



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:05 PM UTC

PDB ID : 7M50 / pdb_00007m50
Title : Crystallographic structure of a cubic crystal form of STMV grown from ammonium sulfate
Authors : McPherson, A.
Deposited on : 2021-03-22
Resolution : 2.31 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
Mogul : 2022.3.0, CSD as543be (2022)
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

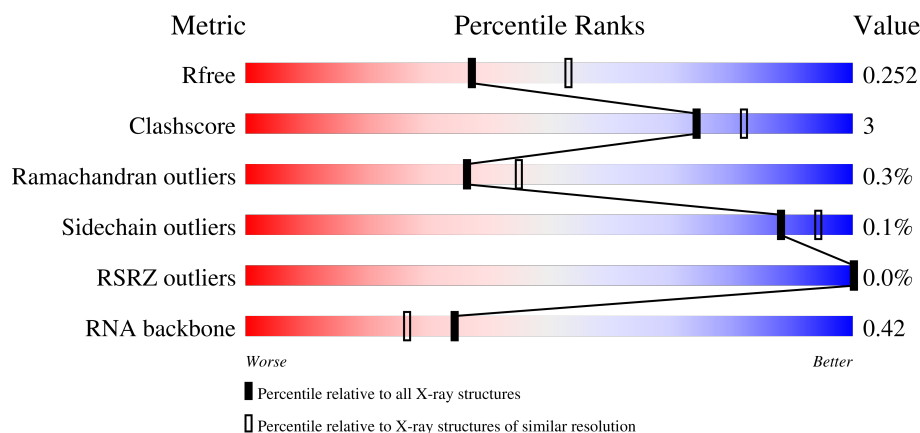
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.31 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.



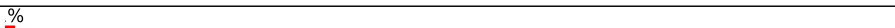
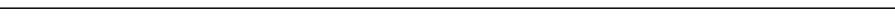
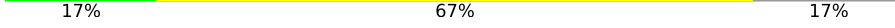
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Ramachandran outliers	187476	8303 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)
RNA backbone	3983	1116 (2.64-2.00)

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	159	 86% 6% 8%
1	B	159	 84% 6% 10%
1	C	159	 85% 5% 10%
1	D	159	 84% 6% 9%

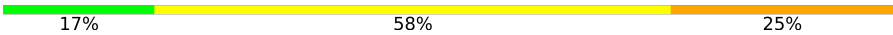
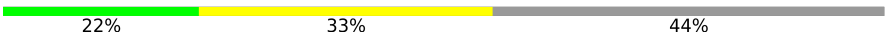
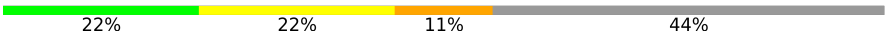
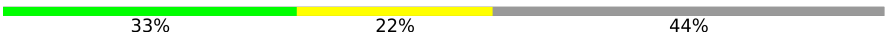
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Mol	Chain	Length	Quality of chain
1	E	159	
1	F	159	
1	G	159	
1	GG	159	
1	H	159	
1	HH	159	
1	I	159	
1	II	159	
1	J	159	
1	JJ	159	
1	K	159	
1	KK	159	
1	L	159	
1	M	159	
1	N	159	
1	O	159	
2	P	8	
2	TT	8	
2	V	8	
3	S	6	
3	T	6	
4	UU	10	
4	X	10	
4	bb	10	
5	Y	7	

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Mol	Chain	Length	Quality of chain
5	a	7	
6	WW	12	
6	e	12	
7	h	10	
8	i	9	
8	ll	9	
8	qq	9	
9	kk	8	
9	m	8	
9	n	8	

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 29786 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 Coat protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	8	0
			1130	709	201	214	6			
1	B	143	Total	C	N	O	S	0	10	0
			1110	699	197	208	6			
1	C	143	Total	C	N	O	S	0	14	0
			1110	699	197	208	6			
1	D	144	Total	C	N	O	S	0	12	0
			1118	703	199	210	6			
1	E	145	Total	C	N	O	S	0	12	0
			1126	707	200	213	6			
1	F	147	Total	C	N	O	S	0	10	0
			1137	713	202	216	6			
1	G	146	Total	C	N	O	S	0	8	0
			1130	709	201	214	6			
1	H	147	Total	C	N	O	S	0	10	0
			1137	713	202	216	6			
1	I	144	Total	C	N	O	S	0	8	0
			1118	703	199	210	6			
1	J	147	Total	C	N	O	S	0	9	0
			1137	713	202	216	6			
1	K	146	Total	C	N	O	S	0	8	0
			1130	709	201	214	6			
1	L	157	Total	C	N	O	S	0	6	0
			1215	761	221	227	6			
1	M	144	Total	C	N	O	S	0	7	0
			1118	703	199	210	6			
1	N	144	Total	C	N	O	S	0	3	0
			1118	703	199	210	6			
1	O	150	Total	C	N	O	S	0	17	0
			1163	728	209	220	6			
1	GG	144	Total	C	N	O	S	0	12	0
			1118	703	199	210	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	HH	144	Total	C	N	O	S	0	9	0
			1118	703	199	210	6			
1	II	143	Total	C	N	O	S	0	9	0
			1110	699	197	208	6			
1	JJ	147	Total	C	N	O	S	0	13	0
			1137	713	202	216	6			
1	KK	145	Total	C	N	O	S	0	9	0
			1126	707	200	213	6			

- Molecule 2 is a RNA chain called RNA (5'-R(P*AP*AP*AP*AP*AP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	7	Total	C	N	O	P	0	7	0
			154	70	35	42	7			
2	V	7	Total	C	N	O	P	0	6	0
			154	70	35	42	7			
2	TT	7	Total	C	N	O	P	0	6	0
			154	70	35	42	7			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	S	5	Total	C	N	O	P	0	5	0
			110	50	25	30	5			
3	T	5	Total	C	N	O	P	0	5	0
			110	50	25	30	5			

- Molecule 4 is a RNA chain called RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	X	8	Total	C	N	O	P	0	8	0
			176	80	40	48	8			
4	UU	8	Total	C	N	O	P	0	6	0
			176	80	40	48	8			
4	bb	9	Total	C	N	O	P	0	8	0
			198	90	45	54	9			

- Molecule 5 is a RNA chain called RNA (5'-R(P*UP*UP*UP*UP*UP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	Y	6	Total 132	C 60	N 30	O 36	P 6	0	6	0
5	a	6	Total 132	C 60	N 30	O 36	P 6	0	6	0

- Molecule 6 is a RNA chain called RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	e	12	Total 240	C 108	N 24	O 96	P 12	0	10	0
6	WW	10	Total 200	C 90	N 20	O 80	P 10	0	2	0

- Molecule 7 is a RNA chain called RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*U))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	h	8	Total	C	N	O	P	0	7	0
			160	72	16	64	8			

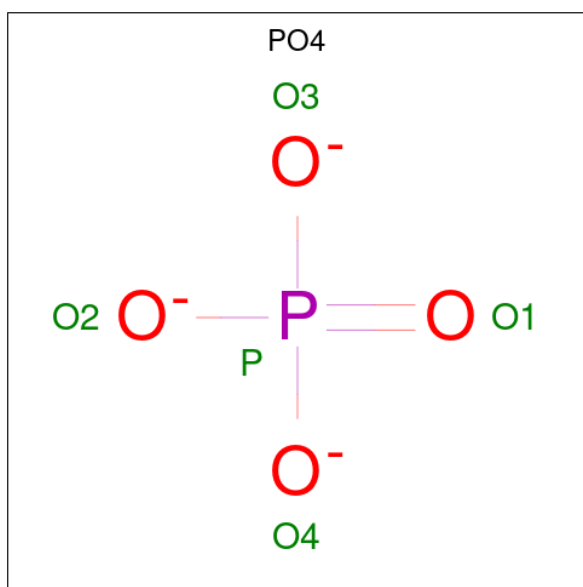
- Molecule 8 is a RNA chain called RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*U))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	i	5	Total 100	C 45	N 10	O 40	P 5	0	5	0
8	ll	5	Total 100	C 45	N 10	O 40	P 5	0	5	0
8	qq	5	Total 100	C 45	N 10	O 40	P 5	0	4	0

- Molecule 9 is a RNA chain called RNA (5'-R(P*UP*UP*UP*UP*UP*UP*U))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	m	7	Total 140	C 63	N 14	O 56	P 7	0	7	0
9	n	7	Total 140	C 63	N 14	O 56	P 7	0	7	0
9	kk	6	Total 120	C 54	N 12	O 48	P 6	0	6	0

- Molecule 10 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	O	P	0	0
			5	4	1		
10	H	1	Total	O	P	0	0
			5	4	1		
10	L	1	Total	O	P	0	0
			5	4	1		
10	HH	1	Total	O	P	0	0
			5	4	1		
10	II	1	Total	O	P	0	0
			5	4	1		
10	R	1	Total	O	P	0	0
			5	4	1		

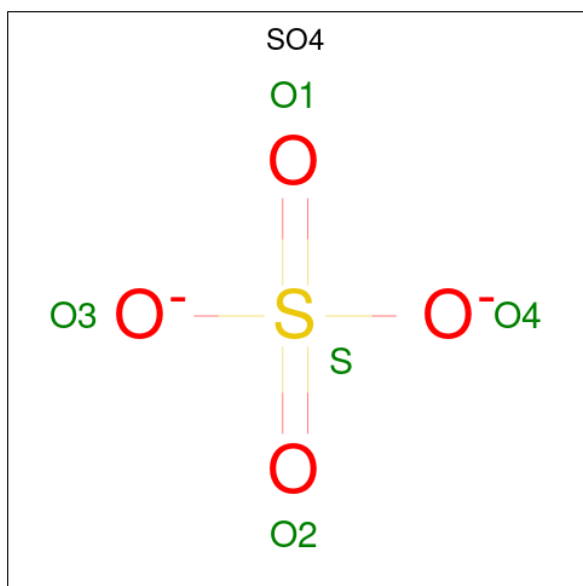
- Molecule 11 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	1	Total	Mg	0	0
			1	1		
11	G	1	Total	Mg	0	0
			1	1		
11	H	1	Total	Mg	0	0
			1	1		
11	JJ	2	Total	Mg	0	0
			2	2		

- Molecule 12 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	C	1	Total Cl 1 1	0	0

- Molecule 13 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
13	F	1	Total O S 5 4 1	0	0
13	M	1	Total O S 5 4 1	0	0
13	M	1	Total O S 5 4 1	0	0
13	O	1	Total O S 5 4 1	0	0
13	JJ	1	Total O S 5 4 1	0	0
13	KK	1	Total O S 5 4 1	0	0

- Molecule 14 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	A	128	Total O 128 128	0	0
14	B	149	Total O 149 149	0	1
14	C	118	Total O 118 118	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	D	177	Total 177	O 177	0	1
14	E	110	Total 110	O 110	0	0
14	F	127	Total 127	O 127	0	0
14	G	153	Total 153	O 153	0	0
14	H	149	Total 149	O 149	0	0
14	I	105	Total 105	O 105	0	0
14	J	155	Total 155	O 155	0	0
14	K	198	Total 198	O 198	0	0
14	L	138	Total 138	O 138	0	0
14	M	150	Total 150	O 150	0	1
14	N	96	Total 96	O 96	0	0
14	O	252	Total 252	O 252	0	0
14	P	39	Total 39	O 39	0	0
14	S	29	Total 29	O 29	0	0
14	T	22	Total 22	O 22	0	0
14	V	61	Total 61	O 61	0	0
14	X	36	Total 36	O 36	0	0
14	Y	55	Total 55	O 55	0	0
14	a	67	Total 67	O 67	0	0
14	e	193	Total 193	O 193	0	0
14	h	64	Total 64	O 64	0	0

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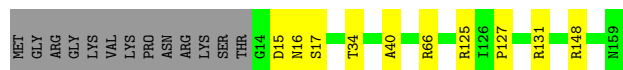
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	i	17	Total 17	O 17	0	0
14	m	39	Total 39	O 39	0	0
14	n	48	Total 48	O 48	0	0
14	GG	127	Total 127	O 127	0	1
14	HH	188	Total 188	O 188	0	0
14	II	271	Total 271	O 271	0	0
14	JJ	125	Total 125	O 125	0	0
14	KK	200	Total 200	O 200	0	0
14	TT	46	Total 46	O 46	0	0
14	UU	96	Total 96	O 96	0	0
14	WW	76	Total 76	O 76	0	0
14	bb	212	Total 212	O 212	0	0
14	kk	65	Total 65	O 65	0	0
14	ll	24	Total 24	O 24	0	0
14	qq	13	Total 13	O 13	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: Coat protein

Chain A:  86% 6% 8%



- Molecule 1: Coat protein

Chain B:  84% 6% 10%



- Molecule 1: Coat protein

Chain C:  85% 5% 10%



- Molecule 1: Coat protein

Chain D:  84% 6% 9%



- Molecule 1: Coat protein

Chain E:  83% 8% 9%

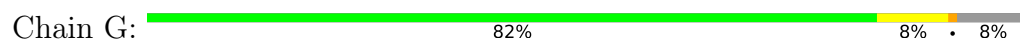


- Molecule 1: Coat protein

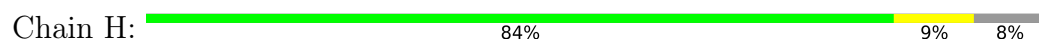
Chain F:  87% 5% 8%



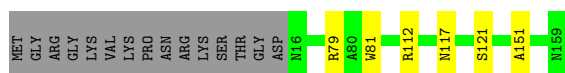
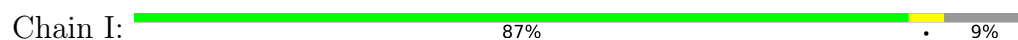
- Molecule 1: Coat protein



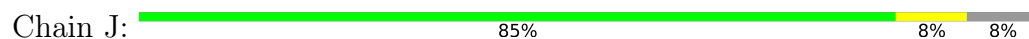
- Molecule 1: Coat protein



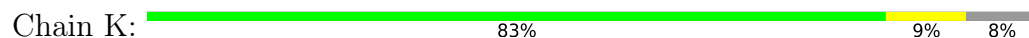
- Molecule 1: Coat protein



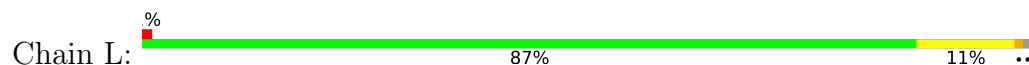
- Molecule 1: Coat protein



- Molecule 1: Coat protein

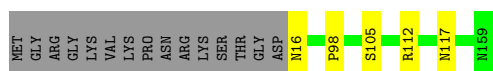


- Molecule 1: Coat protein



- Molecule 1: Coat protein





- Molecule 1: Coat protein

Chain N: 85% 5% 9%



- Molecule 1: Coat protein

Chain O: 89% 6% 6%



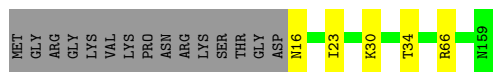
- Molecule 1: Coat protein

Chain GG: 84% 6% 9%



- Molecule 1: Coat protein

Chain HH: 87% 0% 9%



- Molecule 1: Coat protein

Chain II: 81% 8% 10%



- Molecule 1: Coat protein

Chain JJ: 87% 6% 8%



- Molecule 1: Coat protein

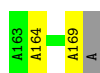
Chain KK: 88% 0% 9%



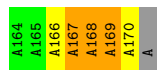
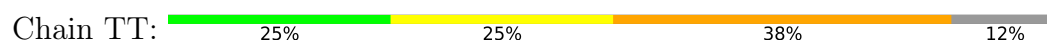
- Molecule 2: RNA (5'-R(P*AP*AP*AP*AP*AP*AP*AP*A)-3')



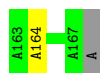
- Molecule 2: RNA (5'-R(P*AP*AP*AP*AP*AP*AP*AP*A)-3')



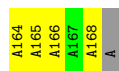
- Molecule 2: RNA (5'-R(P*AP*AP*AP*AP*AP*AP*AP*A)-3')



- Molecule 3: RNA (5'-R(P*AP*AP*AP*AP*AP*A)-3')



- Molecule 3: RNA (5'-R(P*AP*AP*AP*AP*AP*A)-3')

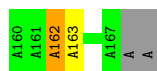


- Molecule 4: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')



- Molecule 4: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')

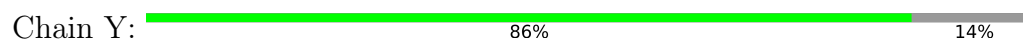




- Molecule 4: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')



- Molecule 5: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*U)-3')



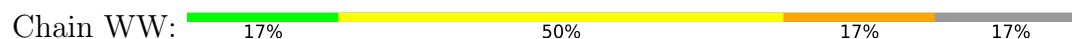
- Molecule 5: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*U)-3')



- Molecule 6: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')



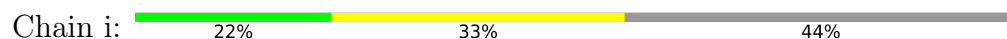
- Molecule 6: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')

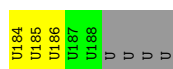


- Molecule 7: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')

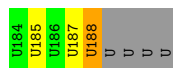


- Molecule 8: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*U)-3')

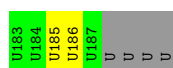
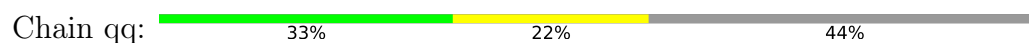




- Molecule 8: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')



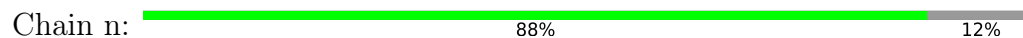
- Molecule 8: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')



- Molecule 9: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')



- Molecule 9: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')



- Molecule 9: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 3	Depositor
Cell constants a, b, c, α , β , γ	234.05Å 234.05Å 234.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	74.01 – 2.31 74.01 – 2.31	Depositor EDS
% Data completeness (in resolution range)	86.0 (74.01-2.31) 82.9 (74.01-2.31)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	0.18	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.71 (at 2.32Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.246 , 0.271 0.232 , 0.252	Depositor DCC
R_{free} test set	1990 reflections (1.25%)	wwPDB-VP
Wilson B-factor (Å ²)	3.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 17.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.28$, $\langle L^2 \rangle = 0.13$	Xtriage
Estimated twinning fraction	0.147 for l,-k,h	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	29786	wwPDB-VP
Average B, all atoms (Å ²)	9.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.96% 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: MG, SO4, PO4, CL

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	0.08	0/1154	0.24	0/1572
1	B	0.07	0/1134	0.23	0/1545
1	C	0.07	0/1134	0.22	0/1545
1	D	0.07	0/1142	0.23	0/1556
1	E	0.07	0/1150	0.24	0/1567
1	F	0.08	0/1161	0.25	0/1582
1	G	0.12	0/1154	0.34	2/1572 (0.1%)
1	GG	0.07	0/1142	0.23	0/1556
1	H	0.08	0/1161	0.25	0/1582
1	HH	0.06	0/1142	0.22	0/1556
1	I	0.07	0/1142	0.23	0/1556
1	II	0.08	0/1134	0.30	0/1545
1	J	0.08	0/1161	0.24	0/1582
1	JJ	0.07	0/1161	0.23	0/1582
1	K	0.07	0/1154	0.24	0/1572
1	KK	0.08	0/1150	0.23	0/1567
1	L	0.07	0/1240	0.24	0/1685
1	M	0.08	0/1142	0.24	0/1556
1	N	0.10	0/1142	0.26	0/1556
1	O	0.07	0/1187	0.23	0/1615
2	P	0.10	0/174	0.24	0/269
2	TT	0.11	0/174	0.83	2/269 (0.7%)
2	V	0.08	0/174	0.18	0/269
3	S	0.09	0/124	0.24	0/191
3	T	0.07	0/124	0.15	0/191
4	UU	0.14	0/199	0.32	0/308
4	X	0.10	0/199	0.32	0/308
4	bb	0.06	0/224	0.14	0/347
5	Y	0.06	0/149	0.17	0/230
5	a	0.13	0/149	0.31	0/230
6	WW	0.14	0/219	0.39	0/336
6	e	0.15	0/263	0.52	0/404

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
7	h	0.10	0/175	0.33	0/268
8	i	0.12	0/109	0.36	0/166
8	ll	0.08	0/109	0.19	0/166
8	qq	0.11	0/109	0.31	0/166
9	kk	0.08	0/131	0.19	0/200
9	m	0.06	0/153	0.13	0/234
9	n	0.05	0/153	0.12	0/234
All	All	0.08	0/26198	0.26	4/36235 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	TT	169[A]	A	OP1-P-O3'	-8.75	81.74	108.00
2	TT	169[A]	A	OP2-P-O3'	-7.86	84.41	108.00
1	G	16	ASN	CA-C-N	5.84	132.21	121.70
1	G	16	ASN	C-N-CA	5.84	132.21	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1130	0	1117	6	0
1	B	1110	0	1103	6	0
1	C	1110	0	1099	4	1
1	D	1118	0	1103	8	0
1	E	1126	0	1108	11	0
1	F	1137	0	1120	6	0
1	G	1130	0	1115	12	0
1	GG	1118	0	1103	6	0
1	H	1137	0	1118	10	0
1	HH	1118	0	1109	5	0
1	I	1118	0	1106	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	II	1110	0	1105	11	0
1	J	1137	0	1122	9	0
1	JJ	1137	0	1118	7	0
1	K	1130	0	1115	10	0
1	KK	1126	0	1110	4	0
1	L	1215	0	1216	13	0
1	M	1118	0	1109	4	0
1	N	1118	0	1112	7	0
1	O	1163	0	1146	8	0
2	P	154	0	78	3	0
2	TT	154	0	78	2	0
2	V	154	0	78	1	0
3	S	110	0	56	0	0
3	T	110	0	56	2	0
4	UU	176	0	89	2	0
4	X	176	0	89	1	0
4	bb	198	0	100	3	0
5	Y	132	0	67	0	0
5	a	132	0	67	1	0
6	WW	200	0	95	1	0
6	e	240	0	121	4	0
7	h	160	0	81	2	0
8	i	100	0	51	2	0
8	ll	100	0	51	2	0
8	qq	100	0	51	2	0
9	kk	120	0	61	1	0
9	m	140	0	71	3	0
9	n	140	0	71	0	0
10	B	5	0	0	0	0
10	H	5	0	0	0	0
10	HH	5	0	0	0	0
10	II	5	0	0	0	0
10	L	5	0	0	0	0
10	R	5	0	0	0	0
11	B	1	0	0	0	0
11	G	1	0	0	0	0
11	H	1	0	0	0	0
11	JJ	2	0	0	0	0
12	C	1	0	0	0	0
13	F	5	0	0	0	0
13	JJ	5	0	0	0	0
13	KK	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	M	10	0	0	0	0
13	O	5	0	0	0	0
14	A	128	0	0	2	0
14	B	149	0	0	0	0
14	C	118	0	0	0	0
14	D	177	0	0	1	0
14	E	110	0	0	1	0
14	F	127	0	0	0	0
14	G	153	0	0	2	0
14	GG	127	0	0	0	0
14	H	149	0	0	2	0
14	HH	188	0	0	0	0
14	I	105	0	0	1	0
14	II	271	0	0	1	0
14	J	155	0	0	1	0
14	JJ	125	0	0	2	0
14	K	198	0	0	0	0
14	KK	200	0	0	0	0
14	L	138	0	0	0	0
14	M	150	0	0	1	0
14	N	96	0	0	0	0
14	O	252	0	0	1	0
14	P	39	0	0	0	0
14	S	29	0	0	0	0
14	T	22	0	0	0	0
14	TT	46	0	0	1	0
14	UU	96	0	0	1	0
14	V	61	0	0	1	0
14	WW	76	0	0	1	0
14	X	36	0	0	0	0
14	Y	55	0	0	0	0
14	a	67	0	0	0	0
14	bb	212	0	0	1	0
14	e	193	0	0	0	0
14	h	64	0	0	0	0
14	i	17	0	0	0	0
14	kk	65	0	0	0	0
14	ll	24	0	0	0	0
14	m	39	0	0	1	0
14	n	48	0	0	0	0
14	qq	13	0	0	0	0
All	All	29786	0	23765	140	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 3.

All (140) 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:P:166[A]:A:N6	6:e:183[A]:U:O4	2.18	0.76
9:m:186[A]:U:OP1	14:m:201:HOH:O	2.13	0.66
1:H:36:THR:OG1	9:kk:186[A]:U:O2'	2.13	0.65
1:L:12:SER:OG	1:L:16:ASN:O	2.15	0.65
8:ll:187[A]:U:H2'	8:ll:188[A]:U:H4'	1.79	0.63
1:JJ:79[A]:ARG:NH2	14:JJ:304:HOH:O	2.33	0.61
4:UU:162[A]:A:N6	14:UU:202:HOH:O	2.33	0.60
1:N:81:TRP:O	1:N:151:ALA:N	2.35	0.59
1:G:38:VAL:O	14:WW:201:HOH:O	2.17	0.58
1:G:48:GLN:NE2	14:G:306:HOH:O	2.37	0.58
1:G:16:ASN:N	1:G:17:SER:HB3	2.17	0.58
1:D:25:ALA:N	14:D:205:HOH:O	2.40	0.55
1:E:125:ARG:NH1	14:E:204:HOH:O	2.39	0.55
1:G:125:ARG:NH2	14:G:307:HOH:O	2.39	0.55
2:TT:168[A]:A:N6	14:TT:201:HOH:O	2.41	0.54
4:bb:162[A]:A:N6	14:bb:202:HOH:O	2.40	0.54
7:h:186[A]:U:O2'	7:h:187[A]:U:OP1	2.24	0.53
4:X:168[A]:A:O2'	4:X:169[A]:A:OP1	2.27	0.53
1:O:14:GLY:HA2	8:i:184[A]:U:H4'	1.92	0.52
1:J:148:ARG:NH1	14:J:207:HOH:O	2.42	0.52
1:H:13:THR:OG1	1:H:14:GLY:N	2.34	0.51
1:II:19:VAL:HG22	1:II:20:VAL:H	1.76	0.51
6:e:179:U:H5''	6:e:180:U:H5	1.76	0.51
1:B:159:ASN:HD22	1:N:19:VAL:HG12	1.76	0.50
1:J:105:SER:HB3	1:J:107:GLU:OE1	2.12	0.50
1:L:36:THR:OG1	5:a:165[A]:A:O2'	2.30	0.50
1:I:112:ARG:NH2	14:I:209:HOH:O	2.44	0.49
1:I:79:ARG:NH2	9:m:188[A]:U:OP1	2.46	0.49
1:KK:57:VAL:HG22	1:KK:140:VAL:HG13	1.93	0.49
1:A:34:THR:HB	1:O:66:ARG:NH2	2.28	0.48
1:J:30:LYS:NZ	1:L:121:SER:O	2.41	0.48
1:E:129:ASN:O	1:II:132:GLN:NE2	2.43	0.48
1:GG:112:ARG:NH2	1:KK:44:GLU:OE2	2.36	0.48
2:V:164[A]:A:O2'	14:V:201:HOH:O	2.12	0.48
1:H:79:ARG:NH2	14:H:605:HOH:O	2.45	0.48
1:O:14:GLY:HA3	8:i:184[A]:U:P	2.54	0.48
1:L:44:GLU:OE2	1:M:112:ARG:NH2	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:81:TRP:N	1:N:151:ALA:O	2.48	0.47
1:D:30:LYS:NZ	1:JJ:121:SER:O	2.36	0.47
8:qq:185[A]:U:H2'	8:qq:186[A]:U:H6	1.80	0.47
1:GG:74:THR:OG1	1:GG:133:ASN:OD1	2.33	0.47
4:bb:164[A]:A:H2'	4:bb:165[A]:A:C8	2.50	0.46
1:JJ:44:GLU:OE2	1:KK:112:ARG:NH1	2.36	0.46
1:O:72:SER:N	14:O:502:HOH:O	2.48	0.46
1:K:117:ASN:HB3	1:L:116:ILE:HG12	1.98	0.46
1:GG:127:PRO:O	1:GG:131:ARG:HG3	2.16	0.46
1:H:121:SER:O	1:HH:30:LYS:NZ	2.41	0.46
1:E:34:THR:HB	1:F:66:ARG:NH2	2.31	0.46
1:GG:36:THR:OG1	8:ll:185[A]:U:O2'	2.31	0.46
1:II:81:TRP:O	1:II:151:ALA:N	2.41	0.45
8:qq:185[A]:U:H2'	8:qq:186[A]:U:C6	2.51	0.45
1:L:48:GLN:HB2	1:L:51:ILE:HG22	1.97	0.45
1:K:54:LYS:NZ	1:K:142:GLU:OE2	2.46	0.45
1:L:11:LYS:HE2	1:L:21:THR:HB	1.97	0.45
1:E:30:LYS:NZ	1:G:121:SER:O	2.35	0.45
1:H:44:GLU:OE2	1:I:112:ARG:NH2	2.40	0.45
1:J:66:ARG:NH2	1:K:34:THR:HB	2.31	0.45
3:T:165[A]:A:H2'	3:T:166[A]:A:C8	2.52	0.45
6:e:189[A]:U:H2'	6:e:190[A]:U:O4'	2.16	0.45
1:E:15:ASP:HB3	1:E:16:ASN:H	1.63	0.45
1:N:44:GLU:OE2	1:O:112:ARG:NH2	2.47	0.45
1:II:129:ASN:O	1:II:132:GLN:NE2	2.34	0.45
1:M:117:ASN:HB3	1:N:116:ILE:HG12	1.97	0.45
1:J:34:THR:HB	1:K:66:ARG:NH2	2.31	0.45
1:D:34:THR:HB	1:II:66:ARG:NH2	2.32	0.44
1:M:16:ASN:N	14:M:311:HOH:O	2.50	0.44
1:K:105:SER:HB2	1:K:107:GLU:OE1	2.18	0.44
1:D:127:PRO:O	1:D:131:ARG:HG3	2.17	0.44
1:K:16:ASN:HB2	9:m:184[A]:U:H5'	1.99	0.44
1:C:84:LEU:HD12	1:C:148:ARG:HB2	2.00	0.44
4:UU:162[A]:A:H2'	4:UU:163[A]:A:H8	1.83	0.44
1:E:16:ASN:HD21	1:F:125:ARG:HH22	1.64	0.44
1:H:148:ARG:NH2	14:H:607:HOH:O	2.50	0.44
1:G:34:THR:HB	1:HH:66:ARG:NH2	2.33	0.44
1:GG:116:ILE:HG12	1:KK:117:ASN:HB3	2.00	0.44
1:H:74:THR:OG1	1:H:133:ASN:OD1	2.31	0.43
1:H:127:PRO:O	1:H:131:ARG:HG3	2.18	0.43
1:A:148:ARG:NH2	14:A:205:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105[A]:SER:HB3	1:B:107:GLU:OE1	2.18	0.43
1:JJ:127:PRO:O	1:JJ:131:ARG:HG3	2.18	0.43
1:I:81:TRP:O	1:I:151:ALA:N	2.45	0.43
1:O:127:PRO:O	1:O:131:ARG:HG3	2.19	0.43
1:G:125:ARG:HB2	1:HH:23:ILE:HD11	2.00	0.43
1:G:66:ARG:NH2	1:HH:34:THR:HB	2.33	0.43
1:II:84:LEU:HG	1:II:149[A]:GLN:HB2	2.01	0.43
1:A:66:ARG:NH2	1:O:34:THR:HB	2.33	0.43
1:A:40:ALA:HB2	6:e:186[A]:U:H4'	2.01	0.43
1:G:127:PRO:O	1:G:131:ARG:HG3	2.19	0.43
1:II:129:ASN:OD1	14:II:301:HOH:O	2.22	0.43
1:D:83:GLN:HB3	1:D:149:GLN:HB3	2.01	0.42
1:F:81:TRP:CD1	1:F:121:SER:HB3	2.54	0.42
1:I:81:TRP:CD1	1:I:121:SER:HB3	2.54	0.42
1:F:44:GLU:OE2	1:G:112:ARG:NH1	2.42	0.42
1:F:127:PRO:O	1:F:131:ARG:HG3	2.19	0.42
4:bb:163[A]:A:H2'	4:bb:164[A]:A:C8	2.54	0.42
1:C:81:TRP:O	1:C:151:ALA:N	2.41	0.42
1:K:112:ARG:NH2	1:O:44:GLU:OE2	2.47	0.42
6:WW:173:U:O2'	6:WW:174:U:O5'	2.36	0.42
1:B:127:PRO:O	1:B:131:ARG:HG3	2.18	0.42
1:E:127:PRO:O	1:E:131:ARG:HG3	2.19	0.42
1:J:81:TRP:CD1	1:J:121:SER:HB3	2.55	0.42
7:h:186[A]:U:HO2'	7:h:187[A]:U:P	2.43	0.42
1:E:66:ARG:NH2	1:F:34:THR:HB	2.34	0.42
1:K:14:GLY:C	1:K:16:ASN:H	2.28	0.42
1:L:99:LEU:HD11	1:L:142:GLU:HG3	2.02	0.42
1:JJ:112:ARG:NH2	14:JJ:307:HOH:O	2.53	0.42
1:B:68:ASP:OD1	1:B:68:ASP:N	2.53	0.42
1:B:117:ASN:HB3	1:C:116:ILE:HG12	2.00	0.42
1:D:117:ASN:HB3	1:E:116:ILE:HG12	2.02	0.42
1:I:117:ASN:HB3	1:J:116:ILE:HG12	2.01	0.42
1:J:99:LEU:HG	1:J:141:CYS:HA	2.01	0.41
1:JJ:97:LYS:HB3	1:JJ:142:GLU:HB2	2.00	0.41
1:A:127:PRO:O	1:A:131:ARG:HG3	2.20	0.41
1:K:127:PRO:O	1:K:131:ARG:HG3	2.20	0.41
1:L:6:VAL:O	1:L:7:LYS:HE2	2.20	0.41
1:N:121:SER:O	1:N:122:VAL:HG12	2.20	0.41
1:C:127:PRO:O	1:C:131:ARG:HG3	2.20	0.41
1:E:54:LYS:NZ	1:E:142:GLU:OE2	2.47	0.41
1:H:40:ALA:HB2	2:TT:167[A]:A:H4'	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:84:LEU:HG	1:G:149:GLN:HB2	2.03	0.41
1:H:99:LEU:HG	1:H:141:CYS:HA	2.03	0.41
1:D:66:ARG:NH2	1:II:34:THR:HB	2.35	0.41
1:K:81:TRP:CD1	1:K:121:SER:HB3	2.56	0.41
1:A:125:ARG:NH2	14:A:206:HOH:O	2.53	0.41
1:D:33:PRO:O	1:II:39:ARG:NH2	2.54	0.41
1:G:81:TRP:O	1:G:151:ALA:N	2.46	0.41
1:J:21:THR:HG22	2:P:165[A]:A:H5''	2.03	0.41
1:L:11:LYS:HG2	1:L:21:THR:HG21	2.03	0.41
1:L:127:PRO:O	1:L:131:ARG:HG3	2.21	0.41
1:M:98:PRO:O	1:M:105:SER:OG	2.33	0.41
1:II:117:ASN:HB3	1:JJ:116:ILE:HG12	2.02	0.41
1:N:84:LEU:HD12	1:N:148:ARG:HB2	2.02	0.41
3:T:164[A]:A:H5'	1:HH:16:ASN:HB2	2.03	0.41
1:L:84:LEU:HG	1:L:149:GLN:HB3	2.03	0.40
2:P:164[A]:A:H2'	2:P:165[A]:A:C8	2.56	0.40
1:II:17:SER:OG	1:II:18:ASN:N	2.53	0.40
1:B:94:VAL:HG13	1:B:122:VAL:HG12	2.02	0.40
1:L:11:LYS:O	1:L:13:THR:N	2.55	0.40
1:GG:81:TRP:CD1	1:GG:121:SER:HB3	2.57	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:C:49[A]:SER:OG	1:C:49[A]:SER:OG[2_675]	2.14	0.06

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	144/159 (91%)	135 (94%)	6 (4%)	3 (2%)	5 4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	141/159 (89%)	134 (95%)	6 (4%)	1 (1%)	18	22
1	C	141/159 (89%)	134 (95%)	7 (5%)	0	100	100
1	D	142/159 (89%)	132 (93%)	10 (7%)	0	100	100
1	E	143/159 (90%)	136 (95%)	7 (5%)	0	100	100
1	F	145/159 (91%)	133 (92%)	12 (8%)	0	100	100
1	G	144/159 (91%)	135 (94%)	8 (6%)	1 (1%)	18	22
1	GG	142/159 (89%)	133 (94%)	9 (6%)	0	100	100
1	H	145/159 (91%)	137 (94%)	8 (6%)	0	100	100
1	HH	142/159 (89%)	135 (95%)	7 (5%)	0	100	100
1	I	142/159 (89%)	134 (94%)	8 (6%)	0	100	100
1	II	141/159 (89%)	129 (92%)	11 (8%)	1 (1%)	18	22
1	J	145/159 (91%)	135 (93%)	10 (7%)	0	100	100
1	JJ	145/159 (91%)	137 (94%)	8 (6%)	0	100	100
1	K	144/159 (91%)	133 (92%)	11 (8%)	0	100	100
1	KK	143/159 (90%)	137 (96%)	6 (4%)	0	100	100
1	L	155/159 (98%)	137 (88%)	16 (10%)	2 (1%)	9	10
1	M	142/159 (89%)	132 (93%)	10 (7%)	0	100	100
1	N	142/159 (89%)	131 (92%)	10 (7%)	1 (1%)	18	22
1	O	148/159 (93%)	138 (93%)	9 (6%)	1 (1%)	18	22
All	All	2876/3180 (90%)	2687 (93%)	179 (6%)	10 (0%)	36	45

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	N	122	VAL
1	II	129	ASN
1	L	12	SER
1	L	16	ASN
1	A	17	SER
1	B	37	TRP
1	G	17	SER
1	O	12	SER
1	A	15	ASP
1	A	16	ASN

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	129/140 (92%)	129 (100%)	0	100	100
1	B	127/140 (91%)	127 (100%)	0	100	100
1	C	127/140 (91%)	127 (100%)	0	100	100
1	D	128/140 (91%)	128 (100%)	0	100	100
1	E	129/140 (92%)	129 (100%)	0	100	100
1	F	130/140 (93%)	130 (100%)	0	100	100
1	G	129/140 (92%)	129 (100%)	0	100	100
1	GG	128/140 (91%)	127 (99%)	1 (1%)	73	85
1	H	130/140 (93%)	130 (100%)	0	100	100
1	HH	128/140 (91%)	128 (100%)	0	100	100
1	I	128/140 (91%)	128 (100%)	0	100	100
1	II	127/140 (91%)	127 (100%)	0	100	100
1	J	130/140 (93%)	130 (100%)	0	100	100
1	JJ	130/140 (93%)	130 (100%)	0	100	100
1	K	129/140 (92%)	129 (100%)	0	100	100
1	KK	129/140 (92%)	129 (100%)	0	100	100
1	L	138/140 (99%)	138 (100%)	0	100	100
1	M	128/140 (91%)	128 (100%)	0	100	100
1	N	128/140 (91%)	127 (99%)	1 (1%)	73	85
1	O	133/140 (95%)	133 (100%)	0	100	100
All	All	2585/2800 (92%)	2583 (100%)	2 (0%)	88	94

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	N	122	VAL
1	GG	51	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	ASN
1	B	159	ASN
1	D	48	GLN
1	G	16	ASN
1	I	117	ASN
1	J	117	ASN
1	M	32	ASN
1	M	139	ASN
1	M	149	GLN
1	O	48	GLN
1	O	89	ASN
1	O	133	ASN
1	O	159	ASN
1	GG	48	GLN
1	GG	89	ASN
1	II	48	GLN
1	JJ	32	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	P	6/8 (75%)	0	0
2	TT	6/8 (75%)	5 (83%)	0
2	V	6/8 (75%)	1 (16%)	0
3	S	4/6 (66%)	1 (25%)	0
3	T	4/6 (66%)	1 (25%)	0
4	UU	7/10 (70%)	1 (14%)	0
4	X	7/10 (70%)	1 (14%)	1 (14%)
4	bb	8/10 (80%)	3 (37%)	0
5	Y	5/7 (71%)	0	0
5	a	5/7 (71%)	2 (40%)	0
6	WW	10/12 (83%)	7 (70%)	2 (20%)
6	e	11/12 (91%)	7 (63%)	0
7	h	7/10 (70%)	3 (42%)	0
8	i	4/9 (44%)	2 (50%)	0
8	ll	4/9 (44%)	1 (25%)	0
8	qq	4/9 (44%)	0	0
9	kk	5/8 (62%)	1 (20%)	0
9	m	6/8 (75%)	1 (16%)	0
9	n	6/8 (75%)	0	0

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
All	All	115/165 (69%)	37 (32%)	3 (2%)

All (37) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	S	164[A]	A
3	T	168[A]	A
2	V	169[A]	A
4	X	169[A]	A
5	a	168[A]	A
5	a	169[A]	A
6	e	180	U
6	e	181[A]	U
6	e	182[A]	U
6	e	183[A]	U
6	e	184[A]	U
6	e	185[A]	U
6	e	186[A]	U
7	h	184[A]	U
7	h	186[A]	U
7	h	187[A]	U
8	i	185[A]	U
8	i	186[A]	U
9	m	188[A]	U
2	TT	166[A]	A
2	TT	167[A]	A
2	TT	168[A]	A
2	TT	169[A]	A
2	TT	170[A]	A
4	UU	162[A]	A
6	WW	166[A]	U
6	WW	167[A]	U
6	WW	168	U
6	WW	171	U
6	WW	172	U
6	WW	173	U
6	WW	174	U
4	bb	161[A]	A
4	bb	163[A]	A
4	bb	168[A]	A
9	kk	184[A]	U
8	ll	188[A]	U

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	X	168[A]	A
6	WW	165	U
6	WW	173	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 6 are monoatomic - leaving 12 for Mogul analysis.

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	$\# Z > 2$
13	SO4	JJ	203	-	4,4,4	0.23	0	6,6,6	0.07	0
10	PO4	HH	401	-	4,4,4	0.97	0	6,6,6	0.46	0
10	PO4	B	701	-	4,4,4	0.97	0	6,6,6	0.46	0
10	PO4	H	501	-	4,4,4	0.97	0	6,6,6	0.45	0
13	SO4	F	201	-	4,4,4	0.24	0	6,6,6	1.26	0
13	SO4	M	202	-	4,4,4	0.24	0	6,6,6	0.14	0
13	SO4	M	201	-	4,4,4	0.24	0	6,6,6	0.10	0
10	PO4	R	1	-	4,4,4	0.95	0	6,6,6	0.47	0
13	SO4	KK	201	-	4,4,4	0.24	0	6,6,6	1.57	1 (16%)
10	PO4	L	901	-	4,4,4	0.97	0	6,6,6	0.46	0
13	SO4	O	201	-	4,4,4	0.24	0	6,6,6	0.08	0
10	PO4	II	201	-	4,4,4	0.95	0	6,6,6	1.14	0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	KK	201	SO4	O3-S-O1	3.22	126.37	109.56

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 [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	146/159 (91%)	-1.69	0 100 100	4, 7, 11, 19	0
1	B	143/159 (89%)	-1.68	0 100 100	4, 6, 9, 11	0
1	C	143/159 (89%)	-1.67	0 100 100	3, 4, 7, 19	0
1	D	144/159 (90%)	-1.73	0 100 100	2, 4, 7, 14	0
1	E	145/159 (91%)	-1.71	0 100 100	4, 5, 8, 15	0
1	F	147/159 (92%)	-1.67	0 100 100	4, 8, 10, 17	0
1	G	146/159 (91%)	-1.70	0 100 100	3, 6, 10, 12	0
1	GG	144/159 (90%)	-1.69	0 100 100	2, 3, 7, 14	0
1	H	147/159 (92%)	-1.70	0 100 100	5, 7, 10, 18	0
1	HH	144/159 (90%)	-1.69	0 100 100	3, 4, 7, 15	0
1	I	144/159 (90%)	-1.70	0 100 100	7, 10, 12, 19	0
1	II	143/159 (89%)	-1.69	0 100 100	2, 4, 7, 12	0
1	J	147/159 (92%)	-1.65	0 100 100	8, 10, 13, 22	0
1	JJ	147/159 (92%)	-1.72	0 100 100	2, 4, 8, 18	0
1	K	146/159 (91%)	-1.64	0 100 100	7, 9, 13, 23	0
1	KK	145/159 (91%)	-1.73	0 100 100	2, 3, 7, 16	0
1	L	157/159 (98%)	-1.58	1 (0%) 85 86	9, 11, 17, 22	0
1	M	144/159 (90%)	-1.65	0 100 100	10, 11, 15, 21	0
1	N	144/159 (90%)	-1.65	0 100 100	6, 10, 14, 16	0
1	O	150/159 (94%)	-1.62	0 100 100	5, 8, 13, 26	0
2	P	7/8 (87%)	-1.77	0 100 100	16, 18, 23, 28	0
2	TT	7/8 (87%)	-1.32	0 100 100	17, 20, 24, 29	0
2	V	7/8 (87%)	-1.87	0 100 100	14, 17, 22, 23	0
3	S	5/6 (83%)	-1.86	0 100 100	15, 16, 19, 21	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
3	T	5/6 (83%)	-2.14	0 100 100	16, 17, 18, 19	0
4	UU	8/10 (80%)	-1.51	0 100 100	17, 18, 26, 26	0
4	X	8/10 (80%)	-1.76	0 100 100	21, 22, 26, 26	0
4	bb	9/10 (90%)	-1.53	0 100 100	14, 19, 25, 27	0
5	Y	6/7 (85%)	-1.71	0 100 100	20, 22, 23, 23	0
5	a	6/7 (85%)	-1.06	0 100 100	20, 23, 29, 31	0
6	WW	10/12 (83%)	-1.18	0 100 100	16, 24, 29, 30	0
6	e	12/12 (100%)	-1.47	0 100 100	19, 23, 28, 28	0
7	h	8/10 (80%)	-1.61	0 100 100	14, 17, 24, 26	0
8	i	5/9 (55%)	-1.90	0 100 100	15, 15, 19, 23	0
8	ll	5/9 (55%)	-1.88	0 100 100	16, 16, 22, 23	0
8	qq	5/9 (55%)	-1.71	0 100 100	14, 15, 22, 22	0
9	kk	6/8 (75%)	-1.56	0 100 100	17, 19, 23, 28	0
9	m	7/8 (87%)	-1.68	0 100 100	17, 20, 28, 28	0
9	n	7/8 (87%)	-1.60	0 100 100	19, 20, 24, 25	0
All	All	3049/3345 (91%)	-1.67	1 (0%) 100 100	2, 7, 16, 31	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	6	VAL	2.7

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](#)

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(Å ²)	Q<0.9
13	SO4	M	202	5/5	0.89	0.07	31,31,34,39	5
13	SO4	M	201	5/5	0.95	0.14	35,40,44,45	5
13	SO4	O	201	5/5	0.97	0.10	30,30,31,34	5
11	MG	G	201	1/1	0.98	0.19	14,14,14,14	0
10	PO4	R	1	5/5	0.98	0.07	42,49,49,49	5
13	SO4	KK	201	5/5	0.98	0.10	24,28,29,30	5
11	MG	JJ	202	1/1	0.99	0.04	4,4,4,4	0
11	MG	B	702	1/1	0.99	0.37	4,4,4,4	0
10	PO4	II	201	5/5	0.99	0.10	23,26,26,33	5
11	MG	H	502	1/1	0.99	0.04	7,7,7,7	0
13	SO4	JJ	203	5/5	0.99	0.05	9,9,12,18	5
11	MG	JJ	201	1/1	0.99	0.07	6,6,6,6	0
13	SO4	F	201	5/5	1.00	0.04	11,12,13,14	5
10	PO4	HH	401	5/5	1.00	0.02	2,2,2,2	0
10	PO4	B	701	5/5	1.00	0.02	3,4,4,4	0
10	PO4	H	501	5/5	1.00	0.03	7,7,7,7	0
10	PO4	L	901	5/5	1.00	0.02	9,10,11,15	0
12	CL	C	201	1/1	1.00	0.02	13,13,13,13	0

6.5 Other polymers

There are no such residues in this entry.