



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 03:30 PM GMT

PDB ID : 3KE6  
Title : The crystal structure of the RsbU and RsbW domains of Rv1364c from Mycobacterium tuberculosis  
Authors : King-Scott, J.; Panjikar, S.; Tucker, P.A.  
Deposited on : 2009-10-24  
Resolution : 2.60 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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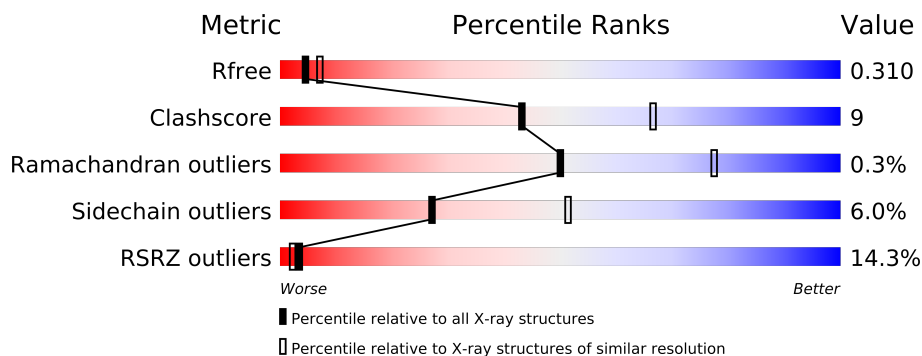
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.60 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))
$R_{free}$	66092	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	399	
1	B	399	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5304 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Protein Rv1364c/MT1410.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	1	0
			2618	1641	459	507	11			
1	B	348	Total	C	N	O	S	0	1	0
			2568	1605	453	500	10			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-27	MET	-	EXPRESSION TAG	UNP Q11034
A	-26	LYS	-	EXPRESSION TAG	UNP Q11034
A	-25	HIS	-	EXPRESSION TAG	UNP Q11034
A	-24	HIS	-	EXPRESSION TAG	UNP Q11034
A	-23	HIS	-	EXPRESSION TAG	UNP Q11034
A	-22	HIS	-	EXPRESSION TAG	UNP Q11034
A	-21	HIS	-	EXPRESSION TAG	UNP Q11034
A	-20	HIS	-	EXPRESSION TAG	UNP Q11034
A	-19	PRO	-	EXPRESSION TAG	UNP Q11034
A	-18	MET	-	EXPRESSION TAG	UNP Q11034
A	-17	SER	-	EXPRESSION TAG	UNP Q11034
A	-16	ASP	-	EXPRESSION TAG	UNP Q11034
A	-15	TYR	-	EXPRESSION TAG	UNP Q11034
A	-14	ASP	-	EXPRESSION TAG	UNP Q11034
A	-13	ILE	-	EXPRESSION TAG	UNP Q11034
A	-12	PRO	-	EXPRESSION TAG	UNP Q11034
A	-11	THR	-	EXPRESSION TAG	UNP Q11034
A	-10	THR	-	EXPRESSION TAG	UNP Q11034
A	-9	GLU	-	EXPRESSION TAG	UNP Q11034
A	-8	ASN	-	EXPRESSION TAG	UNP Q11034
A	-7	LEU	-	EXPRESSION TAG	UNP Q11034
A	-6	TYR	-	EXPRESSION TAG	UNP Q11034
A	-5	PHE	-	EXPRESSION TAG	UNP Q11034
A	-4	GLN	-	EXPRESSION TAG	UNP Q11034
A	-3	GLY	-	EXPRESSION TAG	UNP Q11034

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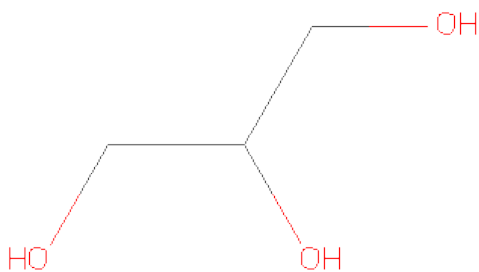
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ALA	-	EXPRESSION TAG	UNP Q11034
A	-1	MET	-	EXPRESSION TAG	UNP Q11034
A	0	VAL	-	EXPRESSION TAG	UNP Q11034
B	-27	MET	-	EXPRESSION TAG	UNP Q11034
B	-26	LYS	-	EXPRESSION TAG	UNP Q11034
B	-25	HIS	-	EXPRESSION TAG	UNP Q11034
B	-24	HIS	-	EXPRESSION TAG	UNP Q11034
B	-23	HIS	-	EXPRESSION TAG	UNP Q11034
B	-22	HIS	-	EXPRESSION TAG	UNP Q11034
B	-21	HIS	-	EXPRESSION TAG	UNP Q11034
B	-20	HIS	-	EXPRESSION TAG	UNP Q11034
B	-19	PRO	-	EXPRESSION TAG	UNP Q11034
B	-18	MET	-	EXPRESSION TAG	UNP Q11034
B	-17	SER	-	EXPRESSION TAG	UNP Q11034
B	-16	ASP	-	EXPRESSION TAG	UNP Q11034
B	-15	TYR	-	EXPRESSION TAG	UNP Q11034
B	-14	ASP	-	EXPRESSION TAG	UNP Q11034
B	-13	ILE	-	EXPRESSION TAG	UNP Q11034
B	-12	PRO	-	EXPRESSION TAG	UNP Q11034
B	-11	THR	-	EXPRESSION TAG	UNP Q11034
B	-10	THR	-	EXPRESSION TAG	UNP Q11034
B	-9	GLU	-	EXPRESSION TAG	UNP Q11034
B	-8	ASN	-	EXPRESSION TAG	UNP Q11034
B	-7	LEU	-	EXPRESSION TAG	UNP Q11034
B	-6	TYR	-	EXPRESSION TAG	UNP Q11034
B	-5	PHE	-	EXPRESSION TAG	UNP Q11034
B	-4	GLN	-	EXPRESSION TAG	UNP Q11034
B	-3	GLY	-	EXPRESSION TAG	UNP Q11034
B	-2	ALA	-	EXPRESSION TAG	UNP Q11034
B	-1	MET	-	EXPRESSION TAG	UNP Q11034
B	0	VAL	-	EXPRESSION TAG	UNP Q11034

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total 2	Mn 2	0	0
4	A	2	Total 2	Mn 2	0	0

- Molecule 5 is water.

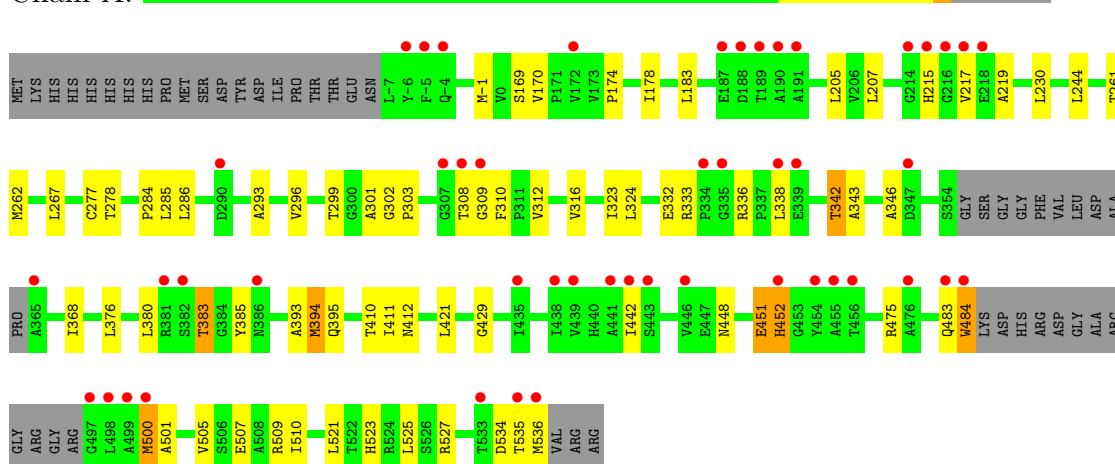
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	51	Total 51	O 51	0	0
5	B	52	Total 52	O 52	0	0

### 3 Residue-property plots

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.

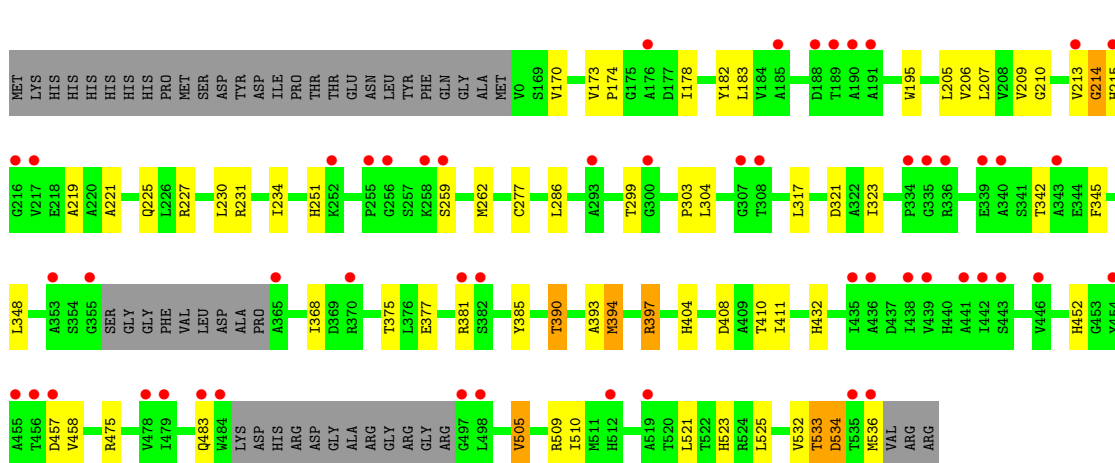
#### • Molecule 1: Protein Rv1364c/MT1410

Chain A:



#### • Molecule 1: Protein Rv1364c/MT1410

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	100.12Å 100.12Å 169.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.63 – 2.60 19.61 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.2 (19.63-2.60) 99.2 (19.61-2.60)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.82 (at 2.59Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.208 , 0.267 0.277 , 0.310	Depositor DCC
$R_{free}$ test set	1356 reflections (5.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	51.6	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 33.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 26968 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	5304	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.31% of the height of the origin peak. No significant pseudotranslation is detected.*

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



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MN, 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	0.57	0/2668	0.74	2/3637 (0.1%)
1	B	0.54	0/2613	0.73	2/3563 (0.1%)
All	All	0.56	0/5281	0.74	4/7200 (0.1%)

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

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	475	ARG	NE-CZ-NH2	-7.42	116.59	120.30
1	A	475	ARG	NE-CZ-NH2	-6.88	116.86	120.30
1	B	475	ARG	NE-CZ-NH1	6.86	123.73	120.30
1	A	475	ARG	NE-CZ-NH1	5.99	123.30	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	452	HIS	Peptide

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2618	0	2588	54	0
1	B	2568	0	2536	38	0
2	A	5	0	0	0	0
3	A	6	0	8	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	51	0	0	3	0
5	B	52	0	0	4	0
All	All	5304	0	5132	91	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (91) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:302:GLY:H	1:A:308:THR:HG21	1.31	0.91
1:B:230:LEU:HD22	1:B:262:MET:CE	2.04	0.88
1:B:206:VAL:HG11	1:B:234:ILE:HD13	1.59	0.83
1:B:214:GLY:O	1:B:219:ALA:HB2	1.82	0.80
1:B:207:LEU:HD11	1:B:394:MET:HG2	1.64	0.78
1:A:410:THR:HG22	1:A:412:ASN:H	1.49	0.77
1:A:230:LEU:HD22	1:A:262:MET:CE	2.15	0.77
1:A:285:LEU:HD21	1:A:293:ALA:HB1	1.65	0.77
1:A:302:GLY:N	1:A:308:THR:HG21	2.01	0.76
1:A:267:LEU:HD22	1:A:394:MET:CE	2.17	0.75
1:B:230:LEU:HD22	1:B:262:MET:HE3	1.74	0.70
1:B:213:VAL:HG12	5:B:30:HOH:O	1.91	0.68
1:A:183:LEU:HD11	1:A:376:LEU:HD21	1.73	0.68
1:A:505:VAL:HG22	1:A:523:HIS:ND1	2.09	0.67
1:A:207:LEU:HD11	1:A:394:MET:HG2	1.77	0.66
1:B:221:ALA:O	1:B:225:GLN:NE2	2.29	0.64
1:A:451:GLU:OE1	1:B:432:HIS:NE2	2.30	0.64
1:A:285:LEU:HD11	1:A:346:ALA:HA	1.81	0.63
1:B:214:GLY:O	1:B:219:ALA:CB	2.47	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:509:ARG:O	1:A:510:ILE:HD12	1.98	0.63
1:A:333:ARG:HD2	1:A:383:THR:HG21	1.79	0.62
1:A:267:LEU:HD22	1:A:394:MET:HE1	1.81	0.62
1:B:521:LEU:HD12	1:B:521:LEU:O	2.00	0.61
1:B:505:VAL:HG22	1:B:523:HIS:ND1	2.16	0.60
1:B:183:LEU:HD23	1:B:385:TYR:CE2	2.37	0.60
1:A:368:ILE:HD11	1:A:393:ALA:HB1	1.86	0.58
1:A:285:LEU:HD23	1:A:286:LEU:N	2.20	0.57
1:B:521:LEU:C	1:B:521:LEU:HD12	2.24	0.57
1:B:404[A]:HIS:NE2	5:B:84:HOH:O	2.27	0.56
1:A:525:LEU:O	5:A:48:HOH:O	2.18	0.56
1:B:505:VAL:HG22	1:B:523:HIS:CG	2.41	0.56
1:B:178:ILE:HD11	1:B:205:LEU:HD21	1.88	0.55
1:B:525:LEU:O	5:B:24:HOH:O	2.18	0.55
1:A:410:THR:CG2	1:A:411:ILE:N	2.72	0.53
1:A:278:THR:H	1:A:299:THR:CG2	2.23	0.51
1:B:509:ARG:O	1:B:510:ILE:HD13	2.11	0.51
1:B:390:THR:HG22	5:B:36:HOH:O	2.11	0.50
1:A:521:LEU:HD12	1:A:521:LEU:O	2.12	0.50
1:B:230:LEU:HD22	1:B:262:MET:HE2	1.89	0.50
1:A:230:LEU:HD22	1:A:262:MET:HE2	1.94	0.50
1:A:308:THR:HG23	1:A:309:GLY:O	2.13	0.49
1:A:267:LEU:HD22	1:A:394:MET:HE2	1.91	0.49
1:B:182:TYR:C	1:B:183:LEU:HD12	2.33	0.49
1:A:380:LEU:HA	1:A:383:THR:HG22	1.94	0.48
1:A:452:HIS:C	1:A:483:GLN:HB2	2.34	0.48
1:A:333:ARG:HH11	1:A:383:THR:HG21	1.79	0.48
1:A:183:LEU:N	1:A:183:LEU:HD12	2.29	0.48
1:A:174:PRO:HA	5:A:53:HOH:O	2.14	0.48
1:A:284:PRO:O	1:A:296:VAL:HG22	2.13	0.48
1:A:278:THR:H	1:A:299:THR:HG22	1.79	0.47
1:A:448:ASN:O	1:A:452:HIS:HB2	2.15	0.47
1:A:261:THR:HG22	1:A:303:PRO:HB3	1.97	0.47
1:B:195:TRP:CZ3	1:B:210:GLY:HA3	2.50	0.47
1:A:521:LEU:C	1:A:521:LEU:HD12	2.36	0.46
1:A:342:THR:HG22	1:A:343:ALA:N	2.31	0.46
1:A:178:ILE:CD1	1:A:205:LEU:HD21	2.46	0.46
1:B:377:GLU:OE2	1:B:536:MET:HE1	2.16	0.46
1:B:532:VAL:HG13	1:B:536:MET:SD	2.56	0.46
1:A:410:THR:HG22	1:A:411:ILE:N	2.31	0.46
1:A:278:THR:OG1	1:A:299:THR:HG22	2.15	0.45
1:B:206:VAL:CG1	1:B:234:ILE:HD13	2.40	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:178:ILE:HD11	1:A:205:LEU:HD21	1.98	0.45
1:B:227:ARG:O	1:B:231:ARG:HD2	2.17	0.45
1:A:215:HIS:HA	1:A:219:ALA:HB3	1.99	0.45
1:B:214:GLY:C	1:B:219:ALA:HB2	2.36	0.44
1:A:-1:MET:CG	1:A:169:SER:O	2.65	0.44
1:A:395:GLN:NE2	5:A:76:HOH:O	2.46	0.44
1:B:368:ILE:HD11	1:B:393:ALA:HB1	1.98	0.44
1:B:259:SER:HA	1:B:303:PRO:HB2	1.98	0.44
1:B:505:VAL:HG11	1:B:521:LEU:HB2	1.99	0.44
1:A:421:LEU:HD21	1:A:442:ILE:HD12	2.00	0.44
1:A:215:HIS:CD2	1:A:215:HIS:C	2.91	0.43
1:A:484:TRP:CE3	1:A:484:TRP:HA	2.53	0.43
1:B:215:HIS:N	1:B:219:ALA:HB2	2.33	0.43
1:B:317:LEU:HD22	1:B:321:ASP:HB3	1.99	0.43
1:B:174:PRO:O	1:B:397:ARG:HD2	2.19	0.43
1:A:484:TRP:HA	1:A:484:TRP:HE3	1.84	0.42
1:A:323:ILE:C	1:A:324:LEU:HD12	2.39	0.42
1:A:332:GLU:OE2	1:A:338:LEU:HD21	2.19	0.42
1:A:244:LEU:HD13	1:A:310:PHE:CD1	2.54	0.42
1:A:316:VAL:O	1:A:316:VAL:HG13	2.18	0.42
1:B:408:ASP:HB3	1:B:410:THR:HG23	2.02	0.42
1:A:509:ARG:C	1:A:510:ILE:HD12	2.39	0.42
1:B:277:CYS:HA	1:B:299:THR:HG21	2.02	0.42
1:A:429:GLY:O	1:A:527:ARG:NH1	2.53	0.42
1:B:452:HIS:O	1:B:483:GLN:HA	2.20	0.41
1:A:299:THR:HG23	1:A:301:ALA:H	1.85	0.41
1:A:500:MET:HE3	1:A:501:ALA:HB2	2.02	0.41
1:B:345:PHE:HE1	1:B:375:THR:HG21	1.85	0.41
1:A:277:CYS:HB2	1:A:312:VAL:HG12	2.02	0.41
1:B:533:THR:HA	1:B:534:ASP:HA	1.85	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	349/399 (88%)	338 (97%)	11 (3%)	0	100	100
1	B	343/399 (86%)	327 (95%)	14 (4%)	2 (1%)	33	63
All	All	692/798 (87%)	665 (96%)	25 (4%)	2 (0%)	50	77

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	214	GLY
1	B	457	ASP

### 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	268/304 (88%)	254 (95%)	14 (5%)	32	59
1	B	263/304 (86%)	245 (93%)	18 (7%)	22	43
All	All	531/608 (87%)	499 (94%)	32 (6%)	27	51

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

Mol	Chain	Res	Type
1	A	170	VAL
1	A	217	VAL
1	A	336	ARG
1	A	342	THR
1	A	383	THR
1	A	385	TYR
1	A	394	MET
1	A	451	GLU
1	A	484	TRP
1	A	500	MET
1	A	507	GLU
1	A	534	ASP
1	A	535	THR
1	A	536	MET
1	B	170	VAL

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Mol	Chain	Res	Type
1	B	173	VAL
1	B	209	VAL
1	B	251	HIS
1	B	286	LEU
1	B	304	LEU
1	B	323	ILE
1	B	342	THR
1	B	348	LEU
1	B	381	ARG
1	B	390	THR
1	B	394	MET
1	B	397	ARG
1	B	411	ILE
1	B	458	VAL
1	B	505	VAL
1	B	533	THR
1	B	534	ASP

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	253	GLN
1	A	395	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	A	1	-	4,4,4	0.92	0	6,6,6	0.19	0
3	GOL	A	540	-	5,5,5	0.63	0	5,5,5	0.59	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	A	1	-	-	0/0/0/0	0/0/0/0
3	GOL	A	540	-	-	0/4/4/4	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.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	354/399 (88%)	0.74	48 (13%) 4 3	46, 53, 69, 88	0
1	B	348/399 (87%)	0.86	53 (15%) 3 2	47, 54, 64, 90	0
All	All	702/798 (87%)	0.80	101 (14%) 3 2	46, 54, 67, 90	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	484	TRP	7.4
1	B	190	ALA	6.2
1	A	484	TRP	6.1
1	B	191	ALA	6.0
1	A	215	HIS	5.6
1	B	215	HIS	5.5
1	A	535	THR	5.4
1	A	365	ALA	5.3
1	B	457	ASP	5.2
1	B	335	GLY	4.7
1	B	188	ASP	4.7
1	A	217	VAL	4.6
1	A	-5	PHE	4.6
1	B	217	VAL	4.4
1	B	535	THR	4.3
1	A	536	MET	4.2
1	A	456	THR	4.2
1	A	218	GLU	4.2
1	B	382	SER	4.1
1	B	189	THR	4.1
1	B	365	ALA	3.7
1	A	216	GLY	3.7
1	B	339	GLU	3.7
1	B	497	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	382	SER	3.6
1	B	252	LYS	3.5
1	B	255	PRO	3.5
1	A	334	PRO	3.4
1	A	188	ASP	3.4
1	A	-6	TYR	3.3
1	A	439	VAL	3.4
1	B	355	GLY	3.3
1	A	443	SER	3.3
1	B	307	GLY	3.3
1	A	307	GLY	3.3
1	A	190	ALA	3.2
1	B	442	ILE	3.2
1	B	308	THR	3.2
1	A	338	LEU	3.2
1	B	536	MET	3.2
1	A	381	ARG	3.2
1	B	340	ALA	3.1
1	B	456	THR	3.1
1	A	214	GLY	3.1
1	B	300	GLY	3.1
1	B	498	LEU	3.0
1	A	454	TYR	3.0
1	B	334	PRO	3.0
1	A	335	GLY	3.0
1	B	370	ARG	3.0
1	B	454	TYR	2.9
1	A	455	ALA	2.9
1	A	191	ALA	2.9
1	A	386	ASN	2.9
1	A	441	ALA	2.8
1	B	483	GLN	2.8
1	B	381	ARG	2.8
1	A	339	GLU	2.8
1	B	443	SER	2.7
1	A	498	LEU	2.7
1	A	189	THR	2.7
1	A	446	VAL	2.7
1	A	442	ILE	2.6
1	B	455	ALA	2.6
1	B	336	ARG	2.6
1	B	439	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	483	GLN	2.6
1	B	479	ILE	2.5
1	B	259	SER	2.5
1	A	-4	GLN	2.5
1	A	497	GLY	2.4
1	A	438	ILE	2.4
1	A	499	ALA	2.4
1	B	441	ALA	2.4
1	B	478	VAL	2.4
1	B	293	ALA	2.4
1	A	187	GLU	2.4
1	A	308	THR	2.4
1	B	438	ILE	2.4
1	B	216	GLY	2.4
1	B	446	VAL	2.3
1	B	435	ILE	2.3
1	A	309	GLY	2.3
1	A	476	ALA	2.2
1	A	452	HIS	2.2
1	B	256	GLY	2.2
1	B	185	ALA	2.2
1	B	258	LYS	2.2
1	A	500	MET	2.2
1	B	512	HIS	2.1
1	A	290	ASP	2.1
1	B	176	ALA	2.1
1	B	343	ALA	2.1
1	B	353	ALA	2.1
1	A	347	ASP	2.1
1	A	172	VAL	2.1
1	B	519	ALA	2.1
1	B	213	VAL	2.1
1	B	436	ALA	2.1
1	A	533	THR	2.0
1	A	435	ILE	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

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	MN	A	2	1/1	0.17	0.73	73,73,73,73	0
4	MN	A	541	1/1	0.09	-1.13	69,69,69,69	0
3	GOL	A	540	6/6	0.17	-2.69	36,42,50,54	0
4	MN	B	3	1/1	0.04	-2.72	67,67,67,67	0
2	SO4	A	1	5/5	0.10	-2.89	42,43,46,47	0
4	MN	B	4	1/1	0.12	-3.10	72,72,72,72	0

### 6.5 Other polymers

There are no such residues in this entry.