



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 28, 2016 – 04:55 AM EST

PDB ID : 5HAA  
Title : Crystal structure of mpy-RNase J, an archaeal RNase J from Methanobrevibacter  
psychrophilus R15  
Authors : Feng, N.; Li, D.F.  
Deposited on : 2015-12-30  
Resolution : 2.90 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.

---

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
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) : rb-20028442

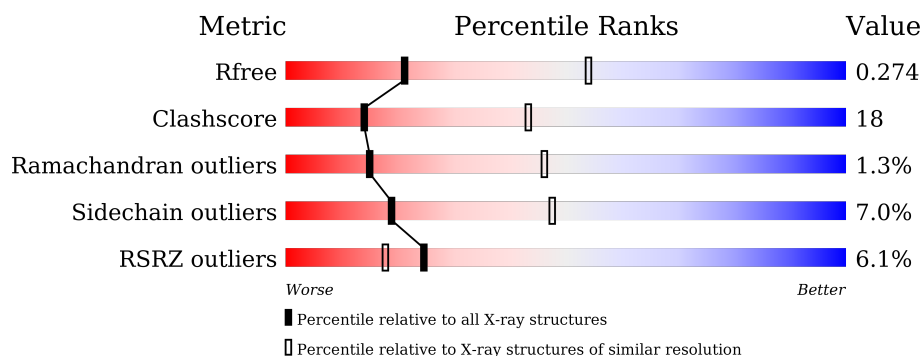
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	470	<div> <div>5%</div> <div> <div></div> <div>57%</div> <div>34%</div> <div>5%</div> <div>5%</div> </div> </div>
1	B	470	<div> <div>7%</div> <div> <div></div> <div>60%</div> <div>35%</div> <div>••</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	ZN	A	501	-	-	-	X
3	SO4	A	504	-	-	-	X
3	SO4	A	506	-	-	X	-
3	SO4	B	504	-	-	-	X
3	SO4	B	507	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7113 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 Ribonuclease J.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	447	Total	C	N	O	S	0	0	0
			3482	2192	608	662	20			
1	B	460	Total	C	N	O	S	0	0	0
			3580	2249	630	681	20			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	initiating methionine	UNP K4MAF9
A	-20	GLY	-	expression tag	UNP K4MAF9
A	-19	SER	-	expression tag	UNP K4MAF9
A	-18	SER	-	expression tag	UNP K4MAF9
A	-17	HIS	-	expression tag	UNP K4MAF9
A	-16	HIS	-	expression tag	UNP K4MAF9
A	-15	HIS	-	expression tag	UNP K4MAF9
A	-14	HIS	-	expression tag	UNP K4MAF9
A	-13	HIS	-	expression tag	UNP K4MAF9
A	-12	HIS	-	expression tag	UNP K4MAF9
A	-11	SER	-	expression tag	UNP K4MAF9
A	-10	SER	-	expression tag	UNP K4MAF9
A	-9	GLY	-	expression tag	UNP K4MAF9
A	-8	LEU	-	expression tag	UNP K4MAF9
A	-7	VAL	-	expression tag	UNP K4MAF9
A	-6	PRO	-	expression tag	UNP K4MAF9
A	-5	ARG	-	expression tag	UNP K4MAF9
A	-4	GLY	-	expression tag	UNP K4MAF9
A	-3	SER	-	expression tag	UNP K4MAF9
A	-2	HIS	-	expression tag	UNP K4MAF9
A	-1	MET	-	expression tag	UNP K4MAF9
A	0	ALA	-	expression tag	UNP K4MAF9
A	1	SER	-	expression tag	UNP K4MAF9
B	-21	MET	-	initiating methionine	UNP K4MAF9
B	-20	GLY	-	expression tag	UNP K4MAF9

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	SER	-	expression tag	UNP K4MAF9
B	-18	SER	-	expression tag	UNP K4MAF9
B	-17	HIS	-	expression tag	UNP K4MAF9
B	-16	HIS	-	expression tag	UNP K4MAF9
B	-15	HIS	-	expression tag	UNP K4MAF9
B	-14	HIS	-	expression tag	UNP K4MAF9
B	-13	HIS	-	expression tag	UNP K4MAF9
B	-12	HIS	-	expression tag	UNP K4MAF9
B	-11	SER	-	expression tag	UNP K4MAF9
B	-10	SER	-	expression tag	UNP K4MAF9
B	-9	GLY	-	expression tag	UNP K4MAF9
B	-8	LEU	-	expression tag	UNP K4MAF9
B	-7	VAL	-	expression tag	UNP K4MAF9
B	-6	PRO	-	expression tag	UNP K4MAF9
B	-5	ARG	-	expression tag	UNP K4MAF9
B	-4	GLY	-	expression tag	UNP K4MAF9
B	-3	SER	-	expression tag	UNP K4MAF9
B	-2	HIS	-	expression tag	UNP K4MAF9
B	-1	MET	-	expression tag	UNP K4MAF9
B	0	ALA	-	expression tag	UNP K4MAF9
B	1	SER	-	expression tag	UNP K4MAF9

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

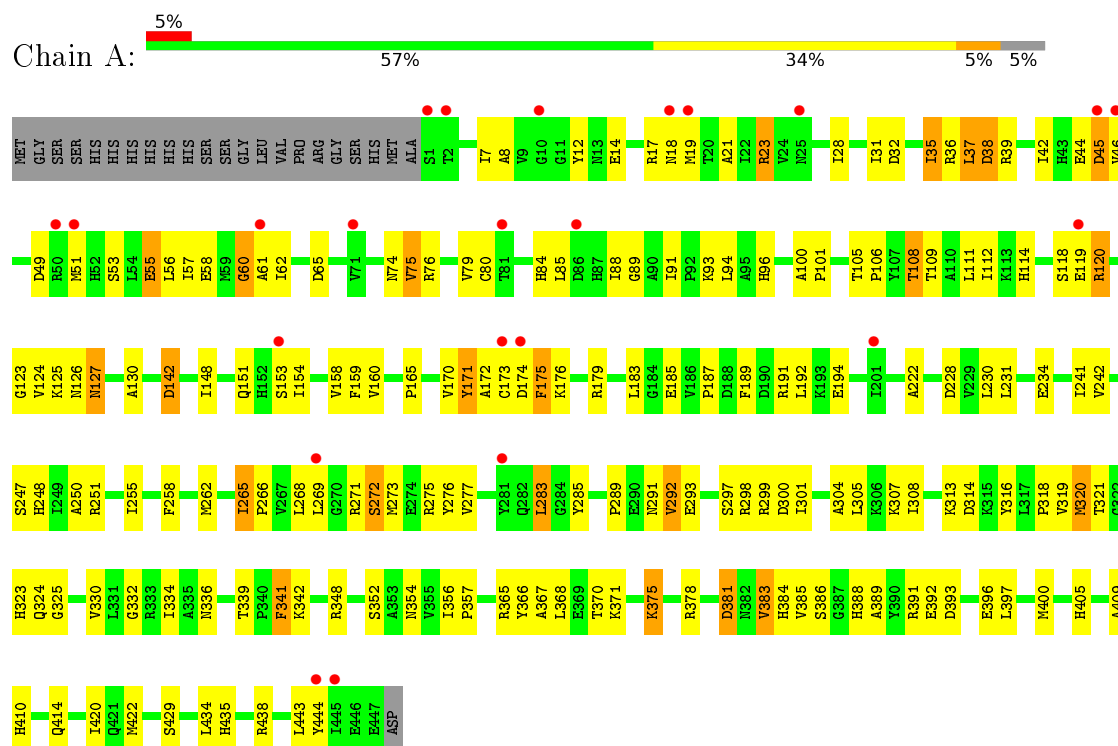
- Molecule 4 is water.

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

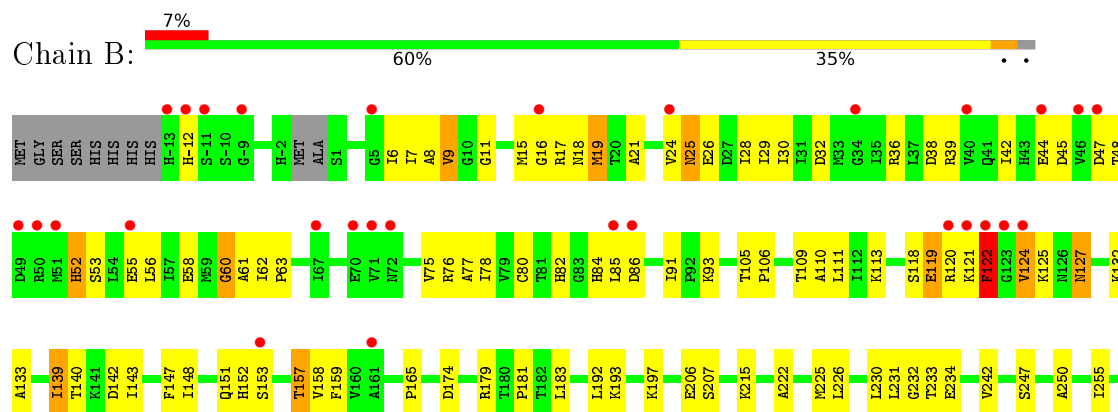
### 3 Residue-property plots [i](#)

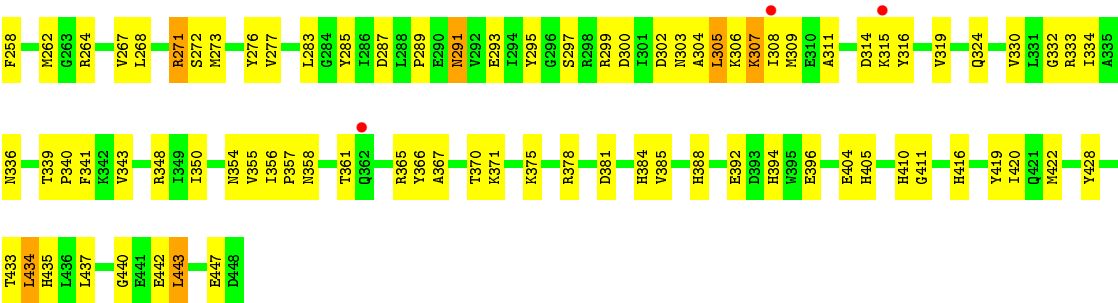
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $RSRZ > 2$ ). 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: Ribonuclease J



#### • Molecule 1: Ribonuclease J







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	162.51Å 162.51Å 180.66Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.03 – 2.90 39.03 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.1 (39.03-2.90) 98.2 (39.03-2.90)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.95 (at 2.90Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1980)	Depositor
R, $R_{free}$	0.237 , 0.274 0.236 , 0.274	Depositor DCC
$R_{free}$ test set	1992 reflections (6.42%)	DCC
Wilson B-factor (Å <sup>2</sup> )	92.8	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 67.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7113	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	110.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.34 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.3398e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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, 0.333 respectively for untwinned datasets, and 0.375, 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: ZN, 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	0.31	0/3544	0.62	1/4798 (0.0%)
1	B	0.37	0/3645	0.66	5/4932 (0.1%)
All	All	0.34	0/7189	0.64	6/9730 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	124	VAL	CG1-CB-CG2	6.88	121.91	110.90
1	B	305	LEU	CB-CG-CD2	-5.82	101.11	111.00
1	B	122	PHE	CB-CG-CD1	-5.55	116.92	120.80
1	A	443	LEU	CA-CB-CG	5.33	127.56	115.30
1	B	124	VAL	N-CA-C	-5.22	96.90	111.00
1	B	305	LEU	CA-CB-CG	5.03	126.87	115.30

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	3482	0	3488	145	0
1	B	3580	0	3577	122	0
2	A	2	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2	0	0	0	0
3	A	20	0	0	4	0
3	B	25	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	7113	0	7065	259	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 18.

All (259) 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:39:ARG:HA	1:A:42:ILE:HD13	1.56	0.85
1:B:305:LEU:HA	1:B:308:ILE:HG22	1.60	0.84
1:B:305:LEU:HD21	1:B:341:PHE:H	1.46	0.81
1:A:179:ARG:NH1	1:B:232:GLY:O	2.15	0.80
1:B:76:ARG:NH1	1:B:139:ILE:O	2.16	0.78
1:B:16:GLY:HA2	1:B:86:ASP:HB2	1.64	0.78
1:B:152:HIS:HB3	1:B:157:THR:HG21	1.64	0.77
1:A:332:GLY:O	1:A:336:ASN:ND2	2.18	0.77
1:A:56:LEU:O	1:A:60:GLY:N	2.15	0.75
1:A:151:GLN:NE2	1:A:185:GLU:O	2.20	0.74
1:A:348:ARG:NH2	1:B:396:GLU:OE1	2.22	0.73
1:A:28:ILE:HB	1:A:75:VAL:HA	1.70	0.72
1:B:343:VAL:HG13	1:B:375:LYS:HG2	1.72	0.71
1:A:334:ILE:HG12	1:A:339:THR:HG21	1.70	0.71
1:A:365:ARG:NH2	1:A:381:ASP:OD1	2.22	0.70
1:B:38:ASP:OD1	1:B:358:ASN:ND2	2.24	0.70
1:A:298:ARG:NH2	3:A:506:SO4:O4	2.26	0.69
1:A:7:ILE:HG23	1:A:21:ALA:HB3	1.75	0.69
1:B:7:ILE:HD13	1:B:442:GLU:HG3	1.73	0.68
1:A:171:TYR:HE1	1:A:173:CYS:HB2	1.59	0.68
1:A:185:GLU:OE1	1:A:185:GLU:N	2.26	0.68
1:B:142:ASP:HB3	1:B:165:PRO:HD3	1.73	0.68
1:B:140:THR:O	1:B:142:ASP:N	2.27	0.67
1:B:183:LEU:HD12	1:B:250:ALA:HA	1.74	0.67
1:B:332:GLY:O	1:B:336:ASN:ND2	2.27	0.67
1:A:96:HIS:HB2	1:A:124:VAL:HG23	1.77	0.67
1:B:306:LYS:HD3	1:B:340:PRO:HB3	1.76	0.67
1:B:365:ARG:NH2	1:B:381:ASP:OD1	2.28	0.66

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:ARG:NH2	3:A:506:SO4:O1	2.29	0.65
1:A:151:GLN:OE1	1:A:187:PRO:HD3	1.97	0.65
1:A:297:SER:OG	1:A:300:ASP:OD2	2.13	0.65
1:A:23:ARG:HH21	1:A:444:TYR:HE1	1.45	0.65
1:A:124:VAL:H	1:A:125:LYS:HD2	1.61	0.65
1:A:187:PRO:HG2	1:A:189:PHE:CE1	2.32	0.64
1:B:420:ILE:HG12	1:B:434:LEU:HD21	1.80	0.64
1:A:23:ARG:HB2	1:A:28:ILE:HD13	1.79	0.64
1:A:228:ASP:O	1:B:215:LYS:NZ	2.31	0.63
1:A:354:ASN:HB3	1:A:384:HIS:HD2	1.63	0.63
1:B:324:GLN:NE2	1:B:361:THR:OG1	2.29	0.63
1:B:15:MET:HG2	1:B:411:GLY:HA2	1.81	0.63
1:B:56:LEU:O	1:B:60:GLY:N	2.23	0.62
1:B:356:ILE:HB	1:B:361:THR:HG21	1.81	0.62
1:A:273:MET:O	1:A:277:VAL:HG22	2.00	0.62
1:A:120:ARG:NH2	1:A:123:GLY:O	2.32	0.62
1:A:183:LEU:HD12	1:A:250:ALA:HA	1.81	0.61
1:A:123:GLY:HA3	1:A:125:LYS:HD3	1.82	0.61
1:A:142:ASP:HB2	1:A:165:PRO:HD3	1.82	0.60
1:A:142:ASP:N	1:A:142:ASP:OD2	2.25	0.60
1:B:6:ILE:HB	1:B:443:LEU:HD23	1.83	0.60
1:A:251:ARG:HD3	1:A:352:SER:OG	2.01	0.59
1:B:110:ALA:HA	1:B:113:LYS:HE2	1.84	0.59
1:A:272:SER:HB2	1:A:321:THR:HG21	1.85	0.59
1:B:109:THR:O	1:B:113:LYS:HG2	2.03	0.58
1:B:267:VAL:HG22	1:B:293:GLU:HB2	1.86	0.58
1:A:111:LEU:HD11	1:A:276:TYR:CE1	2.39	0.58
1:A:101:PRO:HB3	1:A:127:ASN:ND2	2.18	0.58
1:A:258:PHE:O	1:A:262:MET:HG2	2.04	0.58
1:B:437:LEU:HD11	1:B:443:LEU:HD22	1.85	0.58
1:A:153:SER:O	1:A:154:ILE:HD13	2.04	0.57
1:B:226:LEU:HB3	1:B:258:PHE:HE2	1.68	0.57
1:A:405:HIS:HD2	1:A:435:HIS:NE2	2.03	0.57
1:B:305:LEU:HG	1:B:341:PHE:HB2	1.86	0.57
1:A:44:GLU:H	1:A:44:GLU:CD	2.07	0.57
1:A:148:ILE:HB	1:A:159:PHE:HB2	1.87	0.56
1:A:269:LEU:HD13	1:A:301:ILE:HG23	1.87	0.56
1:B:52:HIS:HA	1:B:93:LYS:HZ1	1.71	0.56
1:A:53:SER:O	1:A:57:ILE:HD12	2.06	0.56
1:A:159:PHE:CZ	1:A:175:PHE:HD2	2.24	0.56
1:B:258:PHE:O	1:B:262:MET:HG2	2.06	0.56

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:TYR:O	1:A:370:THR:HG23	2.05	0.56
1:A:84:HIS:NE2	1:A:154:ILE:HD11	2.21	0.56
1:B:273:MET:O	1:B:277:VAL:HG22	2.05	0.56
1:A:79:VAL:HG11	1:A:160:VAL:HG21	1.86	0.55
1:B:366:TYR:O	1:B:370:THR:HG23	2.05	0.55
1:B:302:ASP:OD1	1:B:333:ARG:NH2	2.40	0.55
1:B:334:ILE:HG12	1:B:339:THR:HG21	1.87	0.55
1:B:354:ASN:HB3	1:B:384:HIS:HD2	1.72	0.55
1:B:24:VAL:HB	1:B:143:ILE:HD11	1.88	0.55
1:B:11:GLY:HA3	1:B:18:ASN:HD22	1.72	0.55
1:A:323:HIS:O	1:A:365:ARG:NH1	2.39	0.54
1:B:127:ASN:N	1:B:127:ASN:OD1	2.34	0.54
1:A:304:ALA:O	1:A:308:ILE:HB	2.07	0.54
1:A:293:GLU:OE2	1:A:307:LYS:NZ	2.41	0.54
1:A:289:PRO:O	1:A:292:VAL:HG13	2.08	0.54
1:A:53:SER:OG	1:A:93:LYS:O	2.25	0.54
1:A:325:GLY:HA2	1:A:368:LEU:HD22	1.89	0.54
1:B:334:ILE:HA	1:B:339:THR:HG21	1.88	0.54
1:B:25:ASN:CG	1:B:26:GLU:H	2.05	0.54
1:A:396:GLU:O	1:A:400:MET:HG2	2.08	0.53
1:A:231:LEU:HD22	1:A:262:MET:HB3	1.88	0.53
1:A:391:ARG:HG3	1:A:422:MET:HG3	1.90	0.53
1:B:283:LEU:HD12	1:B:285:TYR:CE2	2.44	0.52
1:A:101:PRO:HB3	1:A:127:ASN:HD21	1.74	0.52
1:A:175:PHE:HE2	1:A:397:LEU:HD22	1.74	0.52
1:A:36:ARG:NE	1:A:38:ASP:OD2	2.39	0.52
1:B:330:VAL:O	1:B:334:ILE:HG13	2.09	0.52
1:B:207:SER:OG	1:B:416:HIS:HD2	1.93	0.52
1:A:389:ALA:HB1	1:A:393:ASP:HB2	1.91	0.52
1:A:273:MET:HG2	1:A:321:THR:HG23	1.92	0.51
1:B:80:CYS:SG	1:B:91:ILE:HD11	2.50	0.51
1:B:394:HIS:CD2	1:B:419:TYR:HE1	2.28	0.51
1:A:23:ARG:HB2	1:A:28:ILE:CD1	2.40	0.51
1:B:314:ASP:N	1:B:314:ASP:OD1	2.43	0.51
1:A:18:ASN:HD21	1:A:410:HIS:CG	2.29	0.51
1:A:175:PHE:CE2	1:A:397:LEU:HD22	2.45	0.51
1:A:342:LYS:HE3	1:A:375:LYS:HZ1	1.76	0.51
1:A:271:ARG:NH2	3:A:506:SO4:S	2.74	0.51
1:B:122:PHE:HE1	1:B:125:LYS:HZ1	1.58	0.51
1:A:189:PHE:CD2	1:B:234:GLU:HG3	2.46	0.51
1:A:17:ARG:HB2	1:A:36:ARG:HB2	1.93	0.51

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:ASN:HB3	1:A:76:ARG:NH1	2.25	0.51
1:B:11:GLY:HA3	1:B:18:ASN:ND2	2.27	0.50
1:A:37:LEU:H	1:A:37:LEU:HD12	1.77	0.50
1:B:206:GLU:OE1	1:B:388:HIS:HD2	1.94	0.50
1:B:55:GLU:O	1:B:58:GLU:HG2	2.12	0.50
1:B:324:GLN:HE22	1:B:356:ILE:H	1.59	0.50
1:B:8:ALA:O	1:B:440:GLY:N	2.40	0.50
1:A:124:VAL:H	1:A:125:LYS:CD	2.24	0.50
1:B:132:LYS:HD3	1:B:133:ALA:H	1.77	0.49
1:B:193:LYS:O	1:B:197:LYS:HG3	2.12	0.49
1:B:304:ALA:O	1:B:307:LYS:HB3	2.12	0.49
1:B:119:GLU:OE1	1:B:120:ARG:N	2.44	0.49
1:B:28:ILE:HB	1:B:75:VAL:HA	1.94	0.49
1:A:283:LEU:HD12	1:A:285:TYR:CE2	2.47	0.49
1:B:428:TYR:HD1	1:B:433:THR:HG22	1.77	0.49
1:B:334:ILE:HA	1:B:339:THR:CG2	2.43	0.49
1:B:305:LEU:HD21	1:B:341:PHE:N	2.21	0.49
1:B:305:LEU:CG	1:B:341:PHE:HB2	2.42	0.49
1:A:242:VAL:HG11	1:A:255:ILE:HD13	1.95	0.49
1:B:355:VAL:HG22	1:B:365:ARG:HD3	1.94	0.49
1:A:109:THR:HG21	1:A:130:ALA:HA	1.94	0.49
1:A:420:ILE:HG23	1:A:434:LEU:HD21	1.94	0.49
1:A:392:GLU:OE1	1:B:378:ARG:HD2	2.13	0.49
1:A:7:ILE:CG2	1:A:21:ALA:HB3	2.42	0.49
1:B:264:ARG:HD3	1:B:316:TYR:O	2.13	0.49
1:A:12:TYR:CE1	1:A:409:ALA:HB1	2.48	0.48
1:A:55:GLU:O	1:A:58:GLU:HG3	2.12	0.48
1:A:313:LYS:HD2	1:A:341:PHE:CE1	2.48	0.48
1:B:255:ILE:HG21	1:B:319:VAL:HG11	1.94	0.48
1:B:309:MET:HG3	1:B:341:PHE:HA	1.94	0.48
1:A:171:TYR:CE1	1:A:173:CYS:HB2	2.43	0.48
1:A:305:LEU:O	1:A:308:ILE:HG22	2.12	0.48
1:B:271:ARG:NH1	1:B:297:SER:HA	2.28	0.48
1:B:305:LEU:HD23	1:B:340:PRO:HB2	1.96	0.48
1:A:19:MET:HE3	1:A:65:ASP:HB3	1.95	0.48
1:B:111:LEU:HD11	1:B:276:TYR:CE1	2.49	0.48
1:B:30:ILE:HB	1:B:78:ILE:HG12	1.94	0.48
1:A:80:CYS:SG	1:A:91:ILE:HD11	2.53	0.48
1:A:265:ILE:HG23	1:A:316:TYR:CD1	2.49	0.47
1:A:305:LEU:HA	1:A:308:ILE:HG22	1.95	0.47
1:B:158:VAL:HG22	1:B:159:PHE:O	2.14	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:GLU:HG2	1:B:179:ARG:HH21	1.79	0.47
1:A:342:LYS:HG3	1:A:375:LYS:HE3	1.97	0.47
1:B:242:VAL:HG22	1:B:350:ILE:HB	1.95	0.47
1:B:305:LEU:HA	1:B:308:ILE:CG2	2.39	0.47
1:B:230:LEU:HB3	1:B:262:MET:HE1	1.97	0.47
1:A:356:ILE:HA	1:A:357:PRO:HD3	1.75	0.47
1:A:14:GLU:CD	1:A:17:ARG:HH21	2.17	0.47
1:B:76:ARG:HD3	1:B:140:THR:OG1	2.15	0.47
1:B:404:GLU:O	1:B:433:THR:HG23	2.15	0.47
1:A:271:ARG:NH2	3:A:506:SO4:O4	2.48	0.47
1:A:354:ASN:HB3	1:A:384:HIS:CD2	2.49	0.47
1:A:158:VAL:HG22	1:A:159:PHE:O	2.15	0.46
1:A:230:LEU:HD12	1:A:258:PHE:HD2	1.79	0.46
1:B:271:ARG:HH12	1:B:297:SER:HA	1.79	0.46
1:A:314:ASP:OD2	1:A:314:ASP:N	2.48	0.46
1:A:174:ASP:OD2	1:A:388:HIS:NE2	2.46	0.46
1:A:323:HIS:CD2	1:A:324:GLN:HG2	2.50	0.46
1:B:311:ALA:O	1:B:315:LYS:NZ	2.49	0.46
1:A:105:THR:HG1	1:A:106:PRO:HD2	1.81	0.45
1:A:334:ILE:HA	1:A:339:THR:HG21	1.98	0.45
1:A:159:PHE:HZ	1:A:175:PHE:HD2	1.64	0.45
1:B:405:HIS:HD2	1:B:435:HIS:NE2	2.14	0.45
1:A:56:LEU:CB	1:A:62:ILE:HG12	2.47	0.45
1:B:7:ILE:HB	1:B:21:ALA:HB3	1.99	0.45
1:B:151:GLN:HE21	1:B:181:PRO:HG2	1.82	0.45
1:A:35:ILE:CD1	1:A:89:GLY:HA3	2.47	0.44
1:B:174:ASP:HA	1:B:206:GLU:OE1	2.16	0.44
1:A:367:ALA:O	1:A:371:LYS:HG3	2.16	0.44
1:A:159:PHE:HE2	1:A:192:LEU:HD21	1.82	0.44
1:A:114:HIS:O	1:A:275:ARG:NH2	2.50	0.44
1:A:51:MET:SD	1:A:55:GLU:HG2	2.58	0.44
1:B:147:PHE:HB3	1:B:158:VAL:HG21	1.99	0.44
1:A:35:ILE:HD12	1:A:85:LEU:HB3	2.00	0.44
1:A:80:CYS:SG	1:A:88:ILE:HG22	2.58	0.44
1:A:334:ILE:HA	1:A:339:THR:CG2	2.48	0.44
1:B:106:PRO:O	1:B:109:THR:OG1	2.31	0.44
1:B:231:LEU:HD23	1:B:262:MET:HE3	1.98	0.44
1:A:222:ALA:HB2	1:A:385:VAL:HG13	1.99	0.44
1:A:268:LEU:HD12	1:A:277:VAL:HG21	1.99	0.44
1:A:96:HIS:O	1:A:124:VAL:HG21	2.17	0.44
1:B:159:PHE:HE2	1:B:192:LEU:HD11	1.83	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:HIS:HD2	1:A:324:GLN:HE21	1.64	0.43
1:A:241:ILE:HD11	1:A:341:PHE:CE2	2.52	0.43
1:A:124:VAL:HB	1:A:125:LYS:HA	2.00	0.43
1:A:414:GLN:OE1	1:A:414:GLN:N	2.46	0.43
1:B:56:LEU:O	1:B:61:ALA:HB3	2.19	0.43
1:A:383:VAL:HG22	1:A:383:VAL:O	2.19	0.43
1:A:176:LYS:NZ	1:A:386:SER:O	2.52	0.43
1:A:7:ILE:HD12	1:A:8:ALA:N	2.34	0.43
1:B:17:ARG:HB2	1:B:36:ARG:HB2	2.01	0.43
1:A:171:TYR:CD1	1:A:171:TYR:C	2.92	0.43
1:A:265:ILE:HA	1:A:266:PRO:HD2	1.76	0.43
1:B:36:ARG:HH12	1:B:39:ARG:NH2	2.17	0.43
1:A:265:ILE:HD13	1:A:291:ASN:HB2	2.00	0.42
1:A:308:ILE:HD13	1:A:318:PRO:HG3	2.01	0.42
1:B:119:GLU:CD	1:B:121:LYS:H	2.22	0.42
1:B:7:ILE:HG13	1:B:21:ALA:HB3	2.01	0.42
1:B:29:ILE:HD12	1:B:77:ALA:HB3	2.02	0.42
1:B:242:VAL:O	1:B:319:VAL:HA	2.19	0.42
1:B:9:VAL:HG13	1:B:19:MET:HB3	2.01	0.42
1:A:31:ILE:HB	1:A:171:TYR:CE2	2.54	0.42
1:A:62:ILE:HD12	1:A:94:LEU:HD21	2.02	0.42
1:A:175:PHE:CE1	1:A:389:ALA:HB3	2.55	0.42
1:A:241:ILE:HD11	1:A:341:PHE:CZ	2.54	0.42
1:B:119:GLU:OE2	1:B:121:LYS:HB3	2.19	0.42
1:A:56:LEU:O	1:A:61:ALA:HB3	2.20	0.42
1:B:308:ILE:HG21	1:B:308:ILE:HD13	1.83	0.42
1:A:191:ARG:HH11	1:A:194:GLU:HG2	1.84	0.42
1:A:36:ARG:HH12	1:A:39:ARG:NH2	2.17	0.42
1:B:367:ALA:O	1:B:371:LYS:HG3	2.20	0.42
1:B:56:LEU:HD13	1:B:93:LYS:HE3	2.01	0.42
1:A:93:LYS:HA	1:A:93:LYS:HD3	1.72	0.42
1:B:289:PRO:HB2	1:B:291:ASN:OD1	2.20	0.42
1:A:108:THR:O	1:A:112:ILE:HD12	2.19	0.42
1:A:305:LEU:HD22	1:A:341:PHE:HD2	1.84	0.42
1:B:293:GLU:HB3	1:B:295:TYR:CZ	2.55	0.42
1:B:230:LEU:HD12	1:B:258:PHE:HD2	1.85	0.42
1:A:396:GLU:OE2	1:B:348:ARG:NH2	2.53	0.41
1:A:45:ASP:OD1	1:A:45:ASP:N	2.52	0.41
1:B:15:MET:HE3	1:B:410:HIS:HE1	1.85	0.41
1:B:174:ASP:OD2	1:B:410:HIS:NE2	2.53	0.41
1:A:378:ARG:NH1	1:B:392:GLU:OE2	2.47	0.41

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:HIS:NE2	1:B:84:HIS:HB2	2.35	0.41
1:B:148:ILE:HB	1:B:159:PHE:HB2	2.03	0.41
1:B:299:ARG:HG3	1:B:300:ASP:OD2	2.21	0.41
1:A:231:LEU:CD2	1:A:262:MET:HB3	2.51	0.41
1:A:251:ARG:O	1:A:255:ILE:HG13	2.19	0.41
1:B:105:THR:OG1	1:B:106:PRO:HD2	2.21	0.41
1:B:268:LEU:HG	1:B:319:VAL:CG2	2.50	0.41
1:B:7:ILE:HD13	1:B:442:GLU:CG	2.45	0.41
1:A:28:ILE:HB	1:A:75:VAL:CA	2.45	0.41
1:B:56:LEU:HD22	1:B:93:LYS:HD3	2.03	0.41
1:B:19:MET:HE2	1:B:63:PRO:HG2	2.03	0.41
1:B:152:HIS:CG	1:B:153:SER:N	2.88	0.41
1:B:9:VAL:HG13	1:B:19:MET:O	2.20	0.41
1:A:100:ALA:O	1:A:126:ASN:HB3	2.20	0.41
1:A:171:TYR:HD1	1:A:172:ALA:N	2.18	0.41
1:A:153:SER:O	1:A:248:HIS:HA	2.21	0.41
1:B:53:SER:O	1:B:62:ILE:HD11	2.21	0.41
1:B:356:ILE:HA	1:B:357:PRO:HD3	1.84	0.40
1:A:230:LEU:HD12	1:A:258:PHE:CD2	2.56	0.40
1:A:320:MET:HB2	1:A:330:VAL:CG1	2.51	0.40
1:A:438:ARG:HD2	1:A:438:ARG:HA	1.95	0.40
1:A:154:ILE:HD12	1:A:247:SER:HB2	2.04	0.40
1:A:56:LEU:HB2	1:A:62:ILE:HG12	2.03	0.40
1:B:222:ALA:HA	1:B:225:MET:HE2	2.04	0.40
1:B:56:LEU:HB2	1:B:62:ILE:HG12	2.04	0.40
1:A:268:LEU:HD22	1:A:319:VAL:HB	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	445/470 (95%)	411 (92%)	29 (6%)	5 (1%)	17	51
1	B	456/470 (97%)	417 (91%)	32 (7%)	7 (2%)	13	42
All	All	901/940 (96%)	828 (92%)	61 (7%)	12 (1%)	15	46

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	ARG
1	B	45	ASP
1	A	118	SER
1	A	45	ASP
1	B	32	ASP
1	B	118	SER
1	A	32	ASP
1	B	25	ASN
1	B	47	ASP
1	B	60	GLY
1	A	60	GLY
1	B	122	PHE

### 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	375/394 (95%)	349 (93%)	26 (7%)	19	48
1	B	386/394 (98%)	359 (93%)	27 (7%)	19	47
All	All	761/788 (97%)	708 (93%)	53 (7%)	19	47

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ARG
1	A	35	ILE
1	A	37	LEU
1	A	38	ASP

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	46	VAL
1	A	49	ASP
1	A	55	GLU
1	A	75	VAL
1	A	108	THR
1	A	119	GLU
1	A	127	ASN
1	A	142	ASP
1	A	170	VAL
1	A	171	TYR
1	A	175	PHE
1	A	265	ILE
1	A	272	SER
1	A	283	LEU
1	A	292	VAL
1	A	299	ARG
1	A	320	MET
1	A	341	PHE
1	A	375	LYS
1	A	381	ASP
1	A	383	VAL
1	A	429	SER
1	B	-12	HIS
1	B	9	VAL
1	B	19	MET
1	B	42	ILE
1	B	44	GLU
1	B	48	THR
1	B	52	HIS
1	B	85	LEU
1	B	119	GLU
1	B	122	PHE
1	B	124	VAL
1	B	127	ASN
1	B	139	ILE
1	B	157	THR
1	B	233	THR
1	B	247	SER
1	B	271	ARG
1	B	272	SER
1	B	287	ASP
1	B	291	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	303	ASN
1	B	307	LYS
1	B	385	VAL
1	B	422	MET
1	B	434	LEU
1	B	443	LEU
1	B	447	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	74	ASN
1	A	82	HIS
1	A	127	ASN
1	A	163	HIS
1	A	209	ASN
1	A	257	GLN
1	A	323	HIS
1	A	405	HIS
1	B	18	ASN
1	B	43	HIS
1	B	82	HIS
1	B	87	HIS
1	B	115	GLN
1	B	209	ASN
1	B	257	GLN
1	B	282	GLN
1	B	324	GLN
1	B	336	ASN
1	B	362	GLN
1	B	388	HIS
1	B	405	HIS
1	B	410	HIS
1	B	416	HIS

### 5.3.3 RNA ⓘ

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 13 ligands modelled in this entry, 4 are monoatomic - leaving 9 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	SO4	A	503	-	4,4,4	0.27	0	6,6,6	0.11	0
3	SO4	A	504	1	4,4,4	0.21	0	6,6,6	0.09	0
3	SO4	A	505	-	4,4,4	0.22	0	6,6,6	0.09	0
3	SO4	A	506	1	4,4,4	0.25	0	6,6,6	0.07	0
3	SO4	B	503	-	4,4,4	0.28	0	6,6,6	0.13	0
3	SO4	B	504	-	4,4,4	0.26	0	6,6,6	0.06	0
3	SO4	B	505	-	4,4,4	0.23	0	6,6,6	0.06	0
3	SO4	B	506	-	4,4,4	0.25	0	6,6,6	0.08	0
3	SO4	B	507	-	4,4,4	0.26	0	6,6,6	0.07	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	SO4	A	503	-	-	0/0/0/0	0/0/0/0
3	SO4	A	504	1	-	0/0/0/0	0/0/0/0
3	SO4	A	505	-	-	0/0/0/0	0/0/0/0
3	SO4	A	506	1	-	0/0/0/0	0/0/0/0
3	SO4	B	503	-	-	0/0/0/0	0/0/0/0
3	SO4	B	504	-	-	0/0/0/0	0/0/0/0

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	B	505	-	-	0/0/0/0	0/0/0/0
3	SO4	B	506	-	-	0/0/0/0	0/0/0/0
3	SO4	B	507	-	-	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.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	506	SO4	4	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	447/470 (95%)	0.39	23 (5%)	32 25	63, 101, 148, 208	0
1	B	460/470 (97%)	0.43	32 (6%)	19 13	63, 109, 168, 246	0
All	All	907/940 (96%)	0.41	55 (6%)	25 18	63, 105, 158, 246	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	44	GLU	7.7
1	B	124	VAL	6.7
1	B	123	GLY	5.5
1	A	46	VAL	5.3
1	B	122	PHE	4.9
1	B	50	ARG	4.7
1	B	71	VAL	4.7
1	A	45	ASP	4.5
1	A	50	ARG	4.4
1	A	86	ASP	4.2
1	B	46	VAL	4.2
1	B	121	LYS	4.1
1	A	445	ILE	4.0
1	A	173	CYS	3.6
1	B	-12	HIS	3.5
1	B	51	MET	3.3
1	A	51	MET	3.3
1	A	201	ILE	3.3
1	A	174	ASP	3.3
1	B	72	ASN	3.1
1	B	70	GLU	3.1
1	B	67	ILE	3.0
1	A	444	TYR	3.0
1	A	18	ASN	3.0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	120	ARG	3.0
1	B	315	LYS	2.9
1	A	119	GLU	2.9
1	B	47	ASP	2.8
1	B	34	GLY	2.7
1	B	-11	SER	2.7
1	B	153	SER	2.6
1	B	55	GLU	2.6
1	B	-9	GLY	2.6
1	A	2	THR	2.5
1	A	153	SER	2.5
1	A	281	TYR	2.5
1	A	1	SER	2.5
1	B	24	VAL	2.4
1	A	25	ASN	2.4
1	A	61	ALA	2.4
1	B	85	LEU	2.4
1	A	19	MET	2.3
1	A	10	GLY	2.3
1	A	71	VAL	2.3
1	B	40	VAL	2.3
1	B	86	ASP	2.3
1	A	81	THR	2.3
1	B	-13	HIS	2.3
1	B	362	GLN	2.2
1	B	161	ALA	2.2
1	B	5	GLY	2.2
1	B	308	ILE	2.1
1	B	49	ASP	2.1
1	B	16	GLY	2.1
1	A	269	LEU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

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
3	SO4	B	507	5/5	0.58	1.02	19.89	428,428,430,431	0
3	SO4	B	504	5/5	0.81	0.43	5.87	167,167,168,168	0
3	SO4	A	504	5/5	0.64	0.38	3.71	137,159,165,165	0
3	SO4	A	506	5/5	0.81	0.28	1.52	130,163,169,170	0
2	ZN	B	501	1/1	0.94	0.35	1.25	139,139,139,139	0
2	ZN	A	501	1/1	0.72	0.40	0.59	152,152,152,152	0
3	SO4	A	503	5/5	0.94	0.21	-0.17	67,78,83,88	0
2	ZN	B	502	1/1	0.96	0.27	-0.18	114,114,114,114	0
3	SO4	B	505	5/5	0.86	0.17	-0.20	128,148,152,153	0
3	SO4	B	503	5/5	0.96	0.18	-0.47	76,87,97,104	0
3	SO4	A	505	5/5	0.86	0.14	-0.72	108,123,132,132	0
2	ZN	A	502	1/1	0.98	0.20	-1.61	93,93,93,93	0
3	SO4	B	506	5/5	0.75	0.25	-	187,187,187,187	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.