MCTDH calculation

Name:

Hamiltonian:

- CH3Quasie_exact
- NOCl
- CH4_rst

Potential Energy Surface:

- on
- off

Type of calculation:

- real - time propagation
- imaginary - time propagation
- Eigenstate calculation
- Thermal fluxe eigenstate calculation

Integrator:

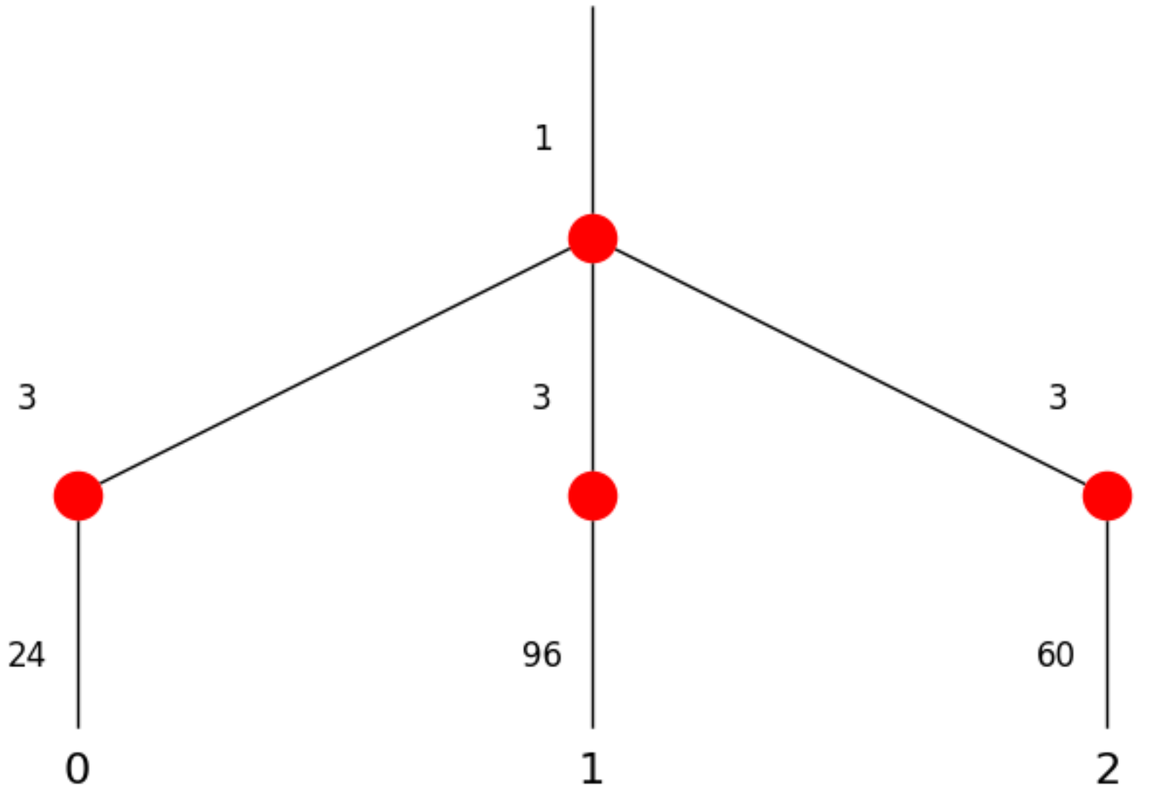
start time: 0 initialtime step: 1

end time: 1985.376 iteration: 100

Scenegraph | Mode

- 1
  - 3
    - 24 0
  - 3
    - 96 1
  - 3

Basis:



Cancel | Save Job | Load... | Start calculation