

Exercise sheet 4

by Robin Heinemann (group 4), Paul Rosendahl (group 4)

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1 Neutrons in the gravitational field

the time-independent Schrödinger equation reads

$$\psi''(z) + \frac{2m}{\hbar^2}(E - V(z))\psi(z) = 0$$

For the gravitational field we can use $V(z) = mgz$ for small changes in z . This results in

$$\psi''(z) + \frac{2m}{\hbar^2}(E - mgz)\psi(z) = 0$$

Using dimensionless coordinate x and energy ε

$$x = \left(\frac{2m^2 g}{\hbar^2} \right)^{1/3} z$$
$$\varepsilon = \left(\frac{2}{g^2 \hbar^2 m} \right)^{1/3} E$$

The Schrödinger equation reads as

$$\psi''(x) + (\varepsilon - x)\psi(x) = 0$$

The Numerov algorithm can be easily used to solve this. The infinite potential for $z < 0$ means the solution must be zero at $z = 0$ and the antisymmetric starting condition must be used.

```
1 reset
2 set xrange[0:5]
3 set yrange[-1000:1000]
4 plot for [idx=0:1] "data" i idx u 1:2 w l title columnheader(1)
```

Listing 1: Gnuplot code for plotting the wave function

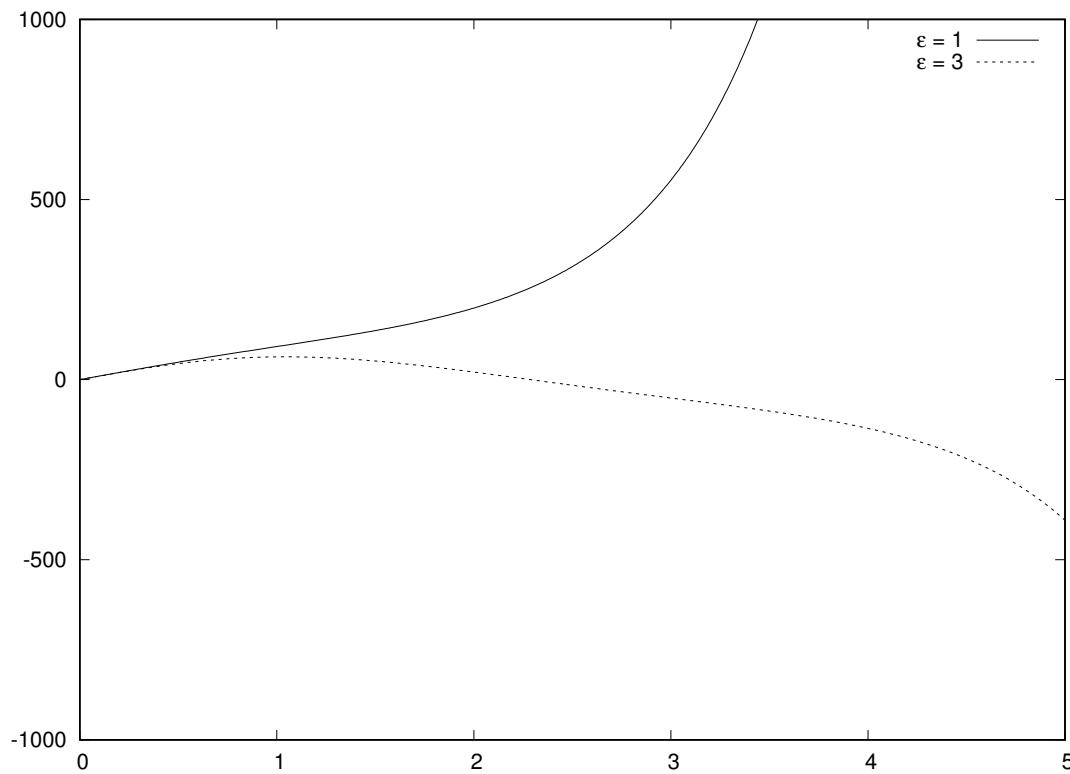


Figure 1: Plot of the wave function for two different ε , one goes towards infinity as x approaches infinity and one goes towards negative infinity

The plot only shows the wave function for small x , to make the structures with small amplitudes visible.

The eigenvalues could be found using a simple linear scan with a stepsize of the accuracy wanted, but is more efficiently implemented with a binary search. This makes it possible to achieve really accurate values for the eigenvalues which just a few calculations of the wavefunction. Table 1 shows the first eight eigenvalues to ten digits accuracy.

2 Rust implementation

```

1 use std::fs::File;
2 use std::path::Path;
3 use std::io::Write;
4
5 fn hermite(n: u64, x: f64) -> f64 {
6     match n {
7         0 => 1.0,
8         1 => 2.0 * x,
9         2 => 4.0 * x * x - 2.0,
10        3 => 8.0 * x * x * x - 12.0 * x,
11        4 => 16.0 * x * x * x * x - 48.0 * x * x + 12.0,
12        _ => 2.0 * x * hermite(n - 1, x) - 2.0 * (n as f64) * hermite(n - 2, x)
13    }
14 }
15

```

```

16 fn numerov(k: &Fn(f64) -> f64, h: f64, n: usize, x0: f64, a: f64, symmetric:
   ↪ bool) -> Vec<f64> {
17     let mut y = Vec::new();
18     let mut i = 2;
19
20     let f = | i | {
21         (1.0 / 12.0) * h * h * k(x0 + (i as f64) * h)
22     };
23
24     if symmetric {
25         y.push(a);
26         let yn = y[0] - h * h * k(0.0) * y[0] / 2.0;
27         y.push(yn);
28     } else {
29         y.push(0.0);
30         y.push(a);
31     }
32
33     let mut kn0 = f(0);
34     let mut kn1 = f(1);
35
36     while i < n {
37         let kn2 = f(i);
38         let yn = 2.0 * (1.0 - 5.0 * kn1) * y[i - 1] - (1.0 + kn0) * y[i - 2];
39
40         y.push(yn / (1.0 + kn2));
41
42         kn0 = kn1;
43         kn1 = kn2;
44         i += 1;
45     }
46
47     y
48 }
49
50 fn harmonic_oscillator(x: f64, eps: f64) -> f64 {
51     2.0 * eps - x * x
52 }
53
54 fn gravitation(x: f64, eps: f64) -> f64 {
55     if x < 0.0 {
56         return 1e308;
57     } else {
58         eps - x
59     }
60 }
61
62 fn factorial(x: u64) -> u64 {
63     let mut f = 1;
64

```

```

65     for n in 1..(x + 1) {
66         f *= n;
67     }
68
69     f
70 }
71
72 fn analytic_harmonic_oscillator(n: u64, x: f64) -> f64 {
73     (- x * x / 2.0).exp() * hermite(n, x) /
74     (2.0_f64.powf(n as f64) * (factorial(n) as f64) *
↪ std::f64::consts::PI.sqrt())
75 }
76
77 fn main() {
78     let path = Path::new("data");
79     let mut file = File::create(&path).unwrap();
80
81
82     for n in 1..3 {
83         let h = 0.01;
84         let ns = (100.0 / h) as u64;
85         let ygrid = numerov(&| x | { gravitation(x, (n as f64) * 2.0 - 1.0) }
86                             , h , ns as usize, 0.0, 1.0, false);
87
88
89         let mut x = 0.0;
90
91
92         file.write_all(format!("{}", {Symbol e}} = {}{}\n", (n as f64) * 2.0 -
↪ 1.0).as_bytes()).unwrap();
93         for y in ygrid {
94             file.write_all(format!("{}", {}{}\n", x, y).as_bytes()).unwrap();
95             x += h;
96         }
97
98         file.write_all("\n\n".as_bytes()).unwrap();
99     }
100
101     let mut e = Vec::new();
102     e.push(find_eigenvalue(0.01, 0.1));
103
104     for i in 0..7 {
105         let last_e = e[i];
106         e.push(find_eigenvalue(last_e, 0.1));
107     }
108
109
110     println!("{}", |$n$| $\backslash$eps_n$|");
111     println!("{}", |-");
112

```

```
113     for i in 0..e.len() {
114         println!("{}", i, e[i]);
115     }
116 }
117
118 fn find_eigenvalue(eps0: f64, delta_eps: f64) -> f64 {
119     let mut eps = eps0;
120     let h = 0.01;
121     let ns = (eps * 10.0 / h) as usize;
122     let ygrid = numerov(&| x | { gravitation(x, eps) }
123         , h , ns as usize, 0.0, 1.0, false);
124
125     let first = ygrid[ns - 1];
126
127     loop {
128         eps += delta_eps;
129         let ns = (eps * 10.0 / h) as usize;
130         let ygrid = numerov(&| x | { gravitation(x, eps) }
131             , h , ns as usize, 0.0, 1.0, false);
132
133         if first * ygrid[ns - 1] < 0.0 { break; }
134     }
135
136
137
138     let mut a = eps - delta_eps;
139     let mut b = eps;
140
141     let mut old_eps = a;
142     let mut old_value = first;
143
144     loop {
145         let c = (a + b) / 2.0;
146
147         let ns = (c * 10.0 / h) as usize;
148         let ygrid = numerov(&| x | { gravitation(x, c) }
149             , h , ns as usize, 0.0, 1.0, false);
150
151         if (old_eps - c).abs() < 1e-10 {
152             break;
153         }
154
155         if ygrid[ns - 1] * old_value > 0.0 {
156             a = c;
157         } else {
158             b = c;
159         }
160
161         old_eps = c;
162     }
```

```

163
164     b
165 }
```

Listing 2: rust implementation of the runge-kutta-4 integrator and application to the three-body problem

Table 1: first eight eigenvalues of the $V(z) = mgz$ potential to ten digits accuracy

n	ε_n
0	2.3381074104830626
1	4.087949443832042
2	5.520559827275572
3	6.786708088442677
4	7.944133584238579
5	9.022650849111365
6	10.040174335651084
7	11.008524295911172

3 Additional notes

All programs written are written using the programming language *rust*. Extra dependencies (*rust crates*) will be listed in a comment in the first line. To get the source files of each program just unzip this *pdf* file. You will find directories for every program in this file. To execute one of the programs run `cargo run` in it's directory. All plots are made with *gnuplot*. This document was written in *org-mode* and converted to *pdf*. The corresponding *org-mode* sources can also be found by unzipping this *pdf* file.