



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:36 AM GMT

PDB ID : 2RE9
Title : Crystal structure of TL1A at 2.1 Å
Authors : Jin, T.C.; Guo, F.; Kim, S.; Howard, A.J.; Zhang, Y.Z.
Deposited on : 2007-09-25
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

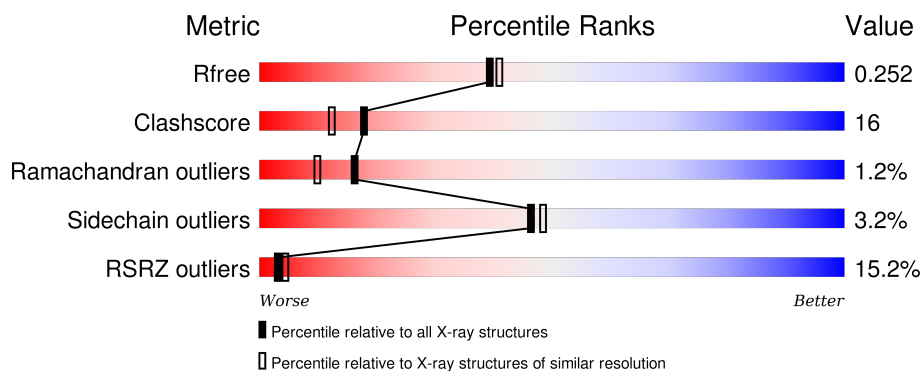
1 Overall quality at a glance

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}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	181	<div> <div>10%</div> <div> <div>66%</div> <div>22%</div> <div>• 9%</div> </div> </div>
1	B	181	<div> <div>17%</div> <div> <div>67%</div> <div>19%</div> <div>• • 9%</div> </div> </div>
1	C	181	<div> <div>14%</div> <div> <div>61%</div> <div>28%</div> <div>• 9%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MG	A	181	-	-	-	X
3	GOL	A	182	-	-	-	X
3	GOL	B	181	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4458 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 TNF superfamily ligand TL1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	164	Total	C	N	O	S	0	0	0
			1317	843	220	247	7			
1	B	164	Total	C	N	O	S	0	0	0
			1311	837	220	247	7			
1	C	164	Total	C	N	O	S	0	0	0
			1317	843	220	247	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	EXPRESSION TAG	UNP Q8NFE9
B	0	GLY	-	EXPRESSION TAG	UNP Q8NFE9
C	0	GLY	-	EXPRESSION TAG	UNP Q8NFE9

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).

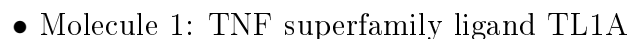


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	155	Total	O	0	0
			155	155		
4	B	185	Total	O	0	0
			185	185		
4	C	160	Total	O	0	0
			160	160		

- Molecule 1: TNF superfamily ligand TL1A



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	114.91Å 114.91Å 119.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.10 47.20 – 2.10	Depositor EDS
% Data completeness (in resolution range)	93.2 (50.00-2.10) 93.5 (47.20-2.10)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.10Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.198 , 0.253 0.199 , 0.252	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	38.5	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 60.0	EDS
Estimated twinning fraction	0.017 for -h,l,k 0.016 for -l,-k,-h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	2 of 46947 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4458	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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, MG

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	0.37	0/1350	0.66	0/1828
1	B	0.39	1/1343 (0.1%)	0.65	0/1819
1	C	0.35	0/1350	0.60	0/1828
All	All	0.37	1/4043 (0.0%)	0.64	0/5475

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	180	LEU	C-OXT	-5.12	1.13	1.23

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	A	1317	0	1292	40	0
1	B	1311	0	1285	39	0
1	C	1317	0	1292	52	0
2	A	1	0	0	0	0
3	A	6	0	8	0	0
3	B	6	0	8	1	0
4	A	155	0	0	2	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	185	0	0	1	0
4	C	160	0	0	3	0
All	All	4458	0	3885	124	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 16.

All (124) 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:114:THR:HG22	1:B:116:SER:H	1.37	0.88
1:B:168:THR:HG22	1:B:169:LYS:HG2	1.59	0.83
1:A:132:GLU:HG2	1:A:137:TRP:HB3	1.63	0.80
1:A:124:LEU:HD21	1:A:147:PHE:HE2	1.50	0.76
1:B:20:ASP:HB2	1:B:23:LYS:HE3	1.67	0.75
1:A:168:THR:HG22	1:A:169:LYS:HG2	1.69	0.73
1:A:172:LYS:HD3	1:B:125:MET:CE	2.19	0.73
1:C:93:GLU:HB2	1:C:95:ARG:HH12	1.54	0.71
1:C:45:ALA:HB3	1:C:66:LYS:HD3	1.74	0.70
1:C:95:ARG:HG3	1:C:100:PRO:HA	1.74	0.70
1:B:128:LYS:HD3	4:B:303:HOH:O	1.92	0.69
1:A:172:LYS:HD3	1:B:125:MET:HE2	1.74	0.69
1:C:111:THR:HG21	1:C:120:PRO:HB2	1.76	0.68
1:A:124:LEU:HD21	1:A:147:PHE:CE2	2.29	0.67
1:C:111:THR:CG2	1:C:120:PRO:HB2	2.26	0.66
1:A:94:ILE:HA	1:A:101:ASN:HD21	1.62	0.64
1:A:128:LYS:HG2	1:C:140:PRO:HD2	1.81	0.62
1:A:59:ASN:O	1:A:60:ARG:HB2	2.00	0.62
1:B:124:LEU:HD13	1:B:147:PHE:HZ	1.68	0.59
1:B:37:GLN:H	1:B:37:GLN:NE2	1.99	0.59
1:A:47:HIS:HB3	1:A:66:LYS:HE3	1.85	0.57
1:A:99:ARG:HD3	1:A:99:ARG:O	2.05	0.57
1:A:117:TYR:H	1:A:118:PRO:HA	1.71	0.56
1:C:118:PRO:HG2	1:C:119:GLU:H	1.71	0.55
1:B:114:THR:HG22	1:B:115:ASP:N	2.21	0.55
1:C:124:LEU:HD13	1:C:147:PHE:HZ	1.72	0.54
1:A:172:LYS:HD3	1:B:125:MET:HE1	1.89	0.54
1:B:111:THR:CG2	1:B:120:PRO:HB2	2.38	0.54
1:C:78:ILE:HD11	1:C:149:LEU:HD11	1.90	0.53
1:B:124:LEU:HD13	1:B:147:PHE:CZ	2.44	0.53
1:A:168:THR:HG21	4:A:279:HOH:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:32:ARG:HH21	1:C:32:ARG:HG2	1.74	0.52
1:B:47:HIS:HB3	1:B:66:LYS:HZ2	1.75	0.52
1:C:99:ARG:HH11	1:C:99:ARG:HG2	1.75	0.52
1:B:37:GLN:NE2	1:B:37:GLN:N	2.58	0.51
1:C:98:GLY:C	1:C:99:ARG:HD2	2.30	0.51
1:C:95:ARG:HD2	1:C:101:ASN:HD22	1.75	0.51
1:C:32:ARG:HD2	4:C:245:HOH:O	2.11	0.51
1:A:94:ILE:HA	1:A:101:ASN:ND2	2.25	0.51
1:C:59:ASN:O	1:C:60:ARG:HB2	2.11	0.51
1:A:28:LEU:HD21	1:A:68:LEU:HD13	1.93	0.50
1:C:95:ARG:HD2	1:C:101:ASN:ND2	2.26	0.49
1:C:90:GLU:H	1:C:90:GLU:CD	2.15	0.49
1:B:47:HIS:HB3	1:B:66:LYS:NZ	2.28	0.49
1:B:40:LYS:HB2	1:B:40:LYS:NZ	2.28	0.48
1:C:110:ILE:HD12	1:C:124:LEU:HD23	1.94	0.48
1:C:67:PHE:CD1	1:C:156:MET:HB3	2.48	0.48
1:C:109:VAL:CG2	1:C:158:ASN:HB2	2.44	0.48
1:C:29:THR:HG21	1:C:170:GLU:HB2	1.96	0.48
1:C:95:ARG:NE	1:C:95:ARG:HA	2.28	0.48
1:A:92:SER:HB3	1:A:132:GLU:O	2.13	0.48
1:A:65:ASN:O	1:A:66:LYS:HB2	2.15	0.47
1:B:90:GLU:N	1:B:90:GLU:OE1	2.48	0.47
1:C:96:GLN:HB2	1:C:99:ARG:HB2	1.96	0.47
1:B:37:GLN:H	1:B:37:GLN:HE21	1.59	0.47
1:B:140:PRO:HD2	1:C:128:LYS:HG2	1.96	0.47
1:B:43:PHE:CD1	1:B:160:SER:HA	2.50	0.47
1:C:20:ASP:HB2	1:C:23:LYS:HE3	1.97	0.47
1:C:168:THR:HG22	1:C:169:LYS:HE2	1.96	0.46
1:C:17:LEU:HD12	1:C:17:LEU:N	2.29	0.46
1:A:31:VAL:CG1	1:A:49:GLU:HB2	2.45	0.46
1:A:140:PRO:HD2	1:B:128:LYS:HE2	1.96	0.46
1:C:65:ASN:O	1:C:66:LYS:HB2	2.14	0.46
1:C:94:ILE:HG23	1:C:94:ILE:O	2.15	0.46
1:C:93:GLU:CB	1:C:95:ARG:HH12	2.24	0.46
1:C:109:VAL:HG12	1:C:125:MET:HG3	1.98	0.46
1:C:43:PHE:HB3	1:C:158:ASN:HB3	1.98	0.45
1:C:60:ARG:NH1	1:C:60:ARG:HB2	2.31	0.45
1:C:47:HIS:NE2	1:C:66:LYS:HE3	2.32	0.45
1:C:32:ARG:NH2	1:C:32:ARG:HG2	2.31	0.45
1:A:117:TYR:HB3	1:A:119:GLU:H	1.81	0.45
1:A:132:GLU:HG2	1:A:137:TRP:CB	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:THR:HG22	1:C:89:SER:O	2.17	0.44
1:C:93:GLU:HB2	1:C:95:ARG:NH1	2.25	0.44
1:C:131:CYS:HA	4:C:285:HOH:O	2.16	0.44
1:B:17:LEU:O	1:B:17:LEU:HD13	2.16	0.44
1:B:102:LYS:HD3	1:B:103:PRO:CD	2.47	0.44
1:C:106:ILE:HG13	1:C:141:ILE:HD11	2.00	0.44
1:A:62:ASN:HB3	1:A:69:LEU:HB3	2.00	0.44
1:B:92:SER:HB3	1:B:132:GLU:O	2.17	0.44
1:B:26:ALA:HB2	1:B:61:MET:HG2	2.00	0.44
1:A:26:ALA:HB2	1:A:61:MET:HG2	2.01	0.43
1:A:95:ARG:HH11	1:A:95:ARG:HG3	1.83	0.43
1:C:34:THR:O	1:C:34:THR:HG23	2.18	0.43
1:B:117:TYR:HD2	1:B:119:GLU:O	2.01	0.43
1:B:40:LYS:HZ3	1:B:40:LYS:HB2	1.84	0.43
1:A:30:VAL:HG22	1:A:31:VAL:N	2.34	0.43
1:B:17:LEU:HD21	4:C:327:HOH:O	2.19	0.43
1:A:111:THR:HG21	1:A:120:PRO:HB3	1.99	0.43
1:B:102:LYS:HD3	1:B:103:PRO:HD2	1.99	0.43
1:B:180:LEU:HD12	1:B:180:LEU:HA	1.85	0.43
1:B:41:ASN:HD22	1:B:41:ASN:N	2.17	0.43
1:B:59:ASN:O	1:B:60:ARG:HB2	2.19	0.43
1:A:119:GLU:HA	1:A:119:GLU:OE1	2.19	0.42
1:A:117:TYR:HB2	1:A:119:GLU:N	2.34	0.42
1:A:111:THR:CG2	1:A:120:PRO:HB3	2.50	0.42
1:B:58:LYS:HB3	3:B:181:GOL:H31	2.01	0.42
1:A:124:LEU:CD2	1:A:147:PHE:CE2	3.02	0.42
1:B:114:THR:CG2	1:B:115:ASP:N	2.82	0.42
1:A:117:TYR:CB	1:A:119:GLU:N	2.83	0.42
1:B:43:PHE:HB3	1:B:159:VAL:O	2.19	0.42
1:C:26:ALA:HB2	1:C:61:MET:HG2	2.02	0.42
1:C:18:ARG:HH21	1:C:18:ARG:HG3	1.84	0.41
1:B:30:VAL:HB	1:B:173:THR:HB	2.02	0.41
1:B:95:ARG:C	1:B:97:ALA:H	2.24	0.41
1:A:119:GLU:HA	1:A:120:PRO:HD3	1.88	0.41
1:A:180:LEU:HA	1:A:180:LEU:HD12	1.92	0.41
1:A:94:ILE:HG21	1:A:100:PRO:HA	2.01	0.41
1:A:109:VAL:HG11	1:A:122:GLN:HG2	2.02	0.41
1:A:72:GLU:OE2	1:C:18:ARG:NH2	2.53	0.41
1:B:20:ASP:O	1:B:23:LYS:HG3	2.21	0.41
1:C:123:LEU:HB3	1:C:147:PHE:CZ	2.55	0.41
1:B:111:THR:HG21	1:B:120:PRO:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99:ARG:HD2	1:C:99:ARG:N	2.36	0.41
1:C:90:GLU:N	1:C:90:GLU:CD	2.75	0.41
1:C:67:PHE:CE1	1:C:156:MET:HB3	2.55	0.41
1:C:31:VAL:CG1	1:C:49:GLU:HB2	2.50	0.41
1:A:95:ARG:HB2	4:A:314:HOH:O	2.21	0.40
1:C:99:ARG:NH1	1:C:99:ARG:HG2	2.36	0.40
1:C:18:ARG:HG3	1:C:18:ARG:NH2	2.37	0.40
1:A:23:LYS:HA	1:A:24:PRO:HD3	1.92	0.40
1:A:117:TYR:CB	1:A:119:GLU:H	2.35	0.40
1:C:117:TYR:HA	1:C:118:PRO:HD2	1.90	0.40
1:C:23:LYS:HA	1:C:24:PRO:HD3	1.98	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)
4:A:309:HOH:O	4:A:309:HOH:O[7_555]	2.00	0.20

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	162/181 (90%)	150 (93%)	8 (5%)	4 (2%)	7	2
1	B	162/181 (90%)	147 (91%)	14 (9%)	1 (1%)	30	24
1	C	162/181 (90%)	149 (92%)	12 (7%)	1 (1%)	30	24
All	All	486/543 (90%)	446 (92%)	34 (7%)	6 (1%)	16	10

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	96	GLN

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Mol	Chain	Res	Type
1	A	97	ALA
1	A	117	TYR
1	B	116	SER
1	A	89	SER
1	C	118	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	147/160 (92%)	143 (97%)	4 (3%)	52	56
1	B	146/160 (91%)	138 (94%)	8 (6%)	27	23
1	C	147/160 (92%)	145 (99%)	2 (1%)	74	80
All	All	440/480 (92%)	426 (97%)	14 (3%)	46	48

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	132	GLU
1	A	155	LEU
1	A	180	LEU
1	B	28	LEU
1	B	37	GLN
1	B	40	LYS
1	B	41	ASN
1	B	124	LEU
1	B	132	GLU
1	B	155	LEU
1	B	180	LEU
1	C	99	ARG
1	C	124	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	42	GLN
1	A	96	GLN
1	A	158	ASN
1	B	37	GLN
1	B	41	ASN
1	B	59	ASN
1	B	139	GLN
1	B	158	ASN
1	C	37	GLN
1	C	59	ASN
1	C	62	ASN
1	C	65	ASN
1	C	139	GLN
1	C	158	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	A	182	-	5,5,5	0.27	0	5,5,5	0.17	0
3	GOL	B	181	-	5,5,5	0.31	0	5,5,5	0.25	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	A	182	-	-	0/4/4/4	0/0/0/0
3	GOL	B	181	-	-	0/4/4/4	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	181	GOL	1	0

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	164/181 (90%)	0.75	19 (11%) 6 9	22, 35, 98, 119	1 (0%)
1	B	164/181 (90%)	1.10	31 (18%) 2 2	24, 36, 108, 122	1 (0%)
1	C	164/181 (90%)	1.15	25 (15%) 3 4	28, 46, 104, 127	0
All	All	492/543 (90%)	1.00	75 (15%) 3 4	22, 39, 104, 127	2 (0%)

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	97	ALA	14.5
1	A	97	ALA	12.2
1	C	94	ILE	11.8
1	A	17	LEU	11.5
1	A	96	GLN	10.9
1	C	98	GLY	10.7
1	C	17	LEU	10.5
1	B	39	PHE	10.0
1	B	94	ILE	9.9
1	B	97	ALA	9.0
1	B	95	ARG	9.0
1	B	38	HIS	8.6
1	C	96	GLN	8.5
1	B	117	TYR	8.2
1	C	95	ARG	8.0
1	A	117	TYR	7.9
1	B	17	LEU	7.4
1	B	96	GLN	7.4
1	B	41	ASN	6.4
1	A	94	ILE	6.3
1	B	91	CYS	6.0
1	A	98	GLY	5.9
1	C	21	GLY	5.8

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Mol	Chain	Res	Type	RSRZ
1	C	99	ARG	5.7
1	C	18	ARG	5.7
1	B	40	LYS	5.7
1	A	20	ASP	5.5
1	A	19	ALA	5.4
1	B	98	GLY	5.3
1	C	36	THR	5.1
1	B	99	ARG	5.1
1	A	95	ARG	5.0
1	B	118	PRO	4.9
1	A	100	PRO	4.9
1	C	92	SER	4.9
1	A	99	ARG	4.8
1	C	118	PRO	4.8
1	B	92	SER	4.8
1	B	37	GLN	4.6
1	C	101	ASN	4.4
1	B	90	GLU	4.3
1	B	93	GLU	4.2
1	C	35	PRO	4.1
1	A	115	ASP	4.0
1	C	20	ASP	3.9
1	B	36	THR	3.9
1	C	90	GLU	3.9
1	B	116	SER	3.8
1	B	42	GLN	3.5
1	B	35	PRO	3.4
1	C	93	GLU	3.3
1	C	89	SER	3.3
1	B	20	ASP	3.2
1	B	115	ASP	3.1
1	C	32	ARG	3.1
1	C	100	PRO	3.0
1	A	93	GLU	3.0
1	B	19	ALA	3.0
1	A	21	GLY	2.9
1	A	91	CYS	2.9
1	A	109	VAL	2.8
1	B	18	ARG	2.5
1	C	88	THR	2.5
1	C	22	ASP	2.5
1	A	18	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	100	PRO	2.5
1	B	119	GLU	2.4
1	C	37	GLN	2.4
1	B	180	LEU	2.4
1	B	133	VAL	2.3
1	C	19	ALA	2.3
1	C	91	CYS	2.3
1	A	102	LYS	2.1
1	B	109	VAL	2.1
1	A	119	GLU	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	GOL	B	181	6/6	0.22	0.46	21.46	84,84,84,85	0
2	MG	A	181	1/1	0.97	0.18	4.97	33,33,33,33	0
3	GOL	A	182	6/6	0.75	0.28	4.88	61,64,65,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.