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# g

+ FQMolecularHamiltonian()
+ FQMolecularHamiltonian()
+ FQMolecularHamiltonian()
+ kinetic()
+ nuclearAttraction()
+ coulombRepulsion()
```

```
+ FQMolecularMagneticHamiltonian()
+ FQMolecularMagneticHamiltonian()
+ paramagnetic()
+ diamagnetic()
```

GQCP::FQMolecularMagnetic