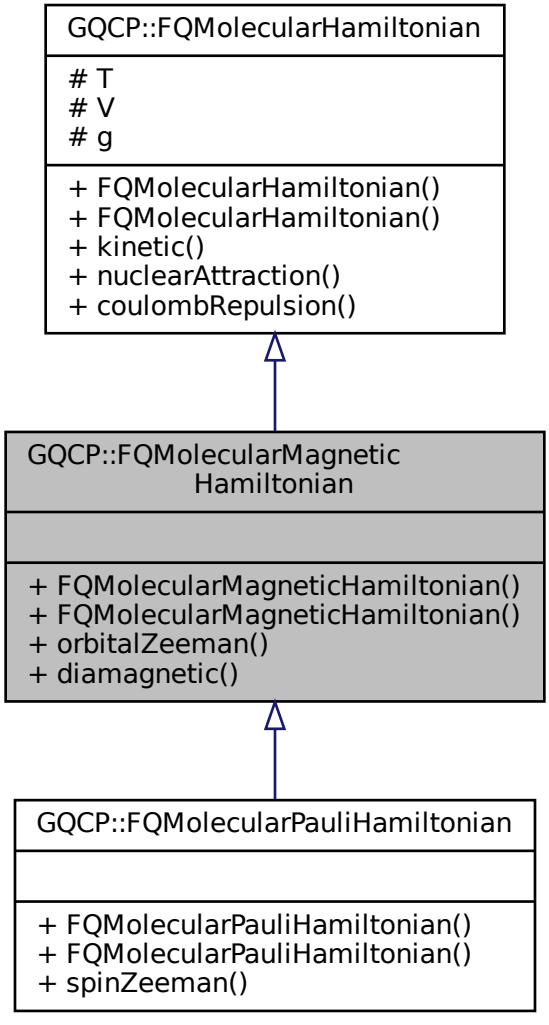
GQCP::FQMolecularHamiltonian
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+ FQMolecularHamiltonian()
+ FQMolecularHamiltonian()
+ kinetic()
+ nuclearAttraction()
+ coulombRepulsion()
GQCP::FQMolecularMagnetic Hamiltonian
+ FQMolecularMagneticHamiltonian()
+ FQMolecularMagneticHamiltonian()
+ orbitalZeeman()
+ diamagnetic()
GQCP::FQMolecularPauliHamiltonian
+ FQMolecularPauliHamiltonian()
+ FQMolecularPauliHamiltonian()
+ spinZeeman()