## GOCP::FOMolecularHamiltonian # T # V # q + FOMolecularHamiltonian() + FOMolecularHamiltonian() + kinetic() + nuclearAttraction() + coulombRepulsion() GOCP::FQMolecularMagnetic Hamiltonian + FQMolecularMagneticHamiltonian() + FOMolecularMagneticHamiltonian() + orbitalZeeman() + diamagnetic() GOCP::FOMolecularPauliHamiltonian

+ FQMolecularPauliHamiltonian() + FOMolecularPauliHamiltonian()

+ spinZeeman()