



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:36 AM UTC

PDB ID : 6O7B / pdb_00006o7b
Title : Crystal structure of Csm1-Csm4 cassette in complex with cA4
Authors : Jia, N.; Patel, D.J.
Deposited on : 2019-03-07
Resolution : 2.40 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

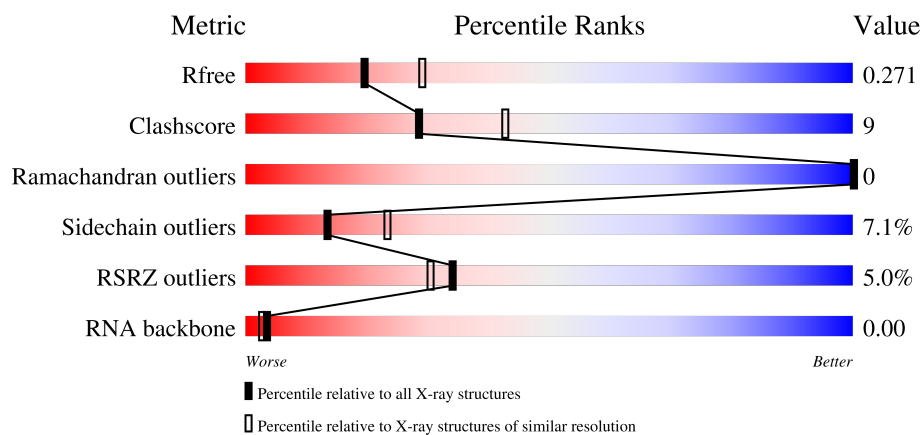
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)
RNA backbone	3983	1155 (2.70-2.10)

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	791	<div> <div>4%</div> <div>68%</div> <div>19%</div> <div>11%</div> </div>
2	B	289	<div> <div>5%</div> <div>75%</div> <div>11%</div> <div>12%</div> </div>
3	C	4	<div> <div>25%</div> <div>75%</div> </div>
3	D	4	<div> <div>25%</div> <div>100%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7884 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	703	Total	C	N	O	S	0	0	0
			5653	3642	971	1024	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	engineered mutation	UNP B6YWB8

- Molecule 2 is a protein called Csm4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	254	Total	C	N	O	S	0	0	0
			2027	1320	334	369	4			

- Molecule 3 is a RNA chain called Cyclic RNA cA4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	4	Total 88	C 40	N 20	O 24	P 4	0	0	0
3	D	4	Total 88	C 40	N 20	O 24	P 4	0	0	0

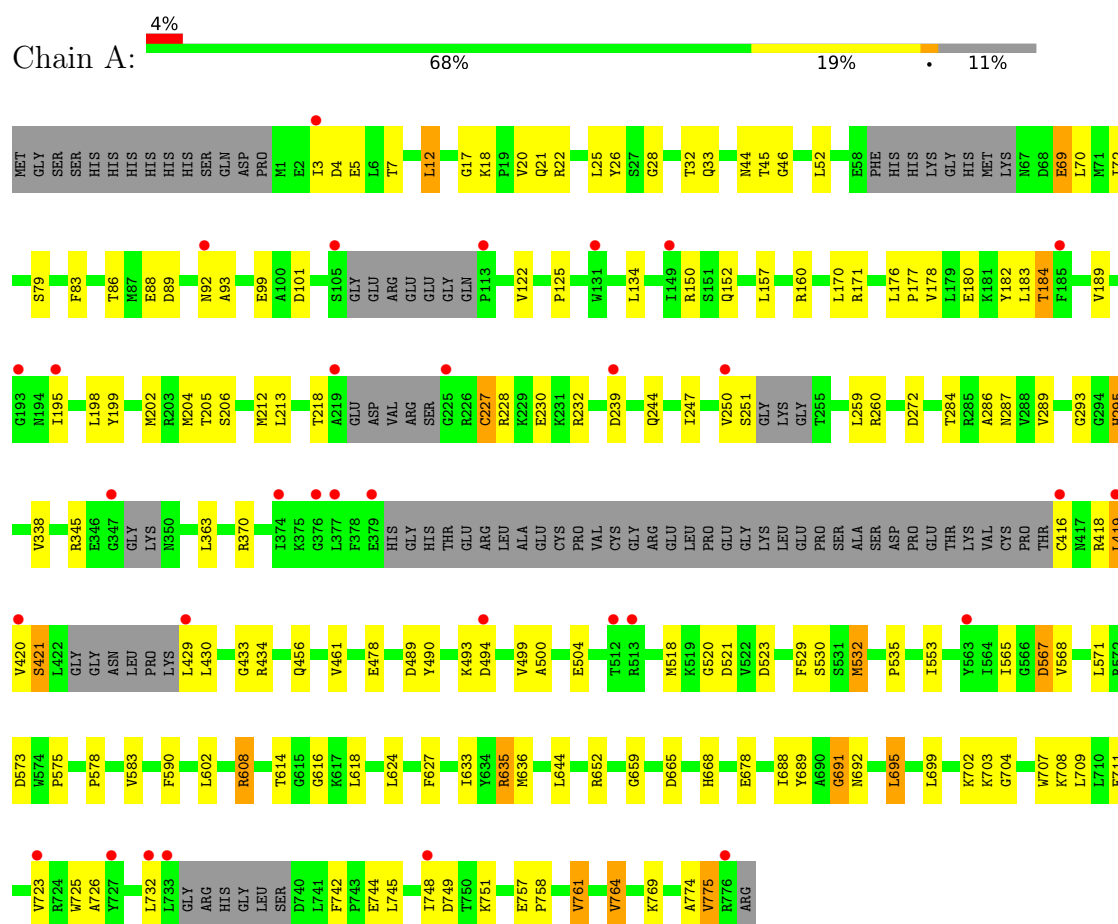
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	21	Total 21	O 21	0	0
4	B	7	Total 7	O 7	0	0

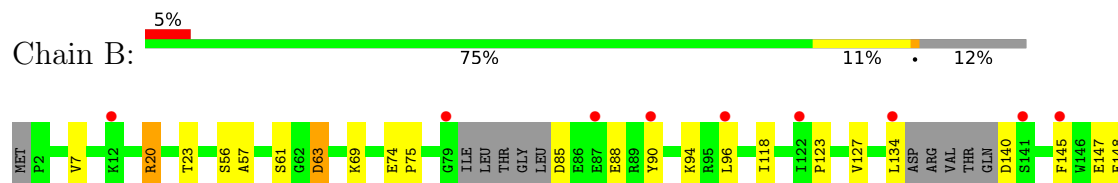
3 Residue-property plots [i](#)

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: CRISPR system single-strand-specific deoxyribonuclease Cas10/Csm1 (subtype III-A)



- Molecule 2: Csm4





● Molecule 3: Cyclic RNA cA4



● Molecule 3: Cyclic RNA cA4



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	156.58Å 156.58Å 187.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.48 – 2.40 49.48 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.48-2.40) 99.4 (49.48-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.23 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.222 , 0.271 0.226 , 0.271	Depositor DCC
R_{free} test set	5165 reflections (9.70%)	wwPDB-VP
Wilson B-factor (Å ²)	67.9	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.6	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	0.95	EDS
Total number of atoms	7884	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.30% 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	0.98	0/5775	1.47	9/7787 (0.1%)
2	B	0.96	0/2078	1.33	0/2811
3	C	2.54	2/99 (2.0%)	3.74	19/152 (12.5%)
3	D	2.37	1/99 (1.0%)	2.51	7/152 (4.6%)
All	All	1.03	3/8051 (0.0%)	1.51	35/10902 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	A	O3'-P	8.22	1.73	1.61
3	C	1	A	C3'-O3'	7.56	1.53	1.42
3	D	3	A	O4'-C1'	-5.04	1.34	1.42

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	A	C1'-O4'-C4'	-15.83	94.07	109.90
3	C	2	A	C1'-O4'-C4'	-11.89	98.01	109.90
3	C	1	A	P-O3'-C3'	11.40	137.30	120.20
3	C	1	A	C3'-C2'-C1'	-10.77	90.53	101.30
3	C	3	A	C1'-O4'-C4'	-10.35	99.55	109.90
3	C	3	A	P-O3'-C3'	9.49	134.44	120.20
3	C	4	A	C1'-O4'-C4'	-9.49	100.21	109.70
3	D	3	A	C1'-O4'-C4'	-9.16	100.54	109.70
3	C	4	A	O4'-C1'-N9	8.56	121.04	108.20
3	C	1	A	O4'-C1'-C2'	-8.35	99.25	107.60
3	C	1	A	N9-C1'-C2'	7.94	123.91	112.00
3	D	2	A	O4'-C1'-N9	-7.59	96.82	108.20
3	C	3	A	O4'-C1'-N9	7.27	119.41	108.50
3	D	4	A	C5'-C4'-C3'	-7.08	104.59	115.20
3	D	1	A	O5'-P-OP1	-6.86	87.42	108.00
3	C	2	A	P-O3'-C3'	6.64	130.15	120.20
3	C	3	A	O4'-C4'-C3'	6.48	110.48	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	A	O5'-P-OP2	-6.22	89.34	108.00
3	C	3	A	O5'-C5'-C4'	-6.02	102.48	111.50
1	A	493	LYS	CA-C-N	5.88	132.01	122.48
1	A	493	LYS	C-N-CA	5.88	132.01	122.48
1	A	703	LYS	CA-C-N	5.87	126.50	119.98
1	A	703	LYS	C-N-CA	5.87	126.50	119.98
3	C	3	A	C8-N9-C1'	-5.68	110.66	127.70
3	D	2	A	O4'-C4'-C3'	-5.65	100.45	106.10
3	D	1	A	C4'-C3'-C2'	-5.61	97.00	102.60
3	C	4	A	O4'-C1'-C2'	-5.56	100.24	105.80
1	A	79	SER	CA-C-N	5.32	125.39	119.32
1	A	79	SER	C-N-CA	5.32	125.39	119.32
1	A	691	GLY	CA-C-O	-5.31	118.50	122.37
1	A	616	GLY	N-CA-C	-5.25	107.55	115.32
3	C	3	A	C5'-C4'-C3'	-5.25	108.12	116.00
3	C	2	A	O5'-C5'-C4'	-5.25	103.63	111.50
1	A	583	VAL	N-CA-C	-5.10	106.71	111.45
3	C	2	A	O4'-C1'-C2'	-5.01	102.59	107.60

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	5653	0	5663	102	0
2	B	2027	0	2025	23	0
3	C	88	0	44	2	0
3	D	88	0	44	19	0
4	A	21	0	0	0	0
4	B	7	0	0	0	0
All	All	7884	0	7776	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 9.

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)
2:B:90:TYR:HD1	2:B:94:LYS:NZ	1.45	1.13
1:A:244:GLN:HG3	3:D:1:A:N7	1.64	1.12
1:A:726:ALA:CB	1:A:748:ILE:HD11	1.98	0.93
2:B:90:TYR:HD1	2:B:94:LYS:HZ1	1.15	0.93
1:A:494:ASP:OD2	1:A:500:ALA:HA	1.68	0.92
2:B:90:TYR:CD1	2:B:94:LYS:NZ	2.38	0.90
1:A:494:ASP:OD1	1:A:500:ALA:HB2	1.74	0.88
1:A:726:ALA:HB2	1:A:748:ILE:HD11	1.64	0.79
2:B:90:TYR:HD1	2:B:94:LYS:HZ3	1.34	0.75
1:A:250:VAL:HG12	1:A:251:SER:N	2.02	0.73
1:A:726:ALA:HB1	1:A:748:ILE:HD11	1.71	0.72
1:A:244:GLN:HG3	3:D:1:A:C8	2.25	0.72
1:A:183:LEU:HD12	1:A:202:MET:HE1	1.74	0.70
2:B:85:ASP:HB2	2:B:88:GLU:OE2	1.92	0.69
1:A:523:ASP:CG	1:A:652:ARG:HB3	2.21	0.66
2:B:85:ASP:HB3	2:B:88:GLU:HB2	1.78	0.65
1:A:284:THR:H	1:A:287:ASN:HD22	1.46	0.63
1:A:627:PHE:HB2	1:A:636:MET:CE	2.29	0.63
1:A:627:PHE:HB2	1:A:636:MET:HE2	1.81	0.62
1:A:571:LEU:HD22	1:A:774:ALA:HB2	1.82	0.62
2:B:173:ALA:HB3	2:B:174:PRO:HD3	1.80	0.61
3:D:1:A:C3'	3:D:2:A:H5'	2.30	0.61
1:A:239:ASP:OD1	1:A:295:HIS:CD2	2.54	0.60
2:B:148:GLU:OE1	2:B:150:ARG:NH2	2.35	0.59
1:A:204:MET:HE1	1:A:289:VAL:HG22	1.84	0.59
2:B:57:ALA:HB1	2:B:157:VAL:HG13	1.85	0.59
1:A:250:VAL:CG1	1:A:251:SER:N	2.65	0.59
2:B:63:ASP:N	2:B:63:ASP:OD1	2.36	0.58
1:A:180:GLU:O	1:A:184:THR:CG2	2.51	0.58
3:D:2:A:H5''	3:D:2:A:N3	2.19	0.58
1:A:295:HIS:CD2	3:D:3:A:H4'	2.41	0.56
3:D:1:A:H4'	3:D:2:A:H5'	1.88	0.56
1:A:26:TYR:HB2	1:A:33:GLN:HG3	1.88	0.56
2:B:214:THR:HG21	2:B:277:THR:HG22	1.87	0.55
1:A:260:ARG:NH2	1:A:489:ASP:O	2.39	0.55
1:A:250:VAL:HG12	1:A:251:SER:H	1.70	0.55
2:B:20:ARG:HA	2:B:56:SER:HB3	1.90	0.54
1:A:635:ARG:HG3	2:B:145:PHE:CE2	2.44	0.53
1:A:250:VAL:CG1	1:A:251:SER:H	2.21	0.53
1:A:101:ASP:HA	1:A:205:THR:HG22	1.89	0.53
1:A:726:ALA:HB1	1:A:748:ILE:CD1	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:LEU:HD22	1:A:83:PHE:CD2	2.44	0.52
1:A:518:MET:HB3	1:A:624:LEU:HD23	1.91	0.52
1:A:695:LEU:HD13	1:A:775:VAL:HG21	1.89	0.52
3:D:1:A:C3'	3:D:2:A:C5'	2.88	0.52
1:A:520:GLY:HA3	1:A:590:PHE:CZ	2.45	0.52
1:A:518:MET:HE2	1:A:602:LEU:HD23	1.91	0.52
2:B:74:GLU:HB2	2:B:75:PRO:HD3	1.92	0.52
2:B:127:VAL:HG22	3:C:2:A:H2'	1.92	0.52
1:A:688:ILE:HA	1:A:699:LEU:CD1	2.38	0.52
3:D:2:A:N3	3:D:2:A:C5'	2.72	0.52
1:A:284:THR:HG23	1:A:286:ALA:HB3	1.92	0.52
3:D:2:A:H4'	3:D:3:A:OP2	2.08	0.52
1:A:12:LEU:HD13	1:A:206:SER:CB	2.41	0.51
1:A:199:TYR:HA	1:A:202:MET:HE2	1.93	0.51
1:A:284:THR:H	1:A:287:ASN:ND2	2.08	0.51
1:A:125:PRO:HG3	1:A:195:ILE:HD11	1.92	0.51
1:A:494:ASP:OD1	1:A:500:ALA:CB	2.52	0.50
1:A:150:ARG:C	1:A:152:GLN:H	2.20	0.50
1:A:659:GLY:O	1:A:769:LYS:NZ	2.37	0.50
1:A:707:TRP:O	1:A:711:GLU:HG3	2.12	0.50
3:D:1:A:C4'	3:D:2:A:H5'	2.42	0.50
1:A:69:GLU:HA	1:A:72:ILE:HD12	1.94	0.50
1:A:247:ILE:CD1	3:D:4:A:C2	2.94	0.50
1:A:430:LEU:HD11	1:A:456:GLN:HA	1.93	0.49
1:A:573:ASP:H	1:A:692:ASN:HD21	1.61	0.49
1:A:26:TYR:H	1:A:33:GLN:NE2	2.10	0.49
1:A:44:ASN:O	1:A:46:GLY:N	2.44	0.49
1:A:204:MET:HE2	1:A:535:PRO:HB3	1.93	0.49
1:A:17:GLY:O	1:A:20:VAL:HG22	2.13	0.48
1:A:293:GLY:O	3:D:3:A:O2'	2.31	0.48
1:A:134:LEU:HD12	1:A:184:THR:HG21	1.95	0.48
1:A:122:VAL:HB	1:A:195:ILE:HG23	1.93	0.48
1:A:518:MET:HE1	1:A:602:LEU:HG	1.95	0.48
1:A:529:PHE:O	1:A:532:MET:HB2	2.13	0.48
1:A:176:LEU:HB2	1:A:177:PRO:HD3	1.95	0.47
1:A:521:ASP:OD2	3:D:2:A:O2'	2.24	0.47
1:A:416:CYS:HA	1:A:419:LEU:HB2	1.96	0.47
1:A:745:LEU:HD11	1:A:764:VAL:HG21	1.95	0.47
1:A:295:HIS:NE2	3:D:3:A:H4'	2.30	0.47
1:A:575:PRO:HG2	1:A:578:PRO:HA	1.96	0.47
1:A:494:ASP:OD2	1:A:504:GLU:OE2	2.33	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:LEU:CD1	1:A:182:TYR:HB3	2.44	0.47
1:A:180:GLU:O	1:A:184:THR:HG22	2.14	0.46
1:A:668:HIS:CE1	1:A:758:PRO:HB3	2.50	0.46
1:A:689:TYR:CE2	1:A:691:GLY:HA2	2.50	0.46
1:A:227:CYS:HA	1:A:232:ARG:HG3	1.97	0.46
1:A:494:ASP:CG	1:A:500:ALA:HA	2.38	0.46
1:A:627:PHE:CB	1:A:636:MET:CE	2.94	0.46
1:A:26:TYR:HD2	1:A:33:GLN:HG3	1.80	0.46
2:B:61:SER:HB2	2:B:123:PRO:HB3	1.98	0.46
1:A:134:LEU:C	1:A:134:LEU:HD23	2.41	0.45
1:A:567:ASP:O	1:A:608:ARG:NH2	2.47	0.45
1:A:709:LEU:HD11	1:A:742:PHE:CD2	2.52	0.45
2:B:85:ASP:CB	2:B:88:GLU:HB2	2.45	0.45
3:D:2:A:OP1	3:D:2:A:C2	2.70	0.45
1:A:18:LYS:O	1:A:22:ARG:HG3	2.16	0.45
1:A:749:ASP:OD1	1:A:751:LYS:HB2	2.16	0.45
1:A:12:LEU:HD13	1:A:206:SER:HB2	1.99	0.45
1:A:272:ASP:OD1	1:A:434:ARG:NH1	2.46	0.45
3:D:2:A:OP1	3:D:2:A:H2	2.00	0.45
1:A:702:LYS:C	1:A:704:GLY:N	2.75	0.44
3:D:1:A:H3'	3:D:2:A:C5'	2.47	0.44
1:A:171:ARG:NH2	1:A:478:GLU:O	2.51	0.44
3:C:3:A:H8	3:C:4:A:OP2	2.01	0.44
1:A:284:THR:HG22	1:A:287:ASN:ND2	2.33	0.44
1:A:180:GLU:O	1:A:184:THR:HG23	2.17	0.44
1:A:433:GLY:HA3	1:A:461:VAL:O	2.19	0.43
1:A:26:TYR:O	1:A:28:GLY:N	2.46	0.43
2:B:85:ASP:CB	2:B:88:GLU:OE2	2.63	0.43
1:A:3:ILE:O	1:A:7:THR:HG23	2.19	0.43
1:A:20:VAL:HG23	1:A:33:GLN:HG2	2.01	0.43
2:B:214:THR:CG2	2:B:277:THR:HB	2.49	0.43
1:A:608:ARG:NH1	1:A:678:GLU:OE1	2.52	0.43
2:B:69:LYS:NZ	2:B:96:LEU:O	2.51	0.42
1:A:363:LEU:HD23	1:A:363:LEU:HA	1.94	0.42
2:B:96:LEU:HD23	2:B:118:ILE:CG1	2.49	0.42
2:B:204:ILE:HG21	2:B:280:LEU:HD13	2.01	0.42
1:A:490:TYR:OH	1:A:578:PRO:HD2	2.19	0.42
1:A:198:LEU:O	1:A:202:MET:HG3	2.20	0.42
1:A:99:GLU:HG2	1:A:212:MET:HE1	2.01	0.41
1:A:3:ILE:HD12	1:A:213:LEU:HD11	2.02	0.41
1:A:565:ILE:O	1:A:608:ARG:HD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:627:PHE:CB	1:A:636:MET:HE3	2.50	0.41
1:A:571:LEU:HD22	1:A:774:ALA:CB	2.48	0.41
3:D:2:A:O4'	3:D:3:A:P	2.78	0.41
1:A:3:ILE:HG22	1:A:89:ASP:HB3	2.00	0.41
3:D:2:A:C4'	3:D:3:A:OP2	2.69	0.41
1:A:25:LEU:HD23	1:A:25:LEU:HA	1.90	0.41
1:A:418:ARG:O	1:A:421:SER:OG	2.38	0.41
1:A:614:THR:HG21	1:A:618:LEU:HB2	2.02	0.41
2:B:7:VAL:O	2:B:158:TYR:HA	2.20	0.41
1:A:521:ASP:HB2	1:A:644:LEU:HD21	2.03	0.41
1:A:92:ASN:OD1	1:A:218:THR:HA	2.21	0.40
1:A:665:ASP:OD1	1:A:668:HIS:HD2	2.04	0.40
1:A:7:THR:HG22	1:A:93:ALA:HA	2.04	0.40
1:A:725:TRP:CZ3	1:A:761:VAL:HG21	2.57	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	685/791 (87%)	652 (95%)	33 (5%)	0	100	100
2	B	244/289 (84%)	231 (95%)	13 (5%)	0	100	100
All	All	929/1080 (86%)	883 (95%)	46 (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	592/664 (89%)	547 (92%)	45 (8%)	12	21
2	B	215/240 (90%)	203 (94%)	12 (6%)	19	33
All	All	807/904 (89%)	750 (93%)	57 (7%)	13	23

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ASP
1	A	5	GLU
1	A	12	LEU
1	A	21	GLN
1	A	32	THR
1	A	45	THR
1	A	69	GLU
1	A	70	LEU
1	A	86	THR
1	A	88	GLU
1	A	160	ARG
1	A	170	LEU
1	A	178	VAL
1	A	184	THR
1	A	189	VAL
1	A	227	CYS
1	A	228	ARG
1	A	230	GLU
1	A	259	LEU
1	A	295	HIS
1	A	338	VAL
1	A	345	ARG
1	A	370	ARG
1	A	419	LEU
1	A	420	VAL
1	A	421	SER
1	A	429	LEU
1	A	499	VAL
1	A	530	SER
1	A	532	MET
1	A	553	ILE
1	A	567	ASP
1	A	568	VAL

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Mol	Chain	Res	Type
1	A	608	ARG
1	A	633	ILE
1	A	635	ARG
1	A	695	LEU
1	A	708	LYS
1	A	723	VAL
1	A	732	LEU
1	A	744	GLU
1	A	757	GLU
1	A	761	VAL
1	A	764	VAL
1	A	775	VAL
2	B	20	ARG
2	B	23	THR
2	B	63	ASP
2	B	134	LEU
2	B	140	ASP
2	B	147	GLU
2	B	152	ARG
2	B	157	VAL
2	B	194	LEU
2	B	213	VAL
2	B	214	THR
2	B	244	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	33	GLN
1	A	287	ASN
1	A	472	ASN
1	A	668	HIS
1	A	675	ASN
1	A	676	HIS
1	A	692	ASN
2	B	30	ASN
2	B	38	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	C	4/4 (100%)	3 (75%)	2 (50%)
3	D	4/4 (100%)	3 (75%)	3 (75%)
All	All	8/8 (100%)	6 (75%)	5 (62%)

All (6) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	C	2	A
3	C	3	A
3	C	4	A
3	D	2	A
3	D	3	A
3	D	4	A

All (5) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	C	1	A
3	C	3	A
3	D	1	A
3	D	2	A
3	D	3	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 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 ⓘ

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	703/791 (88%)	0.45	32 (4%) 37 33	44, 70, 116, 147	0
2	B	254/289 (87%)	0.48	14 (5%) 30 27	48, 70, 123, 164	0
3	C	4/4 (100%)	1.11	1 (25%) 2 1	78, 80, 87, 109	0
3	D	4/4 (100%)	0.96	1 (25%) 2 1	77, 92, 149, 164	0
All	All	965/1088 (88%)	0.46	48 (4%) 34 30	44, 70, 117, 164	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	134	LEU	6.4
1	A	219	ALA	5.9
1	A	225	GLY	5.8
1	A	347	GLY	5.6
1	A	113	PRO	5.3
1	A	92	ASN	5.0
2	B	79	GLY	4.3
1	A	733	LEU	4.1
1	A	250	VAL	4.0
1	A	494	ASP	3.9
2	B	287	GLY	3.8
1	A	420	VAL	3.1
1	A	419	LEU	3.0
2	B	90	TYR	2.9
1	A	732	LEU	2.8
2	B	96	LEU	2.8
1	A	3	ILE	2.6
1	A	429	LEU	2.6
1	A	727	TYR	2.6
2	B	149	ILE	2.5
2	B	122	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	776	ARG	2.5
1	A	195	ILE	2.4
1	A	563	TYR	2.4
2	B	250	GLY	2.4
1	A	748	ILE	2.3
3	D	2	A	2.3
1	A	513	ARG	2.3
1	A	374	ILE	2.3
2	B	231	ARG	2.3
2	B	141	SER	2.2
1	A	239	ASP	2.2
1	A	377	LEU	2.2
1	A	416	CYS	2.2
2	B	145	PHE	2.2
1	A	149	ILE	2.2
1	A	379	GLU	2.2
1	A	131	TRP	2.2
1	A	105	SER	2.1
2	B	12	LYS	2.1
1	A	723	VAL	2.1
1	A	185	PHE	2.1
1	A	376	GLY	2.1
2	B	87	GLU	2.1
1	A	512	THR	2.1
1	A	193	GLY	2.0
3	C	2	A	2.0
2	B	181	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.