



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:58 PM UTC

PDB ID : 7E7A / pdb_00007e7a
Title : Crystal structure of apo ENL YEATS domain T3 mutant
Authors : Li, Y.; Li, H.
Deposited on : 2021-02-25
Resolution : 2.64 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

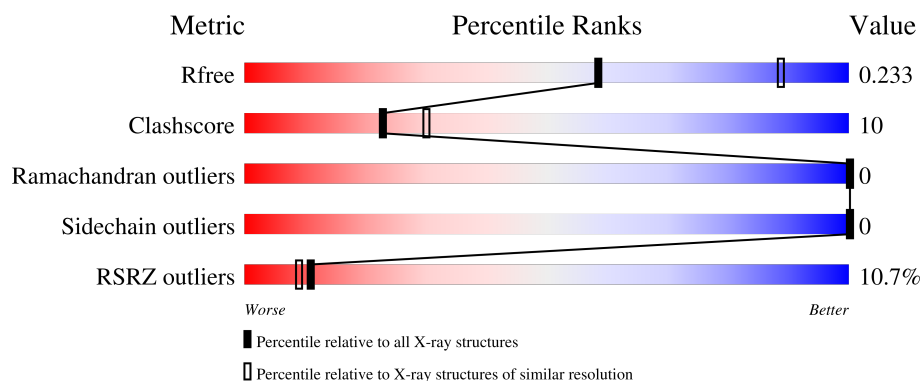
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	155	 6% 70% 17% • 12%
1	B	155	 8% 74% 15% 11%
1	C	155	 11% 70% 19% • 11%
1	D	155	 13% 66% 21% • 12%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4655 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 Protein ENL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	137	Total	C	N	O	S	0	0	0
			1142	738	202	194	8			
1	B	138	Total	C	N	O	S	0	0	0
			1150	743	203	195	9			
1	C	138	Total	C	N	O	S	0	0	0
			1150	743	203	195	9			
1	D	137	Total	C	N	O	S	0	0	0
			1142	738	202	194	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q03111
A	-1	SER	-	expression tag	UNP Q03111
A	0	HIS	-	expression tag	UNP Q03111
A	111	LYS	ASN	engineered mutation	UNP Q03111
A	?	-	PRO	deletion	UNP Q03111
A	?	-	PRO	deletion	UNP Q03111
A	147	HIS	-	expression tag	UNP Q03111
A	148	HIS	-	expression tag	UNP Q03111
A	149	HIS	-	expression tag	UNP Q03111
A	150	HIS	-	expression tag	UNP Q03111
A	151	HIS	-	expression tag	UNP Q03111
A	152	HIS	-	expression tag	UNP Q03111
B	-2	GLY	-	expression tag	UNP Q03111
B	-1	SER	-	expression tag	UNP Q03111
B	0	HIS	-	expression tag	UNP Q03111
B	111	LYS	ASN	engineered mutation	UNP Q03111
B	?	-	PRO	deletion	UNP Q03111
B	?	-	PRO	deletion	UNP Q03111
B	147	HIS	-	expression tag	UNP Q03111
B	148	HIS	-	expression tag	UNP Q03111
B	149	HIS	-	expression tag	UNP Q03111

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Chain	Residue	Modelled	Actual	Comment	Reference
B	150	HIS	-	expression tag	UNP Q03111
B	151	HIS	-	expression tag	UNP Q03111
B	152	HIS	-	expression tag	UNP Q03111
C	-2	GLY	-	expression tag	UNP Q03111
C	-1	SER	-	expression tag	UNP Q03111
C	0	HIS	-	expression tag	UNP Q03111
C	111	LYS	ASN	engineered mutation	UNP Q03111
C	?	-	PRO	deletion	UNP Q03111
C	?	-	PRO	deletion	UNP Q03111
C	147	HIS	-	expression tag	UNP Q03111
C	148	HIS	-	expression tag	UNP Q03111
C	149	HIS	-	expression tag	UNP Q03111
C	150	HIS	-	expression tag	UNP Q03111
C	151	HIS	-	expression tag	UNP Q03111
C	152	HIS	-	expression tag	UNP Q03111
D	-2	GLY	-	expression tag	UNP Q03111
D	-1	SER	-	expression tag	UNP Q03111
D	0	HIS	-	expression tag	UNP Q03111
D	111	LYS	ASN	engineered mutation	UNP Q03111
D	?	-	PRO	deletion	UNP Q03111
D	?	-	PRO	deletion	UNP Q03111
D	147	HIS	-	expression tag	UNP Q03111
D	148	HIS	-	expression tag	UNP Q03111
D	149	HIS	-	expression tag	UNP Q03111
D	150	HIS	-	expression tag	UNP Q03111
D	151	HIS	-	expression tag	UNP Q03111
D	152	HIS	-	expression tag	UNP Q03111

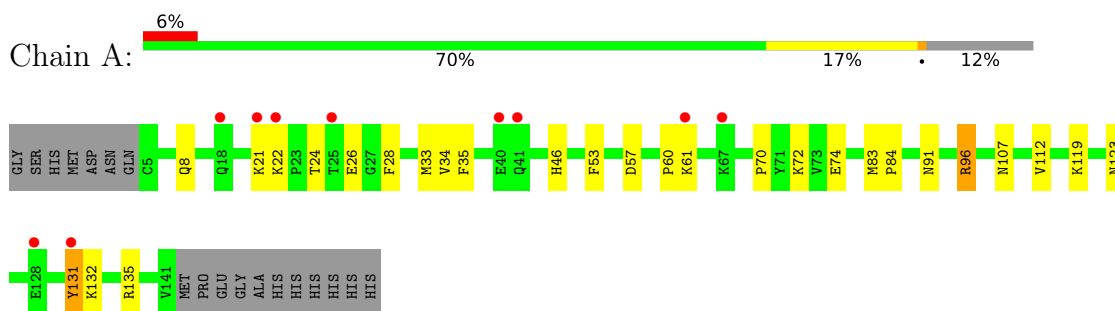
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	23	Total O 23 23	0	0
2	B	20	Total O 20 20	0	0
2	C	13	Total O 13 13	0	0
2	D	15	Total O 15 15	0	0

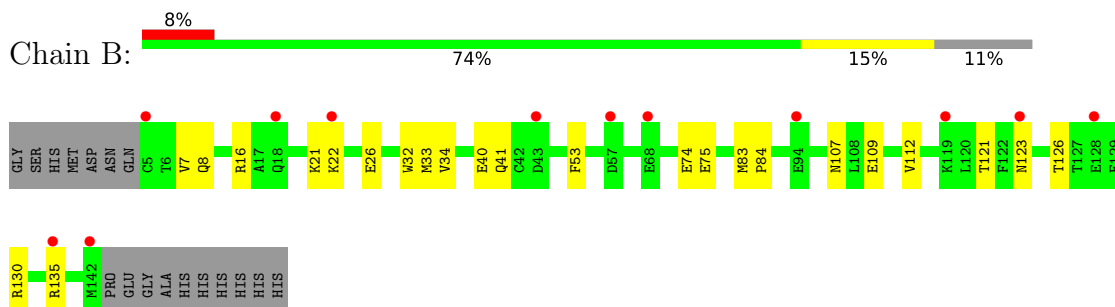
3 Residue-property plots [i](#)

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.

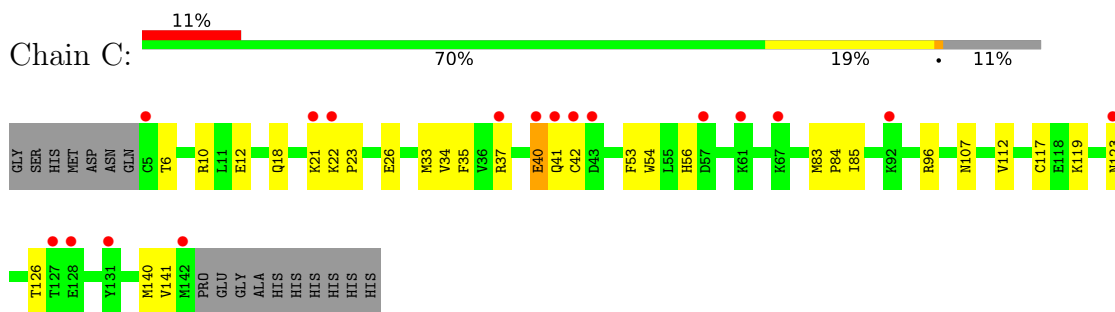
• Molecule 1: Protein ENL



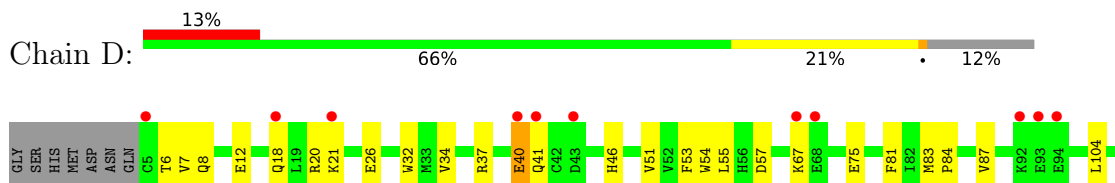
• Molecule 1: Protein ENL

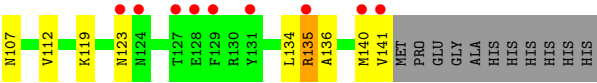


• Molecule 1: Protein ENL



• Molecule 1: Protein ENL





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.32Å 88.26Å 167.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.04 – 2.64 49.04 – 2.64	Depositor EDS
% Data completeness (in resolution range)	96.1 (49.04-2.64) 96.5 (49.04-2.64)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.13 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.213 , 0.230 0.213 , 0.233	Depositor DCC
R_{free} test set	1772 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	50.1	Xtriage
Anisotropy	0.640	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4655	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.39 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.7043e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.41	2/1175 (0.2%)	0.62	2/1585 (0.1%)
1	B	0.26	0/1183	0.56	2/1595 (0.1%)
1	C	0.28	0/1183	0.60	2/1595 (0.1%)
1	D	0.38	1/1175 (0.1%)	0.68	3/1585 (0.2%)
All	All	0.34	3/4716 (0.1%)	0.62	9/6360 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	96	ARG	NE-CZ	7.01	1.40	1.33
1	D	21	LYS	CG-CD	6.36	1.71	1.52
1	A	96	ARG	CD-NE	5.35	1.53	1.46

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	135	ARG	NE-CZ-NH2	-9.56	110.60	119.20
1	C	40	GLU	CA-C-N	7.59	136.04	121.54
1	C	40	GLU	C-N-CA	7.59	136.04	121.54
1	A	131	TYR	CB-CA-C	-7.56	96.87	110.70
1	D	135	ARG	CG-CD-NE	-7.26	96.03	112.00
1	A	131	TYR	CA-CB-CG	6.52	125.63	113.90
1	B	135	ARG	NH1-CZ-NH2	5.61	126.60	119.30
1	D	40	GLU	CA-C-N	5.08	133.96	125.02
1	D	40	GLU	C-N-CA	5.08	133.96	125.02

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	1142	0	1137	24	0
1	B	1150	0	1147	19	0
1	C	1150	0	1147	31	1
1	D	1142	0	1138	29	1
2	A	23	0	0	0	0
2	B	20	0	0	2	0
2	C	13	0	0	3	0
2	D	15	0	0	2	0
All	All	4655	0	4569	95	1

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 10.

All (95) 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:123:ASN:HD21	1:D:123:ASN:HB2	1.25	0.99
1:C:40:GLU:O	1:C:41:GLN:HG3	1.61	0.98
1:C:10:ARG:HE	1:C:119:LYS:HE3	1.27	0.96
1:D:53:PHE:HB3	1:D:83:MET:HE2	1.60	0.82
1:C:33:MET:HE3	1:C:35:PHE:HD2	1.44	0.81
1:D:18:GLN:NE2	2:D:202:HOH:O	2.17	0.77
1:A:34:VAL:HG11	1:A:83:MET:HE1	1.70	0.74
1:C:117:CYS:HB3	1:C:119:LYS:HZ1	1.54	0.73
1:A:46:HIS:CD2	1:A:135:ARG:HD2	2.23	0.73
1:B:123:ASN:ND2	1:D:123:ASN:HB2	2.02	0.72
1:B:84:PRO:HG3	1:D:84:PRO:HG3	1.71	0.72
1:D:7:VAL:HG22	1:D:140:MET:HG2	1.72	0.71
1:C:10:ARG:NE	1:C:119:LYS:HE3	2.04	0.71
1:C:41:GLN:NE2	1:C:42:CYS:SG	2.64	0.71
1:C:12:GLU:OE1	1:C:37:ARG:NH2	2.23	0.70
1:D:40:GLU:HG2	1:D:41:GLN:NE2	2.07	0.69
1:B:53:PHE:HB3	1:B:83:MET:HE3	1.74	0.69
1:C:117:CYS:HB3	1:C:119:LYS:NZ	2.08	0.68
1:C:34:VAL:HG11	1:C:83:MET:HE1	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:GLN:HG2	1:B:121:THR:HG22	1.76	0.66
1:A:84:PRO:HG3	1:C:84:PRO:HG3	1.78	0.66
1:C:119:LYS:O	2:C:201:HOH:O	2.13	0.66
1:C:107:ASN:HB2	1:C:112:VAL:HG23	1.81	0.62
1:B:126:THR:HG23	2:B:202:HOH:O	1.99	0.61
1:C:53:PHE:HB3	1:C:83:MET:HE3	1.82	0.61
1:D:135:ARG:HB3	1:D:135:ARG:CZ	2.30	0.61
1:A:123:ASN:OD1	1:C:123:ASN:ND2	2.34	0.60
1:D:20:ARG:O	2:D:201:HOH:O	2.16	0.60
1:C:33:MET:HE3	1:C:35:PHE:CD2	2.32	0.59
1:A:107:ASN:HB2	1:A:112:VAL:HG23	1.85	0.59
1:D:51:VAL:HG22	1:D:87:VAL:HG13	1.85	0.59
1:C:33:MET:CE	1:C:35:PHE:HD2	2.14	0.59
1:A:33:MET:HG3	1:A:74:GLU:HG2	1.84	0.58
1:A:21:LYS:HD3	1:A:22:LYS:H	1.68	0.58
1:D:12:GLU:OE1	1:D:37:ARG:NH2	2.33	0.57
1:D:107:ASN:HB2	1:D:112:VAL:HG23	1.85	0.57
1:C:10:ARG:HE	1:C:119:LYS:CE	2.11	0.56
1:C:18:GLN:NE2	2:C:202:HOH:O	2.16	0.56
1:C:22:LYS:HD2	1:C:23:PRO:O	2.06	0.55
1:B:21:LYS:HD3	1:C:96:ARG:HH22	1.74	0.53
1:D:12:GLU:HB2	1:D:37:ARG:HG3	1.90	0.53
1:A:131:TYR:O	1:A:132:LYS:C	2.51	0.53
1:D:46:HIS:CE1	1:D:135:ARG:NH2	2.76	0.53
1:B:109:GLU:O	2:B:201:HOH:O	2.19	0.53
1:D:54:TRP:HB2	1:D:84:PRO:HD2	1.91	0.52
1:C:33:MET:CE	1:C:35:PHE:CD2	2.92	0.52
1:B:34:VAL:HG11	1:B:83:MET:HE1	1.92	0.52
1:D:55:LEU:HD23	1:D:83:MET:HE3	1.91	0.51
1:A:131:TYR:O	1:A:135:ARG:N	2.30	0.51
1:D:6:THR:O	1:D:141:VAL:HG23	2.11	0.51
1:D:46:HIS:NE2	1:D:135:ARG:NH2	2.60	0.50
1:A:33:MET:HE3	1:A:35:PHE:HD2	1.76	0.50
1:A:53:PHE:HB3	1:A:83:MET:HE3	1.93	0.50
1:D:26:GLU:OE1	1:D:26:GLU:N	2.43	0.49
1:A:57:ASP:OD2	1:C:56:HIS:NE2	2.45	0.49
1:A:53:PHE:HD2	1:A:83:MET:CE	2.25	0.49
1:A:8:GLN:HB3	1:A:119:LYS:CE	2.42	0.49
1:B:40:GLU:O	1:B:41:GLN:HB2	2.14	0.48
1:C:53:PHE:HD2	1:C:83:MET:CE	2.26	0.48
1:D:107:ASN:CB	1:D:112:VAL:HG23	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:GLU:N	1:C:26:GLU:OE1	2.46	0.48
1:A:91:ASN:O	1:A:96:ARG:HD2	2.14	0.47
1:B:21:LYS:C	1:B:22:LYS:HD3	2.40	0.46
1:D:34:VAL:HG11	1:D:83:MET:HE1	1.98	0.46
1:A:70:PRO:HB2	1:A:72:LYS:HD2	1.98	0.46
1:D:55:LEU:CD2	1:D:83:MET:HE3	2.46	0.46
1:B:33:MET:HG3	1:B:74:GLU:HG2	1.98	0.46
1:B:32:TRP:CZ2	1:B:75:GLU:HB2	2.52	0.45
1:A:26:GLU:OE1	1:A:26:GLU:N	2.40	0.45
1:A:60:PRO:HG2	1:A:61:LYS:HE2	1.99	0.45
1:D:134:LEU:HD21	1:D:140:MET:HE2	1.98	0.45
1:D:81:PHE:CE1	1:D:104:LEU:HD23	2.53	0.44
1:A:24:THR:OG1	1:A:28:PHE:HB2	2.18	0.44
1:C:140:MET:HE2	1:C:140:MET:HB3	1.84	0.44
1:B:7:VAL:HG21	1:B:130:ARG:HG3	1.98	0.44
1:C:6:THR:HB	1:C:141:VAL:CG2	2.48	0.44
1:D:46:HIS:HB2	1:D:136:ALA:HB2	2.00	0.43
1:A:8:GLN:HB3	1:A:119:LYS:HE3	2.00	0.43
1:B:16:ARG:HA	1:B:112:VAL:O	2.19	0.43
1:B:107:ASN:HB2	1:B:112:VAL:HG23	2.01	0.43
1:C:126:THR:HG23	2:C:203:HOH:O	2.18	0.43
1:A:8:GLN:HB3	1:A:119:LYS:HE2	2.02	0.42
1:C:54:TRP:HB2	1:C:84:PRO:HD2	2.00	0.42
1:D:67:LYS:O	1:D:67:LYS:HG2	2.19	0.42
1:A:107:ASN:CB	1:A:112:VAL:HG23	2.48	0.42
1:B:26:GLU:OE2	1:B:26:GLU:N	2.41	0.42
1:D:8:GLN:CD	1:D:119:LYS:HD2	2.44	0.42
1:C:83:MET:HE2	1:C:85:ILE:HD11	2.01	0.42
1:A:46:HIS:NE2	1:A:135:ARG:HD2	2.35	0.41
1:C:21:LYS:HE2	1:C:21:LYS:HB2	1.74	0.41
1:D:57:ASP:OD1	1:D:57:ASP:N	2.47	0.41
1:B:21:LYS:CD	1:C:96:ARG:HH22	2.33	0.41
1:D:32:TRP:CZ2	1:D:75:GLU:HB2	2.55	0.40
1:B:33:MET:CG	1:B:74:GLU:HG2	2.51	0.40
1:A:135:ARG:O	1:A:135:ARG:CD	2.70	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:GLN:OE1	1:D:135:ARG:NH1[1_455]	2.11	0.09

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	135/155 (87%)	131 (97%)	4 (3%)	0	100	100
1	B	136/155 (88%)	133 (98%)	3 (2%)	0	100	100
1	C	136/155 (88%)	134 (98%)	2 (2%)	0	100	100
1	D	135/155 (87%)	134 (99%)	1 (1%)	0	100	100
All	All	542/620 (87%)	532 (98%)	10 (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	126/141 (89%)	126 (100%)	0	100	100
1	B	127/141 (90%)	127 (100%)	0	100	100
1	C	127/141 (90%)	127 (100%)	0	100	100
1	D	126/141 (89%)	126 (100%)	0	100	100
All	All	506/564 (90%)	506 (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 (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	114	HIS
1	C	41	GLN
1	C	45	GLN
1	C	123	ASN

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	137/155 (88%)	0.46	10 (7%) 21 17	32, 49, 72, 84	0
1	B	138/155 (89%)	0.49	12 (8%) 16 12	35, 50, 77, 86	0
1	C	138/155 (89%)	0.60	17 (12%) 8 7	37, 53, 80, 102	0
1	D	137/155 (88%)	0.87	20 (14%) 6 5	40, 58, 87, 103	0
All	All	550/620 (88%)	0.60	59 (10%) 11 9	32, 52, 80, 103	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	142	MET	6.5
1	A	131	TYR	6.5
1	C	41	GLN	5.8
1	D	41	GLN	5.2
1	D	141	VAL	4.0
1	B	142	MET	3.9
1	D	135	ARG	3.8
1	C	57	ASP	3.7
1	D	131	TYR	3.6
1	C	5	CYS	3.6
1	A	41	GLN	3.6
1	D	67	LYS	3.5
1	B	135	ARG	3.5
1	C	40	GLU	3.1
1	D	140	MET	3.1
1	C	123	ASN	3.1
1	D	123	ASN	3.0
1	B	5	CYS	3.0
1	D	129	PHE	2.9
1	D	92	LYS	2.9
1	D	128	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	94	GLU	2.7
1	B	68	GLU	2.7
1	C	43	ASP	2.6
1	A	22	LYS	2.6
1	B	43	ASP	2.5
1	C	37	ARG	2.5
1	C	21	LYS	2.4
1	D	127	THR	2.4
1	B	57	ASP	2.4
1	D	93	GLU	2.4
1	C	42	CYS	2.4
1	D	43	ASP	2.4
1	A	61	LYS	2.4
1	D	18	GLN	2.4
1	D	21	LYS	2.3
1	B	128	GLU	2.3
1	D	94	GLU	2.3
1	B	18	GLN	2.3
1	A	25	THR	2.3
1	B	22	LYS	2.3
1	C	92	LYS	2.3
1	A	67	LYS	2.2
1	D	40	GLU	2.2
1	D	5	CYS	2.2
1	B	123	ASN	2.2
1	A	21	LYS	2.2
1	A	40	GLU	2.2
1	C	131	TYR	2.2
1	D	124	ASN	2.2
1	C	22	LYS	2.2
1	D	68	GLU	2.1
1	B	119	LYS	2.1
1	C	67	LYS	2.1
1	C	61	LYS	2.1
1	A	128	GLU	2.0
1	C	128	GLU	2.0
1	C	127	THR	2.0
1	A	18	GLN	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.