



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 11:05 pm BST

PDB ID : 6O79
Title : Crystal structure of Csm1-Csm4 cassette in complex with cA3
Authors : Jia, N.; Patel, D.J.
Deposited on : 2019-03-07
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

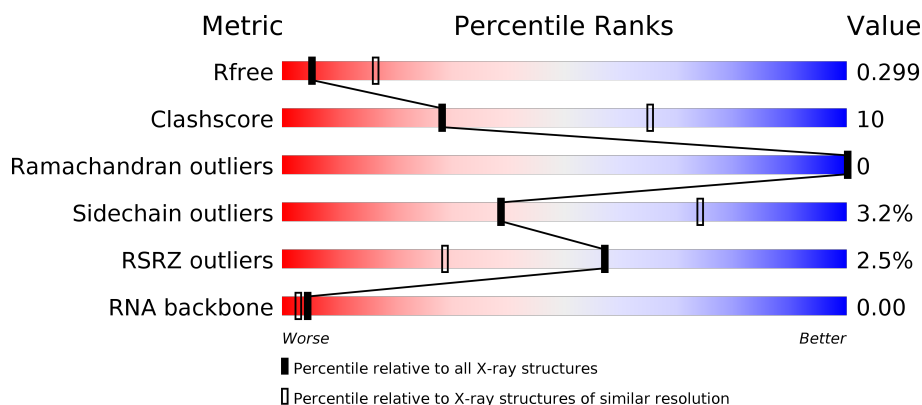
1 Overall quality at a glance

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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)
RNA backbone	3102	1173 (3.30-2.70)

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 on 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	791	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>15%</div> <div>•</div> <div>11%</div> </div> </div>
2	B	289	<div> <div>2%</div> <div> <div></div> <div>67%</div> <div>15%</div> <div>•</div> <div>15%</div> </div> </div>
3	C	3	<div> <div>33%</div> <div>67%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7699 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 system single-strand-specific deoxyribonuclease Cas10/Csm1 (subtype III-A).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	706	Total	C	N	O	S	0	0	0
			5671	3649	975	1031	16			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP B6YWB8
A	-12	GLY	-	expression tag	UNP B6YWB8
A	-11	SER	-	expression tag	UNP B6YWB8
A	-10	SER	-	expression tag	UNP B6YWB8
A	-9	HIS	-	expression tag	UNP B6YWB8
A	-8	HIS	-	expression tag	UNP B6YWB8
A	-7	HIS	-	expression tag	UNP B6YWB8
A	-6	HIS	-	expression tag	UNP B6YWB8
A	-5	HIS	-	expression tag	UNP B6YWB8
A	-4	HIS	-	expression tag	UNP B6YWB8
A	-3	SER	-	expression tag	UNP B6YWB8
A	-2	GLN	-	expression tag	UNP B6YWB8
A	-1	ASP	-	expression tag	UNP B6YWB8
A	0	PRO	-	expression tag	UNP B6YWB8
A	589	ALA	ASP	conflict	UNP B6YWB8

- Molecule 2 is a protein called Csm4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	246	Total	C	N	O	S	0	0	0
			1962	1276	326	356	4			

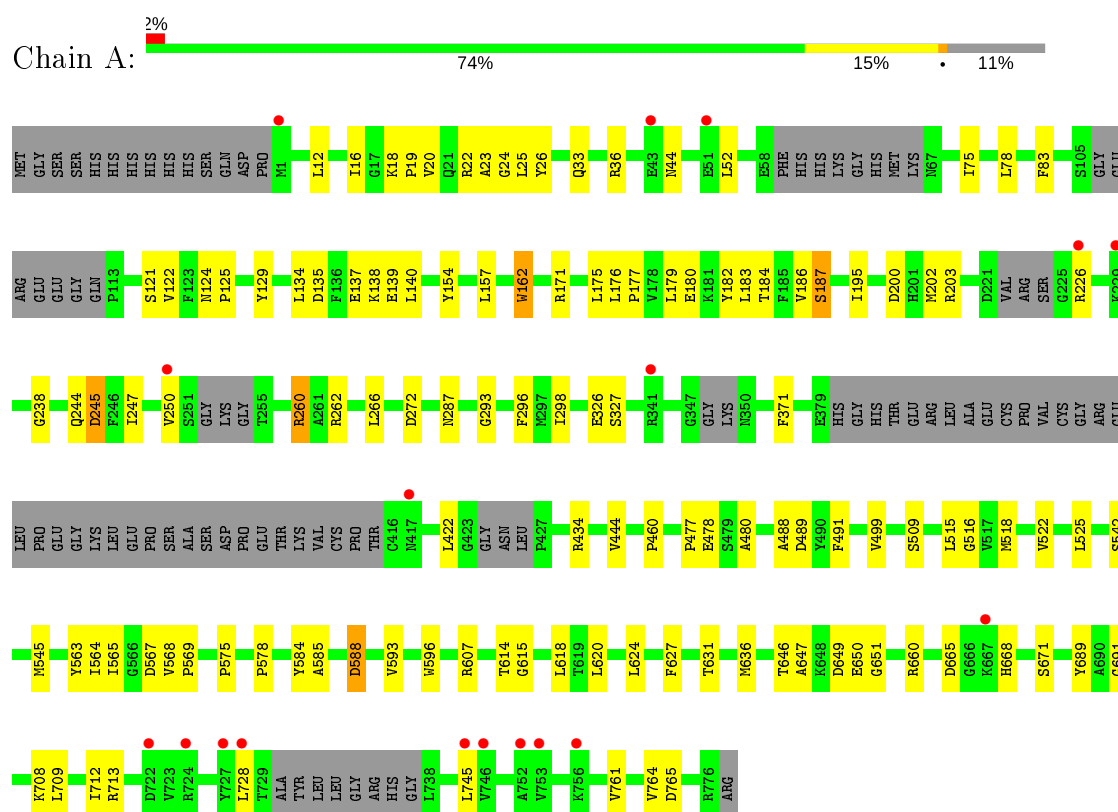
- Molecule 3 is a RNA chain called cyclic RNA cA3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	3	Total	C	N	O	P	0	0	0
			66	30	15	18	3			

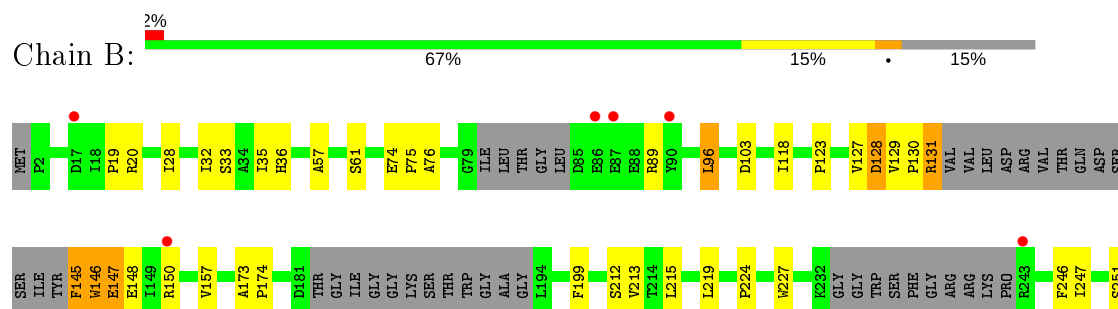
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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 system single-strand-specific deoxyribonuclease Cas10/Csm1 (subtype III-A)



- Molecule 2: Csm4





- Molecule 3: cyclic RNA cA3



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	157.01Å 157.01Å 186.05Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.59 – 3.00 49.54 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.59-3.00) 99.8 (49.54-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.254 , 0.303 0.254 , 0.299	Depositor DCC
R_{free} test set	2707 reflections (9.80%)	wwPDB-VP
Wilson B-factor (Å ²)	93.9	Xtriage
Anisotropy	0.253	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7699	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.97% 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	0.64	0/5793	0.72	0/7809
2	B	0.63	0/2012	0.74	0/2720
3	C	4.46	22/74 (29.7%)	6.45	37/113 (32.7%)
All	All	0.77	22/7879 (0.3%)	0.98	37/10642 (0.3%)

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	A	N7-C5	-11.87	1.32	1.39
3	C	2	A	N7-C5	-9.69	1.33	1.39
3	C	3	A	N7-C5	-8.96	1.33	1.39
3	C	1	A	C5-C6	-8.55	1.33	1.41
3	C	1	A	C2-N3	8.13	1.40	1.33
3	C	2	A	C5-C6	-8.05	1.33	1.41
3	C	3	A	N9-C4	-8.00	1.33	1.37
3	C	2	A	N9-C4	-7.96	1.33	1.37
3	C	3	A	C2-N3	7.75	1.40	1.33
3	C	2	A	C2-N3	7.70	1.40	1.33
3	C	3	A	C5-C6	-7.69	1.34	1.41
3	C	1	A	N9-C4	-7.64	1.33	1.37
3	C	2	A	N1-C2	7.33	1.41	1.34
3	C	1	A	C5-C4	-7.01	1.33	1.38
3	C	3	A	C5-C4	-6.84	1.33	1.38
3	C	3	A	N1-C2	6.75	1.40	1.34
3	C	2	A	N9-C8	-6.72	1.32	1.37
3	C	1	A	N1-C2	6.66	1.40	1.34
3	C	2	A	C5-C4	-6.34	1.34	1.38
3	C	1	A	N9-C8	-6.01	1.32	1.37
3	C	2	A	C3'-C2'	-5.45	1.46	1.52
3	C	3	A	N9-C8	-5.38	1.33	1.37

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	3	A	N1-C2-N3	-22.48	118.06	129.30
3	C	1	A	N1-C2-N3	-20.77	118.91	129.30
3	C	2	A	N1-C2-N3	-20.34	119.13	129.30
3	C	3	A	C2-N3-C4	18.07	119.64	110.60
3	C	1	A	C2-N3-C4	17.08	119.14	110.60
3	C	2	A	C2-N3-C4	16.44	118.82	110.60
3	C	3	A	N7-C8-N9	-12.68	107.46	113.80
3	C	1	A	N7-C8-N9	-11.41	108.09	113.80
3	C	3	A	C1'-O4'-C4'	-11.17	100.97	109.90
3	C	2	A	N7-C8-N9	-11.14	108.23	113.80
3	C	1	A	C5-C6-N6	-10.05	115.66	123.70
3	C	2	A	C5-C6-N6	-9.97	115.73	123.70
3	C	1	A	C4-C5-C6	9.52	121.76	117.00
3	C	3	A	C5-C6-N6	-9.25	116.30	123.70
3	C	1	A	N3-C4-C5	-9.20	120.36	126.80
3	C	1	A	C5-N7-C8	9.11	108.46	103.90
3	C	3	A	C5-C6-N1	9.03	122.22	117.70
3	C	3	A	C5-N7-C8	8.64	108.22	103.90
3	C	2	A	N3-C4-C5	-8.38	120.94	126.80
3	C	2	A	C4-C5-C6	8.36	121.18	117.00
3	C	2	A	C5-C6-N1	8.13	121.77	117.70
3	C	3	A	N3-C4-C5	-8.13	121.11	126.80
3	C	2	A	C8-N9-C4	7.94	108.98	105.80
3	C	1	A	C5-C6-N1	7.85	121.63	117.70
3	C	2	A	C5-N7-C8	7.73	107.77	103.90
3	C	3	A	C8-N9-C4	7.68	108.87	105.80
3	C	1	A	N1-C6-N6	6.84	122.71	118.60
3	C	3	A	C4'-C3'-C2'	-6.68	95.92	102.60
3	C	3	A	C4-C5-C6	6.52	120.26	117.00
3	C	2	A	N1-C6-N6	6.51	122.51	118.60
3	C	2	A	N3-C4-N9	6.21	132.36	127.40
3	C	1	A	N3-C4-N9	5.92	132.14	127.40
3	C	3	A	O5'-C5'-C4'	-5.64	100.98	111.70
3	C	3	A	C4-C5-N7	-5.61	107.90	110.70
3	C	1	A	C4-C5-N7	-5.45	107.98	110.70
3	C	1	A	C8-N9-C4	5.42	107.97	105.80
3	C	1	A	O4'-C1'-N9	5.30	112.44	108.20

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	5671	0	5677	97	0
2	B	1962	0	1953	51	0
3	C	66	0	33	5	0
All	All	7699	0	7663	150	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 10.

All (150) 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)
2:B:130:PRO:CA	2:B:145:PHE:HD1	1.54	1.20
2:B:130:PRO:CA	2:B:145:PHE:CD1	2.33	1.11
2:B:130:PRO:HB3	2:B:145:PHE:CD1	1.86	1.10
2:B:130:PRO:HA	2:B:145:PHE:CB	1.82	1.07
2:B:130:PRO:HA	2:B:145:PHE:CD1	1.92	1.04
2:B:130:PRO:CB	2:B:145:PHE:CD1	2.43	1.00
2:B:262:ARG:HG3	2:B:262:ARG:HH11	1.31	0.95
2:B:127:VAL:O	2:B:147:GLU:HB2	1.68	0.93
2:B:130:PRO:HA	2:B:145:PHE:CG	2.09	0.88
1:A:26:TYR:HD1	1:A:36:ARG:HD2	1.39	0.86
1:A:26:TYR:CD1	1:A:36:ARG:HB3	2.14	0.83
1:A:488:ALA:HB3	1:A:585:ALA:HB3	1.61	0.82
1:A:26:TYR:HB3	1:A:36:ARG:HE	1.43	0.82
2:B:262:ARG:O	2:B:262:ARG:HD2	1.81	0.80
2:B:130:PRO:CB	2:B:145:PHE:HD1	1.86	0.80
2:B:130:PRO:HB3	2:B:145:PHE:CE1	2.17	0.79
2:B:130:PRO:HA	2:B:145:PHE:HB2	1.63	0.78
2:B:128:ASP:OD1	2:B:128:ASP:N	2.16	0.78
2:B:247:ILE:CG2	2:B:251:SER:OG	2.33	0.77
2:B:264:GLU:O	2:B:264:GLU:HG2	1.84	0.76
1:A:26:TYR:CD1	1:A:36:ARG:HD2	2.20	0.76
1:A:24:GLY:O	1:A:33:GLN:NE2	2.20	0.75
1:A:26:TYR:HD1	1:A:36:ARG:CD	2.01	0.73
2:B:127:VAL:C	2:B:128:ASP:OD1	2.28	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:130:PRO:N	2:B:145:PHE:HD1	1.85	0.72
2:B:224:PRO:HG2	2:B:227:TRP:CE3	2.24	0.72
1:A:139:GLU:HG3	1:A:140:LEU:H	1.54	0.72
1:A:522:VAL:HG21	1:A:545:MET:CE	2.20	0.71
1:A:564:ILE:N	1:A:564:ILE:HD12	2.05	0.70
1:A:564:ILE:H	1:A:564:ILE:HD12	1.55	0.70
1:A:23:ALA:HB2	1:A:154:TYR:HB3	1.74	0.69
2:B:130:PRO:HA	2:B:145:PHE:HB3	1.72	0.69
1:A:135:ASP:HB3	1:A:138:LYS:HB2	1.74	0.68
1:A:139:GLU:HG3	1:A:140:LEU:N	2.09	0.68
1:A:250:VAL:HG21	1:A:262:ARG:HH22	1.58	0.67
1:A:293:GLY:HA2	1:A:584:TYR:OH	1.94	0.67
3:C:1:A:O5'	3:C:3:A:O2'	2.10	0.66
2:B:173:ALA:HB3	2:B:174:PRO:HD3	1.77	0.65
1:A:171:ARG:NH2	1:A:478:GLU:O	2.29	0.65
1:A:25:LEU:HD12	1:A:25:LEU:N	2.12	0.64
2:B:148:GLU:OE1	2:B:150:ARG:NH2	2.30	0.64
1:A:26:TYR:HB2	1:A:33:GLN:HG3	1.80	0.63
1:A:134:LEU:HB2	1:A:184:THR:HG21	1.81	0.62
2:B:57:ALA:HB1	2:B:157:VAL:HG13	1.81	0.61
1:A:488:ALA:CB	1:A:585:ALA:HB3	2.30	0.60
1:A:326:GLU:HG2	1:A:326:GLU:O	2.00	0.60
2:B:263:LEU:HD23	2:B:263:LEU:N	2.18	0.59
2:B:262:ARG:HG3	2:B:262:ARG:NH1	2.09	0.59
1:A:564:ILE:CD1	1:A:564:ILE:H	2.15	0.59
1:A:627:PHE:CD2	1:A:636:MET:HG2	2.38	0.58
1:A:660:ARG:HD2	1:A:765:ASP:OD2	2.02	0.58
1:A:26:TYR:CE1	1:A:36:ARG:HB3	2.38	0.58
2:B:57:ALA:HB1	2:B:157:VAL:CG1	2.34	0.58
2:B:247:ILE:HG22	2:B:251:SER:OG	2.04	0.57
1:A:647:ALA:O	1:A:651:GLY:O	2.22	0.57
1:A:491:PHE:HB2	1:A:499:VAL:HG23	1.86	0.56
2:B:96:LEU:HD22	2:B:118:ILE:CG1	2.38	0.54
1:A:565:ILE:HG13	1:A:565:ILE:O	2.06	0.54
2:B:131:ARG:HG3	2:B:131:ARG:O	2.08	0.53
1:A:646:THR:O	1:A:650:GLU:HG3	2.07	0.53
1:A:525:LEU:HD22	1:A:588:ASP:HB3	1.89	0.53
2:B:74:GLU:HB2	2:B:75:PRO:HD3	1.91	0.53
1:A:713:ARG:HG3	1:A:761:VAL:HG12	1.91	0.53
1:A:180:GLU:O	1:A:184:THR:HB	2.09	0.53
2:B:262:ARG:HH11	2:B:262:ARG:CG	2.10	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1:A:H3'	3:C:3:A:O2'	2.09	0.53
1:A:44:ASN:HD22	1:A:162:TRP:HZ2	1.56	0.52
1:A:20:VAL:HG23	1:A:33:GLN:HG2	1.89	0.52
1:A:245:ASP:N	1:A:245:ASP:OD1	2.43	0.52
1:A:122:VAL:HA	1:A:195:ILE:HG23	1.91	0.52
2:B:145:PHE:CD2	2:B:145:PHE:N	2.78	0.52
1:A:179:LEU:HD22	1:A:202:MET:HE1	1.92	0.52
1:A:260:ARG:NH2	1:A:489:ASP:O	2.42	0.52
1:A:631:THR:HB	1:A:636:MET:HE3	1.91	0.51
1:A:745:LEU:HD11	1:A:764:VAL:HG21	1.92	0.51
1:A:22:ARG:HH21	1:A:186:VAL:CG1	2.24	0.51
2:B:262:ARG:NH1	2:B:262:ARG:CG	2.72	0.51
1:A:26:TYR:N	1:A:33:GLN:NE2	2.58	0.51
1:A:647:ALA:HA	1:A:671:SER:OG	2.10	0.50
1:A:25:LEU:N	1:A:25:LEU:CD1	2.74	0.50
1:A:689:TYR:CE2	1:A:691:GLY:HA2	2.47	0.50
1:A:18:LYS:HB2	1:A:19:PRO:HD3	1.93	0.50
1:A:518:MET:HB3	1:A:624:LEU:HD23	1.94	0.49
1:A:22:ARG:NH2	1:A:187:SER:O	2.46	0.49
1:A:708:LYS:O	1:A:712:ILE:HG13	2.13	0.49
1:A:563:TYR:CD2	1:A:563:TYR:C	2.85	0.49
1:A:272:ASP:OD1	1:A:434:ARG:NH1	2.45	0.49
2:B:145:PHE:N	2:B:145:PHE:HD2	2.11	0.49
1:A:614:THR:HG21	1:A:618:LEU:HB2	1.95	0.48
1:A:18:LYS:O	1:A:22:ARG:HG3	2.14	0.48
2:B:96:LEU:N	2:B:96:LEU:HD23	2.29	0.48
1:A:250:VAL:HG11	1:A:262:ARG:HH22	1.78	0.48
1:A:588:ASP:OD1	3:C:1:A:OP1	2.32	0.48
1:A:26:TYR:H	1:A:33:GLN:NE2	2.11	0.48
1:A:477:PRO:HG2	1:A:480:ALA:HB2	1.95	0.48
1:A:607:ARG:HA	1:A:620:LEU:HD12	1.96	0.47
1:A:709:LEU:HA	1:A:712:ILE:HD12	1.96	0.47
1:A:75:ILE:O	1:A:78:LEU:O	2.33	0.47
1:A:125:PRO:HG3	1:A:195:ILE:HD11	1.97	0.47
1:A:26:TYR:CD1	1:A:36:ARG:CD	2.88	0.47
1:A:327:SER:HA	1:A:371:PHE:CE2	2.50	0.47
1:A:522:VAL:HG22	1:A:620:LEU:HD23	1.96	0.46
1:A:121:SER:HA	1:A:129:TYR:CE2	2.51	0.46
1:A:515:LEU:HD23	1:A:636:MET:HE2	1.98	0.46
1:A:247:ILE:CD1	3:C:2:A:C2	2.99	0.46
2:B:215:LEU:HD11	2:B:280:LEU:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:LEU:HD23	1:A:175:LEU:HD23	1.97	0.45
2:B:131:ARG:C	2:B:131:ARG:CD	2.85	0.45
1:A:575:PRO:HG2	1:A:578:PRO:HA	1.98	0.45
2:B:129:VAL:O	2:B:145:PHE:HB2	2.15	0.45
2:B:213:VAL:O	2:B:279:PRO:HA	2.17	0.45
2:B:96:LEU:HD22	2:B:118:ILE:HG13	1.98	0.45
2:B:19:PRO:O	2:B:57:ALA:HB3	2.17	0.45
1:A:157:LEU:CD1	1:A:182:TYR:HB3	2.47	0.45
1:A:238:GLY:HA3	1:A:296:PHE:CZ	2.53	0.44
1:A:26:TYR:HD1	1:A:36:ARG:NE	2.17	0.43
2:B:76:ALA:HB2	2:B:285:PRO:HG3	2.00	0.43
1:A:522:VAL:HG22	1:A:620:LEU:CD2	2.49	0.43
1:A:327:SER:HA	1:A:371:PHE:CZ	2.53	0.43
1:A:745:LEU:HD21	1:A:764:VAL:HG21	1.99	0.43
2:B:35:ILE:HG22	2:B:36:HIS:CD2	2.54	0.43
1:A:287:ASN:O	1:A:298:ILE:HA	2.19	0.43
1:A:183:LEU:HD12	1:A:202:MET:HE1	2.00	0.42
1:A:16:ILE:O	1:A:19:PRO:HD2	2.20	0.42
2:B:263:LEU:HD21	2:B:273:VAL:HG23	2.01	0.42
1:A:137:GLU:H	1:A:137:GLU:HG2	1.50	0.42
1:A:422:LEU:HD22	1:A:444:VAL:HG11	2.01	0.42
1:A:52:LEU:HD22	1:A:83:PHE:CD2	2.54	0.42
1:A:124:ASN:CB	1:A:615:GLY:HA3	2.50	0.41
1:A:122:VAL:CA	1:A:195:ILE:HG23	2.50	0.41
2:B:219:LEU:HG	2:B:246:PHE:CE1	2.54	0.41
2:B:277:THR:O	2:B:279:PRO:HD3	2.20	0.41
1:A:176:LEU:HB2	1:A:177:PRO:HD3	2.01	0.41
1:A:568:VAL:O	1:A:569:PRO:C	2.58	0.41
1:A:200:ASP:HA	1:A:203:ARG:HD3	2.02	0.41
2:B:28:ILE:O	2:B:32:ILE:HG12	2.21	0.41
1:A:266:LEU:HD23	3:C:2:A:C6	2.55	0.41
1:A:515:LEU:HD23	1:A:636:MET:CE	2.51	0.41
1:A:745:LEU:HD21	1:A:764:VAL:CG2	2.50	0.41
2:B:252:ILE:HD12	2:B:282:VAL:HG11	2.02	0.41
2:B:61:SER:HB2	2:B:123:PRO:HG3	2.02	0.41
2:B:146:TRP:N	2:B:146:TRP:CD1	2.89	0.41
1:A:26:TYR:CD1	1:A:36:ARG:CB	2.97	0.40
1:A:460:PRO:HG2	1:A:477:PRO:HD3	2.03	0.40
1:A:712:ILE:HG12	1:A:728:LEU:HD22	2.02	0.40
1:A:509:SER:HB2	1:A:596:TRP:CD1	2.57	0.40
2:B:224:PRO:HG3	2:B:247:ILE:HD11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:SER:OG	1:A:327:SER:O	2.35	0.40
1:A:516:GLY:O	1:A:593:VAL:HA	2.21	0.40
1:A:665:ASP:OD1	1:A:668:HIS:CD2	2.74	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	688/791 (87%)	646 (94%)	42 (6%)	0	100	100
2	B	236/289 (82%)	222 (94%)	14 (6%)	0	100	100
All	All	924/1080 (86%)	868 (94%)	56 (6%)	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	595/664 (90%)	585 (98%)	10 (2%)	60	85
2	B	206/240 (86%)	190 (92%)	16 (8%)	12	42
All	All	801/904 (89%)	775 (97%)	26 (3%)	39	74

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	TRP
1	A	187	SER
1	A	226	ARG
1	A	244	GLN
1	A	245	ASP
1	A	260	ARG
1	A	542	SER
1	A	567	ASP
1	A	588	ASP
1	A	649	ASP
2	B	20	ARG
2	B	33	SER
2	B	89	ARG
2	B	96	LEU
2	B	103	ASP
2	B	128	ASP
2	B	131	ARG
2	B	145	PHE
2	B	146	TRP
2	B	147	GLU
2	B	199	PHE
2	B	212	SER
2	B	262	ARG
2	B	263	LEU
2	B	265	LEU
2	B	271	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	287	ASN
1	A	668	HIS
1	A	675	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	C	3/3 (100%)	1 (33%)	1 (33%)

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	C	3	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	C	1	A

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 carbohydrates 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	706/791 (89%)	0.06	18 (2%) 57 29	65, 98, 154, 192	0
2	B	246/289 (85%)	0.00	6 (2%) 59 30	68, 96, 149, 180	0
3	C	3/3 (100%)	-0.17	0 100 100	91, 91, 98, 131	0
All	All	955/1083 (88%)	0.05	24 (2%) 57 29	65, 98, 153, 192	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	753	VAL	6.0
2	B	86	GLU	4.7
1	A	752	ALA	4.6
2	B	90	TYR	4.4
1	A	1	MET	4.2
1	A	728	LEU	4.1
1	A	250	VAL	4.1
1	A	727	TYR	3.6
1	A	756	LYS	3.4
2	B	243	ARG	3.0
1	A	226	ARG	2.9
1	A	341	ARG	2.5
1	A	229	LYS	2.4
1	A	724	ARG	2.4
1	A	417	ASN	2.3
1	A	667	LYS	2.2
2	B	17	ASP	2.2
1	A	51	GLU	2.2
2	B	150	ARG	2.1
2	B	87	GLU	2.1
1	A	745	LEU	2.0
1	A	43	GLU	2.0
1	A	722	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	746	VAL	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 carbohydrates 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.