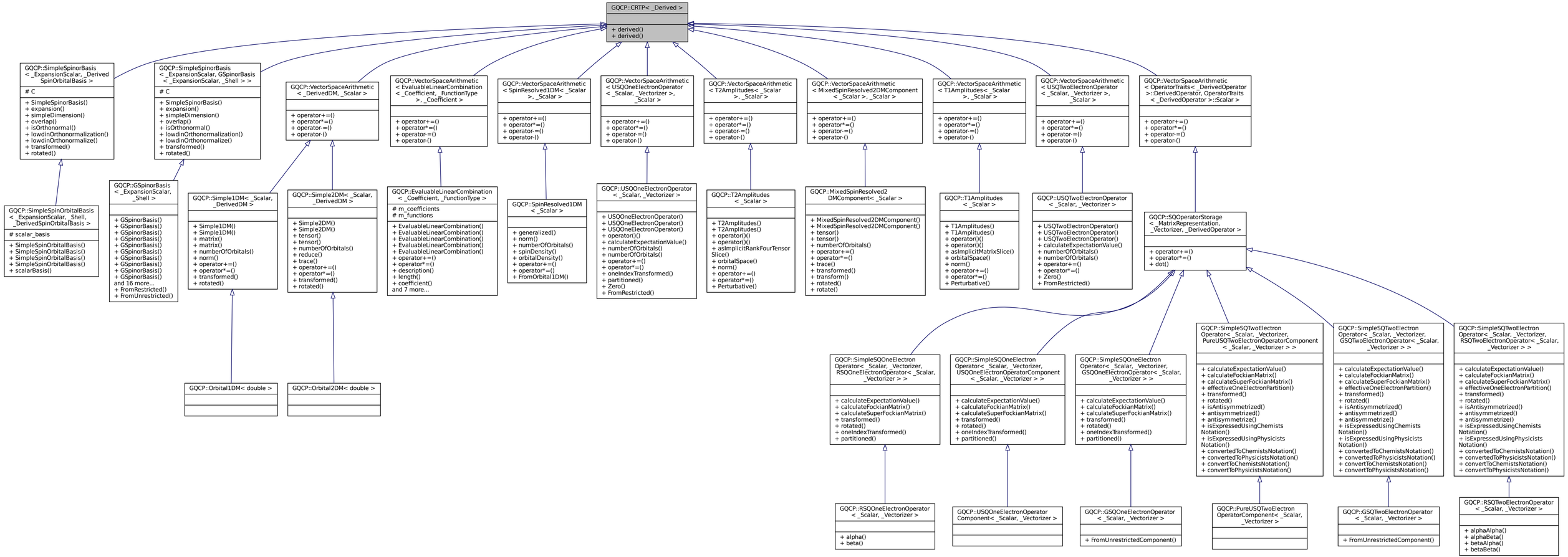
GQCP::CRTP< _Derived >
+ derived()
+ derived()
GQCP::SimpleSpinorBasis < _ExpansionScalar, _DerivedSpinOrbitalBasis >
C
+ SimpleSpinorBasis()
+ expansion()
+ simpleDimension()
+ overlap()
+ isOrthonormal()
+ lowdinOrthonormalization()
+ lowdinOrthonormalize()
+ transformed()
+ rotated()
GQCP::SimpleSpinorBasis < _ExpansionScalar, GSpinorBasis < _ExpansionScalar, _Shell > >
C
+ SimpleSpinorBasis()
+ expansion()
+ simpleDimension()
+ overlap()
+ isOrthonormal()
+ lowdinOrthonormalization()
+ lowdinOrthonormalize()
+ transformed()
+ rotated()
GQCP::VectorSpaceArithmetic < _DerivedDM, _Scalar >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::VectorSpaceArithmetic < EvaluableLinearCombination < _Coefficient, _FunctionType >, _Coefficient >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::VectorSpaceArithmetic < SpinResolved1DM< _Scalar >, _Scalar >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::VectorSpaceArithmetic < USQOneElectronOperator < _Scalar, _Vectorizer >, _Scalar >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::VectorSpaceArithmetic < T2Amplitudes< _Scalar >, _Scalar >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::VectorSpaceArithmetic < MixedSpinResolved2DMComponent < _Scalar >, _Scalar >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::VectorSpaceArithmetic < T1Amplitudes< _Scalar >, _Scalar >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::VectorSpaceArithmetic < USQTwoElectronOperator < _Scalar, _Vectorizer >, _Scalar >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::VectorSpaceArithmetic < OperatorTraits< _DerivedOperator >::DerivedOperator, OperatorTraits < _DerivedOperator >::Scalar >
+ operator+=()
+ operator*=()
+ operator-=()
+ operator-()
GQCP::GSpinorBasis < _ExpansionScalar, _Shell >
+ GSpinorBasis()
+ GSpinorBasis()
+ GSpinorBasis()
+ GSpinorBasis()
+ GSpinorBasis()
+ GSpinorBasis()
+ GSpinorBasis()
+ GSpinorBasis()
+ GSpinorBasis()
+ GSpinorBasis()
and 16 more...
+ FromRestricted()
+ FromUnrestricted()
GQCP::SimpleSpinOrbitalBasis < _ExpansionScalar, _Shell, _DerivedSpinOrbitalBasis >
scalar_basis
+ SimpleSpinOrbitalBasis()
+ SimpleSpinOrbitalBasis()
+ SimpleSpinOrbitalBasis()
+ SimpleSpinOrbitalBasis()
+ scalarBasis()
GQCP::Simple1DM< _Scalar, _DerivedDM >
+ Simple1DM()
+ Simple1DM()
+ matrix()
+ matrix()
+ numberOfOrbitals()
+ norm()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()
GQCP::Simple2DM< _Scalar, _DerivedDM >
+ Simple2DM()
+ Simple2DM()
+ tensor()
+ tensor()
+ numberOfOrbitals()
+ reduce()
+ trace()
+ operator+=()
+ operator*=()
+ transformed()
+ rotated()
GQCP::EvaluableLinearCombination < _Coefficient, _FunctionType >
m_coefficients
m_functions
+ EvaluableLinearCombination()
+ EvaluableLinearCombination()
+ EvaluableLinearCombination()
+ EvaluableLinearCombination()
+ EvaluableLinearCombination()
+ operator+=()
+ operator*=()
+ description()
+ length()
+ coefficient()
and 7 more...
GQCP::SpinResolved1DM < _Scalar >
+ generalized()
+ norm()
+ numberOfOrbitals()
+ spinDensity()
+ orbitalDensity()
+ operator+=()
+ operator*=()
+ FromOrbital1DM()
GQCP::USQOneElectronOperator < _Scalar, _Vectorizer >
+ USQOneElectronOperator()
+ USQOneElectronOperator()
+ USQOneElectronOperator()
+ operator()()
+ calculateExpectationValue()
+ numberOfOrbitals()
+ numberOfOrbitals()
+ operator+=()
+ operator*=()
+ oneIndexTransformed()
+ partitioned()
+ Zero()
+ fromRestricted()
GQCP::T2Amplitudes < _Scalar >
+ T2Amplitudes()
+ T2Amplitudes()
+ operator()()
+ operator()()
+ asImplicitRankFourTensorSlice()
+ orbitalSpace()
+ norm()
+ operator+=()
+ operator*=()
+ Perturbative()
GQCP::MixedSpinResolved2DMComponent< _Scalar >
+ MixedSpinResolved2DMComponent()
+ MixedSpinResolved2DMComponent()
+ tensor()
+ tensor()
+ numberOfOrbitals()
+ operator+=()
+ operator*=()
+ trace()
+ transformed()
+ transform()
+ rotated()
+ rotate()
GQCP::T1Amplitudes < _Scalar >
+ T1Amplitudes()
+ T1Amplitudes()
+ operator()()
+ operator()()
+ asImplicitMatrixSlice()
+ orbitalSpace()
+ norm()
+ operator+=()
+ operator*=()
+ Perturbative()
GQCP::USQTwoElectronOperator < _Scalar, _Vectorizer >
+ USQTwoElectronOperator()
+ USQTwoElectronOperator()
+ USQTwoElectronOperator()
+ calculateExpectationValue()
+ numberOfOrbitals()
+ numberOfOrbitals()
+ operator+=()
+ operator*=()
+ Zero()
+ fromRestricted()
GQCP::SQOperatorStorage < _MatrixRepresentation, _Vectorizer, _DerivedOperator >
+ operator+=()
+ operator*=()
+ dot()
GQCP::Orbital1DM< double >
GQCP::Orbital2DM< double >
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::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, 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, RSQTwoElectronOperator< _Scalar, _Vectorizer > >
+ calculateExpectationValue()
+ calculateFockianMatrix()
+ calculateSuperFockianMatrix()
+ effectiveOneElectronPartition()
+ transformed()
+ rotated()
+ isAntisymmetrized()
+ antisymmetrized()
+ antisymmetrize()
+ isExpressedUsingChemists Notation()
+ isExpressedUsingPhysicists Notation()
+ convertedToChemistsNotation()
+ convertedToPhysicistsNotation()
+ convertToChemistsNotation()
+ convertToPhysicistsNotation()
GQCP::RSQOneElectronOperator < _Scalar, _Vectorizer >
+ alpha()
+ beta()
GQCP::USQOneElectronOperator Component< _Scalar, _Vectorizer >
GQCP::GSQOneElectronOperator < _Scalar, _Vectorizer >
+ FromUnrestrictedComponent()
GQCP::PureUSQTwoElectron OperatorComponent< _Scalar, _Vectorizer >
GQCP::GSQTwoElectronOperator < _Scalar, _Vectorizer >
+ FromUnrestrictedComponent()
GQCP::RSQTwoElectronOperator < _Scalar, _Vectorizer >
+ alphaAlpha()
+ alphaBeta()
+ betaAlpha()
+ betaBeta()