GQCP::FQMolecularHamiltonian # T # V # g + FQMolecularHamiltonian() + FQMolecularHamiltonian() + kinetic() + nuclearAttraction() + coulombRepulsion()

GQCP::FQMolecularMagnetic Hamiltonian

- + FQMolecularMagneticHamiltonian()
- + FOMolecularMagneticHamiltonian()
- + orbitalZeeman()
- + diamagnetic()



GQCP::FQMolecularPauliHamiltonian

- + FOMolecularPauliHamiltonian()
- + FOMolecularPauliHamiltonian()
- + spinZeeman()