



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 8, 2024 – 03:55 AM EDT

PDB ID : 3PHU
Title : OTU Domain of Crimean Congo Hemorrhagic Fever Virus
Authors : Akutsu, M.; Ye, Y.; Virdee, S.; Komander, D.
Deposited on : 2010-11-04
Resolution : 2.20 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

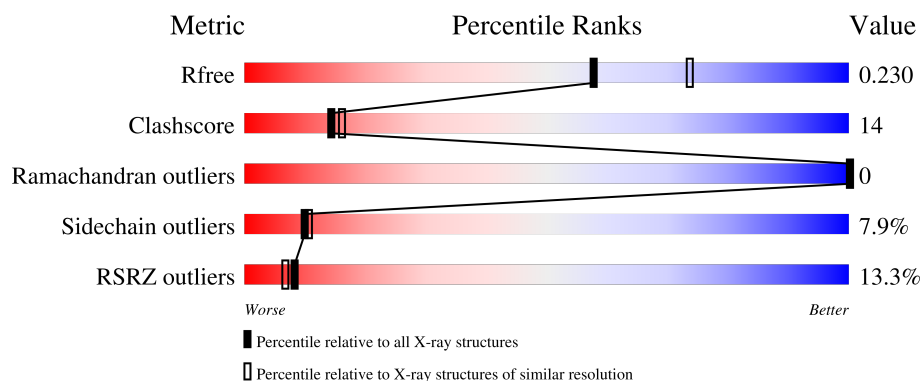
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
R_{free}	164625	5791 (2.20-2.20)
Clashscore	180529	6634 (2.20-2.20)
Ramachandran outliers	177936	6560 (2.20-2.20)
Sidechain outliers	177891	6561 (2.20-2.20)
RSRZ outliers	164620	5791 (2.20-2.20)

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 $\leq 5\%$. 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	219	<div> <div>5%</div> <div>67%</div> <div>14%</div> <div>•</div> <div>16%</div> </div>
1	B	219	<div> <div>17%</div> <div>58%</div> <div>23%</div> <div>•</div> <div>15%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3091 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 RNA-directed RNA polymerase L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	184	Total	C	N	O	S	0	0	0
			1475	933	246	289	7			
1	B	186	Total	C	N	O	S	0	0	0
			1467	927	240	294	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q6TQR6
A	0	PRO	-	expression tag	UNP Q6TQR6
B	-1	GLY	-	expression tag	UNP Q6TQR6
B	0	PRO	-	expression tag	UNP Q6TQR6

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

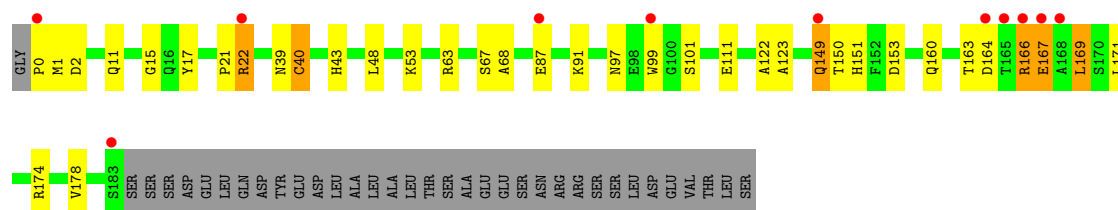
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	91	Total	O	0	0
			91	91		
3	B	52	Total	O	0	0
			52	52		

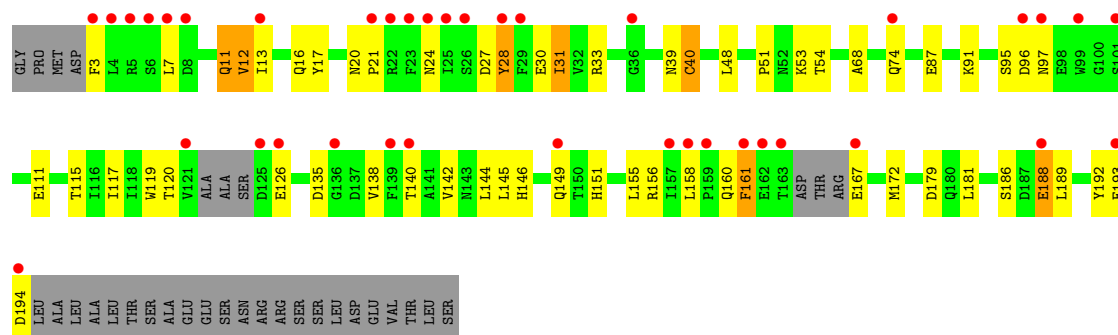
3 Residue-property plots [i](#)

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: RNA-directed RNA polymerase L



• Molecule 1: RNA-directed RNA polymerase L



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	115.99Å 115.99Å 94.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.38 – 2.20 49.38 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.6 (49.38-2.20) 99.7 (49.38-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.4_122)	Depositor
R, R_{free}	0.194 , 0.231 0.193 , 0.230	Depositor DCC
R_{free} test set	1870 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	40.1	Xtriage
Anisotropy	0.353	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3091	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 4.01% 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, CSD

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	0.68	0/1499	0.77	0/2032
1	B	0.61	0/1488	0.69	0/2018
All	All	0.65	0/2987	0.73	0/4050

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	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	A	149	GLN	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	1475	0	1425	37	0
1	B	1467	0	1380	44	0
2	A	6	0	8	0	0
3	A	91	0	0	2	0
3	B	52	0	0	3	0
All	All	3091	0	2813	80	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 14.

All (80) 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:48:LEU:HB3	1:A:166:ARG:NH2	1.49	1.25
1:A:48:LEU:CB	1:A:166:ARG:HH22	1.62	1.12
1:A:48:LEU:HB3	1:A:166:ARG:HH22	0.86	1.00
1:B:31:ILE:HD11	1:B:33:ARG:HD3	1.43	0.99
1:B:39:ASN:HD21	1:B:97:ASN:H	1.21	0.89
1:B:31:ILE:CD1	1:B:33:ARG:HD3	2.06	0.85
1:A:21:PRO:O	1:A:22:ARG:HD2	1.79	0.82
1:A:11:GLN:NE2	1:A:15:GLY:H	1.78	0.80
1:A:11:GLN:HE21	1:A:15:GLY:H	1.26	0.80
1:A:39:ASN:HD21	1:A:97:ASN:H	1.30	0.77
1:B:186:SER:OG	1:B:188:GLU:HG3	1.85	0.76
1:A:48:LEU:O	1:A:166:ARG:NH1	2.20	0.75
1:A:174:ARG:O	1:A:178:VAL:HG23	1.88	0.73
1:B:27:ASP:HB3	1:B:28:TYR:CD2	2.26	0.71
1:B:74:GLN:NE2	3:B:227:HOH:O	2.24	0.69
1:B:160:GLN:HB2	1:B:161:PHE:CE2	2.28	0.69
1:B:31:ILE:HD11	1:B:33:ARG:CD	2.21	0.68
1:B:24:ASN:O	1:B:27:ASP:HB2	1.94	0.67
1:A:166:ARG:NE	3:A:291:HOH:O	2.28	0.67
1:A:167:GLU:H	1:A:167:GLU:CD	1.98	0.66
1:A:48:LEU:CA	1:A:166:ARG:HH22	2.09	0.65
1:A:48:LEU:O	1:A:166:ARG:CZ	2.45	0.65
1:A:68:ALA:HB2	1:A:111:GLU:HB2	1.77	0.64
1:B:179:ASP:OD1	3:B:253:HOH:O	2.15	0.63
1:B:27:ASP:HB3	1:B:28:TYR:HD2	1.66	0.61
1:A:151:HIS:HE1	1:A:153:ASP:OD1	1.82	0.61
1:B:20:ASN:N	1:B:21:PRO:HD3	2.16	0.60
1:B:158:LEU:O	1:B:161:PHE:HD2	1.84	0.59
1:A:48:LEU:O	1:A:166:ARG:NH2	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:0:PRO:HG2	1:A:1:MET:H	1.69	0.57
1:B:115:THR:OG1	1:B:135:ASP:O	2.21	0.57
1:B:53:LYS:HD3	1:B:54:THR:O	2.05	0.56
1:B:161:PHE:N	1:B:161:PHE:CD2	2.74	0.56
1:A:11:GLN:HE21	1:A:15:GLY:N	1.99	0.56
1:A:21:PRO:C	1:A:22:ARG:HD2	2.25	0.55
1:B:117:ILE:HG21	1:B:119:TRP:CE2	2.42	0.55
1:A:22:ARG:HG2	1:A:22:ARG:HH11	1.72	0.55
1:B:3:PHE:N	1:B:28:TYR:HH	2.06	0.54
1:B:28:TYR:CD2	1:B:28:TYR:N	2.76	0.53
1:A:22:ARG:HH11	1:A:22:ARG:CG	2.21	0.53
1:B:39:ASN:HD21	1:B:97:ASN:N	2.00	0.52
1:B:68:ALA:HB2	1:B:111:GLU:HB2	1.92	0.52
1:B:172:MET:HE2	1:B:192:TYR:CE1	2.45	0.52
1:B:51:PRO:HD2	3:B:259:HOH:O	2.09	0.52
1:A:87:GLU:O	1:A:91:LYS:HG3	2.10	0.51
1:A:169:LEU:HD23	1:A:174:ARG:HG2	1.93	0.50
1:B:28:TYR:HD2	1:B:28:TYR:N	2.10	0.49
1:A:48:LEU:CB	1:A:166:ARG:NH2	2.38	0.49
1:A:101:SER:HB3	3:A:308:HOH:O	2.13	0.49
1:A:17:TYR:N	1:A:17:TYR:CD2	2.82	0.47
1:A:0:PRO:HD2	1:A:2:ASP:HB2	1.96	0.47
1:A:68:ALA:CB	1:A:111:GLU:HB2	2.44	0.47
1:B:7:LEU:HA	1:B:7:LEU:HD23	1.56	0.47
1:B:13:ILE:HD11	1:B:16:GLN:NE2	2.31	0.47
1:A:68:ALA:HB2	1:A:111:GLU:CB	2.44	0.46
1:B:27:ASP:HB3	1:B:28:TYR:CE2	2.50	0.46
1:B:12:VAL:HG23	1:B:13:ILE:HG23	1.97	0.46
1:A:163:THR:OG1	1:A:164:ASP:N	2.48	0.46
1:B:146:HIS:NE2	1:B:149:GLN:HA	2.31	0.45
1:A:0:PRO:CG	1:A:1:MET:H	2.28	0.45
1:B:181:LEU:C	1:B:181:LEU:HD12	2.37	0.45
1:B:40:CSD:OD2	1:B:151:HIS:HA	2.16	0.45
1:B:172:MET:CE	1:B:192:TYR:CE1	2.99	0.45
1:B:158:LEU:O	1:B:161:PHE:CD2	2.67	0.44
1:A:40:CSD:HA	1:A:43:HIS:HB2	2.00	0.44
1:B:87:GLU:O	1:B:91:LYS:HG3	2.18	0.44
1:B:48:LEU:HD12	1:B:142:VAL:HG13	1.99	0.43
1:A:99:TRP:CD1	1:A:99:TRP:N	2.84	0.43
1:B:11:GLN:HE21	1:B:17:TYR:HE1	1.66	0.43
1:B:87:GLU:H	1:B:87:GLU:CD	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:GLN:HA	1:B:74:GLN:OE1	2.20	0.42
1:B:193:GLU:O	1:B:194:ASP:CB	2.66	0.42
1:A:122:ALA:O	1:A:123:ALA:HB3	2.20	0.41
1:B:126:GLU:H	1:B:126:GLU:HG2	1.49	0.41
1:B:13:ILE:HD11	1:B:16:GLN:HE21	1.85	0.41
1:A:67:SER:OG	1:A:111:GLU:OE2	2.20	0.41
1:B:20:ASN:N	1:B:21:PRO:CD	2.84	0.40
1:B:145:LEU:HB2	1:B:155:LEU:HD21	2.03	0.40
1:A:0:PRO:CG	1:A:1:MET:N	2.85	0.40
1:A:11:GLN:HG2	1:B:189:LEU:HD23	2.04	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	181/219 (83%)	172 (95%)	9 (5%)	0	100	100
1	B	179/219 (82%)	164 (92%)	15 (8%)	0	100	100
All	All	360/438 (82%)	336 (93%)	24 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	A	159/190 (84%)	149 (94%)	10 (6%)	15	17
1	B	156/190 (82%)	141 (90%)	15 (10%)	7	6
All	All	315/380 (83%)	290 (92%)	25 (8%)	10	11

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ARG
1	A	53	LYS
1	A	63	ARG
1	A	149	GLN
1	A	150	THR
1	A	160	GLN
1	A	166	ARG
1	A	167	GLU
1	A	169	LEU
1	A	171	LEU
1	B	11	GLN
1	B	12	VAL
1	B	28	TYR
1	B	30	GLU
1	B	31	ILE
1	B	95	SER
1	B	96	ASP
1	B	120	THR
1	B	138	VAL
1	B	140	THR
1	B	144	LEU
1	B	156	ARG
1	B	161	PHE
1	B	167	GLU
1	B	188	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	16	GLN
1	A	39	ASN
1	A	97	ASN
1	A	151	HIS

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Mol	Chain	Res	Type
1	B	11	GLN
1	B	16	GLN
1	B	39	ASN
1	B	59	HIS
1	B	97	ASN

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	CSD	B	40	1	4,7,8	2.12	2 (50%)	1,8,10	1.10	0
1	CSD	A	40	1	4,7,8	3.63	1 (25%)	1,8,10	1.90	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
1	CSD	B	40	1	-	2/2/6/8	-
1	CSD	A	40	1	-	0/2/6/8	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	40	CSD	OD1-SG	-6.81	1.41	1.47
1	B	40	CSD	OD1-SG	-3.30	1.44	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	40	CSD	CB-SG	-2.09	1.67	1.79

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	40	CSD	N-CA-CB-SG
1	B	40	CSD	CA-CB-SG-OD1

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	40	CSD	1	0
1	A	40	CSD	1	0

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	A	218	-	5,5,5	0.40	0	5,5,5	1.78	1 (20%)

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	218	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	218	GOL	O1-C1-C2	-3.54	94.42	110.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	218	GOL	O1-C1-C2-C3
2	A	218	GOL	O1-C1-C2-O2

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	183/219 (83%)	0.03	11 (6%) 29 26	24, 40, 76, 108	0
1	B	185/219 (84%)	0.91	38 (20%) 3 3	27, 58, 103, 128	0
All	All	368/438 (84%)	0.47	49 (13%) 8 7	24, 47, 100, 128	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	3	PHE	7.9
1	B	163	THR	6.5
1	A	0	PRO	5.7
1	B	4	LEU	5.3
1	B	194	ASP	5.0
1	B	5	ARG	5.0
1	B	121	VAL	4.7
1	B	7	LEU	4.7
1	A	166	ARG	3.7
1	B	162	GLU	3.7
1	B	149	GLN	3.7
1	A	165	THR	3.6
1	B	167	GLU	3.5
1	B	6	SER	3.4
1	B	22	ARG	3.4
1	A	183	SER	3.4
1	B	96	ASP	3.4
1	B	161	PHE	3.3
1	B	101	SER	3.2
1	B	28	TYR	3.1
1	B	140	THR	3.1
1	A	164	ASP	3.1
1	B	99	TRP	3.0
1	B	193	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	29	PHE	3.0
1	B	36	GLY	2.9
1	B	159	PRO	2.8
1	A	99	TRP	2.8
1	A	167	GLU	2.7
1	B	125	ASP	2.7
1	B	139	PHE	2.7
1	B	8	ASP	2.7
1	A	87	GLU	2.6
1	B	158	LEU	2.5
1	B	188	GLU	2.4
1	B	26	SER	2.4
1	B	126	GLU	2.4
1	B	24	ASN	2.4
1	A	22	ARG	2.3
1	B	97	ASN	2.3
1	B	21	PRO	2.2
1	B	13	ILE	2.2
1	B	25	ILE	2.2
1	A	168	ALA	2.2
1	B	23	PHE	2.2
1	B	157	ILE	2.1
1	A	149	GLN	2.1
1	B	136	GLY	2.1
1	B	74	GLN	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(\AA^2)	Q<0.9
1	CSD	B	40	8/9	0.95	0.09	32,45,69,71	0
1	CSD	A	40	8/9	0.99	0.05	24,28,38,42	0

6.3 Carbohydrates [i](#)

There are no monosaccharides 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(Å ²)	Q<0.9
2	GOL	A	218	6/6	0.83	0.18	19,48,61,65	0

6.5 Other polymers

There are no such residues in this entry.