



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 13, 2022 – 06:10 PM JST

PDB ID : 7ENI
Title : Crystal structure of cas and anti-cas protein complex
Authors : Wang, Y.; Li, X.
Deposited on : 2021-04-17
Resolution : 2.63 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.27
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.27

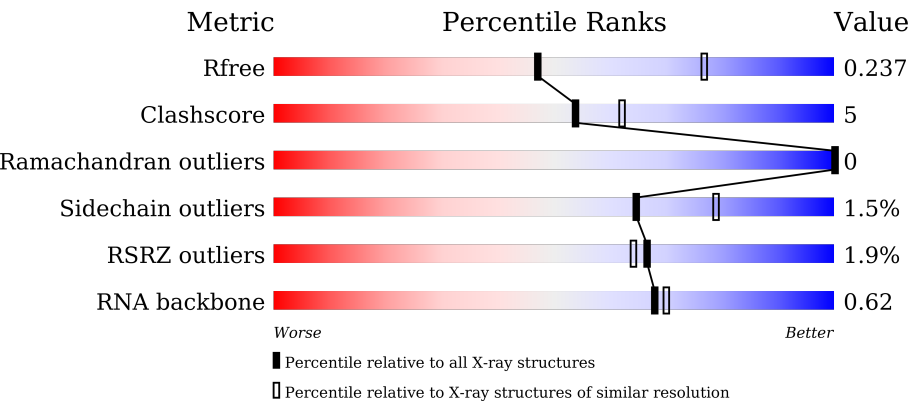
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.63 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}	130704	1426 (2.66-2.62)
Clashscore	141614	1472 (2.66-2.62)
Ramachandran outliers	138981	1446 (2.66-2.62)
Sidechain outliers	138945	1446 (2.66-2.62)
RSRZ outliers	127900	1408 (2.66-2.62)
RNA backbone	3102	1027 (2.96-2.32)

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	1054	<div><div>2%</div><div>86%</div><div>13%</div><div>..</div></div>
2	B	74	<div><div>%</div><div>50%</div><div>35%</div><div>8%</div><div>7%</div></div>
3	C	128	<div><div>4%</div><div>78%</div><div>20%</div><div>..</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11460 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	6	0
			8609	5461	1500	1632	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP J7RUA5

- Molecule 2 is a RNA chain called sgRNA (74-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	69	Total	C	N	O	P	0	0	0
			1441	645	247	480	69			

- Molecule 3 is a protein called AcrIIA13 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	126	Total	C	N	O	S	0	0	0
			1013	639	161	212	1			

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		

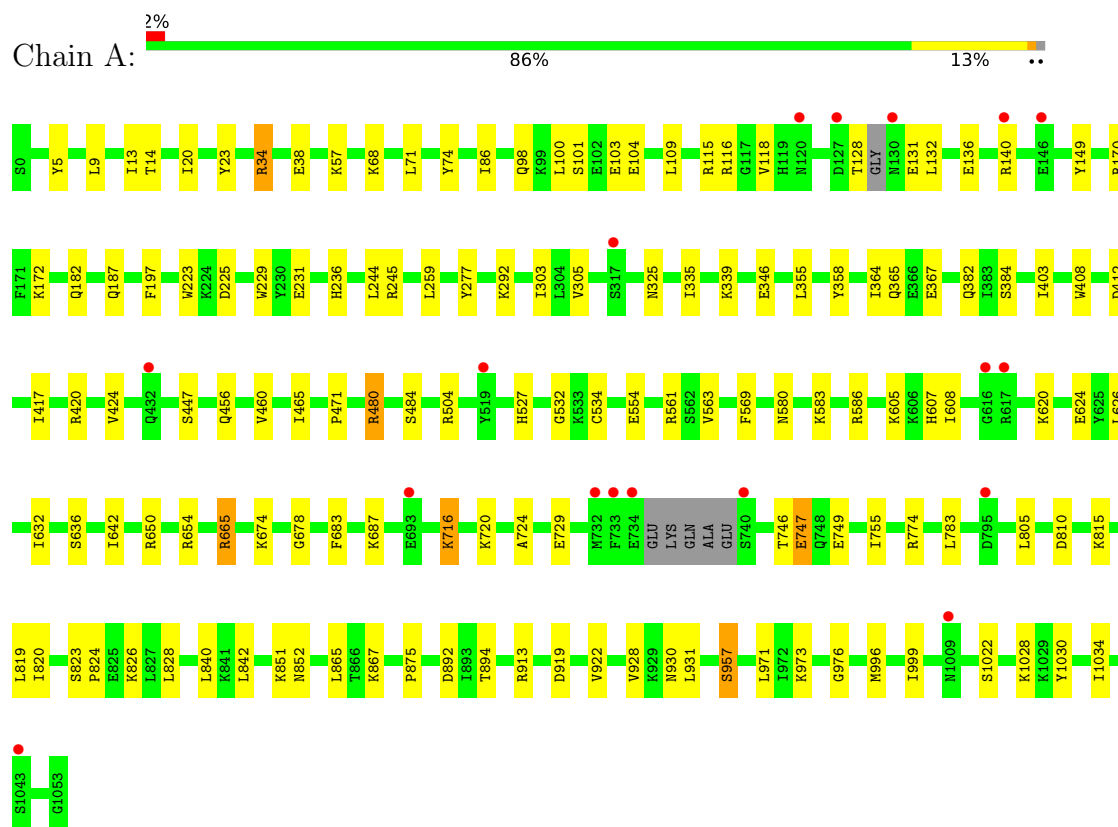
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	263	Total	O	0	0
			263	263		
5	B	74	Total	O	0	0
			74	74		
5	C	15	Total	O	0	0
			15	15		

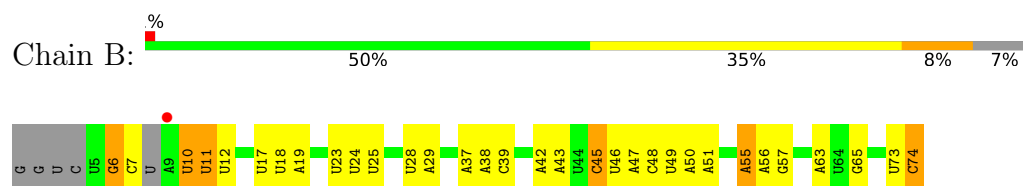
3 Residue-property plots

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.

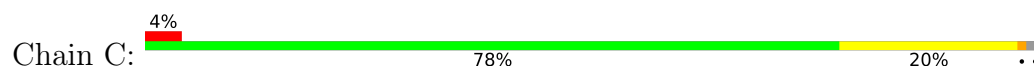
• Molecule 1: CRISPR-associated endonuclease Cas9

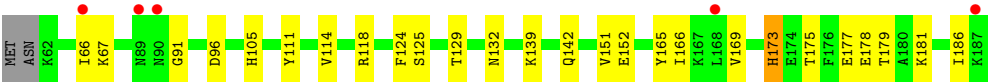


• Molecule 2: sgRNA (74-MER)



• Molecule 3: AcrIIA13 protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.85Å 105.96Å 238.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.24 – 2.63 20.24 – 2.63	Depositor EDS
% Data completeness (in resolution range)	87.1 (20.24-2.63) 87.1 (20.24-2.63)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.63Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.200 , 0.237 0.200 , 0.237	Depositor DCC
R_{free} test set	1995 reflections (3.43%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11460	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.24	0/8770	0.41	0/11807
2	B	0.21	0/1608	0.80	0/2498
3	C	0.25	0/1030	0.47	0/1398
All	All	0.24	0/11408	0.50	0/15703

There are no bond length outliers.

There are no bond angle outliers.

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	8609	0	8609	83	1
2	B	1441	0	726	23	0
3	C	1013	0	927	14	0
4	A	30	0	0	2	0
4	B	15	0	0	0	0
5	A	263	0	0	4	0
5	B	74	0	0	1	0
5	C	15	0	0	0	0
All	All	11460	0	10262	113	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 5.

All (113) 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	1.71	0.91
3:C:175:THR:HB	3:C:178:GLU:HB2	1.66	0.78
1:A:424:VAL:HG22	1:A:665:ARG:HH11	1.62	0.65
3:C:177:GLU:HG2	3:C:181:LYS:HE2	1.80	0.63
1:A:605:LYS:NZ	5:A:1205:HOH:O	2.31	0.63
1:A:867:LYS:HD2	2:B:46:U:H1'	1.81	0.62
1:A:236:HIS:ND1	4:A:1102:PO4:O1	2.30	0.62
1:A:132:LEU:O	1:A:170:ARG:NH2	2.32	0.62
1:A:277:TYR:OH	4:A:1106:PO4:O1	2.19	0.61
1:A:867:LYS:NZ	2:B:45:C:O2	2.36	0.58
3:C:91:GLY:O	3:C:124:PHE:N	2.28	0.58
1:A:913:ARG:HG2	1:A:928:VAL:HB	1.86	0.58
1:A:68:LYS:NZ	5:B:204:HOH:O	2.36	0.58
1:A:922:VAL:HG23	1:A:1022:SER:HA	1.87	0.56
1:A:412:ASP:OD2	1:A:420:ARG:NH2	2.36	0.56
2:B:47:A:H2'	2:B:48:C:C6	2.41	0.56
1:A:840:LEU:HD13	1:A:865:LEU:HD22	1.88	0.56
1:A:115:ARG:NH1	5:A:1217:HOH:O	2.40	0.55
1:A:554:GLU:HG2	1:A:580:ASN:HD22	1.72	0.54
1:A:805:LEU:HD11	1:A:815:LYS:HG2	1.89	0.54
1:A:561:ARG:NH2	1:A:747:GLU:OE1	2.39	0.54
1:A:23:TYR:CZ	1:A:716:LYS:HD2	2.43	0.54
1:A:563:VAL:HG11	1:A:632:ILE:HD12	1.90	0.54
1:A:642:ILE:HD13	1:A:724:ALA:HA	1.91	0.53
3:C:181:LYS:HG2	3:C:186:ILE:HG13	1.91	0.53
3:C:96:ASP:OD2	3:C:118:ARG:NH1	2.43	0.52
1:A:13:ILE:HG13	1:A:14:THR:HG23	1.91	0.52
3:C:66:ILE:O	3:C:67:LYS:HD2	2.08	0.52
1:A:931:LEU:HD22	2:B:56:A:H1'	1.91	0.52
1:A:828:LEU:HB3	2:B:45:C:H5'	1.92	0.52
3:C:129:THR:H	3:C:132:ASN:HB2	1.75	0.52
1:A:810:ASP:O	1:A:852:ASN:ND2	2.27	0.52
3:C:165:TYR:O	3:C:169:VAL:HG22	2.10	0.51
1:A:678:GLY:HA3	5:A:1302:HOH:O	2.10	0.51
1:A:38:GLU:OE1	1:A:484:SER:OG	2.19	0.51
2:B:24:U:H2'	2:B:25:U:C6	2.46	0.51
1:A:867:LYS:HB3	1:A:875:PRO:HD2	1.93	0.51
1:A:292:LYS:O	1:A:325:ASN:HB2	2.11	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:50:A:H2'	2:B:51:A:C8	2.47	0.50
1:A:554:GLU:HG2	1:A:580:ASN:ND2	2.26	0.50
3:C:67:LYS:HE3	3:C:67:LYS:HA	1.94	0.50
1:A:683:PHE:O	1:A:687:LYS:HG2	2.12	0.50
1:A:823:SER:HB2	1:A:826:LYS:HG3	1.93	0.50
1:A:86:ILE:HD12	1:A:104:GLU:HB3	1.94	0.50
1:A:335:ILE:HD13	1:A:382:GLN:HB3	1.94	0.50
1:A:365:GLN:HG3	1:A:384:SER:HB3	1.93	0.49
2:B:6:G:H4'	2:B:7:C:OP2	2.11	0.49
1:A:465:ILE:HG12	1:A:471:PRO:HD3	1.93	0.49
1:A:607:HIS:NE2	5:A:1209:HOH:O	2.35	0.49
3:C:151:VAL:HG12	3:C:152:GLU:HG2	1.94	0.49
1:A:149:TYR:OH	1:A:182:GLN:HG3	2.12	0.49
3:C:139:LYS:HA	3:C:142:GLN:HG2	1.95	0.49
1:A:687:LYS:HB3	1:A:755:ILE:HG23	1.94	0.49
2:B:10:U:O2'	2:B:11:U:OP1	2.28	0.49
1:A:57:LYS:NZ	2:B:55:A:OP1	2.46	0.48
1:A:403:ILE:HG23	1:A:417:ILE:HG21	1.95	0.48
2:B:49:U:H2'	2:B:50:A:C8	2.49	0.48
1:A:5:TYR:CD2	1:A:20:ILE:HG23	2.48	0.48
1:A:223:TRP:CD1	1:A:229:TRP:HB2	2.48	0.48
1:A:674:LYS:HE3	1:A:729:GLU:HG3	1.96	0.47
1:A:828:LEU:HD23	2:B:45:C:H5'	1.95	0.47
1:A:355:LEU:HD23	1:A:364:ILE:HD13	1.97	0.47
2:B:50:A:H2'	2:B:51:A:H8	1.80	0.47
2:B:47:A:H2'	2:B:48:C:H6	1.78	0.47
3:C:105:HIS:HB2	3:C:111:TYR:HE1	1.79	0.47
1:A:98:GLN:O	1:A:187:GLN:NE2	2.49	0.46
1:A:608:ILE:HD12	1:A:626:LEU:HD13	1.97	0.46
1:A:456:GLN:O	1:A:460:VAL:HG23	2.16	0.46
2:B:18:U:H2'	2:B:19:A:C8	2.51	0.46
1:A:74:TYR:CZ	1:A:197:PHE:HB2	2.50	0.46
1:A:23:TYR:CE1	1:A:716:LYS:HD2	2.51	0.45
2:B:17:U:H2'	2:B:18:U:C6	2.51	0.45
1:A:358:TYR:CZ	1:A:367:GLU:HG3	2.52	0.44
1:A:23:TYR:O	1:A:716:LYS:HE2	2.18	0.44
1:A:136:GLU:O	1:A:140:ARG:HD3	2.18	0.44
1:A:527:HIS:CD2	1:A:532:GLY:HA2	2.52	0.44
1:A:231:GLU:HG2	1:A:244:LEU:HD11	1.98	0.44
1:A:820:ILE:O	1:A:824:PRO:HG3	2.18	0.44
2:B:49:U:H2'	2:B:50:A:H8	1.81	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:LYS:HB3	1:A:339:LYS:HE3	1.78	0.44
3:C:114:VAL:HG22	3:C:166:ILE:HD11	1.99	0.43
1:A:259:LEU:HD11	1:A:408:TRP:CE2	2.53	0.43
1:A:100:LEU:HB3	1:A:104:GLU:HG3	2.00	0.43
1:A:34:ARG:HG3	1:A:1034:ILE:HG12	2.00	0.43
1:A:5:TYR:CE2	1:A:471:PRO:HB3	2.54	0.42
1:A:71:LEU:HD21	1:A:109:LEU:HD12	2.00	0.42
1:A:851:LYS:HA	1:A:851:LYS:HD3	1.89	0.42
1:A:996:MET:HB2	1:A:999:ILE:O	2.19	0.42
1:A:819:LEU:HD11	1:A:826:LYS:HE2	2.01	0.42
1:A:919:ASP:O	1:A:922:VAL:HG12	2.19	0.42
1:A:973:LYS:HE3	1:A:976:GLY:HA2	2.02	0.42
1:A:118:VAL:H	1:A:172:LYS:HA	1.84	0.42
2:B:23:U:H2'	2:B:24:U:C6	2.54	0.42
1:A:842:LEU:HD12	1:A:842:LEU:HA	1.95	0.41
1:A:720:LYS:NZ	1:A:749:GLU:OE2	2.41	0.41
1:A:128:THR:C	1:A:131:GLU:H	2.23	0.41
1:A:303:ILE:HG13	1:A:305:VAL:HG22	2.01	0.41
2:B:6:G:O2'	2:B:7:C:OP1	2.36	0.41
2:B:28:U:C2	2:B:29:A:C8	3.09	0.41
1:A:919:ASP:OD1	1:A:957:SER:HB3	2.19	0.41
1:A:892:ASP:OD1	1:A:894:THR:OG1	2.28	0.41
1:A:101:SER:OG	1:A:103[B]:GLU:OE1	2.38	0.41
1:A:583:LYS:HE2	1:A:586:ARG:O	2.20	0.41
3:C:173:HIS:HB2	3:C:179:THR:OG1	2.19	0.40
1:A:650:ARG:O	1:A:654:ARG:HG3	2.21	0.40
1:A:783:LEU:HD23	1:A:930:ASN:HB3	2.03	0.40
2:B:37:A:H2'	2:B:38:A:C8	2.57	0.40
1:A:9:LEU:HD23	1:A:9:LEU:HA	1.89	0.40
1:A:480:ARG:HD3	1:A:480:ARG:HA	1.84	0.40
1:A:620:LYS:O	1:A:624:GLU:HG3	2.21	0.40
1:A:774:ARG:HG3	2:B:74:C:OP1	2.21	0.40
2:B:42:A:C4	2:B:43:A:C8	3.10	0.40
1:A:973:LYS:HB3	1:A:1028:LYS:HB2	2.04	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:346:GLU:OE2	1:A:746:THR:OG1[3_554]	2.09	0.11

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	1048/1054 (99%)	1027 (98%)	21 (2%)	0	100	100
3	C	124/128 (97%)	118 (95%)	6 (5%)	0	100	100
All	All	1172/1182 (99%)	1145 (98%)	27 (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	940/973 (97%)	926 (98%)	14 (2%)	65	79
3	C	109/118 (92%)	107 (98%)	2 (2%)	59	75
All	All	1049/1091 (96%)	1033 (98%)	16 (2%)	65	79

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	ARG
1	A	116	ARG
1	A	225	ASP
1	A	245	ARG
1	A	447	SER
1	A	480	ARG
1	A	504	ARG
1	A	534	CYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	569	PHE
1	A	636	SER
1	A	665	ARG
1	A	716	LYS
1	A	747	GLU
1	A	957	SER
3	C	125	SER
3	C	173	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	958	ASN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	66/74 (89%)	12 (18%)	2 (3%)

All (12) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	6	G
2	B	10	U
2	B	11	U
2	B	12	U
2	B	39	C
2	B	45	C
2	B	55	A
2	B	57	G
2	B	63	A
2	B	65	G
2	B	73	U
2	B	74	C

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	6	G
2	B	10	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	PO4	B	103	-	4,4,4	0.90	0	6,6,6	0.46	0
4	PO4	A	1106	-	4,4,4	0.88	0	6,6,6	0.50	0
4	PO4	A	1102	-	4,4,4	0.91	0	6,6,6	0.48	0
4	PO4	B	101	-	4,4,4	0.88	0	6,6,6	0.45	0
4	PO4	A	1101	-	4,4,4	0.89	0	6,6,6	0.45	0
4	PO4	A	1103	-	4,4,4	0.91	0	6,6,6	0.43	0
4	PO4	A	1104	-	4,4,4	0.90	0	6,6,6	0.44	0
4	PO4	B	102	-	4,4,4	0.89	0	6,6,6	0.43	0
4	PO4	A	1105	-	4,4,4	0.88	0	6,6,6	0.47	0

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1106	PO4	1	0
4	A	1102	PO4	1	0

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	1048/1054 (99%)	-0.32	18 (1%) 70 67	19, 38, 65, 88	0
2	B	69/74 (93%)	-0.65	1 (1%) 75 73	24, 35, 86, 94	0
3	C	126/128 (98%)	-0.13	5 (3%) 38 35	29, 55, 74, 79	0
All	All	1243/1256 (98%)	-0.32	24 (1%) 66 64	19, 39, 67, 94	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	130	ASN	3.6
3	C	168	LEU	3.5
2	B	9	A	3.4
1	A	733	PHE	3.1
1	A	734	GLU	3.0
1	A	120	ASN	2.9
3	C	90	ASN	2.8
1	A	617	ARG	2.7
1	A	740	SER	2.5
1	A	1009	ASN	2.4
3	C	89	ASN	2.3
1	A	616	GLY	2.3
1	A	732	MET	2.3
1	A	146	GLU	2.3
3	C	66	ILE	2.2
1	A	795	ASP	2.2
3	C	187	LYS	2.2
1	A	1043	SER	2.1
1	A	432	GLN	2.1
1	A	317	SER	2.1
1	A	519	TYR	2.1
1	A	140	ARG	2.0
1	A	693	GLU	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	127	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 monosaccharides 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(\AA^2)	Q<0.9
4	PO4	A	1106	5/5	0.86	0.27	55,61,77,89	0
4	PO4	B	103	5/5	0.89	0.31	69,82,85,92	0
4	PO4	B	101	5/5	0.90	0.27	53,64,65,88	0
4	PO4	A	1105	5/5	0.92	0.26	56,56,63,88	0
4	PO4	A	1101	5/5	0.92	0.22	39,58,63,88	0
4	PO4	A	1102	5/5	0.95	0.15	60,65,83,87	0
4	PO4	A	1104	5/5	0.95	0.30	62,63,74,86	0
4	PO4	B	102	5/5	0.97	0.13	29,45,46,56	0
4	PO4	A	1103	5/5	0.97	0.15	43,48,68,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.