



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 08:00 am BST

PDB ID : 6O6X  
Title : Crystal structure of Csm6 W14A/E337A mutant in complex with cA4 by  
cococrystallization  
Authors : Jia, N.; Patel, D.J.  
Deposited on : 2019-03-07  
Resolution : 2.11 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](mailto: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.

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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

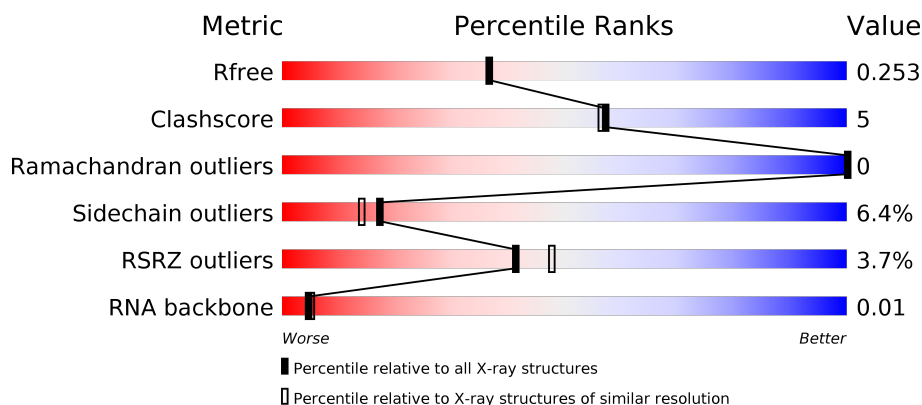
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.11 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	6241 (2.14-2.10)
Clashscore	141614	6778 (2.14-2.10)
Ramachandran outliers	138981	6705 (2.14-2.10)
Sidechain outliers	138945	6706 (2.14-2.10)
RSRZ outliers	127900	6112 (2.14-2.10)
RNA backbone	3102	1013 (2.58-1.66)

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  $\geq 3$ , 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	440	<div> <div>5%</div> <div> <div></div> <div>84%</div> <div>13%</div> <div>..</div> </div> </div>
1	B	440	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>..</div> </div> </div>
2	C	4	<div> <div></div> <div> <div>75%</div> <div>25%</div> </div> </div>
2	D	4	<div> <div></div> <div> <div>75%</div> <div>25%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7126 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 Csm6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	432	Total	C	N	O	S	0	0	0
			3427	2204	582	633	8			
1	B	432	Total	C	N	O	S	0	0	0
			3427	2204	582	633	8			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP B6YWC3
A	0	GLY	-	expression tag	UNP B6YWC3
A	14	ALA	TRP	engineered mutation	UNP B6YWC3
A	337	ALA	GLU	engineered mutation	UNP B6YWC3
A	433	HIS	-	expression tag	UNP B6YWC3
A	434	HIS	-	expression tag	UNP B6YWC3
A	435	HIS	-	expression tag	UNP B6YWC3
A	436	HIS	-	expression tag	UNP B6YWC3
A	437	HIS	-	expression tag	UNP B6YWC3
A	438	HIS	-	expression tag	UNP B6YWC3
B	-1	MET	-	initiating methionine	UNP B6YWC3
B	0	GLY	-	expression tag	UNP B6YWC3
B	14	ALA	TRP	engineered mutation	UNP B6YWC3
B	337	ALA	GLU	engineered mutation	UNP B6YWC3
B	433	HIS	-	expression tag	UNP B6YWC3
B	434	HIS	-	expression tag	UNP B6YWC3
B	435	HIS	-	expression tag	UNP B6YWC3
B	436	HIS	-	expression tag	UNP B6YWC3
B	437	HIS	-	expression tag	UNP B6YWC3
B	438	HIS	-	expression tag	UNP B6YWC3

- Molecule 2 is a RNA chain called Cyclic RNA cA4.

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

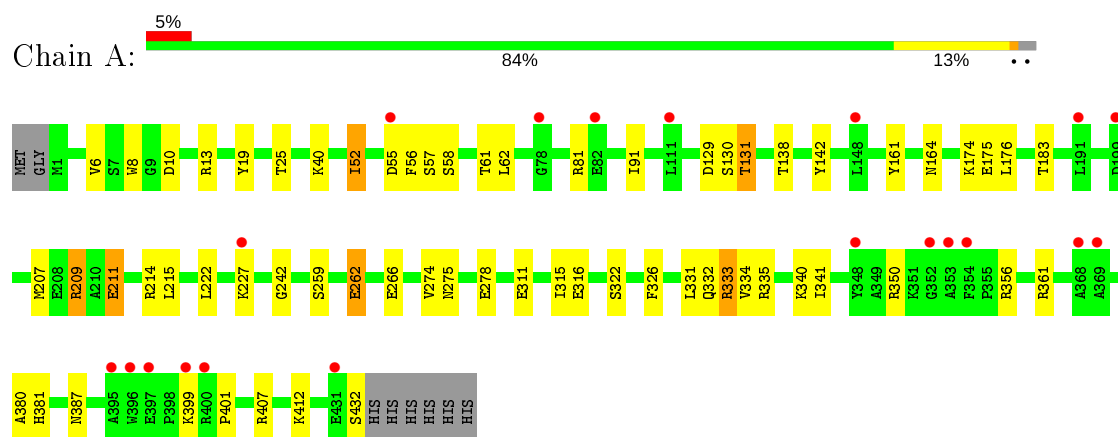
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	38	Total 38	O 38	0	0
3	B	53	Total 53	O 53	0	0
3	C	2	Total 2	O 2	0	0
3	D	3	Total 3	O 3	0	0

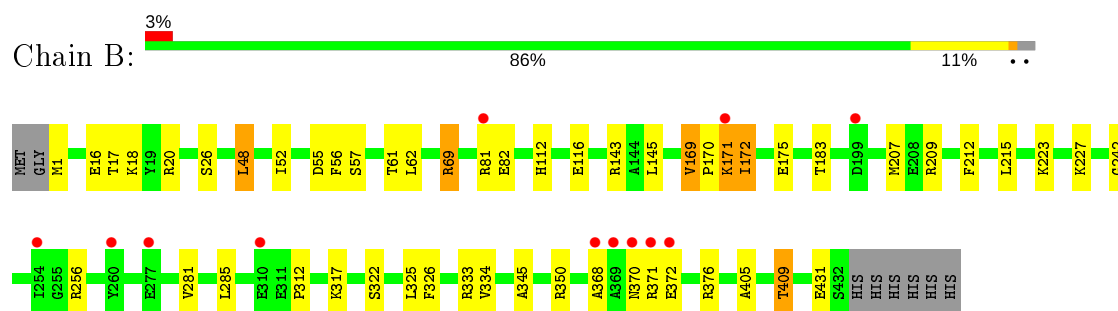
### 3 Residue-property plots [i](#)

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.

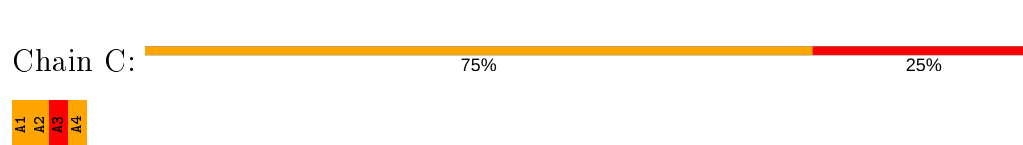
#### • Molecule 1: Csm6



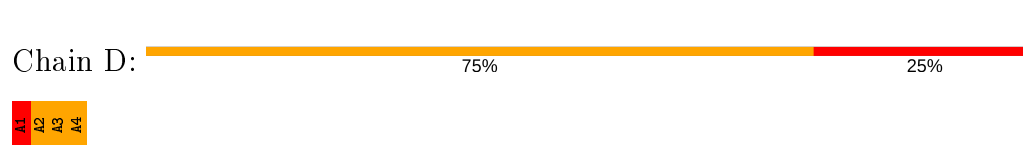
#### • Molecule 1: Csm6



#### • Molecule 2: Cyclic RNA cA4



#### • Molecule 2: Cyclic RNA cA4



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.15Å 115.75Å 161.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.09 – 2.11 49.05 – 2.11	Depositor EDS
% Data completeness (in resolution range)	95.7 (49.09-2.11) 95.7 (49.05-2.11)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.65 (at 2.12Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, $R_{free}$	0.200 , 0.249 0.207 , 0.253	Depositor DCC
$R_{free}$ test set	2838 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.1	Xtriage
Anisotropy	0.566	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 41.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7126	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.70	0/3500	0.84	0/4736
1	B	0.69	0/3500	0.82	0/4736
2	C	4.15	23/99 (23.2%)	6.84	53/152 (34.9%)
2	D	4.05	23/99 (23.2%)	6.39	53/152 (34.9%)
All	All	0.96	46/7198 (0.6%)	1.42	106/9776 (1.1%)

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	A	N7-C5	-10.94	1.32	1.39
2	D	4	A	N7-C5	-10.91	1.32	1.39
2	C	4	A	C2-N3	10.28	1.42	1.33
2	D	3	A	N7-C5	-9.87	1.33	1.39
2	C	1	A	N7-C5	-9.15	1.33	1.39
2	D	2	A	C2-N3	8.96	1.41	1.33
2	D	4	A	N9-C4	-8.67	1.32	1.37
2	D	4	A	C2-N3	8.59	1.41	1.33
2	C	4	A	N7-C5	-8.53	1.34	1.39
2	D	3	A	N1-C2	8.35	1.41	1.34
2	C	2	A	C5-C6	-8.32	1.33	1.41
2	C	2	A	C2-N3	8.22	1.41	1.33
2	D	1	A	N7-C5	-8.15	1.34	1.39
2	D	4	A	N1-C2	8.08	1.41	1.34
2	D	3	A	C2-N3	8.05	1.40	1.33
2	C	4	A	C5-C6	-7.77	1.34	1.41
2	C	4	A	N1-C2	7.46	1.41	1.34
2	C	3	A	N1-C2	7.16	1.40	1.34
2	C	1	A	C2-N3	7.11	1.40	1.33
2	C	3	A	N9-C4	-7.07	1.33	1.37
2	D	1	A	C2-N3	7.05	1.39	1.33
2	C	1	A	C5-C6	-7.05	1.34	1.41
2	C	2	A	N1-C2	7.02	1.40	1.34
2	D	4	A	C5-C6	-6.89	1.34	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	A	N7-C5	-6.82	1.35	1.39
2	C	3	A	C5-C6	-6.80	1.34	1.41
2	D	2	A	N1-C2	6.73	1.40	1.34
2	C	1	A	N1-C2	6.39	1.40	1.34
2	C	4	A	C5-C4	-6.32	1.34	1.38
2	C	3	A	C5-C4	-6.04	1.34	1.38
2	C	3	A	C2-N3	5.98	1.39	1.33
2	C	2	A	O3'-P	5.96	1.68	1.61
2	D	4	A	O4'-C1'	5.96	1.49	1.41
2	D	1	A	C5-C6	-5.68	1.35	1.41
2	D	2	A	P-O5'	5.67	1.65	1.59
2	D	2	A	C5-C4	-5.58	1.34	1.38
2	D	1	A	C5-C4	-5.56	1.34	1.38
2	C	2	A	O4'-C1'	5.53	1.48	1.41
2	C	2	A	N9-C4	-5.51	1.34	1.37
2	D	3	A	C5-C6	-5.47	1.36	1.41
2	D	3	A	C1'-N9	5.46	1.56	1.48
2	D	3	A	O3'-P	5.42	1.67	1.61
2	D	1	A	N9-C4	-5.41	1.34	1.37
2	D	2	A	N9-C8	-5.39	1.33	1.37
2	C	2	A	C5-C4	-5.15	1.35	1.38
2	D	4	A	C8-N7	5.13	1.35	1.31

All (106) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	A	N1-C2-N3	-24.87	116.86	129.30
2	C	4	A	N1-C2-N3	-23.27	117.66	129.30
2	C	1	A	N1-C2-N3	-23.00	117.80	129.30
2	C	2	A	N1-C2-N3	-21.44	118.58	129.30
2	D	1	A	C2-N3-C4	20.86	121.03	110.60
2	D	1	A	N1-C2-N3	-20.58	119.01	129.30
2	C	1	A	C2-N3-C4	20.26	120.73	110.60
2	C	4	A	C2-N3-C4	20.07	120.64	110.60
2	D	3	A	N1-C2-N3	-19.47	119.57	129.30
2	D	2	A	C2-N3-C4	18.85	120.02	110.60
2	C	3	A	C2-N3-C4	18.12	119.66	110.60
2	C	2	A	C2-N3-C4	17.89	119.55	110.60
2	D	3	A	C2-N3-C4	17.48	119.34	110.60
2	D	4	A	N1-C2-N3	-17.22	120.69	129.30
2	D	2	A	N1-C2-N3	-16.29	121.16	129.30
2	C	3	A	N7-C8-N9	-13.83	106.88	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	A	O4'-C1'-N9	13.61	119.08	108.20
2	D	4	A	C2-N3-C4	13.21	117.21	110.60
2	C	4	A	C5-C6-N6	-12.92	113.36	123.70
2	D	2	A	C5-C6-N1	12.60	124.00	117.70
2	D	3	A	C5-N7-C8	12.58	110.19	103.90
2	C	4	A	N7-C8-N9	-12.41	107.59	113.80
2	D	3	A	N7-C8-N9	-11.24	108.18	113.80
2	D	1	A	N9-C4-C5	11.00	110.20	105.80
2	D	2	A	N3-C4-C5	-10.81	119.23	126.80
2	D	4	A	O4'-C1'-N9	-10.75	99.60	108.20
2	D	1	A	N3-C4-C5	-10.60	119.38	126.80
2	D	1	A	C5-N7-C8	10.44	109.12	103.90
2	D	1	A	N7-C8-N9	-10.28	108.66	113.80
2	D	3	A	C4-C5-N7	-10.01	105.70	110.70
2	C	4	A	C5-C6-N1	9.98	122.69	117.70
2	C	4	A	N3-C4-C5	-9.97	119.82	126.80
2	D	1	A	C4-C5-N7	-9.94	105.73	110.70
2	C	4	A	C5-N7-C8	9.89	108.85	103.90
2	C	2	A	C5-C6-N6	-9.87	115.80	123.70
2	C	1	A	N3-C4-C5	-9.75	119.97	126.80
2	C	2	A	N3-C4-C5	-9.72	119.99	126.80
2	C	1	A	C5-C6-N1	9.68	122.54	117.70
2	D	2	A	O4'-C1'-N9	-9.67	100.46	108.20
2	C	1	A	O4'-C1'-N9	9.63	115.90	108.20
2	C	3	A	C5-N7-C8	9.54	108.67	103.90
2	D	3	A	N9-C4-C5	9.45	109.58	105.80
2	C	2	A	C4-C5-C6	9.44	121.72	117.00
2	D	3	A	C4-C5-C6	9.43	121.72	117.00
2	C	3	A	C8-N9-C4	9.41	109.56	105.80
2	C	1	A	N7-C8-N9	-9.35	109.13	113.80
2	D	3	A	N3-C4-C5	-9.35	120.26	126.80
2	D	4	A	C1'-O4'-C4'	-9.24	102.51	109.90
2	C	4	A	N1-C6-N6	8.91	123.94	118.60
2	C	4	A	C4-C5-C6	8.53	121.26	117.00
2	D	4	A	C4-C5-C6	8.51	121.26	117.00
2	D	1	A	C5-C6-N1	8.44	121.92	117.70
2	D	4	A	N9-C4-C5	8.38	109.15	105.80
2	C	1	A	C5-N7-C8	8.35	108.08	103.90
2	D	2	A	C5-C6-N6	-8.34	117.03	123.70
2	D	2	A	N7-C8-N9	-8.32	109.64	113.80
2	C	1	A	C5-C6-N6	-7.84	117.43	123.70
2	D	4	A	C5-C6-N6	-7.80	117.46	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	A	N9-C4-C5	7.76	108.90	105.80
2	D	4	A	O4'-C1'-C2'	-7.75	98.05	105.80
2	C	4	A	N3-C4-N9	7.75	133.60	127.40
2	D	2	A	C4-C5-N7	-7.58	106.91	110.70
2	D	2	A	O4'-C1'-C2'	-7.52	98.28	105.80
2	C	2	A	C5-C6-N1	7.52	121.46	117.70
2	C	3	A	C5-C6-N6	-7.38	117.79	123.70
2	C	2	A	N7-C8-N9	-7.32	110.14	113.80
2	D	1	A	C4-C5-C6	7.27	120.64	117.00
2	C	2	A	C3'-C2'-C1'	-7.12	95.80	101.50
2	D	2	A	C5-N7-C8	7.11	107.46	103.90
2	D	4	A	N3-C4-C5	-6.98	121.92	126.80
2	C	2	A	N1-C6-N6	6.90	122.74	118.60
2	D	2	A	C4-C5-C6	6.88	120.44	117.00
2	C	1	A	N9-C4-C5	6.71	108.48	105.80
2	D	4	A	N7-C8-N9	-6.67	110.46	113.80
2	C	4	A	C8-N9-C4	6.63	108.45	105.80
2	C	3	A	C1'-O4'-C4'	-6.63	104.60	109.90
2	C	3	A	C5-C6-N1	6.55	120.98	117.70
2	D	1	A	C1'-O4'-C4'	-6.53	104.67	109.90
2	D	3	A	C5-C6-N1	6.51	120.95	117.70
2	D	2	A	C1'-O4'-C4'	-6.46	104.73	109.90
2	C	2	A	C5-N7-C8	6.45	107.13	103.90
2	C	4	A	O4'-C1'-N9	6.38	113.31	108.20
2	C	1	A	P-O3'-C3'	6.38	127.36	119.70
2	C	1	A	C4-C5-N7	-6.32	107.54	110.70
2	C	3	A	P-O3'-C3'	6.30	127.26	119.70
2	D	4	A	N9-C1'-C2'	6.23	122.10	114.00
2	C	3	A	N3-C4-C5	-6.05	122.57	126.80
2	C	3	A	C4-C5-N7	-6.01	107.70	110.70
2	C	1	A	C4-C5-C6	5.88	119.94	117.00
2	D	2	A	C6-N1-C2	-5.83	115.10	118.60
2	C	2	A	N9-C4-C5	5.82	108.13	105.80
2	D	4	A	C4-C5-N7	-5.80	107.80	110.70
2	D	4	A	C5-C6-N1	5.73	120.57	117.70
2	D	3	A	C1'-O4'-C4'	-5.70	105.34	109.90
2	C	2	A	O4'-C1'-N9	5.62	112.70	108.20
2	D	4	A	N1-C6-N6	5.62	121.97	118.60
2	C	2	A	N3-C4-N9	5.60	131.88	127.40
2	D	2	A	N3-C4-N9	5.58	131.87	127.40
2	C	3	A	C4-N9-C1'	-5.56	116.29	126.30
2	D	1	A	P-O3'-C3'	5.52	126.32	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	A	C5-C6-N6	-5.51	119.29	123.70
2	D	4	A	C5-N7-C8	5.33	106.56	103.90
2	C	4	A	C1'-O4'-C4'	-5.30	105.66	109.90
2	D	3	A	C5-C6-N6	-5.29	119.47	123.70
2	C	1	A	N3-C4-N9	5.18	131.54	127.40
2	C	2	A	C4'-C3'-C2'	-5.13	97.47	102.60

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	3427	0	3486	41	0
1	B	3427	0	3486	32	0
2	C	88	0	44	1	0
2	D	88	0	44	1	0
3	A	38	0	0	3	0
3	B	53	0	0	1	0
3	C	2	0	0	0	0
3	D	3	0	0	0	0
All	All	7126	0	7060	69	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 5.

All (69) 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:209:ARG:CZ	1:A:209:ARG:HB3	1.89	0.98
1:A:55:ASP:HB3	3:A:616:HOH:O	1.74	0.86
1:A:58:SER:HA	1:A:274:VAL:HG11	1.61	0.82
1:A:56:PHE:HA	1:A:61:THR:HG22	1.64	0.79
1:A:275:ASN:HB3	1:A:278:GLU:OE1	1.86	0.76
1:A:129:ASP:OD1	1:A:131:THR:HG23	1.86	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:ASP:OD1	1:A:131:THR:CG2	2.38	0.72
1:B:405:ALA:O	1:B:409:THR:HG23	1.91	0.71
1:B:207:MET:HE1	1:B:212:PHE:HD1	1.55	0.69
1:A:275:ASN:HB3	1:A:278:GLU:CD	2.15	0.67
1:B:171:LYS:HD2	1:B:171:LYS:O	1.99	0.63
1:B:312:PRO:O	1:B:317:LYS:NZ	2.32	0.61
1:A:40:LYS:HE2	1:A:40:LYS:HA	1.81	0.61
1:B:207:MET:HE3	1:B:212:PHE:HB2	1.85	0.58
1:A:161:TYR:CD2	1:A:183:THR:HG22	2.41	0.56
1:A:56:PHE:HA	1:A:61:THR:CG2	2.33	0.55
1:A:52:ILE:HD12	2:D:1:A:N1	2.22	0.55
1:A:161:TYR:HD2	1:A:183:THR:HG22	1.72	0.54
1:A:332:GLN:OE1	1:B:376:ARG:HG3	2.07	0.54
1:A:209:ARG:CZ	1:A:209:ARG:CB	2.70	0.54
1:B:207:MET:CE	1:B:285:LEU:HD21	2.39	0.53
1:B:18:LYS:HA	1:B:26:SER:O	2.09	0.53
1:A:333:ARG:HG3	1:B:333:ARG:HD3	1.89	0.53
1:A:207:MET:HG3	1:A:211:GLU:HG2	1.90	0.52
1:B:1:MET:N	3:B:501:HOH:O	2.29	0.52
1:B:183:THR:O	1:B:183:THR:HG23	2.09	0.52
1:B:48:LEU:HD12	1:B:48:LEU:H	1.77	0.50
1:B:55:ASP:OD2	1:B:61:THR:HB	2.12	0.49
1:A:322:SER:OG	1:A:334:VAL:HG11	2.13	0.49
1:B:345:ALA:HB1	1:B:409:THR:HG21	1.95	0.48
1:B:112:HIS:CE1	1:B:116:GLU:OE2	2.66	0.48
1:B:20:ARG:HG3	1:B:175:GLU:OE2	2.13	0.48
1:A:242:GLY:HA2	1:A:326:PHE:CE1	2.48	0.47
1:B:48:LEU:HD12	1:B:48:LEU:N	2.29	0.47
1:A:333:ARG:HD2	1:A:380:ALA:O	2.15	0.47
1:A:361:ARG:NH1	1:A:387:ASN:OD1	2.42	0.46
1:B:405:ALA:O	1:B:409:THR:CG2	2.62	0.46
1:A:207:MET:CG	1:A:211:GLU:HG2	2.46	0.46
1:B:345:ALA:CB	1:B:409:THR:HG21	2.46	0.46
1:A:10:ASP:N	3:A:601:HOH:O	2.33	0.46
1:A:262:GLU:O	1:A:266:GLU:HG3	2.16	0.46
1:B:207:MET:HE1	1:B:285:LEU:HD21	1.98	0.46
1:A:129:ASP:CG	1:A:131:THR:HG23	2.36	0.46
1:B:207:MET:CE	1:B:212:PHE:HD1	2.26	0.45
1:B:56:PHE:HE1	1:B:281:VAL:HG23	1.81	0.44
1:A:129:ASP:OD1	1:A:131:THR:HG22	2.14	0.44
1:B:242:GLY:HA2	1:B:326:PHE:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:VAL:CG1	1:A:8:TRP:NE1	2.81	0.44
1:A:315:ILE:HD11	1:A:341:ILE:HG22	1.99	0.43
1:B:322:SER:OG	1:B:334:VAL:HG11	2.18	0.43
1:A:331:LEU:O	1:A:335:ARG:HG3	2.19	0.43
1:A:142:TYR:OH	1:B:143:ARG:HD3	2.19	0.43
1:A:311:GLU:H	1:A:311:GLU:HG3	1.63	0.42
1:A:6:VAL:HG22	1:A:129:ASP:HB3	2.02	0.42
1:A:381:HIS:NE2	2:C:3:A:H1'	2.34	0.42
1:B:207:MET:HE2	1:B:285:LEU:CD2	2.50	0.42
1:B:242:GLY:HA2	1:B:326:PHE:CZ	2.55	0.41
1:A:356:ARG:HD3	3:A:624:HOH:O	2.20	0.41
1:A:275:ASN:CB	1:A:278:GLU:CD	2.87	0.41
1:A:350:ARG:NH1	1:A:401:PRO:HG2	2.36	0.41
1:A:316:GLU:OE1	1:A:407:ARG:HG2	2.21	0.41
1:A:333:ARG:HG3	1:B:333:ARG:HH11	1.85	0.41
1:B:170:PRO:O	1:B:171:LYS:CB	2.68	0.41
1:B:368:ALA:HB1	1:B:371:ARG:HH21	1.86	0.41
1:A:209:ARG:NH1	1:A:209:ARG:HB3	2.32	0.40
1:B:169:VAL:HG22	1:B:172:ILE:HG23	2.03	0.40
1:A:130:SER:HB2	1:A:138:THR:HG21	2.03	0.40
1:A:19:TYR:HA	1:A:176:LEU:O	2.20	0.40
1:B:52:ILE:O	1:B:69:ARG:NH1	2.45	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	
1	A	430/440 (98%)	415 (96%)	15 (4%)	0	100	100
1	B	430/440 (98%)	419 (97%)	11 (3%)	0	100	100
All	All	860/880 (98%)	834 (97%)	26 (3%)	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	369/376 (98%)	345 (94%)	24 (6%)	17	14
1	B	369/376 (98%)	346 (94%)	23 (6%)	18	15
All	All	738/752 (98%)	691 (94%)	47 (6%)	17	14

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	25	THR
1	A	52	ILE
1	A	57	SER
1	A	62	LEU
1	A	81	ARG
1	A	91	ILE
1	A	131	THR
1	A	164	ASN
1	A	174	LYS
1	A	175	GLU
1	A	209	ARG
1	A	211	GLU
1	A	214	ARG
1	A	215	LEU
1	A	222	LEU
1	A	227	LYS
1	A	259	SER
1	A	262	GLU
1	A	333	ARG
1	A	340	LYS
1	A	399	LYS
1	A	412	LYS
1	A	432	SER
1	B	16	GLU

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Mol	Chain	Res	Type
1	B	17	THR
1	B	48	LEU
1	B	57	SER
1	B	62	LEU
1	B	69	ARG
1	B	81	ARG
1	B	82	GLU
1	B	145	LEU
1	B	169	VAL
1	B	171	LYS
1	B	172	ILE
1	B	209	ARG
1	B	215	LEU
1	B	223	LYS
1	B	227	LYS
1	B	256	ARG
1	B	325	LEU
1	B	350	ARG
1	B	370	ASN
1	B	372	GLU
1	B	409	THR
1	B	431	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	112	HIS
1	B	135	ASN
1	B	164	ASN

### 5.3.3 RNA ⓘ

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

All (5) RNA backbone outliers are listed below:

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

All (5) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	1	A
2	C	3	A
2	D	1	A
2	D	2	A
2	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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	432/440 (98%)	0.37	20 (4%) 32 37	35, 48, 77, 121	0
1	B	432/440 (98%)	0.27	12 (2%) 53 59	35, 45, 75, 107	0
2	C	4/4 (100%)	0.17	0 100 100	72, 85, 98, 125	0
2	D	4/4 (100%)	-0.59	0 100 100	40, 43, 49, 52	0
All	All	872/888 (98%)	0.32	32 (3%) 41 48	35, 46, 77, 125	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	369	ALA	7.3
1	B	171	LYS	5.4
1	B	81	ARG	4.9
1	B	371	ARG	4.4
1	A	396	TRP	4.4
1	A	368	ALA	3.5
1	B	199	ASP	3.4
1	B	370	ASN	3.4
1	A	395	ALA	3.0
1	A	353	ALA	2.8
1	B	368	ALA	2.8
1	B	310	GLU	2.7
1	A	399	LYS	2.7
1	A	55	ASP	2.6
1	A	352	GLY	2.6
1	B	260	TYR	2.6
1	A	354	PHE	2.6
1	B	369	ALA	2.5
1	A	227	LYS	2.5
1	B	277	GLU	2.4
1	A	348	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	400	ARG	2.4
1	A	82	GLU	2.4
1	A	111	LEU	2.3
1	A	397	GLU	2.3
1	A	199	ASP	2.2
1	B	254	ILE	2.2
1	A	191	LEU	2.1
1	A	78	GLY	2.1
1	B	372	GLU	2.1
1	A	148	LEU	2.0
1	A	431	GLU	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.