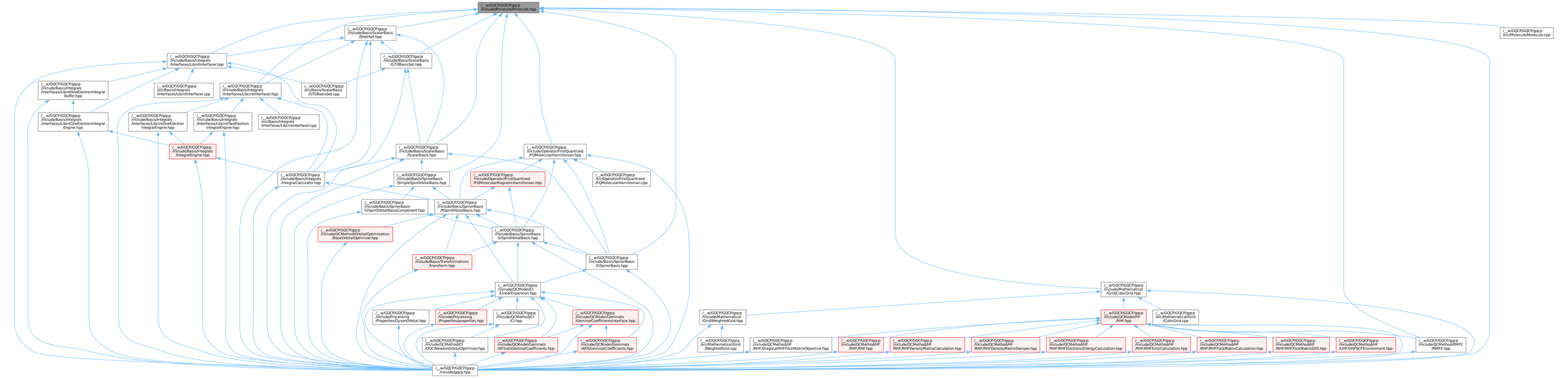
/__w/GQCP/GQCP/gqcp
/include/Molecule/Molecule.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/ScalarBasis
/BaseSet.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Integrals
/Interfaces/LibintInterfacer.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/ScalarBasis
/GTOBasisSet.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Integrals
/Interfaces/LibintOneElectronIntegral
Buffer.hpp
/__w/GQCP/GQCP/gqcp
/src/Basis/Integrals
/Interfaces/LibintInterfacer.cpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Integrals
/Interfaces/LibcintInterfacer.hpp
/__w/GQCP/GQCP/gqcp
/src/Basis/ScalarBasis
/GTOBasisSet.cpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Integrals
/Interfaces/LibintOneElectronIntegral
Engine.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Integrals
/Interfaces/LibintOneElectron
IntegralEngine.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Integrals
/Interfaces/LibcintTwoElectron
IntegralEngine.hpp
/__w/GQCP/GQCP/gqcp
/src/Basis/Integrals
/Interfaces/LibcintInterfacer.cpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Integrals
/IntegralEngine.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/ScalarBasis
/ScalarBasis.hpp
/__w/GQCP/GQCP/gqcp
/include/Operator/FirstQuantized
/FQMolecularHamiltonian.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Integrals
/IntegralCalculator.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/SpinorBasis
/SimpleSpinOrbitalBasis.hpp
/__w/GQCP/GQCP/gqcp
/include/Operator/FirstQuantized
/FQMolecularMagneticHamiltonian.hpp
/__w/GQCP/GQCP/gqcp
/src/Operator/FirstQuantized
/FQMolecularHamiltonian.cpp
/__w/GQCP/GQCP/gqcp
/include/Basis/SpinorBasis
/USpinOrbitalBasisComponent.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/SpinorBasis
/RSpinOrbitalBasis.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/OrbitalOptimization
/BaseOrbitalOptimizer.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/SpinorBasis
/USpinOrbitalBasis.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/Transformations
/transform.hpp
/__w/GQCP/GQCP/gqcp
/include/Basis/SpinorBasis
/GSpinorBasis.hpp
/__w/GQCP/GQCP/gqcp
/include/QCModel/CI
/LinearExpansion.hpp
/__w/GQCP/GQCP/gqcp
/include/Processing
/Properties/DysonOrbital.hpp
/__w/GQCP/GQCP/gqcp
/include/Processing
/Properties/properties.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/CI
/CI.hpp
/__w/GQCP/GQCP/gqcp
/include/QCModel/Geminals
/GeminalCoefficientsInterface.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/CI
/DOCINewtonOrbitalOptimizer.hpp
/__w/GQCP/GQCP/gqcp
/include/QCModel/Geminals
/AP1roGeminalCoefficients.hpp
/__w/GQCP/GQCP/gqcp
/include/QCModel/Geminals
/APIGGeminalCoefficients.hpp
/__w/GQCP/GQCP/gqcp
/include/gqcp.hpp
/__w/GQCP/GQCP/gqcp
/include/Mathematical
/Grid/CubicGrid.hpp
/__w/GQCP/GQCP/gqcp
/src/Molecule/Molecule.cpp
/__w/GQCP/GQCP/gqcp
/Grid/WeightedGrid.hpp
/__w/GQCP/GQCP/gqcp
/include/QCModel/HF
/RHF.hpp
/__w/GQCP/GQCP/gqcp
/src/Mathematical/Grid
/CubicGrid.cpp
/__w/GQCP/GQCP/gqcp
/src/Mathematical/Grid
/WeightedGrid.cpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/RHF/DiagonalRHFFockMatrixObjective.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/RHF/RHF.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/RHF/RHFDensityMatrixCalculation.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/RHF/RHFDensityMatrixDamper.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/RHF/RHFElectronicEnergyCalculation.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/RHF/RHFErrorCalculation.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/RHF/RHFFockMatrixCalculation.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/RHF/RHFFockMatrixDIIS.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/HF
/UHF/UHFSCFEnvironment.hpp
/__w/GQCP/GQCP/gqcp
/include/QCMethod/RMP2
/RMP2.hpp