



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 02:37 PM UTC

PDB ID : 6M0V / pdb_00006m0v
Title : Crsytal structure of streptococcus thermophilus Cas9 in complex with the GGAA PAM
Authors : Zhang, Y.; Zhang, H.; Xu, X.; Wang, Y.; Chen, W.; Wang, Y.; Wu, Z.; Tang, N.; Wang, Y.; Zhao, S.; Gan, J.; Ji, Q.
Deposited on : 2020-02-22
Resolution : 3.00 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

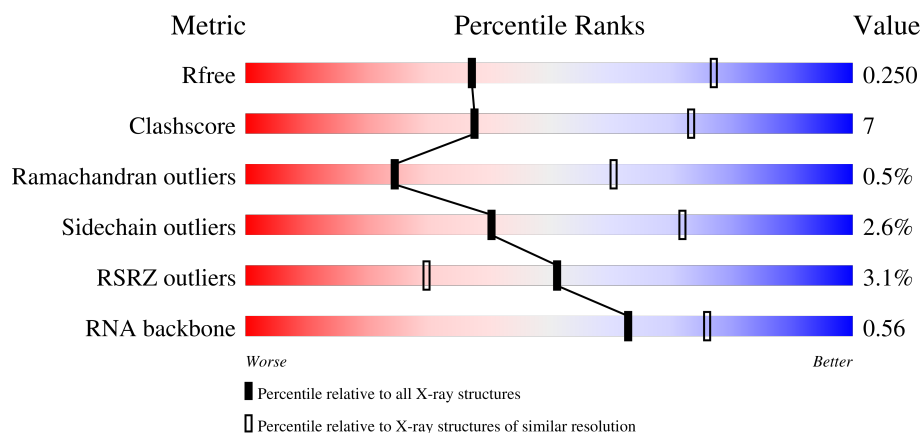
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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)
RNA backbone	3983	1109 (3.20-2.80)

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	B	71	3% 70% 25% .
2	C	28	64% 36%
3	D	8	88% 12%
4	A	1122	3% 73% 17% . 8%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 10482 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 RNA chain called RNA (71-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	71	Total	C	N	O	P	0	0	0
			1500	670	271	488	71			

- Molecule 2 is a DNA chain called DNA (28-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	28	Total	C	N	O	P	0	0	0
			561	272	91	171	27			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	8	Total	C	N	O	P	0	0	0
			166	79	38	42	7			

- Molecule 4 is a protein called CRISPR-associated endonuclease Cas9 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	A	1030	Total	C	N	O	S	0	0	0
			8235	5215	1433	1570	17			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q03LF7
A	1	GLY	-	expression tag	UNP Q03LF7
A	599	ALA	HIS	engineered mutation	UNP Q03LF7

- Molecule 5 is BARIUM ION (CCD ID: BA) (formula: Ba).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	4	Total 4	Ba 4	0	0
5	C	1	Total 1	Ba 1	0	0
5	A	5	Total 5	Ba 5	0	0

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total 1	Mg 1	0	0

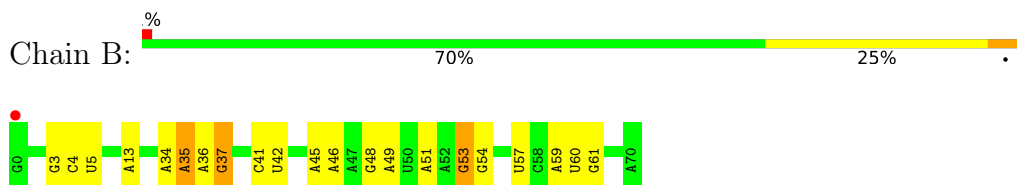
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	6	Total 6	O 6	0	0
7	C	1	Total 1	O 1	0	0
7	A	2	Total 2	O 2	0	0

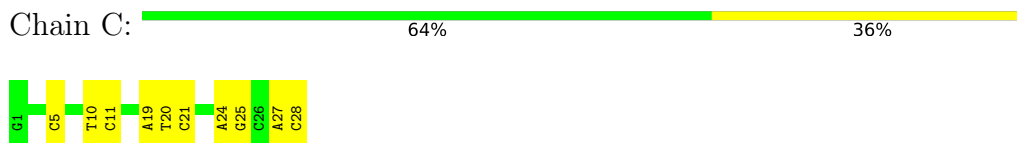
3 Residue-property plots [i](#)

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.

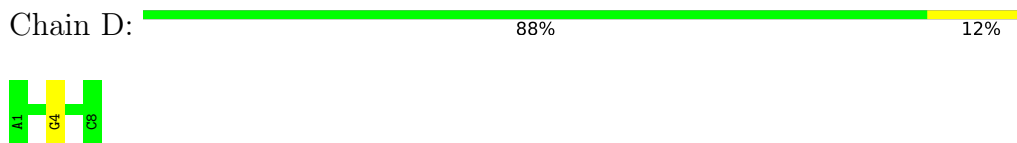
- Molecule 1: RNA (71-MER)



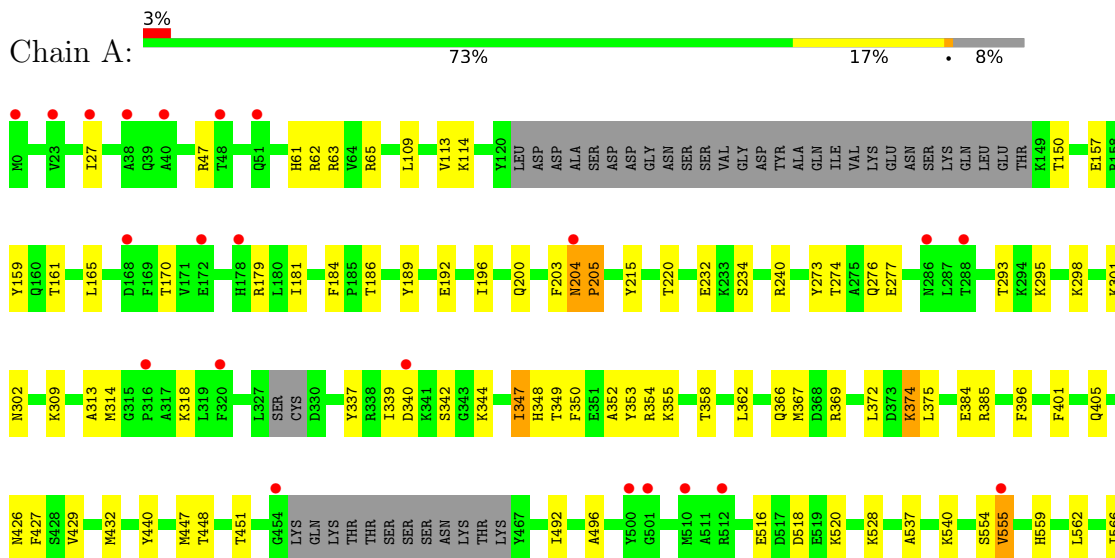
- Molecule 2: DNA (28-MER)

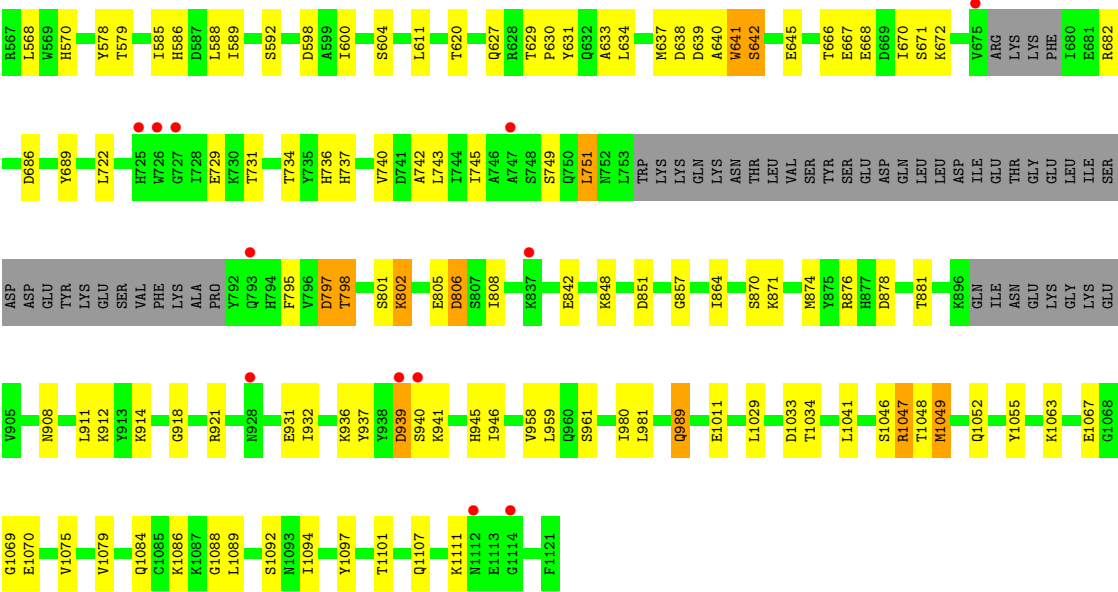


- Molecule 3: DNA (5'-D(*AP*AP*GP*GP*AP*AP*GP*C)-3')



- Molecule 4: CRISPR-associated endonuclease Cas9 1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	321.84Å 75.08Å 70.11Å 90.00° 93.21° 90.00°	Depositor
Resolution (Å)	48.82 – 3.00 48.82 – 3.00	Depositor EDS
% Data completeness (in resolution range)	93.2 (48.82-3.00) 93.2 (48.82-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.213 , 0.257 0.217 , 0.250	Depositor DCC
R_{free} test set	1592 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	57.3	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10482	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BA, MG

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	B	0.25	0/1679	0.36	0/2615
2	C	0.36	0/625	0.52	0/961
3	D	0.69	0/188	0.78	0/289
4	A	0.51	2/8378 (0.0%)	0.54	10/11282 (0.1%)
All	All	0.47	2/10870 (0.0%)	0.52	10/15147 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1046	SER	C-O	-6.61	1.18	1.24
4	A	1070	GLU	CA-C	-5.32	1.46	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1092	SER	N-CA-C	8.19	120.20	111.28
4	A	939	ASP	N-CA-C	6.84	122.02	112.93
4	A	1033	ASP	N-CA-C	-6.59	100.72	110.48
4	A	642	SER	N-CA-C	-6.30	99.75	109.52
4	A	797	ASP	CA-C-N	6.10	129.58	120.31
4	A	797	ASP	C-N-CA	6.10	129.58	120.31
4	A	797	ASP	O-C-N	6.07	128.55	122.12
4	A	751	LEU	N-CA-C	5.79	117.67	111.36
4	A	729	GLU	N-CA-C	-5.32	98.72	108.02
4	A	1070	GLU	N-CA-C	-5.03	103.50	110.35

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	B	1500	0	751	12	0
2	C	561	0	321	9	0
3	D	166	0	90	1	0
4	A	8235	0	8116	127	0
5	A	5	0	0	0	0
5	B	4	0	0	0	0
5	C	1	0	0	0	0
6	A	1	0	0	0	0
7	A	2	0	0	0	0
7	B	6	0	0	0	0
7	C	1	0	0	0	0
All	All	10482	0	9278	137	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 7.

All (137) 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)
4:A:314:MET:HG3	4:A:349:THR:HG23	1.58	0.85
4:A:937:TYR:HE1	4:A:939:ASP:HB2	1.39	0.85
4:A:937:TYR:CE1	4:A:939:ASP:HB2	2.13	0.84
4:A:159:TYR:HB2	4:A:165:LEU:HD21	1.63	0.80
4:A:27:ILE:HD13	4:A:743:LEU:HD13	1.60	0.80
4:A:722:LEU:HD23	4:A:722:LEU:O	1.84	0.77
4:A:347:ILE:HG13	4:A:348:HIS:H	1.52	0.73
4:A:981:LEU:HA	4:A:1011:GLU:HG2	1.72	0.72
4:A:347:ILE:HG13	4:A:348:HIS:N	2.05	0.71
4:A:314:MET:CG	4:A:349:THR:HG23	2.22	0.69
4:A:870:SER:O	4:A:876:ARG:NH2	2.25	0.69
4:A:734:THR:HG23	4:A:1063:LYS:HB3	1.78	0.66
2:C:24:DA:N6	2:C:25:DG:O6	2.30	0.65
2:C:28:DC:H5'	4:A:448:THR:HG22	1.79	0.64
4:A:1075:VAL:HG11	4:A:1094:ILE:HD13	1.81	0.62
4:A:309:LYS:HG2	4:A:353:TYR:CE2	2.34	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:10:DT:H2'	2:C:11:DC:C6	2.37	0.60
4:A:298:LYS:O	4:A:302:ASN:ND2	2.35	0.60
4:A:1047:ARG:NH1	4:A:1052:GLN:O	2.34	0.59
4:A:1075:VAL:HG11	4:A:1094:ILE:CD1	2.32	0.59
1:B:13:A:OP2	4:A:47:ARG:NH2	2.35	0.59
4:A:349:THR:O	4:A:349:THR:HG22	2.02	0.59
4:A:1055:TYR:HA	4:A:1088:GLY:HA2	1.84	0.59
4:A:232:GLU:OE2	4:A:232:GLU:HA	2.03	0.58
4:A:731:THR:HG21	4:A:736:HIS:HD2	1.69	0.56
4:A:639:ASP:O	4:A:640:ALA:HB2	2.05	0.56
4:A:634:LEU:HA	4:A:637:MET:HE3	1.87	0.55
1:B:45:A:H2'	1:B:46:A:C8	2.42	0.55
4:A:585:ILE:HA	4:A:588:LEU:HB3	1.88	0.55
4:A:374:LYS:HB3	4:A:396:PHE:CZ	2.42	0.55
2:C:27:DA:H2'	2:C:28:DC:C6	2.43	0.54
4:A:749:SER:C	4:A:751:LEU:H	2.16	0.54
4:A:806:ASP:OD1	4:A:806:ASP:N	2.40	0.54
4:A:313:ALA:O	4:A:318:LYS:NZ	2.35	0.54
4:A:1029:LEU:O	4:A:1041:LEU:HD12	2.08	0.54
4:A:864:ILE:HG23	4:A:871:LYS:HD3	1.90	0.53
4:A:980:ILE:HD11	4:A:1089:LEU:HB2	1.90	0.53
4:A:642:SER:OG	4:A:645:GLU:HG3	2.07	0.53
4:A:62:ARG:HB3	4:A:113:VAL:HG13	1.91	0.53
4:A:842:GLU:HG2	4:A:941:LYS:HD3	1.91	0.52
4:A:340:ASP:N	4:A:344:LYS:O	2.34	0.52
4:A:851:ASP:O	4:A:857:GLY:HA3	2.10	0.52
1:B:36:A:C5	1:B:37:G:H1'	2.45	0.51
4:A:722:LEU:HD23	4:A:722:LEU:C	2.34	0.51
4:A:349:THR:HG22	4:A:354:ARG:NH1	2.26	0.50
1:B:48:G:OP2	4:A:114:LYS:NZ	2.45	0.50
4:A:204:ASN:N	4:A:205:PRO:HD3	2.27	0.50
4:A:737:HIS:HA	4:A:740:VAL:HG22	1.92	0.50
4:A:181:ILE:HD11	4:A:620:THR:HA	1.94	0.50
4:A:367:MET:HE1	4:A:375:LEU:HD12	1.92	0.50
4:A:921:ARG:NH2	4:A:931:GLU:OE1	2.45	0.49
4:A:562:LEU:HD11	4:A:589:ILE:HG12	1.94	0.49
4:A:731:THR:HG21	4:A:736:HIS:CD2	2.46	0.49
4:A:342:SER:HA	4:A:528:LYS:HD2	1.94	0.49
1:B:60:U:O4	4:A:240:ARG:NH2	2.45	0.49
4:A:682:ARG:NE	4:A:686:ASP:OD2	2.31	0.49
4:A:274:THR:HG21	4:A:429:VAL:HG13	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:568:LEU:HD21	4:A:611:LEU:HD22	1.95	0.48
4:A:914:LYS:HA	4:A:918:GLY:O	2.12	0.48
4:A:633:ALA:O	4:A:637:MET:HG3	2.13	0.48
1:B:49:A:OP2	4:A:63:ARG:NH2	2.43	0.47
4:A:203:PHE:C	4:A:205:PRO:HD3	2.38	0.47
2:C:19:DA:H4'	4:A:689:TYR:CZ	2.49	0.47
4:A:604:SER:HB2	4:A:627:GLN:HE21	1.79	0.47
4:A:537:ALA:HB1	4:A:566:ILE:HG22	1.97	0.47
4:A:742:ALA:HA	4:A:745:ILE:HD12	1.97	0.47
3:D:4:DG:O6	4:A:1086:LYS:NZ	2.45	0.47
4:A:384:GLU:HG2	4:A:385:ARG:H	1.80	0.47
4:A:795:PHE:O	4:A:798:THR:HB	2.15	0.47
4:A:555:VAL:HG22	4:A:589:ILE:CG2	2.45	0.46
1:B:3:G:H2'	1:B:4:C:C6	2.51	0.46
4:A:61:HIS:CD2	4:A:65:ARG:HD3	2.50	0.46
2:C:20:DT:H2'	2:C:21:DC:C6	2.50	0.46
4:A:277:GLU:OE1	4:A:369:ARG:NH1	2.49	0.46
4:A:1011:GLU:CD	4:A:1047:ARG:NH2	2.74	0.46
4:A:276:GLN:CD	4:A:352:ALA:HB3	2.40	0.45
4:A:630:PRO:O	4:A:634:LEU:HB2	2.16	0.45
1:B:53:G:O2'	4:A:989:GLN:HG3	2.16	0.45
4:A:355:LYS:O	4:A:358:THR:OG1	2.31	0.45
4:A:447:MET:HE2	4:A:451:THR:HG23	1.97	0.45
4:A:629:THR:CG2	4:A:670:ILE:HD11	2.46	0.45
1:B:42:U:O2'	4:A:931:GLU:O	2.34	0.45
4:A:492:ILE:O	4:A:496:ALA:N	2.48	0.45
4:A:170:THR:HG21	4:A:638:ASP:OD2	2.17	0.45
4:A:375:LEU:HD23	4:A:375:LEU:HA	1.75	0.45
4:A:578:TYR:O	4:A:641:TRP:HB2	2.17	0.45
4:A:633:ALA:HB1	4:A:637:MET:HE2	1.98	0.45
4:A:874:MET:O	4:A:878:ASP:N	2.49	0.45
4:A:337:TYR:O	4:A:337:TYR:CD1	2.70	0.45
4:A:540:LYS:HD3	4:A:570:HIS:ND1	2.32	0.45
4:A:946:ILE:HB	4:A:959:LEU:HB2	1.99	0.44
2:C:19:DA:H2'	2:C:20:DT:C6	2.52	0.44
4:A:559:HIS:HE1	4:A:592:SER:CB	2.31	0.44
4:A:1069:GLY:H	4:A:1079:VAL:CG2	2.31	0.44
4:A:631:TYR:CD2	4:A:667:GLU:HA	2.53	0.44
4:A:908:ASN:HB3	4:A:911:LEU:HB3	1.99	0.44
4:A:109:LEU:HB3	4:A:215:TYR:CE1	2.53	0.44
4:A:170:THR:OG1	4:A:179:ARG:NE	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:181:ILE:HG13	4:A:620:THR:HG22	2.00	0.44
1:B:4:C:H2'	1:B:5:U:O4'	2.18	0.43
4:A:579:THR:HA	4:A:640:ALA:HA	1.99	0.43
4:A:666:THR:C	4:A:668:GLU:H	2.26	0.43
4:A:945:HIS:HB2	4:A:958:VAL:HB	2.01	0.43
4:A:349:THR:O	4:A:350:PHE:HB2	2.18	0.43
4:A:555:VAL:HG22	4:A:589:ILE:HG23	2.00	0.43
4:A:1049:MET:HE3	4:A:1052:GLN:HG3	2.01	0.43
4:A:276:GLN:OE1	4:A:352:ALA:HB3	2.18	0.43
1:B:41:C:O2'	4:A:874:MET:HA	2.20	0.42
1:B:35:A:OP2	1:B:35:A:H8	2.02	0.42
4:A:722:LEU:C	4:A:722:LEU:CD2	2.92	0.42
2:C:5:DC:N4	4:A:1084:GLN:OE1	2.52	0.42
4:A:427:PHE:HB2	4:A:432:MET:HE3	2.00	0.42
4:A:736:HIS:CD2	4:A:805:GLU:OE1	2.72	0.42
4:A:362:LEU:HD21	4:A:405:GLN:HB3	2.01	0.42
4:A:1097:TYR:CE1	4:A:1111:LYS:HD2	2.55	0.42
4:A:1047:ARG:NH1	4:A:1049:MET:O	2.53	0.42
4:A:196:ILE:O	4:A:200:GLN:HG2	2.20	0.42
4:A:562:LEU:HD22	4:A:566:ILE:HG13	2.01	0.42
4:A:802:LYS:HE3	4:A:802:LYS:HB3	1.90	0.42
4:A:1011:GLU:CD	4:A:1047:ARG:HH22	2.28	0.42
4:A:1069:GLY:H	4:A:1079:VAL:HG23	1.85	0.42
4:A:273:TYR:HB2	4:A:372:LEU:HD23	2.02	0.41
4:A:585:ILE:HD12	4:A:586:HIS:N	2.35	0.41
4:A:157:GLU:O	4:A:161:THR:HG23	2.20	0.41
4:A:516:GLU:O	4:A:520:LYS:N	2.30	0.41
4:A:150:THR:HB	4:A:192:GLU:HB2	2.01	0.41
4:A:301:LYS:NZ	4:A:440:TYR:O	2.49	0.41
4:A:881:THR:HG21	4:A:932:ILE:HD11	2.02	0.41
4:A:848:LYS:HA	4:A:936:LYS:HA	2.02	0.41
2:C:24:DA:H2''	4:A:339:ILE:HG22	2.03	0.41
4:A:184:PHE:HB2	4:A:189:TYR:CE2	2.56	0.41
4:A:293:THR:C	4:A:295:LYS:H	2.28	0.41
4:A:912:LYS:HD3	4:A:912:LYS:HA	1.86	0.41
4:A:186:THR:HG22	4:A:220:THR:HG22	2.03	0.41
4:A:396:PHE:CG	4:A:401:PHE:HE2	2.39	0.41
4:A:518:ASP:OD1	4:A:518:ASP:N	2.54	0.40
4:A:1101:THR:HG22	4:A:1107:GLN:HA	2.04	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
4	A	1016/1122 (91%)	948 (93%)	63 (6%)	5 (0%)	24 60

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	A	204	ASN
4	A	600	ILE
4	A	205	PRO
4	A	1049	MET
4	A	961	SER

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
4	A	874/1000 (87%)	851 (97%)	23 (3%)	40 72

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	A	234	SER
4	A	347	ILE
4	A	366	GLN
4	A	374	LYS
4	A	426	ASN
4	A	554	SER

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Mol	Chain	Res	Type
4	A	555	VAL
4	A	598	ASP
4	A	641	TRP
4	A	671	SER
4	A	672	LYS
4	A	797	ASP
4	A	798	THR
4	A	801	SER
4	A	802	LYS
4	A	806	ASP
4	A	808	ILE
4	A	940	SER
4	A	989	GLN
4	A	1034	THR
4	A	1047	ARG
4	A	1048	THR
4	A	1067	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
4	A	31	ASN
4	A	43	ASN
4	A	85	ASN
4	A	561	GLN
4	A	590	ASN
4	A	593	ASN
4	A	623	GLN
4	A	627	GLN
4	A	736	HIS
4	A	880	GLN
4	A	892	ASN
4	A	944	ASN
4	A	989	GLN
4	A	1039	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	B	69/71 (97%)	9 (13%)	0

All (9) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	B	34	A
1	B	35	A
1	B	37	G
1	B	51	A
1	B	53	G
1	B	54	G
1	B	57	U
1	B	59	A
1	B	61	G

There are no RNA pucker outliers to report.

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 11 ligands modelled in this entry, 11 are monoatomic - leaving 0 for Mogul analysis.

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 [i](#)

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	B	71/71 (100%)	-0.49	1 (1%) 73 51	22, 35, 61, 107	0
2	C	28/28 (100%)	-0.23	0 100 100	27, 36, 75, 78	0
3	D	8/8 (100%)	-0.63	0 100 100	27, 32, 36, 43	0
4	A	1030/1122 (91%)	0.26	34 (3%) 49 28	17, 57, 90, 112	0
All	All	1137/1229 (92%)	0.20	35 (3%) 51 30	17, 55, 90, 112	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	A	286	ASN	3.5
4	A	727	GLY	3.5
4	A	51	GLN	2.9
4	A	204	ASN	2.8
4	A	512	ARG	2.7
4	A	1112	ASN	2.7
4	A	725	HIS	2.7
4	A	726	TRP	2.6
4	A	288	THR	2.6
4	A	555	VAL	2.6
4	A	793	GLN	2.5
4	A	23	VAL	2.5
4	A	837	LYS	2.5
4	A	172	GLU	2.4
4	A	1114	GLY	2.4
4	A	38	ALA	2.3
4	A	500	TYR	2.3
4	A	501	GLY	2.3
4	A	939	ASP	2.3
4	A	928	ASN	2.2
4	A	340	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
4	A	0	MET	2.2
4	A	178	HIS	2.2
4	A	675	VAL	2.2
4	A	316	PRO	2.2
4	A	940	SER	2.1
4	A	454	GLY	2.1
1	B	0	G	2.1
4	A	510	MET	2.1
4	A	168	ASP	2.1
4	A	27	ILE	2.1
4	A	747	ALA	2.0
4	A	40	ALA	2.0
4	A	48	THR	2.0
4	A	320	PHE	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](#)

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
5	BA	C	101	1/1	0.65	0.27	245,245,245,245	0
6	MG	A	1206	1/1	0.79	0.15	41,41,41,41	0
5	BA	B	103	1/1	0.91	0.10	125,125,125,125	0
5	BA	A	1204	1/1	0.93	0.08	121,121,121,121	0
5	BA	A	1201	1/1	0.97	0.04	90,90,90,90	0
5	BA	A	1202	1/1	0.97	0.06	106,106,106,106	0
5	BA	B	104	1/1	0.97	0.07	114,114,114,114	0
5	BA	B	102	1/1	0.97	0.04	85,85,85,85	0
5	BA	A	1205	1/1	0.98	0.03	67,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	BA	A	1203	1/1	0.98	0.09	122,122,122,122	0
5	BA	B	101	1/1	1.00	0.03	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.