



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:39 PM GMT

PDB ID : 4YNA  
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](mailto: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.

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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

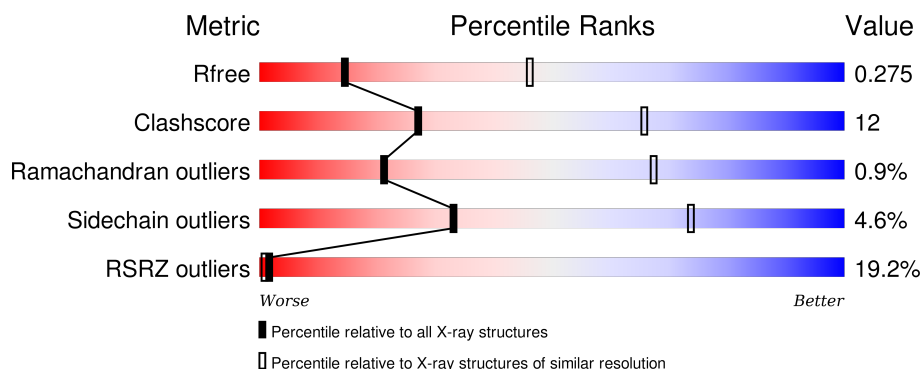
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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  $\geq 3$ , 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	159	<div> <div>2%</div> <div>75%</div> <div>20%</div> <div>• •</div> </div>
1	B	159	<div> <div>%</div> <div>62%</div> <div>26%</div> <div>5%</div> <div>7%</div> </div>
1	C	159	<div> <div>26%</div> <div>71%</div> <div>22%</div> <div>• 6%</div> </div>
1	D	159	<div> <div>43%</div> <div>68%</div> <div>23%</div> <div>• 9%</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 crite-

ria:

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.

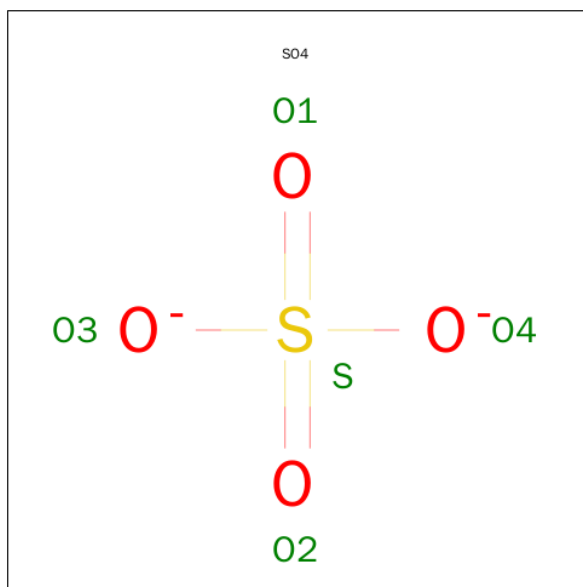
- Molecule 1 is a protein called YfiR.

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 (three-letter code: SO4) (formula: O<sub>4</sub>S).

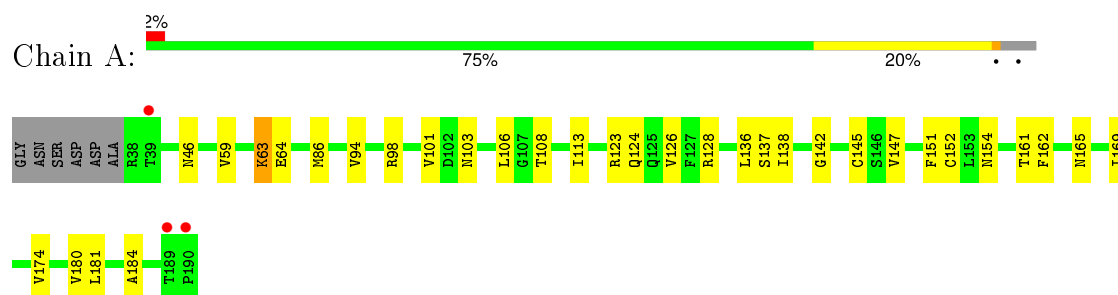


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

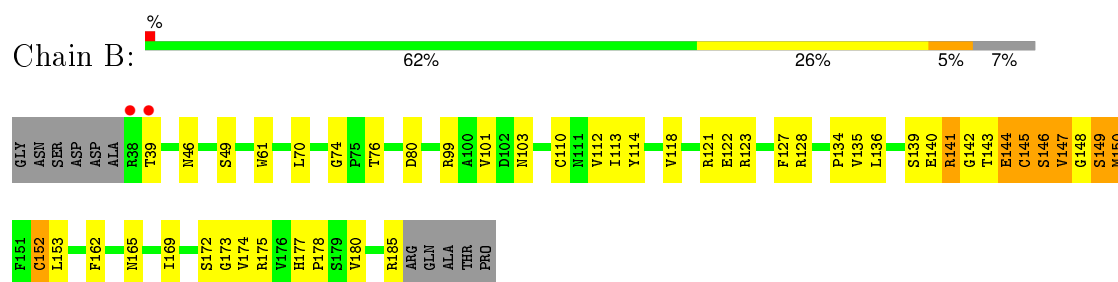
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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.

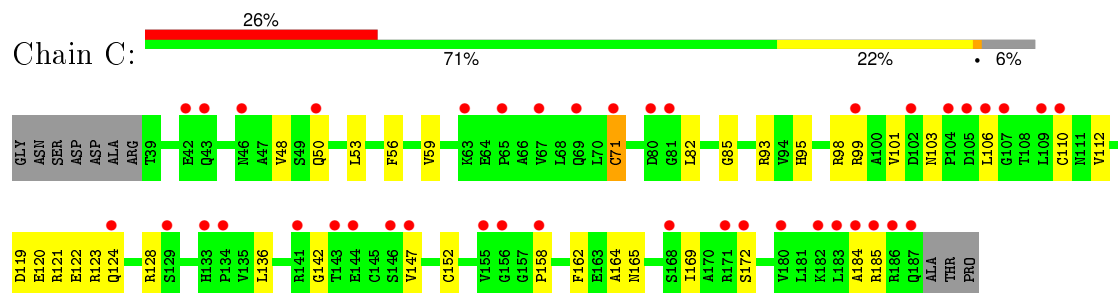
#### • Molecule 1: YfiR



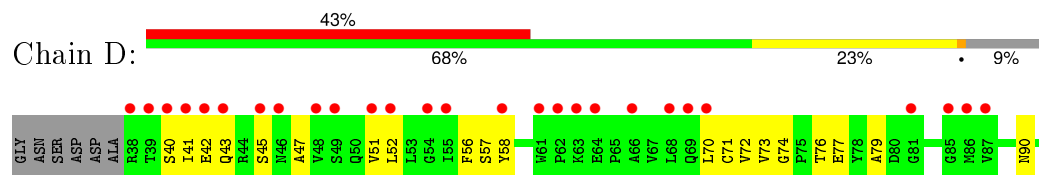
#### • Molecule 1: YfiR

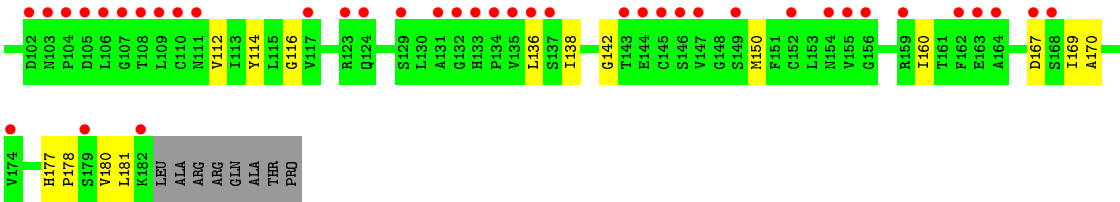


#### • Molecule 1: YfiR



#### • Molecule 1: YfiR





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	119.67Å 119.67Å 85.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.88 – 3.20 34.86 – 3.20	Depositor EDS
% Data completeness (in resolution range)	86.7 (34.88-3.20) 86.8 (34.86-3.20)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.94 (at 3.18Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.246 , 0.281 0.243 , 0.275	Depositor DCC
$R_{free}$ test set	448 reflections (5.04%)	DCC
Wilson B-factor (Å <sup>2</sup> )	60.8	Xtriage
Anisotropy	0.376	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 57.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	0 of 9354 reflections	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	4594	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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: 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	0.61	0/1198	0.76	0/1623
1	B	0.56	0/1158	0.75	1/1568 (0.1%)
1	C	0.60	1/1167 (0.1%)	0.62	0/1580
1	D	0.39	0/1134	0.57	0/1536
All	All	0.55	1/4657 (0.0%)	0.68	1/6307 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	71	CYS	CB-SG	14.18	2.06	1.82

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	152	CYS	CA-CB-SG	-5.21	104.62	114.00

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	1179	0	1194	19	0
1	B	1140	0	1154	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1149	0	1162	29	0
1	D	1116	0	1125	24	0
2	B	10	0	0	3	0
All	All	4594	0	4635	111	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 (111) 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: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:150:MET:HE1	1:B:174:VAL:HG22	1.75	0.67
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
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:D:79:ALA:HB2	1:D:114:TYR:CE2	2.40	0.57
1:B:74:GLY:O	1:B:76:THR:HG23	2.05	0.57
1:B:142:GLY:CA	1:B:152:CYS:HB3	2.33	0.57
1:B:123:ARG:HG3	1:B:123:ARG:HH11	1.69	0.57
1:B:150:MET:HE3	1:B:172:SER:HB3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:GLN:HE21	1:A:128:ARG:HH12	1.53	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:C:59:VAL:HG11	1:C:136:LEU:HD22	1.90	0.54
1:A:123:ARG:HD2	1:A:147:VAL:HB	1.89	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:150:MET:HE3	1:B:172:SER:CB	2.38	0.53
1:B:150:MET:HE1	1:B:174:VAL:CG2	2.39	0.52
1:B:123:ARG:HD2	1:B:147:VAL:CG1	2.32	0.52
1:D:76:THR:OG1	1:D:98:ARG:NH2	2.43	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:D:56:PHE:CE1	1:D:112:VAL:HG21	2.47	0.49
1:B:112:VAL:HG13	1:B:136:LEU:HB3	1.94	0.49
1:B:150:MET:HE1	1:B:172:SER:OG	2.13	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:144:GLU:CD	1:B:144:GLU:H	2.16	0.48
1:B:146:SER:O	1:B:165:ASN:ND2	2.46	0.48
1:B:128:ARG:HG2	1:C:95:HIS:HB2	1.93	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: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
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:B:150:MET:CE	1:B:172:SER:OG	2.64	0.46
1:D:58:TYR:HB3	1:D:177:HIS:HB3	1.98	0.45
1:B:114:TYR:HH	1:B:140:GLU:HG3	1.72	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:CG	1:B:123:ARG:HH11	2.30	0.44
1:A:63:LYS:HE2	1:A:64:GLU:N	2.33	0.44
1:B:128:ARG:HD2	1:C:95:HIS:CD2	2.53	0.43
1:B:113:ILE:HG22	1:B:135:VAL:HG12	1.99	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:B:70:LEU:C	1:B:70:LEU:HD23	2.39	0.43
1:D:51:VAL:HG21	1:D:160:ILE:HG21	2.01	0.42
1:A:162:PHE:CZ	1:A:184:ALA:HB2	2.53	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: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:D:177:HIS:HA	1:D:178:PRO:HD3	1.89	0.41
1:B:145:CYS:O	1:B:145:CYS:SG	2.74	0.41
1:A:180:VAL:HG13	1:A:181:LEU:HD23	2.02	0.41
1:D:56:PHE:CZ	1:D:112:VAL:HG11	2.56	0.41
1:D:47:ALA:O	1:D:51:VAL:HG23	2.21	0.41
1:A:162:PHE:CE2	1:A:184:ALA:HB2	2.56	0.41
1:C:71:CYS:CA	1:C:71:CYS:SG	3.01	0.41
1:D:167:ASP:O	1:D:170:ALA:HB3	2.20	0.41
1:D:112:VAL:HA	1:D:136:LEU:O	2.20	0.41
1:C:56:PHE:CZ	1:C:112:VAL:HG11	2.55	0.41
1:A:103:ASN:ND2	2:B:202:SO4:O4	2.44	0.41
1:B:61:TRP:HD1	1:B:134:PRO:HB2	1.86	0.41
1:A:101:VAL:HG13	1:A:126:VAL:HG23	2.03	0.41
1:C:71:CYS:SG	1:C:99:ARG:HD2	2.60	0.40
1:B:141:ARG:O	1:B:145:CYS:HB3	2.21	0.40
1:C:50:GLN:HA	1:C:53:LEU:HD12	2.03	0.40
1:D:70:LEU:HD21	1:D:114:TYR:HB2	2.04	0.40
1:B:177:HIS:ND1	1:B:178:PRO:HD2	2.36	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	151/159 (95%)	146 (97%)	5 (3%)	0	100	100
1	B	146/159 (92%)	137 (94%)	7 (5%)	2 (1%)	14	57
1	C	147/159 (92%)	138 (94%)	7 (5%)	2 (1%)	14	57
1	D	143/159 (90%)	138 (96%)	4 (3%)	1 (1%)	26	72
All	All	587/636 (92%)	559 (95%)	23 (4%)	5 (1%)	21	67

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%)	124 (96%)	5 (4%)	39	78
1	B	125/133 (94%)	114 (91%)	11 (9%)	12	45
1	C	126/133 (95%)	125 (99%)	1 (1%)	86	96
1	D	123/133 (92%)	117 (95%)	6 (5%)	31	72
All	All	503/532 (94%)	480 (95%)	23 (5%)	33	74

All (23) 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	174	VAL
1	B	99	ARG
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	175	ARG
1	B	180	VAL
1	B	185	ARG
1	C	185	ARG
1	D	40	SER
1	D	43	GLN
1	D	45	SER
1	D	57	SER
1	D	90	ASN
1	D	93	ARG

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

Mol	Chain	Res	Type
1	A	124	GLN
1	B	124	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry

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	0.94	0	6,6,6	0.41	0
2	SO4	B	202	-	4,4,4	0.66	0	6,6,6	0.62	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	SO4	B	201	-	-	0/0/0/0	0/0/0/0
2	SO4	B	202	-	-	0/0/0/0	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.

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	153/159 (96%)	-0.35	3 (1%) 68 54	25, 41, 76, 123	0
1	B	148/159 (93%)	-0.15	2 (1%) 78 65	28, 53, 93, 127	0
1	C	149/159 (93%)	1.63	41 (27%) 1 0	88, 138, 192, 217	0
1	D	145/159 (91%)	2.11	68 (46%) 0 0	82, 134, 175, 192	0
All	All	595/636 (93%)	0.80	114 (19%) 2 1	25, 98, 173, 217	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	156	GLY	8.2
1	C	147	VAL	7.3
1	C	104	PRO	7.0
1	C	171	ARG	6.2
1	C	183	LEU	6.0
1	D	39	THR	6.0
1	C	186	ARG	5.9
1	D	105	ASP	5.7
1	D	46	ASN	5.5
1	D	155	VAL	5.4
1	C	158	PRO	5.4
1	C	182	LYS	5.3
1	D	146	SER	5.2
1	C	146	SER	5.0
1	D	152	CYS	4.9
1	D	129	SER	4.9
1	D	87	VAL	4.9
1	D	143	THR	4.9
1	D	104	PRO	4.8
1	C	124	GLN	4.8
1	D	55	ILE	4.8

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Mol	Chain	Res	Type	RSRZ
1	C	102	ASP	4.7
1	D	111	ASN	4.5
1	C	156	GLY	4.2
1	C	50	GLN	4.2
1	C	71	CYS	4.2
1	D	182	LYS	4.1
1	D	69	GLN	4.1
1	D	110	CYS	4.1
1	D	164	ALA	4.0
1	D	54	GLY	4.0
1	C	180	VAL	4.0
1	D	61	TRP	4.0
1	D	145	CYS	3.9
1	A	190	PRO	3.8
1	D	81	GLY	3.8
1	C	133	HIS	3.7
1	D	52	LEU	3.7
1	C	110	CYS	3.7
1	D	132	GLY	3.7
1	D	136	LEU	3.6
1	C	99	ARG	3.5
1	C	80	ASP	3.4
1	D	159	ARG	3.3
1	D	144	GLU	3.3
1	D	103	ASN	3.3
1	C	69	GLN	3.3
1	D	38	ARG	3.3
1	C	109	LEU	3.2
1	C	143	THR	3.2
1	D	58	TYR	3.2
1	D	147	VAL	3.2
1	D	86	MET	3.1
1	C	42	GLU	3.1
1	C	172	SER	3.1
1	C	46	ASN	3.1
1	C	106	LEU	3.0
1	D	123	ARG	3.0
1	D	137	SER	3.0
1	D	41	ILE	3.0
1	D	43	GLN	3.0
1	D	133	HIS	3.0
1	C	81	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	45	SER	2.9
1	D	135	VAL	2.9
1	B	39	THR	2.9
1	D	108	THR	2.9
1	D	40	SER	2.8
1	B	38	ARG	2.8
1	D	134	PRO	2.8
1	C	107	GLY	2.8
1	C	185	ARG	2.8
1	C	105	ASP	2.8
1	D	168	SER	2.8
1	D	179	SER	2.8
1	D	117	VAL	2.7
1	C	63	LYS	2.7
1	C	67	VAL	2.7
1	C	134	PRO	2.7
1	D	163	GLU	2.7
1	D	102	ASP	2.6
1	D	162	PHE	2.6
1	D	63	LYS	2.6
1	A	189	THR	2.6
1	D	66	ALA	2.6
1	D	167	ASP	2.5
1	D	97	GLU	2.4
1	D	68	LEU	2.4
1	C	187	GLN	2.4
1	D	124	GLN	2.4
1	C	129	SER	2.4
1	D	106	LEU	2.3
1	C	168	SER	2.3
1	C	155	VAL	2.3
1	A	39	THR	2.3
1	D	48	VAL	2.3
1	D	149	SER	2.3
1	D	64	GLU	2.2
1	C	43	GLN	2.2
1	D	42	GLU	2.2
1	D	70	LEU	2.2
1	D	49	SER	2.1
1	D	62	PRO	2.1
1	C	65	PRO	2.1
1	C	184	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	109	LEU	2.1
1	D	154	ASN	2.1
1	D	107	GLY	2.1
1	D	85	GLY	2.0
1	D	131	ALA	2.0
1	C	144	GLU	2.0
1	D	51	VAL	2.0
1	D	174	VAL	2.0
1	C	141	ARG	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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	SO4	B	201	5/5	0.95	0.15	-1.03	54,56,57,59	0
2	SO4	B	202	5/5	0.99	0.13	-1.12	54,55,58,59	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.