Trajectories of S-matrix poles in a new finite-range potential

A. Rácz,¹ P. Salamon,^{1,2} and T. Vertse^{1,2}

¹University of Debrecen, Faculty of Informatics, PO Box 12, H–4010 Debrecen, Hungary ²Institute of Nuclear Research of the Hungarian Academy of Sciences, Debrecen, PO Box 51, H–4001, Hungary (Received 7 June 2011; published 2 September 2011)

The trajectories of S-matrix poles are calculated in the finite-range phenomenological potential introduced recently by Salamon and Vertse [Phys. Rev. C 77, 037302 (2008)] (SV). The potential is similar to a Woods-Saxon (WS) interaction, but it is exactly zero beyond a radius, without any cutoff. The trajectories of the resonance poles in this SV potential are compared to the corresponding trajectories in a cutoff WS potential for l > 0. The dependence on the cutoff radius is demonstrated. The starting points of the trajectories turn out to be related to the average ranges of the two terms in the SV potential.

DOI: 10.1103/PhysRevC.84.037602

PACS number(s): 21.10.Pc, 21.30.-x, 21.60.Cs

Solutions of the Schrödinger equations that belong to poles of the scattering matrix are often used in reaction calculations and also in the description of weakly bound states. Certain pole solutions can be used as single-particle basis functions in a so-called Berggren representation [1], and they are useful in the theoretical description of slightly bound or unbound states in the framework of the Gamow shell model [2] (GSM). The GSM is a powerful tool for the description of the new nuclei recently produced by the radioactive beam facilities. For the numerical calculation of the *S*-matrix poles in the Woods-Saxon (WS) potential, a new method was recently suggested [3].

The WS potential form is often used in reaction calculations. Although the WS potential is considered a finite-range potential, its value becomes zero only at infinite distances. The radial Schrödinger equation with this potential has analytical solutions [4] only for l = 0. Therefore, we have to calculate its solutions (eigenvalues and eigenfunctions) by using numerical methods for the integration. The typical solution method is to integrate the radial equation numerically in an interval $r \in$ $[0, R_{\text{match}}]$ of the radial distance r where the nuclear potential is large, and match the calculated solution at a distance r = R_{match} as in Ref. [9], or at two different distances as in Ref. [3] to the solution of the asymptotic differential equation, which is the Coulomb or the Ricatti-Hankel differential equation. Thus, in this solution we effectively use a modified version of the WS potential in which the potential is cut to zero at a finite distance R_{max} artificially. The cutoff WS potential is the following:

$$V^{\rm WS}(r) = V_0 f^{\rm WS}(r), \tag{1}$$

where the radial shape is

$$f^{\rm WS}(r) = -\begin{cases} \frac{1}{1+e^{\frac{r-R}{a}}}, & \text{if } r < R_{\rm max}\\ 0, & \text{if } r \ge R_{\rm max}. \end{cases}$$
(2)

The solution and, consequently, the resulting pole position does depend on the value of the cutoff radius R_{max} . Therefore, the cutoff radius R_{max} is a parameter of the WS form in Eq. (2), in addition to its depth V_0 , radius R, and diffuseness a. In publications, however, the value of R_{max} is very seldom specified because most physicists think that the cut has a negligible effect on the calculated quantities if R_{max} is large and thus the jump at the cutoff is quite small.

In a recent work by Salamon and Vertse [6] it was discussed that the positions of the broad resonances might be very sensitive to the R_{max} value, and the authors proposed an alternative to the cutoff WS potential. The advantage of the proposed new potential (SV) is that its range is finite in the strict sense since it is exactly zero at a value ρ_{max} . From mathematical point of view the SV potential is nicer than the cutoff WS potential since its derivative exists everywhere in contrast to the latter. Indeed the shape of the SV potential becomes zero smoothly, and even its derivatives become zero at and beyond $r = \rho_{\text{max}}$.

The objective of this Brief Report is to show the implications of the tail of the WS potential for the positions of the *S*-matrix poles as well as their remedy achieved by the SV potential.

The sensitivity to R_{max} can be seen if we follow the position of the S-matrix poles in the complex wave-number plane (k plane) as the strength of the potential V_0 is changed. These curves are the so-called pole trajectories. The shapes of the pole trajectories do depend on the radial shape of the potential concerned. The pole trajectories in a square well/barrier potential had been studied by Nussenzweig [5] over 50 years ago. In a square well the analytical form of the solution is known inside the potential radius R_{sq} as well, and the pole positions are obtained as roots of a simple transcendental equation.

Certain features of the pole trajectories in the square well remain valid for realistic (diffuse) potentials if their range is finite. Other features, however, do depend on the radial shape. The pole trajectories of the cutoff WS form have not been studied as extensively as those of the square well. In a work by Bang *et al.* [7], the trajectories of the resonant poles were given for a WS potential cut off at a fixed R_{max} value, but the R_{max} dependence of the trajectories was not explored.

We now calculate the pole trajectories in cutoff WS potentials with different cutoff radii, and compare them to the similar trajectories in the SV potentials. We present the calculations for l > 0 only since in the l = 0 case we have no

centrifugal barrier to produce narrow resonances, and therefore this case is of reduced importance.

The radial shape of the potentials is fixed, and only the strength is varied along a pole trajectory. For the sake of comparison, we rewrite the original SV potential in Ref. [6] in the form

$$V^{\rm SV}(r) = V_0 f^{\rm SV}(r), \tag{3}$$

where $f^{SV}(r)$ is taken a linear combination:

$$f^{\rm SV}(r,c_1,\rho_0,\rho_1) = f_{\rho_0}(r) + c_1 f'_{\rho_1}(r), \qquad (4)$$

where

$$f_{\rho}(r) = \begin{cases} -e^{\frac{r^2}{r^2 - \rho^2}}, & \text{if } r < \rho \\ 0, & \text{if } r \ge \rho, \end{cases}$$
(5)

and $f'_{\rho}(r)$ is its derivative:

$$f_{\rho}'(r) = \begin{cases} -\frac{2r\rho^2}{(r^2 - \rho^2)^2} e^{\frac{r^2}{r^2 - \rho^2}}, & \text{if } r < \rho \\ 0, & \text{if } r \ge \rho. \end{cases}$$
(6)

Since both terms in Eq. (4) have finite ranges, the SV potential of Eq. (3) has a finite range of $\rho_{\text{max}} = \max\{\rho_0, \rho_1\}$, i.e., $f^{\text{SV}}(r \ge \rho_{\text{max}}) = 0$. With the usual choices $(\rho_0 > \rho_1)$, the range is $\rho_{\text{max}} = \rho_0$.

In order to make the shape of the SV potential similar to the $f^{WS}(r)$ shape, we adjust the range parameters ρ_0 , ρ_1 and the weight c_1 so as to minimize the integral of the squared differences:

$$\Delta(\rho_0, \rho_1, c_1) = \int_0^{\rho_{\max}} [f^{\text{SV}}(r, c_1, \rho_0, \rho_1) - f^{\text{WS}}(r)]^2 dr.$$
(7)

In general, the R_{max} value is large ($\rho_{\text{max}} < R_{\text{max}}$). The fitting fixes the geometrical parameters of the SV potential (ρ_0 , ρ_1 , c_1) so that $f^{\text{SV}}(r)$ depends only on R and a of the WS shape $f^{\text{WS}}(r)$ and does not depend on R_{max} . The quality of the fit can be characterized by the minimum Δ_{min} of Δ .

We consider potentials whose shapes are suitable for describing ${}^{16}\text{O}$ and ${}^{208}\text{Pb}$. We compare the radial shapes of the best-fit SV potentials and the original WS potentials in Fig. 1. One can see that the shapes of the WS forms are reproduced reasonably well both inside and in the surface regions. There are, however, considerable differences in the tail region, where the SV shape goes to zero faster than the WS shape.

In Table I we summarize the geometrical parameters of the WS and SV potentials used in this work. The parameters of the WS potential for oxygen and for lead were taken from Bang *et al.* [7] and from Curutchet *et al.* [8], respectively. The spin-orbit terms were neglected for the sake of simplicity. The SV potential parameters in Table I are the results of the minimization of the deviation Δ in Eq. (7). The values of Δ_{min} are also given in the table.

We can notice that the shape of the SV potential in Eq. (3) has three parameters (c_1, ρ_0, ρ_1) just as the WS shape does (R, a, R_{max}) , but the SV potential has a natural finite range: ρ_{max} . We can observe in Table I that the average values of the two ranges $\bar{\rho}$ show some similarity to the radius parameter *R* of the WS potential, but $\bar{\rho}$ is larger in both cases.

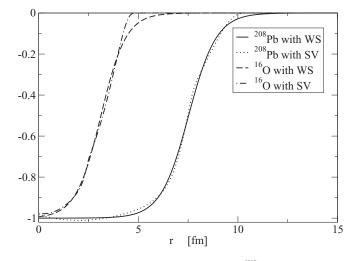


FIG. 1. Comparison of the radial shapes $f^{WS}(r)$ in Eq. (2) and $f^{SV}(r)$ in Eq. (3) for the ¹⁶O and ²⁰⁸Pb nuclei with parameters listed in Table I.

We shall only present trajectories for the decaying resonances since the trajectory of a capturing resonance is simply the mirror image of the trajectory of a decaying resonance.

The pole energies and radial wave functions have been calculated by a modified version of the computer code GAMOW [9]. The node number n starts with n = 1, by convention. This n is the actual radial node number in the limit of large potential depth when the pole becomes a bound-state pole.

For convenience, we start the trajectories with a small V_0 at a small positive strength, $V_0 = 5$ keV. We denote the wave number of the pole belonging to $V_0 = 5$ keV by $k_{n,l}^0$. After finding $k_{n,l}^0$ we increment the V_0 value by small steps repeatedly until the resonance pole reaches the imaginary k axis. For a high l value or for a large node number the imaginary k axis is reached at a very large V_0 value, which is much deeper than the realistic potential well.

Nussenzweig [5] found that for a square well l = 0 the pole trajectories approach the $\operatorname{Re}(k_n) = \frac{n\pi}{R_{sq}}$ lines with the depth approaching to zero. For l = 1 the starting values of the pole trajectories approach the $\operatorname{Re}(k_n) = \frac{(2n-1)\pi}{2R_{sq}}$ values.

Bang *et al.* [7] calculated pole trajectories in a cutoff WS potential for ¹⁶O with a cutoff radius $R_{\text{max}} = 9.4$ fm for l = 0 and l = 1. Figures 5 and 6 of that work suggest that the density of the l = 0 and l = 1 poles are close to the densities of the square well poles belonging to $R_{\text{max}} = R_{\text{sq}}$, where R_{sq} denotes the radius of the square well. But the shapes of the cutoff WS trajectories were quite different from the shapes of the square well trajectories.

TABLE I. Geometrical parameters of the SV and WS potentials and the minimal deviation for nuclei ¹⁶O and ²⁰⁸Pb. All distances are in fm units.

A	c_1	$ ho_0$	$ ho_1$	$\bar{ ho}$	R	а	Δ_{\min}
							5.6×10^{-3} 2.3×10^{-3}

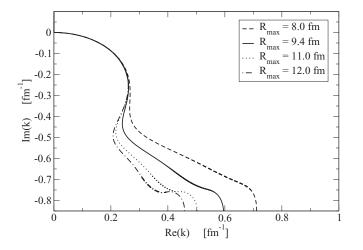


FIG. 2. Trajectories of the n = 2 resonant poles in the complex k plane for l = 1 for WS potentials for ¹⁶O with different values of the cutoff radius R_{max} . The Re $(k) = \frac{3\pi}{2R_{\text{max}}}$ values are 0.59 fm, 0.5 fm, 0.43 fm, and 0.39 fm for increasing R_{max} values, respectively.

In Fig. 2 we present the trajectories of the n = 2, l = 1 resonance pole for ¹⁶O for WS potentials cut off at different distances R_{max} . The curve with $R_{\text{max}} = 9.4$ fm resembles the shape of the n = 2 curve in Fig. 6 of Ref. [7]. It should not be the same since we neglected the spin-orbit term of the potential. We can see that for different R_{max} values the real parts of the starting points $k_{2,1}^0$ are different. The larger the R_{max} value, the smaller $\text{Re}(k_{2,1}^0)$ that is obtained.

We define a sort of pole density at the starting points belonging to $V_0 = 5$ keV as the number of $\operatorname{Re}(k_{n,l}^0)$ values of the same *l*, falling into a $\operatorname{Re}(k)$ interval of unit length. Thus, the smaller the distance between the starting values with node number *n* and *n* + 1, the higher the pole density in a given partial wave *l*. The density of the resonance poles in WS potentials increases with increasing cutoff radii. This is a typical feature of the trajectories calculated in cutoff WS potentials for all *l* and *n* values.



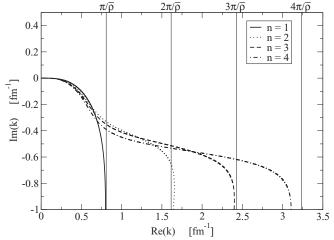


FIG. 4. Trajectories of the resonant poles in the complex k plane for l = 2 for SV potentials for different n values for ¹⁶O.

Now we turn to studying trajectories in the SV potential. In Fig. 3 one can see the trajectories of the l = 1 resonance pole for the SV potential for different n values. The vertical lines in the figure denote the k values for which $\operatorname{Re}(k) = \frac{(2n-1)\pi}{2\bar{o}}$. The starting points of the trajectories with different n values are close to the corresponding vertical lines of the figure. Therefore, we conclude that for the SV potential the average value of the ranges of the two terms plays the same role as the cutoff radius R_{max} of the WS potential. This rule also applies to the l = 2 case, which is shown in Fig. 4. However, in this case *l* is even; therefore the vertical lines correspond to the *k* values with $\operatorname{Re}(k) = \frac{n\pi}{\bar{a}}$. For l = 0 the real parts of the starting $k_{0,n}^0$ values are similar to those of the l = 2 k trajectories [Re $(k_{2,n}^0)$]. The shapes of the trajectories are different, however, because for l = 0, the decaying and capturing resonance poles meet at the imaginary k axis well below the origin.

For ²⁰⁸Pb the pole trajectories in the WS potential depend on the R_{max} value in the same way as observed for ¹⁶O. The

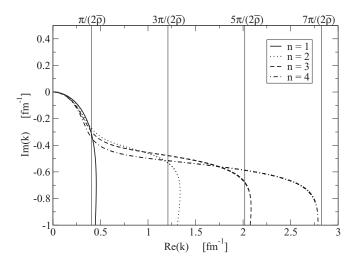


FIG. 3. Trajectories of the resonant poles in the complex k plane for l = 1 for SV potentials for different n values for ¹⁶O.

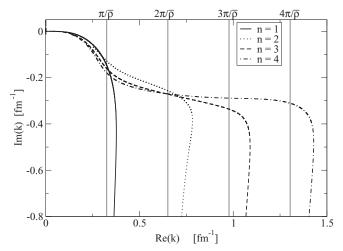


FIG. 5. Trajectories of the resonant poles in the complex k plane in SV potential for l = 2 for different n values for ²⁰⁸Pb.

larger the R_{max} value, the smaller the abscissa of the starting point of the trajectory.

For the SV potential the pole trajectories are very similar to the those for ¹⁶O; therefore we only show the l = 2 case with different *n* values.

In Fig. 5 we show the resonance pole trajectories for l = 2 and for n = 1, 2, 3, 4 values. The curves are similar to the ones in Fig. 4. The starting points of the trajectories $k_{n,2}^0$ are close to the values denoted by vertical lines at $\text{Re}(k) = \frac{n\pi}{\bar{\rho}}$ with $\bar{\rho} = 9.644$ fm.

To sum up, a common feature of the trajectories in the three finite-range potentials (square well, cutoff WS, and SV types) is that the density of the resonant poles is influenced by the range of the potentials. The rule derived by Nussenzweig [5] for square well is that for $V_0 \rightarrow 0$ the distance of the consecutive resonance poles $\operatorname{Re}(k_{n+1,l} - k_{n,l})$ tends to $\frac{\pi}{R_{sq}}$ for any *l* value. Bang *et al.* [7] conjectured that this rule applies for cutoff WS if we replace the R_{sq} by the cutoff radius R_{\max} . They also conjectured that for $V_0 \rightarrow 0$ the l = 0 poles approach the $\operatorname{Re}(k_{n,1}) = \frac{(2n-1)\pi}{2R_{\max}}$ and the l = 1 poles approach the $\operatorname{Re}(k_{n,1}) = \frac{(2n-1)\pi}{2R_{\max}}$. Our results in Fig. 2 support this conjecture qualitatively.

In the SV potential case the shapes of the trajectories are different from that of the WS potential and the starting k value depends on the average of the ranges of the two

terms, rather than the range, i.e., on the value of ρ_{max} . Since $\bar{\rho}$ is close to the radius *R* of the WS form, the distance between the starting Re(*k*) values for consecutive *n* values is governed by the radius of the WS potential to which the parameters of the SV potential were fitted, and not by the cutoff radius R_{max} .

When we calculate the pole energies numerically, we have to take the matching radius R_{match} , or both values of the matching distances used in Ref. [3], larger than the ranges of the potentials (R_{max} and ρ_{max}).

The dependence of the trajectory on the R_{max} value points to the importance of specifying the value of the R_{max} used in the calculations in order to avoid uncertainties in the pole positions. The uncertainty is appreciable for broad resonances.

Contrary to common belief, this strong R_{max} dependence on the cutoff WS potential cannot be diminished by increasing the value of the cutoff radius.

The authors are grateful to R. G. Lovas for valuable discussions. This work was supported by the TÁMOP 4.2.1./B-09/1/KONV-2010-0007/IK/IT project. The project is implemented through the New Hungary Development Plan cofinanced by the European Social Fund and the European Regional Development Fund.

- [1] T. Berggren, Nucl. Phys. A 109, 265 (1968).
- [2] N. Michel, W. Nazarewicz, M. Ploszajczak, and T. Vertse, J. Phys. G: Nucl. Part. Phys. 36, 013101 (2009).
- [3] A. M. Mukhamedzhanov, B. F. Irgaziev, V. Z. Goldberg, Yu. V. Orlov, and I. Qazi, Phys. Rev. C 81, 054314 (2010).
- [4] Gy. Bencze, Commentationes Physico-Mathematicae **31**, 4 (1966).
- [5] H. M. Nussenzveig, Nucl. Phys. 11, 499 (1959).

- [6] P. Salamon and T. Vertse, Phys. Rev. C 77, 037302 (2008).
- [7] J. Bang, S. N. Ershov, F. A. Gareev, and G. S. Kazacha, Nucl. Phys. A **339**, 89 (1980).
- [8] P. Curutchet, T. Vertse, and R. J. Liotta, Phys. Rev. C 39, 1020 (1989).
- [9] T. Vertse, K. F. Pál, and Z. Balogh, Comput. Phys. Commun. 27, 309 (1982).