

First Code Block:

- builds HO basis
- construct free H (H_0) and interaction V in full basis
- split basis into low and high energy (P and Q), P is first n_P states
- computes full H effective for GS up to third order of V (can be used for excited states as well)
- effective energy given by sum of H_1 , H_2 and H_3
- computes relative error to reference value
- does sweep over omega and plots relative error

Second code block:

- builds HO basis
- construct free H (H_0) and interaction V in full basis
- produce an effective matrix acting only on low energy "P space"
- effective put into a qubit space and then pads matrix with zeros to match qubit space size
- compares GS energy from digitized hamiltonian to known reference value
- plots percentage error vs omega