GQCP::QCModel::AP1roG

+ AP1roG()
+ calculateEnergy()
+ calculatePSEJacobian()
+ geminalCoefficients()
+ numberOfElectronPairs()
+ numberOfSpatialOrbitals()
+ orbitalSpace()
+ calculateEnergy()
+ calculatePSECoordinateFunction()
+ calculatePSECoordinateFunctions()
+ calculatePSEJacobian()
+ callablePSEJacobian()
+ callablePSECoordinateFunctions()
+ calculatePSEJacobianElement()