



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:28 PM UTC

PDB ID : 8DL5 / pdb_00008dl5
Title : Crystal structure of PLP-dependent Mannich cyclase LolT
Authors : Gao, J.; Hai, Y.
Deposited on : 2022-07-06
Resolution : 2.10 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

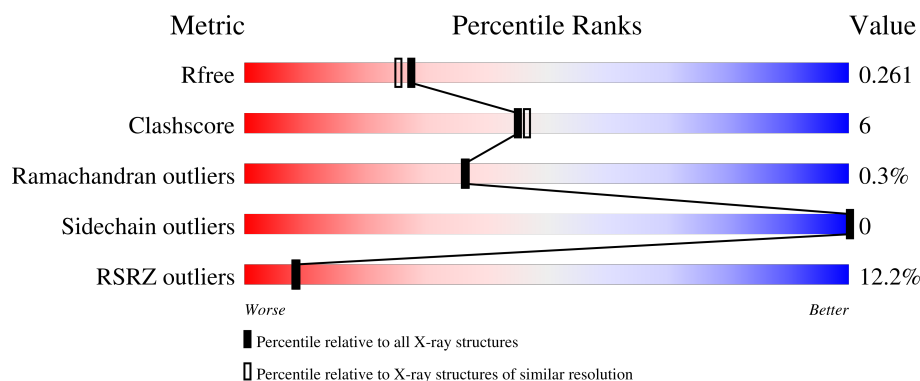
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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))
R_{free}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	496	14% 67% 15% 19%
1	B	496	6% 71% 11% 18%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6655 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 Aminotransferase, class V/Cysteine desulfurase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	P	S	0	0	0
			3212	2057	544	588	1	22			
1	B	408	Total	C	N	O	P	S	0	0	0
			3253	2082	551	597	1	22			

There are 44 discrepancies between the modelled and reference sequences:

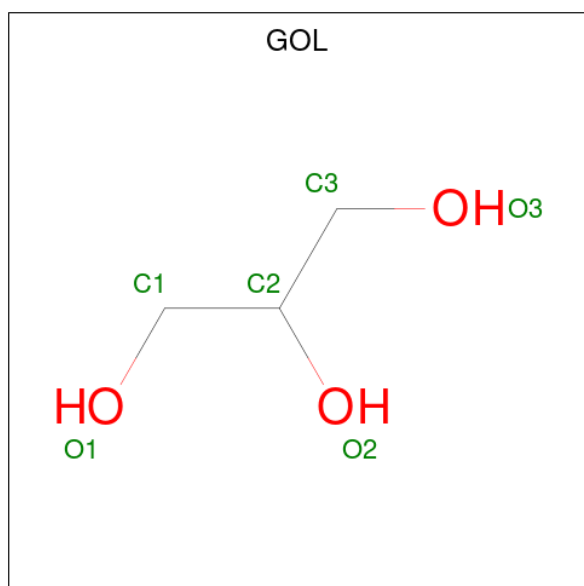
Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP A0A0A2J6G6
A	-20	GLY	-	expression tag	UNP A0A0A2J6G6
A	-19	SER	-	expression tag	UNP A0A0A2J6G6
A	-18	SER	-	expression tag	UNP A0A0A2J6G6
A	-17	HIS	-	expression tag	UNP A0A0A2J6G6
A	-16	HIS	-	expression tag	UNP A0A0A2J6G6
A	-15	HIS	-	expression tag	UNP A0A0A2J6G6
A	-14	HIS	-	expression tag	UNP A0A0A2J6G6
A	-13	HIS	-	expression tag	UNP A0A0A2J6G6
A	-12	HIS	-	expression tag	UNP A0A0A2J6G6
A	-11	SER	-	expression tag	UNP A0A0A2J6G6
A	-10	SER	-	expression tag	UNP A0A0A2J6G6
A	-9	GLY	-	expression tag	UNP A0A0A2J6G6
A	-8	LEU	-	expression tag	UNP A0A0A2J6G6
A	-7	VAL	-	expression tag	UNP A0A0A2J6G6
A	-6	PRO	-	expression tag	UNP A0A0A2J6G6
A	-5	ARG	-	expression tag	UNP A0A0A2J6G6
A	-4	GLY	-	expression tag	UNP A0A0A2J6G6
A	-3	SER	-	expression tag	UNP A0A0A2J6G6
A	-2	HIS	-	expression tag	UNP A0A0A2J6G6
A	-1	MET	-	expression tag	UNP A0A0A2J6G6
A	0	HIS	-	expression tag	UNP A0A0A2J6G6
B	-21	MET	-	expression tag	UNP A0A0A2J6G6
B	-20	GLY	-	expression tag	UNP A0A0A2J6G6
B	-19	SER	-	expression tag	UNP A0A0A2J6G6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	SER	-	expression tag	UNP A0A0A2J6G6
B	-17	HIS	-	expression tag	UNP A0A0A2J6G6
B	-16	HIS	-	expression tag	UNP A0A0A2J6G6
B	-15	HIS	-	expression tag	UNP A0A0A2J6G6
B	-14	HIS	-	expression tag	UNP A0A0A2J6G6
B	-13	HIS	-	expression tag	UNP A0A0A2J6G6
B	-12	HIS	-	expression tag	UNP A0A0A2J6G6
B	-11	SER	-	expression tag	UNP A0A0A2J6G6
B	-10	SER	-	expression tag	UNP A0A0A2J6G6
B	-9	GLY	-	expression tag	UNP A0A0A2J6G6
B	-8	LEU	-	expression tag	UNP A0A0A2J6G6
B	-7	VAL	-	expression tag	UNP A0A0A2J6G6
B	-6	PRO	-	expression tag	UNP A0A0A2J6G6
B	-5	ARG	-	expression tag	UNP A0A0A2J6G6
B	-4	GLY	-	expression tag	UNP A0A0A2J6G6
B	-3	SER	-	expression tag	UNP A0A0A2J6G6
B	-2	HIS	-	expression tag	UNP A0A0A2J6G6
B	-1	MET	-	expression tag	UNP A0A0A2J6G6
B	0	HIS	-	expression tag	UNP A0A0A2J6G6

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	74	Total 74	O 74	0	0
3	B	110	Total 110	O 110	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	77.84Å 77.84Å 298.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.66 – 2.10 37.66 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (37.66-2.10) 94.4 (37.66-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.233 , 0.261 0.234 , 0.261	Depositor DCC
R_{free} test set	3144 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 21.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.048 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6655	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 35.91 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.4240e-04.*

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: GOL, LLP

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	0.20	0/3266	0.44	0/4441
1	B	0.21	0/3307	0.50	3/4494 (0.1%)
All	All	0.20	0/6573	0.47	3/8935 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	459	GLU	N-CA-C	-5.76	105.35	112.90
1	B	459	GLU	CA-C-N	5.52	128.26	120.42
1	B	459	GLU	C-N-CA	5.52	128.26	120.42

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	236	LLP	Mainchain

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	3212	0	3166	46	0
1	B	3253	0	3208	36	0
2	B	6	0	8	0	0
3	A	74	0	0	2	0
3	B	110	0	0	1	0
All	All	6655	0	6382	76	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 6.

All (76) 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:415:SER:OG	1:B:282:GLN:HG2	1.91	0.71
1:A:268:ILE:HD11	1:A:294:ASP:HB3	1.74	0.70
1:A:370:GLU:HG3	1:A:372:ARG:H	1.56	0.68
1:A:240:VAL:HG11	1:A:313:ILE:HD13	1.76	0.66
1:A:293:THR:HG23	1:A:296:GLU:H	1.60	0.66
1:A:137:GLU:OE2	1:B:258:ARG:NH2	2.26	0.66
1:B:293:THR:HG22	1:B:296:GLU:H	1.63	0.62
1:A:456:VAL:O	1:A:460:ILE:HG13	2.01	0.61
1:B:65:TYR:O	1:B:69:VAL:HG22	2.01	0.60
1:B:62:GLN:HB3	1:B:65:TYR:HB3	1.85	0.58
1:B:75:ILE:HD11	1:B:307:ASN:HB3	1.86	0.58
1:B:345:VAL:HG12	1:B:379:VAL:HG11	1.85	0.58
1:A:114:LYS:HB3	1:A:140:PRO:HB2	1.87	0.57
1:A:189:ILE:HG23	1:A:354:LEU:HD23	1.87	0.57
1:A:183:ILE:HB	1:A:213:HIS:HB2	1.86	0.57
1:A:71:GLN:NE2	3:A:503:HOH:O	2.39	0.55
1:A:113:ASP:OD1	1:A:114:LYS:N	2.42	0.53
1:A:230:PHE:HD1	1:A:232:MET:HE2	1.73	0.53
1:A:53:ARG:NH2	1:B:60:GLU:OE2	2.40	0.53
1:A:158:VAL:O	1:A:162:GLU:HG3	2.09	0.53
1:A:293:THR:O	1:A:297:THR:HG23	2.08	0.53
1:A:81:GLU:HB3	1:A:317:LEU:HD12	1.88	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:ILE:HG23	1:A:173:LEU:HB2	1.91	0.53
1:B:107:LEU:HB3	1:B:135:MET:HE1	1.91	0.53
1:B:293:THR:HG21	3:B:617:HOH:O	2.08	0.53
1:A:103:ILE:HD13	1:A:128:ILE:HD11	1.91	0.52
1:A:137:GLU:CD	1:B:258:ARG:HH22	2.16	0.51
1:B:183:ILE:HB	1:B:213:HIS:HB2	1.93	0.51
1:A:428:LEU:HG	1:B:282:GLN:NE2	2.26	0.51
1:B:217:GLN:OE1	1:B:374:CYS:HB2	2.12	0.50
1:B:230:PHE:HD1	1:B:232:MET:HE2	1.77	0.50
1:A:422:HIS:HB3	1:A:452:TRP:CH2	2.47	0.49
1:A:77:GLU:O	1:A:81:GLU:HG3	2.13	0.49
1:A:230:PHE:CD1	1:A:232:MET:HE2	2.48	0.49
1:B:56:LEU:O	1:B:60:GLU:HG2	2.13	0.49
1:B:259:SER:HB2	1:B:267:PHE:HD1	1.78	0.49
1:A:167:GLN:O	1:A:171:GLU:HG3	2.14	0.48
1:B:344:ARG:HE	1:B:451:GLU:CD	2.22	0.47
1:A:75:ILE:HD11	1:A:307:ASN:HB3	1.95	0.47
1:A:189:ILE:HG21	1:A:371:MET:HE3	1.96	0.47
1:A:104:TYR:HD2	1:A:261:ILE:HG21	1.80	0.47
1:A:356:GLU:H	1:A:371:MET:HE2	1.80	0.46
1:B:86:LEU:HD11	1:B:232:MET:HE1	1.97	0.46
1:A:175:PRO:O	1:A:205:VAL:HG22	2.16	0.46
1:B:234:CYS:N	1:B:244:CYS:O	2.48	0.46
1:A:327:GLU:HB3	1:A:331:TYR:CZ	2.50	0.46
1:A:80:ARG:HA	1:A:90:VAL:HG22	1.97	0.45
1:B:404:ASP:O	1:B:407:VAL:HG22	2.16	0.45
1:B:32:GLU:CD	1:B:423:LYS:HD3	2.42	0.45
1:A:43:SER:O	1:A:239:PHE:HA	2.17	0.45
1:A:318:LYS:HE2	1:A:322:GLU:OE1	2.16	0.45
1:A:356:GLU:N	1:A:371:MET:HE2	2.31	0.45
1:B:230:PHE:CD1	1:B:232:MET:HE2	2.52	0.45
1:B:411:VAL:HG13	1:B:428:LEU:HD13	1.98	0.44
1:B:417:THR:O	1:B:421:THR:HB	2.16	0.44
1:A:339:LYS:HA	1:A:372:ARG:NH2	2.32	0.44
1:A:137:GLU:HB3	1:B:270:ARG:NH2	2.32	0.44
1:B:106:VAL:O	1:B:110:ILE:HG12	2.18	0.44
1:B:339:LYS:HA	1:B:372:ARG:NH2	2.32	0.44
1:A:182:THR:OG1	1:A:211:GLY:HA2	2.18	0.44
1:B:268:ILE:HD12	1:B:298:ILE:HD11	2.00	0.44
1:B:278:LEU:HD12	1:B:281:LYS:HG3	2.00	0.43
1:B:49:ARG:O	1:B:53:ARG:HG2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:ARG:HH21	1:A:269:PRO:HA	1.84	0.43
1:A:106:VAL:O	1:A:110:ILE:HG12	2.19	0.43
1:B:381:LEU:C	1:B:383:LEU:H	2.27	0.43
1:A:258:ARG:CZ	1:A:270:ARG:HG3	2.49	0.42
1:A:293:THR:HG21	3:A:516:HOH:O	2.20	0.42
1:B:65:TYR:CZ	1:B:69:VAL:HG21	2.54	0.42
1:B:293:THR:HG22	1:B:295:PHE:N	2.34	0.41
1:A:191:MET:HE3	1:A:192:PRO:HD2	2.03	0.41
1:A:59:LEU:HD22	1:A:308:MET:HE1	2.03	0.41
1:A:89:ARG:HB2	1:A:92:GLU:HG3	2.01	0.41
1:B:425:TRP:O	1:B:426:LEU:HD23	2.21	0.41
1:A:250:PRO:O	1:A:254:GLN:HG3	2.22	0.40
1:B:293:THR:HB	1:B:296:GLU:OE1	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	392/496 (79%)	380 (97%)	11 (3%)	1 (0%)	36	36
1	B	397/496 (80%)	381 (96%)	15 (4%)	1 (0%)	36	36
All	All	789/992 (80%)	761 (96%)	26 (3%)	2 (0%)	36	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	214	SER
1	B	214	SER

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	347/425 (82%)	347 (100%)	0	100	100
1	B	352/425 (83%)	352 (100%)	0	100	100
All	All	699/850 (82%)	699 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	71	GLN
1	A	87	HIS
1	B	51	GLN
1	B	108	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	LLP	B	236	1	23,24,25	2.52	6 (26%)	25,32,34	1.51	4 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	LLP	A	236	1	23,24,25	2.50	7 (30%)	25,32,34	1.57	4 (16%)

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
1	LLP	B	236	1	-	7/16/17/19	0/1/1/1
1	LLP	A	236	1	-	6/16/17/19	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	236	LLP	C4-C4'	7.48	1.62	1.46
1	A	236	LLP	C4-C4'	7.37	1.62	1.46
1	B	236	LLP	C4'-NZ	5.23	1.44	1.27
1	A	236	LLP	C4'-NZ	5.21	1.44	1.27
1	B	236	LLP	C2'-C2	3.67	1.56	1.50
1	A	236	LLP	C2'-C2	3.59	1.56	1.50
1	A	236	LLP	C4-C5	-3.25	1.37	1.42
1	B	236	LLP	C4-C5	-3.02	1.37	1.42
1	A	236	LLP	C6-N1	2.93	1.40	1.34
1	B	236	LLP	C6-N1	2.86	1.40	1.34
1	B	236	LLP	C5'-C5	2.20	1.56	1.50
1	A	236	LLP	C5'-C5	2.15	1.56	1.50
1	A	236	LLP	C4-C3	-2.09	1.37	1.41

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	236	LLP	C4-C4'-NZ	-4.90	101.41	124.04
1	B	236	LLP	C4-C4'-NZ	-4.04	105.41	124.04
1	B	236	LLP	OP4-C5'-C5	2.98	114.95	109.36
1	A	236	LLP	C3-C4-C4'	-2.49	115.91	120.40
1	B	236	LLP	C5'-C5-C6	-2.42	115.42	119.36
1	B	236	LLP	C3-C4-C4'	-2.22	116.40	120.40
1	A	236	LLP	OP4-C5'-C5	2.15	113.39	109.36
1	A	236	LLP	C5-C4-C4'	2.07	124.66	121.47

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	236	LLP	C4-C5-C5'-OP4
1	B	236	LLP	C6-C5-C5'-OP4
1	A	236	LLP	C5-C4-C4'-NZ
1	B	236	LLP	C4-C4'-NZ-CE
1	A	236	LLP	C3-C4-C4'-NZ
1	B	236	LLP	C3-C4-C4'-NZ
1	A	236	LLP	C6-C5-C5'-OP4
1	A	236	LLP	CD-CE-NZ-C4'
1	A	236	LLP	C4-C5-C5'-OP4
1	B	236	LLP	CD-CE-NZ-C4'
1	B	236	LLP	C5-C4-C4'-NZ
1	A	236	LLP	C-CA-CB-CG
1	B	236	LLP	C-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

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	GOL	B	501	-	5,5,5	0.87	0	5,5,5	1.13	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	GOL	B	501	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

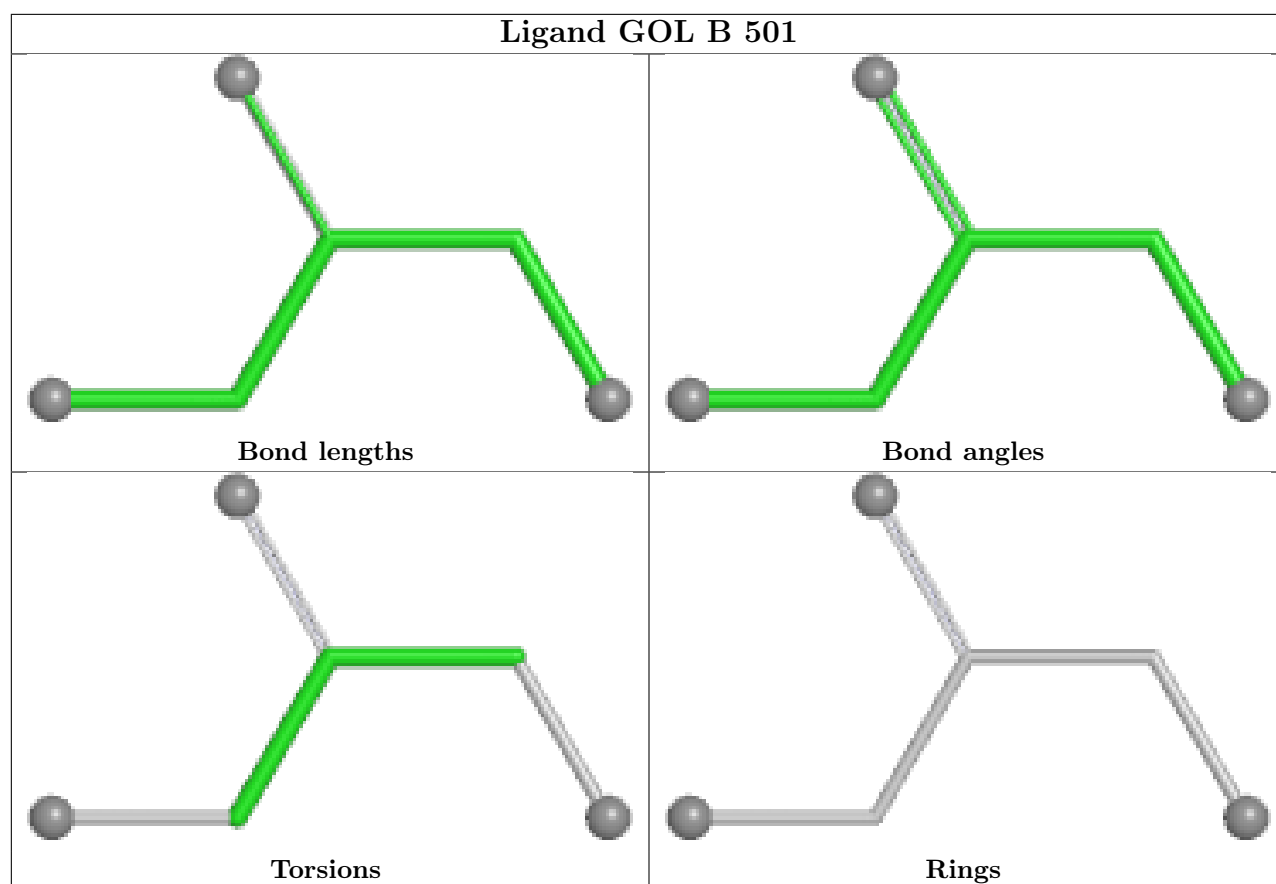
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	402/496 (81%)	1.09	68 (16%) 4 4	26, 44, 67, 85	0
1	B	407/496 (82%)	0.73	31 (7%) 20 21	23, 38, 64, 84	0
All	All	809/992 (81%)	0.91	99 (12%) 8 8	23, 42, 65, 85	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	428	LEU	4.4
1	A	411	VAL	4.2
1	A	414	LEU	4.1
1	B	411	VAL	3.9
1	A	463	THR	3.7
1	A	220	LEU	3.7
1	A	90	VAL	3.7
1	A	457	LEU	3.5
1	A	385	VAL	3.5
1	B	451	GLU	3.4
1	A	407	VAL	3.3
1	B	414	LEU	3.3
1	A	165	VAL	3.3
1	A	199	ALA	3.1
1	A	16	THR	3.1
1	B	385	VAL	3.1
1	A	84	GLN	3.1
1	A	148	VAL	3.1
1	B	465	GLY	3.0
1	A	301	TYR	3.0
1	A	305	SER	2.9
1	A	152	MET	2.8
1	B	138	THR	2.8
1	A	383	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	407	VAL	2.8
1	A	160	ARG	2.7
1	A	450	PHE	2.7
1	A	384	ALA	2.7
1	B	16	THR	2.6
1	B	305	SER	2.6
1	A	454	GLY	2.6
1	A	261	ILE	2.6
1	A	266	GLY	2.6
1	B	384	ALA	2.6
1	A	274	PRO	2.6
1	A	409	PRO	2.6
1	A	461	CYS	2.6
1	A	193	PHE	2.5
1	B	95	PHE	2.5
1	A	316	ALA	2.5
1	A	419	ALA	2.5
1	B	419	ALA	2.5
1	A	437	VAL	2.5
1	A	293	THR	2.5
1	B	282	GLN	2.5
1	B	447	THR	2.5
1	A	166	ALA	2.5
1	B	276	LEU	2.5
1	A	298	ILE	2.5
1	A	451	GLU	2.4
1	A	304	THR	2.4
1	A	452	TRP	2.4
1	B	301	TYR	2.4
1	A	147	THR	2.4
1	B	238	LEU	2.4
1	A	146	VAL	2.4
1	A	40	SER	2.4
1	A	279	TRP	2.4
1	B	459	GLU	2.4
1	B	63	PRO	2.4
1	A	85	LEU	2.3
1	B	445	LEU	2.3
1	A	418	LEU	2.3
1	A	433	GLY	2.3
1	B	452	TRP	2.3
1	A	144	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	135	MET	2.3
1	A	257	ILE	2.3
1	A	435	ILE	2.3
1	A	93	CYS	2.3
1	A	107	LEU	2.3
1	A	162	GLU	2.2
1	A	168	ILE	2.2
1	A	325	GLY	2.2
1	B	340	GLU	2.2
1	A	265	PHE	2.2
1	B	274	PRO	2.2
1	B	462	GLU	2.2
1	A	297	THR	2.2
1	B	308	MET	2.2
1	B	69	VAL	2.2
1	A	198	GLN	2.2
1	A	317	LEU	2.2
1	A	349	LEU	2.2
1	A	292	ALA	2.1
1	A	173	LEU	2.1
1	A	278	LEU	2.1
1	A	315	THR	2.1
1	A	112	PHE	2.1
1	A	106	VAL	2.1
1	B	56	LEU	2.1
1	A	223	GLU	2.1
1	A	205	VAL	2.1
1	B	169	LYS	2.1
1	A	428	LEU	2.1
1	A	132	ILE	2.1
1	A	331	TYR	2.0
1	A	434	TYR	2.0
1	B	442	GLN	2.0

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

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
1	LLP	A	236	24/25	0.93	0.10	30,36,45,51	0
1	LLP	B	236	24/25	0.93	0.11	28,35,40,44	0

6.3 Carbohydrates [i](#)

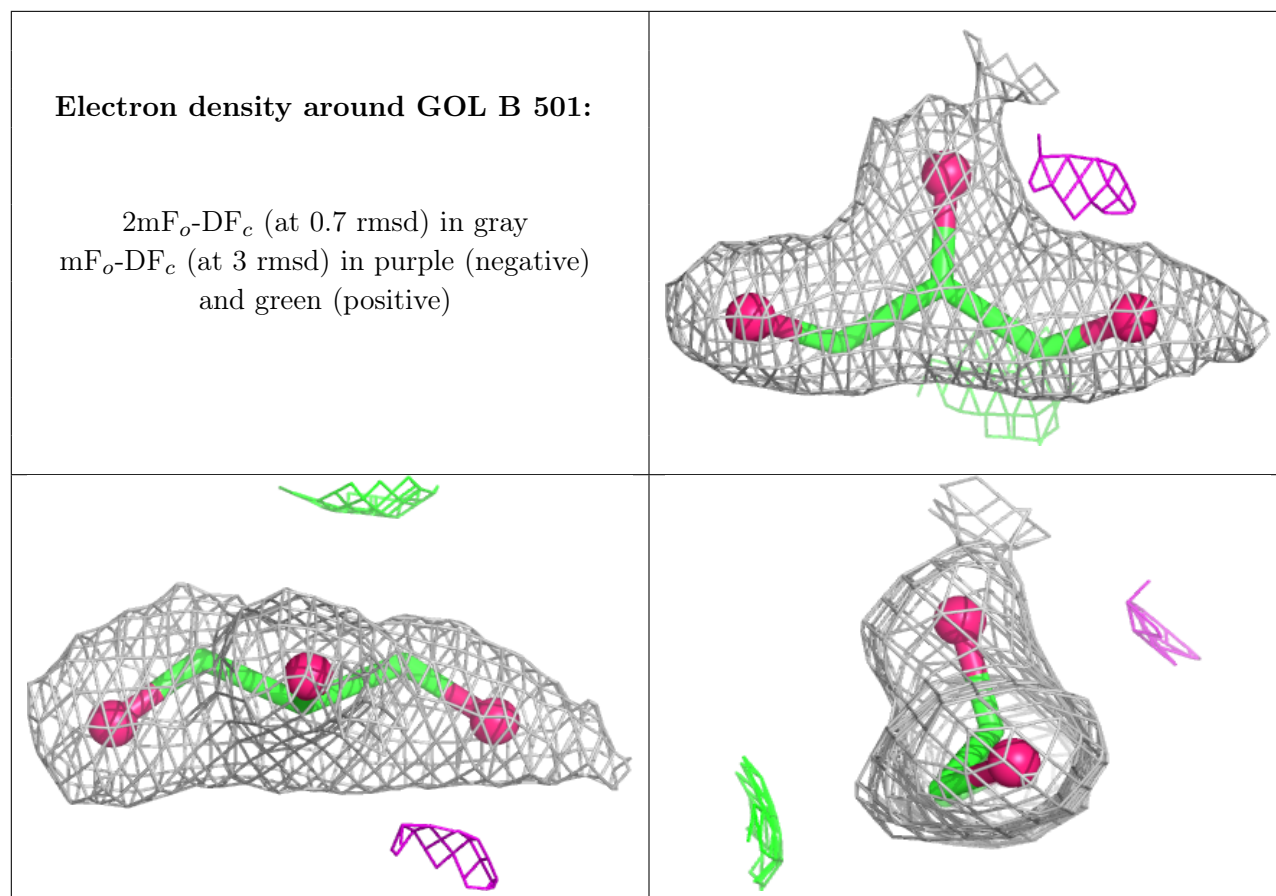
There are no oligosaccharides in this entry.

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	GOL	B	501	6/6	0.87	0.11	44,48,54,55	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.