



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 02:12 PM UTC

PDB ID : 209L / pdb_0000209L
Title : PROTEIN STRUCTURE PLASTICITY EXEMPLIFIED BY INSERTION
AND DELETION MUTANTS IN T4 LYSOZYME
Authors : Vetter, I.R.; Baase, W.A.; Heinz, D.W.; Xiong, J.-P.; Snow, S.; Matthews,
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Deposited on : 1996-09-23
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

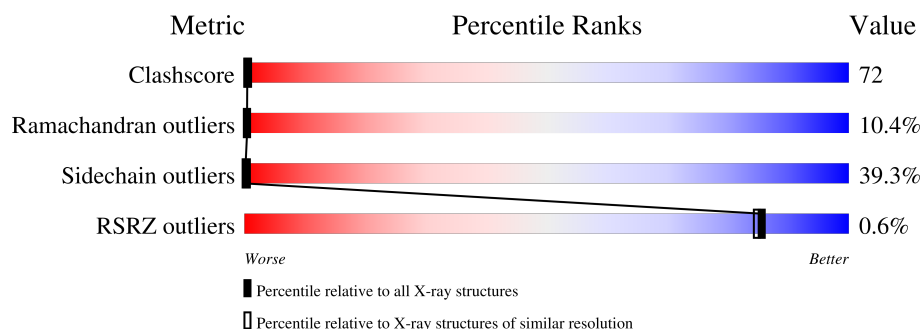
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

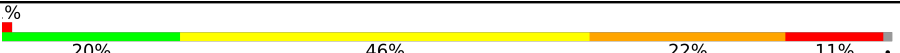
The reported resolution of this entry is 2.70 Å.

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))
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	167	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1328 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 T4 LYSOZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	165	Total	C	N	O	S	0	0	0
			1307	823	238	241	5			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	54	THR	CYS	engineered mutation	UNP P00720
A	73A	ALA	-	insertion	UNP P00720
A	73B	ALA	-	insertion	UNP P00720
A	73C	ALA	-	insertion	UNP P00720
A	97	ALA	CYS	engineered mutation	UNP P00720

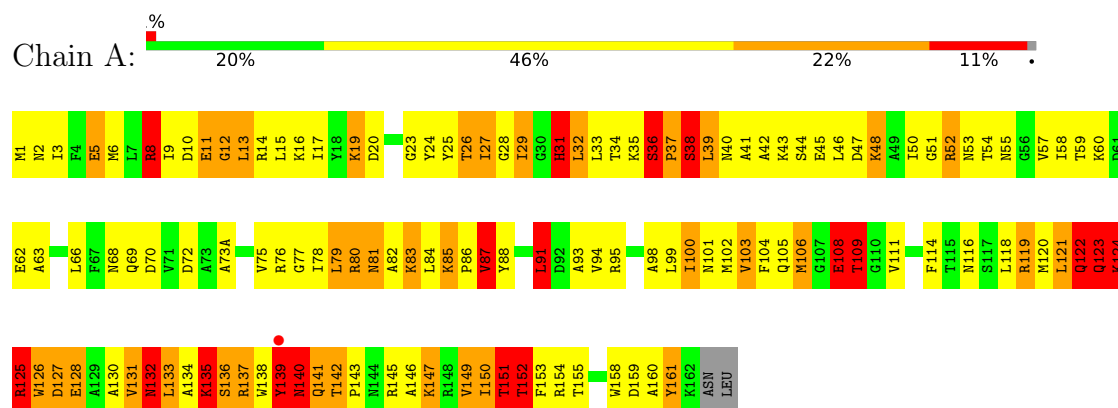
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	21	Total	O	0	0
			21	21		

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: T4 LYSOZYME



4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	95.85Å 95.85Å 117.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	21.37 – 2.70 21.37 – 2.77	Depositor EDS
% Data completeness (in resolution range)	79.5 (21.37-2.70) 85.3 (21.37-2.77)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 2.79Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.179 , (Not available) 0.176 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 222.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	1328	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% 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	1.34	4/1327 (0.3%)	2.09	34/1788 (1.9%)

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	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	35	LYS	C-O	8.17	1.34	1.24
1	A	36	SER	CA-C	7.67	1.62	1.52
1	A	151	THR	CA-C	-6.72	1.43	1.52
1	A	11	GLU	CD-OE2	5.02	1.34	1.25

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	36	SER	CA-C-N	22.55	148.03	119.84
1	A	36	SER	C-N-CA	22.55	148.03	119.84
1	A	36	SER	O-C-N	-14.44	104.71	121.32
1	A	80	ARG	N-CA-C	-9.59	100.52	110.97
1	A	37	PRO	N-CA-CB	9.07	112.77	103.25
1	A	150	ILE	CB-CA-C	-8.67	97.07	111.29
1	A	37	PRO	N-CA-C	8.03	129.00	112.47
1	A	135	LYS	N-CA-C	-7.90	99.36	110.50
1	A	132	ASN	CA-CB-CG	-7.71	104.89	112.60
1	A	87	VAL	N-CA-CB	7.34	121.59	110.58
1	A	8	ARG	N-CA-C	-6.76	103.99	111.36
1	A	36	SER	CA-C-O	6.59	129.19	120.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	31	HIS	CA-CB-CG	-6.55	107.25	113.80
1	A	109	THR	CA-CB-CG2	6.49	121.53	110.50
1	A	108	GLU	N-CA-CB	6.48	121.44	110.49
1	A	17	ILE	N-CA-CB	6.44	118.55	110.05
1	A	109	THR	N-CA-CB	6.43	121.36	110.49
1	A	151	THR	N-CA-C	-6.08	97.84	110.80
1	A	12	GLY	CA-C-N	-5.99	114.36	122.86
1	A	12	GLY	C-N-CA	-5.99	114.36	122.86
1	A	37	PRO	CA-N-CD	-5.87	103.78	112.00
1	A	161	TYR	N-CA-C	-5.83	106.00	113.23
1	A	38	SER	N-CA-C	5.82	117.70	108.67
1	A	68	ASN	CA-CB-CG	-5.82	106.78	112.60
1	A	137	ARG	N-CA-C	-5.80	103.94	111.02
1	A	94	VAL	N-CA-C	5.75	116.40	110.36
1	A	128	GLU	N-CA-C	5.73	118.30	111.71
1	A	124	LYS	N-CA-CB	5.55	121.09	110.82
1	A	150	ILE	CA-CB-CG2	-5.45	101.24	110.50
1	A	139	TYR	CA-CB-CG	-5.40	104.18	113.90
1	A	100	ILE	CB-CA-C	-5.28	105.12	111.88
1	A	91	LEU	O-C-N	5.18	128.81	122.75
1	A	152	THR	N-CA-C	-5.09	105.74	111.28
1	A	36	SER	C-N-CD	-5.04	104.32	125.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	36	SER	Peptide

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	1307	0	1334	190	0
2	A	21	0	0	3	0
All	All	1328	0	1334	190	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 72.

All (190) 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:83:LYS:NZ	1:A:85:LYS:H	1.45	1.15
1:A:13:LEU:HD12	1:A:29:ILE:HD11	1.41	1.02
1:A:3:ILE:HD12	1:A:6:MET:HE2	1.43	1.00
1:A:83:LYS:HZ3	1:A:85:LYS:N	1.63	0.96
1:A:26:THR:HG23	1:A:32:LEU:HA	1.45	0.96
1:A:6:MET:HE3	1:A:101:ASN:HB2	1.50	0.92
1:A:6:MET:CE	1:A:101:ASN:HB2	2.06	0.85
1:A:26:THR:CG2	1:A:32:LEU:HA	2.09	0.82
1:A:79:LEU:HD13	1:A:88:TYR:HB2	1.59	0.82
1:A:83:LYS:HZ3	1:A:85:LYS:H	0.83	0.80
1:A:128:GLU:HA	1:A:131:VAL:HG13	1.65	0.79
1:A:48:LYS:H	1:A:48:LYS:HZ3	1.32	0.77
1:A:98:ALA:O	1:A:102:MET:HG3	1.83	0.77
1:A:50:ILE:HG22	1:A:52:ARG:HG2	1.68	0.76
1:A:79:LEU:HB3	1:A:84:LEU:CD1	2.17	0.74
1:A:5:GLU:O	1:A:9:ILE:HD13	1.88	0.74
1:A:133:LEU:HD23	1:A:134:ALA:N	2.01	0.74
1:A:2:ASN:HB2	2:A:194:HOH:O	1.88	0.74
1:A:106:MET:HE3	1:A:106:MET:HA	1.67	0.74
1:A:83:LYS:HG3	1:A:84:LEU:H	1.53	0.73
1:A:140:ASN:ND2	1:A:140:ASN:H	1.84	0.73
1:A:84:LEU:HB2	1:A:87:VAL:HG13	1.70	0.73
1:A:19:LYS:HA	1:A:24:TYR:O	1.87	0.73
1:A:84:LEU:HB2	1:A:87:VAL:CG1	2.19	0.72
1:A:121:LEU:HD13	1:A:125:ARG:HB3	1.71	0.72
1:A:13:LEU:HD12	1:A:29:ILE:CD1	2.20	0.71
1:A:15:LEU:HD23	1:A:58:ILE:HG13	1.73	0.71
1:A:48:LYS:H	1:A:48:LYS:NZ	1.89	0.70
1:A:79:LEU:HB3	1:A:84:LEU:HD13	1.74	0.70
1:A:3:ILE:HD12	1:A:6:MET:CE	2.21	0.69
1:A:27:ILE:HG13	1:A:28:GLY:H	1.56	0.69
1:A:83:LYS:NZ	1:A:85:LYS:N	2.30	0.69
1:A:6:MET:HE3	1:A:101:ASN:HD22	1.56	0.68
1:A:38:SER:OG	1:A:40:ASN:ND2	2.27	0.68
1:A:140:ASN:H	1:A:140:ASN:HD22	1.41	0.68
1:A:83:LYS:HA	1:A:83:LYS:HE2	1.78	0.66
1:A:102:MET:O	1:A:137:ARG:NH1	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ILE:N	1:A:29:ILE:HD13	2.10	0.66
1:A:151:THR:O	1:A:155:THR:N	2.28	0.66
1:A:44:SER:O	1:A:48:LYS:NZ	2.28	0.66
1:A:141:GLN:HB2	1:A:142:THR:HG22	1.78	0.65
1:A:48:LYS:HZ3	1:A:48:LYS:N	1.95	0.65
1:A:77:GLY:O	1:A:81:ASN:ND2	2.30	0.65
1:A:142:THR:N	1:A:143:PRO:HD3	2.12	0.65
1:A:50:ILE:HD13	1:A:50:ILE:N	2.11	0.65
1:A:8:ARG:HG3	1:A:8:ARG:NH1	2.13	0.64
1:A:26:THR:HG22	1:A:31:HIS:C	2.23	0.64
1:A:26:THR:HG22	1:A:31:HIS:O	1.97	0.63
1:A:128:GLU:O	1:A:132:ASN:HB2	1.97	0.63
1:A:45:GLU:HA	1:A:48:LYS:NZ	2.13	0.63
1:A:50:ILE:CG2	1:A:52:ARG:HG2	2.28	0.63
1:A:48:LYS:HZ3	1:A:48:LYS:CB	2.06	0.63
1:A:27:ILE:HG13	1:A:28:GLY:N	2.14	0.63
1:A:122:GLN:HA	1:A:122:GLN:OE1	1.99	0.63
1:A:3:ILE:HG22	2:A:194:HOH:O	1.98	0.62
1:A:124:LYS:HD3	1:A:124:LYS:C	2.22	0.62
1:A:1:MET:HG3	1:A:2:ASN:N	2.15	0.62
1:A:11:GLU:HB2	1:A:29:ILE:HG22	1.82	0.62
1:A:83:LYS:O	1:A:84:LEU:HD12	2.00	0.61
1:A:85:LYS:N	1:A:86:PRO:HD2	2.15	0.61
1:A:143:PRO:O	1:A:147:LYS:HG3	2.00	0.61
1:A:108:GLU:O	1:A:109:THR:HG22	2.00	0.61
1:A:151:THR:O	1:A:155:THR:HG23	2.00	0.61
1:A:124:LYS:HD3	1:A:125:ARG:N	2.16	0.60
1:A:47:ASP:O	1:A:51:GLY:N	2.31	0.60
1:A:79:LEU:CD1	1:A:88:TYR:HD2	2.15	0.60
1:A:106:MET:HE3	1:A:106:MET:CA	2.32	0.60
1:A:95:ARG:NE	1:A:153:PHE:O	2.35	0.59
1:A:140:ASN:HB2	1:A:141:GLN:HG3	1.85	0.59
1:A:59:THR:H	1:A:62:GLU:HB2	1.68	0.59
1:A:83:LYS:HG3	1:A:84:LEU:N	2.17	0.58
1:A:75:VAL:HG23	1:A:103:VAL:HG11	1.84	0.58
1:A:82:ALA:O	1:A:84:LEU:HG	2.03	0.58
1:A:70:ASP:HB3	1:A:104:PHE:CE2	2.40	0.57
1:A:95:ARG:HH21	1:A:154:ARG:C	2.12	0.57
1:A:84:LEU:O	1:A:87:VAL:HG13	2.05	0.57
1:A:34:THR:CG2	1:A:42:ALA:HB2	2.35	0.57
1:A:8:ARG:HG3	1:A:8:ARG:HH11	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:GLU:HA	1:A:131:VAL:CG1	2.34	0.56
1:A:119:ARG:CZ	1:A:119:ARG:HB3	2.36	0.55
1:A:106:MET:HA	1:A:106:MET:CE	2.34	0.55
1:A:140:ASN:HD22	1:A:140:ASN:N	2.03	0.55
1:A:124:LYS:NZ	1:A:127:ASP:HB3	2.22	0.55
1:A:121:LEU:O	1:A:122:GLN:HB2	2.06	0.55
1:A:146:ALA:O	1:A:150:ILE:HD12	2.06	0.55
1:A:123:GLN:O	1:A:126:TRP:HD1	1.90	0.55
1:A:83:LYS:HA	1:A:83:LYS:CE	2.34	0.54
1:A:151:THR:N	1:A:154:ARG:HG2	2.22	0.54
1:A:141:GLN:C	1:A:143:PRO:HD3	2.32	0.54
1:A:58:ILE:HD12	1:A:63:ALA:HB2	1.90	0.53
1:A:102:MET:HE1	1:A:130:ALA:C	2.33	0.53
1:A:79:LEU:O	1:A:83:LYS:O	2.27	0.53
1:A:73(A):ALA:CB	1:A:100:ILE:HG23	2.38	0.52
1:A:77:GLY:O	1:A:80:ARG:HB3	2.08	0.52
1:A:150:ILE:HG22	1:A:154:ARG:HG2	1.92	0.52
1:A:151:THR:HA	1:A:154:ARG:CG	2.40	0.52
1:A:45:GLU:O	1:A:48:LYS:NZ	2.39	0.52
1:A:6:MET:SD	1:A:158:TRP:HZ3	2.33	0.51
1:A:73(A):ALA:HB3	1:A:100:ILE:HG23	1.91	0.51
1:A:140:ASN:ND2	1:A:140:ASN:N	2.55	0.51
1:A:145:ARG:NH1	1:A:145:ARG:HG3	2.26	0.51
1:A:51:GLY:O	1:A:52:ARG:HB3	2.10	0.50
1:A:151:THR:CA	1:A:154:ARG:HG2	2.41	0.50
1:A:13:LEU:HB2	1:A:29:ILE:HD12	1.93	0.50
1:A:15:LEU:HD22	1:A:58:ILE:O	2.10	0.50
1:A:15:LEU:HD22	1:A:58:ILE:C	2.37	0.50
1:A:28:GLY:C	1:A:29:ILE:HD13	2.37	0.50
1:A:114:PHE:O	1:A:118:LEU:HB2	2.11	0.50
1:A:145:ARG:HG3	1:A:145:ARG:HH11	1.77	0.50
1:A:84:LEU:HD12	1:A:84:LEU:C	2.37	0.49
1:A:85:LYS:N	1:A:86:PRO:CD	2.75	0.49
1:A:45:GLU:C	1:A:48:LYS:HZ1	2.20	0.49
1:A:58:ILE:CD1	1:A:63:ALA:HB2	2.42	0.49
1:A:45:GLU:CA	1:A:48:LYS:NZ	2.76	0.49
1:A:124:LYS:HZ1	1:A:127:ASP:HB3	1.78	0.49
1:A:20:ASP:O	1:A:23:GLY:N	2.42	0.48
1:A:15:LEU:HD23	1:A:58:ILE:CG1	2.42	0.48
1:A:58:ILE:HA	1:A:62:GLU:OE1	2.13	0.48
1:A:6:MET:HE3	1:A:101:ASN:CB	2.32	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ASP:OD1	1:A:161:TYR:HE1	1.95	0.48
1:A:83:LYS:HZ3	1:A:83:LYS:C	2.21	0.48
1:A:138:TRP:HB2	1:A:146:ALA:HB2	1.94	0.48
1:A:55:ASN:N	2:A:179:HOH:O	2.44	0.48
1:A:38:SER:HB3	1:A:40:ASN:HD22	1.79	0.47
1:A:6:MET:HE3	1:A:101:ASN:ND2	2.26	0.47
1:A:34:THR:HG22	1:A:42:ALA:CB	2.45	0.47
1:A:38:SER:O	1:A:41:ALA:N	2.47	0.47
1:A:46:LEU:O	1:A:46:LEU:HD12	2.15	0.47
1:A:76:ARG:HA	1:A:88:TYR:CD2	2.50	0.47
1:A:81:ASN:O	1:A:82:ALA:HB3	2.14	0.47
1:A:8:ARG:HH12	1:A:12:GLY:HA2	1.80	0.46
1:A:82:ALA:HA	1:A:119:ARG:HG2	1.96	0.46
1:A:70:ASP:HB3	1:A:104:PHE:HE2	1.80	0.46
1:A:48:LYS:NZ	1:A:48:LYS:CB	2.76	0.46
1:A:1:MET:O	1:A:2:ASN:HB3	2.16	0.45
1:A:1:MET:CG	1:A:2:ASN:N	2.80	0.45
1:A:25:TYR:HD1	1:A:25:TYR:HA	1.66	0.45
1:A:135:LYS:HB2	1:A:135:LYS:HE3	1.58	0.45
1:A:106:MET:HG2	1:A:111:VAL:HG22	1.98	0.45
1:A:6:MET:HE1	1:A:101:ASN:HB2	1.93	0.45
1:A:149:VAL:O	1:A:152:THR:HB	2.17	0.45
1:A:83:LYS:NZ	1:A:84:LEU:N	2.63	0.45
1:A:79:LEU:CD1	1:A:88:TYR:CD2	2.98	0.45
1:A:133:LEU:HD23	1:A:133:LEU:C	2.42	0.45
1:A:43:LYS:O	1:A:46:LEU:HB3	2.18	0.44
1:A:54:THR:O	1:A:55:ASN:HB3	2.18	0.44
1:A:38:SER:CB	1:A:40:ASN:HD22	2.31	0.44
1:A:45:GLU:HA	1:A:48:LYS:CE	2.47	0.44
1:A:102:MET:HE1	1:A:130:ALA:O	2.18	0.44
1:A:34:THR:HG22	1:A:42:ALA:HA	1.98	0.44
1:A:34:THR:HG22	1:A:42:ALA:HB2	1.98	0.44
1:A:44:SER:C	1:A:48:LYS:HZ2	2.20	0.44
1:A:83:LYS:CG	1:A:84:LEU:N	2.79	0.44
1:A:52:ARG:NH1	1:A:54:THR:HA	2.33	0.43
1:A:136:SER:OG	1:A:137:ARG:N	2.50	0.43
1:A:83:LYS:CE	1:A:83:LYS:CA	2.96	0.43
1:A:151:THR:HG22	1:A:152:THR:N	2.29	0.43
1:A:46:LEU:HD12	1:A:46:LEU:C	2.42	0.43
1:A:139:TYR:O	1:A:143:PRO:HG3	2.19	0.43
1:A:73(A):ALA:HB3	1:A:100:ILE:CG2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:ARG:HH11	1:A:8:ARG:CG	2.32	0.43
1:A:25:TYR:O	1:A:33:LEU:N	2.51	0.42
1:A:76:ARG:O	1:A:80:ARG:N	2.52	0.42
1:A:126:TRP:CE3	1:A:126:TRP:C	2.98	0.42
1:A:102:MET:HA	1:A:138:TRP:CH2	2.54	0.42
1:A:142:THR:N	1:A:143:PRO:CD	2.82	0.42
1:A:48:LYS:HZ3	1:A:48:LYS:CA	2.32	0.42
1:A:91:LEU:N	1:A:91:LEU:HD23	2.34	0.42
1:A:118:LEU:HD12	1:A:118:LEU:HA	1.72	0.42
1:A:12:GLY:O	1:A:29:ILE:HA	2.20	0.42
1:A:132:ASN:HD22	1:A:132:ASN:HA	1.54	0.42
1:A:45:GLU:CA	1:A:48:LYS:HZ2	2.32	0.41
1:A:108:GLU:HG3	1:A:111:VAL:HB	2.01	0.41
1:A:54:THR:HB	1:A:57:VAL:O	2.20	0.41
1:A:114:PHE:O	1:A:118:LEU:N	2.53	0.41
1:A:151:THR:HA	1:A:154:ARG:HG2	2.00	0.41
1:A:103:VAL:O	1:A:106:MET:HB2	2.20	0.41
1:A:119:ARG:CZ	1:A:119:ARG:CB	2.98	0.41
1:A:19:LYS:HD3	1:A:25:TYR:CE1	2.55	0.41
1:A:45:GLU:C	1:A:48:LYS:NZ	2.78	0.41
1:A:125:ARG:C	1:A:127:ASP:N	2.78	0.41
1:A:102:MET:HG2	1:A:149:VAL:HG11	2.02	0.41
1:A:27:ILE:CG1	1:A:28:GLY:N	2.81	0.41
1:A:104:PHE:C	1:A:104:PHE:CD1	2.98	0.41
1:A:119:ARG:C	1:A:121:LEU:N	2.79	0.41
1:A:151:THR:HA	1:A:154:ARG:HG3	2.03	0.41
1:A:102:MET:CE	1:A:133:LEU:CD2	2.99	0.40
1:A:116:ASN:C	1:A:118:LEU:N	2.79	0.40
1:A:83:LYS:NZ	1:A:83:LYS:C	2.79	0.40
1:A:119:ARG:CB	1:A:119:ARG:NH1	2.84	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	163/167 (98%)	126 (77%)	20 (12%)	17 (10%)	0 0

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	37	PRO
1	A	109	THR
1	A	122	GLN
1	A	125	ARG
1	A	139	TYR
1	A	151	THR
1	A	31	HIS
1	A	39	LEU
1	A	52	ARG
1	A	136	SER
1	A	160	ALA
1	A	36	SER
1	A	93	ALA
1	A	140	ASN
1	A	108	GLU
1	A	123	GLN
1	A	149	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	135/137 (98%)	82 (61%)	53 (39%)	0 0

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLU
1	A	8	ARG

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Mol	Chain	Res	Type
1	A	13	LEU
1	A	14	ARG
1	A	16	LYS
1	A	19	LYS
1	A	26	THR
1	A	27	ILE
1	A	29	ILE
1	A	32	LEU
1	A	36	SER
1	A	38	SER
1	A	39	LEU
1	A	48	LYS
1	A	53	ASN
1	A	60	LYS
1	A	66	LEU
1	A	69	GLN
1	A	72	ASP
1	A	78	ILE
1	A	79	LEU
1	A	81	ASN
1	A	83	LYS
1	A	85	LYS
1	A	87	VAL
1	A	91	LEU
1	A	99	LEU
1	A	103	VAL
1	A	105	GLN
1	A	106	MET
1	A	108	GLU
1	A	109	THR
1	A	119	ARG
1	A	120	MET
1	A	121	LEU
1	A	122	GLN
1	A	123	GLN
1	A	124	LYS
1	A	125	ARG
1	A	126	TRP
1	A	127	ASP
1	A	131	VAL
1	A	132	ASN
1	A	133	LEU

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Mol	Chain	Res	Type
1	A	135	LYS
1	A	139	TYR
1	A	140	ASN
1	A	141	GLN
1	A	142	THR
1	A	147	LYS
1	A	151	THR
1	A	152	THR
1	A	159	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	68	ASN
1	A	81	ASN
1	A	101	ASN
1	A	132	ASN
1	A	140	ASN

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

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	165/167 (98%)	-0.42	1 (0%) 85 85	12, 41, 73, 91	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	139	TYR	3.5

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.