An Algebraic Approach to Reflectionless Potentials in One Dimension

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Abstract
We develop algebraic methods to find the eigenenergies and eigenstates of reflectionless potentials in one dimension.

1 Introduction
A few interesting problems in wave mechanics have exact solutions in terms of simple functions. The best known examples – the harmonic oscillator and the hydrogen atom – teach us so much about the structure of quantum systems that they are firmly established in the syllabi of elementary courses. Another class of one-dimensional potential problems also have exact solutions in terms of simple functions. The potentials are inverse hyperbolic cosines,

\[ v_\ell(x) = -\ell(\ell + 1 \sech^2 x) \]

where \( \ell \) is any positive integer \( (\ell = 0, 1, 2, \ldots) \). Both the bound states and scattering states can be found analytically for these potentials in terms of elementary functions. In fact this is the only example (other than step potentials and \( \delta \)-functions) I know of where the scattering states can be found by elementary means. These potentials have remarkable properties including bound states at zero energy, and reflectionless scattering. The latter means that a particle incident on the potential is transmitted with unit probability, albeit with an interaction-dependent phase. As a result they are known as “reflectionless potentials”.
The Schrödinger equation for the harmonic oscillator and the Coulomb potential can be either by the more-or-less standard analysis of differential equations, or by algebraic methods. The algebraic solution to the harmonic oscillator using raising and lowering operators can be found in any textbook. The algebraic solution to the hydrogen atom using the commutation relations of the “Runge-Lenz” vector and the angular momentum is treated in some texts [1].

Eq. (1) can also be solved by a direct attack on the differential equation. The approach can be found in Ref. [2]. The solutions are expressed in terms of confluent hypergeometric functions that reduce to elementary functions when the strength of the potential is \( \ell (\ell + 1) \). On the other hand, the eigenstates of reflectionless potentials can be found very easily using operator methods very similar to those that are used to solve the harmonic oscillator in elementary quantum mechanics texts. This does not appear to be very widely known. In this short paper, I will develop the operator solution to reflectionless potentials [3].

2 General Formalism

We begin with the one-dimensional Schrödinger equation,

\[
\left[ -\frac{\hbar^2}{2m} \frac{d^2}{d\xi^2} - V_0 \operatorname{sech}^2 (b \xi) \right] \psi (\xi) = E \psi (\xi) . \tag{2}
\]

For convenience we scale out the dimensionful quantities by defining \( x = b \xi, v_0 = 2mV_0 b^2 / \hbar^2, \) and \( k^2 = 2m v_0 E / \hbar^2, \) so for \( v_0 = \ell (\ell + 1), \)

\[
\mathcal{H}_\ell \psi (x) = \left[ p^2 - \ell (\ell + 1) \operatorname{sech}^2 x \right] \psi (x) = k^2 \psi (x) \tag{3}
\]

where \( p = -i \frac{d}{dx} \). \( k^2 \leq 0 \) corresponds to bound states and \( k^2 > 0 \) corresponds to scattering. Bound states should have normalizable wavefunctions, \( \int dx |\psi (x)|^2 < \infty \), and scattering states should be defined in terms of incoming, transmitted, and reflected waves. I will show that \( \mathcal{H}_\ell \) has \( \ell \) bound states and also exhibit explicit wave functions for the bound and scattering states of \( \mathcal{H}_\ell \).

In an analogy to the harmonic oscillator we introduce operators

\[
a_\ell = p - i \ell \tanh x \\
a_\ell^\dagger = p + i \ell \tanh x . \tag{4}
\]
Using the canonical commutator between $p$ and $x$, $[x, p] = i$, it is easy to show that

$$A_\ell \equiv a_\ell^\dagger a_\ell = p^2 + \ell^2 - \ell (\ell + 1) \text{sech}^2 x$$
$$B_\ell \equiv a_\ell a_\ell^\dagger = p^2 + \ell^2 - \ell (\ell - 1) \text{sech}^2 x.$$  \hspace{1cm} (5)

First we look for the ground state – a normalizable state annihilated by $a_\ell$. We define the state $|0\rangle_\ell$ by the equation

$$a_\ell |0\rangle_\ell = 0$$

or

$$\left(-i \frac{d}{dx} - i \ell \tanh x\right) \psi_{0\ell}(x) = 0$$

where

$$\psi_{0\ell}(x) = \langle x |0\rangle_\ell$$  \hspace{1cm} (6)

which has the normalizable solution

$$\psi_{0\ell}(x) = N_\ell \text{sech}^\ell (x).$$  \hspace{1cm} (7)

Since $\psi_{0\ell}$ is normalizable it is a bound state. Since it has no nodes, a standard theorem on the one-dimensional Schrödinger equation guarantees it is the ground state.

Now consider the relation between the operators $A_\ell$, $B_\ell$, and $H_\ell$. Comparing eqs. (3) and (5),

$$A_\ell = H_\ell + \ell^2$$
$$B_\ell = H_{\ell-1} + \ell^2.$$  \hspace{1cm} (8)

Suppose $\psi$ is an eigenstate of $A_\ell$,

$$A_\ell |\psi\rangle = \alpha |\psi\rangle.$$  \hspace{1cm} (9)

Then it is also an eigenstate of $B_\ell$ with the same eigenvalue, $\alpha$, as shown by the following algebra:

$$a_\ell \{ A_\ell |\psi\rangle \} = \alpha a_\ell |\psi\rangle$$
$$= \{ a_\ell a_\ell^\dagger \} a_\ell |\psi\rangle = B_\ell a_\ell |\psi\rangle.$$  \hspace{1cm} (10)

The only exception to this is the state $|0\rangle_\ell$, because $a_\ell |0\rangle_\ell = 0$. So $|0\rangle_\ell$ is an eigenstate of $A_\ell$ with eigenvalue $\alpha = 0$, which has no corresponding eigenstate of $B_\ell$.

Eq. (8) enables us to turn this into a statement about the Hamiltonians, $H_\ell$: $H_{\ell-1}$ and $H_\ell$ must share the same spectrum except for the single state $|0\rangle_\ell$. These simple results allow us to construct the eigenstates and eigenenergies of all the Hamiltonians.

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\(^1\ell = 0\) is a special, very simple, case that is treated in the next section.
3 Eigenstates and Eigenenergies

The easiest way to see how the information of the preceding section enables us to solve for eigenstates and eigenenergies is to start with $\ell = 0$, then consider $\ell = 1$, and so on until the pattern becomes obvious.

3.1 $\ell = 0$

For $\ell = 0$, $\mathcal{H}_0 = p^2$. This is a free particle. We know the eigenstates, $|k\rangle_0$. They are labeled by the wave number $k$, and the subscript, 0, which refers to $\ell = 0$,

$$\psi_0(k, x) \equiv \langle x | k \rangle_0 = \exp ikx .$$

(11)

The corresponding eigenenergies are $E(k) = k^2$. According to our operator analysis, there should be a state, $|0\rangle_0$, determined by $a_0 |0\rangle_0 = 0$, or $\frac{d}{dx} \psi_0(0, x) = 0$. The solution is simply a constant, $\psi_0(0, x) = \text{const}$.

3.2 $\ell = 1$

For $\ell = 1$ the results become nontrivial. The Hamiltonian is

$$\mathcal{H}_1 = p^2 - 2 \text{sech}^2 x .$$

(12)

According to our work in Section 2, the spectrum of $\mathcal{H}_1$ is identical to that of $\mathcal{H}_0$ except for the state $|0\rangle_1$. Thus we have established that $\mathcal{H}_1$ has a continuum of eigenstates with $E = k^2$.

The $\ell = 1$ ground state is determined by $\mathcal{A}_1 |0\rangle_1 = 0$. Using eq. (8), $\mathcal{H}_1 = \mathcal{A}_1 + 1$, we find the ground-state energy,

$$\mathcal{H}_1 |0\rangle_1 = - |0\rangle_1 .$$

(13)

So $\ell = 1$ has a bound state with $E^{(0)}_1 = -1$. The spectrum of $\mathcal{H}_1$ is now complete; a bound state at $E = -1$ and a continuum $E = k^2$. It is shown in Fig. 1 along with other values of $\ell$.

The wavefunctions of the eigenstates can be constructed using methods quite similar to those used for the harmonic oscillator. The ground state is easy; from eq. (7) we have

$$\psi^{(0)}_1(x) = \langle x | 0 \rangle_1 = N_1 \text{sech} x .$$

(14)
Figure 1: Energy levels in reflectionless potentials.

To construct the continuum eigenstates, consider the state obtained by acting with $a_1^\dagger$ on the continuum eigenstates of $\mathcal{H}_0$,

$$|k\rangle_1 \equiv a_1^\dagger |k\rangle_0 .$$

The action of $\mathcal{H}_1$ on these states can be related to the $\ell = 0$ problem as follows:

$$\mathcal{H}_1 |k\rangle_1 = (A_1 - 1)a_1^\dagger |k\rangle_0$$
$$= a_1^\dagger B_1 |k\rangle_0 - a_1^\dagger |k\rangle_0$$
$$= (k^2 + 1)a_1^\dagger |k\rangle_0 - a_1^\dagger |k\rangle_0$$
$$= k^2 a_1^\dagger |k\rangle_0 = k^2 |k\rangle_1 .$$

Thus $|k\rangle_1$ is an eigenstate of $\mathcal{H}_1$ with eigenvalue $k^2$.

The continuum state wavefunctions are given by

$$\psi_1(k, x) = \langle x | k\rangle_1 = \langle x | a_1^\dagger |k\rangle_0$$
$$= (-i \frac{d}{dx} + i \tanh x) \exp ikx$$
$$= (k + i \tanh x) \exp ikx .$$

To interpret the continuum states we have to relate them to the usual parameter-
ORIZATION OF SCATTERING IN ONE DIMENSION,
\[ \lim_{x \to -\infty} \psi(k, x) = e^{ikx} + R(k)e^{-ikx} \]
\[ \lim_{x \to \infty} \psi(k, x) = T(k)e^{ikx}. \] (18)

When we take the appropriate limits of eq. (17),
\[ \lim_{x \to -\infty} \psi_1(k, x) = (k - i)e^{ikx} \]
\[ \lim_{x \to \infty} \psi_1(k, x) = (k + i)e^{ikx} \] (19)
and compare with eq. (18) we find
\[ R(k) = 0 \]
\[ T(k) = \frac{k + i}{k - i} \] (20)

As promised, the reflection coefficient vanishes, and the transmission coefficient is a pure phase,
\[ T(k) = \exp\left(2i \tan^{-1}(1/k)\right). \] (21)

This completes the construction for \( \ell = 1 \).

### 3.3 \( \ell = 2 \)

Armed with the methods developed for \( \ell = 1 \), we can construct the solution for \( \ell = 2 \) more quickly. The Hamiltonian is
\[ \mathcal{H}_2 = p^2 - 6 \text{sech}^2 x. \] (22)

According to our general result, the spectrum of \( \mathcal{H}_2 \) coincides with that of \( \mathcal{H}_1 \) except for the ground state, \( |0\rangle_2 \). So there must be two bound states. One with energy \( E = -1 \) is obtained by acting with \( a_2^\dagger \) on \( |0\rangle_1 \), with energy \( E = -1 \), and wavefunction
\[ \psi_2^{(1)}(x) = \langle x | a_2^\dagger | 0 \rangle_1 \]
\[ \propto (p + 2i \tanh x) \text{sech} x \]
\[ \propto \tanh x / \cosh x. \] (23)

Note that this wavefunction is antisymmetric in \( x \to -x \) as we would expect for the first excited state in a one-dimensional potential. The ground-state energy is
determined to be $E_2^{(0)} = -4$ by following an argument analogous to eq. (17). Its wavefunction is given by eq. (7),
\[
\psi_2^{(0)}(x) = \langle x | 0 \rangle_2 = N_2 \text{sech}^2 x.
\]

Finally, the continuum state wavefunctions are constructed by following a procedure analogous to the $\ell = 1$ case. In short,
\[
\psi_2(k, x) = \langle x | a_2^\dagger | k \rangle_1 \\
= (p + 2i \tanh x)(k + i \tanh x) \exp ikx \\
= (1 + k^2 + 3ik \tanh x - 3 \tanh^2 x) \exp ikx.
\]

Comparison with the definition of transmission and reflection coefficients gives
\[
R_2(k) = 0 \\
T_2(k) = \frac{(k + i)(k + 2i)}{(k - i)(k - 2i)} \\
= \exp 2i \left( \tan^{-1} \left( \frac{1}{k} \right) + \tan^{-1} \left( \frac{2}{k} \right) \right).
\]

Clearly we have outlined a method that can be extended to arbitrary $\ell$. The explicit expressions for the wavefunctions are not as interesting as the spectrum and the transmission coefficients.

- A sequence of bound states beginning at $E_\ell^{(0)} = -\ell^2$ and continuing with $E_\ell^{(j)} = -(\ell - j)^2$ until $j = \ell$ and $E_\ell^{(\ell)} = 0$.
- The scattering is reflectionless, and the transmission coefficient is given by
\[
T_\ell(k) = \exp \left( 2i \sum_{j=1}^{\ell} \tan^{-1} \left( \frac{j}{\ell} \right) \right).
\]

4 Discussion

Many interesting features of scattering theory are nicely illustrated by the bound states and transmission coefficients of reflectionless potentials. A full discussion would lead us far afield, so we simply quote some of the most important results:

1. The transmission coefficient, $T_\ell(k)$, has a pole at every value of $k$ at which the potential $\ell(\ell + 1) \text{sech}^2 x$ has a bound state. For $\ell = 1$ we see a pole at $k = i$. For $\ell = 2$ we see poles at $k = i$ and $k = 2i$. 
2. In addition to the bound states at \( k = i, 2i, 3i, \ldots \), the potential \( \ell(\ell + 1) \text{sech}^2 x \) has a bound state at zero energy. The solution to the Schrödinger equation at \( k^2 = 0 \) must become asymptotic to a straight line, \( \psi_\ell(0, x) \to A + Bx \) as \( x \to \pm \infty \). When the slope \( (B) \) of the straight line vanishes, the system is said to possess a bound state at zero energy. The name is justified by the fact that making the potential infinitesimally deeper (and the problem no longer exactly solvable) gives a state bound by an infinitesimal binding energy. Bound states at zero energy are very special to reflectionless potentials.

3. If we parameterize \( T_\ell(k) \) in terms of a phase shift, \( T_\ell(k) = \exp(2i \delta_\ell(k)) \), then it is easy to show that the difference between the phase shift at \( k = 0 \) and \( k \to \infty \) counts \( \pi \) times the number of bound states, with the bound state at zero energy counting as \( \frac{1}{2} \). This result, known as Levinson’s theorem, holds for arbitrary potentials in three dimensions as well as one dimension.

In summary, reflectionless potentials form a simple and versatile laboratory for studying the properties of bound states and scattering.

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**References**


[3] I learned these methods in conversation with Jeffrey Goldstone, who claims they are well known.