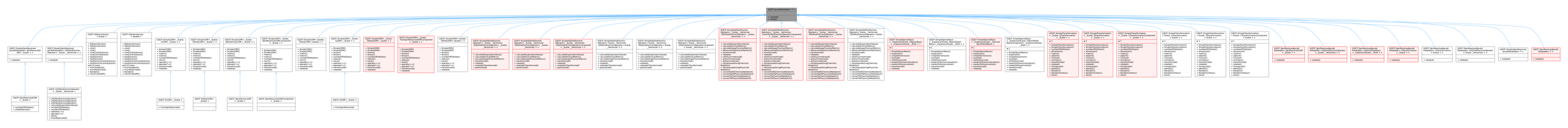

GQCP::JacobiRotatable< T >

+ rotated()
+ rotate()

GQCP::DoublySpinResolved
JacobiRotatable< SpinResolved2
DM< _Scalar > >

+ rotated()

GQCP::DoublySpinResolved
JacobiRotatable< USQTwoElectron
Operator< _Scalar, _Vectorizer > >

+ rotated()

GQCP::SQHamiltonian
< Scalar >

+ SQHamiltonian()
+ SQHamiltonian()
+ core()
+ core()
+ coreContributions()
+ coreContributions()
+ twoElectron()
+ twoElectron()
+ twoElectronContributions()
+ twoElectronContributions()
and 14 more...
+ FromHubbard()
+ FromRestricted()
+ Random()
+ FromFCIDUMP()

GQCP::SQHamiltonian
< double >

+ SQHamiltonian()
+ SQHamiltonian()
+ core()
+ core()
+ coreContributions()
+ coreContributions()
+ twoElectron()
+ twoElectron()
+ twoElectronContributions()
+ twoElectronContributions()
and 14 more...
+ FromHubbard()
+ FromRestricted()
+ Random()
+ FromFCIDUMP()

GQCP::Simple1DM< _Scalar,
G1DM< _Scalar > >

+ Simple1DM()
+ Simple1DM()
+ matrix()
+ matrix()
+ numberOfOrbitals()
+ norm()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::Simple1DM< _Scalar,
Orbital1DM< _Scalar > >

+ Simple1DM()
+ Simple1DM()
+ matrix()
+ matrix()
+ numberOfOrbitals()
+ norm()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::Simple1DM< _Scalar,
SpinDensity1DM< _Scalar > >

+ Simple1DM()
+ Simple1DM()
+ matrix()
+ matrix()
+ numberOfOrbitals()
+ norm()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::Simple1DM< _Scalar,
SpinResolved1DMComponent
< _Scalar > >

+ Simple1DM()
+ Simple1DM()
+ matrix()
+ matrix()
+ numberOfOrbitals()
+ norm()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::Simple1DM< double,
Orbital1DM< double > >

+ Simple1DM()
+ Simple1DM()
+ matrix()
+ matrix()
+ numberOfOrbitals()
+ norm()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::Simple2DM< _Scalar,
G2DM< _Scalar > >

+ Simple2DM()
+ Simple2DM()
+ tensor()
+ tensor()
+ numberOfOrbitals()
+ reduce()
+ trace()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::Simple2DM< _Scalar,
Orbital2DM< _Scalar > >

+ Simple2DM()
+ Simple2DM()
+ tensor()
+ tensor()
+ numberOfOrbitals()
+ reduce()
+ trace()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::Simple2DM< _Scalar,
PureSpinResolved2DMComponent
< _Scalar > >

+ Simple2DM()
+ Simple2DM()
+ tensor()
+ tensor()
+ numberOfOrbitals()
+ reduce()
+ trace()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::Simple2DM< double,
Orbital2DM< double > >

+ Simple2DM()
+ Simple2DM()
+ tensor()
+ tensor()
+ numberOfOrbitals()
+ reduce()
+ trace()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()

GQCP::SimpleSQOneElectron
Operator< _Scalar, _Vectorizer,
GSQOneElectronOperator< _Scalar,
_Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ transformed()
+ rotated()
+ oneIndexTransformed()
+ partitioned()

GQCP::SimpleSQOneElectron
Operator< _Scalar, _Vectorizer,
RSQOneElectronOperator< _Scalar,
_Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ transformed()
+ rotated()
+ oneIndexTransformed()
+ partitioned()

GQCP::SimpleSQOneElectron
Operator< _Scalar, _Vectorizer,
USQOneElectronOperatorComponent
< _Scalar, _Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ transformed()
+ rotated()
+ oneIndexTransformed()
+ partitioned()

GQCP::SimpleSQOneElectron
Operator< Scalar, _Vectorizer,
GSQOneElectronOperator< Scalar,
_Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ transformed()
+ rotated()
+ oneIndexTransformed()
+ partitioned()

GQCP::SimpleSQOneElectron
Operator< Scalar, _Vectorizer,
RSQOneElectronOperator< Scalar,
_Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ transformed()
+ rotated()
+ oneIndexTransformed()
+ partitioned()

GQCP::SimpleSQOneElectron
Operator< Scalar, _Vectorizer,
USQOneElectronOperatorComponent
< Scalar, _Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ transformed()
+ rotated()
+ oneIndexTransformed()
+ partitioned()

GQCP::SimpleSQTwoElectron
Operator< _Scalar, _Vectorizer,
GSQTwoElectronOperator< _Scalar,
_Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ effectiveOneElectronPartition()
+ transformed()
+ rotated()
+ isAntisymmetrized()
+ antisymmetrized()
+ antisymmetrize()
+ isExpressedUsingChemists
Notation()
+ isExpressedUsingPhysicists
Notation()
+ convertedToChemistsNotation()
+ convertedToPhysicistsNotation()
+ convertToChemistsNotation()
+ convertToPhysicistsNotation()

GQCP::SimpleSQTwoElectron
Operator< _Scalar, _Vectorizer,
PureUSQTwoElectronOperatorComponent
< _Scalar, _Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ effectiveOneElectronPartition()
+ transformed()
+ rotated()
+ isAntisymmetrized()
+ antisymmetrized()
+ antisymmetrize()
+ isExpressedUsingChemists
Notation()
+ isExpressedUsingPhysicists
Notation()
+ convertedToChemistsNotation()
+ convertedToPhysicistsNotation()
+ convertToChemistsNotation()
+ convertToPhysicistsNotation()

GQCP::SimpleSQTwoElectron
Operator< _Scalar, _Vectorizer,
RSQTwoElectronOperator< _Scalar,
_Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ effectiveOneElectronPartition()
+ transformed()
+ rotated()
+ isAntisymmetrized()
+ antisymmetrized()
+ antisymmetrize()
+ isExpressedUsingChemists
Notation()
+ isExpressedUsingPhysicists
Notation()
+ convertedToChemistsNotation()
+ convertedToPhysicistsNotation()
+ convertToChemistsNotation()
+ convertToPhysicistsNotation()

GQCP::SimpleSQTwoElectron
Operator< Scalar, _Vectorizer,
GSQTwoElectronOperator< Scalar,
_Vectorizer > >

+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ effectiveOneElectronPartition()
+ transformed()
+ rotated()
+ isAntisymmetrized()
+ antisymmetrized()
+ antisymmetrize()
+ isExpressedUsingChemists
Notation()
+ isExpressedUsingPhysicists
Notation()
+ convertedToChemistsNotation()
+ convertedToPhysicistsNotation()
+ convertToChemistsNotation()
+ convertToPhysicistsNotation()

GQCP::SimpleSpinorBasis
< _ExpansionScalar, GSpinorBasis
< _ExpansionScalar, _Shell > >

C

+ SimpleSpinorBasis()
+ expansion()
+ simpleDimension()
+ overlap()
+ isOrthonormal()
+ lowdinOrthonormalization()
+ lowdinOrthonormalize()
+ transformed()
+ rotated()

GQCP::SimpleSpinorBasis
< _ExpansionScalar, RSpinOrbital
Basis< _ExpansionScalar, _Shell > >

C

+ SimpleSpinorBasis()
+ expansion()
+ simpleDimension()
+ overlap()
+ isOrthonormal()
+ lowdinOrthonormalization()
+ lowdinOrthonormalize()
+ transformed()
+ rotated()

GQCP::SimpleSpinorBasis
< _ExpansionScalar, _Derived
SpinOrbitalBasis >

C

+ SimpleSpinorBasis()
+ expansion()
+ simpleDimension()
+ overlap()
+ isOrthonormal()
+ lowdinOrthonormalization()
+ lowdinOrthonormalize()
+ transformed()
+ rotated()

GQCP::SimpleSpinorBasis
< _ExpansionScalar, USpinOrbital
BasisComponent< _ExpansionScalar,
_Shell > >

C

+ SimpleSpinorBasis()
+ expansion()
+ simpleDimension()
+ overlap()
+ isOrthonormal()
+ lowdinOrthonormalization()
+ lowdinOrthonormalize()
+ transformed()
+ rotated()

GQCP::SimpleTransformation
< _Scalar, GTransformation
< _Scalar > >

T

+ SimpleTransformation()
+ SimpleTransformation()
+ numberOfOrbitals()
+ dimension()
+ matrix()
+ adjoint()
+ inverse()
+ isUnitary()
+ transformed()
+ rotated()
+ FromJacobi()
+ Identity()
+ Random()
+ RandomUnitary()
+ Zero()

GQCP::SimpleTransformation
< _Scalar, RTransformation
< _Scalar > >

T

+ SimpleTransformation()
+ SimpleTransformation()
+ numberOfOrbitals()
+ dimension()
+ matrix()
+ adjoint()
+ inverse()
+ isUnitary()
+ transformed()
+ rotated()
+ FromJacobi()
+ Identity()
+ Random()
+ RandomUnitary()
+ Zero()

GQCP::SimpleTransformation
< _Scalar, UTransformationComponent
< _Scalar > >

T

+ SimpleTransformation()
+ SimpleTransformation()
+ numberOfOrbitals()
+ dimension()
+ matrix()
+ adjoint()
+ inverse()
+ isUnitary()
+ transformed()
+ rotated()
+ FromJacobi()
+ Identity()
+ Random()
+ RandomUnitary()
+ Zero()

GQCP::SimpleTransformation
< Scalar, GTransformation
< Scalar > >

T

+ SimpleTransformation()
+ SimpleTransformation()
+ numberOfOrbitals()
+ dimension()
+ matrix()
+ adjoint()
+ inverse()
+ isUnitary()
+ transformed()
+ rotated()
+ FromJacobi()
+ Identity()
+ Random()
+ RandomUnitary()
+ Zero()

GQCP::SimpleTransformation
< Scalar, RTransformation
< Scalar > >

T

+ SimpleTransformation()
+ SimpleTransformation()
+ numberOfOrbitals()
+ dimension()
+ matrix()
+ adjoint()
+ inverse()
+ isUnitary()
+ transformed()
+ rotated()
+ FromJacobi()
+ Identity()
+ Random()
+ RandomUnitary()
+ Zero()

GQCP::SimpleTransformation
< Scalar, UTransformationComponent
< Scalar > >

T

+ SimpleTransformation()
+ SimpleTransformation()
+ numberOfOrbitals()
+ dimension()
+ matrix()
+ adjoint()
+ inverse()
+ isUnitary()
+ transformed()
+ rotated()
+ FromJacobi()
+ Identity()
+ Random()
+ RandomUnitary()
+ Zero()

GQCP::SpinResolvedJacobi
Rotatable< SpinResolved1DM
< _Scalar > >

+ rotated()

GQCP::SpinResolvedJacobi
Rotatable< USQOneElectronOperator
< _Scalar, _Vectorizer > >

+ rotated()

GQCP::SpinResolvedJacobi
Rotatable< USpinOrbitalBasis
< _ExpansionScalar, _Shell > >

+ rotated()

GQCP::SpinResolvedJacobi
Rotatable< UTransformation
< _Scalar > >

+ rotated()

GQCP::SpinResolvedJacobi
Rotatable< UTransformation
< Scalar > >

+ rotated()

GQCP::SpinResolvedJacobi
Rotatable< USQOneElectronOperator
< Scalar, _Vectorizer > >

+ rotated()

GQCP::DoublySpinResolved
JacobiRotatable< T >

+ rotated()

GQCP::SpinResolvedJacobi
Rotatable< T >

+ rotated()

GQCP::SpinResolved2DM
< _Scalar >

+ numberOfOrbitals()
+ orbitalDensity()

GQCP::USQTwoElectronOperator
< _Scalar, _Vectorizer >

+ USQTwoElectronOperator()
+ USQTwoElectronOperator()
+ USQTwoElectronOperator()
+ calculateExpectationValue()
+ numberOfOrbitals()
+ numberOfOrbitals()
+ operator+=()
+ operator*=()
+ Zero()
+ FromRestricted()

GQCP::G1DM< _Scalar >

+ FromSpinResolved()

GQCP::Orbital1DM<
_Scalar >

GQCP::SpinDensity1DM
< _Scalar >

GQCP::SpinResolved1DMComponent
< _Scalar > GQCP::G2DM< _Scalar >

+ FromSpinResolved()