



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3CLK / pdb_00003clk
Title : Crystal structure of a transcription regulator from *Lactobacillus plantarum*
Authors : Sugadev, R.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center
for Structural Genomics (NYSGXRC)
Deposited on : 2008-03-19
Resolution : 2.08 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

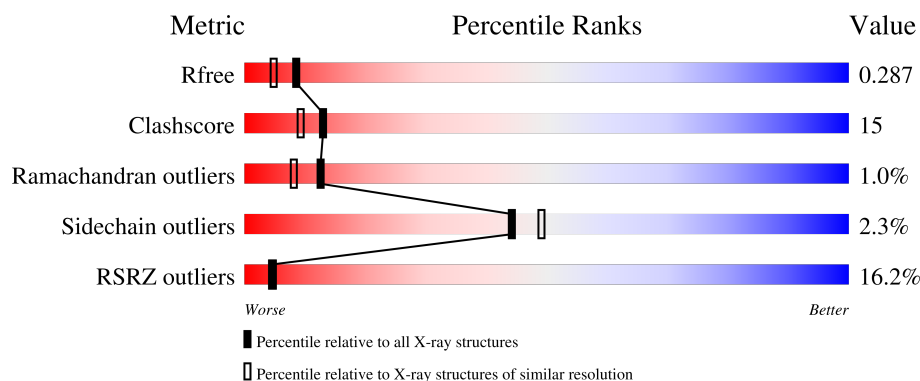
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	290	16% 62% 21% • 13%
1	B	290	12% 67% 21% • 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4153 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 Transcription regulator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	Se	0	0	0
			1932	1223	330	371	2	6			
1	B	262	Total	C	N	O	S	Se	0	0	0
			2025	1278	344	395	2	6			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	54	MSE	-	expression tag	UNP Q88S23
A	55	SER	-	expression tag	UNP Q88S23
A	56	LEU	-	expression tag	UNP Q88S23
A	336	GLU	-	expression tag	UNP Q88S23
A	337	GLY	-	expression tag	UNP Q88S23
A	338	HIS	-	expression tag	UNP Q88S23
A	339	HIS	-	expression tag	UNP Q88S23
A	340	HIS	-	expression tag	UNP Q88S23
A	341	HIS	-	expression tag	UNP Q88S23
A	342	HIS	-	expression tag	UNP Q88S23
A	343	HIS	-	expression tag	UNP Q88S23
B	54	MSE	-	expression tag	UNP Q88S23
B	55	SER	-	expression tag	UNP Q88S23
B	56	LEU	-	expression tag	UNP Q88S23
B	336	GLU	-	expression tag	UNP Q88S23
B	337	GLY	-	expression tag	UNP Q88S23
B	338	HIS	-	expression tag	UNP Q88S23
B	339	HIS	-	expression tag	UNP Q88S23
B	340	HIS	-	expression tag	UNP Q88S23
B	341	HIS	-	expression tag	UNP Q88S23
B	342	HIS	-	expression tag	UNP Q88S23
B	343	HIS	-	expression tag	UNP Q88S23

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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	83	Total	O	0	0
			83	83		
3	B	107	Total	O	0	0
			107	107		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.84Å 72.18Å 104.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.79 – 2.08 45.79 – 2.08	Depositor EDS
% Data completeness (in resolution range)	98.2 (45.79-2.08) 98.2 (45.79-2.08)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.17 (at 2.07Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.240 , 0.286 0.240 , 0.287	Depositor DCC
R_{free} test set	1962 reflections (5.82%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4153	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% 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

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	0.36	0/1952	0.95	9/2624 (0.3%)
1	B	0.38	0/2049	0.88	6/2760 (0.2%)
All	All	0.37	0/4001	0.91	15/5384 (0.3%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	189	TYR	CA-C-N	9.48	131.70	119.84
1	A	189	TYR	C-N-CA	9.48	131.70	119.84
1	A	289	ILE	N-CA-C	-8.64	94.63	107.51
1	B	289	ILE	N-CA-C	-7.32	96.61	107.51
1	A	191	TYR	N-CA-C	6.88	121.42	112.89
1	A	286	LEU	N-CA-C	6.41	119.06	109.25
1	B	286	LEU	N-CA-C	6.09	119.45	109.76
1	B	188	GLN	N-CA-C	5.94	119.63	112.38
1	B	287	THR	N-CA-C	-5.75	101.77	110.28
1	A	190	PRO	N-CA-C	5.71	124.23	112.47
1	B	184	ALA	N-CA-C	5.70	118.50	109.50
1	A	288	SER	N-CA-C	5.53	117.47	109.07
1	B	89	LYS	N-CA-C	-5.45	106.80	113.50
1	A	287	THR	N-CA-C	-5.30	102.44	110.28
1	A	179	ARG	N-CA-C	-5.02	108.89	114.62

There are no chirality outliers.

There are no planarity outliers.

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	1932	0	1951	60	0
1	B	2025	0	2026	62	0
2	A	6	0	8	0	0
3	A	83	0	0	8	0
3	B	107	0	0	4	0
All	All	4153	0	3985	119	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 15.

All (119) 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:248:MSE:SE	1:B:278:MSE:HE3	2.22	0.90
1:A:62:ASN:HA	1:A:120:MSE:HE3	1.56	0.88
1:A:112:LEU:O	1:A:116:GLU:HG2	1.74	0.86
1:A:111:LEU:HD12	1:A:111:LEU:H	1.45	0.81
1:A:67:VAL:O	1:A:127:ILE:HD11	1.84	0.78
1:A:133:ASN:HA	1:A:136:LEU:HD13	1.66	0.78
1:A:134:LEU:HD23	1:A:134:LEU:H	1.55	0.71
1:A:68:VAL:HG12	1:A:125:LEU:HB2	1.75	0.69
1:B:262:ILE:HG23	1:B:267:ASP:OD2	1.95	0.67
1:B:280:LYS:HE3	1:B:328:ARG:HD3	1.79	0.65
1:B:305:HIS:O	1:B:308:VAL:HG22	1.98	0.64
1:B:112:LEU:O	1:B:116:GLU:HB2	1.98	0.63
1:B:303:GLN:HB3	1:B:315:ILE:HD13	1.79	0.63
1:A:136:LEU:HD12	1:A:136:LEU:H	1.62	0.63
1:A:187:ASP:HB2	3:A:523:HOH:O	2.00	0.62
1:B:107:GLN:HE22	1:B:130:THR:H	1.47	0.61
1:A:158:ILE:HD13	1:A:299:THR:HG22	1.83	0.61
1:A:68:VAL:O	1:A:98:TYR:HA	2.00	0.61
1:B:235:LYS:HG3	1:B:260:PHE:HB3	1.81	0.61
1:B:251:ILE:HG12	1:B:282:THR:HG21	1.83	0.60
1:B:280:LYS:O	1:B:285:GLN:NE2	2.35	0.60
1:B:125:LEU:CD2	1:B:147:LEU:HD12	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:THR:HA	1:A:243:ILE:HD11	1.83	0.59
1:B:132:ASP:HA	1:B:135:GLN:HG2	1.84	0.59
1:B:190:PRO:HG2	3:B:345:HOH:O	2.02	0.58
1:B:280:LYS:HD3	1:B:285:GLN:NE2	2.18	0.58
1:B:254:LEU:HD23	1:B:254:LEU:O	2.05	0.57
3:A:573:HOH:O	1:B:283:ARG:HG3	2.05	0.57
1:B:280:LYS:HD3	1:B:285:GLN:CD	2.30	0.57
1:A:137:LEU:HD22	1:A:144:TYR:CD2	2.40	0.57
1:A:110:ALA:HB2	3:A:574:HOH:O	2.06	0.56
1:A:304:ILE:O	1:A:308:VAL:HG22	2.04	0.56
1:B:280:LYS:HE3	1:B:328:ARG:CD	2.36	0.56
1:A:136:LEU:HD12	1:A:136:LEU:N	2.21	0.55
1:B:149:MSE:HA	3:B:442:HOH:O	2.06	0.55
1:A:278:MSE:CE	1:B:278:MSE:HE1	2.37	0.55
1:B:273:ILE:O	1:B:274:ASP:HB2	2.05	0.55
1:A:62:ASN:HA	1:A:120:MSE:CE	2.33	0.55
1:B:283:ARG:NH2	1:B:285:GLN:HE21	2.06	0.54
1:B:223:TYR:HB2	1:B:248:MSE:HE3	1.90	0.54
1:B:191:TYR:O	1:B:195:LYS:HG2	2.08	0.54
1:A:64:ILE:HD13	1:A:304:ILE:HD12	1.90	0.54
1:B:170:THR:HA	1:B:243:ILE:HD11	1.90	0.53
1:B:107:GLN:NE2	1:B:130:THR:H	2.06	0.53
1:A:111:LEU:O	1:A:115:ILE:HG12	2.10	0.52
1:B:254:LEU:HD23	1:B:254:LEU:C	2.35	0.52
1:A:136:LEU:H	1:A:136:LEU:CD1	2.22	0.52
1:B:108:LYS:CB	1:B:133:ASN:HD21	2.23	0.52
1:B:283:ARG:NH2	1:B:285:GLN:NE2	2.57	0.52
1:B:273:ILE:O	1:B:274:ASP:CB	2.58	0.51
1:A:117:ARG:O	1:A:119:VAL:HG13	2.11	0.50
1:B:201:LYS:HD3	1:B:211:ILE:HD13	1.93	0.50
1:A:158:ILE:CD1	1:A:299:THR:HG22	2.42	0.50
1:A:278:MSE:HE1	1:B:278:MSE:HE1	1.93	0.49
1:B:125:LEU:O	1:B:127:ILE:HD12	2.12	0.49
1:B:280:LYS:HE3	1:B:328:ARG:HE	1.77	0.49
1:A:69:SER:O	1:A:70:SER:C	2.55	0.49
1:A:273:ILE:O	1:A:274:ASP:CB	2.61	0.49
1:B:205:LYS:HE3	1:B:211:ILE:HG13	1.95	0.48
1:A:265:PRO:HA	1:A:268:LEU:O	2.12	0.48
1:B:105:GLU:HA	1:B:105:GLU:OE1	2.13	0.48
1:B:125:LEU:HD21	1:B:147:LEU:HD12	1.95	0.48
1:A:111:LEU:H	1:A:111:LEU:CD1	2.19	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:MSE:SE	1:B:278:MSE:HE1	2.64	0.48
1:A:111:LEU:HD12	1:A:111:LEU:N	2.20	0.47
1:B:283:ARG:HH22	1:B:285:GLN:NE2	2.12	0.47
1:A:63:VAL:HG11	1:A:117:ARG:HH12	1.79	0.47
1:B:280:LYS:HE3	1:B:328:ARG:NE	2.29	0.47
1:B:132:ASP:O	1:B:135:GLN:HG3	2.15	0.47
1:A:67:VAL:HG12	1:A:127:ILE:HD13	1.97	0.46
1:B:108:LYS:HA	1:B:133:ASN:ND2	2.30	0.46
1:A:158:ILE:HD13	1:A:299:THR:CG2	2.45	0.46
1:A:251:ILE:HG12	1:A:282:THR:HG21	1.97	0.46
1:A:117:ARG:HH11	1:A:117:ARG:HG2	1.81	0.46
1:A:117:ARG:HD2	3:A:559:HOH:O	2.16	0.46
1:B:265:PRO:HA	1:B:268:LEU:O	2.15	0.46
1:B:188:GLN:HG3	1:B:197:LEU:HD22	1.97	0.46
1:A:69:SER:O	1:A:70:SER:O	2.34	0.45
1:A:115:ILE:HD12	1:A:140:SER:HB3	1.99	0.45
1:B:143:PRO:HG3	1:B:313:ASN:O	2.16	0.45
3:A:513:HOH:O	1:B:283:ARG:HD2	2.15	0.45
1:A:186:ILE:HG22	1:A:187:ASP:O	2.17	0.45
1:A:68:VAL:O	1:A:68:VAL:HG23	2.17	0.45
1:A:86:GLU:OE1	1:A:89:LYS:HD2	2.17	0.45
1:B:159:SER:O	1:B:320:PHE:HA	2.16	0.45
1:A:125:LEU:CD2	1:A:147:LEU:HD12	2.47	0.44
1:A:124:LEU:N	1:A:124:LEU:HD12	2.33	0.44
1:A:221:TYR:CD2	1:A:245:ALA:HB1	2.53	0.44
1:A:189:TYR:HA	1:A:190:PRO:HD3	1.74	0.44
1:B:79:ILE:HG21	1:B:125:LEU:HD22	1.98	0.44
1:A:163:GLU:HA	1:A:195:LYS:O	2.18	0.44
1:B:107:GLN:NE2	1:B:129:LEU:HA	2.32	0.44
1:B:157:PHE:C	1:B:157:PHE:CD1	2.95	0.43
1:A:133:ASN:N	3:A:538:HOH:O	2.51	0.43
1:B:316:VAL:HG12	3:B:409:HOH:O	2.17	0.43
1:A:143:PRO:HD2	3:A:563:HOH:O	2.17	0.43
1:A:172:LEU:O	1:A:172:LEU:HD23	2.19	0.43
1:A:149:MSE:N	1:A:159:SER:HB2	2.33	0.43
1:B:254:LEU:C	1:B:254:LEU:CD2	2.92	0.43
1:B:280:LYS:CE	1:B:328:ARG:HE	2.31	0.43
3:A:543:HOH:O	1:B:283:ARG:HG3	2.19	0.43
1:B:171:ASN:HD22	1:B:174:ILE:HD12	1.84	0.43
1:A:133:ASN:CA	1:A:136:LEU:HD13	2.43	0.42
1:B:108:LYS:HB2	1:B:133:ASN:HD21	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:VAL:C	1:A:127:ILE:HD11	2.42	0.42
1:A:308:VAL:O	1:A:308:VAL:HG23	2.20	0.42
1:B:75:PHE:O	1:B:79:ILE:HG13	2.20	0.42
1:B:172:LEU:C	1:B:172:LEU:HD23	2.45	0.42
1:B:136:LEU:HD13	1:B:136:LEU:C	2.45	0.42
1:B:266:LYS:HB2	3:B:428:HOH:O	2.20	0.42
1:A:71:VAL:HG11	1:A:75:PHE:CE2	2.56	0.41
1:A:74:ASN:OD1	1:A:74:ASN:N	2.44	0.41
1:A:247:ASP:HA	1:A:272:SER:OG	2.20	0.41
1:A:75:PHE:CG	1:A:76:ALA:N	2.89	0.41
1:A:273:ILE:O	1:A:274:ASP:CG	2.64	0.41
1:A:278:MSE:HE2	1:A:278:MSE:HA	2.03	0.41
1:A:288:SER:O	1:A:325:PRO:HA	2.20	0.41
1:B:107:GLN:HE22	1:B:130:THR:N	2.16	0.40
1:B:117:ARG:NH1	1:B:118:PRO:HD2	2.35	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	239/290 (82%)	227 (95%)	8 (3%)	4 (2%)	7	3
1	B	256/290 (88%)	248 (97%)	7 (3%)	1 (0%)	30	28
All	All	495/580 (85%)	475 (96%)	15 (3%)	5 (1%)	12	8

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	70	SER
1	A	139	SER
1	A	274	ASP

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Mol	Chain	Res	Type
1	B	274	ASP
1	A	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	212/239 (89%)	206 (97%)	6 (3%)	38	42
1	B	222/239 (93%)	218 (98%)	4 (2%)	51	58
All	All	434/478 (91%)	424 (98%)	10 (2%)	44	49

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	HIS
1	A	98	TYR
1	A	136	LEU
1	A	137	LEU
1	A	154	ASP
1	A	265	PRO
1	B	84	GLN
1	B	116	GLU
1	B	127	ILE
1	B	283	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	84	GLN
1	A	88	HIS
1	A	138	GLN
1	A	171	ASN
1	A	213	GLN

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Mol	Chain	Res	Type
1	A	236	ASN
1	A	285	GLN
1	A	295	GLN
1	B	74	ASN
1	B	78	GLN
1	B	84	GLN
1	B	93	ASN
1	B	107	GLN
1	B	133	ASN
1	B	138	GLN
1	B	171	ASN
1	B	236	ASN
1	B	256	GLN
1	B	285	GLN
1	B	302	GLN
1	B	313	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](#)

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	500	-	5,5,5	0.30	0	5,5,5	0.40	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	500	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	GOL	O1-C1-C2-C3
2	A	500	GOL	C1-C2-C3-O3
2	A	500	GOL	O1-C1-C2-O2
2	A	500	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	245/290 (84%)	0.94	47 (19%) 3 3	11, 32, 59, 65	0
1	B	256/290 (88%)	0.62	34 (13%) 7 7	12, 28, 53, 64	0
All	All	501/580 (86%)	0.78	81 (16%) 4 5	11, 29, 57, 65	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	98	TYR	5.7
1	A	191	TYR	5.4
1	A	138	GLN	5.1
1	A	189	TYR	4.8
1	A	98	TYR	4.7
1	A	140	SER	4.5
1	A	115	ILE	4.5
1	A	111	LEU	4.4
1	B	88	HIS	3.8
1	A	118	PRO	3.7
1	A	109	HIS	3.7
1	B	91	GLY	3.6
1	A	63	VAL	3.5
1	A	308	VAL	3.5
1	A	134	LEU	3.5
1	A	136	LEU	3.5
1	B	75	PHE	3.5
1	A	70	SER	3.4
1	A	117	ARG	3.3
1	A	77	GLN	3.3
1	A	68	VAL	3.2
1	A	307	SER	3.2
1	B	280	LYS	3.2
1	A	143	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	139	SER	3.1
1	B	117	ARG	3.0
1	B	135	GLN	2.9
1	B	69	SER	2.9
1	B	118	PRO	2.8
1	B	314	ARG	2.8
1	A	128	ALA	2.8
1	A	137	LEU	2.8
1	A	110	ALA	2.8
1	A	71	VAL	2.7
1	B	189	TYR	2.7
1	A	91	GLY	2.6
1	A	88	HIS	2.6
1	A	119	VAL	2.6
1	A	306	GLN	2.5
1	B	308	VAL	2.5
1	B	92	TYR	2.5
1	B	309	LYS	2.5
1	A	61	SER	2.5
1	B	109	HIS	2.5
1	A	211	ILE	2.5
1	B	74	ASN	2.4
1	B	313	ASN	2.4
1	A	190	PRO	2.4
1	A	309	LYS	2.4
1	B	311	GLY	2.4
1	A	87	ALA	2.4
1	B	136	LEU	2.4
1	B	78	GLN	2.4
1	B	141	ASP	2.4
1	B	119	VAL	2.4
1	B	84	GLN	2.3
1	B	90	ASN	2.3
1	A	116	GLU	2.3
1	B	267	ASP	2.3
1	B	139	SER	2.3
1	A	127	ILE	2.3
1	B	205	LYS	2.2
1	A	135	GLN	2.2
1	B	211	ILE	2.2
1	B	324	ASN	2.2
1	B	194	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	205	LYS	2.2
1	A	315	ILE	2.2
1	A	114	ALA	2.2
1	A	72	ARG	2.2
1	A	259	SER	2.2
1	A	112	LEU	2.1
1	A	154	ASP	2.1
1	A	317	SER	2.1
1	B	281	ILE	2.1
1	A	97	VAL	2.1
1	B	105	GLU	2.0
1	B	106	GLU	2.0
1	A	144	TYR	2.0
1	A	146	PHE	2.0
1	B	116	GLU	2.0

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 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	A	500	6/6	0.85	0.12	27,32,34,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.