## Quasipotential approach to scattering theory: Spectral analysis of the vertex function

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The vertex function of the quasipotential in scattering equations is spectral analyzed in terms of the strength eigenstates, and a criterion for its effectiveness established. A simple variational procedure is given to construct the strength eigenstates, by which vertex functions may be evaluated.

NUCLEAR REACTIONS Quasipotential method, convergence of the iteration series.

In a recent paper,<sup>1</sup> we have derived a quasipotential (QP) approach which can be used to solve a class of scattering equations with divergent iteration kernels. It is a direct extension of the earlier work by Sasakawa<sup>2</sup> and Austern,<sup>3</sup> combined with the quasiparticle approach of Weinberg.<sup>4</sup> The vertex function  $\varphi_t$ , which appears in the QP of Ref. 1, was chosen rather arbitrarily, and its effectiveness was examined numerically for a simple form. A more systematic way of constructing  $\varphi_t$  is desirable and, in this paper, we report both on a spectral analysis of  $\varphi_t$  in terms of the strength eigenfunctions and a minimum criterion.

(1) Following Ref. 1, we consider a simple potential scattering of the form

$$D_0|u\rangle = V|u\rangle \quad , \tag{1}$$

where  $D_0 = E - H_0$ , and  $H_0$  may contain some distortion potentials, especially in the case of multichannel scattering. The corresponding integral equation is given by

$$|u\rangle = |u_0\rangle + G_0 V |u\rangle \quad , \tag{2}$$

where

$$D_0|u_0\rangle = 0, \quad D_0G_0 = 1$$
 (3)

The quasipotential approach of Ref. 1 is obtained by a replacement of V in (1) and (2) by  $V_s$  defined by

$$V_{s}(u_{t},\varphi_{t}) = V|u_{t}\rangle \frac{1}{\langle \varphi_{t}^{\dagger}|V|u_{t}\rangle} \langle \varphi_{t}^{\dagger}|V|.$$
(4)

In particular, an iterative procedure is obtained when  $|u_t\rangle$  is chosen to be the *i*th iterated solution of the equation

$$[D_0 - V_s(u_i, \varphi_i)]|u_{i+1}\rangle = 0 \quad . \tag{5}$$

The vertex function  $\varphi_t$  is still arbitrary, but the convergence property of (5) depends *critically* on the choice of  $\varphi_t$ . A simple choice  $|\varphi_t\rangle = |u_0\rangle$  was seen

to result in the Sasakawa-Austern procedure,<sup>2,3</sup> while the form  $|\varphi_t\rangle = f|u_0\rangle$ , with  $f = e^{-ar}$  and *a* a nonlinear adjustable parameter, gave a strongly convergent result in the case of  $p + {}^{16}O$  scattering.

As is well known,<sup>4,5</sup> the convergence property of an iteration series can be studied conveniently in terms of the strength eigenfunctions  $\{\varphi_n\}$  generated by the iteration kernel  $K_0 \equiv G_0 V$ , as

$$K_0 | \varphi_n \rangle = \lambda_n | \varphi_n \rangle$$
 with  $\langle \varphi_n^{\dagger} | V | \varphi_m \rangle = \delta_{nm}$ , (6)

where  $\varphi_n$  are regular at r = 0 and satisfy a purely outgoing-wave boundary condition beyond some  $r > r_c$ . Thus  $K_0$  may be expanded as

$$K_0 = \sum_{m} |\varphi_n\rangle \lambda_n \langle \varphi_n^{\dagger} | V \quad . \tag{7}$$

The scattering function  $|u\rangle$  is given by

$$|u\rangle = |u_0\rangle + \sum a_n |\varphi_n\rangle \quad , \tag{8}$$

with  $a_n = \lambda_n b_n / (1 - \lambda_n)$  and  $b_n = \langle \varphi_n^{\dagger} | V | u_0 \rangle$ . The usual Born series results when  $(1 - \lambda_n)^{-1}$  is expanded in powers of  $\lambda_n$ ; obviously, the series would diverge if any  $\lambda_n$  in (8) would be outside the unit circle, i.e.,  $|\lambda_n| \ge 1$ .

The iteration series obtained from (5) with (4) may be summarized for the *i*th iteration as follows<sup>1</sup>:

$$|\chi_i\rangle = K_0|u_{i-1}\rangle \quad , \tag{9a}$$

$$y_{i} = \frac{\langle \varphi_{t}^{\dagger} | V | U_{0} \rangle}{\langle \varphi_{t} | V | u_{i-1} \rangle - \langle \varphi_{t}^{\dagger} | V | \chi_{i} \rangle} , \qquad (9b)$$

and thus the wave function is given by

$$|u_i\rangle = |u_0\rangle + y_i|\chi_i\rangle \quad . \tag{9c}$$

If the series converges, we expect that

$$|u_i\rangle \rightarrow |u\rangle, \quad y_i \rightarrow 1$$
, (10)

and, after the *i*th iteration,

$$|u_{i}\rangle = |u_{0}\rangle + \sum_{n=1}^{\infty} \lambda_{n} b_{n} |\varphi_{n}\rangle \left(1 + \lambda_{n} y_{i-1} + \lambda_{n}^{2} y_{i-1} y_{i-2} + \cdots + \lambda_{n}^{i-1} \prod_{j=1}^{i-1} y_{j}\right) y_{i} , \qquad (11)$$

$$\frac{27}{413}$$

which is to be compared with the *i*th iterated Born series obtained when all the *y*'s are set equal to unity. Before discussing the role of the vertex function  $\varphi_i$ , we note the following properties: (a) From (8) and (11), the  $\varphi_n$  content of  $|u\rangle$  is determined essentially by the overlap  $b_n = \langle \varphi_n^{\dagger} | V | u_0 \rangle$ , so that, for those  $\varphi_m$ 's with  $b_m = 0$ , no convergence problem arises; (b) the convergence (or divergence) of the iteration series depends on the factor in (11), inside the square bracket. Although eventually  $y_i \rightarrow 1$  as  $i \rightarrow \infty$ , all the *y*'s conspire in such a way that (11) is fundamentally different from the Born series. The precise way in which this occurs will be illustrated below. [The Born series is obtained by setting all  $y_i = 1$  in Eq. (11).]

$$|u_{1}\rangle = |u_{0}\rangle + \sum_{n \neq 1}^{\infty} \lambda_{n} b_{n} |\varphi_{n}\rangle \frac{1}{1 - \lambda_{1}} + a_{1} |\varphi_{1}\rangle ,$$
  

$$|u_{2}\rangle = |u_{0}\rangle + \sum_{n \neq 1}^{\infty} \lambda_{n} b_{n} |\varphi_{n}\rangle \left[1 + \frac{\lambda_{n}}{1 - \lambda_{1}}\right] + a_{1} |\varphi_{1}\rangle ,$$
  

$$|u_{3}\rangle = |u_{0}\rangle + \sum_{n \neq 1}^{\infty} \lambda_{n} b_{n} |\varphi_{n}\rangle \left[1 + \lambda_{n} + \lambda_{n}^{2} \frac{1}{1 - \lambda_{1}}\right] + a_{1} |\varphi_{1}\rangle, \text{ etc.}$$

The above result reveals many salient features of the QP approach: (i) After the *first* iteration, the  $\varphi_1$  part of  $u_i$  has converged completely, and the subsequent iterations do not modify this part of u; (ii) the  $\varphi_n$  parts (n > 1) of  $u_i$  are obtained exactly as in the Born series, *except* for the (i - 1)th term which is multiplied by  $y_1$ . (This may be improved by allowing  $\varphi_i$  to contain  $\varphi_n$ 's other than  $\varphi_1$ .) (iii)  $y_i \rightarrow 1$  immediately after the first iteration, although the overall  $u_i$  has still not converged. Therefore  $y_i \rightarrow 1$  is *not* a reliable indicator of the convergence.

(3) Next, consider the case in which  $|\lambda_1|, |\lambda_2| > 1$ , while  $|\lambda_n| < 1$  for  $n \ge 3$ , and try for  $\varphi_t$  a form

$$|\varphi_t\rangle = c_1|\varphi_1\rangle + c_2|\varphi_2\rangle \quad , \tag{15}$$

with  $c_1$  and  $c_2$  to be determined below.

 $y_1 = \frac{c_1 b_1 + c_2 b_2}{c_1 b_1 (1 - \lambda_1) + c_2 b a_2 (1 - \lambda_2)}$ 

This case is of special interest, because our QP is still a *single* term of rank 1, while the conventional quasipotential should be of rank 2. In fact, we have shown<sup>6</sup> that, when  $\varphi_i$  is chosen properly, the weaker form of rank 1 is *sufficient* to guarantee the convergence of the  $u_i$  series. The form (15) may also be used, even when  $|\lambda_2| < 1$ , to improve the overall convergence. A simple algebraic manipulation yields for the first iteration

and

$$|u_{1}\rangle = |u_{0}\rangle + \sum_{n \neq 1, 2}^{\infty} \lambda_{n} b_{n} |\varphi_{n}\rangle y_{1} + \{\lambda_{1} b_{1} |\varphi_{1}\rangle y_{1} + \lambda_{2} b_{2} |\varphi_{2}\rangle y_{2}\} .$$
(16)

On the other hand, the result (14) with one  $\varphi_1$  sug-

$$|\varphi_t\rangle = c_1|\varphi_1\rangle \quad . \tag{12}$$

[Of course, the QP form (4) is independent of the overall normalization of  $\varphi_t$ .] Then, it is a simple matter to show that

$$y_1 = \frac{1}{1 - \lambda_1}, \quad y_n = 1, \quad n > 1$$
, (13)

and  $|u_i\rangle$  after the *i*th iteration are explicitly given by

(14)

gests that we would like to have, in (16),

$$\{ \} \rightarrow a_1 | \varphi_1 \rangle + a_2 | \varphi_2 \rangle \quad , \tag{17}$$

where  $a_n = \lambda_n b_n / (1 - \lambda_n)$ . (This choice is, of course, not unique but is good enough for our purpose.) Obviously, this is not possible no matter how  $c_1$  and  $c_2$ in  $y_1$  are adjusted. Therefore, for the present case, we need at least two iterations. We have, for i = 2and from (11),

$$|u_{2}\rangle = |u_{0}\rangle + \sum_{n \neq 1, 2}^{\infty'} \lambda_{n} b_{n} |\varphi_{n}\rangle (1 + \lambda_{n} y_{1}) y_{2}$$
  
+  $\{\lambda_{1} b_{1} |\varphi_{1}\rangle (1 + \lambda_{1} y_{1}) y_{2}$   
+  $\lambda_{2} b_{2} |\varphi_{2}\rangle (1 + \lambda_{2} y_{1}) y_{2}\}$ . (18)

It is now possible to adjust the parameters  $c_1$  and  $c_2$ in  $\varphi_r$  of (15) such that the term { } of (18) satisfies (17). That is,

$$\frac{c_1}{c_2} = -\frac{\lambda_2 b_2}{\lambda_1 b_1} \quad . \tag{19}$$

Thus, aside from the overall constant and with  $b_n = \langle \varphi_n^{\dagger} | V | u_0 \rangle$ ,

$$|\varphi_{t}\rangle = \frac{\lambda_{1}}{b_{1}}|\varphi_{1}\rangle - \frac{\lambda_{2}}{b_{2}}|\varphi_{2}\rangle \quad .$$
<sup>(20)</sup>

The choice (20) gives the following iteration series:

$$y_{1} = \frac{1}{1 - (\lambda_{1} + \lambda_{2})} ,$$
  

$$y_{2} = \frac{1 - (\lambda_{1} + \lambda_{2})}{(1 - \lambda_{1})(1 - \lambda_{2})} ,$$
(21)  

$$y_{n} = 1, \quad n \ge 3 .$$

and

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$$|u_{1}\rangle = |u_{0}\rangle + \sum_{\substack{n \neq 1, 2}}^{\infty} \lambda_{n} b_{n} |\varphi_{n}\rangle y_{1} + (\lambda_{1} b_{1} |\varphi_{1}\rangle y_{1} + \lambda_{2} b_{2} |\varphi_{2}\rangle y_{1}) ,$$

$$|u_{2}\rangle = |u_{0}\rangle + \sum_{\substack{n \neq 1, 2}}^{\infty} \lambda_{n} b_{n} |\varphi_{n}\rangle (1 + \lambda_{n} y_{1}) y_{2} + (a_{1} |\varphi_{1}\rangle + a_{2} |\varphi_{2}\rangle) , \qquad (22)$$

$$|u_{3}\rangle = |u_{0}\rangle + \sum_{\substack{n \neq 1, 2}}^{\infty} \lambda_{n} b_{n} |\varphi_{n}\rangle (1 + \lambda_{n} y_{2} + \lambda_{n}^{2} y_{1} y_{2}) + (a_{1} |\varphi_{1}\rangle + a_{2} |\varphi_{2}\rangle), \text{ etc.} ,$$

so that, after the second iteration,  $|u_i\rangle$  will contain the correct  $\varphi_1$  and  $\varphi_2$  components, as we expected. The other components will be just as in the Born series, except for the last two terms, which are multiplied by  $y_2$  and  $y_1y_2$ , respectively.

The above procedure could be carried further for the  $\varphi_t$  with more than two  $\varphi_n$ 's. This proves that (5) with (4) will converge for all cases if  $\varphi_t$  is chosen properly. Even if some  $|\lambda_n| < 1$  for which  $c_n \neq 0$ , the iteration series will be made to converge faster. On the other hand,  $\varphi_t$  need not contain all the troublesome  $\varphi_n$ 's in precisely the form (20), but should include "enough" of them so that the series converges, in the sense of (20), either by cancellations or by reduction in strength.

In summary, we have shown that the rank 1 form (4) for  $V_s$  and the resulting iteration by (5) will converge with a proper choice of  $\varphi_t$  even when a finite number of  $\varphi$ 's occur with  $|\lambda_n| > 1$ . What is not yet clear is whether a simple and systematic way to generate a useful form of  $\varphi_t$  can be found. We have not been able to derive such a method, although a variational principle for the  $\varphi_n$ 's can be formulated.<sup>7</sup> From a practical point of view, it is highly desirable to avoid, if possible, the use of exact  $\varphi_n$ 's in the con-

struction of  $V_s$ , especially when the scattering system is complex and many channels are open. Since the main requirement on the role of  $V_s$  is to *reduce* the strength of those  $\lambda_n$ 's with  $|\lambda_n| \ge 1$ , such that the remaining part of these states in the iteration series is convergent,<sup>4</sup> often a crude approximation to  $\varphi_n$  may be sufficient. For example, from (6), we have

$$VG_0 V |\varphi_n\rangle = \lambda_n V |\varphi_n\rangle \quad , \tag{23}$$

in which case<sup>1</sup> a simultaneous diagonalization of the matrices  $\langle \varphi_{mt} | VG_0 V | \varphi_{nt} \rangle$  and  $\langle \varphi_{mt} | V | \varphi_{nt} \rangle$  will provide a set  $\{\varphi_{nt}\}$  with  $\{\lambda_{nt}\}$ . Alternatively,<sup>7</sup>

$$D_0 |\varphi_n\rangle = \frac{1}{\lambda_n} V |\varphi_n\rangle \quad , \tag{24}$$

in which case the matrices to be diagonalized are  $\langle \varphi_{mt} | D_0 | \varphi_{nt} \rangle$  and  $\langle \varphi_{mt} | V | \varphi_{nt} \rangle$ . This provides an approximate set  $\{\varphi_{nt}\}$ .

Other iterative procedures in which a rank 1 QP of the form (4) is introduced will be discussed elsewhere.<sup>6</sup>

The author would like to thank G. Rawitscher for several useful discussions.

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