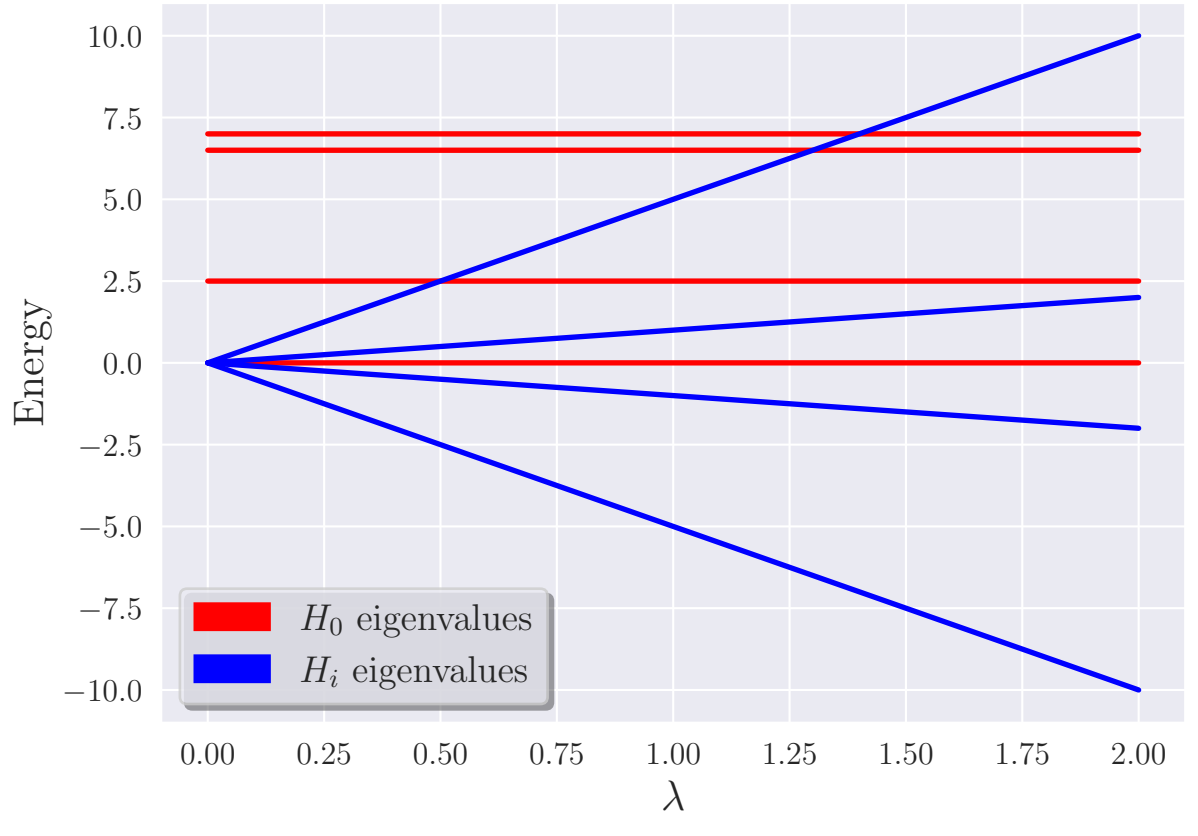
Energy eigenvalues of H_i and H_0 as a function of interaction st
10.0
7.5
5.0
2.5
0.0
−2.5
−5.0
−7.5
−10.0
Energy
H_0 eigenvalues
H_i eigenvalues
0.00
0.25
0.50
0.75
1.00
1.25
1.50
1.75
2.00
λ