

Pseudopotential Units (Schmidt et al.)

$$\hat{H}_S = -\frac{1}{2\mu_1} \Delta r_1 - \frac{1}{2\mu_2} \Delta r_2 - \frac{1}{m_3} \nabla_{r_1} \cdot \nabla_{r_2} + \underbrace{q_1 q_3 V_k(r)}_{q_1 q_3} + q_2 q_3 V_k(r_2) + q_1 q_2 V_k(|r_1 - r_2|) \quad (1)$$

2D:

Schmidt-Keldysh potential:

$$V_k(r) = \frac{\pi}{2r_0} \left(H_0 \left(\frac{r}{r_0} \right) - Y_0 \left(\frac{r}{r_0} \right) \right)$$

- ignores $q_1 q_3$
- and sets $K_e = 1$

this is Gaussian Units

Schmidt au = $10^{-22} \text{ eV} / (\text{m/V})^2$
type

$$V_{k_e} = \begin{cases} V_0 & r < r^* \\ -\frac{\alpha}{2} \left(\frac{dV_k(r)}{dr} \right)^2 & r > r^* \end{cases}$$

Where

$$\alpha = 52 \times 10^3 \times 10^{-22} \left(\frac{\text{eV}}{\text{m}^2} \right)^2$$

au

Schmidt:

$$[\frac{J \sqrt{J}}{m^2}] = [\frac{J \cdot J^2}{C^2 m^2}] = [\frac{J}{C^2 m^2}]$$

This doesn't make sense...

X Schmidt ref. [75]

$$\text{MoS}_2: \alpha = 4.6 \times 10^{-18} \text{ eV m}^2 \text{ V}^{-2}$$

fold. Importantly, the polarizability increases by an order of magnitude with increased screening. Hence, at $\kappa = 10$ the polarizability reaches values in the range of $\alpha = 45 \times 10^{-18} \text{ eV}(\text{m/V})^2$ to $70 \times 10^{-18} \text{ eV}(\text{m/V})^2$. Focusing on freely suspended ($\kappa = 1$) MoS_2 and WSe_2 , the polarizabilities are $\alpha = 4.6 \times 10^{-18} \text{ eV}(\text{m/V})^2$ and $\alpha = 6.3 \times 10^{-18} \text{ eV}(\text{m/V})^2$.

The au. of polarisability is

NIST units $\alpha = 52 \times 10^3 \times 1.64877727436 \times 10^{-41} \text{ C}^2 \text{ m}^2 \text{ J}^{-1}$

```
>>> 52e3*1.64877727436e-41
8.573641826672e-37 C^2 m^2 J^-1
```

Units

When $r > r^*$:

$$\therefore -\frac{\alpha}{2} \left(\frac{dV_k(r)}{dr} \right)^2 = \left[\frac{C^2}{J} \right] \left[\frac{J^2}{m^2 C^2} \right] = \left[\frac{J}{C^2} \right]$$

This makes sense

$$q_1 q_3 V_k(r)$$

$$\text{ACTUAL } V_k(r) = \frac{\pi k_e}{(e_1 + e_2) r_0} \left(H_0 \left(\frac{r}{r_0} \right) - Y_0 \left(\frac{r}{r_0} \right) \right), \quad k_e = \frac{1}{4\pi e_0} \text{ N m}^2 \text{ C}^{-2}$$

The units of $V_k(r)$:

$$\left[\frac{N \text{ m}^2 \text{ C}^{-2}}{m} \right] \rightarrow \frac{q_1 q_2}{C^2} \frac{m}{C^2} = [J]$$

$$V_k \text{ actual: } \left[\frac{Nm}{C^2} \right]$$

Note:

$$1 \text{ J} = 6.242 \times 10^{18} \text{ eV}$$

$$[\text{VJ}] = \left[\frac{J}{C} \right] \left[\frac{J}{A \cdot S} \right]$$



Fundamental Physical Constants

atomic unit of electric polarizability $e^2 a_0^2 / E_h$

Numerical value $1.64877727436 \times 10^{-41} \text{ C}^2 \text{ m}^2 \text{ J}^{-1}$

Standard uncertainty $0.0000000050 \times 10^{-41} \text{ C}^2 \text{ m}^2 \text{ J}^{-1}$

Relative standard uncertainty 3.0×10^{-10}

Concise form $1.64877727436(50) \times 10^{-41} \text{ C}^2 \text{ m}^2 \text{ J}^{-1}$

Finding Bound States -

- Direct diagonalisation - (computationally expensive)

(2D S-wave
Radial Schrödinger equation)

$$\frac{1}{r} \frac{d}{dr} \left(r \frac{dR}{dr} \right) + \left(k^2 - \frac{2\mu V(r)}{\hbar^2} \right) R = 0$$

k is wavenumber: $\Psi_{in} = e^{ikr}$

$$u_\ell(r) \equiv r R(k, r)$$

for $r \rightarrow 0$ the asymptotic solution
 $u_\ell(r \gg 0) \propto r^{\ell m}$

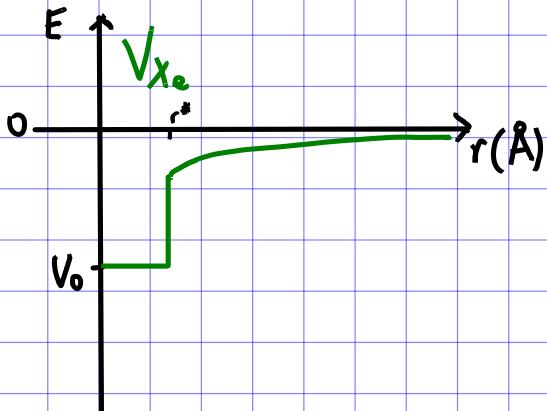
$$\frac{1}{r} \frac{d}{dr} \left(r \frac{dR}{dr} \right) + \left(k^2 - \frac{2\mu V(r)}{\hbar^2} \right) R = 0$$

$$\therefore \frac{1}{r} \frac{d}{dr} \left(r \frac{dR}{dr} \right) - \frac{2\mu V(r) R}{\hbar^2} = -k^2 R$$

$$\therefore \frac{1}{r} \left(\frac{d}{dr} \frac{dR}{dr} + r \frac{d^2 R}{dr^2} \right) - \frac{2\mu V(r) R}{\hbar^2} = -k^2 R$$

$$\therefore \frac{d^2 R}{dr^2} + \frac{1}{r} \frac{dR}{dr} - \frac{2\mu V(r) R}{\hbar^2} = -k^2 R$$

(AKA: Bessel's eqn)



Note: $E = \frac{\hbar^2 k^2}{2\mu}$

$$\therefore \frac{2\mu E}{\hbar^2} = k^2$$

Numerics ($A \vec{R} = \lambda \vec{R}$)

We can write our ODE as a linear system:

$$\left[\begin{bmatrix} \frac{d^2}{dr^2} \\ \frac{d}{dr} \end{bmatrix} + \begin{bmatrix} \frac{1}{r} \\ 0 \end{bmatrix} \right] \left[\begin{bmatrix} \frac{d}{dr} \\ V(r) \end{bmatrix} \right] \begin{bmatrix} 1 \\ R \\ 1 \end{bmatrix} = -k^2 \begin{bmatrix} 1 \\ R \\ 1 \end{bmatrix}$$

where $\begin{bmatrix} \frac{d^2}{dr^2} \end{bmatrix} = \begin{bmatrix} ? & & & \\ & 1 & -2 & 1 \\ & & 1 & -2 & 1 \\ & & & 1 & -2 & 1 \\ & & & & 1 & -2 & 1 \\ & & & & & 1 & -2 & 1 \end{bmatrix}$?? BCS

$\begin{bmatrix} \frac{d}{dr} \end{bmatrix} \begin{bmatrix} \frac{d}{dr} \end{bmatrix} = \begin{bmatrix} ? & & & \\ & 1 & & & \\ & & 1 & & \\ & & & 1 & & \\ & & & & 1 & & \\ & & & & & 1 & & \\ & & & & & & \ddots & \\ & & & & & & & ? ? \end{bmatrix}$

(Diagonal) $\begin{bmatrix} V(r) \end{bmatrix} = \begin{bmatrix} V_{x_e}(r_0) \\ V_{x_e}(r_0+\Delta r) \\ \vdots \\ V_{x_e}(r_{j-1}) \\ V_{x_e}(r_j) \\ \vdots \\ V_{x_e}(r_{N-1}) \end{bmatrix} = \begin{bmatrix} V_0 \\ -V_0 \\ \vdots \\ -V_0 \\ -\frac{1}{2}(\frac{V_0}{\hbar^2} E) \end{bmatrix}$

using centered differences (2nd order)

$$\frac{dR}{dr} \approx \frac{R_{j+1} - 2R_j + R_{j-1}}{\Delta r^2}$$

$$\frac{d^2R}{dr^2} \approx \frac{R_{j+1} - 2R_j + R_{j-1}}{\Delta r^2}$$

if $\vec{A}\vec{R} = \lambda\vec{R}$:

$$\vec{A} = \begin{bmatrix} ?? & ?? \\ 1 & -2 & 1 \\ ?? & ?? \end{bmatrix} + \begin{bmatrix} \frac{1}{r} & & \\ & \frac{1}{r} & \\ & & \frac{1}{r} \end{bmatrix} - \frac{2\mu}{\hbar^2} \begin{bmatrix} V_0 & & \\ & -V_{xe} & \\ & & -\frac{\alpha}{2} \frac{d}{dr} V_{xe} \end{bmatrix}$$

$$\vec{A} = \frac{d^2}{dr^2} + \left(\frac{1}{r} \odot \frac{d}{dr} \right) - \frac{2\mu}{\hbar^2} V_{xe}$$

FINDING Boundary conditions $(l=0)$, $-\frac{2\mu E}{\hbar^2} = k^2$

$$\text{ODE: } \frac{d^2 R}{dr^2} + \frac{1}{r} \frac{dR}{dr} - \frac{2\mu V_{xe}(r)}{\hbar^2} R = -k^2 R$$

$$R'' + \frac{1}{r} R' = \left(-k^2 + \frac{2\mu V_{xe}}{\hbar^2} \right) R$$

As $r \rightarrow 0$, $\frac{1}{r} R'$ dominates

$$\therefore \frac{1}{r} R' = 0 \Rightarrow R' = 0 \quad (\text{Neumann BC, zero derivative BC})$$

$R(r) = C$, C is constant

$$\begin{bmatrix} 0 & ? & 0 \end{bmatrix} \quad S$$

$$\frac{d^2 R}{dr^2} \approx \frac{-2R_j + 2R_{j+1}}{\Delta r^2} \quad \text{Even Symmetry}$$

Because this is ghost point

Matrix coefficients at the boundary

$$\frac{dR}{dr} \approx \frac{R_{j+1} - R_{j-1}}{2\Delta r} = \text{zero}$$

NOT WORKING

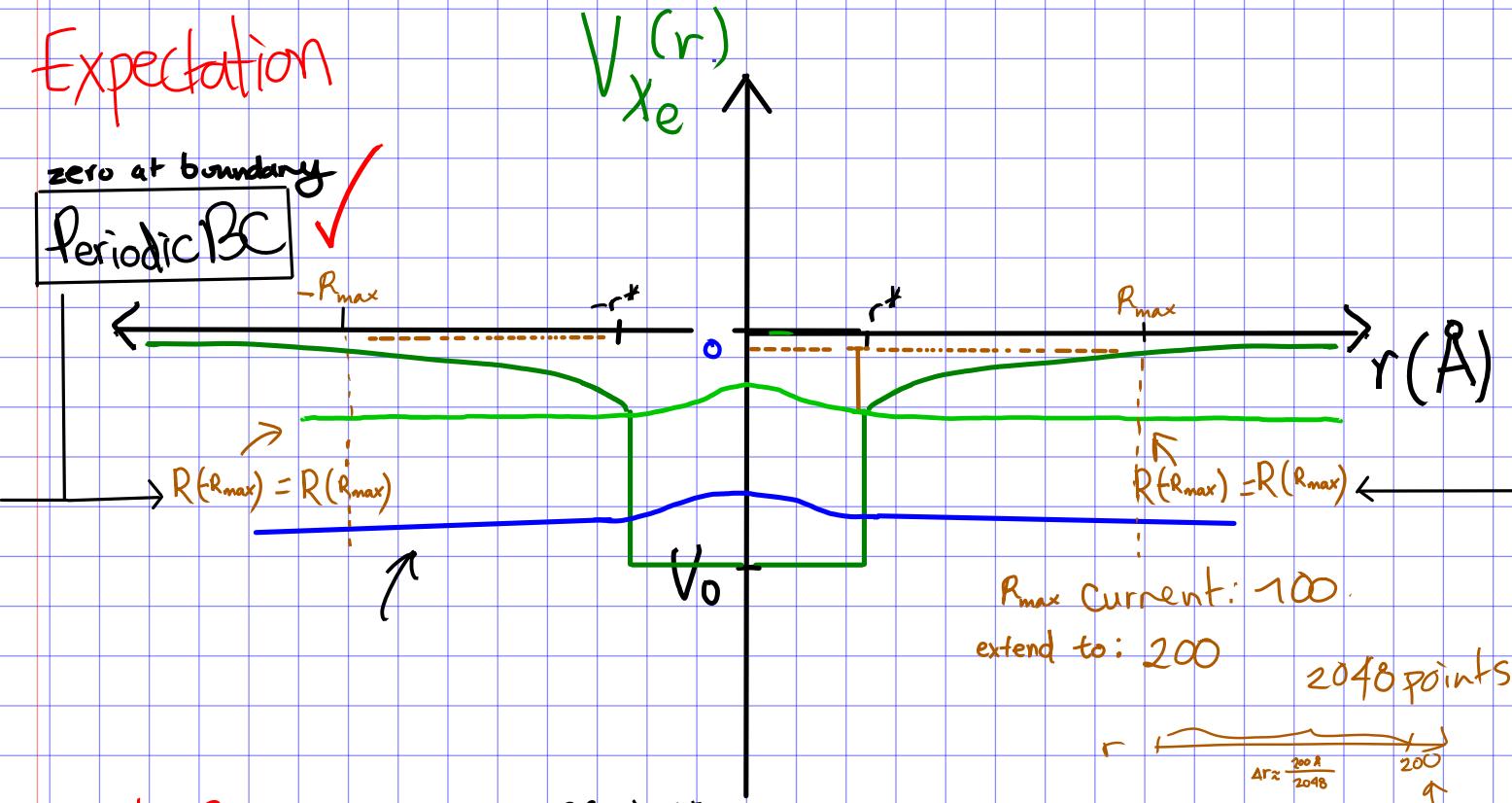
TRY: EVEN extension
TRY: \approx periodic BCs.

Even extension to Schmidt $V_{Xe}(r)$

Expectation

zero at boundary ✓

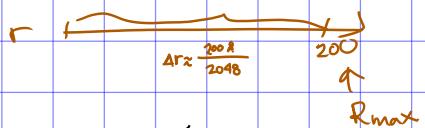
Periodic BC



R_{max} current: 100.

extend to: 200

2048 points



FINDING Boundary conditions ($\ell=0$), $-\frac{2\pi E}{\hbar^2} = k^2$

ODE: $\frac{d^2}{dr^2} + \frac{1}{r} \frac{df}{dr} - \frac{2\mu V_{Xe}(r)}{\hbar^2} R = -k^2 R$ ↗ CHANGE of variables $u(r) = \sqrt{r} R(r)$

① extending r- array to larger R_{max} ✓

② change $r = \sqrt{x^2 + y^2}$ ✓
for $|x|$

modify in

Matrix:

$$A = \frac{1}{\Delta r} D2 + V_{Xe}$$

where $\Delta r = r[1] - r[0]$

[eigenvectors, unitary] = np.linalg.eigh(A)

↑
of A

$$U = \begin{bmatrix} u_{11} & u_{12} & \dots & u_{1N} \\ u_{21} & u_{22} & \dots & u_{2N} \\ u_{31} & u_{32} & \dots & u_{3N} \\ \vdots & \vdots & \ddots & \vdots \\ u_{N1} & u_{N2} & \dots & u_{NN} \end{bmatrix}$$

columns
of U
are eigenvectors
of A.

eventually decided against:

✗ periodic BCS

✗ extension

$$R = \frac{u(r)}{r}$$

Change of variables

20) radial Schrödinger equation ($l=0$)

$$E\psi = -\frac{\hbar^2}{2\mu} \left(\frac{1}{r} \frac{d}{dr} r \frac{d}{dr} \right) \psi + V(r) \psi$$

$$\text{using } \psi = u(r) = \sqrt{r} R(r) \leftarrow$$

$$E = \frac{k^2 \hbar^2}{2\mu} \quad \therefore -k^2 = \frac{2\mu E}{\hbar^2}$$

$$\begin{aligned} \therefore -\frac{k^2 u(r)}{\sqrt{r}} &= \left(\frac{1}{r} \frac{d}{dr} r \frac{d}{dr} \right) \frac{u}{\sqrt{r}} - \frac{2\mu V(r) u}{\hbar^2} \\ &= \frac{1}{r} \frac{d}{dr} r \left(\frac{du}{dr} r^{-1/2} - \frac{u}{2} r^{-3/2} \right) - \frac{2\mu V(r) u}{\hbar^2} \\ &= \frac{1}{r} \frac{d}{dr} \left(\frac{du}{dr} r^{1/2} - \frac{u}{2} r^{-1/2} \right) - \frac{2\mu V(r) u}{\hbar^2} \\ &= \frac{1}{r} \left(\frac{d^2 u}{dr^2} r^{1/2} + \frac{1}{2} \frac{du}{dr} r^{-1/2} - \frac{1}{2} \left(\frac{du}{dr} r^{-1/2} - \frac{u}{2} r^{-3/2} \right) \right) - \frac{2\mu V(r) u}{\hbar^2} \\ &= \frac{d^2 u}{dr^2} r^{-1/2} + \frac{1}{2} \frac{du}{dr} r^{-3/2} - \frac{1}{2} \left(\frac{du}{dr} r^{-3/2} - \frac{u}{2} r^{-5/2} \right) - \frac{2\mu V(r) u}{\hbar^2} \\ &= \frac{1}{Nr} \frac{d^2 u}{dr^2} + \frac{1}{2} \frac{du}{dr} \frac{1}{Nr} - \frac{1}{2} \frac{du}{dr} \frac{1}{Nr} + \frac{u}{4} r^{-5/2} - \frac{2\mu V(r) u}{\hbar^2} \\ \therefore -\frac{k^2 u(r)}{\sqrt{r}} &= \frac{1}{Nr} \frac{d^2 u}{dr^2} + \frac{u}{4} r^{-5/2} - \frac{2\mu V(r) u}{\hbar^2} \end{aligned}$$

$$\therefore -k^2 u(r) = \left(\frac{d^2 u(r)}{dr^2} + \frac{u(r)}{4r^2} - \frac{2\mu V(r) u}{\hbar^2} \right)$$

↑
inverse order approximation.

Schmidt matrix

Tests

- No BCS ✓
- High order approximation to derivative ✓
- zero BCS : ($r \rightarrow 0$, $r \rightarrow R_{\max}$) ✓

FINITE (Mixed) Differences ($\sim 4^{\text{th}}$ order accuracy)

Finite Difference Equations

$$\frac{\partial^{(2)} f}{\partial x^{(2)}} \approx \frac{35f(x+0h) - 104f(x+1h) + 114f(x+2h) - 56f(x+3h) + 11f(x+4h)}{12h^2}$$

$$\frac{\partial^{(2)} f}{\partial x^{(2)}} \approx \frac{11f(x-1h) - 20f(x+0h) + 6f(x+1h) + 4f(x+2h) - 1f(x+3h)}{12h^2}$$

using 4th order centered finite differences

$$\frac{d^2 u(r)}{dr^2} \approx \frac{-1/12 u_{j-2} + 4/3 u_{j-1} - 5/2 u_j + 4/3 u_{j+1} - 1/12 u_{j+2}}{\Delta r^2}$$

$$\frac{\partial^{(2)} f}{\partial x^{(2)}} \approx \frac{-1f(x-3h) + 4f(x-2h) + 6f(x-1h) - 20f(x+0h) + 11f(x+1h)}{12h^2}$$

$$\frac{\partial^{(2)} f}{\partial x^{(2)}} \approx \frac{11f(x-4h) - 56f(x-3h) + 114f(x-2h) - 104f(x-1h) + 35f(x+0h)}{12h^2}$$

	0	1	2	3	4	\dots	$N-1$
0	$\frac{35}{12}$	$-\frac{104}{12}$	$\frac{114}{12}$	$-\frac{56}{12}$	$\frac{11}{12}$	0	Custom
1	$\frac{11}{12}$	$-\frac{20}{12}$	$\frac{6}{12}$	$\frac{1}{12}$	$-\frac{1}{12}$	0	Custom
2	$-\frac{1}{12}$	$\frac{4}{3}$	$-\frac{5}{2}$	$\frac{4}{3}$	$-\frac{1}{12}$	0	\vdots
3	0	$-\frac{1}{12}$	$\frac{4}{3}$	$-\frac{5}{2}$	$\frac{4}{3}$	$-\frac{1}{12}$	\dots
\vdots	\vdots	\vdots	\vdots	\vdots	\vdots	\vdots	\vdots
$N-3$	\dots	$-\frac{1}{12}$	$\frac{4}{3}$	$-\frac{5}{2}$	$\frac{4}{3}$	$-\frac{1}{12}$	0
$N-2$	\dots	\dots	$-\frac{1}{12}$	$\frac{4}{3}$	$-\frac{5}{2}$	$\frac{4}{3}$	$-\frac{1}{12}$
$N-1$	0	\dots	0	$-\frac{1}{12}$	$\frac{4}{3}$	$-\frac{5}{2}$	$\frac{4}{3}$
							Custom
							Custom

Then $A = \frac{1}{\Delta r^2}$ second-D + Schmidt-matrix

A appears to be non-Hermitian due to second-D

This is likely why I get complex phases

Python3-i

```
ana@ana-XPS-13-9343:~/scatterbrain$ python3 -i scatterbrain.py
/usr/lib/python3/dist-packages/numpy/core/_asarray.py:85: ComplexWarning: casting to complex from real
  g complex values to real discards the imaginary part
  return array(a, dtype, copy=False, order=order)
eigenenergy 0: +2.01-0.00j meV
eigenenergy 1: -2.65+3.34j meV
eigenenergy 2: -2.65-3.34j meV
eigenenergy 3: +14.77+11.90j meV
eigenenergy 4: +14.77-11.90j meV
eigenenergy 5: +60.05+20.82j meV
eigenenergy 6: +60.05-20.82j meV
eigenenergy 7: -24.37+0.00j meV
>>> E = -eigenvalues * hbar**2 / (2 * M_red) / meV
>>> for item in np.sort_complex(E)[:100]:
    Schmidt_m = print(item)
... 1
```

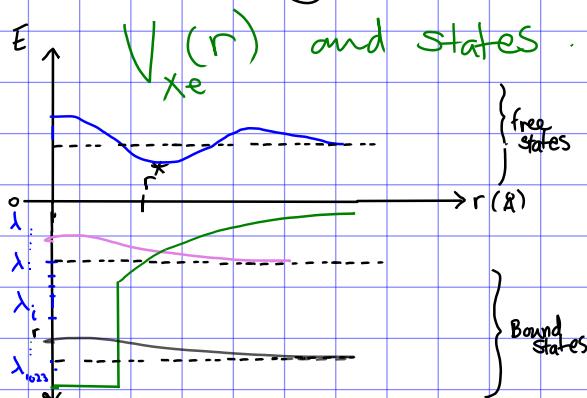
→ lots of complex energy eigenstates as $N \rightarrow$ larger.



Deduced against non-centered differences on the Boundaries!

Result

Direct diagonalisation method



Binding energy stuff

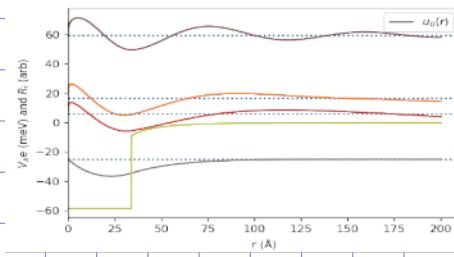
$$\text{CONVERSION: } \frac{1 \text{ J}}{6.242 \times 10^{-21}} = 1 \text{ meV}$$

ENERGY from eigenvalues

$$\lambda_i = k_i^2 = \frac{2 M_{\text{red}} E_i}{\hbar^2}$$

$$\text{GROUND STATE: } -\frac{k_o^2 \hbar^2}{2 M_{\text{red}}} = E_o \quad (\because E_o \text{ in meV})$$

expected $E_o \sim 31.7 \text{ meV}$ (schmidt et al. Fig. 6.)



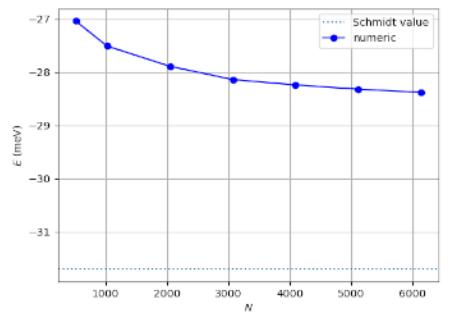
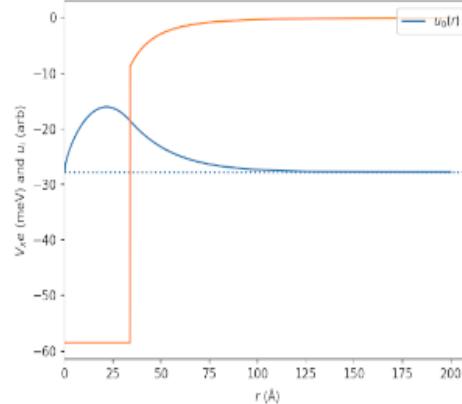
$$N = 6144 \\ r = np.linspace(1e-6, 200, N)$$

Imposing zero BCS $\begin{cases} r \rightarrow 0 : u(r) = 0 \\ r \rightarrow R_{\max} : u(r) = 0 \end{cases}$

A Matrix is Hermitian

$$N = 3072$$

$$E_o = 28.14 \text{ meV}$$



as we increase N value converges $\sim -29 \text{ meV}$
 * Computer cannot handle more points

The screenshot shows a MATLAB IDE interface. On the left, a code editor displays C++ code for generating a sparse matrix. The code uses loops to set elements of a matrix `second` based on conditions involving indices `i` and `j`. It includes comments for 'SPLITTING MATRIX' and 'ELIMINATING MATRIX'. On the right, there are two windows: a 'Matrix Visualization' window showing a 4x4 grid with colored cells (dark purple, yellow, green) representing non-zero values, and a 'Command Window' showing the output of the code execution.

```
function [second] = spalloc(D, i, j)
    % SPLITTING MATRIX
    if (i < j)
        second[i, j] = -1 / 12;
        second[i, i] = 1 / 12;
        second[i, i+1] = 1 / 12;
        second[i+1, i] = 1 / 12;
        second[i+1, i+1] = 1 / 12;
        second[i+1, i+2] = 1 / 12;
        second[i+2, i+1] = 1 / 12;
        second[i+2, i+2] = 1 / 12;
        second[i+2, i+3] = 1 / 12;
        second[i+3, i+2] = 1 / 12;
        second[i+3, i+3] = 1 / 12;
        second[i+3, i+4] = 1 / 12;
    else
        second[i, j] = -1 / 12;
        second[i, i] = 1 / 12;
        second[i, i-1] = 1 / 12;
        second[i-1, i] = 1 / 12;
        second[i-1, i-1] = 1 / 12;
        second[i-1, i-2] = 1 / 12;
        second[i-2, i-1] = 1 / 12;
        second[i-2, i-2] = 1 / 12;
        second[i-2, i-3] = 1 / 12;
        second[i-3, i-2] = 1 / 12;
        second[i-3, i-3] = 1 / 12;
        second[i-3, i-4] = 1 / 12;
    end
    % ELIMINATING MATRIX
    if (i > j)
        second[i, j] = -1 / 3;
        second[i, i] = 5 / 3;
        second[i, i-1] = 1 / 3;
        second[i-1, i] = 1 / 3;
        second[i-1, i-1] = 1 / 3;
        second[i-1, i-2] = 1 / 3;
        second[i-2, i-1] = 1 / 3;
        second[i-2, i-2] = 1 / 3;
        second[i-2, i-3] = 1 / 3;
        second[i-3, i-2] = 1 / 3;
        second[i-3, i-3] = 1 / 3;
        second[i-3, i-4] = 1 / 3;
    else
        second[i, j] = -1 / 3;
        second[i, i] = 5 / 3;
        second[i, i+1] = 1 / 3;
        second[i+1, i] = 1 / 3;
        second[i+1, i+1] = 1 / 3;
        second[i+1, i+2] = 1 / 3;
        second[i+2, i+1] = 1 / 3;
        second[i+2, i+2] = 1 / 3;
        second[i+2, i+3] = 1 / 3;
        second[i+3, i+2] = 1 / 3;
        second[i+3, i+3] = 1 / 3;
        second[i+3, i+4] = 1 / 3;
    end
end

second_matrix = spalloc(4, 4);
second_matrix[1, 1] = 1 / 12;
second_matrix[1, 2] = -1 / 12;
second_matrix[1, 3] = 1 / 12;
second_matrix[1, 4] = 1 / 12;
second_matrix[2, 1] = -1 / 12;
second_matrix[2, 2] = 1 / 12;
second_matrix[2, 3] = 1 / 12;
second_matrix[2, 4] = 1 / 12;
second_matrix[3, 1] = 1 / 12;
second_matrix[3, 2] = 1 / 12;
second_matrix[3, 3] = 1 / 12;
second_matrix[3, 4] = -1 / 12;
second_matrix[4, 1] = 1 / 12;
second_matrix[4, 2] = 1 / 12;
second_matrix[4, 3] = -1 / 12;
second_matrix[4, 4] = 1 / 12;
```

BUG: else statement applied to
The non-defined cases: $j=2, j=N-2$

fixed:

only plot states that satisfy

$$E = -(\text{eigenvalues} * \hbar^2) / 2 M_{\text{red}}$$

iterate over eigenvalues
if not $v_{-0} \leq E[i] \leq 0$:
 continue

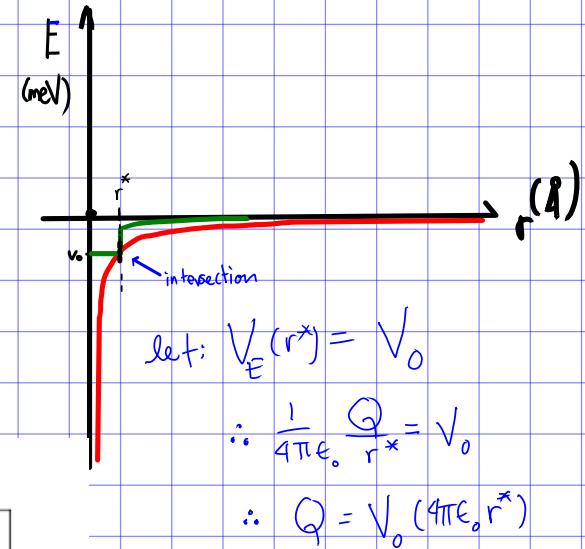
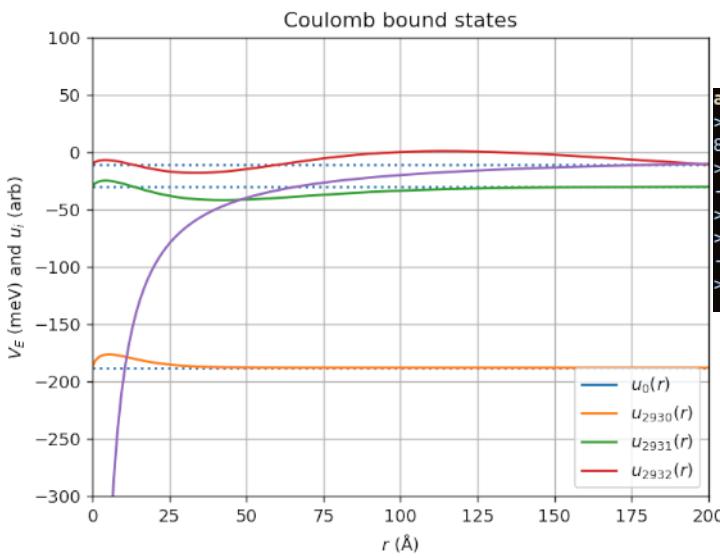
plot: {
| Cote
| plt.show()

Testing Diagonalisation

Coulomb potential

$$V_E(r) = \frac{1}{4\pi\epsilon_0} \cdot \frac{Q}{r}$$

Result



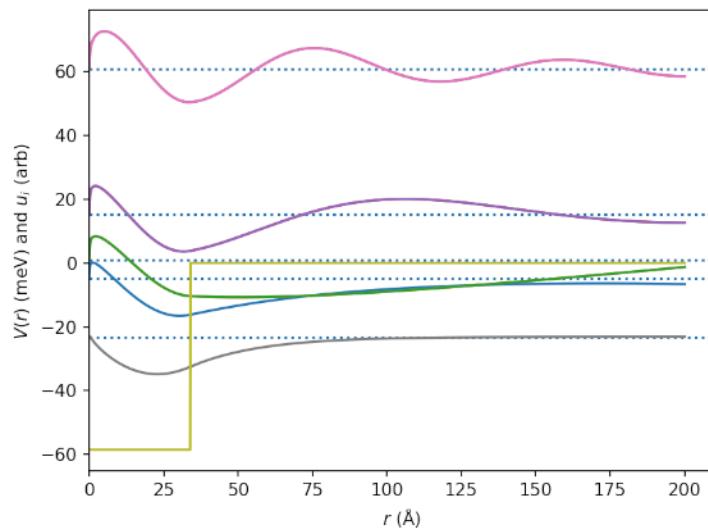
```
ana@ana-XPS-13-9343:~/Desktop/proj$ python3 -i TESTING_D2.py
>>> epsilon
8.854e-12
>>> v_0
-9.37273308899999e-21
>>> Q = v_0 * 4 * np.pi * epsilon * r_star
>>> Q
-3.5456393493840546e-39
>>>
```

Eigenenergies do not satisfy:

$$E_n \propto \frac{1}{(2n-1)^2}$$

where n is eigenstate index.

Diagonalisation using square well and free BCS



Scipy.integrate.solve_ivp(...)

$$\frac{dy}{dt} = f(t, y)$$

- t is 1-D independent variable (scalar)
- y is N-D vector valued $\leftarrow \text{ndarray}$ $\begin{cases} (n,) \\ (\text{shape options}) \\ (n,k) \end{cases}$ (vectorized=True)
- $f(t, y)$ - determines the DE
- t-span is 2-tuple of floats $[t_0, t_f]$ (solver starts @ $t_0 \rightarrow t_f$)
- y_0 is IC. array-like, shape (n,)
- t_eval \hookrightarrow times at which to store the computed Solution
Must be sorted and lie within t-span

return np.shape(arry_like(y))=(n, K)

i.e. each column in this corresponds to a single column in y

return

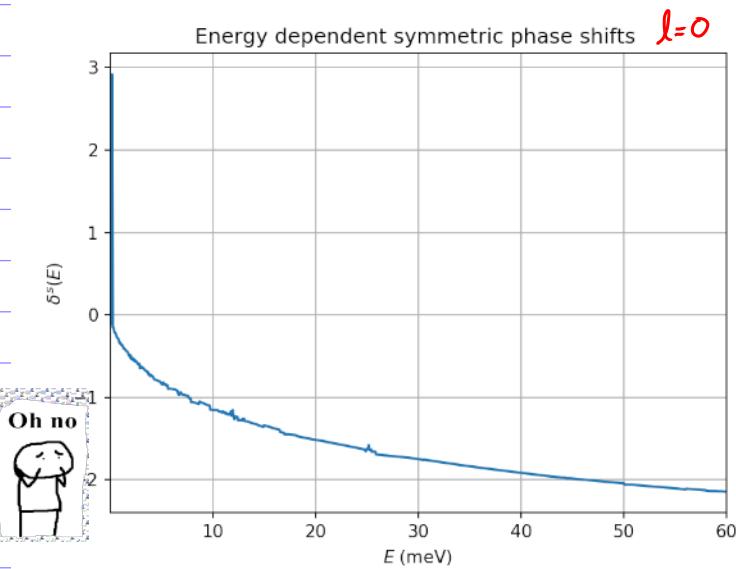
t : ndarray, shape (n_points,)

y : ndarray, shape (n_points,)

sol: Found Solution at t

Defaults

- method = 'RK45'
- t_eval = None
- dense_output = False
- vectorised = False
- options ...



```
#####
## ODE solver SOLVE IVP
delta_0 = [r[0], 1]
delta = []
for i, k_i in enumerate(k):
    # print(i)
    solution_ODE_31 = solve_ivp(
        lambda r, delta : ODE_31(r, delta, k_i),
        [r[0], r[-1]], delta_0, t_eval=r
    )
    delta.append(solution_ODE_31.y[0, -1] - np.pi)

# print(f'THIS IS SOLVE IVP {delta=}\n')
print(f'{np.shape(delta)=}\n')

plt.plot(E / meV, delta)
plt.axis(xmin=E_min / meV, xmax=E_max / meV)
plt.xlabel(r'$E$ (meV)')
plt.ylabel(r'$\delta^s(E)$')
plt.title('Energy dependent symmetric phase shifts ')
plt.grid(True)
plt.savefig('solveivp_symmetric_phase_shifts.png')
plt.show()
#####
```

—Shooting Method—

The essence of the shooting method is to guess a complete \vec{z} at one endpoint, use the relationship for $\frac{d\vec{z}}{dx}$ to propagate a solution to $\vec{z}(x)$ over to the other endpoint, and then compare how close the propagated solution is to known solution in the second boundary condition. You then update your guess and repeat the process to converge the propagated solution to the true solution at the other boundary.

ODE

$$\frac{du^2}{dr^2} = -\frac{u}{4r^2} + \frac{2u}{\hbar^2} V_{xe}(r) u - k^2 u$$

Boundary conditions
(for $r=0$)

$$\vec{y}_0 = \begin{bmatrix} \sqrt{r[0]} \\ \frac{1}{2\sqrt{r[0]}} \end{bmatrix}$$

asymptotic
solution
as $r \rightarrow 0$

1st derivative

We must recast our ODE as a couple of first order ODEs

$$\text{let } \omega = \frac{du}{dr}$$

$$\text{then } \frac{d\omega}{dr} = \frac{d^2u}{dr^2} = -\frac{u}{4r^2} + \frac{2u}{\hbar^2} V_{xe}(r) u - k^2 u$$

Then $\vec{y} = \begin{bmatrix} \omega \\ \frac{d\omega}{dr} \end{bmatrix} = \begin{bmatrix} \frac{du}{dr} \\ -\frac{u}{4r^2} + \frac{2u}{\hbar^2} V_{xe}(r) u - k^2 u \end{bmatrix} = f(r, \vec{y})$

single vector ODE:

Shooting Method

ODE : $\frac{d^2u}{dr^2} + \frac{u}{4r^2} - \frac{2\mu}{\hbar^2} V(r) u + k^2 u = 0$

$$\therefore \frac{d^2u}{dr^2} + \frac{u}{4r^2} - \frac{2\mu V(r)}{\hbar^2} u + k^2 u = 0$$

$$\frac{d^2u}{dr^2} = \left(-\frac{1}{4r^2} + \frac{2\mu V(r)}{\hbar^2} - k^2 \right) u$$

pseudo-potential

$$-\frac{\hbar^2}{2\mu} \frac{d^2u}{dr^2} + V(r) u = E u$$

$$\frac{d^2u}{dr^2} - \frac{2\mu}{\hbar^2} V(r) u = -\frac{2\mu E}{\hbar^2} u$$

$$\frac{d^2u}{dr^2} = \left(\frac{2\mu}{\hbar^2} V(r) - \frac{2\mu E}{\hbar^2} \right) u$$

BCS solve analytically

when: $r \rightarrow 0$, $V_{x_0} = V_0$

$$\frac{d^2u(r)}{dr^2} + \frac{u(r)}{4r^2} = 0$$

asymptotic solution

as $r \rightarrow R_{\max}$
we wait
for shooting method

$$u(r) = C_1 \sqrt{r} + C_2 \sqrt{r} / \ln(r)$$

↓ Throw away

use as asymptotic sol

$$y_0 = \begin{bmatrix} \sqrt{r[0]} \\ \frac{1}{2} r[0]^{-\frac{1}{2}} \end{bmatrix}$$

Initialise: k argument (Guess ①)

1. If $V_0 = -58.5 \text{ meV}$

$$\text{Then } E_{\min} = V_0 = \frac{\hbar^2 k^2}{2 M_{\text{red}}}$$

$$\therefore k^2 = \frac{2 M_{\text{red}} E_{\min}}{\hbar^2}$$

$$E_{\max} = 0$$

if $E > 0$:

$$E_{\text{free}} = \frac{\hbar^2 k^2}{2\mu}$$

$k \in \mathbb{R}$

if $E < 0$:

$$\text{let } k = ik$$

$$E_{\text{bound}} = -\frac{\hbar^2 k^2}{2\mu}$$

CODE + Algorithm

```

# Bound states search:
# SHOOTING method
# ENERGY
# E = -30 * meV #[J]

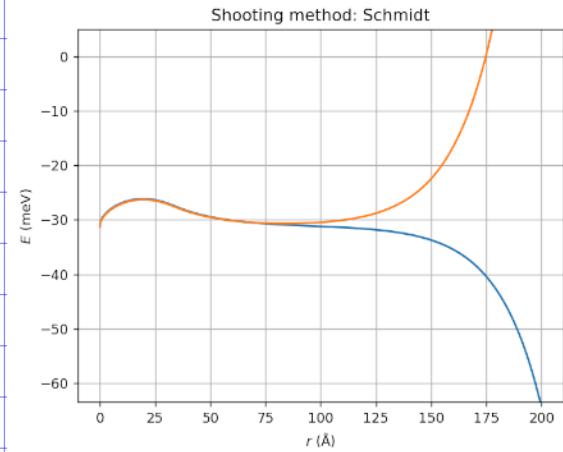
# ODE
def ODE(r, y, E):
    u, w = y
    result = np.zeros_like(y)
    if abs(u) > 1e100: # if blowing up return zero derivative
        return result
    else:
        result[0] = w
        result[1] = (- 1 / (4 * r**2) + 2 * M_red * (V_Xe(r) - E) / hbar**2) * u
        return result

# ICS
y_0 = [np.sqrt(r[0]), 1 / (2 * np.sqrt(r[0]))]
#ODEINT solver
for E in np.linspace(-31.15 * meV, -31.3 * meV, 2):
    solution_ODE = solve_ivp(lambda r, y: ODE(r, y, E), [r[0], r[-1]], y_0, t_eval=r)
    u = solution_ODE.y[0]
    plt.plot(r / angstrom, (u / abs(u[r < r_star]).max() * 5) + E / meV)

plt.axis(ymin=v_0 / meV - 5, ymax=5)
plt.ylabel(r'$E$ (meV)')
plt.xlabel(r'$(\text{\AA})$')
plt.title('Shooting method: Schmidt')
plt.grid(True)
plt.show()

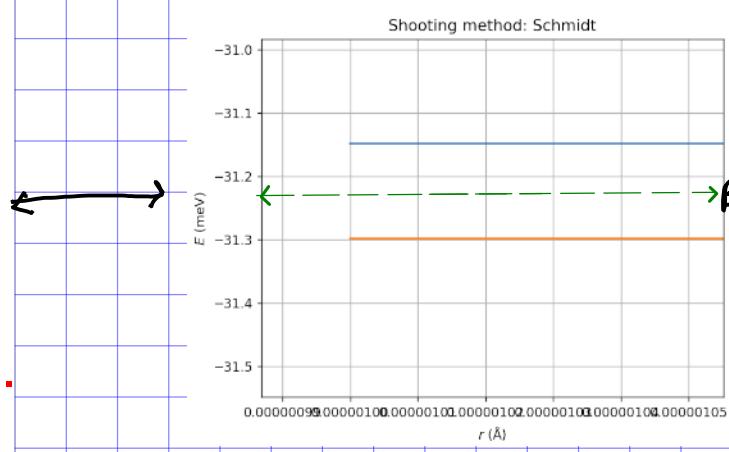
```

Plot:



① Initial guess
 $E_{\text{max}} = -31.15 \text{ meV}$ UB
 $E_{\text{min}} = -31.3 \text{ meV}$ LB

ZOOM!



② Bisection
creates new LB/UB

③ Repeat

BISECTION
new_bound = γ

$$\beta - \alpha = \gamma$$

① set γ as new LB/UB

Automating

```

if sign == -1:
    append to list
if sign == 1:
    append to list

```

-(EL15) Shooting Method -

if: $E = \text{bound state energy}$

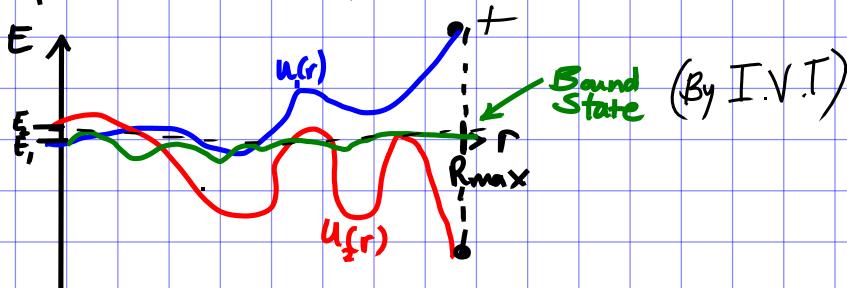
then: $u(R_{\max}) = 0$

\therefore if $E \neq \text{bound state energy}$

then: $u(R_{\max}) \neq 0$

say you have E_1, E_2

for E_1 : $u(R_{\max}) > 0$ and E_2 : $u(R_{\max}) < 0$



$$E \quad Energies = [E_1, E_2, E_3, \dots, E_{1024}] \\ Signs = []$$

$$E_2 - E_1 = \Delta E < E_2$$

1. Divide the potential depth into 1024 energy steps

(1)

Solve ODE
... store solution $\text{sign}(u(R_{\max}))$

(2) search signs for sign flips

$[+, +, +, -, +, +, -, - \dots]$ use these to extract

$[E_1, E_2, E_3, E_4, E_5, E_6, \dots]$ \leftarrow corresponding E_i

(3) construct a list $E_{\text{flip}} = [[E_3, E_4], [E_5, E_6], \dots]$
each of the pairs corresponds to $[E_{\text{upper}}, E_{\text{lower}}]$

The energy bounds for an eigenenergy

we do this until we have scanned the E_{flip} list of Bounds entirely

ITERATIVELY -

④ using E_{low} and E_{up} :

Solve ODE using $E_{\text{low}} + E_{\text{up}} = E_{\text{mid}}$
obtain $\text{sign}(u(R_{\max}))$

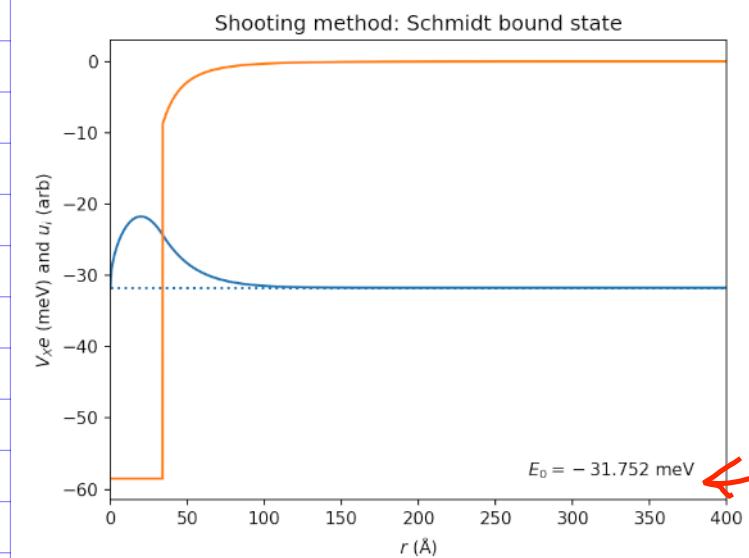
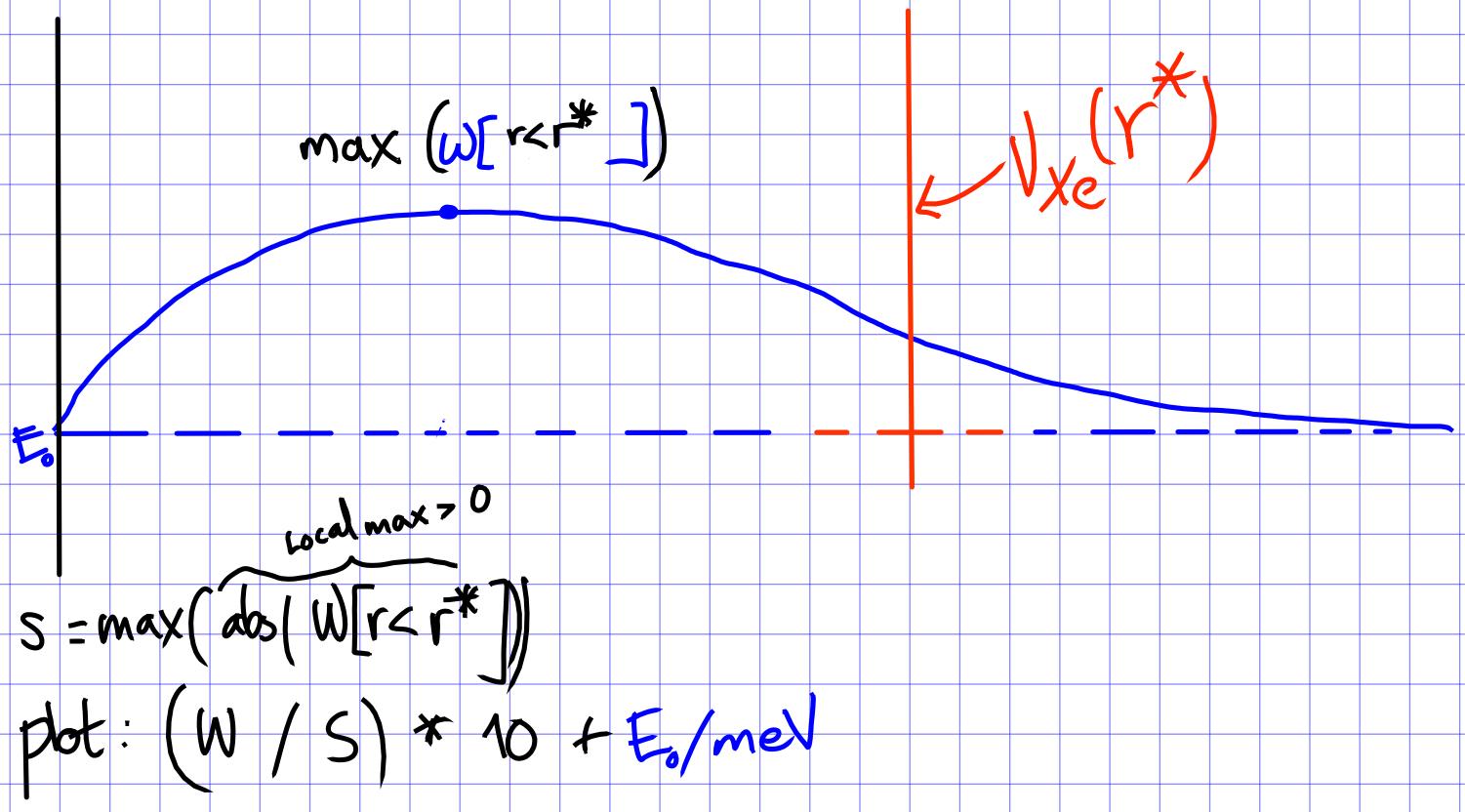
if $\text{sign}(u(R_{\max}))$ for $E_{\text{mid}} = \text{sign}(u(R_{\max}))$ for E_{low}

$$E_{\text{low}} = E_{\text{mid}}$$

else:

$$E_{\text{up}} = E_{\text{mid}}$$

while $\text{abs}(E_{\text{low}} - E_{\text{up}}) > 1e-6$:



E_0 agrees with
 Schmidt et al.