



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:38 AM UTC

PDB ID : 4YNA / pdb_00004yna
Title : Oxidized YfiR
Authors : Xu, M.; Jiang, T.
Deposited on : 2015-03-09
Resolution : 3.20 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	:	4-5-2 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

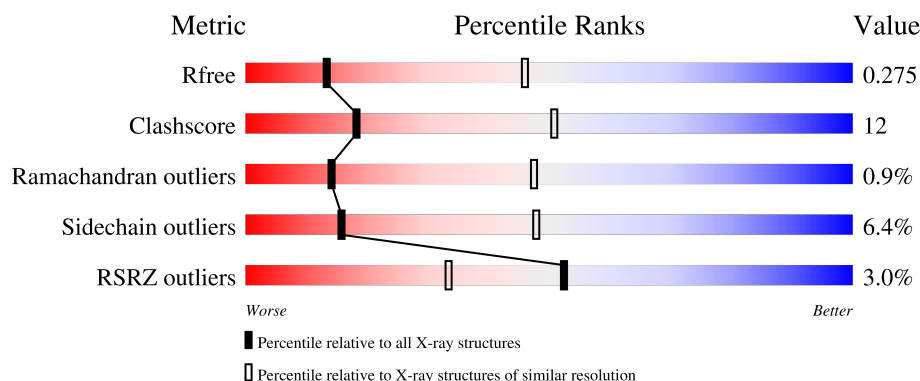
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	159	 72% 23% • •
1	B	159	 61% 26% 5% • 7%
1	C	159	 6% 69% 23% • 6%
1	D	159	 6% 67% 24% • 9%

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	SO4	B	202	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4594 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 YfiR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	0	0
			1179	732	226	215	6			
1	B	148	Total	C	N	O	S	0	0	0
			1140	709	217	208	6			
1	C	149	Total	C	N	O	S	0	0	0
			1149	714	219	210	6			
1	D	145	Total	C	N	O	S	0	0	0
			1116	694	211	205	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	32	GLY	-	expression tag	UNP Q9I4L4
A	33	ASN	-	expression tag	UNP Q9I4L4
A	34	SER	-	expression tag	UNP Q9I4L4
B	32	GLY	-	expression tag	UNP Q9I4L4
B	33	ASN	-	expression tag	UNP Q9I4L4
B	34	SER	-	expression tag	UNP Q9I4L4
C	32	GLY	-	expression tag	UNP Q9I4L4
C	33	ASN	-	expression tag	UNP Q9I4L4
C	34	SER	-	expression tag	UNP Q9I4L4
D	32	GLY	-	expression tag	UNP Q9I4L4
D	33	ASN	-	expression tag	UNP Q9I4L4
D	34	SER	-	expression tag	UNP Q9I4L4

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



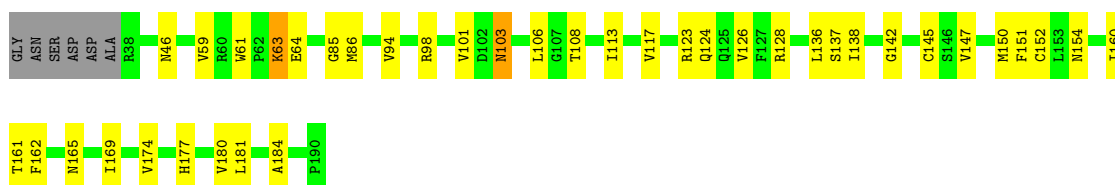
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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.

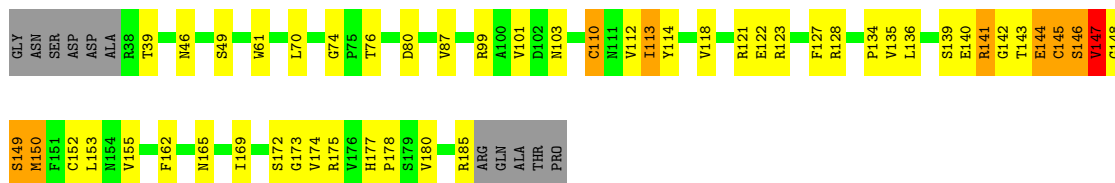
• Molecule 1: YfiR

Chain A: 



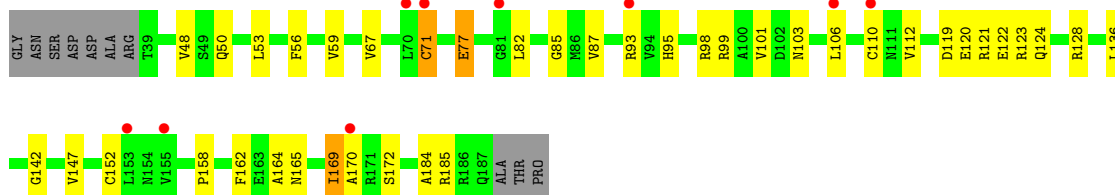
• Molecule 1: YfiR

Chain B: 



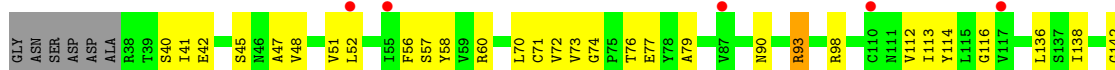
• Molecule 1: YfiR

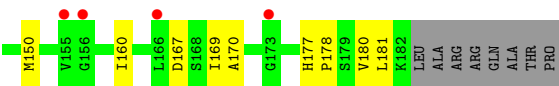
Chain C: 



• Molecule 1: YfiR

Chain D: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	119.67Å 119.67Å 85.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.88 – 3.20 34.88 – 3.20	Depositor EDS
% Data completeness (in resolution range)	86.7 (34.88-3.20) 86.9 (34.88-3.20)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.94 (at 3.18Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.246 , 0.281 0.243 , 0.275	Depositor DCC
R_{free} test set	450 reflections (4.17%)	wwPDB-VP
Wilson B-factor (Å ²)	60.8	Xtriage
Anisotropy	0.376	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 62.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4594	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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

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	0.84	0/1198	1.10	8/1623 (0.5%)
1	B	0.78	0/1158	1.02	2/1568 (0.1%)
1	C	0.66	4/1167 (0.3%)	0.86	0/1580
1	D	0.55	0/1134	0.82	0/1536
All	All	0.72	4/4657 (0.1%)	0.96	10/6307 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	71	CYS	CB-SG	7.61	2.06	1.81
1	C	169	ILE	CA-C	6.33	1.61	1.52
1	C	170	ALA	N-CA	5.93	1.53	1.46
1	C	170	ALA	CA-C	-5.79	1.45	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	110	CYS	N-CA-C	6.77	120.08	108.02
1	A	85	GLY	N-CA-C	-6.58	105.70	111.95
1	A	61	TRP	CA-C-N	6.47	125.97	119.24
1	A	61	TRP	C-N-CA	6.47	125.97	119.24
1	A	177	HIS	CA-C-N	6.39	126.61	119.32
1	A	177	HIS	C-N-CA	6.39	126.61	119.32
1	A	103	ASN	CA-C-N	5.84	126.35	120.04
1	A	103	ASN	C-N-CA	5.84	126.35	120.04
1	A	160	ILE	CB-CA-C	-5.55	104.59	111.08
1	B	147	VAL	N-CA-C	-5.15	102.21	109.63

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	1179	0	1194	21	0
1	B	1140	0	1154	46	0
1	C	1149	0	1162	30	0
1	D	1116	0	1125	25	0
2	B	10	0	0	3	0
All	All	4594	0	4635	112	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 12.

All (112) 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:71:CYS:CB	1:C:71:CYS:SG	2.06	1.42
1:B:114:TYR:OH	1:B:140:GLU:HG3	1.61	0.98
1:A:86:MET:HB2	1:A:94:VAL:HB	1.64	0.79
1:B:148:GLY:CA	1:C:93:ARG:HH21	1.98	0.75
1:B:123:ARG:HD2	1:B:147:VAL:HG11	1.69	0.75
1:B:142:GLY:HA2	1:B:152:CYS:HB3	1.70	0.72
1:B:150:MET:HE1	1:B:172:SER:OG	1.88	0.72
1:A:142:GLY:HA2	1:A:152:CYS:HB3	1.70	0.72
1:C:71:CYS:CB	1:C:110:CYS:SG	2.79	0.71
1:A:145:CYS:N	1:A:152:CYS:SG	2.64	0.70
1:C:121:ARG:HG3	1:C:122:GLU:N	2.07	0.68
1:B:123:ARG:CD	1:B:147:VAL:CG1	2.72	0.67
1:B:148:GLY:HA2	1:C:93:ARG:HH21	1.57	0.67
1:B:39:THR:HB	2:B:202:SO4:O1	1.95	0.67
1:B:123:ARG:HD3	1:B:147:VAL:CG1	2.25	0.66
1:D:71:CYS:O	1:D:114:TYR:N	2.28	0.66
1:A:138:ILE:HG12	1:A:151:PHE:HB2	1.78	0.65
1:B:103:ASN:ND2	2:B:201:SO4:O3	2.32	0.63
1:C:103:ASN:O	1:C:106:LEU:HG	2.00	0.62
1:B:123:ARG:HD3	1:B:147:VAL:HG13	1.82	0.62
1:B:165:ASN:O	1:B:169:ILE:HG13	2.00	0.61
1:D:112:VAL:HG22	1:D:136:LEU:HB3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:123:ARG:HD2	1:C:147:VAL:HB	1.84	0.59
1:C:119:ASP:OD1	1:C:121:ARG:HG2	2.03	0.59
1:C:164:ALA:HB1	1:C:169:ILE:HD11	1.86	0.58
1:B:74:GLY:O	1:B:76:THR:HG23	2.05	0.57
1:D:79:ALA:HB2	1:D:114:TYR:CE2	2.40	0.57
1:B:142:GLY:CA	1:B:152:CYS:HB3	2.33	0.57
1:B:123:ARG:HH11	1:B:123:ARG:HG3	1.69	0.57
1:A:124:GLN:HE21	1:A:128:ARG:HH12	1.53	0.56
1:B:150:MET:HE1	1:B:174:VAL:HG22	1.88	0.56
1:C:165:ASN:O	1:C:169:ILE:HG13	2.06	0.56
1:A:165:ASN:O	1:A:169:ILE:HG13	2.07	0.55
1:A:123:ARG:HD2	1:A:147:VAL:HB	1.89	0.54
1:C:59:VAL:HG11	1:C:136:LEU:HD22	1.90	0.54
1:C:142:GLY:HA2	1:C:152:CYS:HB3	1.89	0.54
1:B:127:PHE:CE1	1:B:148:GLY:HA3	2.44	0.53
1:B:148:GLY:HA2	1:C:93:ARG:NH2	2.23	0.53
1:B:144:GLU:H	1:B:144:GLU:CD	2.16	0.53
1:B:123:ARG:HD2	1:B:147:VAL:CG1	2.32	0.52
1:B:150:MET:HE1	1:B:172:SER:CB	2.39	0.52
1:D:76:THR:OG1	1:D:98:ARG:NH2	2.43	0.52
1:A:150:MET:HE3	1:A:169:ILE:HA	1.91	0.52
1:B:139:SER:OG	1:B:152:CYS:SG	2.67	0.52
1:C:162:PHE:CZ	1:C:184:ALA:HB2	2.45	0.52
1:B:150:MET:CE	1:B:172:SER:CB	2.89	0.51
1:C:124:GLN:HG2	1:C:128:ARG:NH1	2.25	0.51
1:C:169:ILE:O	1:C:172:SER:OG	2.22	0.51
1:C:98:ARG:CZ	1:D:77:GLU:HA	2.41	0.50
1:B:139:SER:HB3	1:B:149:SER:OG	2.11	0.50
1:D:73:VAL:O	1:D:116:GLY:N	2.42	0.50
1:B:112:VAL:HG13	1:B:136:LEU:HB3	1.94	0.49
1:D:56:PHE:CE1	1:D:112:VAL:HG21	2.47	0.49
1:B:153:LEU:HD23	1:B:162:PHE:HB3	1.94	0.49
1:A:59:VAL:HG21	1:A:136:LEU:HD22	1.95	0.48
1:B:128:ARG:HG2	1:C:95:HIS:HB2	1.93	0.48
1:B:146:SER:O	1:B:165:ASN:ND2	2.46	0.48
1:C:120:GLU:HG3	1:C:123:ARG:HH21	1.79	0.48
1:B:101:VAL:HG11	1:B:122:GLU:HB3	1.96	0.47
1:B:114:TYR:CZ	1:B:140:GLU:HG3	2.46	0.47
1:B:70:LEU:C	1:B:70:LEU:HD23	2.39	0.47
1:B:46:ASN:O	1:B:49:SER:HB2	2.14	0.47
1:B:121:ARG:NH1	1:D:42:GLU:OE2	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:101:VAL:HG11	1:C:122:GLU:HB3	1.97	0.46
1:B:110:CYS:O	1:B:135:VAL:HG22	2.16	0.46
1:D:72:VAL:HG13	1:D:76:THR:HG21	1.97	0.46
1:D:58:TYR:HB3	1:D:177:HIS:HB3	1.98	0.45
1:A:86:MET:HE3	1:A:94:VAL:HG11	1.99	0.45
1:B:150:MET:HE3	1:B:169:ILE:HA	1.97	0.45
1:D:52:LEU:HD11	1:D:138:ILE:HG21	1.99	0.45
1:D:180:VAL:HG13	1:D:181:LEU:HD23	1.99	0.45
1:D:150:MET:HB3	1:D:169:ILE:HG12	1.97	0.45
1:C:98:ARG:HB2	1:D:41:ILE:HG21	2.00	0.44
1:D:93:ARG:H	1:D:93:ARG:HG2	1.59	0.44
1:B:123:ARG:HH11	1:B:123:ARG:CG	2.30	0.44
1:C:77:GLU:HG3	1:D:74:GLY:C	2.43	0.44
1:A:63:LYS:HE2	1:A:64:GLU:N	2.33	0.44
1:B:113:ILE:HG22	1:B:135:VAL:HG12	1.99	0.43
1:B:128:ARG:HD2	1:C:95:HIS:CD2	2.53	0.43
1:C:48:VAL:HG12	1:C:82:LEU:HD21	2.00	0.43
1:A:154:ASN:HB3	1:A:161:THR:OG1	2.18	0.43
1:A:103:ASN:O	1:A:106:LEU:HG	2.19	0.43
1:A:98:ARG:HD2	1:B:80:ASP:OD2	2.18	0.43
1:B:123:ARG:CD	1:B:147:VAL:HG13	2.46	0.43
1:B:172:SER:OG	1:B:173:GLY:N	2.52	0.43
1:D:74:GLY:O	1:D:76:THR:HG23	2.19	0.43
1:A:162:PHE:CZ	1:A:184:ALA:HB2	2.53	0.42
1:D:51:VAL:HG21	1:D:160:ILE:HG21	2.01	0.42
1:A:123:ARG:HD2	1:A:147:VAL:CG2	2.49	0.42
1:C:119:ASP:OD1	1:C:121:ARG:CG	2.65	0.42
1:B:150:MET:CE	1:B:172:SER:OG	2.64	0.42
1:A:136:LEU:HD12	1:A:137:SER:N	2.34	0.42
1:C:121:ARG:CG	1:C:122:GLU:N	2.80	0.42
1:D:177:HIS:O	1:D:180:VAL:HG12	2.19	0.42
1:A:180:VAL:HG13	1:A:181:LEU:HD23	2.02	0.41
1:B:145:CYS:O	1:B:145:CYS:SG	2.74	0.41
1:D:177:HIS:HA	1:D:178:PRO:HD3	1.89	0.41
1:A:162:PHE:CE2	1:A:184:ALA:HB2	2.56	0.41
1:D:47:ALA:O	1:D:51:VAL:HG23	2.21	0.41
1:D:56:PHE:CZ	1:D:112:VAL:HG11	2.56	0.41
1:C:71:CYS:SG	1:C:71:CYS:CA	3.01	0.41
1:D:167:ASP:O	1:D:170:ALA:HB3	2.20	0.41
1:C:56:PHE:CZ	1:C:112:VAL:HG11	2.55	0.41
1:D:112:VAL:HA	1:D:136:LEU:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:ASN:ND2	2:B:202:SO4:O4	2.44	0.41
1:A:101:VAL:HG13	1:A:126:VAL:HG23	2.03	0.41
1:B:61:TRP:HD1	1:B:134:PRO:HB2	1.86	0.41
1:B:141:ARG:O	1:B:145:CYS:HB3	2.21	0.40
1:C:71:CYS:SG	1:C:99:ARG:HD2	2.60	0.40
1:C:50:GLN:HA	1:C:53:LEU:HD12	2.03	0.40
1:B:177:HIS:ND1	1:B:178:PRO:HD2	2.36	0.40
1:D:70:LEU:HD21	1:D:114:TYR:HB2	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	151/159 (95%)	146 (97%)	5 (3%)	0	100	100
1	B	146/159 (92%)	137 (94%)	7 (5%)	2 (1%)	9	39
1	C	147/159 (92%)	138 (94%)	7 (5%)	2 (1%)	9	39
1	D	143/159 (90%)	138 (96%)	4 (3%)	1 (1%)	18	52
All	All	587/636 (92%)	559 (95%)	23 (4%)	5 (1%)	14	47

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	146	SER
1	B	149	SER
1	C	85	GLY
1	D	142	GLY
1	C	158	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	129/133 (97%)	123 (95%)	6 (5%)	23	57
1	B	125/133 (94%)	111 (89%)	14 (11%)	6	25
1	C	126/133 (95%)	122 (97%)	4 (3%)	34	65
1	D	123/133 (92%)	115 (94%)	8 (6%)	15	47
All	All	503/532 (94%)	471 (94%)	32 (6%)	16	48

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	63	LYS
1	A	108	THR
1	A	113	ILE
1	A	117	VAL
1	A	174	VAL
1	B	87	VAL
1	B	99	ARG
1	B	113	ILE
1	B	118	VAL
1	B	141	ARG
1	B	143	THR
1	B	144	GLU
1	B	145	CYS
1	B	147	VAL
1	B	150	MET
1	B	155	VAL
1	B	175	ARG
1	B	180	VAL
1	B	185	ARG
1	C	67	VAL
1	C	77	GLU
1	C	87	VAL
1	C	185	ARG

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Mol	Chain	Res	Type
1	D	40	SER
1	D	45	SER
1	D	48	VAL
1	D	57	SER
1	D	60	ARG
1	D	90	ASN
1	D	93	ARG
1	D	113	ILE

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

Mol	Chain	Res	Type
1	A	124	GLN
1	A	133	HIS
1	B	90	ASN
1	B	124	GLN
1	B	133	HIS
1	C	43	GLN
1	C	69	GLN
1	C	165	ASN
1	D	90	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](#)

2 ligands 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	$\# Z > 2$
2	SO4	B	201	-	4,4,4	0.67	0	6,6,6	0.73	0
2	SO4	B	202	-	4,4,4	0.52	0	6,6,6	0.46	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.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	201	SO4	1	0
2	B	202	SO4	2	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	153/159 (96%)	-0.34	0	100	100	25, 41, 76, 123	0
1	B	148/159 (93%)	-0.09	0	100	100	28, 53, 93, 127	0
1	C	149/159 (93%)	0.79	9 (6%)	27	17	88, 138, 192, 217	0
1	D	145/159 (91%)	0.72	9 (6%)	26	17	82, 134, 175, 192	0
All	All	595/636 (93%)	0.26	18 (3%)	52	33	25, 98, 173, 217	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	71	CYS	4.2
1	C	110	CYS	3.8
1	D	117	VAL	3.4
1	D	166	LEU	2.8
1	C	170	ALA	2.6
1	C	153	LEU	2.6
1	D	55	ILE	2.5
1	D	87	VAL	2.4
1	D	110	CYS	2.4
1	C	106	LEU	2.3
1	D	156	GLY	2.2
1	D	52	LEU	2.2
1	C	70	LEU	2.1
1	C	93	ARG	2.1
1	D	173	GLY	2.1
1	C	155	VAL	2.0
1	D	155	VAL	2.0
1	C	81	GLY	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	SO4	B	201	5/5	0.93	0.09	54,56,57,59	0
2	SO4	B	202	5/5	0.97	0.05	54,55,58,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.