



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:59 AM 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 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

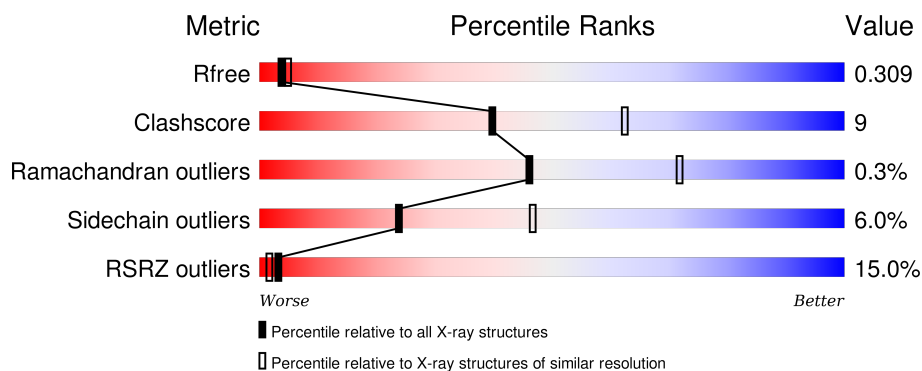
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

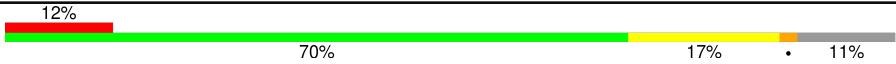

The reported resolution of this entry is 2.60 Å.

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))
$R_{free}$	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (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. 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	399	 12% 70% 17% • 11%
1	B	399	 14% 71% 15% • 13%

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5304 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 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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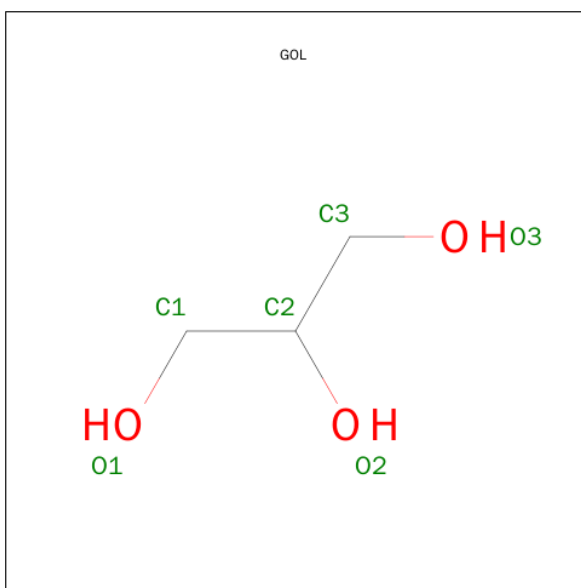
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ALA	-	EXPRESSION TAG	UNP Q11034
A	-1	MET	-	EXPRESSION TAG	UNP Q11034
A	0	VAL	-	EXPRESSION TAG	UNP Q11034
B	-27	MET	-	EXPRESSION TAG	UNP Q11034
B	-26	LYS	-	EXPRESSION TAG	UNP Q11034
B	-25	HIS	-	EXPRESSION TAG	UNP Q11034
B	-24	HIS	-	EXPRESSION TAG	UNP 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

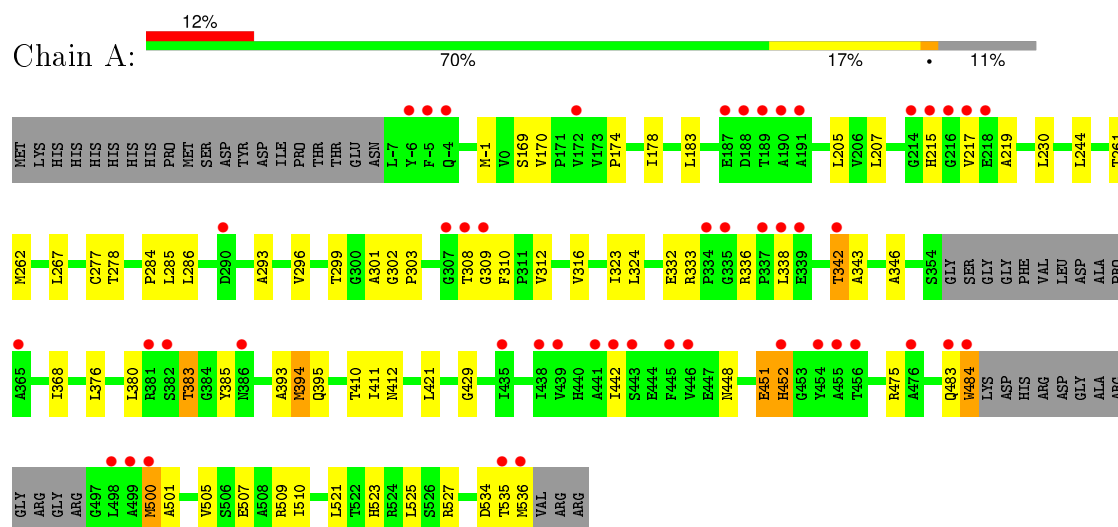
- Molecule 5 is water.

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

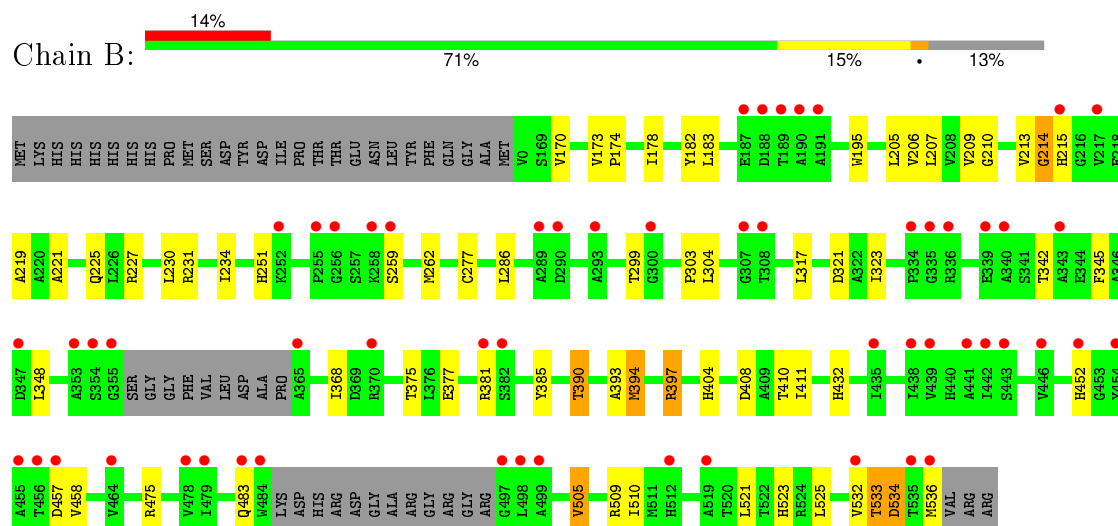
### 3 Residue-property plots [i](#)

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: Protein Rv1364c/MT1410



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



## 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 0.275 , 0.309	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 , 37.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	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.

<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

### 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 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	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

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 9.

All (91) 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: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	Interatomic distance (Å)	Clash overlap (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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:VAL:CG1	1:B:234:ILE:HD13	2.40	0.45
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%)	30	56
All	All	692/798 (87%)	665 (96%)	25 (4%)	2 (0%)	46	72

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%)	29	54
1	B	263/304 (86%)	245 (93%)	18 (7%)	20	39
All	All	531/608 (87%)	499 (94%)	32 (6%)	24	47

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

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Mol	Chain	Res	Type
1	A	534	ASP
1	A	535	THR
1	A	536	MET
1	B	170	VAL
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 [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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	0.91	0	6,6,6	0.20	0
3	GOL	A	540	-	5,5,5	0.64	0	5,5,5	0.58	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.

No monomer is involved in short contacts.

## 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	354/399 (88%)	0.77	48 (13%)  	46, 53, 69, 88	0
1	B	348/399 (87%)	0.90	57 (16%)  	47, 54, 64, 90	0
All	All	702/798 (87%)	0.83	105 (14%)  	46, 54, 67, 90	0

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	484	TRP	7.5
1	A	484	TRP	6.2
1	B	190	ALA	6.0
1	B	191	ALA	6.0
1	A	535	THR	5.8
1	B	215	HIS	5.7
1	B	457	ASP	5.6
1	A	215	HIS	5.4
1	B	188	ASP	5.1
1	A	365	ALA	5.0
1	B	335	GLY	4.9
1	A	217	VAL	4.8
1	B	189	THR	4.7
1	A	456	THR	4.6
1	A	218	GLU	4.5
1	B	217	VAL	4.4
1	B	382	SER	4.3
1	A	188	ASP	4.2
1	B	535	THR	4.2
1	B	252	LYS	4.1
1	A	-5	PHE	4.1
1	A	191	ALA	4.1
1	B	365	ALA	3.9
1	A	536	MET	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	382	SER	3.7
1	B	340	ALA	3.7
1	B	307	GLY	3.6
1	B	355	GLY	3.6
1	B	255	PRO	3.6
1	B	536	MET	3.6
1	A	-6	TYR	3.6
1	B	456	THR	3.5
1	B	334	PRO	3.4
1	A	307	GLY	3.4
1	A	381	ARG	3.4
1	A	190	ALA	3.4
1	B	339	GLU	3.4
1	A	439	VAL	3.4
1	A	443	SER	3.3
1	B	300	GLY	3.3
1	B	497	GLY	3.3
1	B	442	ILE	3.3
1	A	454	TYR	3.3
1	A	216	GLY	3.2
1	A	335	GLY	3.2
1	A	334	PRO	3.2
1	A	338	LEU	3.2
1	B	454	TYR	3.1
1	A	446	VAL	3.1
1	B	308	THR	3.0
1	A	441	ALA	3.0
1	B	455	ALA	2.9
1	B	381	ARG	2.9
1	A	455	ALA	2.8
1	B	443	SER	2.8
1	B	439	VAL	2.8
1	B	483	GLN	2.8
1	A	214	GLY	2.7
1	A	339	GLU	2.7
1	B	370	ARG	2.7
1	B	498	LEU	2.7
1	A	499	ALA	2.7
1	A	189	THR	2.7
1	B	479	ILE	2.7
1	A	483	GLN	2.6
1	A	438	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	500	MET	2.6
1	A	498	LEU	2.6
1	A	-4	GLN	2.6
1	A	386	ASN	2.5
1	A	442	ILE	2.5
1	B	438	ILE	2.5
1	B	446	VAL	2.5
1	A	308	THR	2.5
1	A	187	GLU	2.5
1	B	259	SER	2.5
1	B	478	VAL	2.5
1	B	532	VAL	2.5
1	B	441	ALA	2.4
1	A	337	PRO	2.4
1	A	342	THR	2.4
1	B	336	ARG	2.4
1	A	309	GLY	2.4
1	B	435	ILE	2.4
1	A	172	VAL	2.3
1	B	354	SER	2.3
1	B	289	ALA	2.3
1	B	512	HIS	2.3
1	A	435	ILE	2.3
1	B	293	ALA	2.2
1	A	476	ALA	2.2
1	B	353	ALA	2.2
1	B	347	ASP	2.2
1	B	452	HIS	2.2
1	A	290	ASP	2.2
1	B	258	LYS	2.2
1	B	187	GLU	2.2
1	B	290	ASP	2.1
1	B	519	ALA	2.1
1	B	256	GLY	2.1
1	B	343	ALA	2.1
1	B	464	VAL	2.1
1	B	499	ALA	2.1
1	A	445	PHE	2.0
1	A	452	HIS	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( $\text{\AA}^2$ )	Q<0.9
4	MN	A	2	1/1	0.93	0.18	0.59	73,73,73,73	0
4	MN	A	541	1/1	0.90	0.10	-1.16	69,69,69,69	0
4	MN	B	4	1/1	0.88	0.14	-1.61	72,72,72,72	0
3	GOL	A	540	6/6	0.93	0.17	-2.60	36,42,50,54	0
2	SO4	A	1	5/5	0.98	0.11	-2.96	42,43,46,47	0
4	MN	B	3	1/1	0.94	0.04	-3.01	67,67,67,67	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.