



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 11:47 PM BST

PDB ID : 6O78
Title : Crystal structure of Csm1-Csm4 cassette in complex with pppApApA
Authors : Jia, N.; Patel, D.J.
Deposited on : 2019-03-07
Resolution : 2.80 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 : 2.13.1
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) : 2.13.1

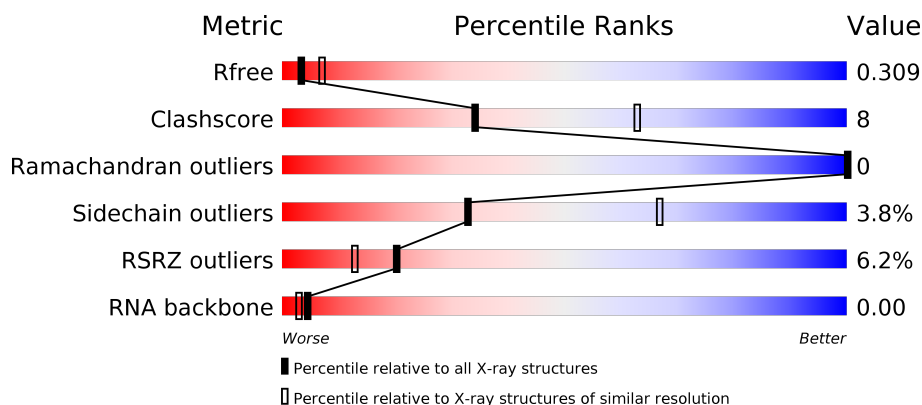
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)
RNA backbone	3102	1227 (3.10-2.50)

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>5%</div> <div> <div></div> <div>78%</div> <div>10%</div> <div>•</div> <div>12%</div> </div> </div>
2	B	289	<div> <div>6%</div> <div> <div></div> <div>67%</div> <div>16%</div> <div>•</div> <div>16%</div> </div> </div>
3	E	3	<div> <div></div> <div> <div>67%</div> <div>33%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7660 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	700	Total	C	N	O	S	0	0	0
			5631	3627	965	1024	15			

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	244	Total	C	N	O	S	0	0	0
			1952	1274	322	352	4			

- Molecule 3 is a RNA chain called RNA (5'-D*(ATP))-R(P*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	3	Total	C	N	O	P	0	0	0
			75	30	15	25	5			

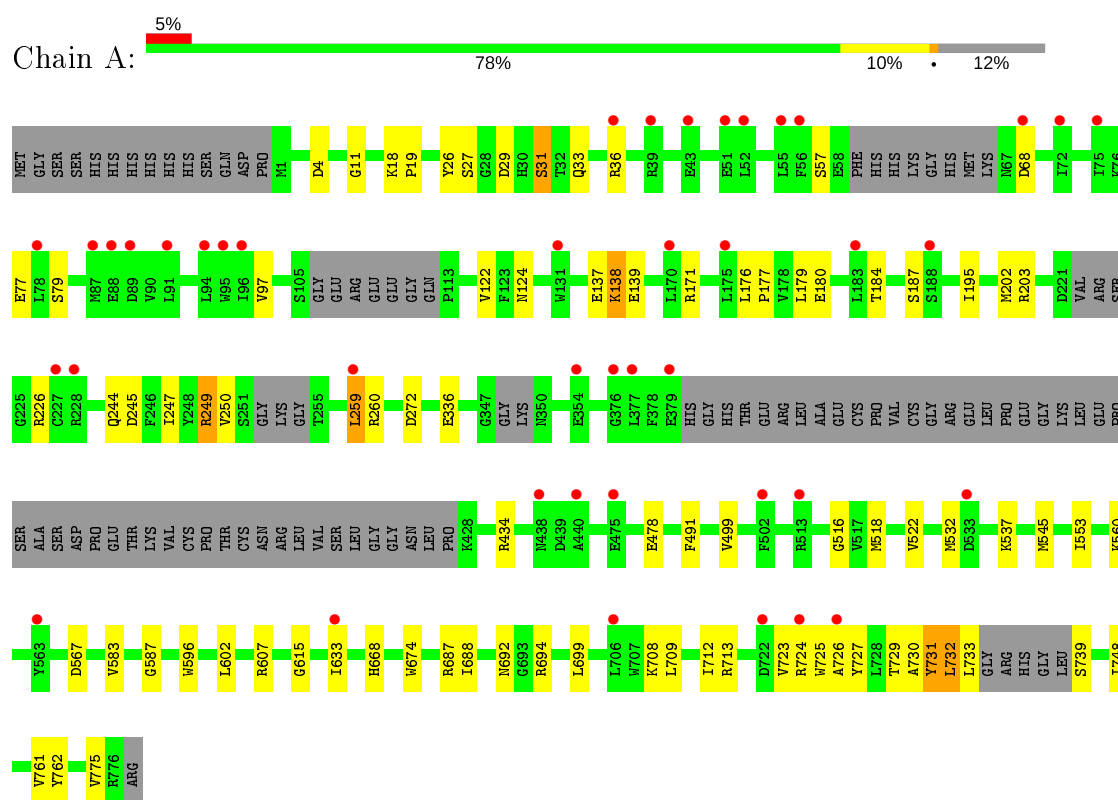
- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mn	0	0
			2	2		

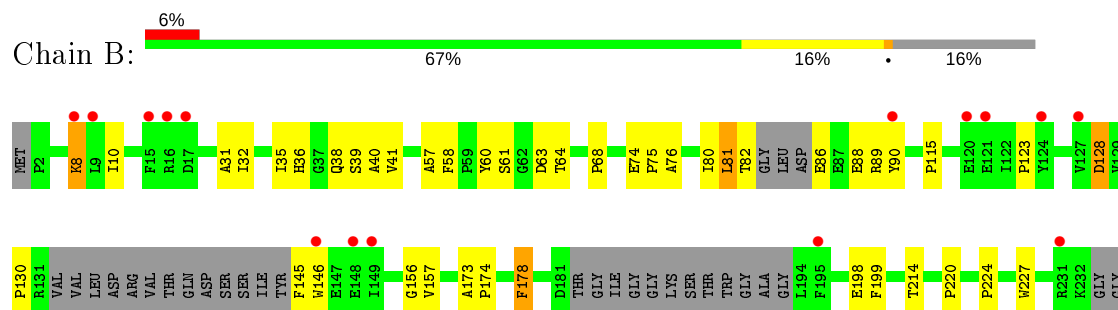
3 Residue-property plots [i](#)

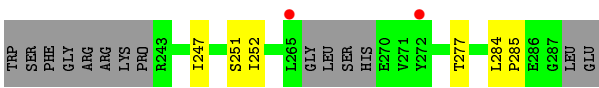
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: RNA (5'-D(* (ATP))-R(P*AP*A)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	154.18Å 154.18Å 182.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.69 – 2.80 48.64 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.69-2.80) 100.0 (48.64-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.257 , 0.304 0.258 , 0.309	Depositor DCC
R_{free} test set	1559 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	97.5	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 63.4	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	0.93	EDS
Total number of atoms	7660	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.66% 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: MN, ATP

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	0.65	0/5754	0.73	0/7759
2	B	0.63	0/2000	0.71	0/2703
3	E	4.09	14/49 (28.6%)	6.23	22/74 (29.7%)
All	All	0.72	14/7803 (0.2%)	0.89	22/10536 (0.2%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	3	A	N7-C5	-9.61	1.33	1.39
3	E	2	A	N7-C5	-9.55	1.33	1.39
3	E	2	A	C2-N3	8.01	1.40	1.33
3	E	2	A	C5-C6	-7.69	1.34	1.41
3	E	3	A	C5-C6	-7.47	1.34	1.41
3	E	3	A	C2-N3	7.41	1.40	1.33
3	E	2	A	N9-C4	-7.38	1.33	1.37
3	E	3	A	N9-C4	-7.36	1.33	1.37
3	E	3	A	N1-C2	6.89	1.40	1.34
3	E	2	A	N1-C2	6.87	1.40	1.34
3	E	2	A	C5-C4	-6.28	1.34	1.38
3	E	3	A	C5-C4	-5.92	1.34	1.38
3	E	3	A	N9-C8	-5.56	1.33	1.37
3	E	2	A	N9-C8	-5.40	1.33	1.37

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	2	A	N1-C2-N3	-21.91	118.34	129.30
3	E	3	A	N1-C2-N3	-20.99	118.81	129.30
3	E	2	A	C2-N3-C4	17.88	119.54	110.60
3	E	3	A	C2-N3-C4	17.25	119.22	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	2	A	N7-C8-N9	-11.47	108.07	113.80
3	E	3	A	N7-C8-N9	-10.74	108.43	113.80
3	E	2	A	C5-C6-N6	-9.15	116.38	123.70
3	E	2	A	C5-C6-N1	8.76	122.08	117.70
3	E	2	A	C5-N7-C8	8.49	108.14	103.90
3	E	2	A	N3-C4-C5	-8.47	120.87	126.80
3	E	3	A	N3-C4-C5	-8.46	120.88	126.80
3	E	3	A	C5-N7-C8	8.37	108.08	103.90
3	E	3	A	C5-C6-N6	-8.15	117.18	123.70
3	E	3	A	C5-C6-N1	8.12	121.76	117.70
3	E	3	A	C1'-O4'-C4'	-8.02	103.48	109.90
3	E	3	A	C4-C5-C6	7.50	120.75	117.00
3	E	2	A	C4-C5-C6	7.21	120.61	117.00
3	E	3	A	C4-C5-N7	-6.35	107.53	110.70
3	E	2	A	C8-N9-C4	6.20	108.28	105.80
3	E	3	A	N9-C4-C5	5.97	108.19	105.80
3	E	2	A	C4-C5-N7	-5.61	107.89	110.70
3	E	2	A	N3-C4-N9	5.15	131.52	127.40

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	5631	0	5632	49	0
2	B	1952	0	1960	80	0
3	E	75	0	34	1	0
4	A	2	0	0	0	0
All	All	7660	0	7626	129	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 8.

All (129) 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:8:LYS:CE	2:B:198:GLU:HB2	1.44	1.42
2:B:8:LYS:HE3	2:B:198:GLU:CB	1.62	1.30
2:B:60:TYR:HB3	2:B:156:GLY:H	1.06	1.15
2:B:8:LYS:HE2	2:B:198:GLU:HB2	1.34	1.03
2:B:60:TYR:HB3	2:B:156:GLY:N	1.75	1.00
2:B:8:LYS:HE3	2:B:198:GLU:HB2	1.01	0.99
2:B:8:LYS:HG3	2:B:198:GLU:O	1.62	0.99
2:B:81:LEU:HD12	2:B:81:LEU:H	1.25	0.97
2:B:60:TYR:HB2	2:B:156:GLY:C	1.85	0.96
2:B:8:LYS:HB2	2:B:60:TYR:CE1	2.01	0.95
2:B:8:LYS:HE3	2:B:198:GLU:CA	1.96	0.95
1:A:250:VAL:HG22	1:A:633:ILE:HD11	1.51	0.91
1:A:26:TYR:HB2	1:A:33:GLN:HG3	1.53	0.91
2:B:60:TYR:CB	2:B:156:GLY:C	2.44	0.85
2:B:31:ALA:O	2:B:35:ILE:HG12	1.77	0.84
2:B:8:LYS:CE	2:B:198:GLU:CB	2.32	0.84
2:B:60:TYR:CB	2:B:156:GLY:O	2.29	0.81
2:B:80:ILE:HG21	2:B:115:PRO:HA	1.63	0.79
2:B:60:TYR:HB2	2:B:156:GLY:O	1.82	0.79
2:B:8:LYS:HE3	2:B:198:GLU:N	1.97	0.79
2:B:86:GLU:HA	2:B:89:ARG:HG2	1.64	0.78
2:B:32:ILE:HD13	2:B:35:ILE:HD11	1.66	0.76
2:B:35:ILE:CG2	2:B:178:PHE:CE1	2.71	0.74
2:B:35:ILE:CG2	2:B:178:PHE:CD1	2.72	0.72
2:B:8:LYS:O	2:B:8:LYS:HD2	1.88	0.72
2:B:60:TYR:N	2:B:156:GLY:O	2.24	0.71
1:A:725:TRP:O	1:A:729:THR:HG23	1.91	0.70
1:A:250:VAL:CG2	1:A:633:ILE:HD11	2.21	0.69
2:B:8:LYS:HB2	2:B:60:TYR:HE1	1.52	0.69
2:B:80:ILE:CG2	2:B:115:PRO:HA	2.24	0.68
2:B:8:LYS:HZ3	2:B:10:ILE:HG13	1.58	0.67
2:B:224:PRO:HG2	2:B:227:TRP:CE3	2.30	0.67
2:B:35:ILE:HG21	2:B:178:PHE:CD1	2.29	0.66
2:B:61:SER:HB2	2:B:123:PRO:HG2	1.78	0.66
2:B:60:TYR:HB3	2:B:156:GLY:CA	2.26	0.66
1:A:138:LYS:HD2	1:A:138:LYS:O	1.97	0.65
1:A:775:VAL:O	1:A:775:VAL:HG12	1.96	0.65
2:B:8:LYS:CB	2:B:60:TYR:HE1	2.10	0.65
2:B:86:GLU:CA	2:B:89:ARG:HG2	2.28	0.64
2:B:61:SER:HB2	2:B:123:PRO:CG	2.28	0.64
2:B:8:LYS:NZ	2:B:10:ILE:CG1	2.60	0.63
2:B:60:TYR:HB3	2:B:156:GLY:C	2.19	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:8:LYS:HZ2	2:B:10:ILE:HG12	1.64	0.62
2:B:35:ILE:HG22	2:B:178:PHE:CE1	2.35	0.62
2:B:8:LYS:NZ	2:B:10:ILE:HG12	2.14	0.61
2:B:8:LYS:HB2	2:B:60:TYR:CZ	2.35	0.61
2:B:86:GLU:O	2:B:90:TYR:N	2.34	0.61
1:A:139:GLU:OE1	1:A:139:GLU:HA	1.99	0.60
1:A:723:VAL:HG12	1:A:726:ALA:HB3	1.82	0.60
1:A:522:VAL:HG21	1:A:545:MET:HE1	1.84	0.60
1:A:712:ILE:HD11	1:A:732:LEU:HG	1.83	0.59
1:A:137:GLU:O	1:A:560:LYS:NZ	2.36	0.58
2:B:81:LEU:N	2:B:81:LEU:HD12	2.04	0.57
2:B:8:LYS:NZ	2:B:10:ILE:HG13	2.20	0.57
2:B:8:LYS:CB	2:B:60:TYR:CE1	2.81	0.57
1:A:723:VAL:O	1:A:724:ARG:C	2.43	0.57
1:A:491:PHE:HB2	1:A:499:VAL:HG23	1.87	0.57
1:A:723:VAL:HG11	1:A:748:ILE:HD11	1.86	0.57
2:B:38:GLN:O	2:B:41:VAL:N	2.33	0.57
1:A:272:ASP:OD1	1:A:434:ARG:NH1	2.38	0.56
1:A:712:ILE:HG21	1:A:729:THR:HG22	1.86	0.56
1:A:124:ASN:HB2	1:A:615:GLY:HA3	1.87	0.56
2:B:8:LYS:HZ3	2:B:10:ILE:CG1	2.19	0.55
1:A:733:LEU:HB3	1:A:739:SER:HA	1.89	0.55
1:A:692:ASN:O	1:A:694:ARG:NH1	2.40	0.54
1:A:171:ARG:NH2	1:A:478:GLU:O	2.41	0.54
1:A:26:TYR:CE2	1:A:36:ARG:O	2.61	0.54
2:B:32:ILE:HA	2:B:35:ILE:CG1	2.38	0.53
2:B:130:PRO:HA	2:B:145:PHE:CB	2.40	0.52
1:A:31:SER:HB3	1:A:57:SER:O	2.10	0.51
1:A:731:TYR:CD1	1:A:731:TYR:C	2.84	0.51
1:A:26:TYR:H	1:A:33:GLN:NE2	2.09	0.50
2:B:252:ILE:HG21	2:B:284:LEU:HD21	1.93	0.50
2:B:178:PHE:CD2	2:B:178:PHE:C	2.86	0.49
1:A:688:ILE:HA	1:A:699:LEU:CD1	2.42	0.49
1:A:712:ILE:HG21	1:A:729:THR:CG2	2.42	0.49
2:B:224:PRO:HG3	2:B:247:ILE:HD11	1.93	0.49
2:B:247:ILE:CG2	2:B:251:SER:HB2	2.43	0.49
1:A:775:VAL:O	1:A:775:VAL:CG1	2.61	0.48
2:B:74:GLU:HB2	2:B:75:PRO:HD3	1.95	0.48
2:B:57:ALA:HB1	2:B:157:VAL:CG1	2.44	0.48
2:B:76:ALA:HB2	2:B:285:PRO:HG3	1.96	0.48
2:B:146:TRP:HA	2:B:146:TRP:CE3	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:81:LEU:CD1	2:B:81:LEU:H	2.05	0.48
1:A:179:LEU:HD22	1:A:202:MET:HE1	1.95	0.48
1:A:245:ASP:O	1:A:249:ARG:HB2	2.14	0.47
1:A:587:GLY:C	3:E:1:ATP:HN61	2.18	0.47
2:B:38:GLN:O	2:B:39:SER:C	2.53	0.47
2:B:8:LYS:HE3	2:B:198:GLU:H	1.77	0.46
1:A:708:LYS:O	1:A:712:ILE:HG13	2.15	0.46
2:B:8:LYS:HZ2	2:B:10:ILE:CG1	2.26	0.46
2:B:38:GLN:O	2:B:40:ALA:N	2.49	0.46
1:A:723:VAL:CG1	1:A:726:ALA:HB3	2.46	0.45
2:B:63:ASP:OD1	2:B:63:ASP:N	2.37	0.45
1:A:26:TYR:HB2	1:A:33:GLN:CG	2.36	0.45
2:B:32:ILE:HA	2:B:35:ILE:HG12	1.97	0.45
2:B:173:ALA:HB3	2:B:174:PRO:HD3	1.99	0.45
2:B:214:THR:HG22	2:B:277:THR:HB	1.99	0.45
2:B:57:ALA:HB1	2:B:157:VAL:HG13	2.00	0.44
1:A:11:GLY:HA2	1:A:97:VAL:HG13	1.99	0.44
1:A:516:GLY:HA3	1:A:596:TRP:HA	2.00	0.44
2:B:80:ILE:HG22	2:B:80:ILE:O	2.17	0.44
1:A:713:ARG:HG3	1:A:761:VAL:HG12	2.00	0.44
1:A:607:ARG:HG3	1:A:674:TRP:CE2	2.52	0.43
2:B:220:PRO:HB3	2:B:247:ILE:HD12	2.00	0.43
1:A:180:GLU:O	1:A:184:THR:HB	2.19	0.43
2:B:60:TYR:HB2	2:B:157:VAL:N	2.30	0.43
2:B:58:PHE:CE1	2:B:68:PRO:HD3	2.53	0.43
1:A:176:LEU:N	1:A:177:PRO:CD	2.82	0.42
1:A:532:MET:HE3	1:A:537:LYS:HB3	2.01	0.42
2:B:60:TYR:HB3	2:B:156:GLY:O	2.14	0.42
2:B:247:ILE:HG23	2:B:251:SER:CB	2.49	0.42
1:A:122:VAL:HA	1:A:195:ILE:HG23	2.02	0.42
1:A:259:LEU:HD23	1:A:583:VAL:HG13	2.02	0.41
1:A:18:LYS:N	1:A:19:PRO:CD	2.83	0.41
2:B:130:PRO:HA	2:B:145:PHE:HB3	2.02	0.41
2:B:8:LYS:HZ1	2:B:198:GLU:HG3	1.85	0.41
2:B:247:ILE:CG2	2:B:251:SER:CB	2.98	0.41
1:A:518:MET:HE2	1:A:602:LEU:HD23	2.02	0.41
2:B:35:ILE:HG23	2:B:178:PHE:CE1	2.56	0.41
2:B:60:TYR:CB	2:B:156:GLY:CA	2.94	0.41
2:B:88:GLU:OE1	2:B:88:GLU:HA	2.21	0.41
1:A:668:HIS:HB3	1:A:762:TYR:CZ	2.56	0.41
1:A:727:TYR:O	1:A:730:ALA:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:732:LEU:HA	1:A:732:LEU:HD22	1.92	0.41
2:B:128:ASP:OD1	2:B:128:ASP:N	2.53	0.41
1:A:723:VAL:O	1:A:725:TRP:N	2.54	0.41
2:B:8:LYS:HD2	2:B:8:LYS:C	2.42	0.41
1:A:709:LEU:HA	1:A:712:ILE:HD12	2.03	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	684/791 (86%)	645 (94%)	39 (6%)	0	100	100
2	B	232/289 (80%)	219 (94%)	13 (6%)	0	100	100
All	All	916/1080 (85%)	864 (94%)	52 (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	589/664 (89%)	567 (96%)	22 (4%)	34	68
2	B	206/240 (86%)	198 (96%)	8 (4%)	32	66
All	All	795/904 (88%)	765 (96%)	30 (4%)	33	67

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ASP
1	A	27	SER
1	A	29	ASP
1	A	31	SER
1	A	68	ASP
1	A	77	GLU
1	A	79	SER
1	A	138	LYS
1	A	187	SER
1	A	203	ARG
1	A	226	ARG
1	A	244	GLN
1	A	247	ILE
1	A	249	ARG
1	A	259	LEU
1	A	260	ARG
1	A	336	GLU
1	A	553	ILE
1	A	567	ASP
1	A	687	ARG
1	A	731	TYR
1	A	732	LEU
2	B	8	LYS
2	B	36	HIS
2	B	64	THR
2	B	81	LEU
2	B	82	THR
2	B	128	ASP
2	B	178	PHE
2	B	199	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	287	ASN
1	A	472	ASN
1	A	668	HIS
1	A	675	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	E	1/3 (33%)	1 (100%)	0

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	E	3	A

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 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 ⓘ

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	700/791 (88%)	0.42	42 (6%)	21	14	74, 108, 152, 175	0
2	B	244/289 (84%)	0.44	17 (6%)	16	9	77, 116, 162, 187	0
3	E	2/3 (66%)	0.10	0	100	100	93, 93, 93, 140	0
All	All	946/1083 (87%)	0.42	59 (6%)	20	13	74, 109, 155, 187	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	68	ASP	5.2
1	A	87	MET	4.7
1	A	39	ARG	4.6
2	B	146	TRP	4.5
1	A	91	LEU	4.4
2	B	120	GLU	4.0
1	A	440	ALA	3.8
1	A	227	CYS	3.6
2	B	265	LEU	3.6
1	A	563	TYR	3.6
1	A	56	PHE	3.5
2	B	16	ARG	3.4
1	A	377	LEU	3.3
2	B	8	LYS	3.2
1	A	75	ILE	3.2
1	A	722	ASP	3.1
2	B	17	ASP	3.1
1	A	379	GLU	3.1
2	B	148	GLU	3.0
1	A	96	ILE	3.0
2	B	121	GLU	2.9
1	A	170	LEU	2.8
1	A	51	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	175	LEU	2.8
1	A	706	LEU	2.7
2	B	90	TYR	2.7
1	A	513	ARG	2.7
1	A	354	GLU	2.7
1	A	72	ILE	2.7
1	A	228	ARG	2.7
1	A	36	ARG	2.6
1	A	52	LEU	2.6
1	A	188	SER	2.6
2	B	127	VAL	2.6
1	A	78	LEU	2.5
1	A	88	GLU	2.5
1	A	131	TRP	2.5
1	A	724	ARG	2.5
2	B	124	TYR	2.4
1	A	533	ASP	2.4
1	A	183	LEU	2.4
1	A	95	TRP	2.3
2	B	149	ILE	2.3
1	A	726	ALA	2.3
1	A	259	LEU	2.3
1	A	633	ILE	2.3
2	B	195	PHE	2.2
2	B	272	TYR	2.2
2	B	15	PHE	2.2
2	B	9	LEU	2.2
1	A	376	GLY	2.2
2	B	231	ARG	2.1
1	A	55	LEU	2.1
1	A	43	GLU	2.1
1	A	94	LEU	2.0
1	A	475	GLU	2.0
1	A	502	PHE	2.0
1	A	89	ASP	2.0
1	A	438	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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	MN	A	801	1/1	0.95	0.06	130,130,130,130	0
4	MN	A	802	1/1	0.97	0.08	102,102,102,102	0

6.5 Other polymers [i](#)

There are no such residues in this entry.