



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:19 PM UTC

PDB ID : 1FHE / pdb_00001fhe
Title : GLUTATHIONE TRANSFERASE (FH47) FROM FASCIOLA HEPATICA
Authors : Rossjohn, J.; Parker, M.W.
Deposited on : 1997-07-24
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

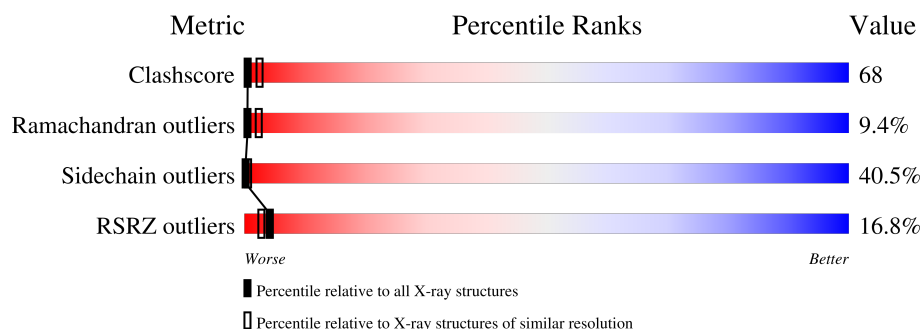
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	217	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GSH	A	300	X	X	-	-

2 Entry composition [i](#)

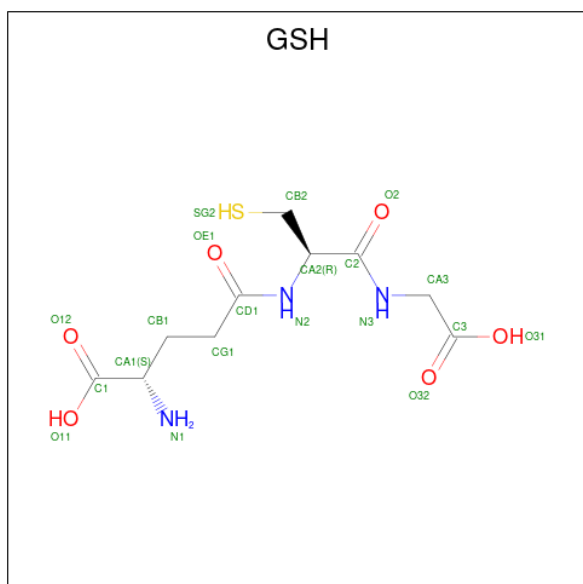
There are 3 unique types of molecules in this entry. The entry contains 1810 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called GLUTATHIONE TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1760	1139	293	316	12			

- Molecule 2 is GLUTATHIONE (CCD ID: GSH) (formula: C₁₀H₁₇N₃O₆S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			20	10	3	6	1		

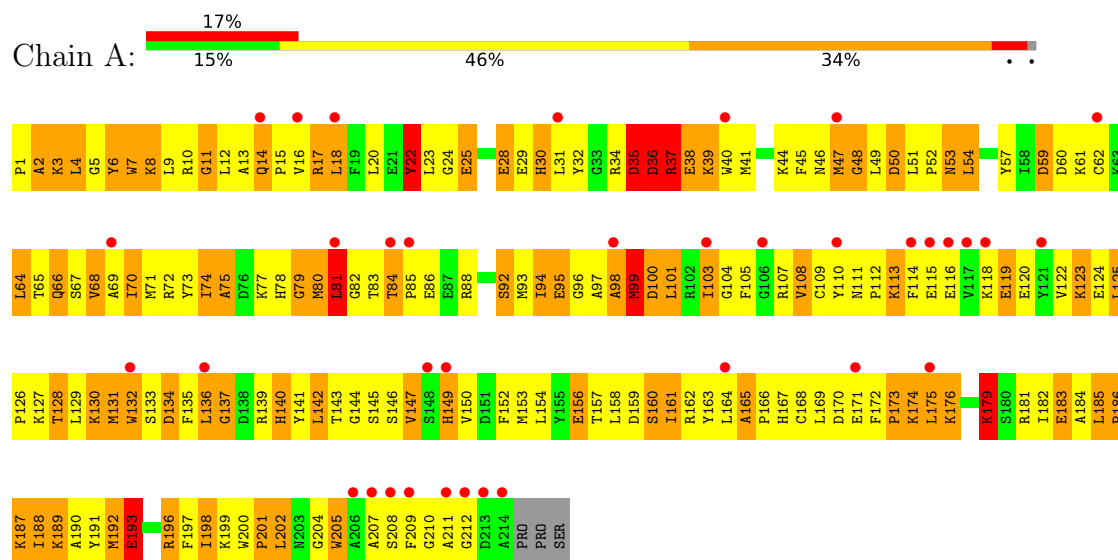
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total	O	0	0
			30	30		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: GLUTATHIONE TRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	158.04Å 158.04Å 74.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.00 30.00 – 3.04	Depositor EDS
% Data completeness (in resolution range)	94.0 (30.00-3.00) 90.5 (30.00-3.04)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.00 (at 3.05Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.237 , 0.354 0.331 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	92.5	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 64.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	1810	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.61% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GSH

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.34	0/1808	0.97	13/2433 (0.5%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	185	LEU	CA-C-N	7.72	129.50	119.84
1	A	185	LEU	C-N-CA	7.72	129.50	119.84
1	A	37	ARG	N-CA-C	-7.20	102.84	111.69
1	A	36	ASP	N-CA-C	6.98	119.81	110.88
1	A	35	ASP	N-CA-C	-6.08	105.95	113.19
1	A	75	ALA	N-CA-C	-5.97	106.13	113.41
1	A	136	LEU	N-CA-C	-5.81	105.90	112.87
1	A	98	ALA	N-CA-C	-5.49	106.42	113.23
1	A	60	ASP	N-CA-C	-5.43	105.52	111.82
1	A	99	MET	N-CA-C	-5.25	107.01	113.41
1	A	143	THR	N-CA-C	5.21	118.77	112.93
1	A	22	TYR	N-CA-C	-5.12	105.60	111.07
1	A	1	PRO	N-CA-CB	5.00	108.50	103.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1760	0	1744	239	0
2	A	20	0	16	5	0
3	A	30	0	0	2	0
All	All	1810	0	1760	239	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 68.

All (239) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:VAL:HG23	1:A:71:MET:HG3	1.45	0.99
1:A:165:ALA:H	1:A:166:PRO:HD3	1.26	0.97
1:A:158:LEU:HD11	1:A:175:LEU:HD13	1.45	0.97
1:A:207:ALA:HB3	1:A:211:ALA:HB2	1.47	0.95
1:A:44:LYS:HE3	1:A:52:PRO:HB2	1.49	0.95
1:A:18:LEU:HB3	1:A:152:PHE:CD2	2.09	0.87
1:A:53:ASN:HD21	2:A:300:GSH:HG13	1.38	0.86
1:A:57:TYR:HB2	1:A:70:ILE:HD13	1.57	0.85
1:A:112:PRO:HB2	1:A:113:LYS:HD3	1.57	0.85
1:A:22:TYR:CE2	1:A:147:VAL:HB	2.13	0.84
1:A:18:LEU:HB3	1:A:152:PHE:HD2	1.41	0.84
1:A:2:ALA:HB1	1:A:57:TYR:HE1	1.44	0.82
1:A:158:LEU:O	1:A:161:ILE:HG13	1.78	0.81
1:A:9:LEU:HD22	1:A:204:GLY:HA2	1.63	0.80
1:A:22:TYR:HE2	1:A:147:VAL:HB	1.46	0.80
1:A:2:ALA:HB1	1:A:57:TYR:CE1	2.18	0.77
1:A:133:SER:OG	1:A:175:LEU:HG	1.83	0.77
1:A:165:ALA:N	1:A:166:PRO:HD3	1.97	0.76
1:A:165:ALA:H	1:A:166:PRO:CD	1.99	0.76
1:A:7:TRP:HA	1:A:32:TYR:H	1.51	0.75
1:A:8:LYS:HG2	1:A:31:LEU:HD13	1.70	0.73
1:A:57:TYR:HB2	1:A:70:ILE:HG23	1.70	0.72
1:A:14:GLN:HB3	1:A:156:GLU:HG3	1.71	0.72
1:A:36:ASP:HB3	1:A:39:LYS:HB3	1.70	0.72
1:A:10:ARG:HG2	1:A:198:ILE:HG22	1.72	0.71
1:A:68:VAL:HA	1:A:71:MET:HE3	1.74	0.70
1:A:105:PHE:O	1:A:108:VAL:HG22	1.92	0.70
1:A:124:GLU:O	1:A:128:THR:HG23	1.92	0.70
1:A:18:LEU:HG	1:A:188:ILE:CD1	2.21	0.70
1:A:196:ARG:HD2	3:A:317:HOH:O	1.91	0.69
1:A:130:LYS:O	1:A:134:ASP:HB2	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:MET:O	1:A:157:THR:HG23	1.93	0.69
1:A:72:ARG:HD2	1:A:81:LEU:HD22	1.74	0.68
1:A:66:GLN:HB2	1:A:69:ALA:HB3	1.76	0.68
1:A:164:LEU:HB2	1:A:209:PHE:CE1	2.30	0.66
1:A:40:TRP:CH2	1:A:44:LYS:HA	2.30	0.66
1:A:85:PRO:N	1:A:88:ARG:HH12	1.94	0.66
1:A:122:VAL:O	1:A:126:PRO:HD3	1.95	0.66
1:A:168:CYS:O	1:A:171:GLU:HB3	1.96	0.66
1:A:124:GLU:HA	1:A:127:LYS:HG3	1.78	0.65
1:A:200:TRP:CD2	1:A:201:PRO:HA	2.31	0.65
1:A:7:TRP:O	1:A:9:LEU:N	2.31	0.64
1:A:9:LEU:CD2	1:A:204:GLY:HA2	2.27	0.64
1:A:71:MET:O	1:A:74:ILE:HG22	1.98	0.64
1:A:165:ALA:N	1:A:166:PRO:CD	2.57	0.64
1:A:187:LYS:HD3	1:A:190:ALA:HB3	1.80	0.63
1:A:65:THR:HA	1:A:70:ILE:HD11	1.81	0.63
1:A:99:MET:O	1:A:103:ILE:HB	1.98	0.62
1:A:49:LEU:O	1:A:52:PRO:HD3	1.99	0.62
1:A:65:THR:O	1:A:66:GLN:HB2	1.98	0.62
1:A:207:ALA:HB3	1:A:211:ALA:CB	2.26	0.62
1:A:52:PRO:O	1:A:53:ASN:HB2	1.97	0.62
1:A:170:ASP:HA	1:A:176:LYS:CD	2.30	0.62
1:A:18:LEU:HD12	1:A:191:TYR:CE2	2.35	0.61
1:A:81:LEU:O	1:A:88:ARG:HG3	2.01	0.61
1:A:7:TRP:CE3	1:A:32:TYR:HB3	2.36	0.60
1:A:97:ALA:HA	1:A:100:ASP:HB2	1.81	0.60
1:A:130:LYS:HA	1:A:172:PHE:CZ	2.36	0.60
1:A:53:ASN:CG	1:A:54:LEU:H	2.09	0.60
1:A:169:LEU:HG	1:A:176:LYS:HE3	1.82	0.60
1:A:46:ASN:C	1:A:48:GLY:H	2.09	0.60
1:A:139:ARG:HE	1:A:144:GLY:HA2	1.68	0.59
1:A:36:ASP:HB3	1:A:39:LYS:CB	2.31	0.59
1:A:132:TRP:HA	1:A:132:TRP:CE3	2.37	0.59
1:A:53:ASN:HD21	2:A:300:GSH:CG1	2.15	0.58
1:A:147:VAL:HG21	1:A:185:LEU:HD11	1.86	0.58
1:A:18:LEU:HD12	1:A:191:TYR:HE2	1.68	0.58
1:A:10:ARG:O	1:A:11:GLY:C	2.47	0.58
1:A:14:GLN:OE1	1:A:14:GLN:HA	2.04	0.58
1:A:156:GLU:O	1:A:160:SER:HB2	2.04	0.58
1:A:129:LEU:HD22	1:A:175:LEU:HD12	1.85	0.57
1:A:8:LYS:O	1:A:9:LEU:HD23	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:TRP:HA	1:A:32:TYR:N	2.18	0.57
1:A:7:TRP:CD1	1:A:7:TRP:H	2.20	0.57
1:A:65:THR:O	1:A:66:GLN:CB	2.51	0.57
1:A:119:GLU:HG3	1:A:120:GLU:N	2.18	0.57
1:A:122:VAL:HA	1:A:125:LEU:HB2	1.87	0.57
1:A:88:ARG:HH11	1:A:88:ARG:HB2	1.70	0.56
1:A:209:PHE:CD1	1:A:210:GLY:N	2.73	0.56
1:A:44:LYS:O	1:A:44:LYS:HG3	2.05	0.56
1:A:149:HIS:CE1	1:A:150:VAL:HG22	2.40	0.56
1:A:181:ARG:O	1:A:184:ALA:HB3	2.04	0.56
1:A:92:SER:HA	1:A:95:GLU:CG	2.35	0.56
1:A:127:LYS:O	1:A:130:LYS:HB3	2.05	0.56
1:A:92:SER:O	1:A:95:GLU:HG3	2.06	0.56
1:A:92:SER:HA	1:A:95:GLU:HG3	1.87	0.55
1:A:188:ILE:O	1:A:192:MET:HG3	2.06	0.55
1:A:10:ARG:HG2	1:A:198:ILE:CG2	2.35	0.55
1:A:59:ASP:HB2	1:A:62:CYS:O	2.05	0.55
1:A:191:TYR:C	1:A:193:GLU:H	2.15	0.54
1:A:64:LEU:HD23	1:A:69:ALA:O	2.08	0.54
1:A:94:ILE:CG2	1:A:150:VAL:HG11	2.38	0.54
1:A:48:GLY:C	1:A:49:LEU:HD23	2.33	0.53
1:A:53:ASN:ND2	2:A:300:GSH:HG13	2.15	0.53
1:A:132:TRP:CZ2	1:A:154:LEU:HD22	2.44	0.53
1:A:53:ASN:OD1	2:A:300:GSH:HG12	2.09	0.53
1:A:167:HIS:CE1	1:A:168:CYS:HB2	2.44	0.52
1:A:44:LYS:HG2	1:A:45:PHE:CE2	2.43	0.52
1:A:57:TYR:CB	1:A:70:ILE:HD13	2.34	0.52
1:A:161:ILE:C	1:A:163:TYR:H	2.17	0.52
1:A:104:GLY:C	1:A:105:PHE:HD1	2.18	0.52
1:A:13:ALA:O	1:A:17:ARG:HB2	2.10	0.52
1:A:34:ARG:C	1:A:36:ASP:H	2.18	0.52
1:A:72:ARG:HH11	1:A:81:LEU:HD22	1.75	0.52
1:A:75:ALA:HB3	1:A:81:LEU:HD21	1.91	0.52
1:A:10:ARG:HB2	1:A:202:LEU:HA	1.90	0.52
1:A:39:LYS:HD2	1:A:39:LYS:O	2.10	0.51
1:A:57:TYR:HB2	1:A:70:ILE:CG2	2.38	0.51
1:A:74:ILE:HD13	1:A:74:ILE:O	2.10	0.51
1:A:84:THR:CG2	1:A:86:GLU:HG2	2.41	0.51
1:A:170:ASP:HA	1:A:176:LYS:HD2	1.91	0.51
1:A:35:ASP:OD1	1:A:35:ASP:O	2.28	0.51
1:A:95:GLU:HA	1:A:98:ALA:HB3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:GLU:C	1:A:126:PRO:HD2	2.35	0.51
1:A:71:MET:HE1	1:A:153:MET:HE2	1.92	0.51
1:A:46:ASN:C	1:A:48:GLY:N	2.68	0.51
1:A:6:TYR:O	1:A:31:LEU:HA	2.11	0.50
1:A:74:ILE:O	1:A:74:ILE:HG23	2.12	0.50
1:A:185:LEU:HD23	1:A:186:PRO:HD2	1.94	0.50
1:A:9:LEU:HD11	1:A:205:TRP:CZ3	2.47	0.50
1:A:159:ASP:HA	1:A:179:LYS:HE2	1.93	0.50
1:A:209:PHE:C	1:A:211:ALA:H	2.20	0.50
1:A:94:ILE:HB	1:A:150:VAL:HG11	1.93	0.50
1:A:53:ASN:ND2	1:A:54:LEU:H	2.10	0.49
1:A:16:VAL:O	1:A:20:LEU:HG	2.12	0.49
1:A:3:LYS:O	1:A:57:TYR:HD1	1.96	0.49
1:A:5:GLY:O	1:A:6:TYR:HB2	2.12	0.49
1:A:66:GLN:O	1:A:67:SER:C	2.56	0.49
1:A:88:ARG:HB2	1:A:88:ARG:NH1	2.28	0.49
1:A:200:TRP:CG	1:A:201:PRO:HA	2.48	0.48
1:A:70:ILE:HG22	1:A:70:ILE:O	2.12	0.48
1:A:57:TYR:CB	1:A:70:ILE:HG23	2.43	0.48
1:A:78:HIS:O	1:A:79:GLY:C	2.57	0.48
1:A:190:ALA:HA	1:A:193:GLU:CD	2.39	0.48
1:A:34:ARG:HG3	1:A:205:TRP:CD2	2.48	0.48
1:A:49:LEU:O	1:A:51:LEU:N	2.47	0.48
1:A:152:PHE:CE2	1:A:188:ILE:HD11	2.48	0.48
1:A:167:HIS:CG	1:A:168:CYS:N	2.82	0.48
1:A:32:TYR:CD2	1:A:54:LEU:HD21	2.48	0.48
1:A:125:LEU:N	1:A:126:PRO:HD2	2.29	0.48
1:A:158:LEU:C	1:A:160:SER:N	2.71	0.48
1:A:3:LYS:HG3	1:A:28:GLU:HB3	1.95	0.48
1:A:62:CYS:SG	1:A:64:LEU:HD13	2.54	0.48
1:A:111:ASN:OD1	1:A:114:PHE:HB3	2.14	0.48
1:A:182:ILE:O	1:A:188:ILE:HG21	2.14	0.47
1:A:18:LEU:HG	1:A:188:ILE:HD11	1.96	0.47
1:A:119:GLU:CG	1:A:120:GLU:N	2.77	0.47
1:A:7:TRP:HB3	1:A:32:TYR:C	2.40	0.47
1:A:47:MET:HE3	1:A:49:LEU:HD11	1.96	0.47
1:A:129:LEU:HA	1:A:132:TRP:HB2	1.96	0.47
1:A:192:MET:HE2	1:A:192:MET:HB3	1.82	0.47
1:A:22:TYR:CD1	1:A:22:TYR:C	2.93	0.46
1:A:49:LEU:HD13	1:A:65:THR:CG2	2.45	0.46
1:A:85:PRO:CD	1:A:88:ARG:HH12	2.27	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:LYS:HD3	1:A:187:LYS:HA	1.58	0.46
1:A:114:PHE:HD1	1:A:114:PHE:H	1.62	0.46
1:A:135:PHE:C	1:A:137:GLY:H	2.23	0.46
1:A:132:TRP:CZ3	1:A:142:LEU:HD11	2.50	0.46
1:A:147:VAL:HG21	1:A:185:LEU:CD1	2.45	0.46
1:A:38:GLU:HA	1:A:41:MET:HB2	1.97	0.46
1:A:135:PHE:C	1:A:137:GLY:N	2.69	0.46
1:A:35:ASP:OD1	1:A:35:ASP:C	2.59	0.46
1:A:68:VAL:HA	1:A:71:MET:CE	2.43	0.46
1:A:204:GLY:O	1:A:211:ALA:HB1	2.16	0.46
1:A:23:LEU:HD11	1:A:80:MET:HE1	1.98	0.45
1:A:161:ILE:C	1:A:163:TYR:N	2.72	0.45
1:A:34:ARG:C	1:A:36:ASP:N	2.74	0.45
1:A:188:ILE:HG22	1:A:189:LYS:N	2.31	0.45
1:A:80:MET:C	1:A:82:GLY:H	2.25	0.45
1:A:110:TYR:HA	1:A:207:ALA:HB1	1.97	0.45
1:A:141:TYR:CD2	1:A:147:VAL:HG13	2.51	0.45
1:A:120:GLU:O	1:A:123:LYS:HB3	2.16	0.45
1:A:192:MET:HA	1:A:197:PHE:CD2	2.52	0.45
1:A:128:THR:O	1:A:132:TRP:HB2	2.16	0.45
1:A:109:CYS:O	1:A:207:ALA:HB1	2.16	0.45
1:A:7:TRP:HZ3	1:A:37:ARG:HA	1.82	0.45
1:A:53:ASN:ND2	2:A:300:GSH:CG1	2.78	0.45
1:A:158:LEU:C	1:A:160:SER:H	2.24	0.45
1:A:38:GLU:C	1:A:40:TRP:N	2.74	0.45
1:A:49:LEU:O	1:A:52:PRO:CD	2.65	0.45
1:A:160:SER:HA	1:A:202:LEU:HD11	1.98	0.45
1:A:57:TYR:CD2	1:A:70:ILE:HG23	2.52	0.44
1:A:190:ALA:O	1:A:193:GLU:HG3	2.17	0.44
1:A:53:ASN:O	1:A:54:LEU:HD12	2.17	0.44
1:A:173:PRO:O	1:A:174:LYS:C	2.60	0.44
1:A:44:LYS:HE3	1:A:52:PRO:CB	2.35	0.44
1:A:183:GLU:HG3	1:A:192:MET:SD	2.56	0.44
1:A:103:ILE:C	1:A:105:PHE:H	2.26	0.44
1:A:7:TRP:CA	1:A:32:TYR:HB2	2.48	0.44
1:A:123:LYS:HE2	1:A:127:LYS:CE	2.47	0.44
1:A:94:ILE:C	1:A:96:GLY:N	2.74	0.44
1:A:125:LEU:N	1:A:126:PRO:CD	2.81	0.44
1:A:10:ARG:HB3	1:A:14:GLN:HG2	2.00	0.43
1:A:72:ARG:HD2	1:A:81:LEU:CD2	2.46	0.43
1:A:84:THR:O	1:A:85:PRO:C	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:LEU:HB3	3:A:305:HOH:O	2.19	0.43
1:A:25:GLU:OE1	1:A:74:ILE:HD11	2.18	0.43
1:A:114:PHE:CD1	1:A:114:PHE:N	2.86	0.43
1:A:23:LEU:HD11	1:A:80:MET:SD	2.58	0.43
1:A:140:HIS:HB3	1:A:141:TYR:CD1	2.54	0.43
1:A:164:LEU:HD22	1:A:209:PHE:CG	2.54	0.43
1:A:110:TYR:HA	1:A:207:ALA:CB	2.49	0.43
1:A:7:TRP:CZ3	1:A:37:ARG:HA	2.53	0.42
1:A:10:ARG:HD2	1:A:156:GLU:OE2	2.19	0.42
1:A:3:LYS:HD3	1:A:30:HIS:CD2	2.53	0.42
1:A:28:GLU:CD	1:A:29:GLU:H	2.25	0.42
1:A:132:TRP:HA	1:A:132:TRP:HE3	1.81	0.42
1:A:83:THR:O	1:A:88:ARG:NH1	2.53	0.42
1:A:4:LEU:HD21	1:A:16:VAL:HG11	2.02	0.42
1:A:66:GLN:O	1:A:70:ILE:HG13	2.18	0.42
1:A:84:THR:C	1:A:88:ARG:HH12	2.27	0.42
1:A:130:LYS:O	1:A:130:LYS:HE2	2.20	0.42
1:A:202:LEU:CD2	1:A:210:GLY:HA3	2.50	0.42
1:A:50:ASP:OD1	1:A:50:ASP:N	2.47	0.42
1:A:103:ILE:HD13	1:A:103:ILE:HA	1.77	0.42
1:A:204:GLY:O	1:A:205:TRP:C	2.63	0.41
1:A:160:SER:HA	1:A:202:LEU:CD1	2.50	0.41
1:A:201:PRO:HB2	1:A:202:LEU:H	1.67	0.41
1:A:103:ILE:HG22	1:A:104:GLY:N	2.35	0.41
1:A:113:LYS:CD	1:A:113:LYS:N	2.84	0.41
1:A:166:PRO:HA	1:A:169:LEU:HB3	2.02	0.41
1:A:4:LEU:HD11	1:A:16:VAL:HG13	2.03	0.41
1:A:59:ASP:HB3	1:A:62:CYS:H	1.86	0.41
1:A:126:PRO:O	1:A:130:LYS:N	2.53	0.41
1:A:14:GLN:CB	1:A:15:PRO:HD3	2.51	0.41
1:A:49:LEU:O	1:A:50:ASP:C	2.64	0.41
1:A:103:ILE:C	1:A:105:PHE:N	2.79	0.41
1:A:130:LYS:HA	1:A:130:LYS:HD2	1.89	0.41
1:A:201:PRO:HG3	1:A:212:GLY:O	2.21	0.41
1:A:72:ARG:O	1:A:75:ALA:HB3	2.20	0.41
1:A:164:LEU:HB2	1:A:209:PHE:CZ	2.56	0.41
1:A:131:MET:O	1:A:131:MET:HG2	2.19	0.40
1:A:4:LEU:HD21	1:A:16:VAL:CG1	2.52	0.40
1:A:7:TRP:HA	1:A:32:TYR:HB2	2.03	0.40
1:A:14:GLN:HG3	1:A:156:GLU:CD	2.46	0.40
1:A:85:PRO:N	1:A:88:ARG:NH1	2.66	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:HIS:CE1	1:A:150:VAL:CG2	3.03	0.40
1:A:53:ASN:CG	1:A:54:LEU:N	2.74	0.40
1:A:114:PHE:O	1:A:118:LYS:N	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	212/217 (98%)	154 (73%)	38 (18%)	20 (9%)	0 2

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	ALA
1	A	8	LYS
1	A	24	GLY
1	A	186	PRO
1	A	11	GLY
1	A	50	ASP
1	A	53	ASN
1	A	66	GLN
1	A	137	GLY
1	A	165	ALA
1	A	173	PRO
1	A	6	TYR
1	A	101	LEU
1	A	81	LEU
1	A	179	LYS
1	A	192	MET
1	A	193	GLU
1	A	79	GLY

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Mol	Chain	Res	Type
1	A	48	GLY
1	A	201	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	185/188 (98%)	110 (60%)	75 (40%)	0 0

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	4	LEU
1	A	7	TRP
1	A	14	GLN
1	A	17	ARG
1	A	18	LEU
1	A	22	TYR
1	A	25	GLU
1	A	28	GLU
1	A	30	HIS
1	A	35	ASP
1	A	36	ASP
1	A	37	ARG
1	A	38	GLU
1	A	39	LYS
1	A	47	MET
1	A	54	LEU
1	A	59	ASP
1	A	61	LYS
1	A	64	LEU
1	A	68	VAL
1	A	70	ILE
1	A	73	TYR
1	A	74	ILE

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Mol	Chain	Res	Type
1	A	77	LYS
1	A	80	MET
1	A	81	LEU
1	A	84	THR
1	A	92	SER
1	A	93	MET
1	A	94	ILE
1	A	95	GLU
1	A	99	MET
1	A	100	ASP
1	A	101	LEU
1	A	103	ILE
1	A	107	ARG
1	A	108	VAL
1	A	113	LYS
1	A	115	GLU
1	A	116	GLU
1	A	119	GLU
1	A	123	LYS
1	A	125	LEU
1	A	128	THR
1	A	130	LYS
1	A	131	MET
1	A	132	TRP
1	A	134	ASP
1	A	136	LEU
1	A	140	HIS
1	A	142	LEU
1	A	145	SER
1	A	146	SER
1	A	147	VAL
1	A	149	HIS
1	A	156	GLU
1	A	160	SER
1	A	161	ILE
1	A	162	ARG
1	A	174	LYS
1	A	175	LEU
1	A	176	LYS
1	A	179	LYS
1	A	183	GLU
1	A	187	LYS

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Mol	Chain	Res	Type
1	A	188	ILE
1	A	189	LYS
1	A	193	GLU
1	A	196	ARG
1	A	198	ILE
1	A	199	LYS
1	A	202	LEU
1	A	205	TRP
1	A	208	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	46	ASN
1	A	53	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GSH	A	300	-	18,19,19	2.35	8 (44%)	21,24,24	2.75	11 (52%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GSH	A	300	-	1/1/6/8	9/24/24/24	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	300	GSH	O32-C3	4.93	1.38	1.22
2	A	300	GSH	CG1-CD1	-3.68	1.44	1.51
2	A	300	GSH	O2-C2	3.55	1.30	1.23
2	A	300	GSH	O12-C1	3.49	1.32	1.22
2	A	300	GSH	OE1-CD1	3.23	1.29	1.23
2	A	300	GSH	CB2-SG2	-2.67	1.76	1.81
2	A	300	GSH	O11-C1	-2.54	1.22	1.30
2	A	300	GSH	CB1-CG1	-2.23	1.46	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300	GSH	CB1-CG1-CD1	6.76	128.14	113.06
2	A	300	GSH	CB2-CA2-N2	4.52	117.74	111.28
2	A	300	GSH	OE1-CD1-CG1	-4.11	114.57	122.02
2	A	300	GSH	O31-C3-O32	-3.57	114.16	123.33
2	A	300	GSH	CB2-CA2-C2	3.40	116.90	109.50
2	A	300	GSH	O2-C2-CA2	-3.32	113.53	120.48
2	A	300	GSH	O31-C3-CA3	3.19	124.96	112.81
2	A	300	GSH	CA3-N3-C2	2.40	127.43	121.38
2	A	300	GSH	OE1-CD1-N2	2.37	126.97	122.95
2	A	300	GSH	CG1-CB1-CA1	-2.25	108.69	113.86
2	A	300	GSH	CA2-N2-CD1	2.01	126.72	121.68

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	300	GSH	CA1

All (9) torsion outliers are listed below:

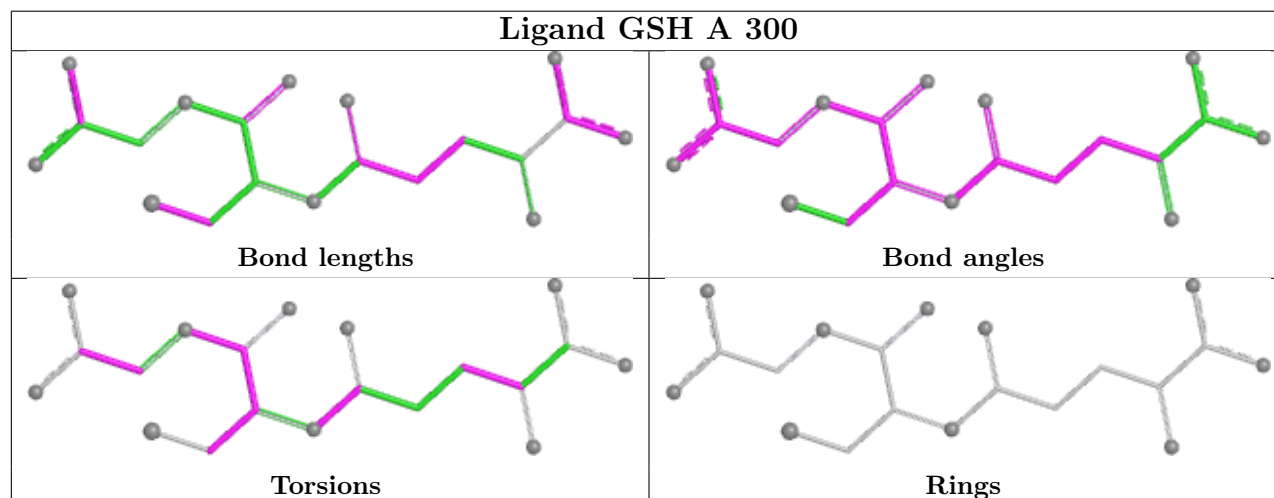
Mol	Chain	Res	Type	Atoms
2	A	300	GSH	N2-CA2-CB2-SG2
2	A	300	GSH	C2-CA2-CB2-SG2
2	A	300	GSH	O2-C2-N3-CA3
2	A	300	GSH	OE1-CD1-N2-CA2
2	A	300	GSH	O31-C3-CA3-N3
2	A	300	GSH	N1-CA1-CB1-CG1
2	A	300	GSH	N3-C2-CA2-N2
2	A	300	GSH	O32-C3-CA3-N3
2	A	300	GSH	C1-CA1-CB1-CG1

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	300	GSH	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	214/217 (98%)	0.87	36 (16%) 4 3	7, 41, 83, 100	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	214	ALA	7.3
1	A	208	SER	6.0
1	A	115	GLU	5.7
1	A	207	ALA	4.3
1	A	110	TYR	3.7
1	A	117	VAL	3.7
1	A	85	PRO	3.6
1	A	212	GLY	3.6
1	A	118	LYS	3.5
1	A	116	GLU	3.4
1	A	40	TRP	3.3
1	A	209	PHE	3.1
1	A	171	GLU	3.1
1	A	114	PHE	2.9
1	A	62	CYS	2.8
1	A	84	THR	2.6
1	A	106	GLY	2.6
1	A	69	ALA	2.5
1	A	149	HIS	2.5
1	A	18	LEU	2.5
1	A	98	ALA	2.4
1	A	121	TYR	2.4
1	A	148	SER	2.4
1	A	132	TRP	2.3
1	A	164	LEU	2.3
1	A	213	ASP	2.3
1	A	103	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	211	ALA	2.3
1	A	47	MET	2.2
1	A	16	VAL	2.2
1	A	14	GLN	2.1
1	A	206	ALA	2.1
1	A	81	LEU	2.1
1	A	136	LEU	2.1
1	A	175	LEU	2.1
1	A	31	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

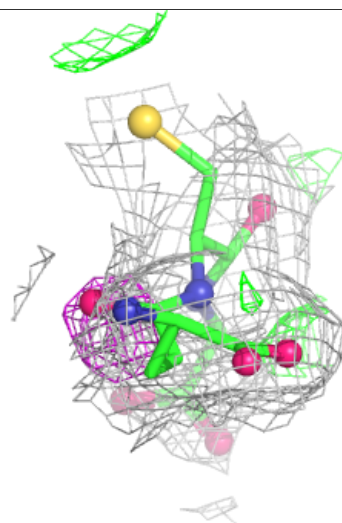
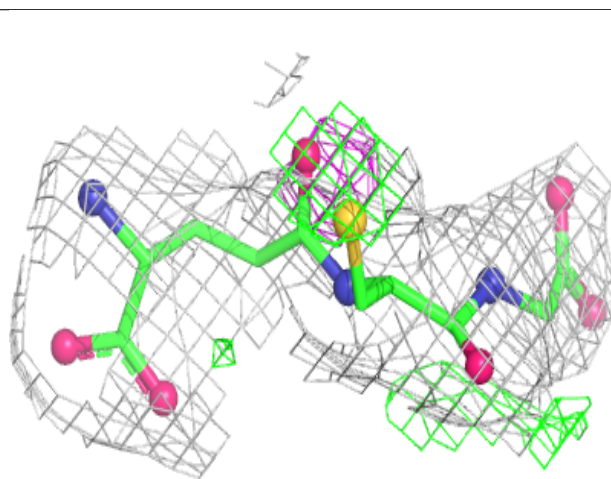
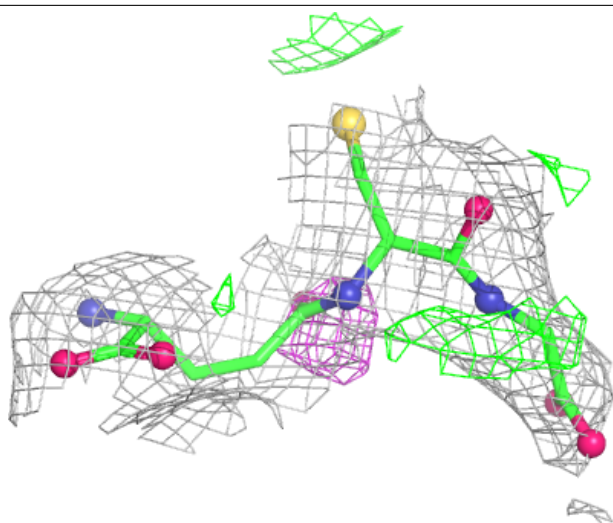
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GSH	A	300	20/20	0.70	0.14	24,24,24,24	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around GSH A 300:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.