

eikonal_s.f takes input from a user specified filename with the format below. It calculates the eikonal S-matrix $S(b)$ according to one of a number of complex potential forms that are specified as either: calc, chap, read, gaus (see below). The name of the S-matrix output file is prompted for upon running the code.

Output in S-matrix file is:

```
no_points  b_step
b  Re.S(b)  Im.S(b)  |S(b)|
.....
```

```
%-----
0.  1.  6.  12.
80.0
0.05  601
0.  30.  0.1
1.25
calc
37.4  1.20  0.75
calc
10.0  1.30  0.60
0.0  1.30  0.60

%-----
% Potentials --> eikonal S-matrix S(b): input data
%-----
% projectile : target:  z1, mass1  z2, mass2
% incident energy per nucleon in lab: elab
% integration step and number of steps:  drx,nx
% impact parameter range and step for S(b): bmin,bmax,bstep
% recommend 0.  30.  0.1 for input to other codes
% Coulomb radius parameter: rc
% real potential options: calc, chap, read, gaus
% use chapel hill global set if chap used
% provide volume parameters vr,rr,ar if calc is used
% potential numbers must be in file if read is used
% (input file format is produced by jlmP.f)
% provide volume parameters vr,rr if gaus is used
% imag potential options: calc, chap, read, gaus
% use chapel hill global set if chap used
% provide volume parameters wv,riv,aiv if calc is used
% provide surface parameters ws,ris,ais if calc is used
% potential numbers must be in file if read is used
% (input file format is produced by jlmP.f)
% provide volume parameters wv,riv if gaus is used
%-----
% NB: all radius parameters are multiplied by target mass2**(1/3)
% unless the projectile mass (mass1) is four or greater: i.e.
% a13=mass2**0.33333333333333d0
% if(mass1.gt.4.d0) a13=a13+mass1**0.33333333333333d0
%-----

calc
37.4  1.20  0.75
calc
10.0  1.30  0.60
0.0  1.30  0.60

chap
```