



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:56 PM UTC

PDB ID : 7YW1 / pdb_00007yw1
Title : crystal structure of UBE2O
Authors : Fu, Z.; Zhu, W.; Huang, H.
Deposited on : 2022-08-20
Resolution : 3.27 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

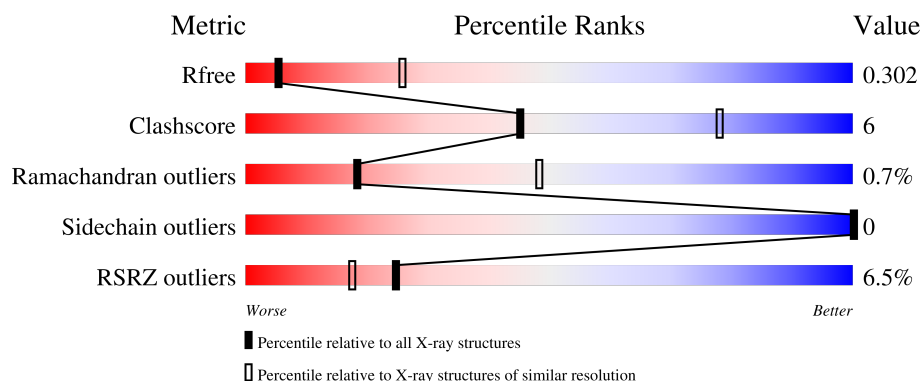
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.27 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	1303 (3.30-3.26)
Clashscore	190562	1354 (3.30-3.26)
Ramachandran outliers	187476	1334 (3.30-3.26)
Sidechain outliers	187428	1333 (3.30-3.26)
RSRZ outliers	180081	1303 (3.30-3.26)

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	928	5% 66% 12% 22%
1	B	928	5% 60% 9% 30%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 10032 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.

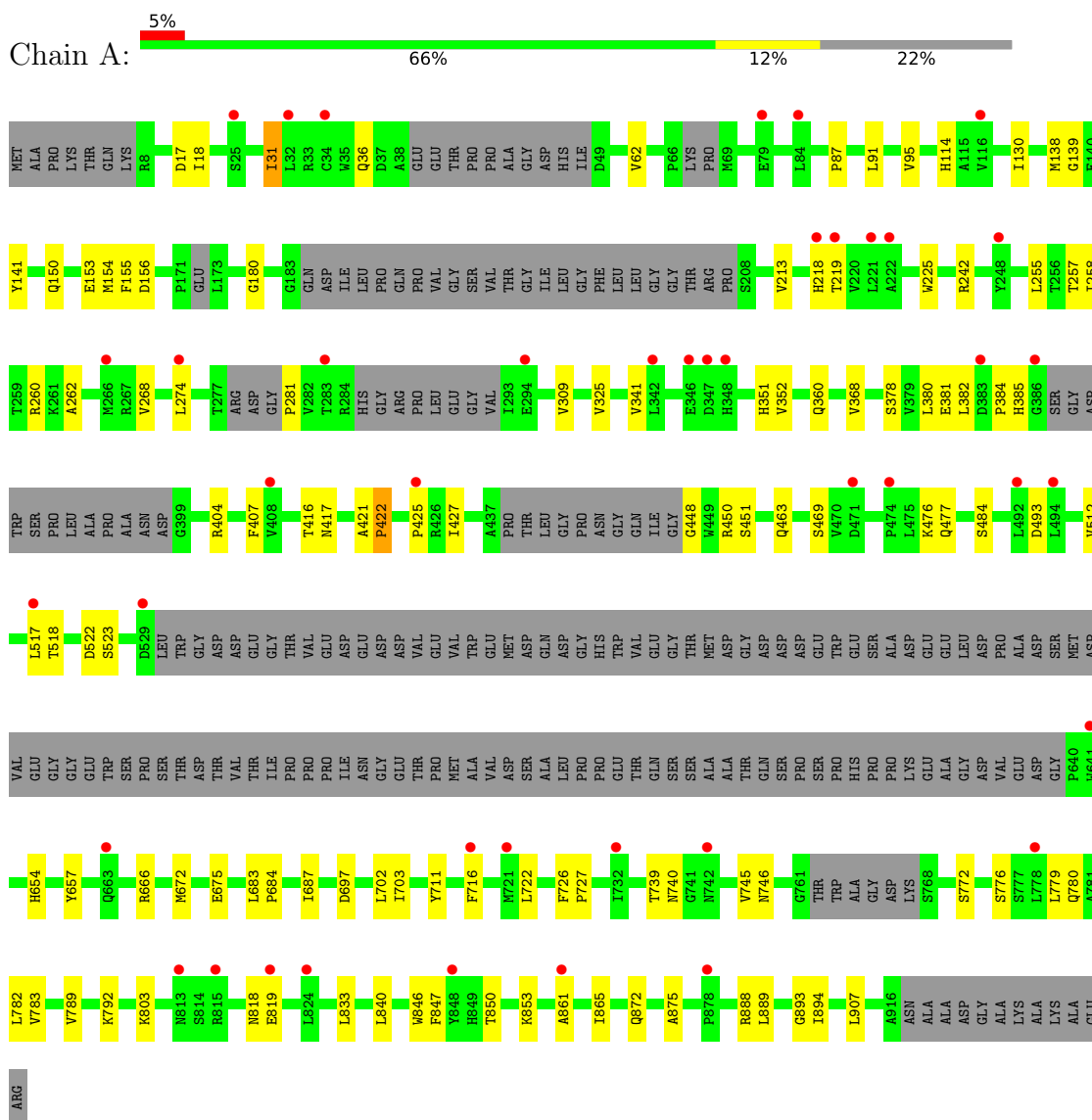
- Molecule 1 is a protein called (E3-independent) E2 ubiquitin-conjugating enzyme UBE2O.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	723	Total	C	N	O	S	0	0	0
			5350	3390	932	1014	14			
1	B	648	Total	C	N	O	S	0	0	0
			4682	2969	806	897	10			

3 Residue-property plots

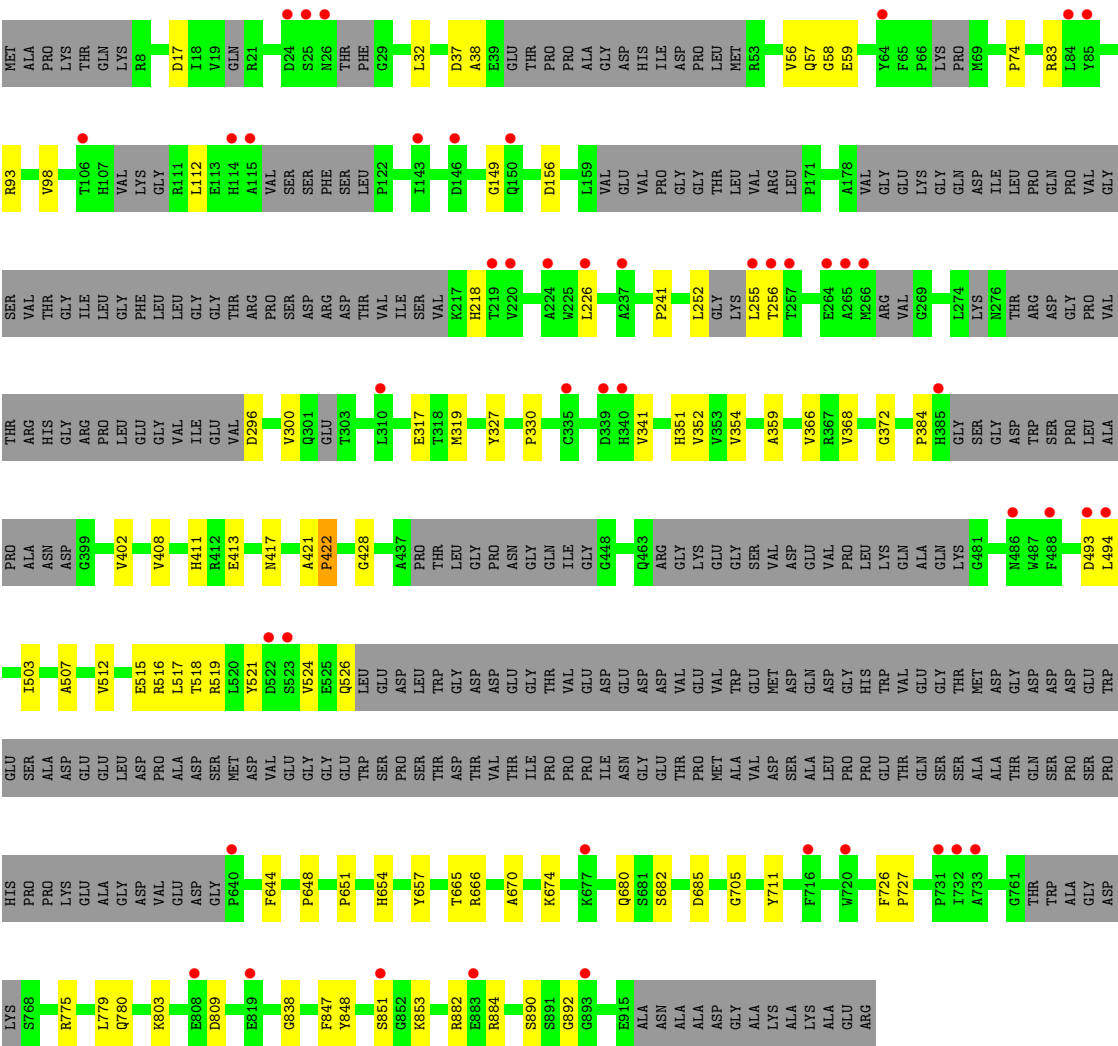
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: (E3-independent) E2 ubiquitin-conjugating enzyme UBE2O



- Molecule 1: (E3-independent) E2 ubiquitin-conjugating enzyme UBE2O





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	115.32Å 132.53Å 155.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.04 – 3.27 54.04 – 3.27	Depositor EDS
% Data completeness (in resolution range)	76.6 (54.04-3.27) 76.6 (54.04-3.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 3.26Å)	Xtriage
Refinement program	PHENIX 1.17.1-3660	Depositor
R, R_{free}	0.249 , 0.300 (Not available) , 0.302	Depositor DCC
R_{free} test set	1351 reflections (3.62%)	wwPDB-VP
Wilson B-factor (Å ²)	87.2	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 96.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	10032	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.09	0/5459	0.27	0/7457
1	B	0.09	0/4770	0.26	0/6524
All	All	0.09	0/10229	0.27	0/13981

There are no bond length outliers.

There are no bond angle outliers.

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	5350	0	4977	63	0
1	B	4682	0	4224	52	0
All	All	10032	0	9201	113	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 6.

All (113) 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:378:SER:HB3	1:A:381:GLU:HG3	1.83	0.61
1:A:384:PRO:O	1:A:385:HIS:ND1	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:644:PHE:HB3	1:B:680:GLN:HG2	1.83	0.60
1:A:114:HIS:HD2	1:A:281:PRO:HB2	1.67	0.59
1:B:32:LEU:H	1:B:32:LEU:HD23	1.68	0.59
1:A:404:ARG:NH2	1:A:493:ASP:OD1	2.36	0.58
1:A:666:ARG:HH21	1:B:526:GLN:HA	1.68	0.58
1:A:36:GLN:HA	1:A:803:LYS:HD3	1.84	0.58
1:B:682:SER:HB2	1:B:779:LEU:HD21	1.86	0.58
1:A:31:ILE:H	1:A:62:VAL:HG23	1.70	0.57
1:A:687:ILE:HG23	1:A:703:ILE:HG12	1.87	0.57
1:A:255:LEU:HB2	1:A:476:LYS:HG2	1.86	0.56
1:A:865:ILE:HD12	1:A:894:ILE:HG23	1.88	0.56
1:A:463:GLN:HG3	1:A:469:SER:HB3	1.88	0.55
1:B:408:VAL:HG12	1:B:519:ARG:HA	1.87	0.55
1:A:833:LEU:HB2	1:A:907:LEU:HD13	1.88	0.55
1:A:257:THR:HG23	1:A:476:LYS:HB3	1.88	0.54
1:A:351:HIS:ND1	1:A:422:PRO:HG3	2.23	0.54
1:A:352:VAL:HG12	1:A:368:VAL:HA	1.90	0.54
1:A:780:GLN:HA	1:A:783:VAL:HG22	1.90	0.53
1:B:17:ASP:HB2	1:B:83:ARG:HA	1.91	0.53
1:A:425:PRO:HB2	1:A:518:THR:HG23	1.89	0.53
1:B:56:VAL:HG13	1:B:58:GLY:H	1.74	0.53
1:B:411:HIS:NE2	1:B:518:THR:OG1	2.37	0.52
1:A:153:GLU:OE1	1:A:450:ARG:NH1	2.42	0.52
1:A:87:PRO:HD2	1:A:262:ALA:HA	1.92	0.51
1:A:792:LYS:O	1:A:818:ASN:ND2	2.43	0.51
1:A:448:GLY:O	1:A:451:SER:OG	2.28	0.51
1:B:59:GLU:HA	1:B:74:PRO:HA	1.93	0.51
1:B:156:ASP:HA	1:B:218:HIS:HA	1.92	0.50
1:B:352:VAL:HB	1:B:366:VAL:HB	1.92	0.50
1:B:341:VAL:HG12	1:B:384:PRO:HB3	1.93	0.50
1:B:354:VAL:HA	1:B:366:VAL:HG12	1.93	0.50
1:B:413:GLU:OE2	1:B:516:ARG:NH1	2.45	0.49
1:A:18:ILE:HG12	1:A:268:VAL:HG21	1.94	0.49
1:A:225:TRP:O	1:A:242:ARG:NH2	2.39	0.49
1:A:341:VAL:HG21	1:A:382:LEU:HD13	1.95	0.49
1:B:654:HIS:HB3	1:B:657:TYR:HB2	1.94	0.49
1:A:421:ALA:HB1	1:A:422:PRO:HD2	1.96	0.48
1:B:417:ASN:N	1:B:515:GLU:O	2.46	0.48
1:B:351:HIS:CE1	1:B:422:PRO:HG3	2.49	0.48
1:A:155:PHE:O	1:A:219:THR:OG1	2.31	0.48
1:B:93:ARG:NH1	1:B:317:GLU:OE2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:MET:HE3	1:A:154:MET:HE2	1.96	0.47
1:A:872:GLN:HA	1:A:888:ARG:HH21	1.79	0.47
1:A:746:ASN:ND2	1:A:789:VAL:O	2.47	0.47
1:A:17:ASP:N	1:A:17:ASP:OD1	2.45	0.47
1:A:697:ASP:HA	1:A:722:LEU:HD12	1.97	0.47
1:A:141:TYR:HB2	1:A:258:ILE:HB	1.96	0.47
1:A:260:ARG:HG2	1:A:407:PHE:CE2	2.49	0.47
1:A:683:LEU:HD23	1:A:684:PRO:HD2	1.95	0.47
1:B:890:SER:C	1:B:892:GLY:H	2.23	0.47
1:A:522:ASP:OD1	1:A:523:SER:N	2.48	0.47
1:A:779:LEU:HA	1:A:782:LEU:HD12	1.97	0.47
1:B:37:ASP:HA	1:B:803:LYS:HD2	1.96	0.46
1:A:156:ASP:OD1	1:A:156:ASP:N	2.44	0.46
1:A:380:LEU:HD12	1:A:893:GLY:HA2	1.96	0.46
1:B:512:VAL:HG21	1:B:517:LEU:HD21	1.97	0.46
1:B:428:GLY:HA3	1:B:521:TYR:HB3	1.97	0.46
1:B:775:ARG:NH1	1:B:780:GLN:OE1	2.48	0.46
1:B:56:VAL:HG22	1:B:57:GLN:H	1.80	0.46
1:A:130:ILE:HG22	1:A:274:LEU:HD23	1.97	0.46
1:B:685:ASP:N	1:B:685:ASP:OD1	2.48	0.46
1:A:846:TRP:HA	1:A:850:THR:HB	1.98	0.45
1:B:352:VAL:HG12	1:B:368:VAL:HA	1.97	0.45
1:B:705:GLY:HA3	1:B:711:TYR:O	2.17	0.45
1:B:411:HIS:O	1:B:516:ARG:NH2	2.42	0.45
1:B:368:VAL:O	1:B:372:GLY:N	2.44	0.45
1:B:149:GLY:HA2	1:B:226:LEU:HG	1.98	0.45
1:B:112:LEU:HB2	1:B:300:VAL:HA	1.99	0.45
1:B:56:VAL:H	1:B:59:GLU:HG2	1.82	0.44
1:A:360:GLN:HE22	1:A:875:ALA:HA	1.82	0.44
1:A:260:ARG:HG2	1:A:407:PHE:HE2	1.82	0.44
1:B:93:ARG:HG2	1:B:319:MET:HE1	2.00	0.44
1:B:847:PHE:O	1:B:851:SER:OG	2.35	0.44
1:A:711:TYR:HE1	1:A:745:VAL:HG22	1.82	0.44
1:B:651:PRO:HB3	1:B:838:GLY:HA3	1.99	0.44
1:B:503:ILE:HD11	1:B:507:ALA:HA	2.00	0.44
1:A:180:GLY:H	1:A:213:VAL:HB	1.83	0.43
1:A:666:ARG:NH2	1:B:526:GLN:HA	2.33	0.43
1:B:882:ARG:C	1:B:884:ARG:H	2.27	0.43
1:A:512:VAL:HG21	1:A:517:LEU:HD21	2.01	0.43
1:B:421:ALA:O	1:B:422:PRO:C	2.61	0.43
1:A:416:THR:OG1	1:A:417:ASN:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:861:ALA:O	1:A:865:ILE:HG12	2.19	0.43
1:B:726:PHE:CG	1:B:727:PRO:HA	2.54	0.43
1:B:848:TYR:HA	1:B:853:LYS:HB2	2.00	0.43
1:A:726:PHE:CG	1:A:727:PRO:HA	2.54	0.42
1:A:711:TYR:HB3	1:A:716:PHE:CZ	2.53	0.42
1:A:772:SER:H	1:A:776:SER:HB2	1.85	0.42
1:A:819:GLU:HG3	1:A:889:LEU:HD23	2.01	0.42
1:B:38:ALA:H	1:B:803:LYS:HD3	1.84	0.42
1:A:421:ALA:O	1:A:422:PRO:C	2.62	0.42
1:A:427:ILE:HD13	1:A:427:ILE:HA	1.95	0.42
1:B:327:TYR:CZ	1:B:330:PRO:HA	2.55	0.42
1:A:91:LEU:HD21	1:A:309:VAL:HG13	2.02	0.41
1:B:515:GLU:N	1:B:515:GLU:OE1	2.53	0.41
1:B:252:LEU:HD23	1:B:252:LEU:H	1.85	0.41
1:B:809:ASP:OD1	1:B:809:ASP:N	2.39	0.41
1:B:493:ASP:OD1	1:B:494:LEU:N	2.53	0.41
1:A:95:VAL:HG13	1:A:325:VAL:HG11	2.03	0.41
1:A:139:GLY:O	1:A:150:GLN:NE2	2.54	0.41
1:A:672:MET:HA	1:A:675:GLU:HB2	2.02	0.41
1:A:702:LEU:HD22	1:A:840:LEU:HD13	2.03	0.41
1:B:665:THR:OG1	1:B:666:ARG:N	2.50	0.41
1:B:670:ALA:O	1:B:674:LYS:HG2	2.20	0.41
1:A:847:PHE:O	1:A:853:LYS:HB2	2.21	0.41
1:B:98:VAL:HG12	1:B:359:ALA:HB1	2.03	0.40
1:B:255:LEU:HB2	1:B:256:THR:H	1.67	0.40
1:A:654:HIS:HB3	1:A:657:TYR:HB2	2.03	0.40
1:B:296:ASP:OD1	1:B:296:ASP:N	2.54	0.40
1:A:156:ASP:HB3	1:A:218:HIS:HA	2.03	0.40
1:A:739:THR:HG22	1:A:740:ASN:H	1.86	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	701/928 (76%)	635 (91%)	62 (9%)	4 (1%)	21	51
1	B	611/928 (66%)	554 (91%)	52 (8%)	5 (1%)	16	44
All	All	1312/1856 (71%)	1189 (91%)	114 (9%)	9 (1%)	18	48

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	31	ILE
1	A	422	PRO
1	B	422	PRO
1	B	524	VAL
1	B	402	VAL
1	A	484	SER
1	A	477	GLN
1	B	648	PRO
1	B	241	PRO

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	321/786 (41%)	321 (100%)	0	100	100
1	B	438/786 (56%)	438 (100%)	0	100	100
All	All	759/1572 (48%)	759 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	360	GLN
1	A	734	HIS
1	A	740	ASN
1	A	786	GLN
1	B	301	GLN

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Mol	Chain	Res	Type
1	B	323	GLN
1	B	329	ASN
1	B	340	HIS

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](#)

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	723/928 (77%)	0.55	43 (5%)	28 19	39, 80, 118, 143	0
1	B	648/928 (69%)	0.64	46 (7%)	22 16	48, 91, 142, 174	0
All	All	1371/1856 (73%)	0.59	89 (6%)	25 17	39, 84, 132, 174	0

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	335	CYS	5.0
1	A	221	LEU	4.6
1	B	494	LEU	4.6
1	B	808	GLU	4.3
1	B	143	ILE	4.1
1	B	256	THR	4.0
1	A	529	ASP	4.0
1	B	677	LYS	3.9
1	A	348	HIS	3.8
1	A	848	TYR	3.8
1	B	819	GLU	3.7
1	B	219	THR	3.5
1	B	493	ASP	3.5
1	B	237	ALA	3.4
1	A	383	ASP	3.3
1	A	248	TYR	3.3
1	A	342	LEU	3.2
1	B	716	PHE	3.2
1	B	720	TRP	3.1
1	B	339	ASP	3.1
1	A	222	ALA	3.1
1	B	488	PHE	3.0
1	A	494	LEU	3.0
1	A	32	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	224	ALA	2.9
1	A	732	ILE	2.9
1	B	150	GLN	2.9
1	A	79	GLU	2.9
1	A	425	PRO	2.9
1	B	115	ALA	2.9
1	B	731	PRO	2.8
1	A	815	ARG	2.8
1	B	25	SER	2.8
1	A	742	ASN	2.7
1	A	517	LEU	2.7
1	A	641	TRP	2.7
1	A	721	MET	2.6
1	A	283	THR	2.6
1	A	778	LEU	2.6
1	A	408	VAL	2.6
1	A	386	GLY	2.6
1	A	346	GLU	2.5
1	A	474	PRO	2.5
1	A	716	PHE	2.5
1	B	893	GLY	2.5
1	A	347	ASP	2.5
1	A	492	LEU	2.5
1	B	523	SER	2.5
1	A	813	ASN	2.5
1	A	266	MET	2.4
1	A	218	HIS	2.4
1	B	265	ALA	2.4
1	B	385	HIS	2.4
1	A	471	ASP	2.4
1	B	264	GLU	2.4
1	B	220	VAL	2.4
1	B	851	SER	2.3
1	B	733	ALA	2.3
1	A	663	GLN	2.3
1	B	640	PRO	2.3
1	A	34	CYS	2.3
1	B	64	TYR	2.3
1	A	84	LEU	2.3
1	B	883	GLU	2.3
1	B	114	HIS	2.3
1	B	486	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	732	ILE	2.2
1	A	861	ALA	2.2
1	A	878	PRO	2.2
1	B	226	LEU	2.2
1	B	266	MET	2.2
1	A	219	THR	2.2
1	B	257	THR	2.2
1	B	26	ASN	2.2
1	A	819	GLU	2.2
1	B	24	ASP	2.1
1	B	85	TYR	2.1
1	A	294	GLU	2.1
1	A	116	VAL	2.1
1	B	106	THR	2.1
1	B	255	LEU	2.1
1	B	340	HIS	2.1
1	B	522	ASP	2.1
1	B	84	LEU	2.0
1	A	824	LEU	2.0
1	B	310	LEU	2.0
1	A	25	SER	2.0
1	B	146	ASP	2.0
1	A	274	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.