



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:34 PM UTC

PDB ID : 7EL1 / pdb_00007el1
Title : Structure of a protein from bacteria
Authors : Liu, H.; Zhu, Y.; Huang, Z.
Deposited on : 2021-04-07
Resolution : 2.22 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

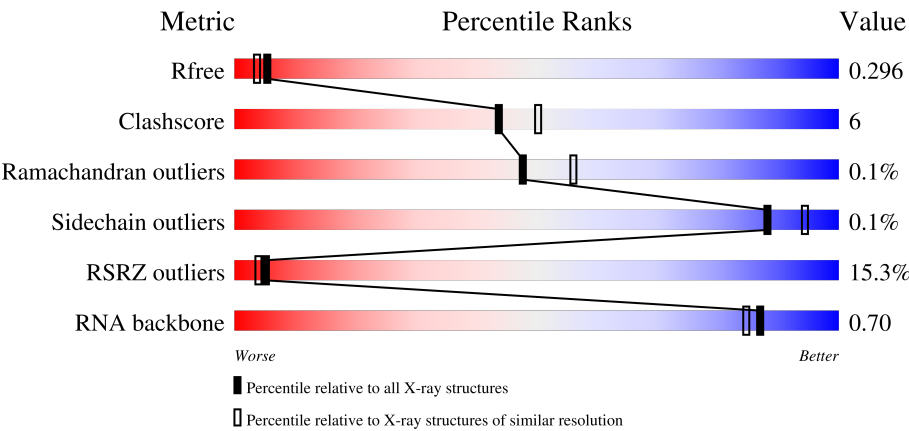
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)
RNA backbone	3983	1098 (2.54-1.90)

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	1053	17% 85% 12% .
2	B	73	% 81% 15% .
3	C	28	4% 93% 7%
4	D	8	100%

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Mol	Chain	Length	Quality of chain
5	E	100	5% 79% 18% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11584 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 CRISPR-associated endonuclease Cas9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1028	Total	C	N	O	S	0	0	1
			8369	5320	1460	1577	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	580	ALA	ASN	engineered mutation	UNP J7RUA5
A	946	ALA	CYS	engineered mutation	UNP J7RUA5

- Molecule 2 is a RNA chain called RNA (73-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	73	Total	C	N	O	P	0	0	0
			1545	689	279	504	73			

- Molecule 3 is a DNA chain called DNA (28-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	28	Total	C	N	O	P	0	0	0
			561	270	99	165	27			

- Molecule 4 is a DNA chain called DNA (5'-D(*TP*TP*GP*AP*AP*TP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	8	Total	C	N	O	P	0	0	0
			164	80	31	46	7			

- Molecule 5 is a protein called 100AA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	98	Total 819	C 526	N 134	O 157	S 2	0	0	0

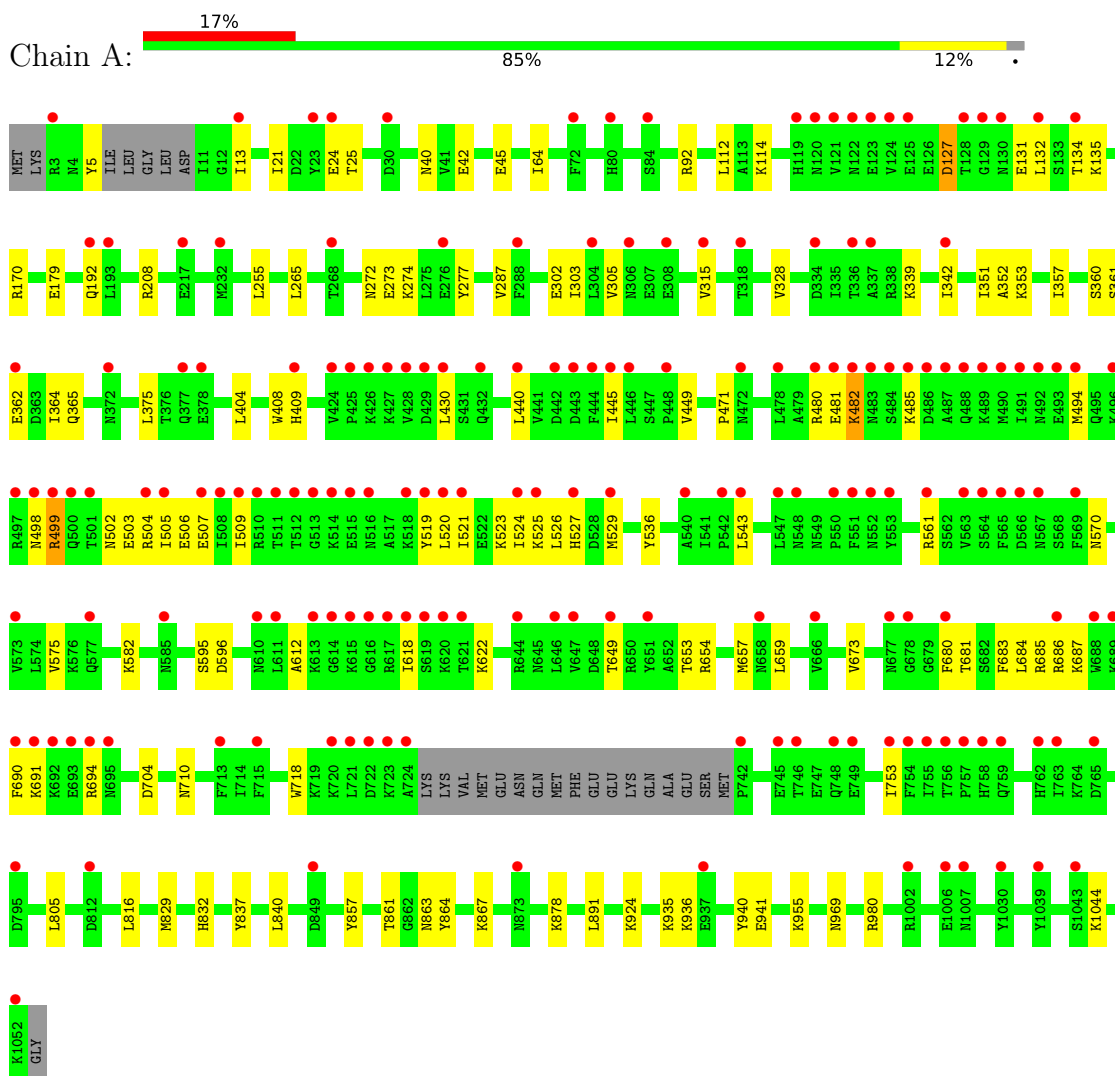
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	73	Total 73	O 73	0	0
6	B	35	Total 35	O 35	0	0
6	C	5	Total 5	O 5	0	0
6	D	9	Total 9	O 9	0	0
6	E	4	Total 4	O 4	0	0

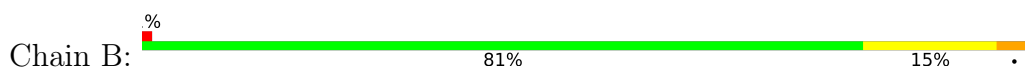
3 Residue-property plots

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: CRISPR-associated endonuclease Cas9

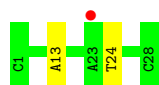
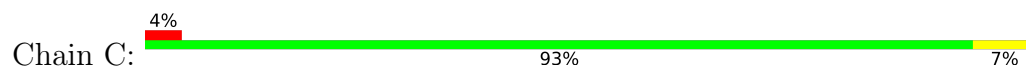


• Molecule 2: RNA (73-MER)





- Molecule 3: DNA (28-MER)



- Molecule 4: DNA (5'-D(*TP*TP*GP*AP*AP*TP*AP*G)-3')



There are no outlier residues recorded for this chain.

- Molecule 5: 100AA



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	330.58Å 105.22Å 68.55Å 90.00° 92.71° 90.00°	Depositor
Resolution (Å)	50.00 – 2.22 50.00 – 2.22	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-2.22) 99.8 (50.00-2.22)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 2.22Å)	Xtriage
Refinement program	PHENIX v1.11	Depositor
R, R_{free}	0.213 , 0.242 (Not available) , 0.296	Depositor DCC
R_{free} test set	5675 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	54.9	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.011 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11584	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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

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	0.43	0/8523	0.64	3/11480 (0.0%)
2	B	0.37	0/1729	0.56	1/2692 (0.0%)
3	C	0.48	0/627	0.72	0/963
4	D	0.59	0/184	0.81	0/283
5	E	0.56	1/836 (0.1%)	0.69	2/1120 (0.2%)
All	All	0.44	1/11899 (0.0%)	0.64	6/16538 (0.0%)

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 (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	30	LYS	C-N	11.07	1.40	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	46	LYS	CD-CE-NZ	-9.12	82.72	111.90
1	A	499	ARG	N-CA-C	8.59	120.34	110.97
1	A	878	LYS	CD-CE-NZ	-7.61	87.55	111.90
2	B	57	G	O5'-P-OP2	-5.16	92.53	108.00
1	A	208	ARG	NE-CZ-NH2	5.11	123.80	119.20
5	E	46	LYS	CB-CG-CD	-5.06	99.67	111.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	480	ARG	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	8369	0	8382	120	0
2	B	1545	0	773	6	0
3	C	561	0	317	2	0
4	D	164	0	93	0	0
5	E	819	0	801	12	0
6	A	73	0	0	2	0
6	B	35	0	0	0	0
6	C	5	0	0	0	0
6	D	9	0	0	0	0
6	E	4	0	0	1	0
All	All	11584	0	10366	134	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 (134) 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:520:LEU:O	1:A:524:ILE:HD12	1.55	1.06
1:A:657:MET:HE1	1:A:673:VAL:HG12	1.32	1.05
1:A:134:THR:HG22	1:A:170:ARG:HH11	1.28	0.96
1:A:520:LEU:HD13	1:A:524:ILE:HD11	1.48	0.95
1:A:521:ILE:HA	1:A:524:ILE:CD1	1.97	0.94
1:A:504:ARG:HA	1:A:504:ARG:HH21	1.36	0.88
1:A:861:THR:HG23	1:A:863:ASN:H	1.40	0.87
1:A:504:ARG:HH22	1:A:507:GLU:HB3	1.40	0.86
1:A:521:ILE:HA	1:A:524:ILE:HD12	1.57	0.85
1:A:657:MET:CE	1:A:673:VAL:HG12	2.11	0.81
1:A:504:ARG:HH22	1:A:507:GLU:CB	1.95	0.80
1:A:690:PHE:HB3	1:A:694:ARG:HH22	1.48	0.79
1:A:499:ARG:O	1:A:503:GLU:HG2	1.83	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:ILE:HD11	1:A:305:VAL:HG22	1.67	0.76
1:A:134:THR:HG22	1:A:170:ARG:NH1	2.01	0.75
1:A:521:ILE:HA	1:A:524:ILE:HD13	1.67	0.75
5:E:30:LYS:HD3	5:E:60:VAL:HG12	1.69	0.75
1:A:502:ASN:HA	1:A:505:ILE:HG13	1.70	0.73
1:A:272:ASN:HD22	1:A:273:GLU:N	1.89	0.71
1:A:92:ARG:HD2	1:A:179:GLU:OE2	1.89	0.71
1:A:504:ARG:HA	1:A:504:ARG:NH2	2.05	0.71
1:A:520:LEU:C	1:A:524:ILE:HD12	2.16	0.70
1:A:595:SER:HB3	5:E:8:SER:HB3	1.77	0.66
1:A:690:PHE:HB3	1:A:694:ARG:NH2	2.10	0.66
1:A:521:ILE:CA	1:A:524:ILE:HD12	2.25	0.66
1:A:653:THR:O	1:A:657:MET:HG2	1.96	0.64
5:E:47:ILE:HD11	5:E:82:GLY:HA2	1.80	0.64
1:A:303:ILE:CD1	1:A:305:VAL:HG22	2.28	0.64
1:A:360:SER:O	1:A:364:ILE:HG12	1.98	0.64
1:A:328:VAL:HG11	1:A:352:ALA:HB2	1.80	0.64
1:A:64:ILE:HD11	1:A:114:LYS:HB2	1.79	0.63
1:A:685:ARG:HH11	1:A:694:ARG:NH1	1.97	0.62
1:A:520:LEU:O	1:A:524:ILE:CD1	2.40	0.62
1:A:857:TYR:O	1:A:861:THR:HG22	1.98	0.62
1:A:40:ASN:HD21	1:A:485:LYS:HB2	1.64	0.62
1:A:690:PHE:C	1:A:694:ARG:HH22	2.09	0.61
1:A:867:LYS:NZ	2:B:45:C:O2	2.33	0.60
1:A:685:ARG:NH2	1:A:704:ASP:OD1	2.34	0.59
1:A:131:GLU:OE1	1:A:132:LEU:HD23	2.02	0.59
1:A:504:ARG:NH2	1:A:507:GLU:HB3	2.14	0.59
1:A:525:LYS:O	1:A:529:MET:HG2	2.03	0.59
1:A:612:ALA:HA	1:A:618:ILE:HG22	1.83	0.58
1:A:494:MET:HE3	1:A:505:ILE:CD1	2.32	0.58
1:A:339:LYS:HA	1:A:342:ILE:HG22	1.84	0.58
1:A:303:ILE:C	1:A:303:ILE:HD12	2.29	0.58
1:A:92:ARG:HD3	1:A:112:LEU:HG	1.87	0.57
1:A:685:ARG:HB3	1:A:690:PHE:HB2	1.86	0.57
1:A:680:PHE:O	1:A:683:PHE:HB3	2.04	0.56
1:A:936:LYS:HD2	1:A:941:GLU:OE1	2.05	0.56
1:A:13:ILE:HD12	1:A:449:VAL:HG11	1.88	0.56
1:A:265:LEU:O	1:A:274:LYS:NZ	2.28	0.55
1:A:561:ARG:HB2	1:A:686:ARG:NH1	2.22	0.55
1:A:494:MET:HE1	1:A:509:ILE:HG12	1.90	0.54
1:A:829:MET:HE3	1:A:837:TYR:HA	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:685:ARG:HH11	1:A:694:ARG:CZ	2.21	0.54
1:A:526:LEU:HD23	1:A:529:MET:SD	2.48	0.54
1:A:362:GLU:H	1:A:362:GLU:CD	2.15	0.54
1:A:867:LYS:HD2	2:B:46:U:H1'	1.89	0.54
5:E:66:ARG:HD2	5:E:68:ASP:OD2	2.07	0.54
5:E:70:ASN:OD1	5:E:72:ILE:N	2.41	0.53
1:A:277:TYR:CD1	1:A:409:HIS:CD2	2.96	0.53
1:A:649:THR:O	1:A:654:ARG:NH1	2.42	0.52
1:A:691:LYS:O	1:A:694:ARG:CZ	2.57	0.52
1:A:681:THR:O	1:A:685:ARG:HG3	2.08	0.52
1:A:519:TYR:O	1:A:523:LYS:HG3	2.11	0.51
1:A:502:ASN:HA	1:A:505:ILE:CG1	2.38	0.51
1:A:864:TYR:CE2	6:A:1138:HOH:O	2.62	0.50
1:A:287:VAL:HG23	1:A:302:GLU:HG3	1.93	0.50
5:E:39:LYS:O	5:E:43:GLU:HG3	2.12	0.50
1:A:718:TRP:CD1	1:A:753:ILE:HD11	2.46	0.50
1:A:21:ILE:HD13	1:A:710:ASN:OD1	2.11	0.49
1:A:303:ILE:HD13	1:A:305:VAL:HG13	1.94	0.49
1:A:683:PHE:CZ	1:A:687:LYS:HE2	2.47	0.49
1:A:498:ASN:O	1:A:502:ASN:HB2	2.12	0.49
1:A:505:ILE:HD12	1:A:506:GLU:N	2.28	0.49
1:A:502:ASN:HD22	1:A:505:ILE:HD11	1.78	0.49
1:A:805:LEU:HD12	1:A:816:LEU:HD23	1.93	0.49
1:A:525:LYS:NZ	1:A:529:MET:HE3	2.27	0.49
2:B:23:U:H2'	2:B:24:U:C6	2.47	0.49
1:A:351:ILE:HD11	1:A:375:LEU:HD11	1.94	0.48
1:A:277:TYR:CD1	1:A:409:HIS:HD2	2.32	0.48
2:B:12:G:H2'	2:B:13:C:C6	2.48	0.48
1:A:127:ASP:OD1	1:A:127:ASP:N	2.46	0.48
1:A:690:PHE:CB	1:A:694:ARG:HH22	2.22	0.48
1:A:680:PHE:CZ	1:A:684:LEU:HD11	2.48	0.48
1:A:829:MET:O	1:A:832:HIS:O	2.32	0.48
5:E:60:VAL:HG22	6:E:203:HOH:O	2.13	0.48
1:A:303:ILE:CD1	1:A:303:ILE:C	2.87	0.47
1:A:5:TYR:CE2	1:A:471:PRO:HB3	2.49	0.47
1:A:273:GLU:HG3	1:A:274:LYS:HG2	1.97	0.47
1:A:40:ASN:ND2	1:A:485:LYS:HB2	2.28	0.47
1:A:24:GLU:HG2	1:A:25:THR:H	1.79	0.46
1:A:1044:LYS:NZ	6:A:1101:HOH:O	2.33	0.46
1:A:924:LYS:NZ	1:A:955:LYS:O	2.44	0.46
1:A:829:MET:HE1	1:A:840:LEU:HD23	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:430:LEU:HD21	1:A:440:LEU:HD13	1.98	0.46
1:A:255:LEU:HD23	1:A:404:LEU:HD21	1.97	0.46
1:A:361:SER:O	1:A:365:GLN:HG3	2.16	0.46
1:A:64:ILE:HD11	1:A:114:LYS:CB	2.44	0.45
1:A:570:ASN:HA	1:A:622:LYS:HD3	1.99	0.45
1:A:969:ASN:OD1	1:A:980:ARG:HD2	2.17	0.44
1:A:502:ASN:ND2	1:A:505:ILE:HD11	2.32	0.44
1:A:536:TYR:CD2	1:A:575:VAL:HG11	2.52	0.44
1:A:891:LEU:HD21	2:B:56:A:H4'	2.00	0.44
1:A:42:GLU:HA	1:A:45:GLU:HB2	2.01	0.43
1:A:440:LEU:HD22	1:A:659:LEU:HD13	2.00	0.43
1:A:683:PHE:HD2	1:A:684:LEU:HD12	1.84	0.43
1:A:272:ASN:ND2	1:A:274:LYS:H	2.16	0.43
1:A:482:LYS:HD3	1:A:482:LYS:HA	1.86	0.43
1:A:582:LYS:HE3	1:A:596:ASP:O	2.18	0.43
1:A:481:GLU:O	1:A:482:LYS:HG2	2.18	0.43
1:A:494:MET:HE3	1:A:505:ILE:HD13	1.99	0.43
1:A:408:TRP:HB3	1:A:409:HIS:CD2	2.54	0.42
1:A:582:LYS:HB2	1:A:582:LYS:HE2	1.74	0.42
5:E:70:ASN:OD1	5:E:70:ASN:C	2.62	0.42
1:A:445:ILE:HD12	1:A:445:ILE:HA	1.84	0.42
2:B:62:A:H4'	2:B:63:A:O5'	2.19	0.42
5:E:5:LYS:HG2	5:E:23:VAL:HG22	2.00	0.42
1:A:192:GLN:CD	1:A:192:GLN:H	2.28	0.42
1:A:526:LEU:HA	1:A:529:MET:CG	2.49	0.42
5:E:17:TYR:CD1	5:E:42:ILE:HG13	2.54	0.42
1:A:40:ASN:HD21	1:A:485:LYS:CB	2.31	0.41
1:A:527:HIS:CE1	1:A:543:LEU:HG	2.56	0.41
1:A:502:ASN:HD22	1:A:505:ILE:CG1	2.33	0.41
1:A:135:LYS:HG3	3:C:13:DA:H5''	2.02	0.41
1:A:829:MET:HE1	1:A:840:LEU:CD2	2.50	0.41
1:A:24:GLU:HG2	1:A:25:THR:N	2.35	0.41
1:A:342:ILE:HG21	1:A:342:ILE:HD13	1.85	0.41
5:E:89:LEU:O	5:E:93:LYS:NZ	2.52	0.41
1:A:353:LYS:O	1:A:357:ILE:HG12	2.21	0.41
1:A:829:MET:HE2	1:A:829:MET:HB3	1.87	0.41
5:E:69:LYS:HB3	5:E:69:LYS:HE2	1.79	0.41
1:A:315:VAL:HG12	3:C:24:DT:H2''	2.03	0.40
1:A:935:LYS:HB2	1:A:940:TYR:CE1	2.56	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	1024/1053 (97%)	987 (96%)	36 (4%)	1 (0%)	48	56
5	E	96/100 (96%)	94 (98%)	2 (2%)	0	100	100
All	All	1120/1153 (97%)	1081 (96%)	38 (3%)	1 (0%)	48	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	482	LYS

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	856/970 (88%)	855 (100%)	1 (0%)	88	94
5	E	91/94 (97%)	91 (100%)	0	100	100
All	All	947/1064 (89%)	946 (100%)	1 (0%)	88	94

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	127	ASP

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	40	ASN
1	A	272	ASN
1	A	325	ASN
1	A	377	GLN
1	A	557	HIS
1	A	570	ASN
1	A	585	ASN
1	A	700	HIS
1	A	832	HIS
1	A	835	GLN
1	A	885	ASN
1	A	888	ASN
1	A	1037	ASN
5	E	84	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	71/73 (97%)	7 (9%)	1 (1%)

All (7) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	38	A
2	B	39	C
2	B	55	A
2	B	57	G
2	B	62	A
2	B	63	A
2	B	65	G

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	62	A

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 [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	1026/1053 (97%)	1.07	182 (17%) 4 3	36, 65, 103, 146	0
2	B	73/73 (100%)	-0.04	1 (1%) 73 72	42, 54, 112, 120	0
3	C	28/28 (100%)	0.05	1 (3%) 46 43	46, 56, 82, 83	0
4	D	8/8 (100%)	-0.53	0 100 100	41, 45, 50, 52	0
5	E	98/100 (98%)	0.81	5 (5%) 33 30	53, 69, 81, 82	0
All	All	1233/1262 (97%)	0.95	189 (15%) 5 4	36, 65, 101, 146	0

All (189) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	124	VAL	7.2
1	A	614	GLY	6.8
1	A	446	LEU	6.7
1	A	122	ASN	6.5
1	A	481	GLU	6.4
1	A	647	VAL	6.2
1	A	491	ILE	6.0
1	A	505	ILE	5.5
1	A	129	GLY	5.4
1	A	440	LEU	5.3
1	A	499	ARG	5.3
1	A	615	LYS	5.2
1	A	551	PHE	5.1
1	A	649	THR	5.0
1	A	724	ALA	4.9
1	A	489	LYS	4.8
1	A	432	GLN	4.7
1	A	498	ASN	4.7
1	A	130	ASN	4.5
1	A	618	ILE	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	757	PRO	4.4
1	A	585	ASN	4.3
1	A	753	ILE	4.2
1	A	428	VAL	4.1
1	A	123	GLU	4.1
1	A	485	LYS	4.1
1	A	132	LEU	4.1
1	A	617	ARG	4.1
1	A	565	PHE	4.1
1	A	511	THR	4.0
1	A	513	GLY	4.0
1	A	490	MET	3.9
1	A	482	LYS	3.9
1	A	715	PHE	3.9
1	A	484	SER	3.9
1	A	680	PHE	3.8
1	A	755	ILE	3.8
1	A	426	LYS	3.8
1	A	425	PRO	3.8
1	A	508	ILE	3.8
1	A	430	LEU	3.7
1	A	496	LYS	3.7
1	A	692	LYS	3.7
1	A	742	PRO	3.7
1	A	409	HIS	3.7
1	A	125	GLU	3.7
1	A	483	ASN	3.6
1	A	514	LYS	3.6
1	A	494	MET	3.6
1	A	443	ASP	3.6
1	A	763	ILE	3.6
1	A	492	ASN	3.5
1	A	445	ILE	3.5
1	A	487	ALA	3.5
1	A	543	LEU	3.5
1	A	444	PHE	3.4
1	A	759	GLN	3.4
1	A	691	LYS	3.3
1	A	529	MET	3.3
1	A	569	PHE	3.3
1	A	518	LYS	3.2
1	A	765	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	849	ASP	3.2
1	A	548	ASN	3.2
1	A	758	HIS	3.2
1	A	480	ARG	3.1
1	A	694	ARG	3.1
1	A	80	HIS	3.1
1	A	542	PRO	3.1
1	A	509	ILE	3.1
1	A	507	GLU	3.1
1	A	693	GLU	3.1
1	A	722	ASP	3.1
1	A	713	PHE	3.0
1	A	510	ARG	3.0
1	A	620	LYS	3.0
1	A	72	PHE	3.0
1	A	666	VAL	3.0
1	A	521	ILE	3.0
1	A	561	ARG	3.0
1	A	651	TYR	3.0
1	A	308	GLU	2.9
1	A	550	PRO	2.9
1	A	23	TYR	2.9
1	A	678	GLY	2.9
1	A	232	MET	2.9
1	A	342	ILE	2.9
1	A	493	GLU	2.8
1	A	567	ASN	2.8
1	A	695	ASN	2.8
1	A	24	GLU	2.8
1	A	424	VAL	2.8
1	A	519	TYR	2.8
1	A	553	TYR	2.8
1	A	427	LYS	2.7
1	A	690	PHE	2.7
1	A	754	PHE	2.7
1	A	795	ASP	2.7
1	A	1043	SER	2.7
1	A	488	GLN	2.7
1	A	1030	TYR	2.7
1	A	512	THR	2.7
1	A	577	GLN	2.6
1	A	217	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	611	LEU	2.6
1	A	276	GLU	2.6
1	A	128	THR	2.6
1	A	478	LEU	2.6
1	A	504	ARG	2.6
1	A	644	ARG	2.6
1	A	1002	ARG	2.6
1	A	336	THR	2.6
1	A	448	PRO	2.6
1	A	646	LEU	2.5
1	A	540	ALA	2.5
1	A	658	ASN	2.5
1	A	304	LEU	2.5
1	A	721	LEU	2.5
1	A	486	ASP	2.5
1	A	525	LYS	2.5
1	A	563	VAL	2.5
1	A	686	ARG	2.5
1	A	1007	ASN	2.5
1	A	762	HIS	2.5
1	A	566	ASP	2.4
2	B	35	G	2.4
1	A	3	ARG	2.4
1	A	748	GLN	2.4
3	C	23	DA	2.4
1	A	192	GLN	2.4
1	A	745	GLU	2.4
1	A	1006	GLU	2.4
1	A	1039	TYR	2.3
1	A	616	GLY	2.3
5	E	93	LYS	2.3
1	A	756	THR	2.3
1	A	564	SER	2.3
1	A	120	ASN	2.3
1	A	516	ASN	2.3
1	A	121	VAL	2.3
1	A	315	VAL	2.3
1	A	547	LEU	2.3
1	A	613	LYS	2.3
1	A	621	THR	2.3
1	A	746	THR	2.3
1	A	193	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	1052	LYS	2.3
1	A	610	ASN	2.3
5	E	17	TYR	2.3
1	A	134	THR	2.3
1	A	119	HIS	2.3
1	A	268	THR	2.2
1	A	318	THR	2.2
1	A	689	LYS	2.2
1	A	472	ASN	2.2
1	A	677	ASN	2.2
1	A	937	GLU	2.2
1	A	720	LYS	2.2
1	A	723	LYS	2.2
1	A	812	ASP	2.2
1	A	337	ALA	2.2
1	A	524	ILE	2.1
1	A	378	GLU	2.1
1	A	501	THR	2.1
1	A	306	ASN	2.1
1	A	362	GLU	2.1
1	A	429	ASP	2.1
1	A	515	GLU	2.1
1	A	688	TRP	2.1
1	A	497	ARG	2.1
1	A	619	SER	2.1
1	A	13	ILE	2.1
1	A	372	ASN	2.1
1	A	377	GLN	2.1
5	E	71	GLY	2.1
5	E	39	LYS	2.1
1	A	442	ASP	2.1
1	A	500	GLN	2.1
5	E	89	LEU	2.1
1	A	873	ASN	2.1
1	A	288	PHE	2.1
1	A	84	SER	2.0
1	A	749	GLU	2.0
1	A	527	HIS	2.0
1	A	520	LEU	2.0
1	A	573	VAL	2.0
1	A	30	ASP	2.0
1	A	334	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	552	ASN	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.