



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 06:52 PM EDT

PDB ID : 6A4L / pdb_00006a4l
Title : AcrR from Mycobacterium tuberculosis
Authors : Kang, S.M.
Deposited on : 2018-06-20
Resolution : 2.80 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

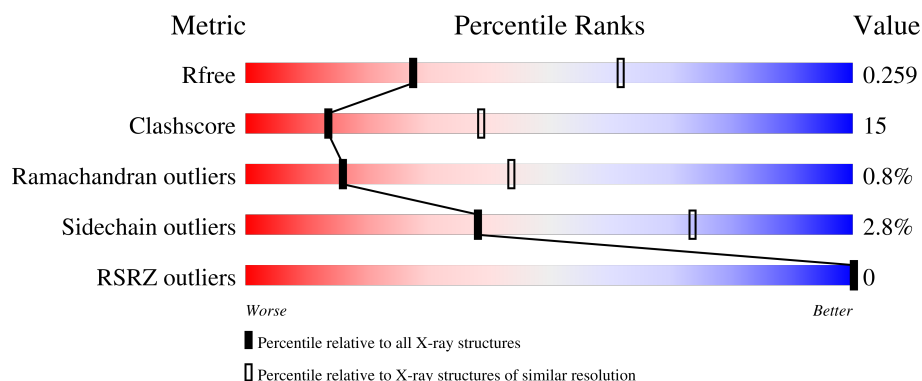
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	230	
1	B	230	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3211 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 TetR family transcriptional regulator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	Se	0	0	0
			1534	967	273	288	1	5			
1	B	198	Total	C	N	O	S	Se	0	0	0
			1556	979	278	293	1	5			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	initiating methionine	UNP A0A045J2D2
A	-18	GLY	-	expression tag	UNP A0A045J2D2
A	-17	SER	-	expression tag	UNP A0A045J2D2
A	-16	SER	-	expression tag	UNP A0A045J2D2
A	-15	HIS	-	expression tag	UNP A0A045J2D2
A	-14	HIS	-	expression tag	UNP A0A045J2D2
A	-13	HIS	-	expression tag	UNP A0A045J2D2
A	-12	HIS	-	expression tag	UNP A0A045J2D2
A	-11	HIS	-	expression tag	UNP A0A045J2D2
A	-10	HIS	-	expression tag	UNP A0A045J2D2
A	-9	SER	-	expression tag	UNP A0A045J2D2
A	-8	SER	-	expression tag	UNP A0A045J2D2
A	-7	GLY	-	expression tag	UNP A0A045J2D2
A	-6	LEU	-	expression tag	UNP A0A045J2D2
A	-5	VAL	-	expression tag	UNP A0A045J2D2
A	-4	PRO	-	expression tag	UNP A0A045J2D2
A	-3	ARG	-	expression tag	UNP A0A045J2D2
A	-2	GLY	-	expression tag	UNP A0A045J2D2
A	-1	SER	-	expression tag	UNP A0A045J2D2
A	0	HIS	-	expression tag	UNP A0A045J2D2
B	-19	MSE	-	initiating methionine	UNP A0A045J2D2
B	-18	GLY	-	expression tag	UNP A0A045J2D2
B	-17	SER	-	expression tag	UNP A0A045J2D2
B	-16	SER	-	expression tag	UNP A0A045J2D2
B	-15	HIS	-	expression tag	UNP A0A045J2D2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP A0A045J2D2
B	-13	HIS	-	expression tag	UNP A0A045J2D2
B	-12	HIS	-	expression tag	UNP A0A045J2D2
B	-11	HIS	-	expression tag	UNP A0A045J2D2
B	-10	HIS	-	expression tag	UNP A0A045J2D2
B	-9	SER	-	expression tag	UNP A0A045J2D2
B	-8	SER	-	expression tag	UNP A0A045J2D2
B	-7	GLY	-	expression tag	UNP A0A045J2D2
B	-6	LEU	-	expression tag	UNP A0A045J2D2
B	-5	VAL	-	expression tag	UNP A0A045J2D2
B	-4	PRO	-	expression tag	UNP A0A045J2D2
B	-3	ARG	-	expression tag	UNP A0A045J2D2
B	-2	GLY	-	expression tag	UNP A0A045J2D2
B	-1	SER	-	expression tag	UNP A0A045J2D2
B	0	HIS	-	expression tag	UNP A0A045J2D2

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		

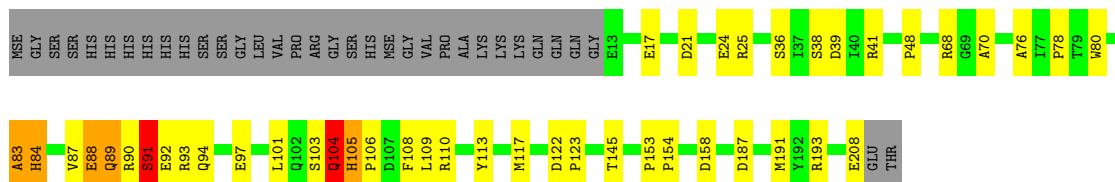
- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Na	0	0
			2	2		
3	B	1	Total	Na	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		
4	B	47	Total	O	0	0
			47	47		

- Molecule 1: TetR family transcriptional regulator



A83	A84	G85	G86	H87	E88	G89	R90	S91	E92	R93	G94	L95	T96	E97	L98	V99	S100	L101	G102	S103	G104	H105	P106	D107	F108	L109	R110	Q121	D122	V128	G174	V175	Y176	D184	D187	M191	T210							
SE	GLY	SE	SE	HIS	HIS	HIS	HIS	HIS	SE	SE	GLY	LEU	VAL	PRO	ARG	GLY	SE	HIS	MSE	GLY	VAL	PRO	ALA	LYS	LYS	GLN	GLN	GLY	R14	D21	R25	T29	G31	G32	A33	A34	R41	P48	K58	E59	A63	E67	Q71	F72

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	118.75Å 118.75Å 93.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.05 – 2.80 45.05 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (45.05-2.80) 95.8 (45.05-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.71 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.213 , 0.259 0.222 , 0.259	Depositor DCC
R_{free} test set	949 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	0.004	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 13.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.458 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3211	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.69% 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: SO4, NA

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	0.44	1/1560 (0.1%)	0.97	9/2109 (0.4%)
1	B	0.52	2/1582 (0.1%)	0.99	9/2138 (0.4%)
All	All	0.48	3/3142 (0.1%)	0.98	18/4247 (0.4%)

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	87	VAL	CA-CB	5.97	1.61	1.55
1	A	83	ALA	CA-C	5.51	1.58	1.53
1	B	86	PRO	CA-C	5.28	1.58	1.53

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	89	GLN	CA-CB-CG	9.01	132.13	114.10
1	B	86	PRO	CA-C-O	-7.35	112.21	123.16
1	A	105	HIS	CA-C-N	-7.17	110.88	119.84
1	A	105	HIS	C-N-CA	-7.17	110.88	119.84
1	B	85	GLY	N-CA-C	-7.03	97.99	112.34
1	A	88	GLU	CA-C-N	6.93	134.18	121.70
1	A	88	GLU	C-N-CA	6.93	134.18	121.70
1	B	30	LYS	N-CA-C	6.65	124.96	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	91	SER	N-CA-C	6.60	119.81	111.69
1	A	88	GLU	CA-CB-CG	6.01	126.12	114.10
1	A	89	GLN	CB-CA-C	6.00	121.50	110.10
1	A	89	GLN	CA-C-N	5.83	132.19	121.70
1	A	89	GLN	C-N-CA	5.83	132.19	121.70
1	B	88	GLU	N-CA-C	5.65	120.69	111.37
1	B	104	GLN	N-CA-C	5.40	117.78	111.02
1	B	109	LEU	N-CA-C	5.37	120.23	111.37
1	B	93	ARG	N-CA-C	-5.34	106.82	113.55
1	B	87	VAL	CB-CA-C	-5.29	106.41	111.44

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	104	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	1534	0	1521	49	0
1	B	1556	0	1545	44	0
2	A	10	0	0	1	0
2	B	15	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
4	A	46	0	0	4	0
4	B	47	0	0	4	0
All	All	3211	0	3066	90	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 (90) 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:88:GLU:HB2	1:A:89:GLN:HB3	1.17	1.14
1:B:29:THR:O	4:B:401:HOH:O	1.83	0.95
1:A:89:GLN:HG2	1:A:90:ARG:C	1.95	0.91
1:A:88:GLU:CB	1:A:89:GLN:HB3	2.01	0.89
1:A:83:ALA:HB1	1:A:84:HIS:HA	1.57	0.86
1:A:24:GLU:OE2	4:A:401:HOH:O	1.94	0.86
1:B:105:HIS:HB3	1:B:107:ASP:HB2	1.58	0.85
1:B:184:ASP:OD1	4:B:402:HOH:O	1.94	0.85
1:B:96:THR:O	1:B:100:SER:N	2.15	0.78
1:B:88:GLU:HG2	1:B:92:GLU:H	1.48	0.78
1:A:89:GLN:CG	1:A:90:ARG:HA	2.16	0.76
1:A:89:GLN:CD	1:A:90:ARG:HA	2.12	0.74
1:A:83:ALA:CB	1:A:84:HIS:HA	2.18	0.74
1:A:88:GLU:HB2	1:A:89:GLN:CB	2.08	0.74
1:B:174:GLY:O	4:B:404:HOH:O	2.06	0.73
1:A:208:GLU:O	4:A:402:HOH:O	2.06	0.73
1:A:17:GLU:OE1	1:A:68:ARG:NH1	2.23	0.72
1:B:31:GLY:N	4:B:401:HOH:O	2.21	0.72
1:B:84:HIS:CG	1:B:85:GLY:H	2.08	0.71
1:A:89:GLN:HG3	1:A:92:GLU:CB	2.21	0.70
1:A:87:VAL:HA	1:A:90:ARG:HG3	1.74	0.70
1:A:193:ARG:HG2	1:A:193:ARG:HH11	1.57	0.70
1:B:93:ARG:HA	1:B:96:THR:OG1	1.93	0.69
1:B:84:HIS:CG	1:B:85:GLY:N	2.58	0.68
1:B:93:ARG:O	1:B:97:GLU:N	2.15	0.68
1:A:104:GLN:HA	1:A:106:PRO:HD3	1.79	0.65
1:A:89:GLN:HG2	1:A:90:ARG:HA	1.79	0.64
1:A:89:GLN:HG2	1:A:91:SER:N	2.13	0.63
1:B:85:GLY:CA	1:B:90:ARG:HH21	2.12	0.63
1:A:89:GLN:HG2	1:A:90:ARG:CA	2.28	0.62
1:A:93:ARG:O	1:A:97:GLU:N	2.23	0.62
1:B:88:GLU:HG3	1:B:92:GLU:HB2	1.81	0.62
1:A:145:THR:O	4:A:405:HOH:O	2.16	0.62
1:B:88:GLU:HG2	1:B:92:GLU:N	2.15	0.61
1:A:89:GLN:HG3	1:A:92:GLU:CG	2.31	0.61
1:A:89:GLN:HG3	1:A:92:GLU:HG2	1.82	0.60
1:B:41:ARG:HD3	1:B:48:PRO:HD3	1.84	0.58
1:B:30:LYS:O	1:B:34:ALA:HB3	2.04	0.58
1:A:78:PRO:O	1:A:94:GLN:NE2	2.37	0.58
1:A:117:MSE:HE2	1:B:110:ARG:HG3	1.85	0.58
1:B:105:HIS:HB3	1:B:106:PRO:HA	1.86	0.57
1:A:110:ARG:NH1	2:A:301:SO4:O4	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:TRP:CZ2	1:A:90:ARG:HB2	2.41	0.56
1:A:113:TYR:O	1:A:117:MSE:HG3	2.06	0.55
1:A:193:ARG:HG2	1:A:193:ARG:NH1	2.21	0.54
1:A:89:GLN:HG3	1:A:92:GLU:HB2	1.90	0.53
1:B:106:PRO:HD3	1:B:110:ARG:HB3	1.91	0.53
1:B:106:PRO:CG	1:B:110:ARG:HB3	2.39	0.52
1:B:88:GLU:CG	1:B:92:GLU:HB2	2.39	0.52
1:A:70:ALA:HB2	1:A:108:PHE:CE1	2.45	0.52
1:B:85:GLY:HA2	1:B:90:ARG:HH21	1.74	0.51
1:A:104:GLN:CA	1:A:106:PRO:HD3	2.42	0.50
1:B:106:PRO:HG3	1:B:110:ARG:HB3	1.93	0.50
1:A:89:GLN:CG	1:A:92:GLU:HG2	2.42	0.50
1:A:88:GLU:CA	1:A:89:GLN:HB3	2.43	0.49
1:A:117:MSE:HE1	1:B:110:ARG:HA	1.95	0.49
1:B:110:ARG:HB2	1:B:176:TYR:OH	2.13	0.49
1:A:80:TRP:CH2	1:A:90:ARG:HB2	2.48	0.48
1:A:89:GLN:HG3	1:A:92:GLU:N	2.28	0.48
1:A:87:VAL:HG12	1:A:89:GLN:O	2.14	0.48
1:A:158:ASP:O	4:A:406:HOH:O	2.20	0.48
1:B:88:GLU:N	1:B:89:GLN:O	2.47	0.47
1:B:83:ALA:HB1	1:B:93:ARG:NH2	2.30	0.47
1:B:84:HIS:ND1	1:B:85:GLY:O	2.46	0.47
1:A:41:ARG:HD3	1:A:48:PRO:HD3	1.96	0.47
1:B:91:SER:O	1:B:95:LEU:N	2.33	0.47
1:B:83:ALA:O	1:B:90:ARG:NE	2.48	0.46
1:B:121:GLN:HG3	1:B:122:ASP:H	1.80	0.46
1:A:38:SER:OG	1:A:41:ARG:NH2	2.49	0.45
1:B:84:HIS:CE1	1:B:87:VAL:HG22	2.52	0.45
1:B:187:ASP:O	1:B:191:MSE:HG3	2.17	0.45
1:B:102:GLN:HA	1:B:104:GLN:H	1.82	0.44
1:A:21:ASP:O	1:A:25:ARG:HG3	2.17	0.44
1:B:88:GLU:CG	1:B:92:GLU:H	2.24	0.43
1:A:153:PRO:HA	1:A:154:PRO:HD3	1.89	0.43
1:B:73:PHE:CD2	1:B:108:PHE:HE1	2.36	0.43
1:B:59:GLU:H	1:B:59:GLU:CD	2.26	0.43
1:B:106:PRO:CD	1:B:110:ARG:HB3	2.48	0.43
1:A:36:SER:HB2	1:A:39:ASP:H	1.83	0.43
1:B:96:THR:HA	1:B:99:VAL:HB	2.01	0.43
1:A:106:PRO:O	1:A:109:LEU:N	2.33	0.42
1:B:32:TYR:CZ	1:B:58:LYS:HE2	2.55	0.42
1:B:21:ASP:O	1:B:25:ARG:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:GLU:O	1:B:71:GLN:HG3	2.21	0.41
1:A:76:ALA:HB3	1:A:101:LEU:HD11	2.02	0.41
1:A:117:MSE:HE3	1:B:176:TYR:HE2	1.85	0.40
1:A:122:ASP:HA	1:A:123:PRO:HD3	1.93	0.40
1:A:103:SER:O	1:A:103:SER:OG	2.33	0.40
1:A:187:ASP:O	1:A:191:MSE:HG3	2.21	0.40
1:B:63:ALA:HB1	1:B:128:VAL:HG21	2.03	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	194/230 (84%)	178 (92%)	14 (7%)	2 (1%)	12	38
1	B	196/230 (85%)	182 (93%)	13 (7%)	1 (0%)	24	55
All	All	390/460 (85%)	360 (92%)	27 (7%)	3 (1%)	16	44

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	GLN
1	A	91	SER
1	B	89	GLN

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	161/183 (88%)	158 (98%)	3 (2%)	50	81
1	B	164/183 (90%)	158 (96%)	6 (4%)	30	65
All	All	325/366 (89%)	316 (97%)	9 (3%)	38	73

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	HIS
1	A	91	SER
1	A	105	HIS
1	B	14	ARG
1	B	59	GLU
1	B	88	GLU
1	B	89	GLN
1	B	90	ARG
1	B	104	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	HIS
1	B	146	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	$\# Z > 2$
2	SO4	B	302	-	4,4,4	0.26	0	6,6,6	0.20	0
2	SO4	B	303	-	4,4,4	0.23	0	6,6,6	0.11	0
2	SO4	B	301	-	4,4,4	0.25	0	6,6,6	0.14	0
2	SO4	A	302	-	4,4,4	0.24	0	6,6,6	0.13	0
2	SO4	A	301	-	4,4,4	0.25	0	6,6,6	0.12	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
2	A	301	SO4	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	191/230 (83%)	-1.44	0 100 100	14, 34, 97, 120	0
1	B	193/230 (83%)	-1.41	0 100 100	14, 35, 92, 124	0
All	All	384/460 (83%)	-1.43	0 100 100	14, 35, 94, 124	0

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
3	NA	B	304	1/1	0.97	0.05	45,45,45,45	0
3	NA	A	304	1/1	0.98	0.07	61,61,61,61	0
2	SO4	B	301	5/5	0.99	0.11	61,63,94,103	0
2	SO4	B	302	5/5	0.99	0.13	56,60,68,92	0
2	SO4	B	303	5/5	0.99	0.10	49,61,89,95	0
3	NA	A	303	1/1	0.99	0.04	33,33,33,33	0
2	SO4	A	301	5/5	0.99	0.09	71,75,102,118	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	302	5/5	0.99	0.09	39,60,77,104	0

6.5 Other polymers [i](#)

There are no such residues in this entry.