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