



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:50 PM UTC

PDB ID : 5XWM / pdb_00005xwm
Title : human ERp44 zinc-bound form
Authors : Watanabe, S.; Harayama, M.; Inaba, K.
Deposited on : 2017-06-30
Resolution : 2.45 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

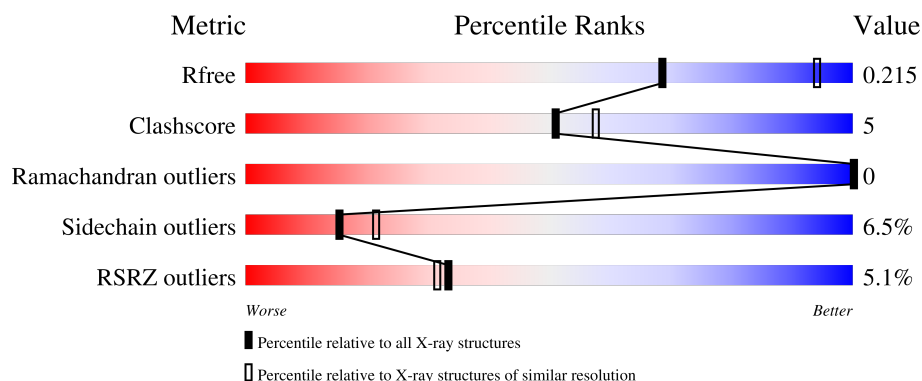
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	383	4% 73% 16% • 10%
1	B	383	4% 78% 13% • 8%
1	C	383	3% 78% 11% • 9%
1	D	383	7% 81% 14% • •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11949 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 Endoplasmic reticulum resident protein 44.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	0	1	0
			2757	1749	469	524	15			
1	B	354	Total	C	N	O	S	0	0	0
			2875	1825	491	545	14			
1	C	348	Total	C	N	O	S	0	0	0
			2823	1795	480	534	14			
1	D	368	Total	C	N	O	S	0	0	0
			2949	1869	505	561	14			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP Q9BS26
A	-4	SER	-	expression tag	UNP Q9BS26
A	-3	HIS	-	expression tag	UNP Q9BS26
A	-2	MET	-	expression tag	UNP Q9BS26
A	-1	ALA	-	expression tag	UNP Q9BS26
A	0	SER	-	expression tag	UNP Q9BS26
B	-5	GLY	-	expression tag	UNP Q9BS26
B	-4	SER	-	expression tag	UNP Q9BS26
B	-3	HIS	-	expression tag	UNP Q9BS26
B	-2	MET	-	expression tag	UNP Q9BS26
B	-1	ALA	-	expression tag	UNP Q9BS26
B	0	SER	-	expression tag	UNP Q9BS26
C	-5	GLY	-	expression tag	UNP Q9BS26
C	-4	SER	-	expression tag	UNP Q9BS26
C	-3	HIS	-	expression tag	UNP Q9BS26
C	-2	MET	-	expression tag	UNP Q9BS26
C	-1	ALA	-	expression tag	UNP Q9BS26
C	0	SER	-	expression tag	UNP Q9BS26
D	-5	GLY	-	expression tag	UNP Q9BS26
D	-4	SER	-	expression tag	UNP Q9BS26
D	-3	HIS	-	expression tag	UNP Q9BS26

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	MET	-	expression tag	UNP Q9BS26
D	-1	ALA	-	expression tag	UNP Q9BS26
D	0	SER	-	expression tag	UNP Q9BS26

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Zn	0	0
			3	3		
2	B	3	Total	Zn	0	0
			3	3		
2	C	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Cl	0	0
			2	2		
3	B	2	Total	Cl	0	0
			2	2		
3	C	2	Total	Cl	0	0
			2	2		
3	D	2	Total	Cl	0	0
			2	2		

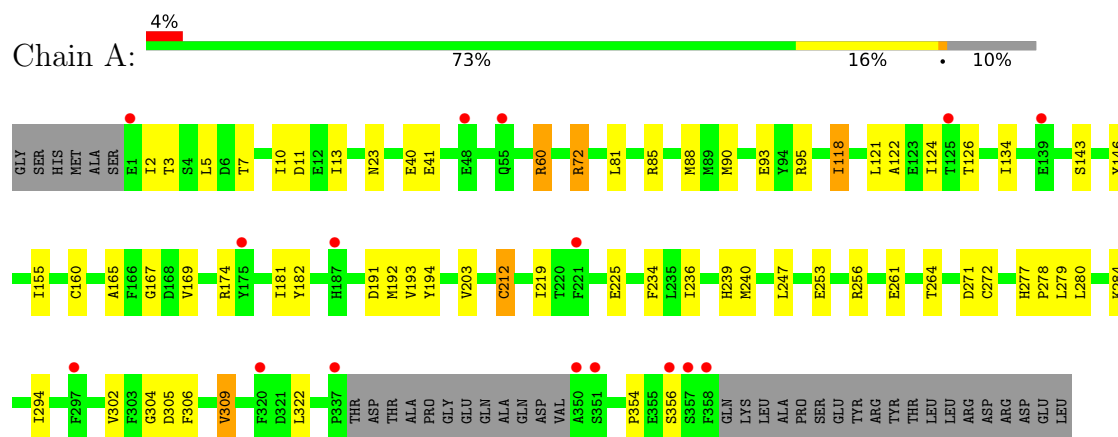
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	42	Total	O	0	0
			42	42		
4	B	152	Total	O	0	0
			152	152		
4	C	155	Total	O	0	0
			155	155		
4	D	178	Total	O	0	0
			178	178		

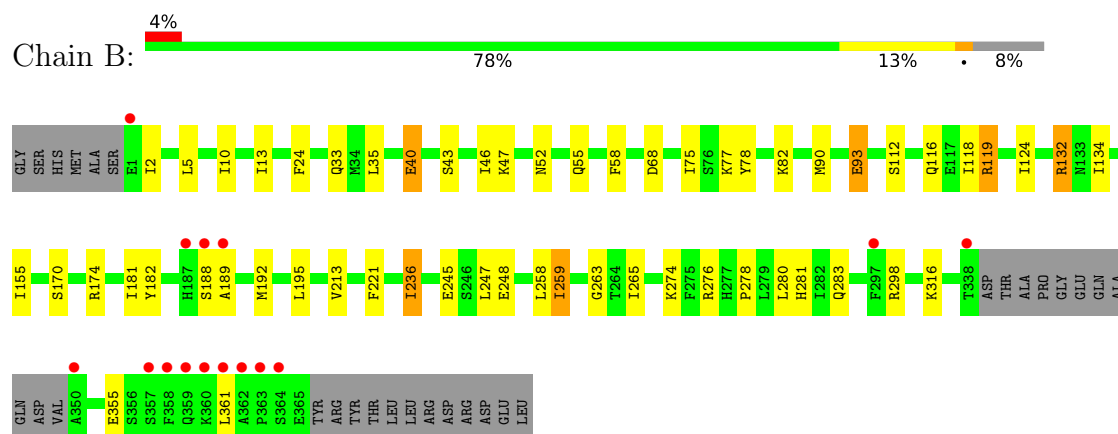
3 Residue-property plots [i](#)

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.

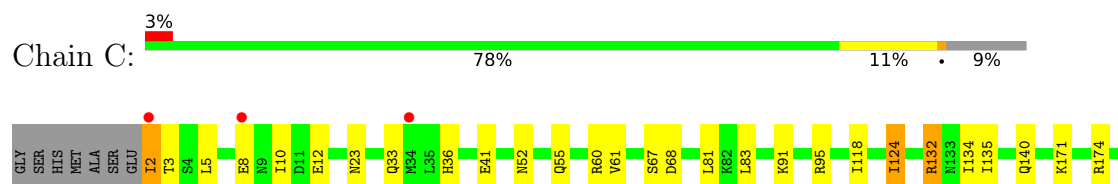
- Molecule 1: Endoplasmic reticulum resident protein 44

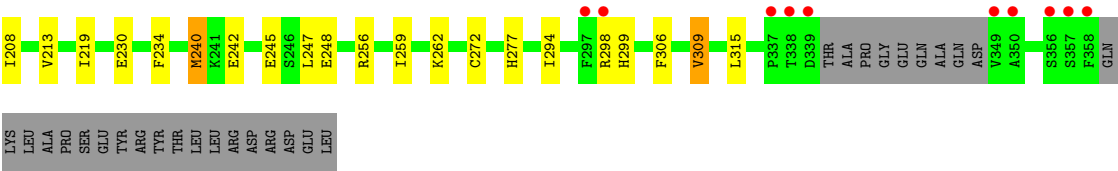


- Molecule 1: Endoplasmic reticulum resident protein 44

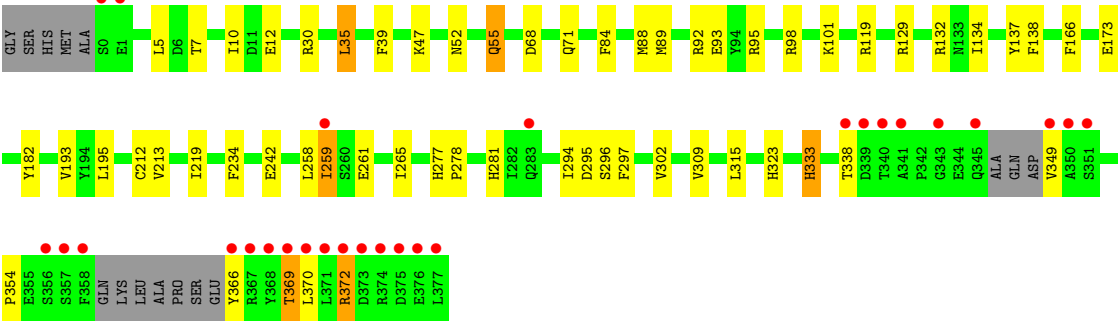
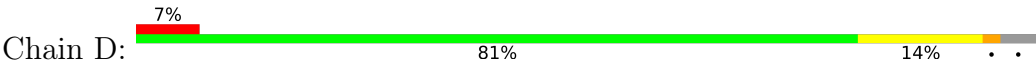


- Molecule 1: Endoplasmic reticulum resident protein 44





● Molecule 1: Endoplasmic reticulum resident protein 44



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	175.15Å 175.15Å 407.81Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.06 – 2.45 44.06 – 2.45	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.06-2.45) 99.9 (44.06-2.45)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.59 (at 2.45Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.182 , 0.214 0.183 , 0.215	Depositor DCC
R_{free} test set	6621 reflections (4.61%)	wwPDB-VP
Wilson B-factor (Å ²)	41.9	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 64.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11949	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.90% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, CL

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	0.31	0/2824	0.48	0/3833
1	B	0.35	0/2945	0.55	0/3984
1	C	0.37	0/2892	0.56	0/3915
1	D	0.37	0/3018	0.62	2/4087 (0.0%)
All	All	0.35	0/11679	0.55	2/15819 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	212	CYS	CA-C-N	-6.30	117.85	123.33
1	D	212	CYS	C-N-CA	-6.30	117.85	123.33

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	2757	0	2587	36	0
1	B	2875	0	2761	30	0
1	C	2823	0	2703	26	0
1	D	2949	0	2798	31	0
2	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	3	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	2	0	0	1	0
3	B	2	0	0	2	0
3	C	2	0	0	1	0
3	D	2	0	0	2	0
4	A	42	0	0	0	0
4	B	152	0	0	2	0
4	C	155	0	0	0	0
4	D	178	0	0	4	0
All	All	11949	0	10849	116	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 5.

All (116) 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:52:ASN:HB2	1:B:55:GLN:HG3	1.69	0.76
1:A:253:GLU:HG2	1:A:309:VAL:HG12	1.69	0.75
1:A:118:ILE:HD11	1:A:165:ALA:HB2	1.72	0.71
1:A:306:PHE:O	1:A:309:VAL:HG23	1.92	0.69
1:A:90[A]:MET:HE1	1:A:155:ILE:HA	1.76	0.67
1:D:281:HIS:HD1	1:D:366:TYR:HA	1.59	0.66
1:C:2:ILE:HD11	1:C:60:ARG:HD2	1.78	0.66
1:B:278:PRO:HG3	1:D:372:ARG:HG2	1.77	0.65
1:C:240:MET:HE3	1:C:272:CYS:SG	2.37	0.65
1:D:369:THR:O	1:D:369:THR:OG1	2.15	0.64
1:C:33:GLN:HA	1:C:36:HIS:HD2	1.62	0.64
1:D:242:GLU:H	1:D:242:GLU:CD	2.07	0.62
1:D:261:GLU:HG2	4:D:529:HOH:O	1.99	0.62
1:A:122:ALA:O	1:A:126:THR:HG22	2.00	0.61
1:D:88:MET:SD	4:D:676:HOH:O	2.57	0.60
1:B:90:MET:HE1	1:B:155:ILE:HG23	1.84	0.59
1:A:294:ILE:HG21	1:A:322:LEU:HD22	1.84	0.58
1:B:2:ILE:HG13	1:B:40:GLU:HG2	1.86	0.58
1:B:78:TYR:O	1:D:333:HIS:HE1	1.87	0.58
1:A:121:LEU:O	1:A:124:ILE:HG12	2.03	0.57
3:B:405:CL:CL	3:D:404:CL:CL	2.96	0.57
1:C:306:PHE:O	1:C:309:VAL:HG22	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:ARG:NH2	1:C:248:GLU:OE1	2.36	0.56
1:C:118:ILE:HD13	1:C:124:ILE:HD11	1.88	0.56
1:A:181:ILE:HD11	1:A:191:ASP:OD2	2.05	0.55
1:A:23:ASN:HA	1:A:81:LEU:HD23	1.87	0.55
1:A:277:HIS:HE2	1:C:277:HIS:CD2	2.14	0.54
1:D:134:ILE:HD13	1:D:182:TYR:HA	1.90	0.54
1:A:88:MET:HE2	1:A:90[B]:MET:HE1	1.89	0.54
1:B:5:LEU:HD13	1:B:10:ILE:HG13	1.91	0.53
1:B:283:GLN:HG2	1:B:361:LEU:CB	2.38	0.53
1:B:258:LEU:HD21	1:B:316:LYS:HA	1.91	0.52
1:B:82:LYS:HD3	1:B:93:GLU:HB2	1.90	0.52
1:D:5:LEU:HD13	1:D:10:ILE:HG13	1.90	0.52
1:A:2:ILE:HG22	1:A:60:ARG:HG2	1.93	0.50
1:A:354:PRO:O	1:C:95:ARG:NH1	2.37	0.49
1:D:265:ILE:HD11	1:D:323:HIS:CE1	2.48	0.49
1:C:2:ILE:O	1:C:2:ILE:HG13	2.12	0.49
1:D:35:LEU:HD22	1:D:39:PHE:HB2	1.93	0.49
3:A:405:CL:CL	3:C:404:CL:CL	3.04	0.49
1:D:84:PHE:CE2	1:D:89:MET:HB2	2.48	0.49
1:B:118:ILE:HD13	1:B:124:ILE:HD11	1.95	0.49
1:D:234:PHE:HB2	1:D:295:ASP:HB3	1.95	0.49
1:B:77:LYS:HD2	1:D:354:PRO:HB2	1.95	0.48
1:A:160:CYS:HB2	1:A:212:CYS:SG	2.52	0.48
1:A:277:HIS:NE2	1:C:277:HIS:CD2	2.80	0.48
1:C:259:ILE:O	1:C:262:LYS:HB2	2.14	0.48
1:D:281:HIS:ND1	1:D:366:TYR:HA	2.27	0.48
1:B:259:ILE:HD12	1:B:259:ILE:HA	1.79	0.48
1:C:5:LEU:HB2	1:C:61:VAL:HG22	1.96	0.48
1:A:277:HIS:O	1:A:280:LEU:HB3	2.13	0.47
1:A:272:CYS:HB2	1:A:279:LEU:HD11	1.95	0.47
1:C:171:LYS:HA	1:C:174:ARG:NH1	2.30	0.47
1:B:248:GLU:OE1	1:C:132:ARG:NH2	2.48	0.47
3:B:404:CL:CL	3:D:403:CL:CL	3.08	0.46
1:A:302:VAL:HG12	1:A:304:GLY:H	1.81	0.46
1:B:281:HIS:HD1	1:D:370:LEU:H	1.62	0.46
1:C:134:ILE:HD13	1:C:208:ILE:HG23	1.98	0.46
1:C:5:LEU:HD13	1:C:10:ILE:HG13	1.98	0.46
1:C:298:ARG:HG2	1:C:299:HIS:CD2	2.51	0.46
1:A:239:HIS:HE1	1:A:247:LEU:HD22	1.81	0.46
1:B:170:SER:OG	1:B:174:ARG:HD3	2.16	0.46
1:D:129:ARG:HA	1:D:129:ARG:HD2	1.74	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:ARG:HE	1:A:90[A]:MET:HE2	1.80	0.45
1:A:121:LEU:HD12	1:A:121:LEU:H	1.80	0.45
1:A:11:ASP:OD1	1:A:72:ARG:NH2	2.49	0.45
1:A:284:LYS:HD3	1:A:284:LYS:HA	1.81	0.45
1:A:277:HIS:HD2	1:A:280:LEU:HD23	1.81	0.45
1:C:23:ASN:HA	1:C:81:LEU:HD23	1.97	0.45
1:B:188:SER:OG	1:B:189:ALA:N	2.50	0.45
1:D:281:HIS:HD1	1:D:366:TYR:CA	2.25	0.45
1:B:182:TYR:HD2	1:B:192:MET:HE3	1.81	0.45
1:B:119:ARG:HA	1:B:119:ARG:HD2	1.75	0.45
1:B:118:ILE:CD1	1:B:124:ILE:HD11	2.47	0.45
1:B:24:PHE:CD2	1:B:75:ILE:HG13	2.52	0.45
1:D:349:VAL:N	4:D:507:HOH:O	2.50	0.45
1:A:121:LEU:HD11	1:A:169:VAL:O	2.17	0.45
1:C:68:ASP:OD1	1:C:68:ASP:N	2.50	0.45
1:A:194:TYR:OH	1:A:203:VAL:HG12	2.17	0.44
1:C:52:ASN:HB2	1:C:55:GLN:CD	2.41	0.44
1:C:91:LYS:HB3	1:C:230:GLU:HG2	1.98	0.44
1:D:259:ILE:HD13	1:D:259:ILE:HA	1.76	0.44
1:A:3:THR:HG21	1:A:13:ILE:HD13	2.00	0.44
1:A:143:SER:HB2	1:A:146:TYR:H	1.82	0.43
1:A:277:HIS:HB3	1:A:278:PRO:HD3	2.00	0.43
1:D:68:ASP:OD1	1:D:68:ASP:N	2.51	0.43
1:B:47:LYS:HA	1:B:47:LYS:HD2	1.83	0.43
1:D:234:PHE:O	1:D:294:ILE:HA	2.19	0.43
1:A:167:GLY:O	1:A:174:ARG:NH2	2.50	0.43
1:A:240:MET:HE3	1:A:240:MET:HB3	1.88	0.42
1:A:261:GLU:OE1	1:A:264:THR:OG1	2.35	0.42
1:D:138:PHE:O	1:D:166:PHE:HA	2.19	0.42
1:C:315:LEU:HD12	1:C:315:LEU:HA	1.80	0.42
1:D:137:TYR:CE2	1:D:173:GLU:HG3	2.54	0.42
1:C:83:LEU:HD12	1:C:83:LEU:HA	1.87	0.42
1:D:333:HIS:CD2	1:D:333:HIS:H	2.37	0.42
1:D:258:LEU:HD11	1:D:315:LEU:HG	2.00	0.42
1:A:5:LEU:HD13	1:A:10:ILE:HD12	2.01	0.42
1:B:236:ILE:HD13	1:B:236:ILE:HG21	1.80	0.41
1:A:134:ILE:HG12	1:A:182:TYR:HD1	1.86	0.41
1:A:234:PHE:O	1:A:294:ILE:HA	2.20	0.41
1:B:43:SER:HB3	1:B:58:PHE:CD2	2.55	0.41
1:B:263:GLY:N	4:B:501:HOH:O	2.15	0.41
1:B:68:ASP:OD1	1:B:68:ASP:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:52:ASN:HB2	1:D:55:GLN:HG2	2.02	0.41
1:D:92:ARG:HD2	4:D:665:HOH:O	2.19	0.41
1:C:124:ILE:HG12	1:C:135:ILE:HD13	2.03	0.41
1:D:296:SER:O	1:D:297:PHE:HB2	2.20	0.41
1:B:90:MET:HE2	4:B:578:HOH:O	2.21	0.41
1:B:181:ILE:HD12	1:B:181:ILE:HG23	1.83	0.41
1:C:36:HIS:CE1	1:C:60:ARG:HH21	2.39	0.41
1:A:253:GLU:OE2	1:A:256:ARG:NH1	2.54	0.41
1:B:298:ARG:O	1:D:369:THR:HG21	2.20	0.41
1:B:221:PHE:HZ	1:B:274:LYS:O	2.04	0.40
1:C:234:PHE:O	1:C:294:ILE:HA	2.22	0.40
1:D:277:HIS:HB3	1:D:278:PRO:HD3	2.02	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	343/383 (90%)	332 (97%)	11 (3%)	0	100	100
1	B	350/383 (91%)	344 (98%)	6 (2%)	0	100	100
1	C	344/383 (90%)	338 (98%)	6 (2%)	0	100	100
1	D	362/383 (94%)	350 (97%)	12 (3%)	0	100	100
All	All	1399/1532 (91%)	1364 (98%)	35 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	296/344 (86%)	278 (94%)	18 (6%)	17	24
1	B	316/344 (92%)	295 (93%)	21 (7%)	15	20
1	C	310/344 (90%)	293 (94%)	17 (6%)	19	28
1	D	318/344 (92%)	294 (92%)	24 (8%)	12	16
All	All	1240/1376 (90%)	1160 (94%)	80 (6%)	15	21

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	THR
1	A	40	GLU
1	A	41	GLU
1	A	60	ARG
1	A	72	ARG
1	A	93	GLU
1	A	95	ARG
1	A	118	ILE
1	A	192	MET
1	A	193	VAL
1	A	212	CYS
1	A	219	ILE
1	A	225	GLU
1	A	236	ILE
1	A	271	ASP
1	A	305	ASP
1	A	309	VAL
1	A	356	SER
1	B	13	ILE
1	B	33	GLN
1	B	35	LEU
1	B	40	GLU
1	B	46	ILE
1	B	93	GLU
1	B	112	SER
1	B	116	GLN
1	B	119	ARG
1	B	132	ARG

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Mol	Chain	Res	Type
1	B	134	ILE
1	B	195	LEU
1	B	213	VAL
1	B	236	ILE
1	B	245	GLU
1	B	247	LEU
1	B	259	ILE
1	B	265	ILE
1	B	276	ARG
1	B	280	LEU
1	B	355	GLU
1	C	2	ILE
1	C	3	THR
1	C	8	GLU
1	C	12	GLU
1	C	41	GLU
1	C	67	SER
1	C	124	ILE
1	C	132	ARG
1	C	140	GLN
1	C	213	VAL
1	C	219	ILE
1	C	240	MET
1	C	242	GLU
1	C	245	GLU
1	C	247	LEU
1	C	256	ARG
1	C	309	VAL
1	D	7	THR
1	D	12	GLU
1	D	30	ARG
1	D	35	LEU
1	D	47	LYS
1	D	55	GLN
1	D	71	GLN
1	D	93	GLU
1	D	95	ARG
1	D	98	ARG
1	D	101	LYS
1	D	119	ARG
1	D	132	ARG
1	D	193	VAL

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Mol	Chain	Res	Type
1	D	195	LEU
1	D	213	VAL
1	D	219	ILE
1	D	259	ILE
1	D	302	VAL
1	D	309	VAL
1	D	333	HIS
1	D	338	THR
1	D	369	THR
1	D	372	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	97	GLN
1	A	179	ASN
1	A	323	HIS
1	B	55	GLN
1	B	66	HIS
1	C	36	HIS
1	C	52	ASN
1	C	97	GLN
1	C	140	GLN
1	C	187	HIS
1	C	206	ASN
1	C	257	GLN
1	D	55	GLN

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

Of 18 ligands modelled in this entry, 18 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	346/383 (90%)	0.18	16 (4%) 37 36	25, 74, 108, 135	1 (0%)
1	B	354/383 (92%)	-0.29	15 (4%) 40 40	27, 52, 104, 192	0
1	C	348/383 (90%)	-0.38	13 (3%) 45 45	25, 51, 105, 134	0
1	D	368/383 (96%)	-0.19	28 (7%) 20 17	25, 50, 119, 169	0
All	All	1416/1532 (92%)	-0.17	72 (5%) 33 31	25, 57, 109, 192	1 (0%)

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	377	LEU	19.6
1	B	363	PRO	10.0
1	B	362	ALA	9.5
1	D	368	TYR	9.0
1	D	375	ASP	9.0
1	A	350	ALA	8.6
1	C	339	ASP	7.7
1	D	0	SER	6.6
1	B	364	SER	6.6
1	A	358	PHE	6.5
1	D	349	VAL	6.2
1	B	361	LEU	6.2
1	D	371	LEU	6.1
1	D	1	GLU	5.7
1	C	358	PHE	5.6
1	D	366	TYR	5.6
1	D	376	GLU	5.5
1	D	367	ARG	5.2
1	B	360	LYS	4.9
1	A	337	PRO	4.8
1	A	221	PHE	4.7

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Mol	Chain	Res	Type	RSRZ
1	B	188	SER	4.2
1	B	338	THR	4.1
1	D	341	ALA	3.9
1	C	349	VAL	3.9
1	D	357	SER	3.8
1	B	187	HIS	3.8
1	B	358	PHE	3.7
1	D	358	PHE	3.7
1	C	357	SER	3.7
1	A	175	TYR	3.6
1	D	374	ARG	3.6
1	A	357	SER	3.6
1	B	350	ALA	3.5
1	B	189	ALA	3.5
1	C	350	ALA	3.4
1	D	339	ASP	3.3
1	B	1	GLU	3.3
1	D	370	LEU	3.2
1	D	340	THR	3.2
1	C	34	MET	3.2
1	A	297	PHE	3.0
1	B	359	GLN	3.0
1	A	1	GLU	3.0
1	A	320	PHE	2.9
1	C	338	THR	2.9
1	D	369	THR	2.9
1	D	372	ARG	2.9
1	D	345	GLN	2.8
1	C	337	PRO	2.8
1	D	373	ASP	2.7
1	D	343	GLY	2.7
1	A	351	SER	2.7
1	D	259	ILE	2.6
1	D	356	SER	2.5
1	A	187	HIS	2.5
1	B	297	PHE	2.4
1	C	297	PHE	2.4
1	A	48	GLU	2.4
1	A	125	THR	2.4
1	A	139	GLU	2.4
1	D	351	SER	2.3
1	C	2	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	298	ARG	2.3
1	C	356	SER	2.1
1	D	338	THR	2.1
1	B	357	SER	2.1
1	D	283	GLN	2.1
1	A	356	SER	2.1
1	C	8	GLU	2.0
1	A	55	GLN	2.0
1	D	350	ALA	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(Å ²)	Q<0.9
3	CL	B	405	1/1	0.96	0.09	56,56,56,56	0
3	CL	A	404	1/1	0.97	0.08	58,58,58,58	0
3	CL	D	403	1/1	0.97	0.09	53,53,53,53	0
3	CL	D	404	1/1	0.97	0.06	62,62,62,62	0
3	CL	C	403	1/1	0.98	0.12	60,60,60,60	0
3	CL	C	404	1/1	0.98	0.10	78,78,78,78	0
3	CL	B	404	1/1	0.98	0.10	57,57,57,57	0
3	CL	A	405	1/1	0.98	0.08	75,75,75,75	0
2	ZN	C	401	1/1	0.99	0.07	50,50,50,50	0
2	ZN	D	402	1/1	0.99	0.04	45,45,45,45	0
2	ZN	A	403	1/1	0.99	0.04	53,53,53,53	0
2	ZN	B	401	1/1	0.99	0.03	72,72,72,72	0
2	ZN	B	402	1/1	0.99	0.05	43,43,43,43	0
2	ZN	D	401	1/1	1.00	0.08	46,46,46,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	401	1/1	1.00	0.08	50,50,50,50	0
2	ZN	B	403	1/1	1.00	0.02	48,48,48,48	0
2	ZN	A	402	1/1	1.00	0.03	80,80,80,80	0
2	ZN	C	402	1/1	1.00	0.03	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.