



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:10 PM UTC

PDB ID : 1LJR / pdb_00001ljr
Title : GLUTATHIONE TRANSFERASE (HGST T2-2) FROM HUMAN
Authors : Rossjohn, J.; Mckinstry, W.J.; Oakley, A.J.; Verger, D.; Flanagan, J.; Chelvanayagam, G.; Tan, K.L.; Board, P.G.; Parker, M.W.
Deposited on : 1998-03-08
Resolution : 3.20 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

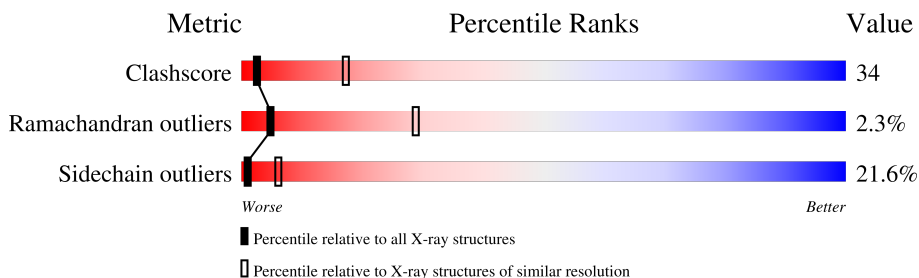
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	244	
1	B	244	

2 Entry composition [i](#)

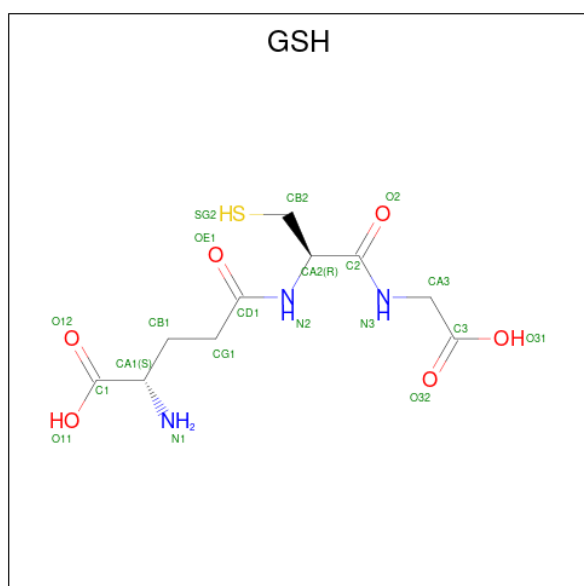
There are 2 unique types of molecules in this entry. The entry contains 3920 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 S-TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	244	Total	C	N	O	S	0	0	0
			1940	1253	331	348	8			
1	B	244	Total	C	N	O	S	0	0	0
			1940	1253	331	348	8			

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



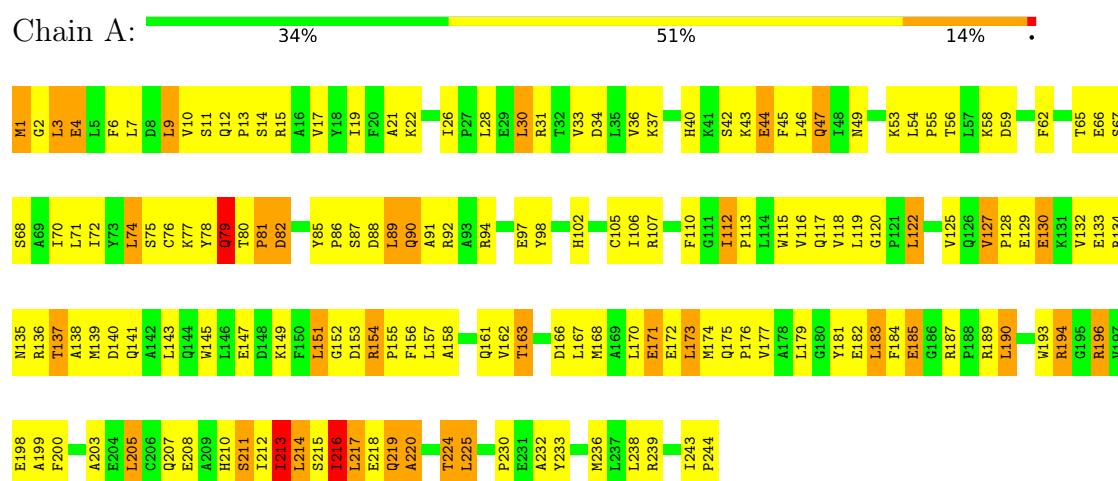
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	B	1	Total	C	N	O	S	0	0
			20	10	3	6	1		

3 Residue-property plots

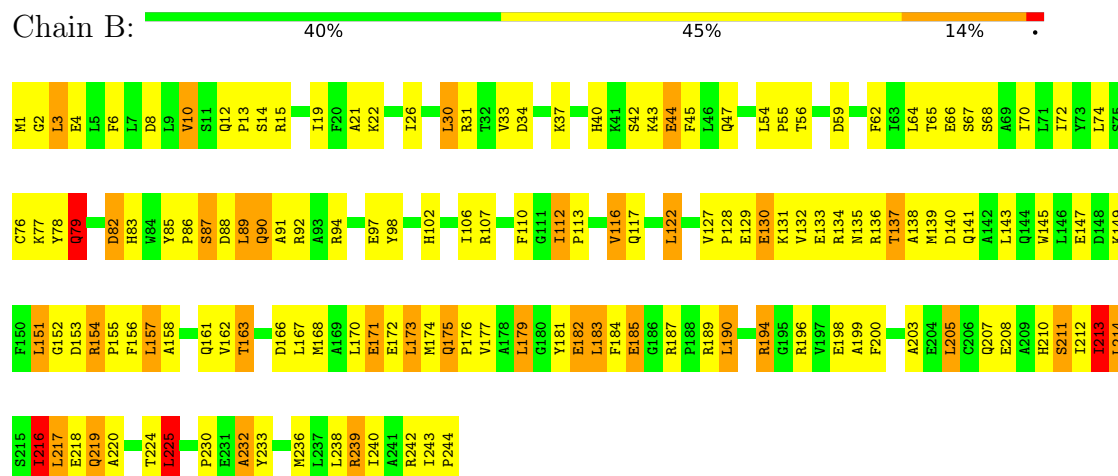
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

• Molecule 1: GLUTATHIONE S-TRANSFERASE



• Molecule 1: GLUTATHIONE S-TRANSFERASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	94.35 Å 94.35 Å 120.17 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 3.20	Depositor
% Data completeness (in resolution range)	96.0 (15.00-3.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.208 , 0.306	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3920	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	0.49	0/1985	0.99	9/2692 (0.3%)
1	B	0.52	0/1985	0.98	9/2692 (0.3%)
All	All	0.50	0/3970	0.98	18/5384 (0.3%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	216	ILE	N-CA-C	6.67	117.37	110.36
1	B	216	ILE	N-CA-C	6.54	117.23	110.36
1	A	10	VAL	N-CA-C	-6.33	104.87	111.58
1	A	213	ILE	N-CA-C	5.84	118.39	111.09
1	A	127	VAL	N-CA-C	5.53	114.37	108.95
1	A	47	GLN	N-CA-C	-5.46	105.88	112.54
1	B	213	ILE	N-CA-C	5.43	117.88	111.09
1	B	87	SER	N-CA-C	-5.28	106.79	113.18
1	B	232	ALA	N-CA-C	-5.27	105.13	112.45
1	A	232	ALA	N-CA-C	-5.17	106.12	112.90
1	B	225	LEU	CA-C-N	5.11	125.09	120.03
1	B	225	LEU	C-N-CA	5.11	125.09	120.03
1	A	219	GLN	N-CA-C	-5.09	105.72	112.34
1	B	47	GLN	N-CA-C	-5.08	106.34	112.54
1	B	10	VAL	N-CA-C	-5.05	106.22	111.58
1	A	220	ALA	N-CA-C	-5.01	105.46	111.03
1	B	219	GLN	N-CA-C	-5.01	105.82	112.34
1	A	105	CYS	N-CA-C	5.01	119.54	113.38

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1940	0	1978	146	0
1	B	1940	0	1978	133	0
2	A	20	0	15	1	0
2	B	20	0	15	0	0
All	All	3920	0	3986	271	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 34.

All (271) 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:B:158:ALA:HB3	1:B:163:THR:HG21	1.31	1.08
1:A:158:ALA:HB3	1:A:163:THR:HG21	1.29	1.05
1:A:112:ILE:HD13	1:A:243:ILE:HG12	1.43	0.98
1:B:112:ILE:HD13	1:B:243:ILE:HG12	1.46	0.96
1:A:33:VAL:HB	1:A:54:LEU:HD11	1.46	0.95
1:B:33:VAL:HB	1:B:54:LEU:HD11	1.49	0.94
1:A:118:VAL:HG12	1:A:119:LEU:HD23	1.51	0.93
1:A:214:LEU:HD23	1:A:214:LEU:H	1.33	0.92
1:B:214:LEU:H	1:B:214:LEU:HD23	1.38	0.86
1:A:214:LEU:H	1:A:214:LEU:CD2	1.89	0.86
1:A:182:GLU:HB3	1:A:185:GLU:HG2	1.57	0.84
1:B:214:LEU:H	1:B:214:LEU:CD2	1.90	0.84
1:B:182:GLU:HB3	1:B:185:GLU:HG2	1.61	0.82
1:B:210:HIS:HB3	1:B:214:LEU:HD21	1.61	0.81
1:A:158:ALA:HB3	1:A:163:THR:CG2	2.11	0.81
1:A:106:ILE:HG22	1:A:172:GLU:HG2	1.62	0.80
1:B:106:ILE:HG22	1:B:172:GLU:HG2	1.63	0.80
1:B:158:ALA:HB3	1:B:163:THR:CG2	2.14	0.77
1:A:175:GLN:HB3	1:A:176:PRO:HD3	1.68	0.76
1:B:232:ALA:O	1:B:236:MET:HG3	1.87	0.75
1:A:177:VAL:HA	1:A:181:TYR:O	1.87	0.74
1:B:12:GLN:HG2	1:B:13:PRO:HD3	1.70	0.74
1:A:210:HIS:HB3	1:A:214:LEU:HD21	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:ALA:HA	1:A:225:LEU:HD22	1.70	0.73
1:B:177:VAL:HA	1:B:181:TYR:O	1.89	0.73
1:A:147:GLU:OE2	1:A:189:ARG:HB2	1.89	0.72
1:B:175:GLN:HB3	1:B:176:PRO:HD3	1.72	0.72
1:B:145:TRP:O	1:B:149:LYS:HB3	1.90	0.71
1:A:154:ARG:HB2	1:A:155:PRO:HD2	1.71	0.71
1:B:219:GLN:HB3	1:B:224:THR:HB	1.72	0.71
1:B:196:ARG:O	1:B:199:ALA:HB3	1.90	0.71
1:A:196:ARG:O	1:A:199:ALA:HB3	1.91	0.71
1:B:130:GLU:CD	1:B:130:GLU:H	2.00	0.70
1:B:154:ARG:HB2	1:B:155:PRO:HD2	1.73	0.70
1:B:175:GLN:O	1:B:179:LEU:HD22	1.91	0.70
1:B:137:THR:O	1:B:141:GLN:HB2	1.92	0.69
1:A:219:GLN:HB3	1:A:224:THR:HB	1.75	0.68
1:A:21:ALA:HA	1:A:26:ILE:HD12	1.75	0.68
1:B:14:SER:HA	1:B:55:PRO:HG3	1.77	0.67
1:A:137:THR:O	1:A:141:GLN:HB2	1.94	0.67
1:B:88:ASP:HB3	1:B:91:ALA:HB3	1.77	0.66
1:A:12:GLN:HG2	1:A:13:PRO:HD3	1.78	0.66
1:A:173:LEU:O	1:A:176:PRO:HD2	1.96	0.66
1:B:214:LEU:CD2	1:B:214:LEU:N	2.59	0.66
1:B:6:PHE:HA	1:B:31:ARG:O	1.96	0.65
1:A:145:TRP:O	1:A:149:LYS:HB3	1.96	0.65
1:B:15:ARG:HB2	1:B:171:GLU:HG3	1.78	0.65
1:B:133:GLU:HA	1:B:136:ARG:HD2	1.78	0.65
1:A:14:SER:HA	1:A:55:PRO:HG3	1.79	0.65
1:B:21:ALA:HA	1:B:26:ILE:HD12	1.77	0.65
1:A:6:PHE:HA	1:A:31:ARG:O	1.97	0.65
1:A:90:GLN:HG3	1:B:62:PHE:CE1	2.32	0.64
1:A:15:ARG:HB2	1:A:171:GLU:HG3	1.80	0.64
1:A:213:ILE:HG12	1:A:214:LEU:N	2.13	0.64
1:A:88:ASP:HB3	1:A:91:ALA:HB3	1.80	0.64
1:A:130:GLU:H	1:A:130:GLU:CD	2.04	0.63
1:B:162:VAL:HG11	1:B:200:PHE:CE2	2.33	0.63
1:A:214:LEU:CD2	1:A:214:LEU:N	2.61	0.63
1:A:12:GLN:CD	1:A:12:GLN:H	2.07	0.63
1:B:187:ARG:HG2	1:B:187:ARG:HH11	1.65	0.62
1:B:173:LEU:O	1:B:176:PRO:HD2	1.98	0.62
1:B:161:GLN:O	1:B:163:THR:HG22	1.99	0.62
1:B:65:THR:O	1:B:66:GLU:HB2	1.98	0.62
1:A:6:PHE:CE2	1:A:31:ARG:HG2	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:SER:OG	1:A:13:PRO:HD2	2.00	0.61
1:B:6:PHE:CE2	1:B:31:ARG:HG2	2.35	0.61
1:A:133:GLU:HA	1:A:136:ARG:HD2	1.82	0.61
1:A:187:ARG:HG2	1:A:187:ARG:HH11	1.66	0.61
1:B:128:PRO:O	1:B:132:VAL:HG23	2.00	0.61
1:B:15:ARG:NH1	1:B:213:ILE:HG21	2.15	0.60
1:A:62:PHE:CE1	1:B:90:GLN:HG3	2.36	0.60
1:B:3:LEU:HD22	1:B:78:TYR:OH	2.02	0.60
1:B:163:THR:HG23	1:B:166:ASP:OD1	2.02	0.59
1:A:182:GLU:HB3	1:A:185:GLU:CG	2.32	0.58
1:A:65:THR:O	1:A:66:GLU:HB2	2.03	0.58
1:B:213:ILE:HG12	1:B:214:LEU:N	2.17	0.58
1:A:163:THR:HG23	1:A:166:ASP:OD1	2.04	0.58
1:B:122:LEU:O	1:B:122:LEU:HD22	2.03	0.58
1:A:116:VAL:HA	1:A:127:VAL:HG21	1.84	0.58
1:B:130:GLU:CD	1:B:130:GLU:N	2.61	0.57
1:B:185:GLU:OE1	1:B:185:GLU:HA	2.04	0.57
1:B:147:GLU:HG3	1:B:190:LEU:HG	1.85	0.57
1:A:220:ALA:HA	1:A:225:LEU:CD2	2.33	0.57
1:A:161:GLN:O	1:A:163:THR:HG22	2.05	0.56
1:A:34:ASP:CG	1:A:37:LYS:HB2	2.31	0.56
1:A:170:LEU:O	1:A:170:LEU:HD12	2.06	0.56
1:A:182:GLU:O	1:A:185:GLU:HG2	2.04	0.56
1:B:1:MET:O	1:B:3:LEU:N	2.37	0.56
1:B:34:ASP:CG	1:B:37:LYS:HB2	2.31	0.56
1:A:162:VAL:HG11	1:A:200:PHE:CE2	2.40	0.56
1:A:98:TYR:OH	1:A:102:HIS:HD2	1.89	0.56
1:A:1:MET:O	1:A:3:LEU:N	2.36	0.56
1:B:182:GLU:O	1:B:185:GLU:HG2	2.05	0.55
1:B:12:GLN:CD	1:B:12:GLN:H	2.13	0.55
1:A:3:LEU:HD22	1:A:78:TYR:OH	2.06	0.55
1:A:110:PHE:CD2	1:A:139:MET:HE3	2.42	0.55
1:A:205:LEU:O	1:A:208:GLU:HG2	2.07	0.55
1:A:193:TRP:HA	1:A:196:ARG:NH1	2.22	0.54
1:B:194:ARG:HH11	1:B:198:GLU:CD	2.14	0.54
1:A:97:GLU:O	1:B:66:GLU:HG3	2.07	0.54
1:A:140:ASP:HA	1:A:143:LEU:HB2	1.89	0.54
1:A:19:ILE:HG21	1:A:167:LEU:HD11	1.89	0.54
1:B:107:ARG:HG3	1:B:107:ARG:HH11	1.72	0.54
1:B:213:ILE:O	1:B:216:ILE:HG22	2.07	0.54
1:A:174:MET:HG3	1:A:210:HIS:NE2	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:PRO:HA	1:A:233:TYR:HB3	1.90	0.54
1:B:19:ILE:HG21	1:B:167:LEU:HD11	1.89	0.54
1:A:185:GLU:OE1	1:A:185:GLU:HA	2.08	0.54
1:B:205:LEU:O	1:B:208:GLU:HG2	2.08	0.54
1:A:76:CYS:HB3	1:B:89:LEU:HD21	1.90	0.54
1:B:155:PRO:HA	1:B:196:ARG:NH1	2.23	0.54
1:B:211:SER:O	1:B:214:LEU:HD23	2.09	0.53
1:A:89:LEU:HD21	1:B:76:CYS:HB3	1.91	0.52
1:B:182:GLU:HB3	1:B:185:GLU:CG	2.37	0.52
1:A:153:ASP:OD1	1:A:154:ARG:N	2.42	0.52
1:A:9:LEU:HD22	1:A:15:ARG:HG2	1.92	0.52
1:A:1:MET:HG2	1:A:3:LEU:HB2	1.92	0.52
1:A:118:VAL:HG12	1:A:119:LEU:N	2.23	0.52
1:B:98:TYR:CE2	1:B:102:HIS:HB2	2.45	0.52
1:A:40:HIS:HA	1:A:45:PHE:CD2	2.46	0.51
1:B:12:GLN:HG2	1:B:13:PRO:CD	2.38	0.51
1:B:110:PHE:CD2	1:B:139:MET:HE3	2.45	0.51
1:A:184:PHE:CE1	1:A:190:LEU:HD13	2.45	0.51
1:A:218:GLU:C	1:A:220:ALA:N	2.66	0.51
1:B:98:TYR:OH	1:B:102:HIS:HD2	1.93	0.51
1:A:6:PHE:CD1	1:A:6:PHE:N	2.78	0.51
1:A:213:ILE:O	1:A:216:ILE:HG22	2.11	0.51
1:A:135:ASN:O	1:A:138:ALA:HB3	2.10	0.51
1:A:136:ARG:HG2	1:A:137:THR:N	2.25	0.51
1:A:107:ARG:HH11	1:A:107:ARG:HG3	1.75	0.51
1:B:236:MET:O	1:B:240:ILE:HG13	2.10	0.51
1:A:210:HIS:O	1:A:212:ILE:N	2.43	0.51
1:B:151:LEU:O	1:B:152:GLY:C	2.53	0.50
1:B:230:PRO:HA	1:B:233:TYR:HB3	1.93	0.50
1:A:130:GLU:CD	1:A:130:GLU:N	2.62	0.50
1:A:217:LEU:O	1:A:220:ALA:HB3	2.11	0.50
1:B:140:ASP:HA	1:B:143:LEU:HB2	1.94	0.50
1:B:143:LEU:HD11	1:B:183:LEU:HB3	1.92	0.50
1:A:151:LEU:O	1:A:152:GLY:C	2.53	0.49
1:B:155:PRO:HB2	1:B:156:PHE:CE1	2.48	0.49
1:A:175:GLN:HB3	1:A:176:PRO:CD	2.41	0.49
1:B:112:ILE:CD1	1:B:243:ILE:HG12	2.31	0.49
1:B:78:TYR:C	1:B:79:GLN:HG3	2.38	0.49
1:A:147:GLU:OE1	1:A:187:ARG:HG2	2.13	0.49
1:B:210:HIS:O	1:B:212:ILE:N	2.45	0.49
1:B:218:GLU:C	1:B:220:ALA:N	2.67	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:HIS:C	1:B:214:LEU:HD21	2.38	0.48
1:B:211:SER:OG	1:B:212:ILE:N	2.46	0.48
1:B:220:ALA:HA	1:B:225:LEU:HD22	1.95	0.48
1:A:119:LEU:HD23	1:A:119:LEU:N	2.27	0.48
1:B:155:PRO:HG2	1:B:156:PHE:HD1	1.77	0.48
1:A:210:HIS:O	1:A:211:SER:C	2.56	0.48
1:A:211:SER:O	1:A:214:LEU:HD23	2.14	0.48
1:A:210:HIS:HB3	1:A:214:LEU:CD2	2.40	0.48
1:B:1:MET:HE1	1:B:59:ASP:OD2	2.14	0.48
1:B:210:HIS:O	1:B:211:SER:C	2.56	0.48
1:B:147:GLU:OE2	1:B:189:ARG:HB2	2.13	0.48
1:B:184:PHE:O	1:B:185:GLU:C	2.57	0.48
1:A:90:GLN:HG3	1:B:62:PHE:CZ	2.49	0.47
1:A:143:LEU:HD22	1:A:190:LEU:HD11	1.96	0.47
1:A:184:PHE:O	1:A:185:GLU:C	2.56	0.47
1:B:6:PHE:CD1	1:B:6:PHE:N	2.82	0.47
1:A:70:ILE:HG23	1:A:74:LEU:HD13	1.95	0.47
1:A:210:HIS:C	1:A:214:LEU:HD21	2.39	0.47
1:B:216:ILE:CG2	1:B:217:LEU:N	2.77	0.47
1:A:155:PRO:HG2	1:A:156:PHE:HD1	1.79	0.47
1:A:116:VAL:HA	1:A:127:VAL:CG2	2.44	0.47
1:B:143:LEU:HD22	1:B:190:LEU:HD11	1.96	0.47
1:B:216:ILE:HG23	1:B:217:LEU:N	2.28	0.47
1:A:122:LEU:HD22	1:A:122:LEU:O	2.15	0.47
1:B:175:GLN:HB3	1:B:176:PRO:CD	2.44	0.47
1:A:113:PRO:HB3	1:A:181:TYR:CE1	2.49	0.47
1:B:42:SER:O	1:B:43:LYS:C	2.58	0.47
1:B:83:HIS:CD2	1:B:162:VAL:HG13	2.50	0.47
1:B:174:MET:HG3	1:B:210:HIS:NE2	2.29	0.47
1:A:78:TYR:C	1:A:79:GLN:HG3	2.40	0.46
1:B:134:ARG:NH1	1:B:244:PRO:O	2.48	0.46
1:B:239:ARG:O	1:B:242:ARG:HG2	2.15	0.46
1:A:15:ARG:NH1	1:A:213:ILE:HG21	2.30	0.46
1:B:116:VAL:HG22	1:B:127:VAL:HB	1.98	0.46
1:A:187:ARG:HB3	1:A:190:LEU:HB2	1.96	0.46
1:B:136:ARG:HG2	1:B:137:THR:N	2.30	0.46
1:A:211:SER:OG	1:A:212:ILE:N	2.49	0.46
1:A:187:ARG:HG2	1:A:187:ARG:NH1	2.29	0.46
1:B:40:HIS:HA	1:B:45:PHE:CD2	2.51	0.46
1:B:154:ARG:HG3	1:B:156:PHE:O	2.16	0.46
1:B:225:LEU:HD12	1:B:225:LEU:HA	1.78	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:SER:O	1:A:218:GLU:HB3	2.16	0.46
1:A:216:ILE:HG23	1:A:217:LEU:N	2.30	0.46
1:B:44:GLU:H	1:B:44:GLU:HG2	1.42	0.45
1:B:131:LYS:HD3	1:B:243:ILE:HG22	1.99	0.45
1:A:42:SER:O	1:A:43:LYS:C	2.59	0.45
1:B:187:ARG:HG2	1:B:187:ARG:NH1	2.29	0.45
1:A:44:GLU:O	1:A:47:GLN:HB2	2.16	0.45
1:B:135:ASN:O	1:B:138:ALA:HB3	2.17	0.45
1:A:72:ILE:HG23	1:A:85:TYR:CE2	2.51	0.45
1:A:78:TYR:O	1:A:79:GLN:C	2.58	0.45
1:B:147:GLU:OE1	1:B:187:ARG:HG2	2.16	0.45
1:A:44:GLU:H	1:A:44:GLU:HG2	1.45	0.45
1:B:6:PHE:HA	1:B:31:ARG:HB3	1.98	0.45
1:A:6:PHE:HA	1:A:31:ARG:HB3	1.98	0.45
1:A:143:LEU:CD1	1:A:183:LEU:HB3	2.47	0.45
1:B:78:TYR:O	1:B:79:GLN:C	2.59	0.45
1:B:210:HIS:HB3	1:B:214:LEU:CD2	2.41	0.45
1:A:155:PRO:HB2	1:A:156:PHE:CE1	2.52	0.45
1:A:98:TYR:CE2	1:A:102:HIS:HB2	2.52	0.45
1:A:115:TRP:HZ3	1:A:236:MET:HG2	1.82	0.44
1:B:72:ILE:HG23	1:B:85:TYR:CE2	2.53	0.44
1:A:134:ARG:NH1	1:A:244:PRO:O	2.50	0.44
1:B:30:LEU:N	1:B:30:LEU:HD23	2.32	0.44
1:A:220:ALA:CA	1:A:225:LEU:HD22	2.45	0.44
1:B:3:LEU:HD13	1:B:59:ASP:OD2	2.17	0.44
1:B:86:PRO:O	1:B:92:ARG:HB2	2.18	0.44
1:B:170:LEU:HD12	1:B:170:LEU:O	2.17	0.44
1:A:162:VAL:HG11	1:A:200:PHE:CZ	2.52	0.44
1:A:82:ASP:OD2	1:A:86:PRO:HA	2.17	0.44
1:A:12:GLN:N	1:A:13:PRO:HD2	2.33	0.44
1:A:194:ARG:HH11	1:A:198:GLU:CD	2.26	0.44
1:A:143:LEU:HD11	1:A:183:LEU:HB3	2.00	0.43
1:A:179:LEU:HD12	1:A:217:LEU:HD23	2.00	0.43
1:A:216:ILE:CG2	1:A:217:LEU:N	2.80	0.43
1:B:113:PRO:HB3	1:B:181:TYR:CE1	2.53	0.43
1:A:128:PRO:O	1:A:132:VAL:HG23	2.17	0.43
1:A:106:ILE:CG2	1:A:172:GLU:HG2	2.39	0.43
1:A:218:GLU:C	1:A:220:ALA:H	2.27	0.43
1:B:162:VAL:HG11	1:B:200:PHE:CZ	2.53	0.43
1:B:3:LEU:HD12	1:B:59:ASP:HB2	2.01	0.43
1:A:154:ARG:HB2	1:A:155:PRO:CD	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:SER:N	1:A:214:LEU:HD21	2.34	0.43
1:A:90:GLN:CG	1:B:62:PHE:CE1	3.02	0.43
1:A:7:LEU:HA	1:A:54:LEU:HD13	2.01	0.42
1:B:89:LEU:HD23	1:B:89:LEU:HA	1.78	0.42
1:A:137:THR:O	1:A:140:ASP:OD1	2.38	0.42
1:A:193:TRP:HA	1:A:196:ARG:HH11	1.83	0.42
1:A:4:GLU:OE2	1:A:58:LYS:HE2	2.20	0.42
1:A:53:LYS:HG2	2:A:245:GSH:O32	2.18	0.42
1:A:167:LEU:O	1:A:171:GLU:HB2	2.20	0.42
1:B:106:ILE:CG2	1:B:172:GLU:HG2	2.40	0.42
1:B:66:GLU:C	1:B:70:ILE:HD12	2.44	0.42
1:B:155:PRO:O	1:B:196:ARG:NH1	2.52	0.42
1:B:167:LEU:O	1:B:171:GLU:HB2	2.19	0.42
1:A:151:LEU:HD23	1:A:151:LEU:HA	1.84	0.42
1:A:203:ALA:O	1:A:207:GLN:HB2	2.19	0.42
1:A:80:THR:O	1:A:81:PRO:C	2.63	0.42
1:A:46:LEU:HA	1:A:49:ASN:O	2.20	0.42
1:A:133:GLU:O	1:A:134:ARG:C	2.62	0.42
1:B:167:LEU:HD12	1:B:167:LEU:HA	1.77	0.42
1:B:190:LEU:HD23	1:B:190:LEU:HA	1.86	0.42
1:A:40:HIS:O	1:A:45:PHE:HD2	2.03	0.41
1:B:3:LEU:CD1	1:B:59:ASP:HB2	2.50	0.41
1:B:157:LEU:H	1:B:157:LEU:HG	1.66	0.41
1:B:217:LEU:O	1:B:220:ALA:HB3	2.20	0.41
1:B:12:GLN:HG3	1:B:172:GLU:CD	2.45	0.41
1:A:21:ALA:HB1	1:A:28:LEU:HD12	2.03	0.41
1:A:140:ASP:OD1	1:A:141:GLN:N	2.53	0.41
1:A:155:PRO:HG2	1:A:156:PHE:CD1	2.55	0.41
1:B:155:PRO:HG2	1:B:156:PHE:CD1	2.56	0.41
1:B:143:LEU:CD1	1:B:183:LEU:HB3	2.51	0.41
1:B:154:ARG:HB2	1:B:155:PRO:CD	2.46	0.41
1:A:3:LEU:HD13	1:A:59:ASP:OD2	2.20	0.41
1:A:17:VAL:HG22	1:A:71:LEU:HG	2.02	0.41
1:A:30:LEU:N	1:A:30:LEU:HD23	2.36	0.41
1:A:86:PRO:O	1:A:92:ARG:HB2	2.20	0.41
1:A:120:GLY:C	1:A:125:VAL:O	2.64	0.41
1:B:1:MET:HE1	1:B:77:LYS:HE2	2.02	0.41
1:B:82:ASP:O	1:B:83:HIS:C	2.63	0.41
1:A:66:GLU:HG3	1:B:97:GLU:O	2.21	0.41
1:B:8:ASP:OD1	1:B:10:VAL:HG23	2.21	0.41
1:A:112:ILE:N	1:A:113:PRO:CD	2.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:MET:HE1	1:A:176:PRO:CG	2.50	0.40
1:B:203:ALA:O	1:B:207:GLN:HB2	2.22	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	242/244 (99%)	210 (87%)	26 (11%)	6 (2%)	4	27
1	B	242/244 (99%)	207 (86%)	30 (12%)	5 (2%)	5	31
All	All	484/488 (99%)	417 (86%)	56 (12%)	11 (2%)	5	29

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	185	GLU
1	A	211	SER
1	B	185	GLU
1	B	211	SER
1	A	3	LEU
1	B	3	LEU
1	A	79	GLN
1	B	79	GLN
1	A	2	GLY
1	B	2	GLY
1	A	81	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	206/206 (100%)	161 (78%)	45 (22%)	1	6
1	B	206/206 (100%)	162 (79%)	44 (21%)	1	6
All	All	412/412 (100%)	323 (78%)	89 (22%)	1	6

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	4	GLU
1	A	9	LEU
1	A	22	LYS
1	A	30	LEU
1	A	36	VAL
1	A	44	GLU
1	A	56	THR
1	A	67	SER
1	A	68	SER
1	A	74	LEU
1	A	75	SER
1	A	77	LYS
1	A	79	GLN
1	A	82	ASP
1	A	87	SER
1	A	89	LEU
1	A	90	GLN
1	A	94	ARG
1	A	112	ILE
1	A	117	GLN
1	A	122	LEU
1	A	129	GLU
1	A	130	GLU
1	A	137	THR
1	A	151	LEU
1	A	154	ARG

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Mol	Chain	Res	Type
1	A	157	LEU
1	A	163	THR
1	A	168	MET
1	A	171	GLU
1	A	173	LEU
1	A	183	LEU
1	A	190	LEU
1	A	194	ARG
1	A	196	ARG
1	A	205	LEU
1	A	213	ILE
1	A	214	LEU
1	A	216	ILE
1	A	217	LEU
1	A	224	THR
1	A	225	LEU
1	A	238	LEU
1	A	239	ARG
1	B	4	GLU
1	B	22	LYS
1	B	30	LEU
1	B	44	GLU
1	B	56	THR
1	B	64	LEU
1	B	67	SER
1	B	68	SER
1	B	74	LEU
1	B	79	GLN
1	B	82	ASP
1	B	87	SER
1	B	89	LEU
1	B	90	GLN
1	B	94	ARG
1	B	112	ILE
1	B	116	VAL
1	B	117	GLN
1	B	122	LEU
1	B	129	GLU
1	B	130	GLU
1	B	137	THR
1	B	151	LEU
1	B	153	ASP

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Mol	Chain	Res	Type
1	B	154	ARG
1	B	157	LEU
1	B	163	THR
1	B	168	MET
1	B	171	GLU
1	B	173	LEU
1	B	175	GLN
1	B	179	LEU
1	B	182	GLU
1	B	183	LEU
1	B	190	LEU
1	B	194	ARG
1	B	205	LEU
1	B	213	ILE
1	B	214	LEU
1	B	216	ILE
1	B	217	LEU
1	B	225	LEU
1	B	238	LEU
1	B	239	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	79	GLN
1	A	83	HIS
1	A	102	HIS
1	A	141	GLN
1	A	144	GLN
1	B	24	ASN
1	B	79	GLN
1	B	83	HIS
1	B	102	HIS
1	B	117	GLN
1	B	141	GLN
1	B	144	GLN

5.3.3 RNA ⓘ

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](#)

2 ligands are 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	B	245	-	18,19,19	0.87	0	21,24,24	0.71	0
2	GSH	A	245	-	18,19,19	0.82	0	21,24,24	0.66	0

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	B	245	-	-	3/24/24/24	-
2	GSH	A	245	-	-	4/24/24/24	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	245	GSH	C2-CA2-CB2-SG2
2	B	245	GSH	C2-CA2-CB2-SG2

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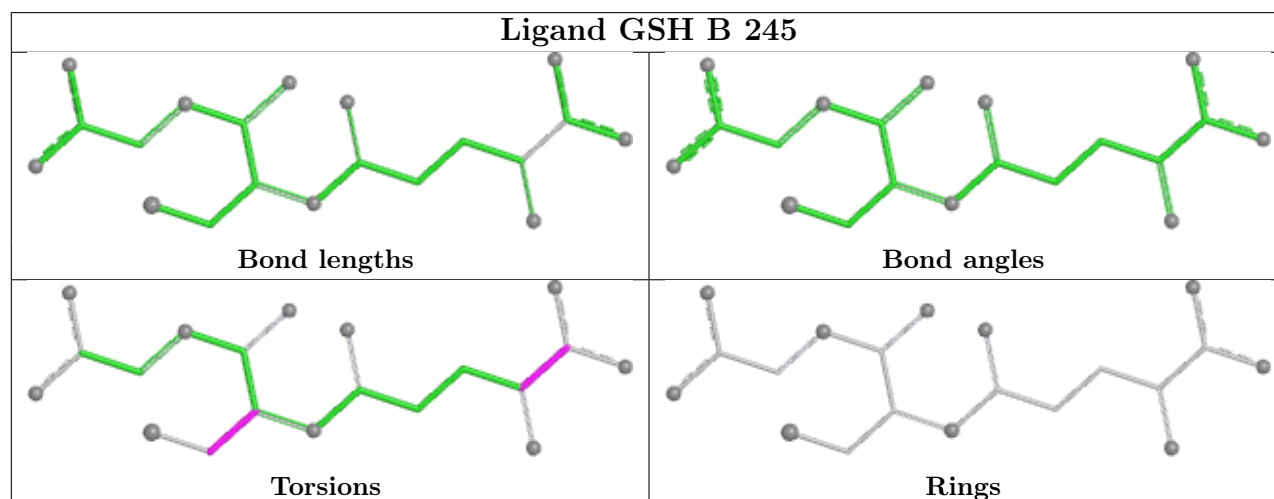
Mol	Chain	Res	Type	Atoms
2	B	245	GSH	N2-CA2-CB2-SG2
2	A	245	GSH	N2-CA2-CB2-SG2
2	B	245	GSH	O11-C1-CA1-N1
2	A	245	GSH	O11-C1-CA1-N1
2	A	245	GSH	O12-C1-CA1-N1

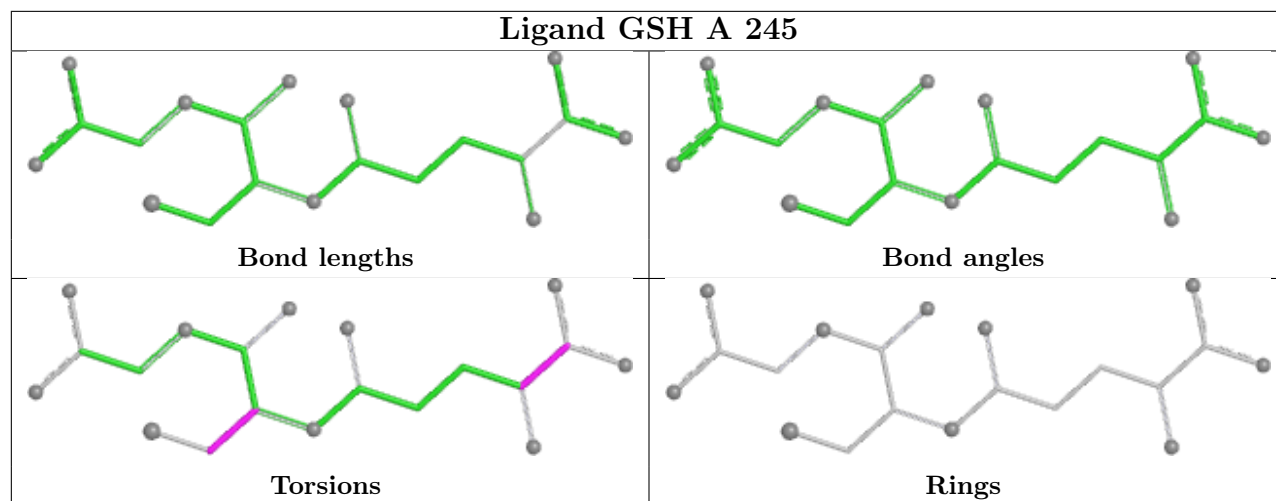
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	245	GSH	1	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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.