## Exercise sheet 6

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## 1 Perturbed quantum mechanical oscillator

The Hamiltonian is given by

$$h = \frac{H}{\hbar\omega} = \left(\frac{1}{2}\Pi^2 + \frac{1}{2}Q^2 + \lambda Q^4\right)$$
$$(h)_{nm} = (h_0)_{nm} + \lambda \left(Q^4\right)_{nm}$$
$$(h_0)_{nm} = \left(n + \frac{1}{2}\right)\delta_{nm}$$

 $Q_{nm}$  is given by

$$Q_{nm} = rac{1}{\sqrt{2}} \Big( \sqrt{n+1} \delta_{n,m-1} + \sqrt{n} \delta_{n,m+1} \Big)$$

For  $Q^2$  we can find

$$\begin{split} (Q^2)_{ik} &= \frac{1}{2}(\sqrt{i+1}\delta_{i,j-1} + \sqrt{i}\delta_{i,j+1})(\sqrt{j+1}\delta_{j,k-1} + \sqrt{j}\delta_{j,k+1}) \\ &= \frac{1}{2}(\sqrt{i+1}\sqrt{j+1}\delta_{i,j-1}\delta_{j,k-1} + \sqrt{i+1}\sqrt{j}\delta_{i,j-1}\delta_{j,k+1} + \sqrt{i}\sqrt{j+1}\delta_{i,j+1}\delta_{j,k-1} + \sqrt{i}\sqrt{j}\delta_{i,j+1}\delta_{j,k+1}) \\ &= \frac{1}{2}(\sqrt{(i+1)(i+2)}\delta_{i,k-2} + (2i+1)\delta_{ik} + \sqrt{i(i-1)}\delta_{i,k+2}) \\ (Q^4)_{ik} &= \frac{1}{4}(\sqrt{(i+1)(i+2)}\delta_{i,j-2} + (2i+1)\delta_{ij} + \sqrt{i(i-1)}\delta_{i,j+2}) \\ &= \frac{1}{4}(\sqrt{(j+1)(j+2)}\delta_{j,k-2} + (2j+1)\delta_{jk} + \sqrt{j(j-1)}\delta_{j,k+2}) \\ &= \frac{1}{4}(\sqrt{(j+1)(j+2)}\sqrt{(i+1)(i+2)}\delta_{i,j-2}\delta_{j,k-2} + (2j+1)\sqrt{(i+1)(i+2)}\delta_{i,j-2}\delta_{jk} \\ &+ \sqrt{j(j-1)}\sqrt{(i+1)(i+2)}\delta_{i,j-2}\delta_{j,k+2} + \sqrt{(j+1)(j+2)}(2i+1)\delta_{ij}\delta_{j,k-2} \\ &+ \sqrt{(j+1)(j+2)}\sqrt{i(i-1)}\delta_{i,j+1}\delta_{j,k-2} + (2j+1)\sqrt{i(i-1)}\delta_{i,j+2}\delta_{jk} \\ &+ \sqrt{j(j-1)}\sqrt{i(i-1)}\delta_{i,j+2}\delta_{j,k+2}) \\ &= \frac{1}{4}(\sqrt{(i+1)(i+2)(i+3)(i+4)}\delta_{i,k-4} + (2i+5)\sqrt{(i+1)(i+2)}\delta_{i,k-2} \\ &+ (2i+1)^2\delta_{i,k} + \sqrt{(i+1)(i+2)}(2i+1)\delta_{i,k+2} \\ &+ (2i+1)^2\delta_{i,k} + \sqrt{i(i-1)(2i+1)}\delta_{i,k+2} \\ &+ i(i-1)\delta_{i,j} + (2i+5)\sqrt{i(i-1)}\delta_{i,k+2} \\ &+ \sqrt{i(i-1)(i-2)(i-3)}\delta_{i,k+4}) \end{split}$$

If the matrix is of size n there are boundary terms for the diagonals, because of the finite size, which results in

$$= \frac{1}{4} (\sqrt{(i+1)(i+2)(i+3)(i+4)} \delta_{i,k-4} + (2i+5) \sqrt{(i+1)(i+2)} \delta_{i,k-2}$$

$$+ (i+1)(i+2) \delta_{i,k} + \sqrt{(i+1)(i+2)} (2i+1) \delta_{i,k-2}$$

$$+ (2i+1)^2 \delta_{i,k} + \sqrt{i(i-1)} (2i+1) \delta_{i,k+2}$$

$$+ i(i-1) \delta_{i,j} + (2i+5) \sqrt{i(i-1)} \delta_{i,k+2}$$

$$+ \sqrt{i(i-1)(i-2)(i-3)} \delta_{i,k+4}$$

$$(i+1)(i+2) (\delta_{i,0} \delta_{k,0} + \delta_{i,1}, \delta_{k,1}) - i(i-1) (\delta_{i,n-1} \delta_{i,n-1} + \delta_{i,n} \delta_{i,n}))$$

```
#include <tred2.c>
#include <tqli.c>
#include <nrutil.c>
#include <stdio.h>
#include <math.h>
```

6

```
double pythag(double a, double b) {
       return sqrt(a * a + b * b);
   }
   void matmul(double ** a, double ** b, double ** c, int n) {
11
       for(int i = 1; i <= n; i++) {</pre>
12
       for(int j = 1; j <= n; j++) {</pre>
13
            c[i][j] = 0.0;
14
            for(int k = 1; k \le n; k++) {
15
            c[i][j] += a[i][k] * b[k][j];
16
       }
18
       }
19
20
21
   void p(double ** m, int n) {
       for(int i = 1; i <= n; i++) {</pre>
23
       for(int j = 1; j <= n; j++) {
24
            printf("%2.5lf ", m[i][j]);
25
       }
26
       printf("\n");
27
       }
28
   }
29
30
   void perturbed_oscillator(int n, double 1) {
31
       double ** m = dmatrix(1, n + 4, 1, n + 4);
32
       double ** m2 = dmatrix(1, n + 4, 1, n + 4);
33
       double ** m3 = dmatrix(1, n + 4, 1, n + 4);
35
       double sqrt2 = sqrt(2.0);
36
37
       for(int i = 1; i < n; i++) {</pre>
38
       m[i + 1][i] = sqrt(i) / sqrt2;
       m[i][i + 1] = sqrt(i) / sqrt2;
42
       matmul(m, m, m2, n);
43
       matmul(m2, m2, m, n);
44
       for(int i = 1; i <= n; i++) {
       m[i][i] += (i - 1 + .5) / 1;
47
48
       for(int j = 1; j \le n; j++) {
            m[i][j] *= 1;
50
            m2[i][j] = m[i][j];
       }
53
       }
54
55
       double * d = malloc(sizeof(double) * (n + 1));
```

```
double * e = malloc(sizeof(double) * (n + 1));
57
        d = d - 1;
58
        e = e - 1;
59
        tred2(m, n, d, e);
        tqli(d, e, n, m);
62
        matmul(m2, m, m3, n);
64
65
        double * eps = malloc(sizeof(double) * n);
        int * idx = malloc(sizeof(int) * n);
        for(int i = 1; i <= n; i++) {
70
        idx[i-1] = i;
71
        }
        for (int i = 1; i < n; i++) {</pre>
74
        int key = idx[i];
75
        int j = i - 1;
76
        while(j \ge 0 \&\& d[idx[j]] > d[key]) {
            idx[j + 1] = idx[j];
            j = j - 1;
80
        }
81
        idx[j + 1] = key;
82
        printf("|-\n");
        printf("| n = %d n", n);
86
        for(int i = 1; i <= 10; i++) {
        int ix = idx[i - 1];
88
        double eps = 0.0;
        for(int j = 1; j <= n; j++) {
91
            eps += fabs(m[j][ix] - m3[j][ix] / d[ix]);
92
        }
93
        printf("| %.15lf | %.15lf \n", d[ix], eps);
95
   }
97
98
    int main() {
99
        printf("| $\\lambda_i$ | $\\eps$ $\n");
100
        for(int i = 15; i <= 30; i += 5) {
101
        perturbed_oscillator(i, .1);
102
        }
103
   }
104
```

Listing 1: rust implementation of gauss elimination for a tridiagonal matrix

Table 1: the lowest ten eigenvalues  $\lambda_i$  and the 1-norm of  $hv - \lambda v$  where v is the eigenvector with eigenvalue  $\lambda$ 

$aes \lambda_i$ and the 1-norm of	
$\lambda_i$	ε\$
n = 15	
0.559146331717481	0.000000000000056
1.769503154752390	0.0000000000000001
3.138630351927867	0.0000000000000002
4.628832924839890	0.00000000000000002
6.219589268571163	0.0000000000000004
7.900514741235316	0.0000000000000003
9.684402683933815	0.000000000000003
11.394960405389993	0.000000000000003
13.144176458068515	0.0000000000000002
16.606417302164019	0.0000000000000004
n = 20	
0.559146327396061	0.000000000000017
1.769502650598558	0.0000000000000023
3.138624310963206	0.0000000000000003
4.628881549200992	0.0000000000000009
6.220291663918584	0.0000000000000004
7.899835172646707	0.0000000000000007
9.658345956233044	0.00000000000000002
11.484717894488245	0.0000000000000006
13.364075642272910	0.0000000000000005
15.422401778971913	0.0000000000000003
n = 25	
0.559146327187369	0.0000000000000233
1.769502644004068	0.0000000000000003
3.138624307573727	0.0000000000000028
4.628882782968839	0.0000000000000006
6.220300826989985	0.0000000000000010
7.899769051837731	0.0000000000000004
9.657851696340218	0.00000000000000009
11.487225950553698	0.0000000000000006
13.381940814696728	0.0000000000000008
15.340927001633565	0.00000000000000006
n = 30	0.0000000000000000000000000000000000000
0.559146327183577	0.000000000000052
1.769502643949268	0.0000000000000000000000000000000000000
3.138624308466358	0.0000000000000000000000000000000000000
4.628882808395193	0.0000000000000000000000000000000000000
6.220300899870113	0.000000000000038
7.899767283848700	0.0000000000000000000000000000000000000
9.657840218111822	0.000000000000023
11.487312780939364	0.0000000000000000000000000000000000000
13.382455649955588	0.000000000000024
15.338741261872123	0.00000000000000004
13.330/412010/2123	0.0000000000000011

Comparing the eigenvalues for different n it is clear that the lowest one is by far the most accurate. For n=20

 $\lambda_0$  is accurate to about nine digits. Also  $\varepsilon$  is really small for all of the values, which shows that these values are actually eigenvalues.