



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2026 – 06:06 PM EDT

PDB ID : 7ENR / pdb_00007enr
Title : Crystal structure of cas and anti-cas protein complex
Authors : Wang, Y.; Li, X.
Deposited on : 2021-04-19
Resolution : 4.21 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

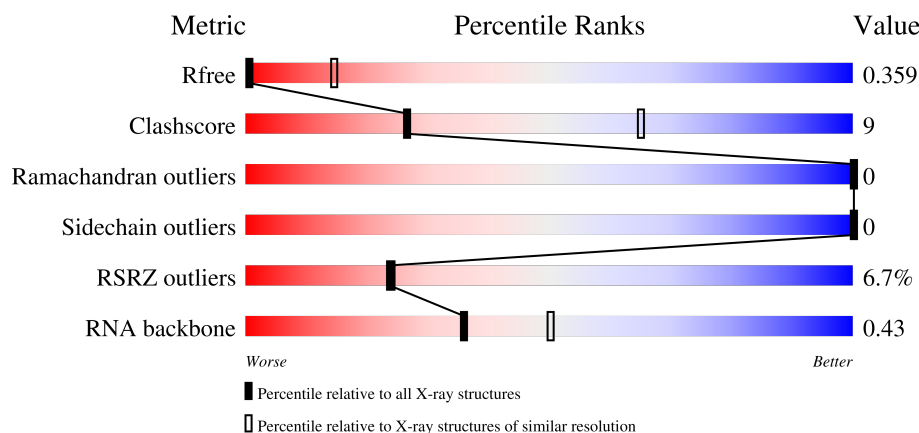
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.21 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	1004 (4.52-3.88)
Clashscore	190562	1019 (4.50-3.90)
Ramachandran outliers	187476	1020 (4.56-3.84)
Sidechain outliers	187428	1006 (4.56-3.84)
RSRZ outliers	180081	1001 (4.52-3.88)
RNA backbone	3983	1026 (5.04-3.10)

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	 6% 81% 13%
2	B	98	 5% 34% 38% 21% 4%
3	C	100	 14% 88% 9%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10498 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	1048	Total	C	N	O	S	0	4	0
			7930	5021	1381	1515	13			

- Molecule 2 is a RNA chain called sgRNA (98-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	92	Total	C	N	O	P	0	0	0
			1929	863	331	643	92			

- Molecule 3 is a protein called AcrIIA14.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	97	Total	C	N	O	0	0	0
			639	399	112	128			

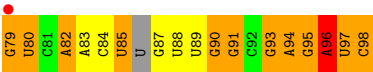
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

Chain A:

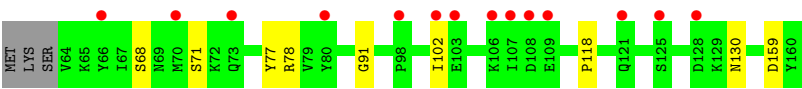
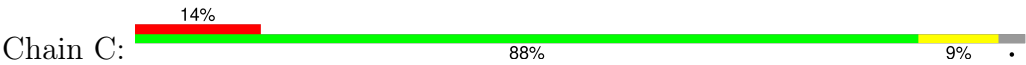
Chain B:

Category	Percentage
Red	5%
Green	34%
Yellow	38%
Orange	21%
Grey	6%

Sequence: G, G, U, U, C, G6, G7, A9, U10, U11, U17, U18, A19, U23, U24, U25, A26, G27, U33, G34, A38, C39, A40, G41, A42, A43, U46, A47, C48, U49, A50, A51, A52, A53, C54, A55, A56, G57, A61, A62, A63, U64, G65, G66, C67, U72, U73, A74, U75, C76, U77, U78



● Molecule 3: AcrIIA14



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	173.80Å 173.80Å 185.89Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.99 – 4.21 38.99 – 4.21	Depositor EDS
% Data completeness (in resolution range)	95.7 (38.99-4.21) 95.5 (38.99-4.21)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 4.28Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.260 , 0.307 0.264 , 0.359	Depositor DCC
R_{free} test set	342 reflections (2.25%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.640	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 1.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.038 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	10498	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% 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	0.14	0/8076	0.34	0/10988
2	B	0.27	0/2152	0.51	1/3343 (0.0%)
3	C	0.11	0/651	0.30	0/897
All	All	0.17	0/10879	0.38	1/15228 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	96	A	C4-N9-C1'	-5.46	109.93	126.30

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	7930	0	7302	132	1
2	B	1929	0	973	53	1
3	C	639	0	451	7	0
All	All	10498	0	8726	177	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 9.

All (177) 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:971:LEU:HB2	1:A:1030:TYR:O	1.69	0.91
2:B:78:C:N3	2:B:95:G:N2	2.30	0.79
1:A:719:LYS:HE3	1:A:742:PRO:HG3	1.64	0.77
2:B:39:C:H2'	2:B:40:A:H8	1.49	0.77
1:A:526:LEU:HD12	1:A:574:LEU:HB3	1.71	0.71
2:B:49:U:H2'	2:B:50:A:H8	1.56	0.70
1:A:105:PHE:HE1	1:A:183:LEU:HD21	1.57	0.68
1:A:608:ILE:HG12	1:A:626:LEU:HD12	1.76	0.67
1:A:275:LEU:O	1:A:411:ASN:ND2	2.29	0.66
1:A:526:LEU:CD1	1:A:574:LEU:H	2.09	0.64
1:A:602:GLU:OE2	3:C:78:ARG:NH2	2.30	0.64
1:A:968:ASN:OD1	1:A:969:ASN:ND2	2.32	0.63
1:A:608:ILE:HG21	1:A:627:LEU:HD21	1.81	0.62
2:B:52:A:H2'	2:B:53:A:H8	1.63	0.62
1:A:674:LYS:HD3	1:A:728:MET:HE2	1.80	0.62
2:B:49:U:H2'	2:B:50:A:C8	2.35	0.62
2:B:42:A:H2'	2:B:43:A:C8	2.35	0.61
1:A:234:MET:HE1	1:A:247:VAL:HG13	1.82	0.61
1:A:689:LYS:O	1:A:764:LYS:NZ	2.33	0.61
1:A:897:TYR:HE1	2:B:53:A:H1'	1.65	0.61
1:A:526:LEU:HD11	1:A:574:LEU:H	1.66	0.61
1:A:897:TYR:HE2	1:A:904:VAL:HG11	1.66	0.61
1:A:582:LYS:O	1:A:586:ARG:NH2	2.31	0.60
1:A:55:ARG:NH2	2:B:67:C:OP1	2.35	0.59
1:A:579:GLU:OE1	1:A:591:TYR:OH	2.21	0.59
1:A:974:ILE:HG22	1:A:1027:ILE:HG12	1.85	0.59
2:B:50:A:H2'	2:B:51:A:C8	2.39	0.58
1:A:857:TYR:O	1:A:861:THR:OG1	2.20	0.58
2:B:48:C:H2'	2:B:49:U:H6	1.69	0.57
1:A:625:TYR:OH	1:A:645:ASN:ND2	2.36	0.57
2:B:50:A:H2'	2:B:51:A:H8	1.69	0.56
1:A:354:ILE:HG21	1:A:368:LEU:HG	1.87	0.56
1:A:1010:ASP:OD2	1:A:1012:ARG:NE	2.39	0.56
2:B:10:U:O2'	2:B:11:U:OP1	2.22	0.56
1:A:66:ARG:HH12	2:B:64:U:H4'	1.69	0.56
1:A:791:THR:OG1	1:A:903:LYS:O	2.23	0.56
1:A:526:LEU:HD12	1:A:574:LEU:CB	2.34	0.56
1:A:558:ILE:HD11	1:A:573:VAL:HG13	1.86	0.56
1:A:5:TYR:HE1	1:A:471:PRO:HB3	1.71	0.56
1:A:529:MET:HE1	2:B:6:G:H1	1.71	0.56
1:A:969:ASN:OD1	1:A:980:ARG:NH2	2.38	0.56
1:A:5:TYR:CE1	1:A:471:PRO:HB3	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:93:G:O2'	2:B:94:A:O4'	2.24	0.55
1:A:536:TYR:CD1	1:A:575:VAL:HG11	2.42	0.55
1:A:720:LYS:HA	1:A:742:PRO:HB2	1.88	0.55
2:B:52:A:H2'	2:B:53:A:C8	2.40	0.55
1:A:92:ARG:HD2	1:A:112:LEU:HG	1.89	0.54
1:A:887:LEU:HD11	1:A:905:VAL:HG21	1.89	0.54
2:B:85:U:H5	2:B:87:G:C6	2.26	0.54
1:A:558:ILE:HG23	1:A:588:PRO:HG2	1.88	0.54
1:A:940:TYR:HB2	1:A:1051:LYS:HB3	1.90	0.54
1:A:134:THR:HG23	1:A:170:ARG:HD2	1.88	0.54
1:A:203:ASP:OD1	1:A:207:THR:OG1	2.25	0.54
1:A:5:TYR:HE2	1:A:468:TYR:HB3	1.73	0.54
1:A:424:VAL:HG13	1:A:662:SER:HB3	1.90	0.53
1:A:118:VAL:H	1:A:172:LYS:HA	1.73	0.53
1:A:177:VAL:HG21	1:A:206:GLU:HB3	1.90	0.53
1:A:89:TYR:O	1:A:93:VAL:HG13	2.09	0.53
1:A:247:VAL:HG23	1:A:250:ALA:HB2	1.90	0.52
1:A:881:LYS:NZ	2:B:46:U:OP2	2.38	0.52
1:A:9:LEU:HB2	1:A:476:ILE:HG12	1.91	0.52
1:A:676:ILE:HD11	1:A:680:PHE:HD2	1.73	0.52
1:A:280:LYS:O	1:A:284:ILE:HG12	2.09	0.52
3:C:71:SER:HA	3:C:159:ASP:HB3	1.90	0.52
1:A:505:ILE:HD11	1:A:524:ILE:HG21	1.90	0.52
1:A:526:LEU:C	1:A:526:LEU:HD13	2.35	0.52
1:A:233:LEU:HD23	1:A:357:ILE:HG22	1.90	0.52
2:B:38:A:C5	2:B:39:C:H1'	2.45	0.52
3:C:68:SER:O	3:C:130:ASN:ND2	2.43	0.52
1:A:642:ILE:HG23	1:A:646:LEU:HD23	1.92	0.52
1:A:558:ILE:HG12	1:A:626:LEU:HD22	1.91	0.52
1:A:942:VAL:HB	1:A:1049:ILE:HG23	1.92	0.52
1:A:827:LEU:HD23	1:A:882:TYR:HB3	1.93	0.51
1:A:244:LEU:HD12	1:A:245:ARG:O	2.11	0.51
1:A:132:LEU:HB3	1:A:170:ARG:HH22	1.75	0.50
1:A:282:GLN:NE2	1:A:302:GLU:OE1	2.45	0.50
2:B:48:C:H2'	2:B:49:U:C6	2.47	0.50
1:A:644:ARG:NH1	1:A:645:ASN:OD1	2.42	0.50
1:A:708:ILE:O	1:A:711:ALA:HB3	2.12	0.50
1:A:739:GLU:OE1	1:A:742:PRO:HD2	2.11	0.50
1:A:456:GLN:NE2	2:B:73:U:OP2	2.45	0.49
3:C:77:TYR:HB2	3:C:102:ILE:HG13	1.94	0.49
2:B:42:A:H2'	2:B:43:A:H8	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:SER:HB2	1:A:397:LEU:HG	1.94	0.49
1:A:314:ARG:NH1	1:A:322:GLU:O	2.45	0.49
1:A:749:GLU:O	1:A:753:ILE:HG12	2.12	0.49
1:A:913:ARG:NH2	1:A:967:TYR:OH	2.46	0.49
2:B:78:C:H42	2:B:94:A:H61	1.60	0.48
1:A:536:TYR:HE1	1:A:573:VAL:HG21	1.77	0.48
2:B:39:C:H2'	2:B:40:A:C8	2.39	0.48
1:A:183:LEU:O	1:A:187:GLN:HB2	2.14	0.48
1:A:601:TYR:HB3	3:C:78:ARG:NH2	2.29	0.48
1:A:88:PRO:O	1:A:92:ARG:HB2	2.14	0.47
1:A:561:ARG:NH1	1:A:567:ASN:OD1	2.47	0.47
1:A:208:ARG:HB2	2:B:17:U:H5'	1.96	0.47
1:A:973:LYS:HD3	1:A:1030:TYR:CE2	2.49	0.47
1:A:744:ILE:HG13	1:A:749:GLU:HB3	1.96	0.47
2:B:77:U:C4	2:B:96:A:N1	2.82	0.47
1:A:5:TYR:OH	1:A:469:GLY:O	2.28	0.47
1:A:676:ILE:HD11	1:A:680:PHE:CD2	2.49	0.47
1:A:820:ILE:HD12	1:A:844:MET:HG3	1.96	0.47
2:B:82:A:C2	2:B:83:A:C2	3.02	0.47
2:B:77:U:O4	2:B:95:G:C6	2.68	0.47
1:A:561:ARG:HD2	1:A:750:TYR:CE1	2.49	0.46
1:A:92:ARG:HG3	1:A:108:ALA:HB1	1.98	0.46
1:A:454:PHE:CZ	1:A:652:ALA:HB1	2.50	0.46
2:B:23:U:H2'	2:B:24:U:H6	1.80	0.46
1:A:879:LYS:NZ	2:B:23:U:OP1	2.49	0.46
1:A:9:LEU:HD11	1:A:461:ILE:HG13	1.97	0.46
2:B:23:U:H2'	2:B:24:U:C6	2.51	0.46
1:A:808:LEU:HD22	1:A:816:LEU:HD22	1.97	0.46
2:B:72:U:O2'	2:B:73:U:OP1	2.31	0.46
1:A:358:TYR:CE2	1:A:367:GLU:HG3	2.51	0.45
1:A:977:GLU:OE1	1:A:979:TYR:OH	2.29	0.45
2:B:41:G:C2	2:B:42:A:N7	2.84	0.45
2:B:33:U:C4	2:B:41:G:N2	2.85	0.45
1:A:568:SER:OG	2:B:6:G:OP1	2.34	0.45
2:B:24:U:H2'	2:B:25:U:C6	2.52	0.45
1:A:394:ASN:HD21	1:A:445:ILE:HD11	1.82	0.44
1:A:184:LEU:HD13	1:A:198:ILE:HG23	1.99	0.44
1:A:897:TYR:CE2	1:A:904:VAL:HG11	2.51	0.44
1:A:39:ALA:O	1:A:449:VAL:HG22	2.18	0.44
1:A:683:PHE:CE1	1:A:687:LYS:HG3	2.51	0.44
2:B:10:U:HO2'	2:B:11:U:P	2.40	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:721:LEU:HD11	1:A:754:PHE:HE1	1.83	0.44
2:B:97:U:O2'	2:B:98:C:OP1	2.32	0.44
1:A:50:LYS:HD3	1:A:50:LYS:HA	1.84	0.43
1:A:527:HIS:ND1	1:A:532:GLY:HA2	2.33	0.43
1:A:649:THR:OG1	2:B:9:A:N1	2.49	0.43
1:A:851:LYS:HD3	1:A:851:LYS:HA	1.90	0.43
1:A:93:VAL:HG12	1:A:151:ALA:HB1	1.99	0.43
1:A:962:PHE:CG	1:A:1047:PRO:HD3	2.54	0.43
1:A:259:LEU:HB3	1:A:407:LEU:HD22	1.98	0.43
1:A:1013:PRO:HA	1:A:1014:PRO:HD3	1.90	0.43
1:A:284:ILE:HA	1:A:288:PHE:HB2	2.01	0.43
1:A:292:LYS:HA	1:A:330:HIS:CD2	2.54	0.43
1:A:820:ILE:HD12	1:A:844:MET:HE3	2.01	0.43
1:A:962:PHE:CE1	1:A:1047:PRO:HB3	2.53	0.43
1:A:5:TYR:CE2	1:A:468:TYR:HB3	2.53	0.43
1:A:115:ARG:HG2	2:B:19:A:P	2.59	0.43
1:A:118:VAL:HA	1:A:208:ARG:HH22	1.84	0.43
1:A:354:ILE:O	1:A:358:TYR:HB2	2.18	0.43
1:A:954:LYS:HB2	1:A:956:ILE:HG13	1.99	0.43
2:B:85:U:C5	2:B:87:G:C6	3.06	0.43
2:B:49:U:C2	2:B:50:A:C8	3.06	0.42
2:B:61:A:H8	2:B:61:A:OP2	2.02	0.42
2:B:82:A:C2	2:B:83:A:N1	2.87	0.42
1:A:63:ARG:HD2	1:A:204:LEU:O	2.19	0.42
1:A:454:PHE:HZ	1:A:652:ALA:HB1	1.85	0.42
1:A:783:LEU:CD2	1:A:930:ASN:HB3	2.50	0.42
1:A:16:VAL:HG23	1:A:36:PHE:CE1	2.54	0.42
1:A:16:VAL:HG23	1:A:36:PHE:HE1	1.84	0.42
1:A:34:ARG:CZ	1:A:1034:ILE:HG13	2.49	0.42
1:A:600:SER:OG	3:C:159:ASP:OD2	2.32	0.42
1:A:840:LEU:HD11	1:A:854:LEU:HD11	2.02	0.42
2:B:77:U:O4	2:B:96:A:N1	2.53	0.42
2:B:18:U:H2'	2:B:19:A:C8	2.55	0.42
1:A:982:ILE:HD11	1:A:995:ASN:HB2	2.02	0.42
1:A:820:ILE:O	1:A:824:PRO:HG3	2.20	0.41
1:A:410:THR:HG21	1:A:420:ARG:HH22	1.86	0.41
1:A:919:ASP:HB2	1:A:960:ALA:HB2	2.01	0.41
1:A:92:ARG:HE	1:A:151:ALA:HB2	1.85	0.41
2:B:82:A:C2	2:B:90:G:C6	3.08	0.41
3:C:91:GLY:O	3:C:118:PRO:HA	2.19	0.41
1:A:566:ASP:O	1:A:571:ASN:ND2	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:559:ILE:HG21	1:A:641:PHE:CE2	2.56	0.41
1:A:879:LYS:HZ3	2:B:23:U:P	2.44	0.41
2:B:74:A:H4'	2:B:75:U:OP1	2.20	0.41
2:B:74:A:P	2:B:74:A:H3'	2.61	0.41
1:A:521:ILE:O	1:A:522:GLU:C	2.63	0.40
2:B:26:A:H2'	2:B:27:G:H8	1.86	0.40
1:A:1045:LYS:O	1:A:1046:HIS:HB2	2.21	0.40
2:B:90:G:C6	2:B:91:G:N7	2.89	0.40
1:A:805:LEU:O	1:A:879:LYS:HA	2.21	0.40
1:A:657:MET:HE1	1:A:673:VAL:HG12	2.03	0.40
2:B:79:G:H2'	2:B:80:U:C6	2.57	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:A:317:SER:OG	2:B:82:A:OP1[5_555]	2.19	0.01

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	1048/1053 (100%)	990 (94%)	58 (6%)	0	100	100
3	C	95/100 (95%)	93 (98%)	2 (2%)	0	100	100
All	All	1143/1153 (99%)	1083 (95%)	60 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	768/972 (79%)	768 (100%)	0	100	100
3	C	44/94 (47%)	44 (100%)	0	100	100
All	All	812/1066 (76%)	812 (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 (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	462	ASN
1	A	633	ASN
1	A	667	ASN
1	A	710	ASN
1	A	773	HIS
1	A	1025	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	88/98 (89%)	31 (35%)	4 (4%)

All (31) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	6	G
2	B	10	U
2	B	11	U
2	B	34	G
2	B	39	C
2	B	47	A
2	B	55	A
2	B	56	A

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Mol	Chain	Res	Type
2	B	57	G
2	B	61	A
2	B	63	A
2	B	65	G
2	B	73	U
2	B	74	A
2	B	75	U
2	B	78	C
2	B	79	G
2	B	80	U
2	B	82	A
2	B	84	C
2	B	85	U
2	B	88	U
2	B	89	U
2	B	90	G
2	B	91	G
2	B	93	G
2	B	94	A
2	B	95	G
2	B	96	A
2	B	97	U
2	B	98	C

All (4) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	10	U
2	B	72	U
2	B	74	A
2	B	97	U

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

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

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	1048/1053 (99%)	0.86	64 (6%)	27 27	6, 39, 64, 86	4 (0%)
2	B	92/98 (93%)	0.68	5 (5%)	31 29	8, 46, 103, 135	0
3	C	97/100 (97%)	1.22	14 (14%)	6 10	50, 76, 104, 124	0
All	All	1237/1251 (98%)	0.88	83 (6%)	24 24	6, 41, 79, 135	4 (0%)

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	536	TYR	5.3
2	B	74	A	4.2
3	C	106	LYS	4.0
1	A	85	GLY	4.0
1	A	585	ASN	3.7
1	A	1033	ASP	3.3
1	A	428	VAL	3.3
1	A	130	ASN	3.2
1	A	844	MET	3.2
1	A	422	LYS	3.2
2	B	34	G	3.2
1	A	251	TYR	3.2
1	A	425	PRO	3.1
1	A	388	GLY	2.9
1	A	308	GLU	2.9
1	A	537	SER	2.9
1	A	111	HIS	2.8
1	A	743	GLU	2.8
1	A	107	ALA	2.8
1	A	629	GLU	2.8
1	A	816	LEU	2.8
1	A	829	MET	2.7
1	A	351	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	379	GLU	2.7
1	A	565	PHE	2.6
3	C	98	PRO	2.6
1	A	516	ASN	2.5
1	A	342	ILE	2.5
1	A	692	LYS	2.5
1	A	754	PHE	2.5
3	C	102	ILE	2.5
1	A	554	GLU	2.5
1	A	600	SER	2.4
3	C	70	MET	2.4
1	A	255	LEU	2.4
3	C	125	SER	2.4
1	A	1021	ALA	2.4
2	B	62	A	2.4
1	A	840	LEU	2.4
1	A	599	ILE	2.3
1	A	635	PHE	2.3
3	C	109	GLU	2.3
3	C	66	TYR	2.3
1	A	505	ILE	2.3
1	A	680	PHE	2.3
2	B	79	G	2.2
1	A	123	GLU	2.2
1	A	573	VAL	2.2
1	A	877	ILE	2.2
1	A	131	GLU	2.2
1	A	392	THR	2.2
1	A	383	ILE	2.2
1	A	961	GLU	2.2
3	C	128	ASP	2.2
3	C	107	ILE	2.2
1	A	341	ILE	2.1
1	A	682	SER	2.1
3	C	73	GLN	2.1
1	A	220	PRO	2.1
2	B	6	G	2.1
3	C	80	TYR	2.1
1	A	500	GLN	2.1
1	A	219	SER	2.1
1	A	297	LYS	2.1
3	C	103	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	445	ILE	2.1
1	A	738	ALA	2.1
1	A	128	THR	2.1
1	A	214	GLY	2.1
1	A	404	LEU	2.1
1	A	880	ILE	2.1
1	A	48	ARG	2.0
1	A	109	LEU	2.0
3	C	108	ASP	2.0
1	A	288	PHE	2.0
1	A	535	LEU	2.0
1	A	909	LEU	2.0
1	A	129	GLY	2.0
1	A	238	THR	2.0
3	C	121	GLN	2.0
1	A	865	LEU	2.0
1	A	95	GLY	2.0
1	A	22	ASP	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.