



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:49 PM UTC

PDB ID : 6ZHT / pdb_00006zht
Title : Uba1-Ubc13 disulfide mediated complex
Authors : Schaefer, A.; Misra, M.; Schindelin, H.
Deposited on : 2020-06-23
Resolution : 2.30 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

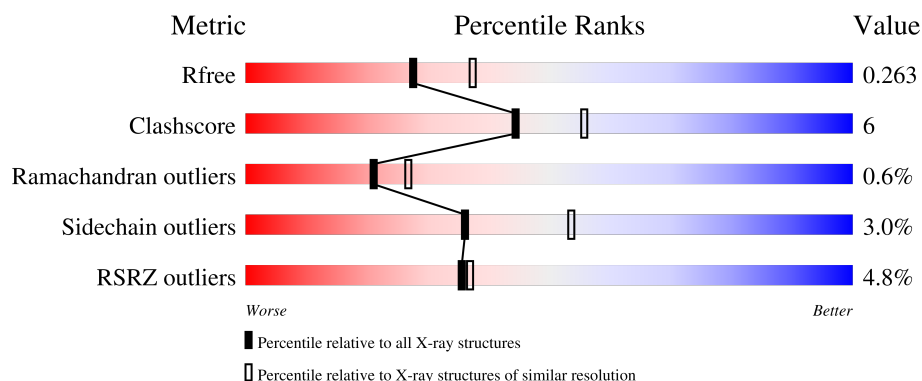
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	C	1001	
2	B	153	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18472 atoms, of which 9042 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 Ubiquitin-activating enzyme E1 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	C	983	Total	C	H	N	O	S	0	8	0
			15478	4958	7696	1288	1513	23			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	24	ALA	-	expression tag	UNP P22515
C	479	ALA	LEU	conflict	UNP P22515

- Molecule 2 is a protein called Ubiquitin-conjugating enzyme E2 13.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	153	Total	C	H	N	O	S	0	2	0
			2485	793	1250	207	231	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP P52490

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		

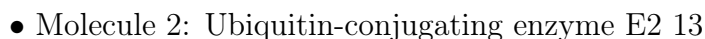
- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	4	Total	Cl	0	0
			4	4		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	287	Total	O	0	1
			288	288		
5	B	48	Total	O	0	1
			49	49		

- Molecule 1: Ubiquitin-activating enzyme E1 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	181.39Å 58.27Å 137.46Å 90.00° 109.72° 90.00°	Depositor
Resolution (Å)	24.71 – 2.30 24.71 – 2.30	Depositor EDS
% Data completeness (in resolution range)	71.4 (24.71-2.30) 71.3 (24.71-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.31Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.206 , 0.263 0.206 , 0.263	Depositor DCC
R_{free} test set	2177 reflections (3.60%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	18472	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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: GOL, CL

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	C	0.17	0/7971	0.39	0/10784
2	B	0.17	0/1271	0.42	0/1730
All	All	0.17	0/9242	0.39	0/12514

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	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	36	THR	Peptide

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	C	7782	7696	7665	98	0
2	B	1235	1250	1244	16	0
3	B	6	8	8	0	0
3	C	66	88	88	0	0
4	C	4	0	0	0	0
5	B	49	0	0	0	0
5	C	288	0	0	2	2
All	All	9430	9042	9005	109	2

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 (109) 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:C:449:MET:HE2	1:C:540:THR:HG21	1.45	0.98
1:C:928:TRP:O	1:C:930:ARG:NH2	2.22	0.73
1:C:25:VAL:HG21	1:C:28:LYS:HZ1	1.58	0.68
1:C:958:SER:OG	2:B:9:ILE:HD11	1.94	0.68
2:B:20:VAL:HG13	2:B:23:ILE:HB	1.76	0.67
1:C:449:MET:HE1	1:C:877:VAL:CG1	2.26	0.65
1:C:416:ARG:NH1	5:C:1201:HOH:O	2.28	0.65
1:C:222:LEU:HD22	1:C:243:VAL:HG11	1.79	0.63
1:C:413:VAL:O	1:C:413:VAL:HG12	1.96	0.63
1:C:869:THR:O	1:C:873:VAL:HG23	1.98	0.63
1:C:933:ILE:HB	1:C:1022:ILE:HD13	1.81	0.62
1:C:968:PHE:CD2	2:B:16:VAL:HG21	2.36	0.60
1:C:966:ALA:HB3	1:C:969:PHE:CD1	2.37	0.60
1:C:994:ILE:HD11	1:C:1001:MET:HE1	1.84	0.59
1:C:728:PHE:O	1:C:732:VAL:HG23	2.03	0.59
1:C:449:MET:HE1	1:C:877:VAL:HG11	1.85	0.58
1:C:227:LEU:CD2	1:C:258:VAL:HG21	2.34	0.58
1:C:440:VAL:HG12	1:C:543:LEU:HD11	1.85	0.58
1:C:995:PRO:O	1:C:996:ALA:HB3	2.04	0.57
1:C:227:LEU:HD22	1:C:258:VAL:HG21	1.86	0.57
1:C:1006:CYS:SG	2:B:9:ILE:HG22	2.44	0.57
1:C:854:LYS:O	1:C:858:ILE:HD12	2.05	0.57
1:C:40:LEU:HD21	1:C:42:LEU:HD21	1.85	0.56
1:C:957:LEU:C	1:C:957:LEU:HD23	2.31	0.56
1:C:137:LYS:NZ	1:C:155:GLU:OE2	2.39	0.56
1:C:558:PHE:O	1:C:930:ARG:NH1	2.39	0.55
1:C:956:MET:HE3	1:C:963:LEU:HD11	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:294:LEU:HD22	1:C:329:LEU:HD22	1.88	0.55
1:C:958:SER:CB	2:B:9:ILE:HD11	2.37	0.55
1:C:28:LYS:HG2	1:C:32:LEU:HD13	1.90	0.54
1:C:294:LEU:HD22	1:C:329:LEU:CD2	2.36	0.54
1:C:25:VAL:HG12	1:C:853:GLN:HG2	1.90	0.53
1:C:71:VAL:HG22	1:C:91:ARG:HG2	1.90	0.53
1:C:163:THR:HG21	1:C:369:LEU:HD21	1.90	0.53
1:C:160:PHE:CZ	1:C:394:LEU:HD21	2.44	0.53
1:C:79:GLN:HG2	1:C:82:LEU:HD12	1.91	0.52
2:B:76:TYR:CE2	2:B:125:VAL:HG13	2.44	0.52
1:C:696:PRO:HD2	1:C:699:ALA:HB2	1.92	0.51
1:C:449:MET:HE2	1:C:540:THR:CG2	2.31	0.51
1:C:25:VAL:HG11	1:C:28:LYS:NZ	2.26	0.51
1:C:78:THR:HA	1:C:455:LEU:HD13	1.93	0.51
1:C:466:ILE:HB	1:C:513:ILE:HD13	1.93	0.51
1:C:807:ASP:OD2	1:C:809:SER:OG	2.27	0.50
1:C:509:LEU:HD23	1:C:513:ILE:HD11	1.93	0.50
1:C:228:PHE:HB3	1:C:240:ILE:HG23	1.93	0.50
2:B:4:LEU:HB3	2:B:5:PRO:HD3	1.94	0.50
1:C:219:LEU:HD23	1:C:247:GLY:C	2.37	0.50
1:C:176:PRO:HD2	1:C:259:LYS:HG2	1.93	0.50
1:C:995:PRO:O	1:C:996:ALA:CB	2.60	0.49
2:B:8:ILE:HG21	2:B:32:LEU:HB3	1.93	0.49
1:C:449:MET:HE1	1:C:877:VAL:HG12	1.93	0.49
1:C:957:LEU:HD23	1:C:958:SER:N	2.28	0.49
1:C:231:GLU:HB3	1:C:239:ARG:HB3	1.95	0.49
1:C:797:GLU:N	1:C:797:GLU:OE2	2.45	0.48
1:C:437:VAL:HG11	1:C:453:TRP:CH2	2.48	0.48
2:B:71:PHE:CD1	2:B:86:ILE:HD11	2.49	0.47
1:C:49:VAL:HG11	1:C:79:GLN:HE22	1.78	0.47
1:C:705:GLU:HB3	1:C:706:PRO:HD2	1.96	0.47
1:C:450:LEU:HB3	1:C:502:VAL:HG11	1.96	0.47
1:C:406:ASN:O	1:C:410:THR:HB	2.14	0.46
1:C:230:VAL:HG13	1:C:238:PHE:CD2	2.50	0.46
1:C:413:VAL:O	1:C:413:VAL:CG1	2.63	0.46
1:C:96:ARG:HD3	1:C:110:VAL:HG23	1.98	0.46
2:B:55:GLU:HG3	2:B:72:LEU:HD11	1.98	0.46
1:C:619:LEU:HD13	5:C:1338:HOH:O	2.16	0.46
1:C:49:VAL:HG11	1:C:79:GLN:OE1	2.16	0.46
1:C:410:THR:CG2	1:C:410:THR:O	2.63	0.45
2:B:71:PHE:HD1	2:B:86:ILE:HD11	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:881:LEU:C	1:C:881:LEU:HD23	2.41	0.45
1:C:227:LEU:HD22	1:C:258:VAL:CG2	2.46	0.45
1:C:958:SER:HB3	2:B:9:ILE:HD11	1.99	0.45
1:C:410:THR:HG23	1:C:422:ALA:HA	1.98	0.45
1:C:631:VAL:HB	1:C:816:LEU:HD12	1.98	0.45
1:C:53:LYS:NZ	1:C:448:GLU:OE1	2.48	0.44
1:C:71:VAL:HG13	1:C:82:LEU:HD13	1.99	0.44
1:C:211:VAL:HG13	1:C:230:VAL:HG21	2.00	0.44
1:C:994:ILE:HG22	1:C:996:ALA:H	1.82	0.44
1:C:615:TRP:CE3	1:C:841:ARG:HD2	2.53	0.44
1:C:49:VAL:HG11	1:C:79:GLN:NE2	2.32	0.44
1:C:964:LEU:HD23	1:C:989:VAL:HG21	2.00	0.44
1:C:150:ARG:NH1	1:C:377:ALA:O	2.51	0.43
1:C:313:THR:HG22	1:C:410:THR:HG21	1.99	0.43
1:C:79:GLN:HG2	1:C:82:LEU:CD1	2.48	0.43
1:C:994:ILE:CD1	1:C:1001:MET:HE1	2.48	0.43
1:C:79:GLN:CG	1:C:82:LEU:HG	2.49	0.43
1:C:499:ALA:CB	1:C:513:ILE:HG21	2.49	0.43
1:C:955:THR:HG23	1:C:967:SER:HB2	2.01	0.42
2:B:95:TRP:HA	2:B:99:LEU:HD12	2.02	0.42
1:C:198:LEU:HD21	1:C:710:GLY:HA2	2.01	0.42
1:C:922:LYS:NZ	1:C:950:GLU:OE1	2.47	0.42
2:B:81:ASP:OD1	2:B:81:ASP:C	2.62	0.42
1:C:557:VAL:HA	1:C:928:TRP:CZ3	2.54	0.42
2:B:43:GLU:O	2:B:44:GLN:HB2	2.20	0.42
1:C:313:THR:HG21	1:C:410:THR:CG2	2.49	0.42
1:C:936:ASP:OD2	1:C:983:THR:OG1	2.32	0.42
1:C:59:GLY:O	1:C:60:VAL:HG13	2.21	0.41
1:C:301:GLN:HE22	1:C:329:LEU:HD12	1.85	0.41
1:C:722:ILE:HG12	1:C:755:VAL:HG13	2.03	0.41
1:C:481:ARG:C	1:C:482[B]:GLN:HG2	2.46	0.41
1:C:482[A]:GLN:HA	1:C:482[A]:GLN:NE2	2.36	0.41
1:C:924:TYR:CD2	1:C:1018:PRO:HG3	2.56	0.41
1:C:977:ARG:HA	1:C:980:LEU:HD12	2.01	0.41
1:C:481:ARG:O	1:C:482[B]:GLN:CB	2.69	0.41
1:C:634:TYR:OH	1:C:806:PRO:O	2.36	0.41
2:B:1[B]:MET:C	2:B:1[B]:MET:SD	3.04	0.41
1:C:50:GLU:HG3	1:C:363:VAL:HG12	2.03	0.40
1:C:313:THR:CG2	1:C:410:THR:HG21	2.51	0.40
1:C:481:ARG:O	1:C:482[A]:GLN:HB3	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:1460:HOH:O	5:C:1460:HOH:O[2_555]	1.45	0.75
5:C:1214:HOH:O	5:C:1424:HOH:O[2_556]	2.13	0.07

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	C	987/1001 (99%)	934 (95%)	46 (5%)	7 (1%)	18	23
2	B	153/153 (100%)	150 (98%)	2 (1%)	1 (1%)	18	23
All	All	1140/1154 (99%)	1084 (95%)	48 (4%)	8 (1%)	21	23

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	193	GLY
1	C	750	ASN
1	C	996	ALA
1	C	220	ASP
2	B	92	LYS
1	C	482[A]	GLN
1	C	482[B]	GLN
1	C	194	THR

5.3.2 Protein sidechains [i](#)

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	C	871/879 (99%)	845 (97%)	26 (3%)	36	53
2	B	136/135 (101%)	131 (96%)	5 (4%)	30	45
All	All	1007/1014 (99%)	976 (97%)	31 (3%)	36	52

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	112	ASP
1	C	130	ASP
1	C	180	GLU
1	C	250	LYS
1	C	251	LYS
1	C	263	LYS
1	C	265	SER
1	C	299	LEU
1	C	307[A]	ASN
1	C	307[B]	ASN
1	C	312	ARG
1	C	329	LEU
1	C	395	GLU
1	C	410	THR
1	C	430	LYS
1	C	517	ILE
1	C	619	LEU
1	C	670	GLU
1	C	747	ASP
1	C	748	ASP
1	C	869	THR
1	C	930	ARG
1	C	961	VAL
1	C	972	LYS
1	C	978	LEU
1	C	999	SER
2	B	17	SER
2	B	20	VAL
2	B	94	ASN
2	B	139	LYS
2	B	150	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	201	ASN
1	C	429	GLN
1	C	493	ASN
1	C	541	ASN
1	C	639	ASN
1	C	690	GLN
1	C	725	ASN
1	C	844	ASN
1	C	878	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](#)

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 for Mogul analysis.

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	$\# Z > 2$
3	GOL	C	1102	-	5,5,5	0.82	0	5,5,5	1.09	0
3	GOL	C	1103	-	5,5,5	0.83	0	5,5,5	1.04	0
3	GOL	C	1104	-	5,5,5	0.83	0	5,5,5	1.01	0
3	GOL	C	1106	-	5,5,5	0.77	0	5,5,5	0.95	0
3	GOL	B	201	-	5,5,5	0.84	0	5,5,5	0.97	0
3	GOL	C	1107	-	5,5,5	0.83	0	5,5,5	1.01	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	C	1108	-	5,5,5	0.81	0	5,5,5	1.03	0
3	GOL	C	1101	-	5,5,5	0.85	0	5,5,5	1.05	0
3	GOL	C	1105	-	5,5,5	0.83	0	5,5,5	1.04	0
3	GOL	C	1109	-	5,5,5	0.83	0	5,5,5	0.94	0
3	GOL	C	1111	-	5,5,5	0.81	0	5,5,5	0.89	0
3	GOL	C	1110	-	5,5,5	0.80	0	5,5,5	0.98	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
3	GOL	C	1102	-	-	1/4/4/4	-
3	GOL	C	1103	-	-	2/4/4/4	-
3	GOL	C	1104	-	-	2/4/4/4	-
3	GOL	C	1106	-	-	4/4/4/4	-
3	GOL	B	201	-	-	1/4/4/4	-
3	GOL	C	1107	-	-	2/4/4/4	-
3	GOL	C	1108	-	-	0/4/4/4	-
3	GOL	C	1101	-	-	2/4/4/4	-
3	GOL	C	1105	-	-	4/4/4/4	-
3	GOL	C	1109	-	-	2/4/4/4	-
3	GOL	C	1111	-	-	4/4/4/4	-
3	GOL	C	1110	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1101	GOL	C1-C2-C3-O3
3	C	1103	GOL	O1-C1-C2-C3
3	C	1105	GOL	O1-C1-C2-C3
3	C	1106	GOL	O1-C1-C2-C3
3	C	1107	GOL	C1-C2-C3-O3
3	C	1109	GOL	O1-C1-C2-O2
3	C	1109	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	C	1111	GOL	O1-C1-C2-C3
3	C	1104	GOL	O2-C2-C3-O3
3	C	1107	GOL	O2-C2-C3-O3
3	C	1111	GOL	O1-C1-C2-O2
3	C	1104	GOL	C1-C2-C3-O3
3	C	1105	GOL	C1-C2-C3-O3
3	C	1106	GOL	C1-C2-C3-O3
3	C	1111	GOL	C1-C2-C3-O3
3	C	1105	GOL	O1-C1-C2-O2
3	C	1106	GOL	O1-C1-C2-O2
3	C	1101	GOL	O2-C2-C3-O3
3	C	1102	GOL	O1-C1-C2-O2
3	C	1103	GOL	O1-C1-C2-O2
3	C	1111	GOL	O2-C2-C3-O3
3	B	201	GOL	O1-C1-C2-O2
3	C	1110	GOL	O2-C2-C3-O3
3	C	1106	GOL	O2-C2-C3-O3
3	C	1105	GOL	O2-C2-C3-O3
3	C	1110	GOL	O1-C1-C2-C3

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	C	983/1001 (98%)	0.14	49 (4%) 34 35	13, 42, 94, 156	4 (0%)
2	B	153/153 (100%)	0.07	5 (3%) 49 51	27, 43, 71, 114	1 (0%)
All	All	1136/1154 (98%)	0.13	54 (4%) 35 37	13, 42, 93, 156	5 (0%)

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	479	ALA	4.8
1	C	776	LEU	4.7
1	C	25	VAL	4.7
1	C	236	PHE	4.6
1	C	191	PRO	4.5
1	C	480	ASN	3.9
1	C	482[A]	GLN	3.9
1	C	24	ALA	3.7
1	C	526[A]	GLU	3.6
1	C	774	ALA	3.5
1	C	31	MET	3.5
1	C	478	ASN	3.5
1	C	30	ALA	3.4
1	C	702	SER	3.3
1	C	863	ILE	3.3
2	B	93	THR	3.3
2	B	94	ASN	3.2
1	C	243	VAL	3.2
1	C	26	LEU	3.2
1	C	638	PRO	3.1
1	C	749	SER	3.0
2	B	152	PRO	3.0
1	C	745	LYS	2.9
1	C	249	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	1[A]	MET	2.9
1	C	751	SER	2.8
1	C	195	VAL	2.8
1	C	199	ASP	2.8
1	C	795	SER	2.7
1	C	996	ALA	2.7
1	C	750	ASN	2.6
1	C	477	SER	2.6
1	C	994	ILE	2.5
1	C	747	ASP	2.5
1	C	189	ILE	2.5
2	B	91	LEU	2.5
1	C	32	LEU	2.4
1	C	194	THR	2.4
1	C	233	LEU	2.4
1	C	746	SER	2.4
1	C	864	PRO	2.3
1	C	490	VAL	2.3
1	C	232	VAL	2.3
1	C	238	PHE	2.3
1	C	235	PRO	2.2
1	C	798	ILE	2.2
1	C	193	GLY	2.1
1	C	481	ARG	2.1
1	C	951	GLY	2.1
1	C	248	GLU	2.1
1	C	993	ASP	2.1
1	C	919	TYR	2.1
1	C	105	TYR	2.0
1	C	221	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands

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
3	GOL	C	1107	6/6	0.68	0.15	53,64,75,75	0
3	GOL	C	1102	6/6	0.78	0.16	62,75,84,95	0
3	GOL	C	1105	6/6	0.81	0.13	44,53,62,63	0
3	GOL	C	1111	6/6	0.81	0.16	54,65,76,79	0
4	CL	C	1113	1/1	0.81	0.25	81,81,81,81	1
3	GOL	C	1110	6/6	0.82	0.18	49,59,65,65	0
3	GOL	B	201	6/6	0.84	0.13	41,52,63,67	0
3	GOL	C	1106	6/6	0.85	0.13	43,57,68,71	0
3	GOL	C	1109	6/6	0.85	0.21	44,56,75,75	0
3	GOL	C	1108	6/6	0.88	0.10	37,50,59,61	0
3	GOL	C	1103	6/6	0.88	0.11	33,50,60,72	0
3	GOL	C	1104	6/6	0.90	0.11	42,56,67,67	0
4	CL	C	1114	1/1	0.92	0.15	64,64,64,64	0
4	CL	C	1115	1/1	0.92	0.08	53,53,53,53	0
3	GOL	C	1101	6/6	0.95	0.10	27,33,37,38	0
4	CL	C	1112	1/1	0.96	0.05	35,35,35,35	0

6.5 Other polymers

There are no such residues in this entry.