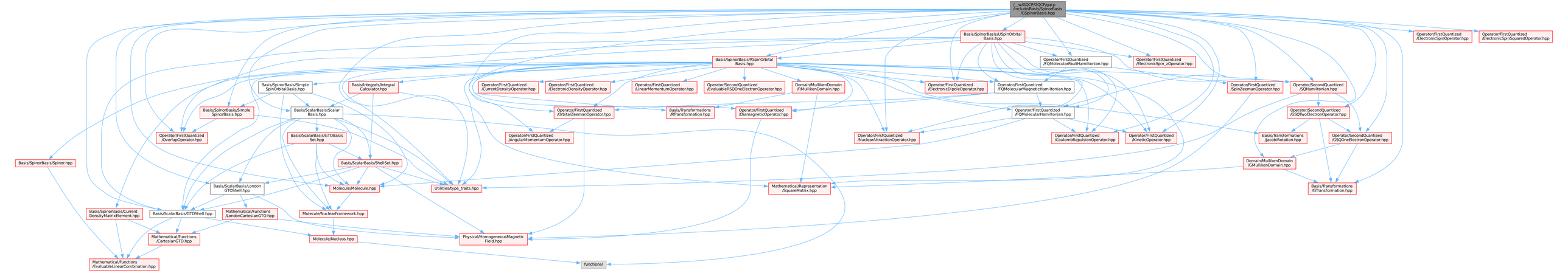
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Basis/SpinorBasis/USpinOrbital
Basis.hpp
Basis/SpinorBasis/RSpinOrbital
Basis.hpp
Operator/FirstQuantized
/FQMolecularPauliHamiltonian.hpp
Operator/FirstQuantized
/ElectronicSpin_zOperator.hpp
Operator/FirstQuantized
/ElectronicSpinOperator.hpp
Operator/FirstQuantized
/ElectronicSpinSquaredOperator.hpp
Basis/SpinorBasis/Simple
SpinOrbitalBasis.hpp
Basis/Integrals/Integral
Calculator.hpp
Operator/FirstQuantized
/CurrentDensityOperator.hpp
Operator/FirstQuantized
/ElectronicDensityOperator.hpp
Operator/FirstQuantized
/LinearMomentumOperator.hpp
Operator/SecondQuantized
/EvaluableRSQOneElectronOperator.hpp
Domain/MullikenDomain
/RMullikenDomain.hpp
Operator/FirstQuantized
/ElectronicDipoleOperator.hpp
Operator/FirstQuantized
/FQMolecularMagneticHamiltonian.hpp
Operator/FirstQuantized
/SpinZeemanOperator.hpp
Operator/SecondQuantized
/SQHamiltonian.hpp
Basis/SpinorBasis/Simple
SpinorBasis.hpp
Basis/ScalarBasis/Scalar
Basis.hpp
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/OrbitalZeemanOperator.hpp
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/RTransformation.hpp
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/FQMolecularHamiltonian.hpp
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/GSQTwoElectronOperator.hpp
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/OverlapOperator.hpp
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Set.hpp
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/AngularMomentumOperator.hpp
Operator/FirstQuantized
/NuclearAttractionOperator.hpp
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/CoulombRepulsionOperator.hpp
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/KineticOperator.hpp
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/JacobiRotation.hpp
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/GSQOneElectronOperator.hpp
Basis/SpinorBasis/Spinor.hpp
Basis/ScalarBasis/ShellSet.hpp
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/GMullikenDomain.hpp
Basis/ScalarBasis/London
GTOShell.hpp
Molecule/Molecule.hpp
Utilities/type_traits.hpp
Mathematical/Representation
/SquareMatrix.hpp
Basis/Transformations
/GTransformation.hpp
Basis/SpinorBasis/Current
DensityMatrixElement.hpp
Basis/ScalarBasis/GTOShell.hpp
Mathematical/Functions
/LondonCartesianGTO.hpp
Molecule/NuclearFramework.hpp
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/CartesianGTO.hpp
Molecule/Nucleus.hpp
Physical/HomogeneousMagnetic
Field.hpp
Mathematical/Functions
/EvaluableLinearCombination.hpp
functional