



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:14 PM GMT

PDB ID : 4O9X  
Title : Crystal Structure of TcdB2-TccC3  
Authors : Meusch, D.; Gatsogiannis, C.; Efremov, R.G.; Lang, A.E.; Hofnagel, O.; Vetter, I.R.; Aktories, K.; Raunser, S.  
Deposited on : 2014-01-03  
Resolution : 2.17 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

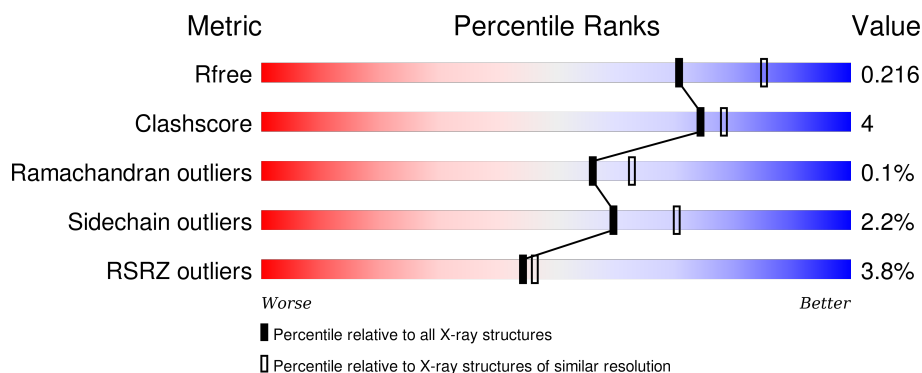
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 2.17 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	5130 (2.20-2.16)
Clashscore	102246	5965 (2.20-2.16)
Ramachandran outliers	100387	5863 (2.20-2.16)
Sidechain outliers	100360	5864 (2.20-2.16)
RSRZ outliers	91569	5142 (2.20-2.16)

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	2191	

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	HG	A	2204	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17581 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 TcdB2, TccC3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	2082	Total	C	N	O	S	0	0	0
			16589	10392	2939	3225	33			

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP Q8GF99
A	2	GLY	-	EXPRESSION TAG	UNP Q8GF99
A	3	SER	-	EXPRESSION TAG	UNP Q8GF99
A	4	SER	-	EXPRESSION TAG	UNP Q8GF99
A	5	HIS	-	EXPRESSION TAG	UNP Q8GF99
A	6	HIS	-	EXPRESSION TAG	UNP Q8GF99
A	7	HIS	-	EXPRESSION TAG	UNP Q8GF99
A	8	HIS	-	EXPRESSION TAG	UNP Q8GF99
A	9	HIS	-	EXPRESSION TAG	UNP Q8GF99
A	10	HIS	-	EXPRESSION TAG	UNP Q8GF99
A	11	SER	-	EXPRESSION TAG	UNP Q8GF99
A	12	SER	-	EXPRESSION TAG	UNP Q8GF99
A	13	GLY	-	EXPRESSION TAG	UNP Q8GF99
A	14	LEU	-	EXPRESSION TAG	UNP Q8GF99
A	15	VAL	-	EXPRESSION TAG	UNP Q8GF99
A	16	PRO	-	EXPRESSION TAG	UNP Q8GF99
A	17	ARG	-	EXPRESSION TAG	UNP Q8GF99
A	18	GLY	-	EXPRESSION TAG	UNP Q8GF99
A	19	SER	-	EXPRESSION TAG	UNP Q8GF99
A	20	HIS	-	EXPRESSION TAG	UNP Q8GF99
A	21	MET	-	EXPRESSION TAG	UNP Q8GF99
A	22	ALA	-	EXPRESSION TAG	UNP Q8GF99
A	23	SER	-	EXPRESSION TAG	UNP Q8GF99
A	24	MET	-	EXPRESSION TAG	UNP Q8GF99
A	25	THR	-	EXPRESSION TAG	UNP Q8GF99
A	26	GLY	-	EXPRESSION TAG	UNP Q8GF99
A	27	GLY	-	EXPRESSION TAG	UNP Q8GF99

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Chain	Residue	Modelled	Actual	Comment	Reference
A	28	GLN	-	EXPRESSION TAG	UNP Q8GF99
A	29	GLN	-	EXPRESSION TAG	UNP Q8GF99
A	30	MET	-	EXPRESSION TAG	UNP Q8GF99
A	31	GLY	-	EXPRESSION TAG	UNP Q8GF99
A	32	ARG	-	EXPRESSION TAG	UNP Q8GF99
A	33	GLY	-	EXPRESSION TAG	UNP Q8GF99
A	34	SER	-	EXPRESSION TAG	UNP Q8GF99
A	1509	PRO	-	LINKER	UNP Q8GF99
A	1510	GLY	-	LINKER	UNP Q8GF99
A	1511	SER	-	LINKER	UNP Q8GF99
A	1512	ARG	-	LINKER	UNP Q8GF99
A	1513	PRO	-	LINKER	UNP Q8GF99

- Molecule 2 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	9	Total Hg 9 9	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	983	Total O 983 983	0	0

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.

- Chain A: 85% 4% 10% 5%

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	232.36Å 232.36Å 143.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.77 – 2.17 29.77 – 2.17	Depositor EDS
% Data completeness (in resolution range)	92.6 (29.77-2.17) 93.1 (29.77-2.17)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.64 (at 2.18Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, $R_{free}$	0.192 , 0.215 0.193 , 0.216	Depositor DCC
$R_{free}$ test set	10883 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	41.9	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 45.6	EDS
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 217656 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	17581	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.91% 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: HG

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	0.42	0/17001	0.59	1/23181 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	709	LEU	CB-CG-CD1	5.08	119.63	111.00

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	16589	0	15934	129	0
2	A	9	0	0	0	0
3	A	983	0	0	18	0
All	All	17581	0	15934	129	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 4.

All (129) 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:1264:PHE:HE2	1:A:1285:CYS:HG	1.07	1.03
1:A:1478:LYS:HG3	1:A:1813:GLU:HG3	1.47	0.96
1:A:487:LEU:HD13	1:A:496:LEU:HD23	1.68	0.75
1:A:1959:GLN:NE2	3:A:2820:HOH:O	2.22	0.72
1:A:436:LYS:HG2	1:A:492:GLY:H	1.55	0.71
1:A:1242:THR:HG22	1:A:1285:CYS:SG	2.31	0.70
1:A:495:GLN:HG3	1:A:496:LEU:HD12	1.74	0.70
1:A:1264:PHE:HE2	1:A:1285:CYS:SG	2.14	0.69
1:A:1692:ARG:NH1	3:A:3192:HOH:O	2.25	0.69
1:A:301:ARG:NH1	1:A:311:PHE:O	2.26	0.67
1:A:1112:THR:HG22	1:A:1157:VAL:HG12	1.77	0.67
1:A:685:PRO:HG2	1:A:688:LEU:HD12	1.76	0.66
1:A:1284:GLN:NE2	3:A:3101:HOH:O	2.32	0.62
1:A:370:ASN:HB2	1:A:382:PHE:HB3	1.79	0.62
1:A:514:ASP:HB2	1:A:515:GLY:HA2	1.80	0.62
1:A:301:ARG:NE	3:A:2931:HOH:O	2.30	0.61
1:A:1835:ARG:HG3	1:A:1848:ARG:HH21	1.64	0.61
1:A:1683:ARG:HG2	1:A:1692:ARG:HG2	1.82	0.61
1:A:1605:ILE:HD12	1:A:1945:SER:HA	1.83	0.60
1:A:1992:LEU:HD21	1:A:2143:THR:HG22	1.83	0.60
1:A:476:SER:HB3	1:A:491:ASN:HB3	1.84	0.59
1:A:824:HIS:HD2	3:A:3000:HOH:O	1.86	0.58
1:A:543:ALA:HB2	1:A:564:ASP:HB2	1.86	0.58
1:A:1493:ASN:ND2	3:A:3167:HOH:O	2.31	0.58
1:A:772:TYR:CE1	1:A:773:LEU:HD22	2.39	0.57
1:A:134:GLU:N	1:A:134:GLU:OE1	2.25	0.57
1:A:729:LYS:HD2	1:A:778:HIS:CD2	2.40	0.57
1:A:1217:LEU:HD23	3:A:3083:HOH:O	2.05	0.56
1:A:2177:MET:HG3	1:A:2178:VAL:HG22	1.88	0.55
1:A:1569:GLN:HG2	1:A:1590:LEU:HD23	1.87	0.55
1:A:219:ALA:HB1	1:A:329:TYR:CD1	2.42	0.55
1:A:742:VAL:HG23	1:A:1382:GLY:HA3	1.89	0.55
1:A:1785:LYS:NZ	1:A:1803:ASP:OD2	2.40	0.54
1:A:1545:ARG:NH2	1:A:1552:PRO:HG3	2.22	0.54
1:A:668:PHE:CE1	1:A:679:PRO:HG3	2.43	0.53
1:A:704:LEU:HA	1:A:778:HIS:HE1	1.73	0.53
1:A:246:CYS:HG	1:A:319:CYS:HG	1.58	0.52
1:A:1266:PRO:HB3	1:A:1285:CYS:SG	2.49	0.52
1:A:749:ARG:NH2	1:A:757:ASP:OD2	2.40	0.52
1:A:1692:ARG:NH2	1:A:1708:GLN:OE1	2.34	0.52
1:A:772:TYR:CD1	1:A:773:LEU:HD22	2.45	0.51
1:A:1381:ASP:HB3	1:A:1387:LEU:HD21	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:802:ARG:NH2	1:A:858:GLU:OE2	2.44	0.51
1:A:1649:GLU:OE1	1:A:1658:ARG:NH2	2.42	0.50
1:A:1176:LYS:HD3	1:A:1185:LYS:HD3	1.93	0.50
1:A:889:ASP:OD2	1:A:892:ARG:NH2	2.41	0.49
1:A:1421:THR:HG22	1:A:1438:PRO:HB3	1.94	0.49
1:A:563:ARG:O	1:A:568:LYS:NZ	2.37	0.49
1:A:301:ARG:NH1	1:A:310:ALA:O	2.42	0.49
1:A:704:LEU:HB3	1:A:706:VAL:HG22	1.94	0.49
1:A:478:ILE:HB	1:A:481:LEU:HD12	1.93	0.49
1:A:1545:ARG:HH21	1:A:1552:PRO:HG3	1.77	0.49
1:A:772:TYR:CD1	1:A:803:GLY:HA3	2.48	0.49
1:A:1042:SER:HB2	1:A:1146:LEU:HD13	1.95	0.49
1:A:751:SER:OG	1:A:773:LEU:HD23	2.13	0.49
1:A:2078:ASP:OD1	1:A:2078:ASP:N	2.35	0.48
1:A:1917:ARG:NH2	3:A:2643:HOH:O	2.36	0.48
1:A:2041:ARG:NH2	3:A:3263:HOH:O	2.46	0.48
1:A:1603:GLU:HG2	1:A:1609:ARG:HG3	1.95	0.48
1:A:704:LEU:HD12	1:A:778:HIS:HE1	1.79	0.47
1:A:1114:PHE:HB2	1:A:1156:TRP:HB2	1.96	0.47
1:A:562:THR:HG22	1:A:568:LYS:HD2	1.96	0.47
1:A:1400:GLN:NE2	1:A:1410:ASN:O	2.47	0.47
1:A:1057:ALA:HA	1:A:1058:GLY:HA2	1.66	0.47
1:A:2115:GLY:HA2	1:A:2135:TYR:CD1	2.50	0.47
1:A:326:ARG:NH2	3:A:2895:HOH:O	2.46	0.47
1:A:1293:TRP:O	1:A:1352:ARG:NH1	2.48	0.47
1:A:1083:GLN:HG2	1:A:1112:THR:CG2	2.46	0.46
1:A:75:LEU:HB3	1:A:91:LEU:HB3	1.97	0.46
1:A:689:ARG:NH1	1:A:691:ASP:OD2	2.47	0.46
1:A:167:ARG:NH2	3:A:2920:HOH:O	2.48	0.46
1:A:1969:HIS:HB3	1:A:1981:TRP:CE2	2.50	0.46
1:A:451:TRP:NE1	1:A:479:PRO:HG2	2.30	0.46
1:A:2024:SER:HA	1:A:2025:ASN:HA	1.63	0.46
1:A:491:ASN:HA	1:A:492:GLY:HA2	1.66	0.45
1:A:834:ARG:NH1	3:A:3004:HOH:O	2.49	0.45
1:A:1553:ASP:OD2	1:A:1917:ARG:HD2	2.15	0.45
1:A:760:ALA:O	1:A:763:LEU:HB2	2.17	0.45
1:A:414:HIS:CD2	1:A:416:GLN:HE22	2.35	0.45
1:A:998:PRO:HD3	3:A:2821:HOH:O	2.16	0.45
1:A:704:LEU:HA	1:A:778:HIS:CE1	2.51	0.45
1:A:809:GLU:HB3	1:A:811:GLU:HG2	1.99	0.45
1:A:652:SER:N	1:A:653:GLY:HA2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:ILE:HG12	1:A:714:PRO:HA	2.00	0.44
1:A:242:ASP:OD2	3:A:2923:HOH:O	2.21	0.44
1:A:1705:GLN:NE2	3:A:3195:HOH:O	2.50	0.44
1:A:1771:SER:HB3	1:A:1784:ILE:HB	1.99	0.44
1:A:406:ASP:N	1:A:406:ASP:OD1	2.44	0.44
1:A:921:TYR:CZ	1:A:957:ARG:HD2	2.52	0.44
1:A:301:ARG:NH2	3:A:2931:HOH:O	2.51	0.44
1:A:2134:ARG:HB2	1:A:2135:TYR:H	1.61	0.43
1:A:468:THR:OG1	1:A:469:TRP:N	2.49	0.43
1:A:454:ARG:HG2	1:A:472:MET:HA	1.98	0.43
1:A:1402:ASN:ND2	1:A:1408:ILE:HD11	2.34	0.43
1:A:869:PHE:CG	1:A:927:ARG:HB2	2.53	0.43
1:A:570:LYS:HD2	1:A:570:LYS:HA	1.76	0.43
1:A:1164:ASP:HB2	1:A:1176:LYS:HB2	2.00	0.43
1:A:1574:ARG:NH2	1:A:1942:SER:O	2.31	0.43
1:A:1714:GLN:NE2	3:A:3200:HOH:O	2.23	0.43
1:A:448:LYS:HA	1:A:449:GLY:HA2	1.41	0.43
1:A:1293:TRP:CE2	1:A:1352:ARG:HD3	2.54	0.42
1:A:867:ALA:O	1:A:951:ARG:NH2	2.53	0.42
1:A:799:ARG:NH1	1:A:819:GLU:OE2	2.52	0.42
1:A:1201:ARG:HG2	1:A:1207:THR:HG22	2.02	0.42
1:A:440:LEU:HG	1:A:752:SER:HB3	2.00	0.42
1:A:686:GLU:HA	1:A:687:GLY:HA2	1.65	0.42
1:A:755:TRP:CE3	1:A:772:TYR:HE2	2.37	0.42
1:A:522:PRO:HD2	1:A:550:LEU:HD21	2.01	0.42
1:A:1655:LYS:HD2	1:A:1655:LYS:HA	1.81	0.42
1:A:435:LEU:HD13	1:A:490:ILE:HD13	2.02	0.42
1:A:620:ARG:NH1	3:A:2979:HOH:O	2.25	0.42
1:A:536:GLN:HG2	1:A:578:ILE:HD11	2.02	0.41
1:A:1293:TRP:CZ2	1:A:1352:ARG:HD3	2.55	0.41
1:A:1286:GLN:HA	1:A:1358:LEU:O	2.21	0.41
1:A:831:ALA:HA	1:A:832:PRO:HD3	1.96	0.41
1:A:958:ILE:HD13	1:A:997:TYR:CE1	2.56	0.41
1:A:1551:ASP:HA	1:A:1552:PRO:HD3	1.97	0.41
1:A:435:LEU:HD23	1:A:435:LEU:HA	1.88	0.41
1:A:1234:LEU:HD23	1:A:1240:PRO:HA	2.03	0.41
1:A:451:TRP:CE2	1:A:479:PRO:HG2	2.56	0.41
1:A:1143:THR:HG23	1:A:1144:ASN:O	2.21	0.41
1:A:772:TYR:N	1:A:772:TYR:CD2	2.89	0.40
1:A:652:SER:CB	1:A:654:PRO:HD2	2.51	0.40
1:A:496:LEU:O	1:A:496:LEU:HD22	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1588:ASN:HA	1:A:1605:ILE:HG12	2.04	0.40
1:A:1785:LYS:HE3	1:A:1785:LYS:HB2	1.90	0.40
1:A:2025:ASN:ND2	1:A:2026:ASN:HB3	2.36	0.40
1:A:1518:ASP:HA	1:A:1519:PRO:HD2	1.95	0.40
1:A:640:ASN:HA	1:A:641:PRO:HD2	1.90	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	2076/2191 (95%)	2012 (97%)	61 (3%)	3 (0%)	56 63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2149	GLY
1	A	496	LEU
1	A	486	SER

### 5.3.2 Protein sidechains [i](#)

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	1795/1889 (95%)	1755 (98%)	40 (2%)	60 70

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	142	ASP
1	A	167	ARG
1	A	204	SER
1	A	253	HIS
1	A	293	TYR
1	A	317	TRP
1	A	374	SER
1	A	420	VAL
1	A	446	GLN
1	A	457	GLN
1	A	464	SER
1	A	477	VAL
1	A	481	LEU
1	A	488	VAL
1	A	496	LEU
1	A	563	ARG
1	A	634	GLN
1	A	665	LEU
1	A	682	LEU
1	A	709	LEU
1	A	771	CYS
1	A	772	TYR
1	A	830	ASN
1	A	884	LEU
1	A	940	TYR
1	A	964	CYS
1	A	970	LEU
1	A	1078	ARG
1	A	1094	THR
1	A	1097	LEU
1	A	1117	SER
1	A	1285	CYS
1	A	1503	VAL
1	A	1601	CYS
1	A	1848	ARG
1	A	1974	ILE
1	A	2025	ASN
1	A	2078	ASP
1	A	2110	GLU
1	A	2159	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	237	GLN
1	A	416	GLN
1	A	1400	GLN
1	A	1750	GLN
1	A	2096	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 9 ligands modelled in this entry, 9 are monoatomic - leaving 0 for Mogul analysis.

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	2082/2191 (95%)	-0.20	80 (3%)	44 46	22, 37, 70, 141	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	492	GLY	26.8
1	A	566	PHE	14.2
1	A	565	GLY	12.6
1	A	564	ASP	10.3
1	A	491	ASN	9.5
1	A	567	ALA	7.8
1	A	1998	GLY	7.2
1	A	569	GLY	6.9
1	A	514	ASP	6.3
1	A	481	LEU	6.3
1	A	2026	ASN	5.7
1	A	2023	ALA	5.6
1	A	2025	ASN	5.5
1	A	568	LYS	5.1
1	A	2024	SER	5.0
1	A	480	SER	4.9
1	A	763	LEU	4.7
1	A	768	THR	4.7
1	A	476	SER	4.6
1	A	1351	PRO	4.5
1	A	1999	ALA	4.5
1	A	1997	LYS	4.5
1	A	1295	PRO	4.5
1	A	832	PRO	4.1
1	A	563	ARG	3.9
1	A	767	GLN	3.9
1	A	479	PRO	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	306	LYS	3.6
1	A	490	ILE	3.6
1	A	773	LEU	3.3
1	A	1655	LYS	3.3
1	A	515	GLY	3.1
1	A	766	GLY	3.1
1	A	772	TYR	3.1
1	A	765	THR	3.1
1	A	506	ARG	3.1
1	A	1518	ASP	3.0
1	A	2078	ASP	3.0
1	A	2045	ASN	3.0
1	A	2077	GLU	2.9
1	A	307	THR	2.9
1	A	543	ALA	2.9
1	A	1409	ILE	2.8
1	A	833	GLU	2.7
1	A	448	LYS	2.7
1	A	482	GLN	2.7
1	A	828	GLN	2.7
1	A	411	HIS	2.7
1	A	2022	GLN	2.6
1	A	493	ASP	2.6
1	A	1996	ASP	2.6
1	A	1403	GLU	2.6
1	A	359	ASN	2.6
1	A	913	ASP	2.6
1	A	1932	GLN	2.6
1	A	495	GLN	2.5
1	A	392	GLY	2.5
1	A	2047	THR	2.5
1	A	496	LEU	2.4
1	A	462	ILE	2.4
1	A	390	GLN	2.4
1	A	510	SER	2.4
1	A	715	HIS	2.4
1	A	414	HIS	2.3
1	A	764	THR	2.3
1	A	1057	ALA	2.3
1	A	2027	ALA	2.3
1	A	709	LEU	2.3
1	A	509	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	1410	ASN	2.2
1	A	513	PRO	2.2
1	A	775	PHE	2.2
1	A	577	ASP	2.2
1	A	562	THR	2.2
1	A	834	ARG	2.1
1	A	1901	SER	2.1
1	A	1902	ASP	2.1
1	A	2097	LEU	2.1
1	A	878	ASP	2.1
1	A	1549	ASN	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	HG	A	2204	1/1	0.39	0.23	6.35	348,348,348,348	0
2	HG	A	2208	1/1	0.98	0.04	-1.72	96,96,96,96	0
2	HG	A	2203	1/1	0.96	0.04	-1.72	113,113,113,113	0
2	HG	A	2202	1/1	0.99	0.04	-1.94	87,87,87,87	0
2	HG	A	2209	1/1	0.82	0.09	-1.95	213,213,213,213	0
2	HG	A	2201	1/1	0.99	0.03	-2.32	62,62,62,62	0
2	HG	A	2206	1/1	0.94	0.04	-2.56	117,117,117,117	0
2	HG	A	2205	1/1	0.98	0.04	-4.07	101,101,101,101	0
2	HG	A	2207	1/1	0.98	0.04	-	60,60,60,60	0



## 6.5 Other polymers [i](#)

There are no such residues in this entry.