



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:35 AM UTC

PDB ID : 5MTZ / pdb_00005mtz
Title : Crystal structure of a long form RNase Z from yeast
Authors : Li de la Sierra-Gallay, I.; Miao, M.; van Tilbeurgh, H.
Deposited on : 2017-01-11
Resolution : 2.99 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

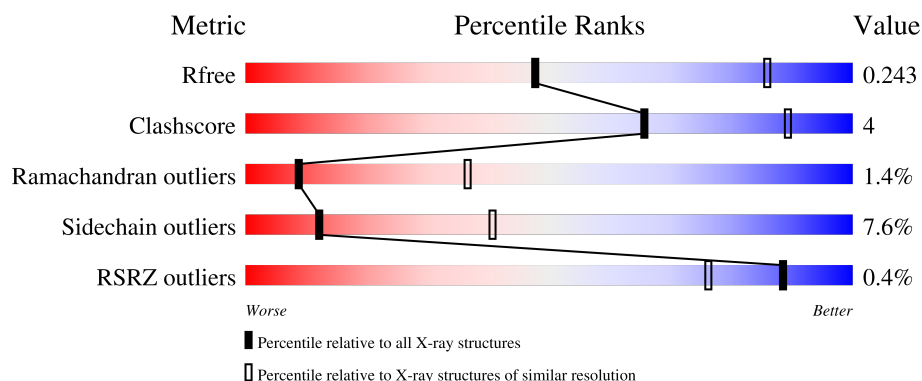
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.99 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	874	
1	B	874	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12472 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 Ribonuclease Z.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	768	Total	C	N	O	S	Se	0	0	0
			6284	4020	1046	1192	8	18			
1	B	755	Total	C	N	O	S	Se	0	0	0
			6174	3954	1031	1163	8	18			

There are 74 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-35	MSE	-	initiating methionine	UNP P36159
A	-34	ALA	-	expression tag	UNP P36159
A	-33	HIS	-	expression tag	UNP P36159
A	-32	HIS	-	expression tag	UNP P36159
A	-31	HIS	-	expression tag	UNP P36159
A	-30	HIS	-	expression tag	UNP P36159
A	-29	HIS	-	expression tag	UNP P36159
A	-28	HIS	-	expression tag	UNP P36159
A	-27	VAL	-	expression tag	UNP P36159
A	-26	GLY	-	expression tag	UNP P36159
A	-25	THR	-	expression tag	UNP P36159
A	-24	GLY	-	expression tag	UNP P36159
A	-23	SER	-	expression tag	UNP P36159
A	-22	ASN	-	expression tag	UNP P36159
A	-21	ASP	-	expression tag	UNP P36159
A	-20	ASP	-	expression tag	UNP P36159
A	-19	ASP	-	expression tag	UNP P36159
A	-18	ASP	-	expression tag	UNP P36159
A	-17	LYS	-	expression tag	UNP P36159
A	-16	SER	-	expression tag	UNP P36159
A	-15	PRO	-	expression tag	UNP P36159
A	-14	ASP	-	expression tag	UNP P36159
A	-13	PRO	-	expression tag	UNP P36159
A	-12	ASN	-	expression tag	UNP P36159
A	-11	TRP	-	expression tag	UNP P36159

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	GLU	-	expression tag	UNP P36159
A	-9	LEU	-	expression tag	UNP P36159
A	-8	VAL	-	expression tag	UNP P36159
A	-7	TYR	-	expression tag	UNP P36159
A	-6	THR	-	expression tag	UNP P36159
A	-5	ALA	-	expression tag	UNP P36159
A	-4	ARG	-	expression tag	UNP P36159
A	-3	LEU	-	expression tag	UNP P36159
A	-2	GLN	-	expression tag	UNP P36159
A	-1	GLU	-	expression tag	UNP P36159
A	0	PHE	-	expression tag	UNP P36159
A	1	MSE	-	expression tag	UNP P36159
B	-35	MSE	-	initiating methionine	UNP P36159
B	-34	ALA	-	expression tag	UNP P36159
B	-33	HIS	-	expression tag	UNP P36159
B	-32	HIS	-	expression tag	UNP P36159
B	-31	HIS	-	expression tag	UNP P36159
B	-30	HIS	-	expression tag	UNP P36159
B	-29	HIS	-	expression tag	UNP P36159
B	-28	HIS	-	expression tag	UNP P36159
B	-27	VAL	-	expression tag	UNP P36159
B	-26	GLY	-	expression tag	UNP P36159
B	-25	THR	-	expression tag	UNP P36159
B	-24	GLY	-	expression tag	UNP P36159
B	-23	SER	-	expression tag	UNP P36159
B	-22	ASN	-	expression tag	UNP P36159
B	-21	ASP	-	expression tag	UNP P36159
B	-20	ASP	-	expression tag	UNP P36159
B	-19	ASP	-	expression tag	UNP P36159
B	-18	ASP	-	expression tag	UNP P36159
B	-17	LYS	-	expression tag	UNP P36159
B	-16	SER	-	expression tag	UNP P36159
B	-15	PRO	-	expression tag	UNP P36159
B	-14	ASP	-	expression tag	UNP P36159
B	-13	PRO	-	expression tag	UNP P36159
B	-12	ASN	-	expression tag	UNP P36159
B	-11	TRP	-	expression tag	UNP P36159
B	-10	GLU	-	expression tag	UNP P36159
B	-9	LEU	-	expression tag	UNP P36159
B	-8	VAL	-	expression tag	UNP P36159
B	-7	TYR	-	expression tag	UNP P36159
B	-6	THR	-	expression tag	UNP P36159

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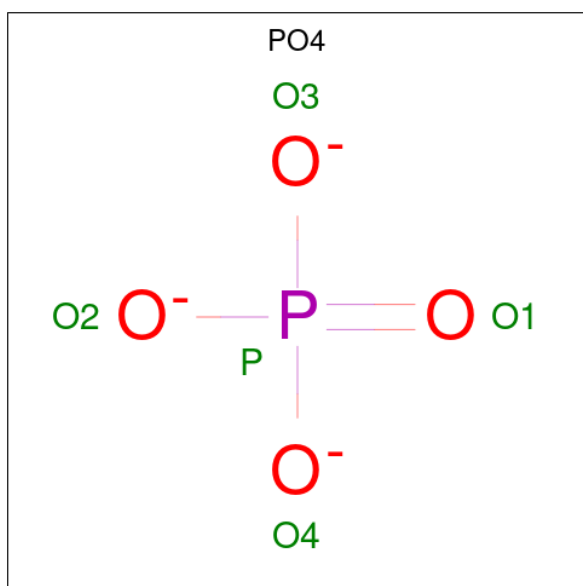
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	ALA	-	expression tag	UNP P36159
B	-4	ARG	-	expression tag	UNP P36159
B	-3	LEU	-	expression tag	UNP P36159
B	-2	GLN	-	expression tag	UNP P36159
B	-1	GLU	-	expression tag	UNP P36159
B	0	PHE	-	expression tag	UNP P36159
B	1	MSE	-	expression tag	UNP P36159

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

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

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).

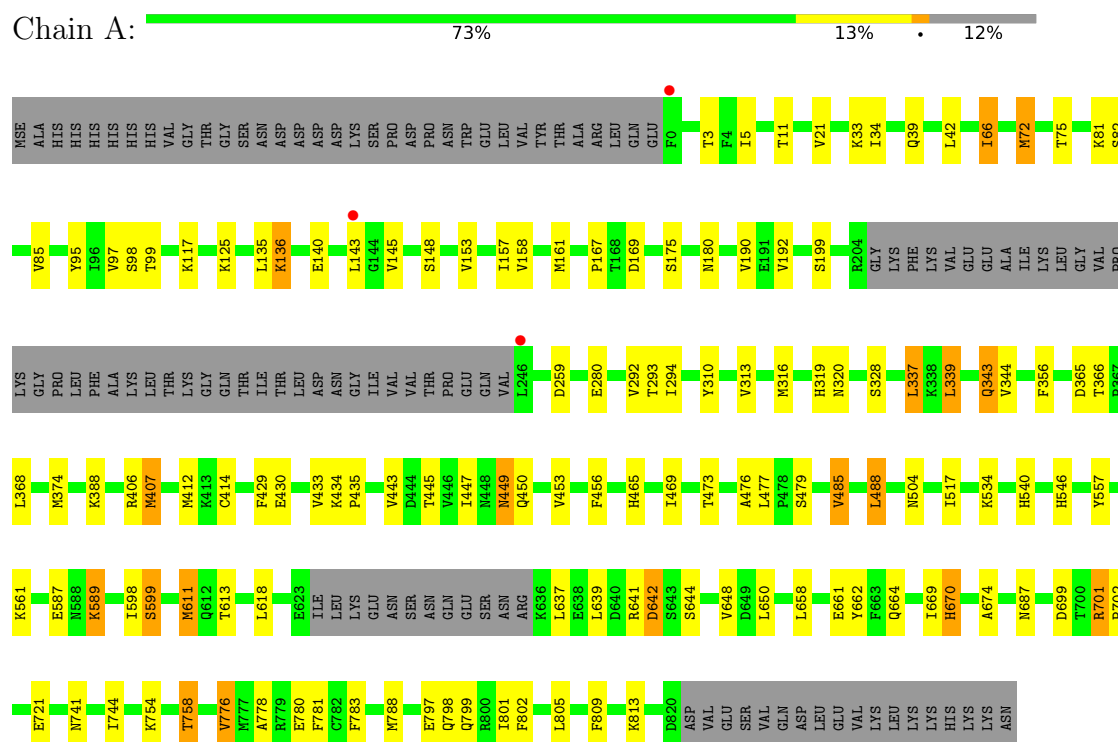


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

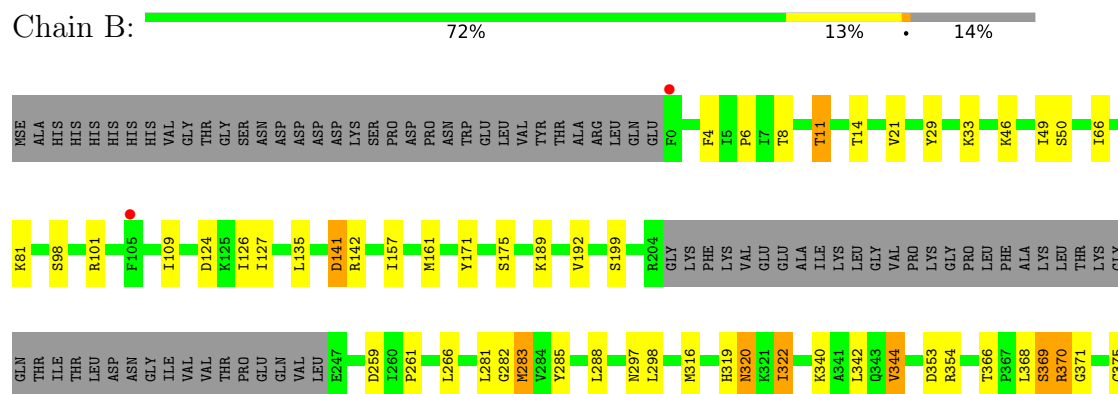
3 Residue-property plots

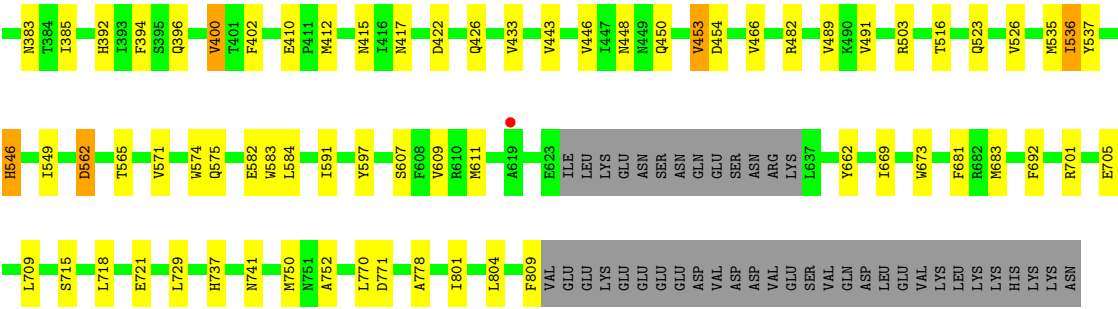
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Ribonuclease Z



• Molecule 1: Ribonuclease Z





4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	136.00Å 136.00Å 115.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.61 – 2.99 49.61 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.61-2.99) 99.8 (49.61-2.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.181 , 0.262 (Not available) , 0.243	Depositor DCC
R_{free} test set	2135 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	85.6	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.038 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12472	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% 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: PO4, ZN

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	0.74	0/6410	0.96	7/8638 (0.1%)
1	B	0.74	0/6300	0.95	5/8491 (0.1%)
All	All	0.74	0/12710	0.96	12/17129 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	433	VAL	N-CA-C	6.60	117.24	111.90
1	B	453	VAL	N-CA-C	6.22	116.82	108.11
1	A	701	ARG	N-CA-C	-5.78	102.11	110.08
1	B	126	ILE	CB-CA-C	5.46	115.76	110.63
1	A	517	ILE	CB-CA-C	-5.46	104.99	111.81
1	A	546	HIS	N-CA-C	5.43	121.67	114.12
1	A	485	VAL	CB-CA-C	-5.34	102.62	111.32
1	A	433	VAL	N-CA-C	5.29	116.56	112.12
1	A	66	ILE	N-CA-C	5.26	118.25	113.20
1	A	587	GLU	N-CA-C	5.12	116.72	111.03
1	B	124	ASP	N-CA-C	-5.08	101.69	109.41
1	B	546	HIS	N-CA-C	5.06	121.65	114.16

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	6284	0	6147	50	0
1	B	6174	0	6048	44	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
All	All	12472	0	12195	93	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 (93) 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:365:ASP:O	1:A:366:THR:HG22	1.78	0.83
1:A:157:ILE:HG22	1:A:161:MSE:HE2	1.65	0.79
1:B:681:PHE:HB2	1:B:683:MSE:CE	2.20	0.72
1:B:681:PHE:HB2	1:B:683:MSE:HE1	1.73	0.71
1:A:343:GLN:HE22	1:A:798:GLN:HE21	1.38	0.70
1:B:283:MSE:HE1	1:B:400:VAL:HG21	1.75	0.68
1:A:98:SER:HA	1:A:161:MSE:HE1	1.78	0.65
1:A:158:VAL:HA	1:A:161:MSE:HE3	1.78	0.64
1:B:98:SER:HA	1:B:161:MSE:HE1	1.82	0.61
1:A:280:GLU:HA	1:A:412:MSE:HE1	1.84	0.59
1:A:445:THR:HA	1:A:449:ASN:HD21	1.66	0.59
1:B:426:GLN:NE2	1:B:448:ASN:OD1	2.35	0.59
1:A:754:LYS:HE2	1:A:780:GLU:OE1	2.03	0.58
1:B:157:ILE:HG22	1:B:161:MSE:HE3	1.85	0.58
1:B:523:GLN:HA	1:B:526:VAL:HG12	1.86	0.58
1:A:153:VAL:HG21	1:A:618:LEU:HD11	1.87	0.57
1:B:718:LEU:HB2	1:B:750:MSE:HE1	1.86	0.57
1:A:343:GLN:NE2	1:A:798:GLN:HE21	2.02	0.56
1:A:788:MSE:HE1	1:A:798:GLN:HE22	1.69	0.56
1:A:741:ASN:HA	1:A:744:ILE:HG22	1.88	0.56
1:B:266:LEU:HD21	1:B:298:LEU:HA	1.88	0.55
1:A:337:LEU:HD11	1:A:447:ILE:HD11	1.89	0.55
1:B:342:LEU:HB3	1:B:801:ILE:HD13	1.89	0.55
1:A:356:PHE:HB2	1:A:368:LEU:HD21	1.88	0.55
1:A:598:ILE:O	1:A:599:SER:CB	2.54	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:536:ILE:HD11	1:B:549:ILE:HG12	1.90	0.54
1:A:799:GLN:HA	1:A:802:PHE:CD1	2.43	0.53
1:A:598:ILE:HD13	1:A:650:LEU:HD23	1.92	0.50
1:B:50:SER:HA	1:B:81:LYS:HE2	1.93	0.50
1:A:33:LYS:HE2	1:A:259:ASP:OD2	2.11	0.50
1:B:11:THR:HG23	1:B:14:THR:OG1	2.12	0.50
1:A:741:ASN:O	1:A:744:ILE:HG22	2.12	0.50
1:B:683:MSE:HE3	1:B:692:PHE:HB3	1.92	0.49
1:A:42:LEU:HD12	1:A:72:MSE:HE1	1.94	0.49
1:B:33:LYS:HE2	1:B:259:ASP:OD2	2.13	0.49
1:A:97:VAL:HG23	1:A:158:VAL:HG22	1.95	0.49
1:A:366:THR:O	1:A:366:THR:HG23	2.13	0.48
1:B:283:MSE:CE	1:B:400:VAL:HG21	2.42	0.48
1:A:406:ARG:O	1:A:407:MSE:C	2.57	0.48
1:A:557:TYR:O	1:A:561:LYS:HB2	2.14	0.48
1:A:473:THR:HA	1:A:758:THR:HG22	1.97	0.47
1:B:340:LYS:O	1:B:344:VAL:N	2.49	0.46
1:A:95:TYR:O	1:A:99:THR:HG23	2.16	0.46
1:A:180:ASN:HD22	1:A:613:THR:HB	1.81	0.46
1:A:456:PHE:CE1	1:A:469:ILE:HD11	2.51	0.46
1:B:575:GLN:HE22	1:B:673:TRP:CD1	2.33	0.46
1:A:670:HIS:CD2	1:A:699:ASP:HB2	2.50	0.46
1:B:157:ILE:HG22	1:B:161:MSE:CE	2.44	0.46
1:A:11:THR:HG21	1:A:328:SER:OG	2.17	0.45
1:A:135:LEU:O	1:A:136:LYS:C	2.59	0.45
1:B:535:MSE:HE3	1:B:537:TYR:HB2	1.99	0.45
1:B:574:TRP:HA	1:B:597:TYR:OH	2.17	0.44
1:A:98:SER:HA	1:A:161:MSE:CE	2.46	0.44
1:A:294:ILE:HD11	1:A:319:HIS:ND1	2.32	0.44
1:A:429:PHE:CD1	1:A:443:VAL:HG22	2.53	0.44
1:A:450:GLN:NE2	1:A:453:VAL:HG11	2.33	0.44
1:B:701:ARG:NH1	1:B:737:HIS:O	2.49	0.44
1:B:6:PRO:HB3	1:B:394:PHE:CZ	2.53	0.43
1:B:282:GLY:HA2	1:B:412:MSE:HE1	2.01	0.43
1:A:669:ILE:HB	1:A:701:ARG:HB3	1.99	0.43
1:B:750:MSE:HE2	1:B:752:ALA:HB2	2.00	0.43
1:A:744:ILE:HD11	1:A:776:VAL:HG13	2.01	0.43
1:B:584:LEU:HD13	1:B:591:ILE:HB	2.01	0.43
1:A:534:LYS:HB3	1:A:658:LEU:HD22	2.01	0.42
1:A:611:MSE:HE2	1:A:639:LEU:HB2	2.01	0.42
1:B:319:HIS:HB3	1:B:322:ILE:HB	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:340:LYS:HG3	1:B:344:VAL:HG12	2.01	0.42
1:B:283:MSE:HE3	1:B:402:PHE:CZ	2.55	0.42
1:A:339:LEU:HD13	1:A:809:PHE:HZ	1.85	0.42
1:B:135:LEU:HG	1:B:192:VAL:HG12	2.01	0.42
1:B:369:SER:O	1:B:371:GLY:N	2.53	0.42
1:B:583:TRP:CD1	1:B:583:TRP:C	2.98	0.42
1:A:642:ASP:OD2	1:B:50:SER:OG	2.38	0.41
1:B:8:THR:HA	1:B:396:GLN:HE21	1.84	0.41
1:B:546:HIS:HB2	1:B:549:ILE:HD12	2.01	0.41
1:A:701:ARG:HG3	1:A:702:PRO:HD2	2.02	0.41
1:B:101:ARG:NH2	1:B:582:GLU:OE1	2.54	0.41
1:B:715:SER:OG	1:B:750:MSE:HE3	2.20	0.41
1:A:598:ILE:O	1:A:599:SER:HB3	2.20	0.41
1:B:741:ASN:HD21	1:B:771:ASP:HB2	1.86	0.41
1:A:488:LEU:HD21	1:A:504:ASN:HB3	2.01	0.41
1:B:415:ASN:HD21	1:B:417:ASN:HB2	1.85	0.41
1:B:450:GLN:HA	1:B:453:VAL:HG12	2.02	0.41
1:A:540:HIS:HB3	1:A:674:ALA:HA	2.03	0.41
1:B:21:VAL:HG22	1:B:29:TYR:HB2	2.03	0.41
1:B:285:TYR:CE1	1:B:316:MSE:HG2	2.56	0.41
1:B:320:ASN:O	1:B:354:ARG:NH1	2.54	0.41
1:A:310:TYR:CD1	1:A:310:TYR:N	2.89	0.40
1:A:776:VAL:HG11	1:A:781:PHE:CE1	2.56	0.40
1:A:783:PHE:O	1:A:788:MSE:HE1	2.20	0.40
1:A:66:ILE:C	1:A:66:ILE:HD12	2.46	0.40
1:A:434:LYS:HB3	1:A:435:PRO:HD3	2.04	0.40
1:B:4:PHE:HB2	1:B:400:VAL:HG13	2.02	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	762/874 (87%)	686 (90%)	66 (9%)	10 (1%)	9	38
1	B	749/874 (86%)	669 (89%)	69 (9%)	11 (2%)	8	35
All	All	1511/1748 (86%)	1355 (90%)	135 (9%)	21 (1%)	9	36

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	141	ASP
1	B	370	ARG
1	A	145	VAL
1	A	167	PRO
1	A	407	MSE
1	A	599	SER
1	A	778	ALA
1	A	136	LYS
1	A	476	ALA
1	B	368	LEU
1	B	562	ASP
1	A	589	LYS
1	A	670	HIS
1	B	142	ARG
1	B	171	TYR
1	B	261	PRO
1	A	797	GLU
1	B	778	ALA
1	B	322	ILE
1	B	109	ILE
1	B	344	VAL

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	707/783 (90%)	651 (92%)	56 (8%)	11	39
1	B	694/783 (89%)	644 (93%)	50 (7%)	13	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1401/1566 (90%)	1295 (92%)	106 (8%)	12	41

All (106) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	5	ILE
1	A	21	VAL
1	A	34	ILE
1	A	39	GLN
1	A	72	MSE
1	A	75	THR
1	A	81	LYS
1	A	82	SER
1	A	85	VAL
1	A	117	LYS
1	A	125	LYS
1	A	140	GLU
1	A	143	LEU
1	A	148	SER
1	A	169	ASP
1	A	175	SER
1	A	190	VAL
1	A	192	VAL
1	A	199	SER
1	A	292	VAL
1	A	293	THR
1	A	313	VAL
1	A	316	MSE
1	A	320	ASN
1	A	337	LEU
1	A	339	LEU
1	A	343	GLN
1	A	344	VAL
1	A	374	MSE
1	A	388	LYS
1	A	414	CYS
1	A	430	GLU
1	A	449	ASN
1	A	465	HIS
1	A	477	LEU
1	A	479	SER

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Mol	Chain	Res	Type
1	A	485	VAL
1	A	488	LEU
1	A	589	LYS
1	A	611	MSE
1	A	637	LEU
1	A	641	ARG
1	A	642	ASP
1	A	644	SER
1	A	648	VAL
1	A	661	GLU
1	A	662	TYR
1	A	664	GLN
1	A	687	ASN
1	A	721	GLU
1	A	758	THR
1	A	776	VAL
1	A	801	ILE
1	A	805	LEU
1	A	813	LYS
1	B	11	THR
1	B	46	LYS
1	B	49	ILE
1	B	66	ILE
1	B	127	ILE
1	B	141	ASP
1	B	175	SER
1	B	189	LYS
1	B	199	SER
1	B	281	LEU
1	B	283	MSE
1	B	288	LEU
1	B	297	ASN
1	B	320	ASN
1	B	353	ASP
1	B	366	THR
1	B	369	SER
1	B	370	ARG
1	B	375	CYS
1	B	383	ASN
1	B	385	ILE
1	B	392	HIS
1	B	400	VAL

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Mol	Chain	Res	Type
1	B	410	GLU
1	B	422	ASP
1	B	443	VAL
1	B	446	VAL
1	B	454	ASP
1	B	466	VAL
1	B	482	ARG
1	B	489	VAL
1	B	491	VAL
1	B	503	ARG
1	B	516	THR
1	B	536	ILE
1	B	562	ASP
1	B	565	THR
1	B	571	VAL
1	B	607	SER
1	B	609	VAL
1	B	611	MSE
1	B	662	TYR
1	B	669	ILE
1	B	705	GLU
1	B	709	LEU
1	B	721	GLU
1	B	729	LEU
1	B	770	LEU
1	B	804	LEU
1	B	809	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	HIS
1	A	79	GLN
1	A	94	ASN
1	A	449	ASN
1	A	455	ASN
1	A	751	ASN
1	A	798	GLN
1	A	799	GLN
1	B	62	ASN
1	B	94	ASN
1	B	251	HIS

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Mol	Chain	Res	Type
1	B	396	GLN
1	B	415	ASN
1	B	449	ASN
1	B	614	GLN
1	B	653	GLN
1	B	664	GLN
1	B	686	ASN
1	B	769	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PO4	B	903	2	4,4,4	0.82	0	6,6,6	1.01	0
3	PO4	A	903	2	4,4,4	1.02	0	6,6,6	0.58	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	750/874 (85%)	-0.38	3 (0%)	88 76	50, 80, 121, 166	0
1	B	737/874 (84%)	-0.34	3 (0%)	88 76	51, 80, 134, 171	0
All	All	1487/1748 (85%)	-0.36	6 (0%)	88 76	50, 80, 128, 171	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	0	PHE	4.1
1	A	246	LEU	3.1
1	B	0	PHE	2.6
1	A	143	LEU	2.6
1	B	105	PHE	2.1
1	B	619	ALA	2.1

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
3	PO4	A	903	5/5	0.95	0.06	77,89,95,102	0
3	PO4	B	903	5/5	0.96	0.05	75,78,85,86	0
2	ZN	B	901	1/1	0.99	0.04	73,73,73,73	0
2	ZN	B	902	1/1	1.00	0.01	68,68,68,68	0
2	ZN	A	902	1/1	1.00	0.02	73,73,73,73	0
2	ZN	A	901	1/1	1.00	0.03	83,83,83,83	0

6.5 Other polymers [i](#)

There are no such residues in this entry.