



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 06:48 PM UTC

PDB ID : 4O9X / pdb_00004o9x
Title : Crystal Structure of TcdB2-TccC3
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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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

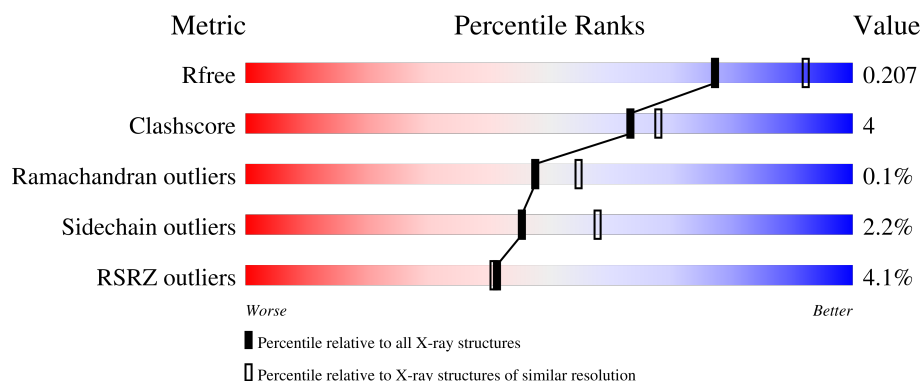
1 Overall quality at a glance

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}	180053	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	

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 (CCD ID: HG) (formula: Hg).

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

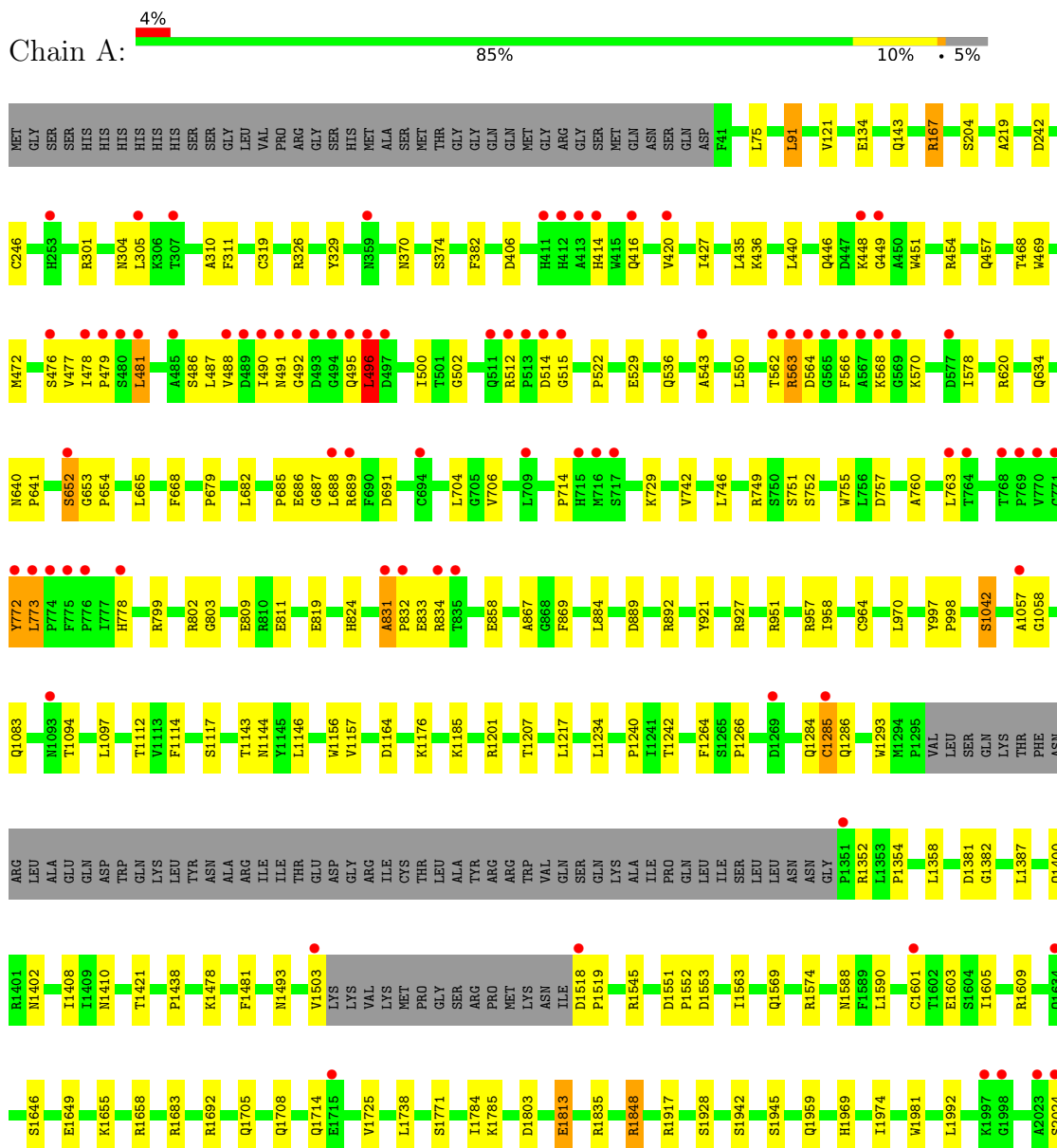
- Molecule 3 is water.

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

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: TcdB2, TccC3





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	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) 93.0 (29.77-2.17)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.18Å)	Xtriage
Refinement program	PHENIX 1.8.1 _1168	Depositor
R, R_{free}	0.192 , 0.215 0.184 , 0.207	Depositor DCC
R_{free} test set	10883 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	41.9	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17581	wwPDB-VP
Average B, all atoms (Å ²)	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.*

¹Intensities estimated from amplitudes.

²Theoretical 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

5.1 Standard geometry

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	0.49	0/17001	0.81	16/23181 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	652	SER	N-CA-C	8.08	121.63	111.69
1	A	512	ARG	CA-C-N	-7.66	112.32	120.66
1	A	512	ARG	C-N-CA	-7.66	112.32	120.66
1	A	2164	ASP	CA-C-N	6.51	126.20	119.56
1	A	2164	ASP	C-N-CA	6.51	126.20	119.56
1	A	831	ALA	CA-C-N	6.48	125.98	119.24
1	A	831	ALA	C-N-CA	6.48	125.98	119.24
1	A	1354	PRO	O-C-N	5.81	123.98	121.31
1	A	1042	SER	N-CA-C	5.79	117.86	109.07
1	A	1928	SER	CA-C-N	5.47	124.66	118.97
1	A	1928	SER	C-N-CA	5.47	124.66	118.97
1	A	502	GLY	CA-C-N	5.43	125.26	119.28
1	A	502	GLY	C-N-CA	5.43	125.26	119.28
1	A	1481	PHE	N-CA-C	5.39	117.38	109.24
1	A	563	ARG	N-CA-C	5.30	118.32	110.48
1	A	1646	SER	N-CA-C	5.23	116.79	108.79

There are no chirality outliers.

There are no planarity outliers.

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	16589	0	15934	131	0
2	A	9	0	0	0	0
3	A	983	0	0	18	0
All	All	17581	0	15934	131	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 (131) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:GLU:OE1	1:A:134:GLU:N	2.25	0.57
1:A:772:TYR:CE1	1:A:773:LEU:HD22	2.39	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:563:ARG:O	1:A:568:LYS:NZ	2.37	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:1545:ARG:NH2	1:A:1552:PRO:HG3	2.22	0.54
1:A:1785:LYS:NZ	1:A:1803:ASP:OD2	2.40	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:749:ARG:NH2	1:A:757:ASP:OD2	2.40	0.52
1:A:1266:PRO:HB3	1:A:1285:CYS:SG	2.49	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
1:A:802:ARG:NH2	1:A:858:GLU:OE2	2.44	0.51
1:A:246:CYS:HG	1:A:319:CYS:HG	1.59	0.50
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: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:1057:ALA:HA	1:A:1058:GLY:HA2	1.66	0.47
1:A:1400:GLN:NE2	1:A:1410:ASN:O	2.47	0.47
1:A:2115:GLY:HA2	1:A:2135:TYR:CD1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:75:LEU:HB3	1:A:91:LEU:HB3	1.97	0.46
1:A:1083:GLN:HG2	1:A:1112:THR:CG2	2.46	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:451:TRP:NE1	1:A:479:PRO:HG2	2.30	0.46
1:A:1969:HIS:HB3	1:A:1981:TRP:CE2	2.50	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:760:ALA:O	1:A:763:LEU:HB2	2.17	0.45
1:A:1553:ASP:OD2	1:A:1917:ARG:HD2	2.15	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
1:A:242:ASP:OD2	3:A:2923:HOH:O	2.21	0.44
1:A:427:ILE:HG12	1:A:714:PRO:HA	2.00	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:OD1	1:A:406:ASP:N	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:454:ARG:HG2	1:A:472:MET:HA	1.98	0.43
1:A:468:THR:OG1	1:A:469:TRP:N	2.49	0.43
1:A:2134:ARG:HB2	1:A:2135:TYR:H	1.61	0.43
1:A:570:LYS:HD2	1:A:570:LYS:HA	1.76	0.43
1:A:869:PHE:CG	1:A:927:ARG:HB2	2.53	0.43
1:A:1402:ASN:ND2	1:A:1408:ILE:HD11	2.34	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:522:PRO:HD2	1:A:550:LEU:HD21	2.01	0.42
1:A:755:TRP:CE3	1:A:772:TYR:HE2	2.37	0.42
1:A:1655:LYS:HA	1:A:1655:LYS:HD2	1.81	0.42
1:A:435:LEU:HD13	1:A:490:ILE:HD13	2.02	0.42
1:A:564:ASP:C	1:A:566:PHE:H	2.28	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:831:ALA:HA	1:A:832:PRO:HD3	1.96	0.41
1:A:1286:GLN:HA	1:A:1358:LEU:O	2.21	0.41
1:A:958:ILE:HD13	1:A:997:TYR:CE1	2.56	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:1551:ASP:HA	1:A:1552:PRO:HD3	1.97	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:652:SER:CB	1:A:654:PRO:HD2	2.51	0.40
1:A:772:TYR:CD2	1:A:772:TYR:N	2.89	0.40
1:A:496:LEU:O	1:A:496:LEU:HD22	2.21	0.40
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:304:ASN:C	1:A:305:LEU:HD12	2.47	0.40
1:A:640:ASN:HA	1:A:641:PRO:HD2	1.90	0.40
1:A:1518:ASP:HA	1:A:1519:PRO:HD2	1.95	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	2076/2191 (95%)	2012 (97%)	61 (3%)	3 (0%)	48	55

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 ⓘ

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%)	45 58

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

Mol	Chain	Res	Type
1	A	91	LEU
1	A	121	VAL
1	A	143	GLN
1	A	167	ARG
1	A	204	SER
1	A	374	SER
1	A	420	VAL
1	A	446	GLN
1	A	457	GLN
1	A	477	VAL
1	A	481	LEU
1	A	488	VAL
1	A	496	LEU
1	A	500	ILE
1	A	529	GLU
1	A	634	GLN
1	A	665	LEU
1	A	682	LEU
1	A	746	LEU
1	A	772	TYR
1	A	773	LEU
1	A	833	GLU

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Mol	Chain	Res	Type
1	A	884	LEU
1	A	964	CYS
1	A	970	LEU
1	A	1094	THR
1	A	1097	LEU
1	A	1117	SER
1	A	1285	CYS
1	A	1503	VAL
1	A	1563	ILE
1	A	1601	CYS
1	A	1725	VAL
1	A	1738	LEU
1	A	1813	GLU
1	A	1848	ARG
1	A	1974	ILE
1	A	2078	ASP
1	A	2110	GLU
1	A	2159	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	358	HIS
1	A	405	GLN
1	A	416	GLN
1	A	824	HIS
1	A	896	ASN
1	A	902	GLN
1	A	977	GLN
1	A	1098	GLN
1	A	1141	GLN
1	A	1221	GLN
1	A	1228	ASN
1	A	1400	GLN
1	A	1454	GLN
1	A	1750	GLN
1	A	1767	GLN
1	A	1832	GLN
1	A	1938	ASN
1	A	2025	ASN
1	A	2037	ASN

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 oligosaccharides 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, 95th 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 ²)	Q<0.9
1	A	2082/2191 (95%)	-0.16	85 (4%) 41 41	22, 37, 70, 141	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	492	GLY	10.8
1	A	496	LEU	6.4
1	A	491	ASN	6.4
1	A	772	TYR	6.2
1	A	490	ILE	6.0
1	A	566	PHE	5.7
1	A	513	PRO	5.6
1	A	567	ALA	5.4
1	A	494	GLY	5.2
1	A	773	LEU	5.1
1	A	2023	ALA	5.0
1	A	1518	ASP	5.0
1	A	564	ASP	4.9
1	A	768	THR	4.7
1	A	479	PRO	4.5
1	A	2045	ASN	4.2
1	A	778	HIS	4.1
1	A	1285	CYS	4.1
1	A	770	VAL	4.0
1	A	771	CYS	4.0
1	A	565	GLY	4.0
1	A	514	ASP	4.0
1	A	774	PRO	3.9
1	A	495	GLN	3.9
1	A	485	ALA	3.7
1	A	481	LEU	3.6
1	A	775	PHE	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	715	HIS	3.4
1	A	563	ARG	3.4
1	A	569	GLY	3.4
1	A	716	MET	3.3
1	A	694	CYS	3.3
1	A	480	SER	3.2
1	A	562	THR	3.2
1	A	515	GLY	3.1
1	A	414	HIS	3.1
1	A	511	GLN	3.1
1	A	831	ALA	3.1
1	A	307	THR	3.0
1	A	1269	ASP	3.0
1	A	2027	ALA	3.0
1	A	488	VAL	2.9
1	A	568	LYS	2.9
1	A	1503	VAL	2.9
1	A	413	ALA	2.8
1	A	489	ASP	2.8
1	A	305	LEU	2.8
1	A	717	SER	2.8
1	A	689	ARG	2.8
1	A	688	LEU	2.8
1	A	776	PRO	2.8
1	A	709	LEU	2.8
1	A	763	LEU	2.7
1	A	493	ASP	2.7
1	A	835	THR	2.7
1	A	359	ASN	2.7
1	A	476	SER	2.7
1	A	764	THR	2.6
1	A	2024	SER	2.6
1	A	832	PRO	2.5
1	A	834	ARG	2.5
1	A	411	HIS	2.4
1	A	412	HIS	2.4
1	A	1998	GLY	2.4
1	A	1715	GLU	2.4
1	A	2025	ASN	2.4
1	A	420	VAL	2.3
1	A	2050	THR	2.3
1	A	769	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	497	ASP	2.3
1	A	1997	LYS	2.3
1	A	478	ILE	2.2
1	A	448	LYS	2.2
1	A	1057	ALA	2.2
1	A	1634	GLN	2.2
1	A	253	HIS	2.2
1	A	652	SER	2.2
1	A	512	ARG	2.2
1	A	543	ALA	2.2
1	A	1601	CYS	2.2
1	A	416	GLN	2.1
1	A	577	ASP	2.1
1	A	1351	PRO	2.0
1	A	1093	ASN	2.0
1	A	449	GLY	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 oligosaccharides 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. The B-factors column lists the minimum, median, 95th 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	B-factors($\text{\AA}^2$)	Q<0.9
2	HG	A	2204	1/1	0.59	0.40	348,348,348,348	0
2	HG	A	2209	1/1	0.88	0.26	213,213,213,213	0
2	HG	A	2206	1/1	0.95	0.11	117,117,117,117	0
2	HG	A	2203	1/1	0.95	0.10	113,113,113,113	0
2	HG	A	2207	1/1	0.97	0.07	60,60,60,60	0
2	HG	A	2205	1/1	0.97	0.15	101,101,101,101	0
2	HG	A	2208	1/1	0.98	0.11	96,96,96,96	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	HG	A	2202	1/1	0.99	0.07	87,87,87,87	0
2	HG	A	2201	1/1	1.00	0.04	62,62,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.