



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 08:44 PM EDT

PDB ID : 7DEJ / pdb_00007dej
Title : Structure of human ORP3 ORD in apo-form
Authors : Tong, J.; Tan, L.; Im, Y.J.
Deposited on : 2020-11-04
Resolution : 2.70 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

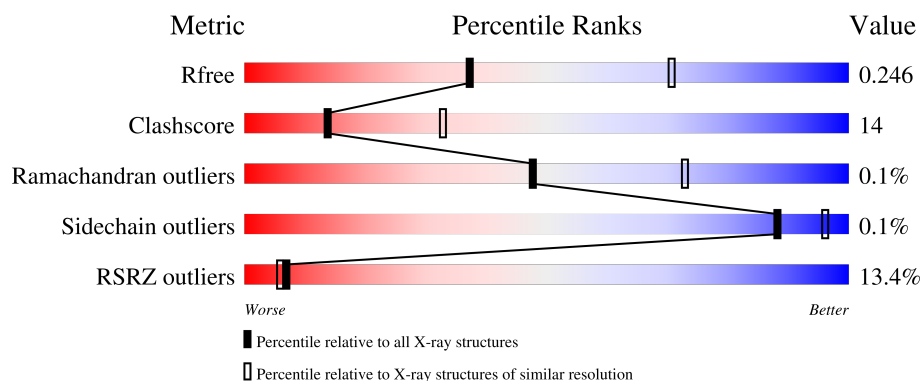
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	388	20% 64% 31% ..
1	B	388	6% 77% 19% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6210 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 Oxysterol-binding protein-related protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	0	0	0
			3072	1955	540	564	13			
1	B	376	Total	C	N	O	S	0	0	0
			3083	1961	544	565	13			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	499	GLY	-	expression tag	UNP Q9H4L5
A	500	SER	-	expression tag	UNP Q9H4L5
A	501	PRO	-	expression tag	UNP Q9H4L5
A	502	GLU	-	expression tag	UNP Q9H4L5
A	503	PHE	-	expression tag	UNP Q9H4L5
A	515	SER	CYS	engineered mutation	UNP Q9H4L5
A	520	SER	CYS	engineered mutation	UNP Q9H4L5
A	?	-	ASP	deletion	UNP Q9H4L5
A	860	GLY	ASP	engineered mutation	UNP Q9H4L5
B	499	GLY	-	expression tag	UNP Q9H4L5
B	500	SER	-	expression tag	UNP Q9H4L5
B	501	PRO	-	expression tag	UNP Q9H4L5
B	502	GLU	-	expression tag	UNP Q9H4L5
B	503	PHE	-	expression tag	UNP Q9H4L5
B	515	SER	CYS	engineered mutation	UNP Q9H4L5
B	520	SER	CYS	engineered mutation	UNP Q9H4L5
B	?	-	ASP	deletion	UNP Q9H4L5
B	860	GLY	ASP	engineered mutation	UNP Q9H4L5

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	O	0	0
			10	10		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	45	Total	O	0	0
			45	45		

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	95.91Å 95.91Å 188.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.44 – 2.70 29.44 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.44-2.70) 98.8 (29.44-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.93 (at 2.72Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.233 , 0.285 0.236 , 0.246	Depositor DCC
R_{free} test set	1417 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	57.6	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 58.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6210	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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

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	0.44	2/3162 (0.1%)	0.77	6/4284 (0.1%)
1	B	0.57	2/3173 (0.1%)	0.82	3/4298 (0.1%)
All	All	0.51	4/6335 (0.1%)	0.80	9/8582 (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	3
1	B	0	1
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	615	CYS	CB-SG	8.84	2.10	1.81
1	B	760	CYS	CB-SG	6.46	2.02	1.81
1	A	760	CYS	CB-SG	5.64	1.99	1.81
1	A	689	CYS	CB-SG	5.47	1.99	1.81

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	658	TRP	CA-C-N	-7.30	114.41	120.98
1	B	658	TRP	C-N-CA	-7.30	114.41	120.98
1	A	614	GLU	CG-CD-OE1	-6.67	103.06	118.40
1	A	853	ARG	CA-C-N	-6.49	112.36	122.67
1	A	853	ARG	C-N-CA	-6.49	112.36	122.67
1	A	614	GLU	N-CA-C	-6.33	97.48	108.69
1	A	689	CYS	CA-CB-SG	5.91	127.99	114.40
1	B	760	CYS	CA-CB-SG	5.78	127.70	114.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	634	PRO	N-CD-CG	5.50	111.46	103.20

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	601	GLY	Peptide
1	A	614	GLU	Sidechain
1	A	883	HIS	Peptide
1	B	858	SER	Peptide

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	3072	0	2959	105	0
1	B	3083	0	2972	70	0
2	A	10	0	0	7	0
2	B	45	0	0	21	0
All	All	6210	0	5931	175	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 14.

All (175) 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:B:760:CYS:SG	1:B:760:CYS:CB	2.02	1.47
1:B:615:CYS:CB	1:B:615:CYS:SG	2.10	1.39
1:B:725:ALA:HB1	2:B:901:HOH:O	1.40	1.21
1:B:558:ARG:HG2	2:B:902:HOH:O	1.52	1.09
1:B:726:LYS:N	2:B:901:HOH:O	1.84	1.08
1:B:760:CYS:SG	2:B:937:HOH:O	2.24	0.95
1:B:725:ALA:CB	2:B:901:HOH:O	2.03	0.90
1:B:726:LYS:O	2:B:901:HOH:O	1.92	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:540:LEU:HD11	1:B:657:PHE:HB2	1.58	0.86
1:B:561:GLU:OE1	2:B:902:HOH:O	1.95	0.84
1:A:513:ARG:HE	1:A:515:SER:HB2	1.40	0.84
1:B:527:SER:HB3	1:B:546:PRO:HG2	1.60	0.84
1:B:735:ILE:HD11	1:B:758:ILE:HD12	1.59	0.84
1:A:832:ILE:O	1:A:836:GLN:NE2	2.11	0.84
1:A:602:SER:O	2:A:901:HOH:O	1.95	0.83
1:A:617:ARG:HH11	1:A:619:ASP:HB3	1.46	0.79
1:B:876:LEU:C	2:B:908:HOH:O	2.27	0.78
1:A:581:ARG:NH2	1:A:679:ASP:OD2	2.17	0.78
1:B:876:LEU:O	2:B:903:HOH:O	2.03	0.77
1:A:716:CYS:SG	2:A:909:HOH:O	2.42	0.77
1:B:617:ARG:NH2	1:B:791:GLU:OE1	2.19	0.74
1:B:879:SER:HB2	2:B:903:HOH:O	1.88	0.74
1:A:615:CYS:SG	2:A:907:HOH:O	2.44	0.74
1:B:724:LYS:O	2:B:904:HOH:O	2.06	0.72
1:A:602:SER:CB	2:A:901:HOH:O	2.38	0.71
1:A:602:SER:HB2	2:A:901:HOH:O	1.89	0.71
1:A:617:ARG:HG2	1:A:620:LYS:H	1.56	0.71
1:B:614:GLU:OE2	2:B:905:HOH:O	2.10	0.70
1:B:686:VAL:HG13	1:B:703:TYR:O	1.92	0.69
1:B:812:PRO:HD2	1:B:832:ILE:HD11	1.73	0.69
1:A:686:VAL:HG22	1:A:703:TYR:O	1.93	0.69
1:A:617:ARG:HD3	1:A:620:LYS:HG3	1.76	0.68
1:B:703:TYR:HB3	1:B:724:LYS:HB2	1.76	0.67
1:A:602:SER:C	2:A:901:HOH:O	2.36	0.66
1:A:853:ARG:NH1	1:A:883:HIS:HB2	2.11	0.66
1:A:602:SER:CA	2:A:901:HOH:O	2.43	0.66
1:B:823:GLU:OE1	2:B:906:HOH:O	2.14	0.66
1:B:617:ARG:HH11	1:B:620:LYS:HG3	1.60	0.66
1:A:653:TRP:CZ3	1:A:666:PRO:HG3	2.31	0.66
1:A:871:GLU:N	1:A:871:GLU:OE1	2.30	0.65
1:A:692:ASN:HB3	1:A:695:SER:HB2	1.79	0.65
1:A:828:GLN:O	1:A:832:ILE:HD12	1.98	0.64
1:B:878:PHE:N	2:B:908:HOH:O	2.32	0.62
1:B:520:SER:HB2	1:B:783:TYR:H	1.64	0.62
1:A:854:PHE:HA	1:A:866:ASN:HD21	1.65	0.61
1:A:635:ILE:HG12	1:A:652:ARG:HB3	1.82	0.60
1:A:785:PHE:HB3	1:A:789:ALA:HB3	1.83	0.60
1:B:758:ILE:HG12	1:B:770:TRP:HE3	1.67	0.60
1:A:871:GLU:H	1:A:871:GLU:CD	2.09	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:758:ILE:HG12	1:A:770:TRP:HE3	1.66	0.59
1:B:725:ALA:HA	2:B:904:HOH:O	2.03	0.58
1:B:562:GLU:OE1	1:B:685:LYS:NZ	2.35	0.58
1:B:837:ARG:HG2	1:B:840:ARG:HH22	1.68	0.58
1:A:697:GLN:HG3	1:A:727:TYR:CE2	2.40	0.57
1:A:625:PHE:CZ	1:A:639:HIS:HB3	2.39	0.56
1:A:513:ARG:HG3	1:A:515:SER:H	1.68	0.56
1:A:758:ILE:HG12	1:A:770:TRP:CE3	2.40	0.56
1:B:548:GLU:HG2	1:B:822:LEU:HD12	1.87	0.55
1:B:528:LEU:HD12	1:B:532:LEU:HD13	1.88	0.55
1:A:627:GLU:OE2	1:A:853:ARG:HB2	2.06	0.55
1:B:590:ILE:HD13	1:B:735:ILE:HD12	1.87	0.55
1:A:673:THR:HG23	1:A:680:HIS:HD2	1.70	0.55
1:A:617:ARG:HH11	1:A:619:ASP:CB	2.19	0.55
1:A:680:HIS:ND1	1:A:711:LEU:HD12	2.21	0.55
1:A:843:LEU:HB2	1:A:848:VAL:HB	1.88	0.55
1:B:616:ILE:HG23	1:B:623:GLN:HG2	1.89	0.55
1:A:528:LEU:HG	1:A:549:LEU:HD13	1.88	0.54
1:A:512:ARG:NH2	1:A:617:ARG:HH12	2.06	0.54
1:A:513:ARG:NE	1:A:515:SER:HB2	2.19	0.54
1:A:561:GLU:OE2	1:A:599:ARG:NH1	2.40	0.54
1:B:583:VAL:HG13	1:B:750:LEU:HD21	1.89	0.54
1:B:579:LEU:HD22	1:B:747:VAL:HG21	1.90	0.53
1:A:579:LEU:HD13	1:A:747:VAL:HG11	1.91	0.52
1:A:805:PRO:HG3	1:A:882:ASP:OD2	2.09	0.52
1:A:579:LEU:CD2	1:A:716:CYS:HB3	2.40	0.51
1:A:812:PRO:HB2	1:A:828:GLN:HG3	1.92	0.51
1:B:735:ILE:HD11	1:B:758:ILE:CD1	2.36	0.51
1:B:693:ILE:O	1:B:694:LEU:HG	2.10	0.51
1:B:561:GLU:OE2	1:B:599:ARG:NH1	2.41	0.51
1:B:723:ILE:CG2	2:B:904:HOH:O	2.57	0.51
1:B:757:SER:O	1:B:758:ILE:HD13	2.12	0.50
1:A:573:ALA:HB2	1:A:674:LEU:HB3	1.93	0.50
1:A:612:THR:HG22	1:A:627:GLU:HG3	1.92	0.50
1:B:701:GLU:OE1	1:B:724:LYS:NZ	2.43	0.50
1:A:552:PRO:O	1:A:553:LEU:HD22	2.12	0.50
1:A:556:LEU:HD23	1:A:636:SER:HB3	1.93	0.50
1:A:799:SER:HA	1:A:802:LEU:HD12	1.92	0.50
1:A:883:HIS:HB3	1:A:884:PRO:HD2	1.92	0.49
1:B:531:ILE:HG22	1:B:532:LEU:HD12	1.95	0.49
1:B:689:CYS:HB2	1:B:703:TYR:OH	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:878:PHE:C	1:B:880:LYS:H	2.20	0.49
1:A:812:PRO:O	1:A:816:PHE:HD1	1.94	0.49
1:A:831:ARG:HH22	1:A:887:TRP:HA	1.77	0.49
1:B:590:ILE:HD13	1:B:735:ILE:CD1	2.42	0.49
1:B:876:LEU:O	2:B:908:HOH:O	2.19	0.49
1:A:653:TRP:HZ3	1:A:666:PRO:HG3	1.75	0.48
1:B:724:LYS:C	2:B:904:HOH:O	2.51	0.48
1:A:553:LEU:HD23	1:A:785:PHE:CZ	2.48	0.48
1:B:579:LEU:CD2	1:B:747:VAL:HG21	2.44	0.48
1:B:811:ARG:HG2	1:B:813:ASP:OD1	2.13	0.48
1:B:516:LEU:HB2	1:B:782:TYR:CD1	2.49	0.48
1:A:532:LEU:HB3	1:A:657:PHE:CD2	2.49	0.48
1:A:608:VAL:HG11	1:A:886:LEU:HD12	1.95	0.48
1:A:871:GLU:HA	1:A:874:LYS:HB2	1.96	0.47
1:B:528:LEU:HD22	1:B:698:ARG:NH1	2.29	0.47
1:B:837:ARG:HG2	1:B:840:ARG:NH2	2.29	0.47
1:A:608:VAL:HG23	1:A:811:ARG:NH1	2.29	0.47
1:A:673:THR:HA	1:A:679:ASP:O	2.14	0.47
1:B:609:LEU:HD11	1:B:850:HIS:CD2	2.50	0.47
1:B:705:GLU:N	2:B:907:HOH:O	2.17	0.47
1:B:826:GLU:O	1:B:830:GLN:HG3	2.14	0.47
1:A:559:LEU:HD11	1:A:651:VAL:HG22	1.96	0.47
1:A:673:THR:HG22	1:A:680:HIS:HA	1.96	0.47
1:A:878:PHE:CE2	1:A:880:LYS:HD2	2.50	0.47
1:B:566:SER:OG	1:B:620:LYS:HE3	2.15	0.47
1:A:697:GLN:HG2	1:A:697:GLN:O	2.15	0.47
1:B:839:ARG:NH2	1:B:885:VAL:H	2.14	0.46
1:A:868:THR:O	1:A:872:LEU:HB2	2.15	0.46
1:A:654:LYS:HB3	1:A:665:VAL:HB	1.96	0.46
1:A:581:ARG:NH2	1:A:679:ASP:CG	2.72	0.46
1:B:726:LYS:HE2	1:B:728:TRP:HE1	1.80	0.46
1:A:697:GLN:HG3	1:A:727:TYR:CZ	2.50	0.46
1:B:528:LEU:HD21	1:B:694:LEU:H	1.79	0.46
1:A:520:SER:HB2	1:A:783:TYR:CD1	2.51	0.46
1:A:882:ASP:N	1:A:882:ASP:OD1	2.47	0.46
1:B:726:LYS:HB3	1:B:728:TRP:NE1	2.31	0.46
1:B:703:TYR:CB	1:B:724:LYS:HB2	2.46	0.45
1:A:709:LYS:HB2	1:A:709:LYS:HE3	1.55	0.45
1:A:556:LEU:CD2	1:A:636:SER:HB3	2.46	0.45
1:A:647:PHE:HD1	1:A:672:VAL:HG22	1.80	0.45
1:A:854:PHE:HA	1:A:866:ASN:ND2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:813:ASP:OD1	1:A:814:GLN:N	2.50	0.45
1:A:555:THR:HG23	1:A:607:PRO:HD3	1.99	0.45
1:A:608:VAL:HG13	1:A:836:GLN:HG2	1.99	0.45
1:A:617:ARG:HD3	1:A:620:LYS:CG	2.46	0.44
1:A:645:PHE:HA	1:A:673:THR:O	2.17	0.44
1:A:675:PRO:C	1:A:677:PHE:H	2.24	0.44
1:A:760:CYS:SG	1:A:761:GLY:N	2.91	0.44
1:B:610:GLY:O	1:B:853:ARG:NH1	2.50	0.44
1:A:675:PRO:C	1:A:677:PHE:N	2.76	0.44
1:A:516:LEU:O	1:A:518:ALA:N	2.51	0.44
1:A:701:GLU:CD	1:A:724:LYS:HG3	2.43	0.44
1:A:623:GLN:O	1:A:640:ALA:HA	2.18	0.44
1:A:573:ALA:HB2	1:A:674:LEU:HD22	2.01	0.43
1:A:563:LEU:HD21	1:A:647:PHE:CD2	2.53	0.43
1:A:574:GLN:NE2	1:A:644:ASN:HD22	2.16	0.43
1:A:693:ILE:HA	1:A:698:ARG:HG2	2.00	0.43
1:A:561:GLU:CD	1:A:599:ARG:HH22	2.27	0.42
1:A:590:ILE:HD13	1:A:735:ILE:HD13	2.01	0.42
1:A:742:ARG:HG3	1:A:743:SER:N	2.34	0.42
1:B:878:PHE:O	1:B:880:LYS:N	2.49	0.42
1:A:638:CYS:SG	1:A:649:GLN:HG2	2.60	0.42
1:B:758:ILE:HG12	1:B:770:TRP:CE3	2.50	0.42
1:A:572:ALA:C	1:A:574:GLN:H	2.28	0.42
1:A:839:ARG:O	1:A:843:LEU:HG	2.20	0.41
1:A:556:LEU:HB3	1:A:626:SER:OG	2.20	0.41
1:A:673:THR:CG2	1:A:680:HIS:CD2	3.03	0.41
1:A:515:SER:HA	1:A:782:TYR:HE2	1.84	0.41
1:A:680:HIS:O	1:A:710:ASN:HA	2.21	0.41
1:A:564:GLU:HA	1:A:788:PHE:CE2	2.56	0.41
1:A:662:MET:HE3	1:A:664:ILE:HD11	2.02	0.41
1:A:750:LEU:HA	1:A:759:TYR:O	2.21	0.41
1:B:558:ARG:CG	2:B:902:HOH:O	2.35	0.41
1:B:679:ASP:HB3	1:B:681:PHE:CE2	2.56	0.41
1:A:649:GLN:OE1	1:A:685:LYS:HE3	2.21	0.41
1:A:813:ASP:HB3	1:A:828:GLN:HB2	2.02	0.41
1:B:581:ARG:HD3	1:B:677:PHE:CE2	2.56	0.41
1:A:673:THR:HG23	1:A:680:HIS:CD2	2.53	0.40
1:B:725:ALA:CA	2:B:904:HOH:O	2.65	0.40
1:A:675:PRO:O	1:A:677:PHE:N	2.54	0.40
1:A:732:ALA:O	1:A:753:LYS:HE3	2.21	0.40
1:A:516:LEU:HG	1:A:517:PRO:HD2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:550:ASN:HA	1:B:603:LYS:O	2.21	0.40
1:A:689:CYS:HB2	1:A:703:TYR:CE1	2.57	0.40
1:B:787:GLN:O	1:B:791:GLU:HG3	2.22	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	373/388 (96%)	342 (92%)	31 (8%)	0	100	100
1	B	374/388 (96%)	351 (94%)	22 (6%)	1 (0%)	36	60
All	All	747/776 (96%)	693 (93%)	53 (7%)	1 (0%)	48	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	859	ASP

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	336/346 (97%)	336 (100%)	0	100	100
1	B	337/346 (97%)	336 (100%)	1 (0%)	86	94

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	673/692 (97%)	672 (100%)	1 (0%)	88	96

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	686	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	535	ASN
1	A	550	ASN
1	A	644	ASN
1	A	671	HIS
1	A	680	HIS
1	A	731	ASN
1	A	733	HIS
1	B	631	HIS
1	B	828	GLN
1	B	850	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	375/388 (96%)	1.22	79 (21%) 2 2	40, 88, 131, 142	0
1	B	376/388 (96%)	0.07	22 (5%) 28 25	27, 48, 106, 131	0
All	All	751/776 (96%)	0.65	101 (13%) 7 6	27, 64, 124, 142	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	541	SER	6.0
1	A	806	THR	5.5
1	B	725	ALA	5.4
1	B	694	LEU	4.6
1	A	799	SER	4.5
1	A	728	TRP	4.4
1	A	887	TRP	4.3
1	A	873	ARG	4.1
1	B	526	ILE	3.9
1	A	886	LEU	3.9
1	B	531	ILE	3.8
1	A	624	PHE	3.7
1	A	842	VAL	3.7
1	A	616	ILE	3.6
1	B	529	TRP	3.5
1	A	613	TYR	3.4
1	A	625	PHE	3.4
1	A	802	LEU	3.4
1	A	876	LEU	3.4
1	A	573	ALA	3.3
1	A	833	GLU	3.3
1	A	578	PRO	3.3
1	A	727	TYR	3.3
1	A	615	CYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	790	LEU	3.2
1	A	878	PHE	3.2
1	B	615	CYS	3.2
1	A	820	GLY	3.2
1	A	522	SER	3.2
1	A	798	SER	3.2
1	A	881	LEU	3.2
1	A	854	PHE	3.1
1	A	805	PRO	3.0
1	A	852	PRO	3.0
1	A	825	ALA	3.0
1	B	751	PHE	3.0
1	B	716	CYS	2.9
1	A	673	THR	2.9
1	A	867	GLY	2.9
1	A	699	TRP	2.8
1	A	671	HIS	2.8
1	B	527	SER	2.8
1	A	803	LEU	2.8
1	A	760	CYS	2.8
1	B	696	GLY	2.8
1	B	760	CYS	2.7
1	B	726	LYS	2.7
1	A	602	SER	2.7
1	A	843	LEU	2.7
1	A	813	ASP	2.7
1	A	817	LEU	2.6
1	A	761	GLY	2.6
1	A	646	VAL	2.6
1	A	623	GLN	2.6
1	B	695	SER	2.5
1	A	680	HIS	2.5
1	B	659	GLY	2.5
1	A	880	LYS	2.5
1	A	804	PRO	2.5
1	A	574	GLN	2.5
1	B	532	LEU	2.5
1	A	640	ALA	2.5
1	A	689	CYS	2.5
1	A	870	LEU	2.5
1	A	696	GLY	2.4
1	A	605	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	863	TRP	2.4
1	A	695	SER	2.4
1	A	558	ARG	2.3
1	A	785	PHE	2.3
1	A	626	SER	2.3
1	A	762	GLY	2.3
1	A	629	VAL	2.3
1	A	795	MET	2.3
1	A	634	PRO	2.3
1	A	650	ASP	2.3
1	A	877	GLY	2.3
1	B	537	GLY	2.3
1	A	865	SER	2.2
1	B	657	PHE	2.2
1	A	577	SER	2.2
1	B	799	SER	2.2
1	A	810	PHE	2.2
1	A	797	PRO	2.2
1	A	610	GLY	2.1
1	A	524	SER	2.1
1	A	521	PRO	2.1
1	B	530	ASN	2.1
1	A	869	TYR	2.1
1	B	543	VAL	2.1
1	A	855	PHE	2.1
1	A	853	ARG	2.0
1	A	711	LEU	2.0
1	A	834	GLN	2.0
1	A	832	ILE	2.0
1	A	676	VAL	2.0
1	A	681	PHE	2.0
1	B	661	SER	2.0
1	A	517	PRO	2.0
1	A	556	LEU	2.0
1	A	763	GLY	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.