

Quantum Scattering Project: Using the Lippmann–Schwinger equation for scattering off a non-local potential

Nathan Jansen

1 Introduction

In high energy physics and the study of quantum mechanical phenomena, particle properties and composition, scattering is an important event to study and understand. This is because the scattering amplitude is something that can be directly measure. It was due to the Rutherford gold foil experiment that an early model of the atom was formed, this model was built on the fact that alpha particles would scatter via Coulombic forces. When shot through a piece of gold foil there was many angles that could only be justified by the idea of a nucleus and the particle participating in an elastic scattering event. Clearly scattering is integral to our understanding of the universe.

Rather than dealing with the wave equation itself, instead the reaction amplitude is solved using the Lippmann–Schwinger equation. The Lippmann–Schwinger equation is an integral form of the Schrodinger equation and constructs a reaction amplitude matrix, the R matrix, on which the diagonal represents an observable in momentum space.

$$R(k', k) = V(k', k) + \frac{2}{\pi} \rho \int_0^\infty dp \frac{(p^2 V(k', p) R(p, k))}{(k_o^2 - p^2)/2\mu}$$

This program employs a method of breaking down the Lippmann–Schwinger integral equation into a system of linear equations to be solved using a matrix.

$$R = V_i + \frac{2}{\pi} \sum_{j=1}^N \frac{k_j^2 V_{ij} R_j w_j}{(k_o^2 - k_j^2)/2\mu} - \frac{2}{\pi} V_i o R_o \sum_{m=1}^N \frac{w_m}{(k_o^2 - k_m^2)/2\mu}$$

Where V is velocity, k_o is observable momentum, k_j is quadrature points, w is weight and R is the R matrix. The weights and denominators are turned into a single vector D and the equation takes on the matrix form:

$$R - DV R = [1 - DV] R = V$$
$$[1 - DV] = F$$

Finally you can solve for the R matrix using matrix inversion

$$[R] = [F]^{-1}[V]$$

2 Error Analysis

For the error analysis we used a known analytical solution. The fitting was very good in some regions but poor in other. The region of good fitting is from 0 to π , where we are seeing resonant behavior, specifically at $\pi/2$ where $\sin = 1$. After we pass region around π the computation begins to fit poorly. The region of best fitting occurs where there is resonance of the system. The error of Gaussian quadrature is expected to follow a power law in respect to N . In the region of 16-20 points error scales as $e = x \log N$ where x is the power. In this case we found that the power of the error is $1/10$. It's lowest value at 19 points of $10e-8$. When the number of points reaches around 30 then the error is highly variable, range from -2 to -5 (log scale). It gains This is because this particular reasons fits well so this is likely a case subtractive cancellation occurring in this particular region.

When doing an analysis with a global error however scales differently. It has a very steep slope in the beginning but then levels out around $10e-5$ on a logarithmic scale. Unfortunately it is a bit too computationally heavy to see any signs of round off for the Gaussian method in terms of solving these matrices. It is a bit too much for my laptop to handle to go too much further than about 200 points. However that far didn't vary much from 50 points (slope in this region of the graph was 0.0019). But as the method of integration makes use of subtracting two summations I would suspect a subtractive cancellation error to occur with a high number of points as well.