GOCP::FOMolecularHamiltonian # T # V # g + FOMolecularHamiltonian() + FOMolecularHamiltonian() + kinetic() + nuclearAttraction() + coulombRepulsion() GOCP::FQMolecularMagnetic Hamiltonian + FQMolecularMagneticHamiltonian() + FOMolecularMagneticHamiltonian() + orbitalZeeman() + diamagnetic() GQCP::FQMolecularPauliHamiltonian + FOMolecularPauliHamiltonian() + FQMolecularPauliHamiltonian() + spinZeeman()