



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:18 AM UTC

PDB ID : 5B7S / pdb_00005b7s
Title : Apo structure of Cysteine Desulfurase from Thermococcus onnurineus NA1
Authors : Ho, T.-H.; Kang, L.W.
Deposited on : 2016-06-08
Resolution : 2.58 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
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

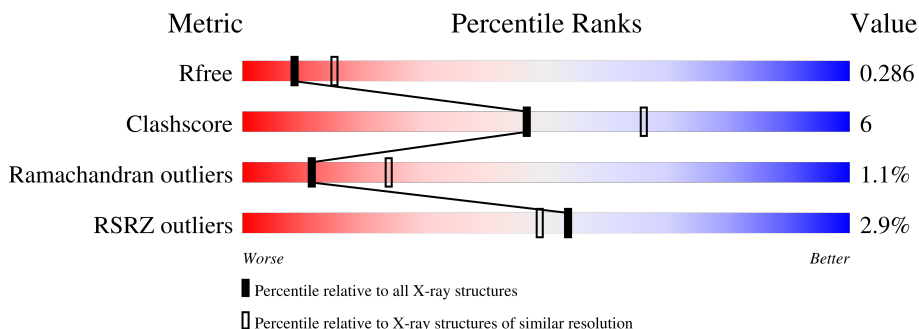
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	419	 3% 79% 13% • 5%
1	B	419	 2% 83% 12% 5%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6264 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 Cysteine desulfurase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	397	Total	C	N	O	P	S	0	0	0
			3111	1982	538	581	1	9			
1	B	400	Total	C	N	O	P	S	0	0	0
			3040	1937	518	574	1	10			

There are 40 discrepancies between the modelled and reference sequences:

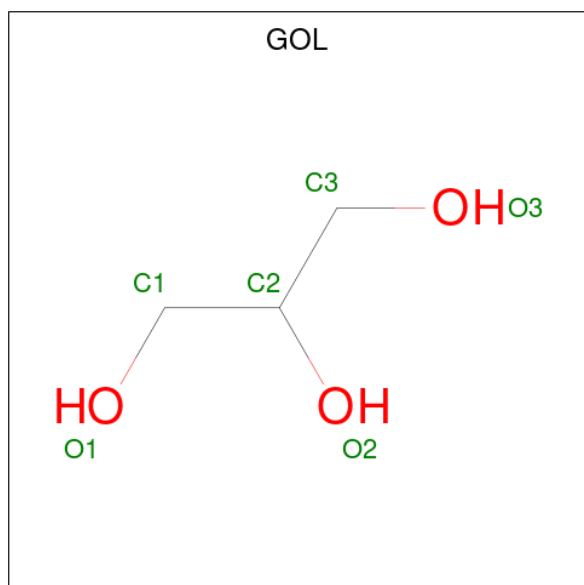
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP B6YT87
A	-18	HIS	-	expression tag	UNP B6YT87
A	-17	HIS	-	expression tag	UNP B6YT87
A	-16	HIS	-	expression tag	UNP B6YT87
A	-15	HIS	-	expression tag	UNP B6YT87
A	-14	HIS	-	expression tag	UNP B6YT87
A	-13	HIS	-	expression tag	UNP B6YT87
A	-12	SER	-	expression tag	UNP B6YT87
A	-11	SER	-	expression tag	UNP B6YT87
A	-10	GLU	-	expression tag	UNP B6YT87
A	-9	ASN	-	expression tag	UNP B6YT87
A	-8	LEU	-	expression tag	UNP B6YT87
A	-7	TYR	-	expression tag	UNP B6YT87
A	-6	PHE	-	expression tag	UNP B6YT87
A	-5	GLN	-	expression tag	UNP B6YT87
A	-4	GLY	-	expression tag	UNP B6YT87
A	-3	HIS	-	expression tag	UNP B6YT87
A	-2	MET	-	expression tag	UNP B6YT87
A	-1	ALA	-	expression tag	UNP B6YT87
A	0	SER	-	expression tag	UNP B6YT87
B	-19	MET	-	expression tag	UNP B6YT87
B	-18	HIS	-	expression tag	UNP B6YT87
B	-17	HIS	-	expression tag	UNP B6YT87
B	-16	HIS	-	expression tag	UNP B6YT87
B	-15	HIS	-	expression tag	UNP B6YT87

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP B6YT87
B	-13	HIS	-	expression tag	UNP B6YT87
B	-12	SER	-	expression tag	UNP B6YT87
B	-11	SER	-	expression tag	UNP B6YT87
B	-10	GLU	-	expression tag	UNP B6YT87
B	-9	ASN	-	expression tag	UNP B6YT87
B	-8	LEU	-	expression tag	UNP B6YT87
B	-7	TYR	-	expression tag	UNP B6YT87
B	-6	PHE	-	expression tag	UNP B6YT87
B	-5	GLN	-	expression tag	UNP B6YT87
B	-4	GLY	-	expression tag	UNP B6YT87
B	-3	HIS	-	expression tag	UNP B6YT87
B	-2	MET	-	expression tag	UNP B6YT87
B	-1	ALA	-	expression tag	UNP B6YT87
B	0	SER	-	expression tag	UNP B6YT87

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

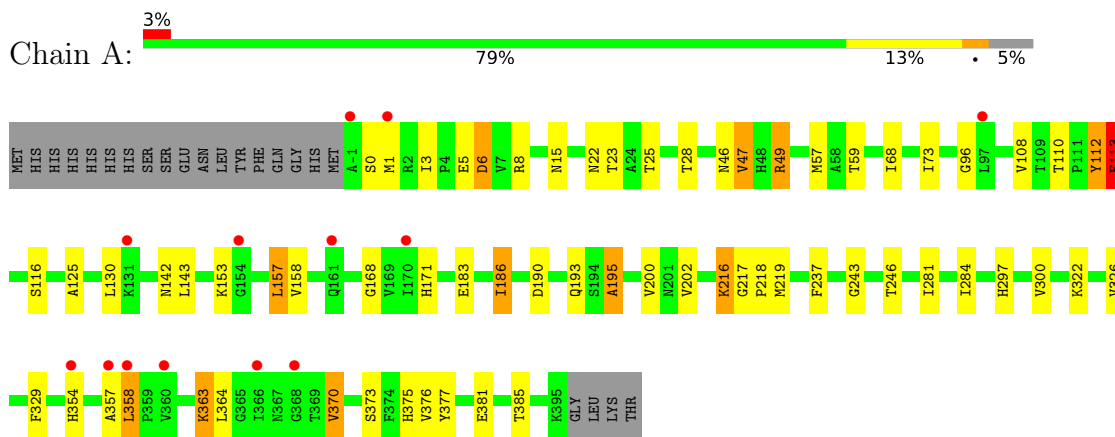
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	57	Total 57	O 57	0	0
3	B	44	Total 44	O 44	0	0

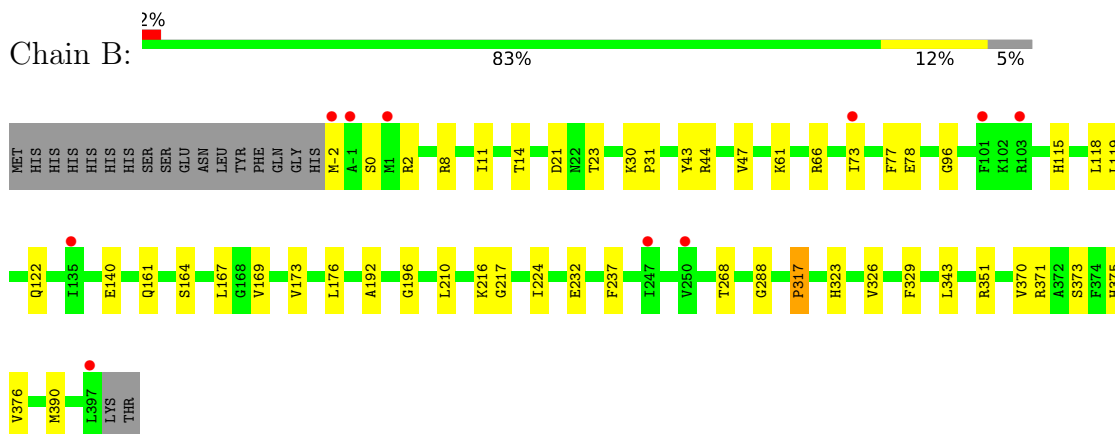
3 Residue-property plots [i](#)

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: Cysteine desulfurase



- Molecule 1: Cysteine desulfurase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.01Å 92.53Å 145.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.36 – 2.58 36.36 – 2.58	Depositor EDS
% Data completeness (in resolution range)	99.2 (36.36-2.58) 99.2 (36.36-2.58)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.20 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.203 , 0.295 (Not available) , 0.286	Depositor DCC
R_{free} test set	1465 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6264	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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

5.1 Standard geometry

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	0.92	2/3153 (0.1%)	1.20	21/4263 (0.5%)
1	B	0.87	0/3081	1.07	6/4187 (0.1%)
All	All	0.90	2/6234 (0.0%)	1.14	27/8450 (0.3%)

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	A	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	6	ASP	N-CA	6.70	1.55	1.46
1	A	46	ASN	N-CA	5.26	1.52	1.45

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	46	ASN	CA-C-N	9.20	138.26	121.70
1	A	46	ASN	C-N-CA	9.20	138.26	121.70
1	A	357	ALA	N-CA-C	-8.10	97.42	109.81
1	A	113	GLU	N-CA-C	-7.75	94.30	110.80
1	A	363	LYS	N-CA-C	7.48	120.08	111.11
1	A	3	ILE	N-CA-C	6.88	114.71	107.76
1	B	2	ARG	N-CA-C	6.87	119.75	109.25
1	A	6	ASP	CA-C-N	6.45	133.32	121.70
1	A	6	ASP	C-N-CA	6.45	133.32	121.70

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	6	ASP	N-CA-C	6.39	120.53	107.69
1	A	47	VAL	CB-CA-C	6.32	123.40	111.40
1	B	192	ALA	N-CA-C	6.23	118.87	111.33
1	A	46	ASN	O-C-N	-5.90	115.45	122.65
1	A	363	LYS	CA-C-N	5.84	132.22	121.70
1	A	363	LYS	C-N-CA	5.84	132.22	121.70
1	A	46	ASN	N-CA-C	-5.64	103.48	110.41
1	B	317	PRO	N-CA-C	5.47	120.17	111.26
1	B	268	THR	CA-C-N	-5.46	114.97	120.21
1	B	268	THR	C-N-CA	-5.46	114.97	120.21
1	A	157	LEU	N-CA-C	5.41	117.41	109.24
1	A	195	ALA	N-CA-C	5.39	117.85	111.33
1	A	46	ASN	CA-C-O	-5.34	115.90	121.99
1	B	169	VAL	CB-CA-C	-5.25	102.67	111.29
1	A	15	ASN	N-CA-C	5.24	119.52	113.18
1	A	186	ILE	N-CA-CB	5.23	116.23	110.53
1	A	116	SER	N-CA-C	5.09	118.26	111.75
1	A	370	VAL	CB-CA-C	5.00	118.03	110.98

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	0	SER	Peptide
1	A	112	TYR	Peptide

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	3111	0	3108	40	0
1	B	3040	0	2940	33	0
2	A	6	0	8	0	0
2	B	6	0	8	0	0
3	A	57	0	0	1	1
3	B	44	0	0	0	1
All	All	6264	0	6064	70	1

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 (70) 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:6:ASP:N	1:A:6:ASP:OD1	2.10	0.85
1:B:23:THR:HG23	1:B:373:SER:CB	2.20	0.71
1:B:23:THR:OG1	1:B:371:ARG:CZ	2.47	0.63
1:B:73:ILE:HD11	1:B:210:LEU:HD22	1.80	0.63
1:B:21:ASP:OD2	1:B:351:ARG:NH1	2.32	0.62
1:A:193:GLN:NE2	1:A:216:LLP:O3	2.32	0.62
1:A:153:LYS:HA	1:A:183:GLU:HG2	1.81	0.61
1:B:161:GLN:HE22	1:B:164:SER:HB2	1.66	0.60
1:A:193:GLN:HE22	1:A:216:LLP:C3	2.16	0.59
1:A:354:HIS:HB2	1:A:358:LEU:HD12	1.86	0.58
1:A:142:ASN:OD1	1:A:322:LYS:HE3	2.04	0.57
1:A:28:THR:CG2	1:B:44:ARG:HD3	2.35	0.57
1:A:23:THR:HG23	1:A:373:SER:CB	2.35	0.56
1:A:73:ILE:O	1:A:202:VAL:HG22	2.06	0.55
1:A:25:THR:OG1	1:B:44:ARG:NH2	2.41	0.53
1:A:193:GLN:HG3	3:A:548:HOH:O	2.08	0.53
1:A:153:LYS:CA	1:A:183:GLU:HG2	2.38	0.52
1:A:375:HIS:CD2	1:A:376:VAL:H	2.27	0.52
1:A:363:LYS:HB3	1:A:364:LEU:HB2	1.92	0.52
1:A:5:GLU:O	1:A:8:ARG:HG3	2.10	0.52
1:A:297:HIS:HA	1:A:300:VAL:HG12	1.93	0.50
1:A:6:ASP:HA	1:A:8:ARG:H	1.75	0.50
1:B:78:GLU:H	1:B:78:GLU:CD	2.19	0.50
1:B:140:GLU:CG	1:B:317:PRO:HA	2.42	0.50
1:B:23:THR:HG23	1:B:373:SER:HB3	1.94	0.50
1:B:343:LEU:HD11	1:B:390:MET:HE3	1.94	0.49
1:B:8:ARG:HB3	1:B:14:THR:HG21	1.95	0.49
1:A:6:ASP:HA	1:A:8:ARG:N	2.27	0.49
1:B:23:THR:HG23	1:B:373:SER:HB2	1.93	0.48
1:B:11:ILE:O	1:B:14:THR:HB	2.14	0.48
1:A:125:ALA:HA	1:A:130:LEU:HB2	1.96	0.47
1:A:157:LEU:HB2	1:A:186:ILE:HG12	1.96	0.47
1:B:217:GLY:O	1:B:375:HIS:HE1	1.96	0.47
1:B:375:HIS:CD2	1:B:376:VAL:H	2.33	0.47
1:B:343:LEU:HD21	1:B:390:MET:HE3	1.97	0.46
1:B:115:HIS:HA	1:B:118:LEU:CD2	2.44	0.46
1:B:164:SER:HG	1:B:167:LEU:H	1.62	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:PHE:CZ	1:B:370:VAL:HG22	2.51	0.46
1:B:-2:MET:HE2	1:B:288:GLY:HA3	1.98	0.46
1:A:22:ASN:OD1	1:A:377:TYR:OH	2.25	0.45
1:A:28:THR:HG22	1:B:44:ARG:HD3	1.98	0.45
1:A:381:GLU:O	1:A:385:THR:HG23	2.16	0.45
1:A:217:GLY:O	1:A:375:HIS:HE1	1.99	0.45
1:B:43:TYR:OH	1:B:61:LYS:HG2	2.17	0.45
1:A:218:PRO:HA	1:A:284:ILE:CD1	2.46	0.45
1:A:96:GLY:HA3	1:A:237:PHE:CE1	2.52	0.44
1:A:195:ALA:HA	1:A:200:VAL:HG11	2.00	0.44
1:B:96:GLY:HA3	1:B:237:PHE:CE1	2.52	0.43
1:B:173:VAL:HA	1:B:176:LEU:HB2	1.99	0.43
1:B:323:HIS:HE1	1:B:326:VAL:O	2.00	0.43
1:A:57:MET:HE3	1:A:57:MET:HA	2.00	0.43
1:A:68:ILE:HG22	1:A:281:ILE:HG12	2.00	0.43
1:A:110:THR:OG1	1:A:113:GLU:OE2	2.32	0.43
1:B:115:HIS:HA	1:B:118:LEU:HD23	2.00	0.43
1:B:161:GLN:HE22	1:B:164:SER:CB	2.30	0.43
1:A:243:GLY:O	1:A:246:THR:HB	2.18	0.43
1:B:119:LEU:HA	1:B:122:GLN:HB2	2.00	0.43
1:A:143:LEU:HD22	1:A:171:HIS:CD2	2.54	0.43
1:A:157:LEU:CB	1:A:186:ILE:HG12	2.49	0.42
1:A:190:ASP:OD2	1:A:216:LLP:N1	2.53	0.42
1:A:219:MET:HE3	1:A:284:ILE:HG13	2.02	0.42
1:A:218:PRO:HA	1:A:284:ILE:HD12	2.01	0.42
1:A:168:GLY:HA3	1:A:326:VAL:HG13	2.02	0.42
1:B:224:ILE:O	1:B:224:ILE:HG23	2.20	0.42
1:A:329:PHE:CZ	1:A:370:VAL:HG22	2.55	0.41
1:A:108:VAL:HG23	1:A:158:VAL:HG13	2.01	0.41
1:B:30:LYS:N	1:B:31:PRO:CD	2.84	0.41
1:B:196:GLY:HA3	1:B:217:GLY:C	2.45	0.41
1:A:49:ARG:HG3	1:A:59:THR:HG21	2.03	0.41
1:B:66:ARG:HD2	1:B:77:PHE:CE1	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:552:HOH:O	3:B:502:HOH:O[4_445]	2.05	0.15

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	394/419 (94%)	372 (94%)	16 (4%)	6 (2%)	8	16
1	B	397/419 (95%)	379 (96%)	15 (4%)	3 (1%)	16	32
All	All	791/838 (94%)	751 (95%)	31 (4%)	9 (1%)	11	24

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	47	VAL
1	A	113	GLU
1	A	358	LEU
1	A	1	MET
1	A	49	ARG
1	A	112	TYR
1	B	0	SER
1	B	47	VAL
1	B	232	GLU

5.3.2 Protein sidechains [i](#)

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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	A	216	1	23,24,25	3.13	5 (21%)	25,32,34	1.98	7 (28%)
1	LLP	B	216	1	23,24,25	2.98	5 (21%)	25,32,34	1.45	3 (12%)

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	A	216	1	-	3/16/17/19	0/1/1/1
1	LLP	B	216	1	-	4/16/17/19	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	216	LLP	C3-C2	10.94	1.52	1.41
1	B	216	LLP	C3-C2	9.60	1.50	1.41
1	A	216	LLP	C4'-NZ	5.61	1.45	1.27
1	B	216	LLP	C4-C5	5.58	1.49	1.42
1	A	216	LLP	C4-C5	5.41	1.49	1.42
1	B	216	LLP	C4'-NZ	5.24	1.44	1.27
1	B	216	LLP	C4-C3	5.16	1.49	1.41
1	A	216	LLP	C4-C3	4.49	1.48	1.41
1	A	216	LLP	C4-C4'	3.82	1.54	1.46
1	B	216	LLP	C4-C4'	3.50	1.54	1.46

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	216	LLP	C3-C4-C5	-4.91	114.33	118.28
1	A	216	LLP	O3-C3-C2	3.95	125.77	117.58
1	B	216	LLP	C4-C3-C2	-3.91	117.94	120.14
1	B	216	LLP	C3-C4-C5	-3.12	115.77	118.28
1	A	216	LLP	C5-C4-C4'	3.04	126.17	121.47
1	A	216	LLP	C2'-C2-C3	2.98	124.28	120.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	216	LLP	OP4-P-OP1	-2.45	99.83	106.44
1	B	216	LLP	C6-N1-C2	2.27	123.31	119.20
1	A	216	LLP	C4-C4'-NZ	-2.11	114.31	124.04
1	A	216	LLP	C6-N1-C2	2.08	122.98	119.20

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	216	LLP	C4-C4'-NZ-CE
1	A	216	LLP	O-C-CA-CB
1	B	216	LLP	C4-C4'-NZ-CE
1	B	216	LLP	O-C-CA-CB
1	B	216	LLP	CG-CD-CE-NZ
1	A	216	LLP	C3-C4-C4'-NZ
1	B	216	LLP	C3-C4-C4'-NZ

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	216	LLP	3	0

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	GOL	A	401	-	5,5,5	0.50	0	5,5,5	0.47	0
2	GOL	B	401	-	5,5,5	0.56	0	5,5,5	0.35	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	A	401	-	-	0/4/4/4	-
2	GOL	B	401	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	GOL	C1-C2-C3-O3
2	B	401	GOL	O2-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

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	396/419 (94%)	0.23	13 (3%) 49 44	28, 50, 82, 108	0
1	B	399/419 (95%)	0.20	10 (2%) 58 54	29, 50, 75, 106	0
All	All	795/838 (94%)	0.22	23 (2%) 53 49	28, 50, 79, 108	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	357	ALA	3.7
1	B	247	ILE	3.6
1	B	-1	ALA	3.3
1	A	1	MET	3.2
1	B	-2	MET	3.1
1	A	358	LEU	2.9
1	A	-1	ALA	2.9
1	A	354	HIS	2.7
1	A	366	ILE	2.7
1	B	1	MET	2.7
1	A	154	GLY	2.7
1	A	161	GLN	2.7
1	A	131	LYS	2.4
1	B	73	ILE	2.4
1	A	170	ILE	2.3
1	B	397	LEU	2.3
1	A	360	VAL	2.3
1	A	97	LEU	2.2
1	B	135	ILE	2.2
1	B	101	PHE	2.2
1	A	368	GLY	2.1
1	B	250	VAL	2.1
1	B	103	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	LLP	A	216	24/25	0.95	0.10	31,56,67,69	0
1	LLP	B	216	24/25	0.95	0.10	34,49,74,77	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	401	6/6	0.81	0.16	73,76,77,77	0
2	GOL	A	401	6/6	0.89	0.10	51,56,64,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.