



wwPDB EM Validation Summary Report ⓘ

Mar 10, 2026 – 11:36 AM UTC

PDB ID : 6K32 / pdb_00006k32
EMDB ID : EMD-9907
Title : RdRp complex
Authors : Li, X.W.
Deposited on : 2019-05-16
Resolution : 3.20 Å(reported)

This is a wwPDB EM Validation Summary 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/EMValidationReportHelp>

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:

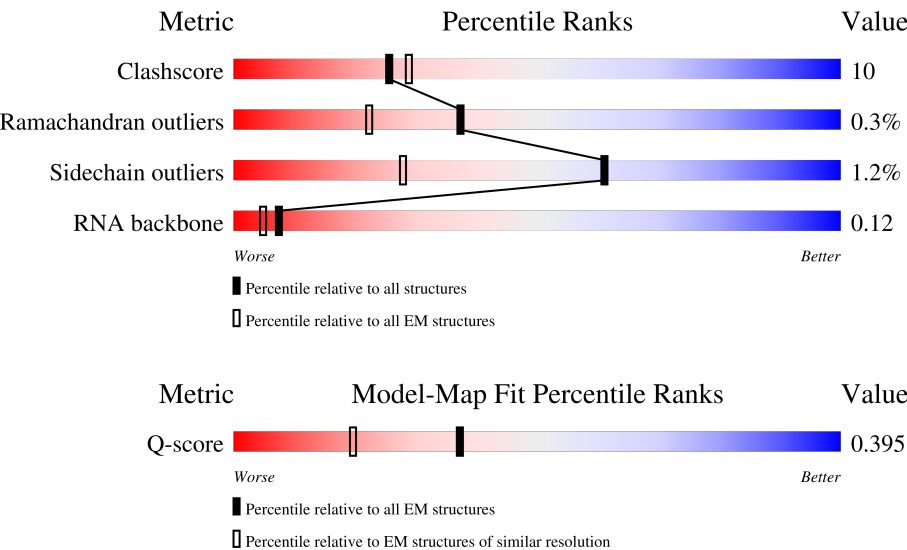
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.20 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.



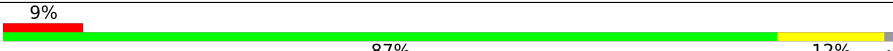
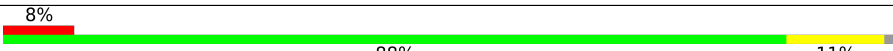
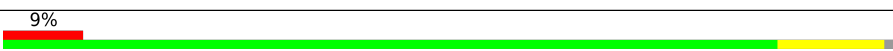
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	15020 (2.70 - 3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1208	
2	B	559	
3	t	5	

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Mol	Chain	Length	Quality of chain
4	p	3	 33% 67%
5	C	1220	 8% 87% 12%
6	D	1205	 9% 87% 12%
6	E	1205	 8% 88% 11%
6	F	1205	 7% 89% 10%
7	G	1226	 9% 87% 12%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
12	UTP	A	1306	-	-	X	-
9	MG7	A	1302	-	-	X	-

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 61364 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 RNA-dependent RNA polymerase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1208	Total	C	N	O	S	0	0
			9652	6124	1668	1824	36		

- Molecule 2 is a protein called Viral structural protein 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	498	Total	C	N	O	S	0	0
			3981	2552	662	749	18		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	t	5	Total	C	N	O	P	0	0
			103	46	14	38	5		

- Molecule 4 is a RNA chain called RNA (5'-D(*(3PO))-R(*AP*G)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
4	p	3	Total	C	N	O	P	0	0
			54	20	10	20	4		

- Molecule 5 is a protein called VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	C	1220	Total	C	N	O	S	0	0
			9622	6072	1676	1836	38		

- Molecule 6 is a protein called VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	D	1197	Total	C	N	O	S	0	0
			9447	5966	1645	1798	38		

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Mol	Chain	Residues	Atoms					AltConf	Trace
6	E	1191	Total	C	N	O	S	0	0
			9402	5937	1637	1791	37		
6	F	1191	Total	C	N	O	S	0	0
			9402	5937	1637	1791	37		

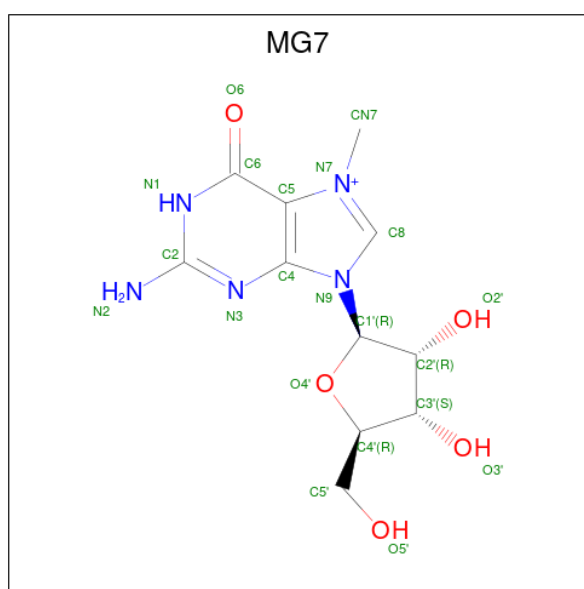
- Molecule 7 is a protein called VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	1218	Total	C	N	O	S	0	0
			9618	6073	1675	1832	38		

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

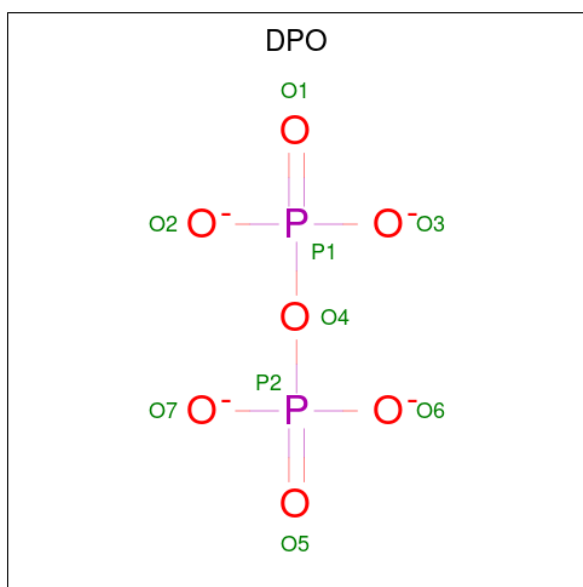
Mol	Chain	Residues	Atoms		AltConf
8	A	2	Total	Mg	0
			2	2	

- Molecule 9 is 7-METHYLGUANOSINE (CCD ID: MG7) (formula: C₁₁H₁₆N₅O₅).



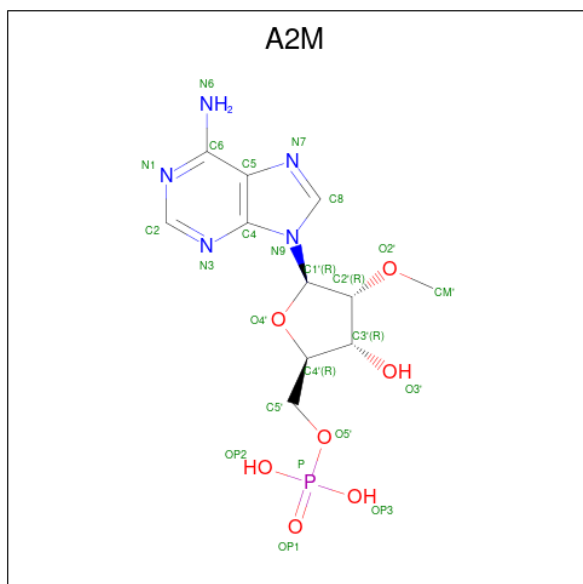
Mol	Chain	Residues	Atoms				AltConf
9	A	1	Total	C	N	O	0
			21	11	5	5	

- Molecule 10 is DIPHOSPHATE (CCD ID: DPO) (formula: O₇P₂).



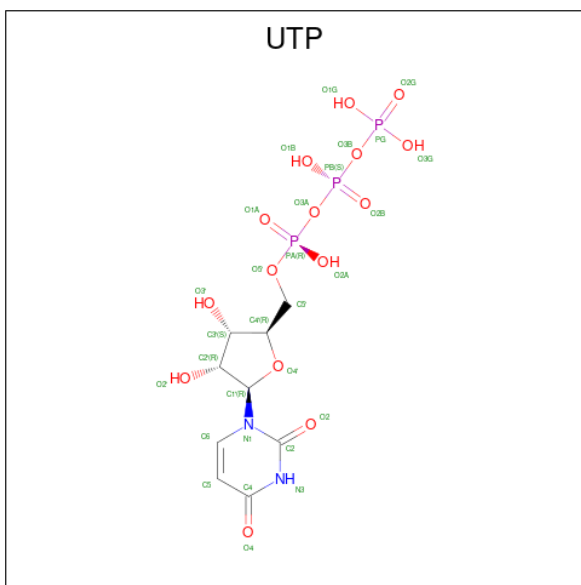
Mol	Chain	Residues	Atoms			AltConf
10	A	1	Total	O	P	0
			8	6	2	

- Molecule 11 is 2'-O-methyladenosine 5'-(dihydrogen phosphate) (CCD ID: A2M) (formula: $C_{11}H_{16}N_5O_7P$).



Mol	Chain	Residues	Atoms					AltConf
11	A	1	Total	C	N	O	P	0
			23	11	5	6	1	

- Molecule 12 is URIDINE 5'-TRIPHOSPHATE (CCD ID: UTP) (formula: $C_9H_{15}N_2O_{15}P_3$).

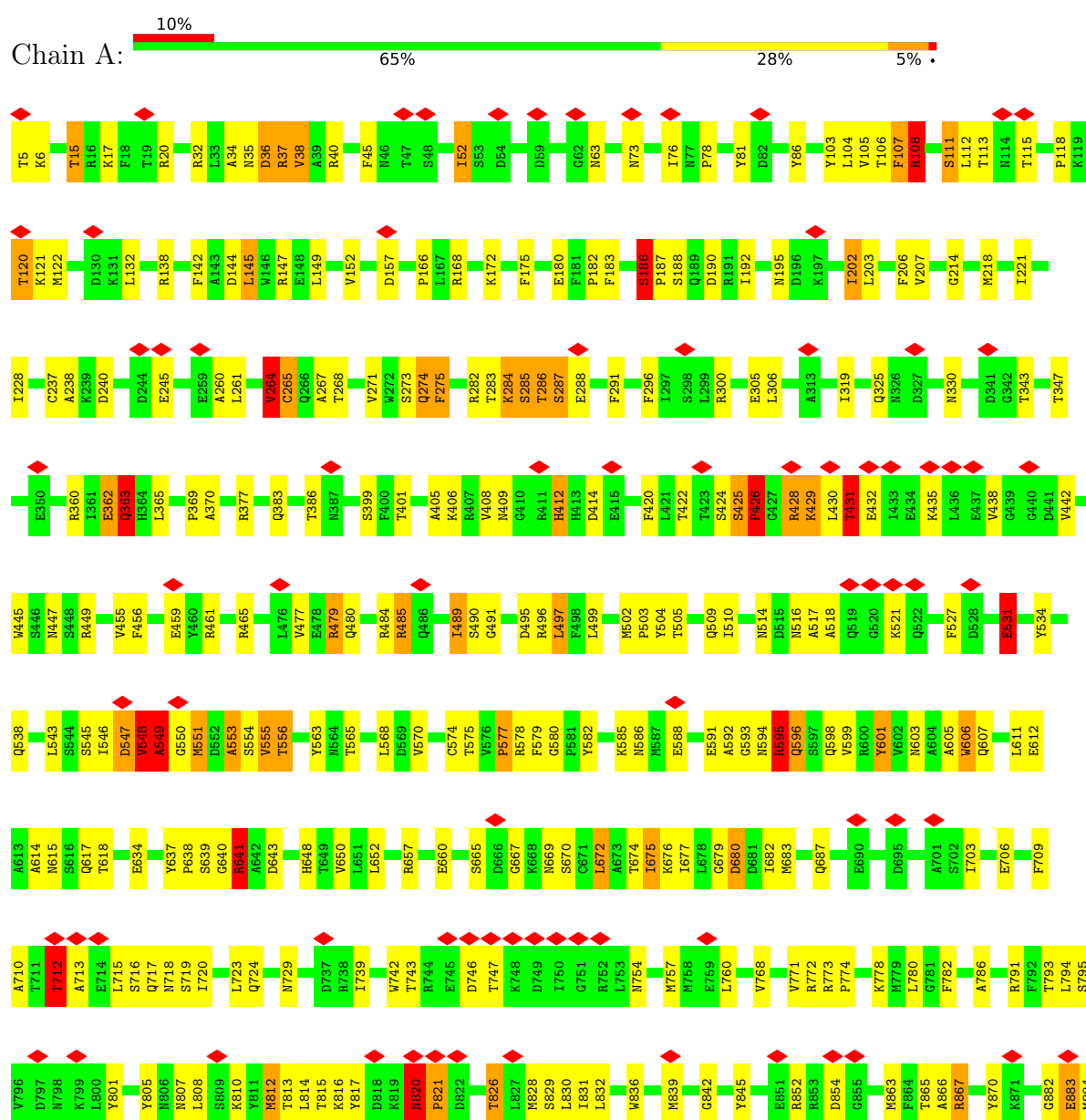


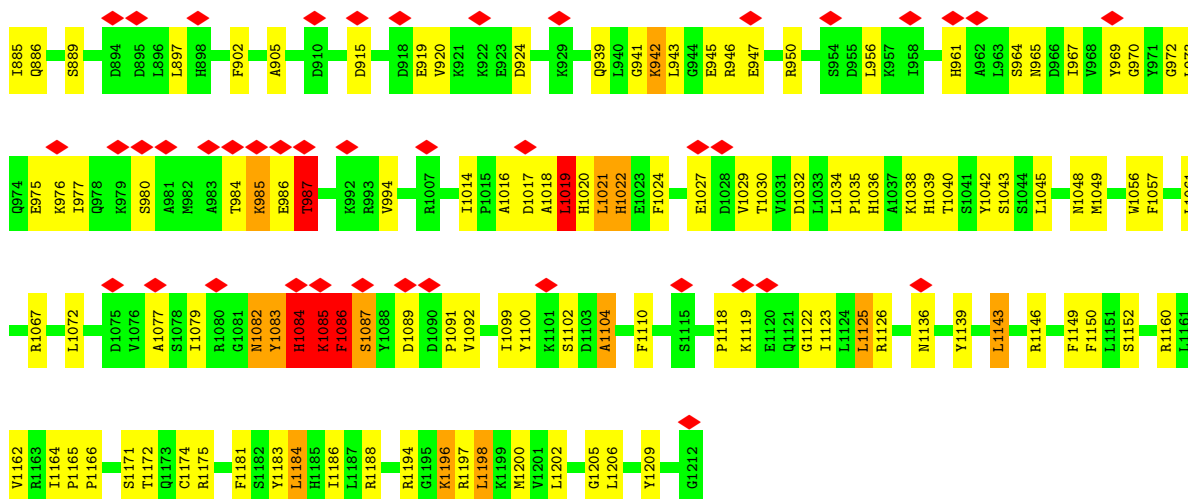
Mol	Chain	Residues	Atoms					AltConf
12	A	1	Total	C	N	O	P	0
			29	9	2	15	3	

3 Residue-property plots

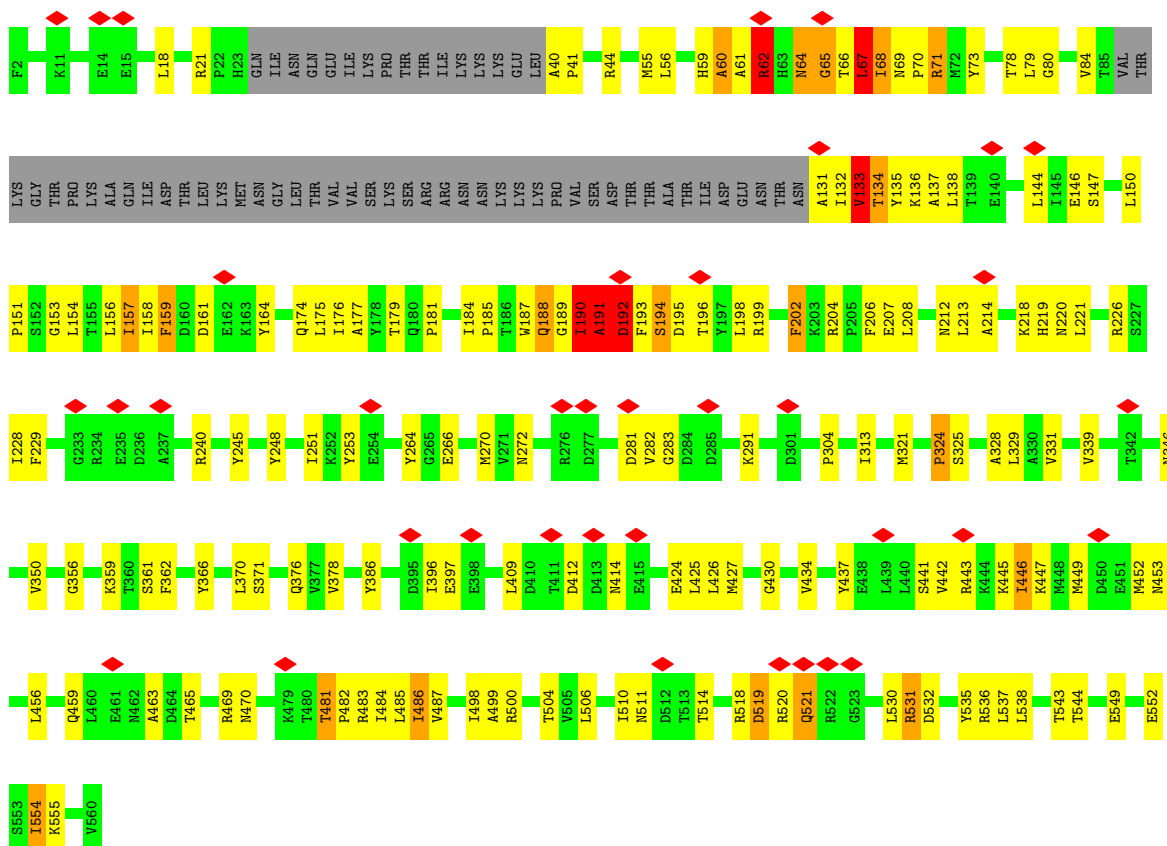
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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: RNA-dependent RNA polymerase





• Molecule 2: Viral structural protein 4



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



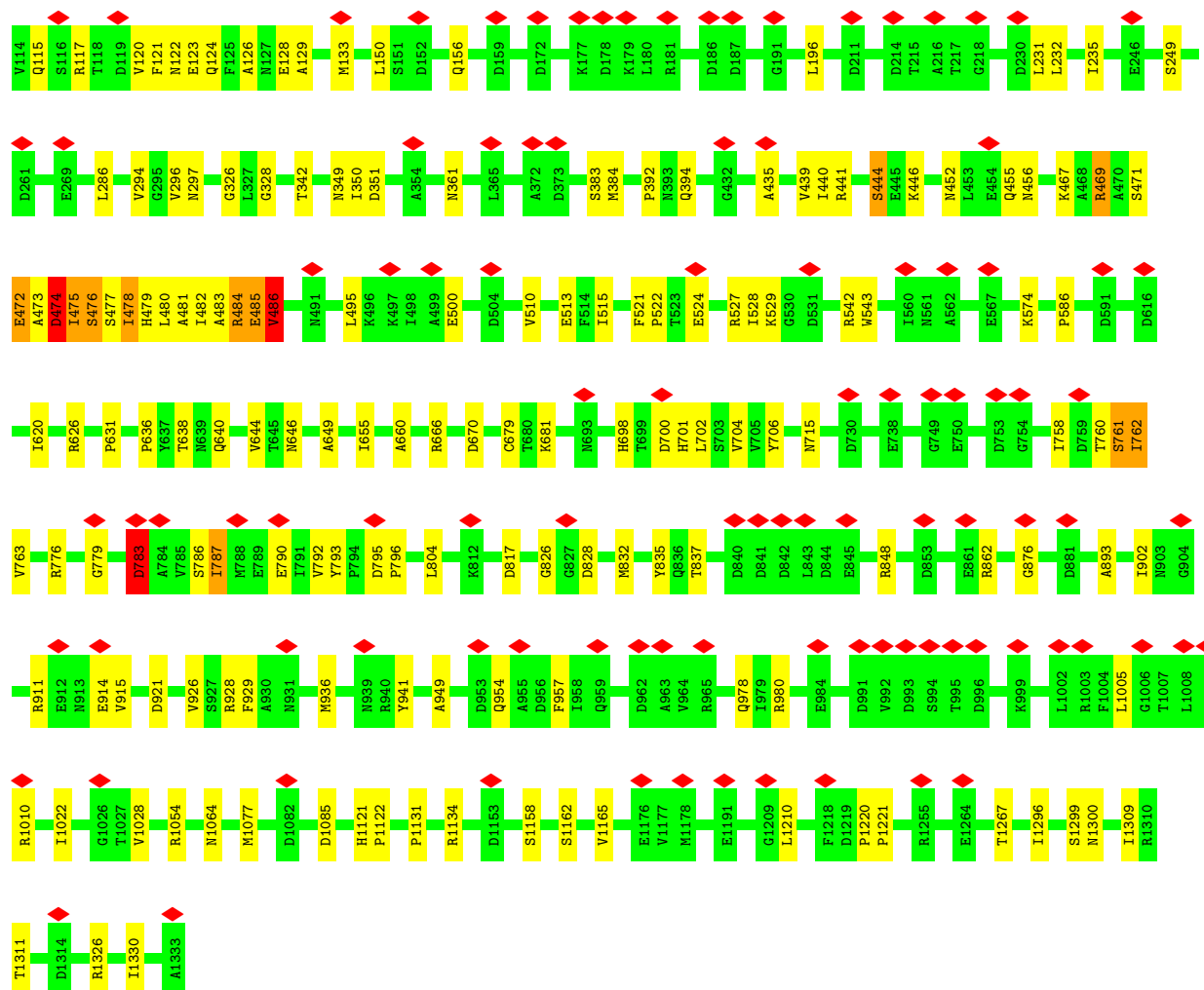
• Molecule 4: RNA (5'-D*(3PO))-R(*AP*G)-3'

Chain p: 



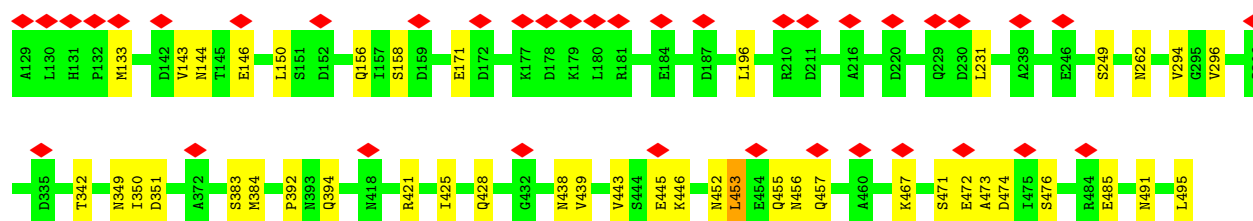
• Molecule 5: VP1

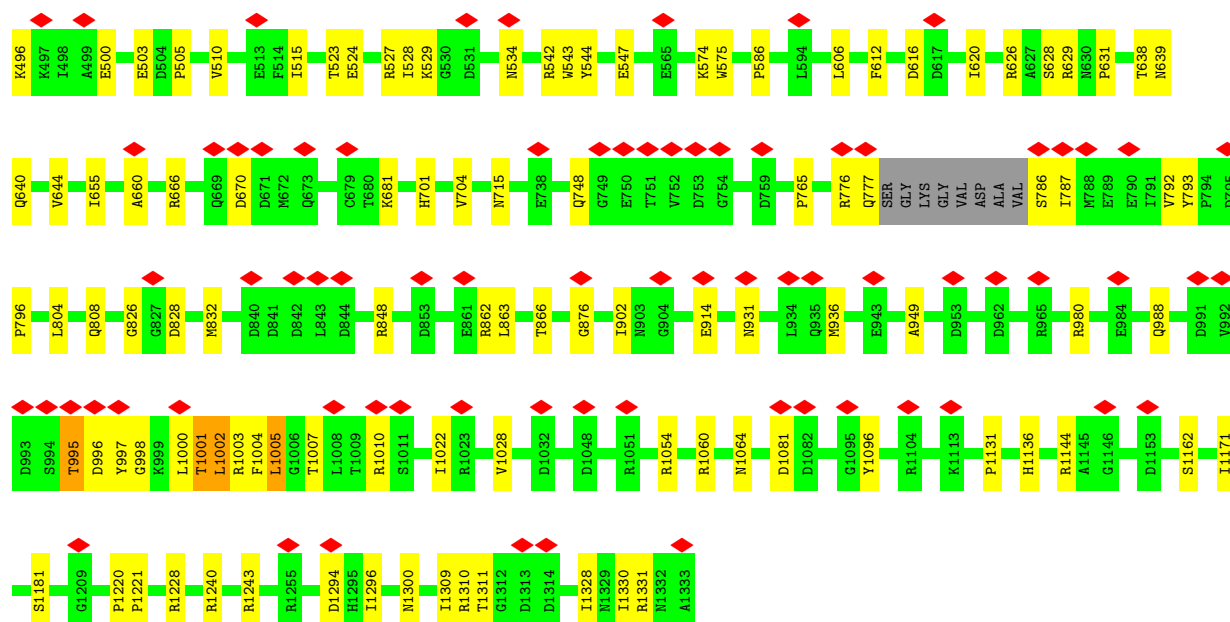
Chain C: 



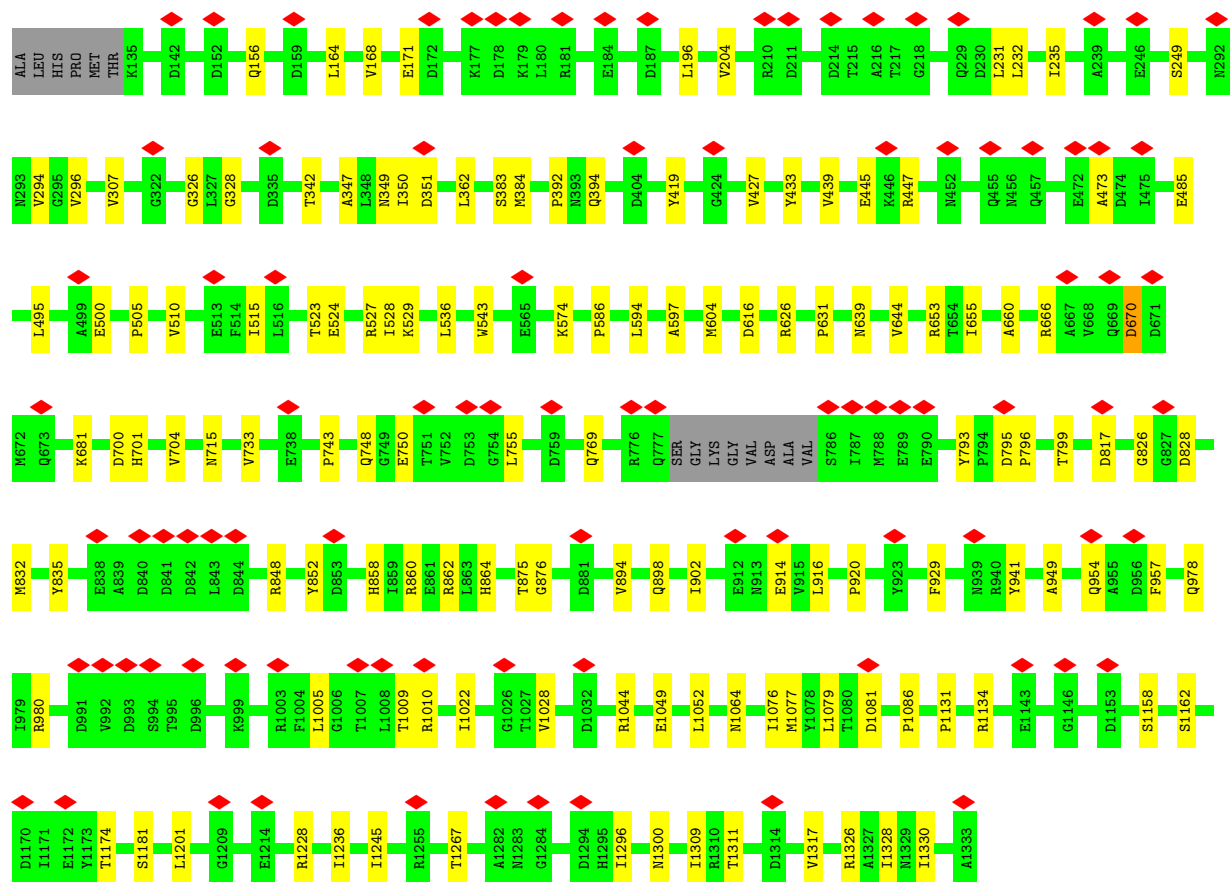
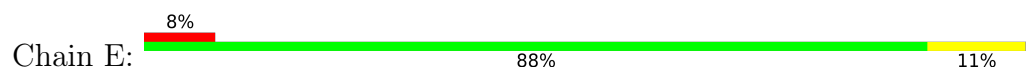
• Molecule 6: VP1

Chain D: 





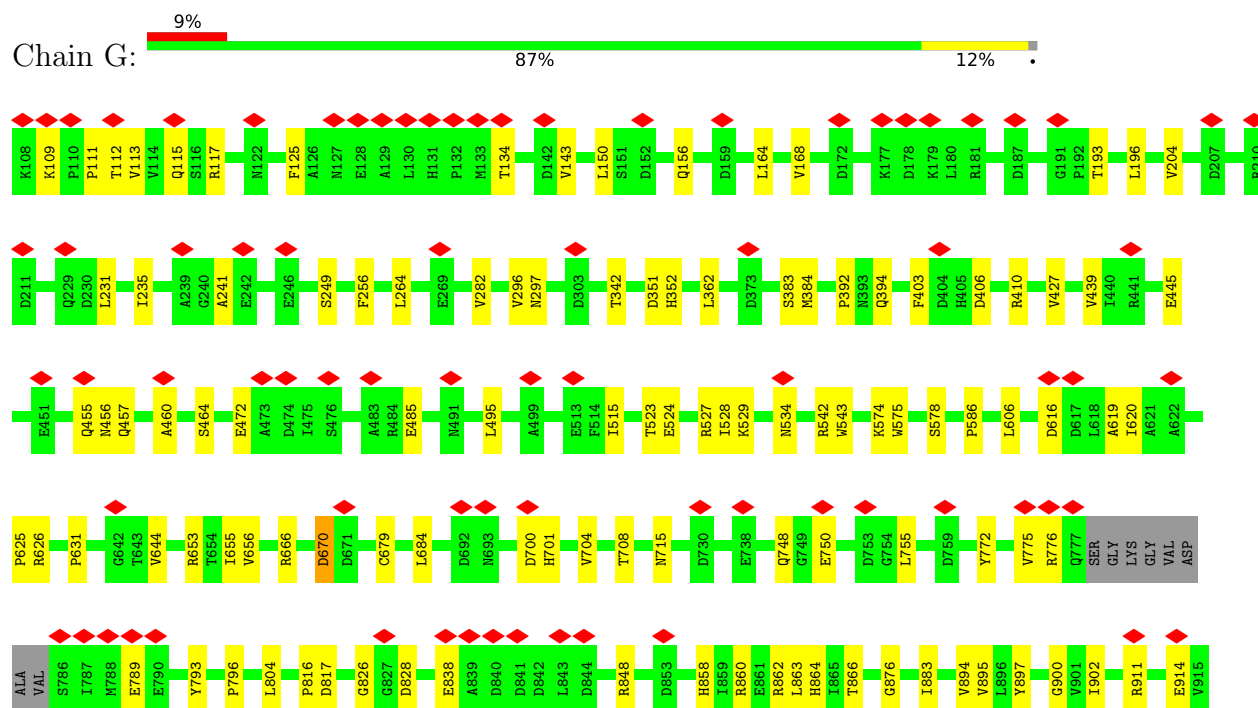
• Molecule 6: VP1

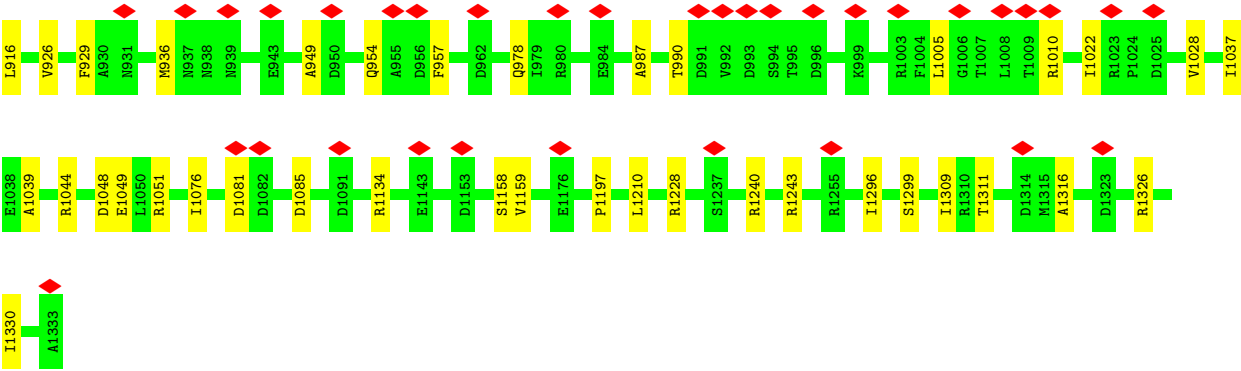


• Molecule 6: VP1



• Molecule 7: VP1





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	100000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TECNAI ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	20	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	161.942	Depositor
Minimum map value	-93.427	Depositor
Average map value	2.237	Depositor
Map value standard deviation	16.144	Depositor
Recommended contour level	25	Depositor
Map size ($\text{\AA}$)	285.75, 285.75, 285.75	wwPDB
Map dimensions	225, 225, 225	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.27, 1.27, 1.27	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: DPO, 3PO, MG, A2M, MG7, UTP

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	0.56	2/9859 (0.0%)	0.88	32/13346 (0.2%)
2	B	0.44	0/4067	0.82	9/5513 (0.2%)
3	t	5.40	22/113 (19.5%)	6.17	33/171 (19.3%)
4	p	1.63	1/46 (2.2%)	1.77	2/71 (2.8%)
5	C	0.32	0/9820	0.58	13/13367 (0.1%)
6	D	0.29	0/9642	0.53	1/13125 (0.0%)
6	E	0.26	0/9595	0.52	4/13060 (0.0%)
6	F	0.26	0/9595	0.51	5/13060 (0.0%)
7	G	0.26	0/9817	0.54	7/13363 (0.1%)
All	All	0.42	25/62554 (0.0%)	0.68	106/85076 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	43
2	B	0	21
5	C	0	7
6	D	0	2
7	G	0	3
All	All	0	76

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1083	TYR	C-N	30.86	1.59	1.33
3	t	1	U	C5'-C4'	-25.16	1.13	1.51
3	t	1	U	C4'-O4'	20.52	1.76	1.45
3	t	1	U	C3'-O3'	16.61	1.67	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	t	2	U	C3'-O3'	15.03	1.64	1.42

The worst 5 of 106 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	t	1	U	C2'-C3'-O3'	-30.26	68.31	113.70
3	t	1	U	C5'-C4'-C3'	27.35	157.03	116.00
3	t	1	U	C4'-C3'-O3'	21.61	145.42	113.00
3	t	2	U	C1'-C2'-O2'	-21.16	76.66	108.40
3	t	1	U	C5'-C4'-O4'	-18.94	81.39	109.80

There are no chirality outliers.

5 of 76 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	120	THR	Peptide
1	A	186	SER	Peptide
1	A	264	VAL	Peptide
1	A	285	SER	Peptide
1	A	52	ILE	Peptide

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	9652	0	9538	563	0
2	B	3981	0	3985	268	0
3	t	103	0	52	73	0
4	p	54	0	21	31	0
5	C	9622	0	9522	172	0
6	D	9447	0	9360	169	0
6	E	9402	0	9314	74	0
6	F	9402	0	9314	77	0
7	G	9618	0	9526	101	0
8	A	2	0	0	0	0
9	A	21	0	5	7	0
10	A	8	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
11	A	23	0	10	5	0
12	A	29	0	9	17	0
All	All	61364	0	60656	1255	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

The worst 5 of 1255 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)
5:C:473:ALA:HB2	5:C:761:SER:CB	1.42	1.49
3:t:2:U:C3'	3:t:2:U:O3'	1.64	1.43
3:t:1:U:O3'	3:t:1:U:C3'	1.67	1.42
1:A:1077:ALA:HB2	1:A:1086:PHE:CE2	1.55	1.39
2:B:185:PRO:HG3	6:D:1000:LEU:CD2	1.51	1.36

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 PDB entries followed by that with respect to all EM entries.

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	1206/1208 (100%)	1091 (90%)	96 (8%)	19 (2%)	7	36
2	B	492/559 (88%)	437 (89%)	50 (10%)	5 (1%)	12	45
5	C	1218/1220 (100%)	1152 (95%)	64 (5%)	2 (0%)	43	73
6	D	1193/1205 (99%)	1127 (94%)	66 (6%)	0	100	100
6	E	1187/1205 (98%)	1120 (94%)	67 (6%)	0	100	100
6	F	1187/1205 (98%)	1124 (95%)	63 (5%)	0	100	100
7	G	1214/1226 (99%)	1133 (93%)	81 (7%)	0	100	100
All	All	7697/7828 (98%)	7184 (93%)	487 (6%)	26 (0%)	37	68

5 of 26 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	187	PRO
1	A	425	SER
1	A	432	GLU
1	A	549	ALA
1	A	555	VAL

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 PDB entries followed by that with respect to all EM entries.

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	1049/1049 (100%)	1000 (95%)	49 (5%)	23	57
2	B	435/492 (88%)	422 (97%)	13 (3%)	36	66
5	C	1063/1063 (100%)	1052 (99%)	11 (1%)	68	80
6	D	1044/1049 (100%)	1038 (99%)	6 (1%)	78	84
6	E	1039/1049 (99%)	1036 (100%)	3 (0%)	86	88
6	F	1039/1049 (99%)	1038 (100%)	1 (0%)	88	91
7	G	1064/1069 (100%)	1063 (100%)	1 (0%)	88	91
All	All	6733/6820 (99%)	6649 (99%)	84 (1%)	61	79

5 of 84 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	446	ILE
5	C	783	ASP
2	B	449	MET
5	C	472	GLU
6	D	995	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 120 such sidechains are listed below:

Mol	Chain	Res	Type
5	C	1234	GLN

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Mol	Chain	Res	Type
7	G	577	GLN
6	D	1015	GLN
7	G	455	GLN
7	G	1285	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	t	4/5 (80%)	3 (75%)	0
4	p	1/3 (33%)	1 (100%)	0
All	All	5/8 (62%)	4 (80%)	0

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	t	2	U
3	t	3	A
3	t	4	C
4	p	3	G

There are no RNA pucker outliers to report.

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 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z > 2$
11	A2M	A	1304	10	22,25,26	3.56	9 (40%)	30,36,39	5.34	17 (56%)
9	MG7	A	1302	10	23,23,23	7.51	11 (47%)	35,35,35	4.90	16 (45%)
10	DPO	A	1303	9,11	3,7,8	1.12	0	6,10,13	1.75	1 (16%)
12	UTP	A	1306	8	29,30,30	2.20	6 (20%)	43,47,47	2.40	13 (30%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	A2M	A	1304	10	-	4/9/27/28	0/3/3/3
9	MG7	A	1302	10	-	3/6/22/22	0/3/3/3
10	DPO	A	1303	9,11	-	0/2/5/6	-
12	UTP	A	1306	8	-	5/22/38/38	0/2/2/2

The worst 5 of 26 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	1302	MG7	O3'-C3'	-18.62	0.96	1.43
9	A	1302	MG7	CN7-N7	-16.86	1.18	1.46
9	A	1302	MG7	O2'-C2'	-13.23	1.10	1.43
9	A	1302	MG7	C3'-C2'	-11.63	1.21	1.53
9	A	1302	MG7	C3'-C4'	-9.75	1.28	1.53

The worst 5 of 47 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	A	1304	A2M	C1'-N9-C8	-16.61	90.24	127.09
9	A	1302	MG7	C1'-N9-C8	-13.74	80.36	126.74
11	A	1304	A2M	C2'-C1'-N9	13.41	135.82	113.75
9	A	1302	MG7	C1'-N9-C4	12.39	163.09	126.49
9	A	1302	MG7	O4'-C4'-C5'	-11.88	84.10	109.22

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

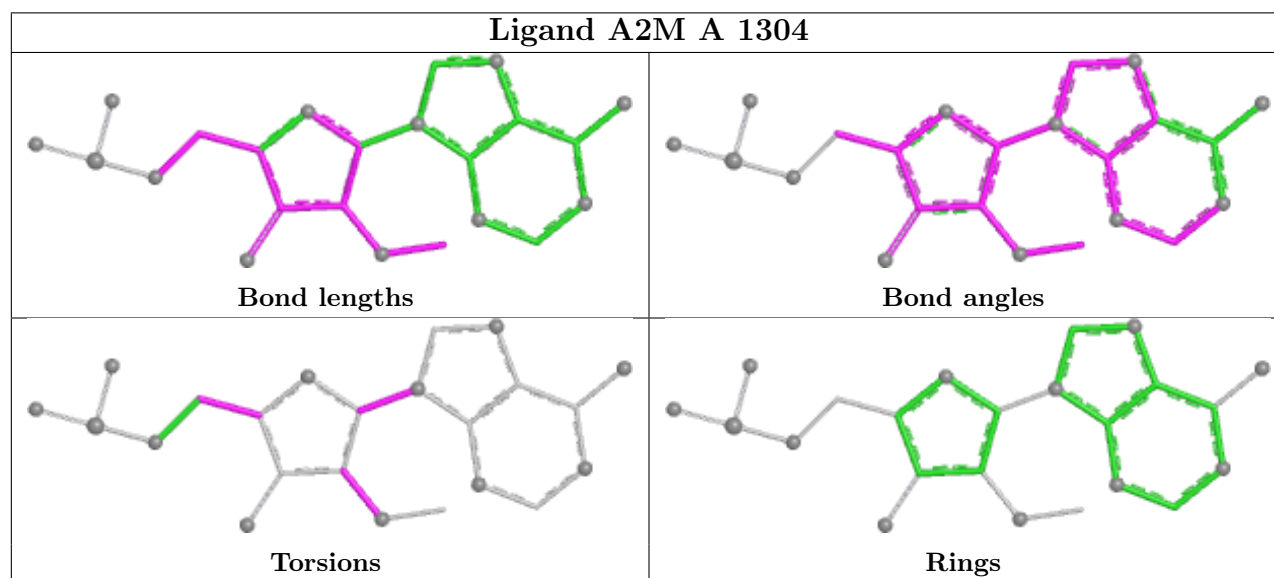
Mol	Chain	Res	Type	Atoms
9	A	1302	MG7	O4'-C1'-N9-C4
11	A	1304	A2M	C2'-C1'-N9-C8
12	A	1306	UTP	C5'-O5'-PA-O1A
12	A	1306	UTP	C5'-O5'-PA-O3A
12	A	1306	UTP	C2'-C1'-N1-C2

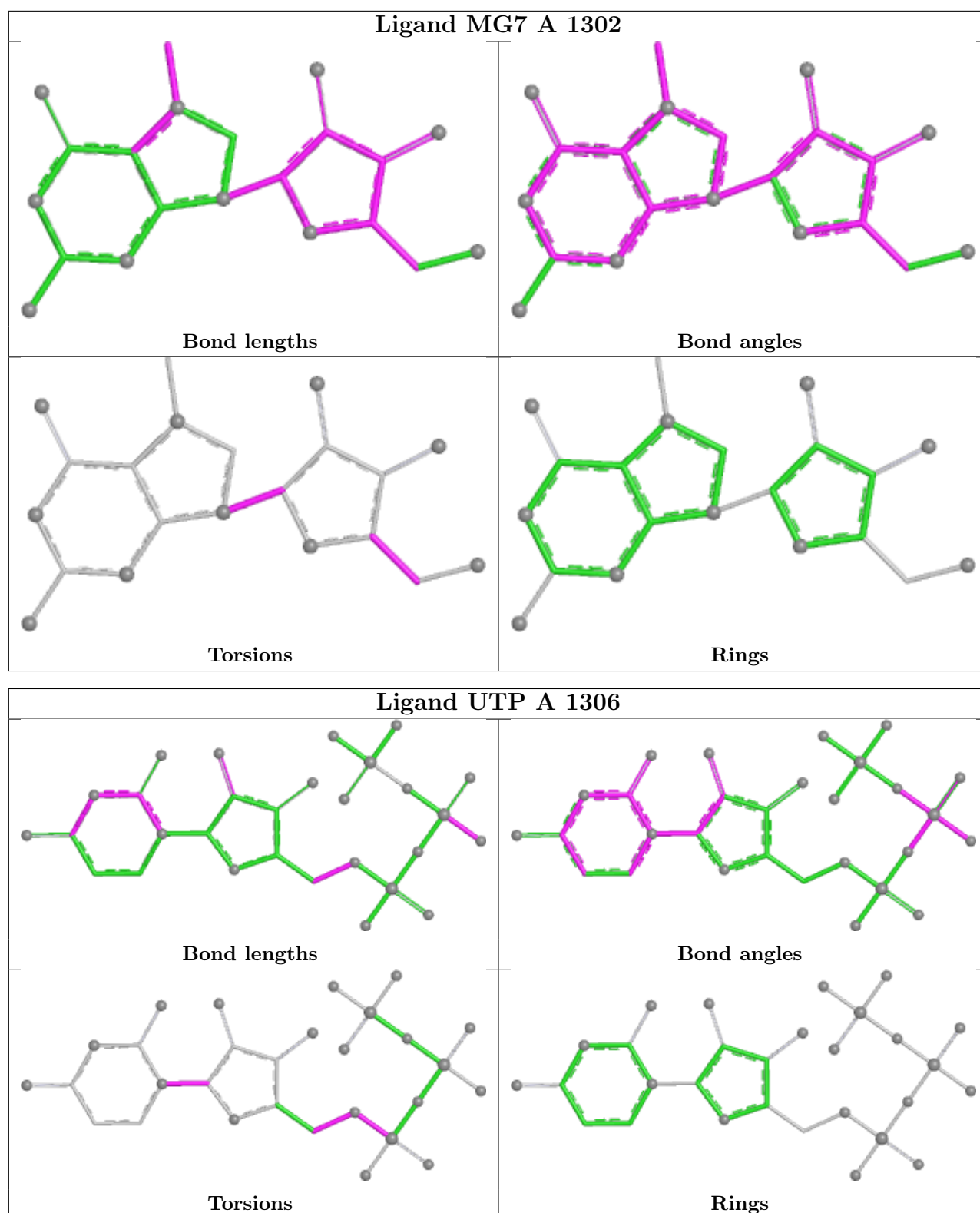
There are no ring outliers.

4 monomers are involved in 30 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	A	1304	A2M	5	0
9	A	1302	MG7	7	0
10	A	1303	DPO	1	0
12	A	1306	UTP	17	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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.

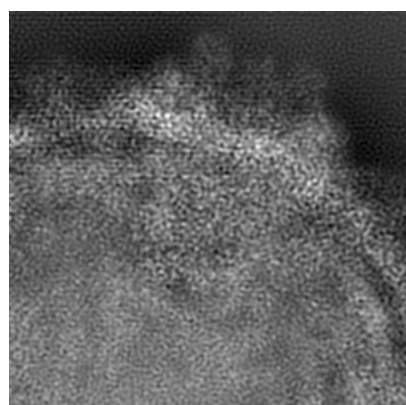
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9907. These allow visual inspection of the internal detail of the map and identification of artifacts.

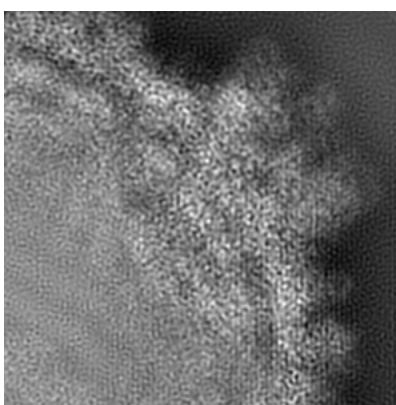
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

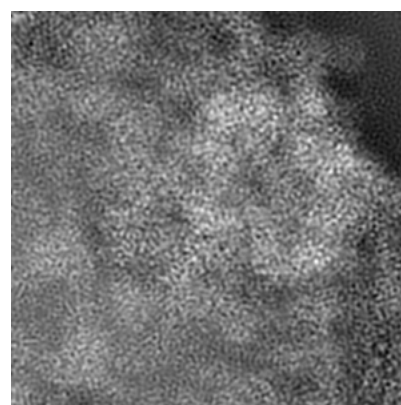
6.1.1 Primary map



X



Y

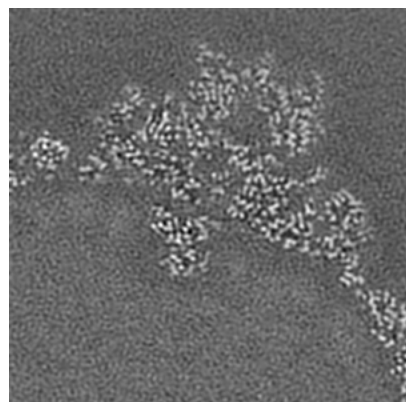


Z

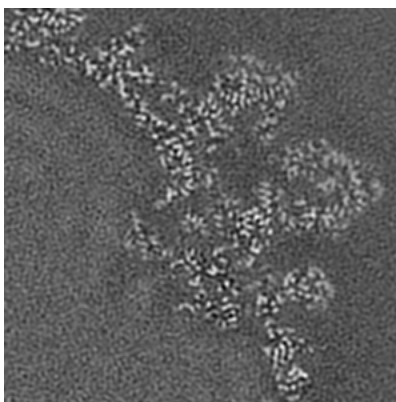
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

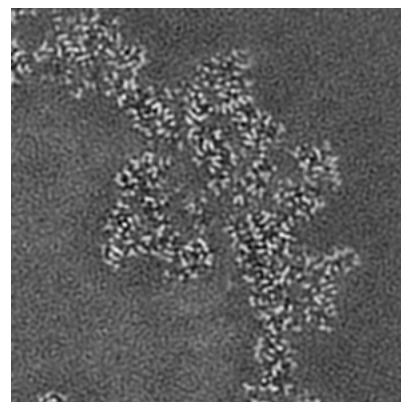
6.2.1 Primary map



X Index: 112



Y Index: 112

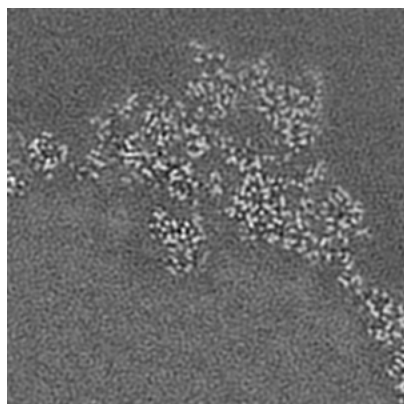


Z Index: 112

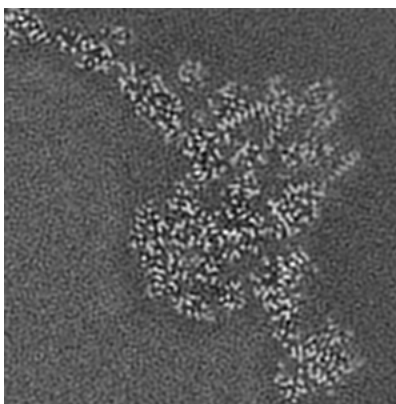
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

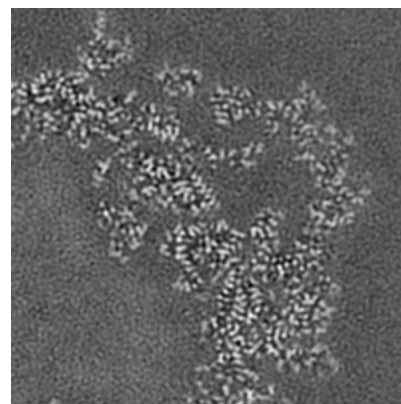
6.3.1 Primary map



X Index: 113



Y Index: 90

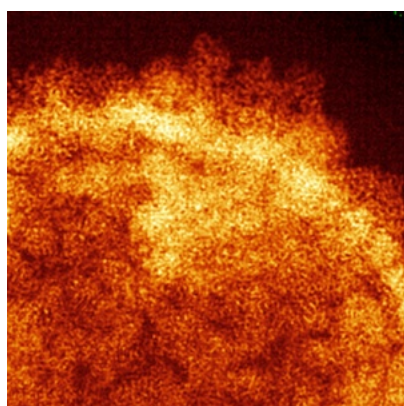


Z Index: 131

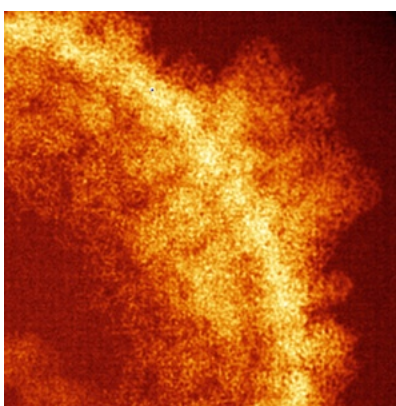
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

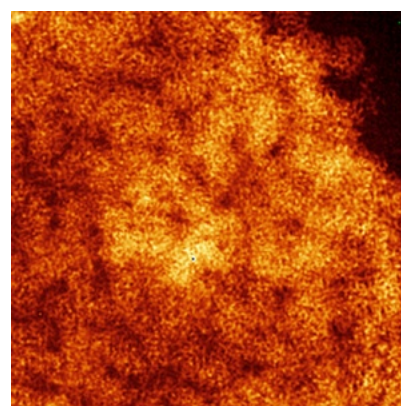
6.4.1 Primary map



X



Y

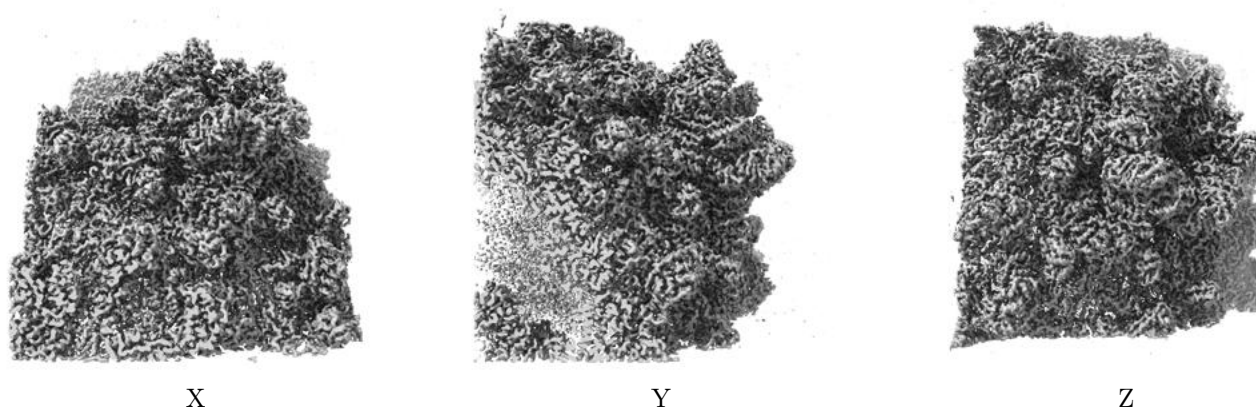


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 25.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

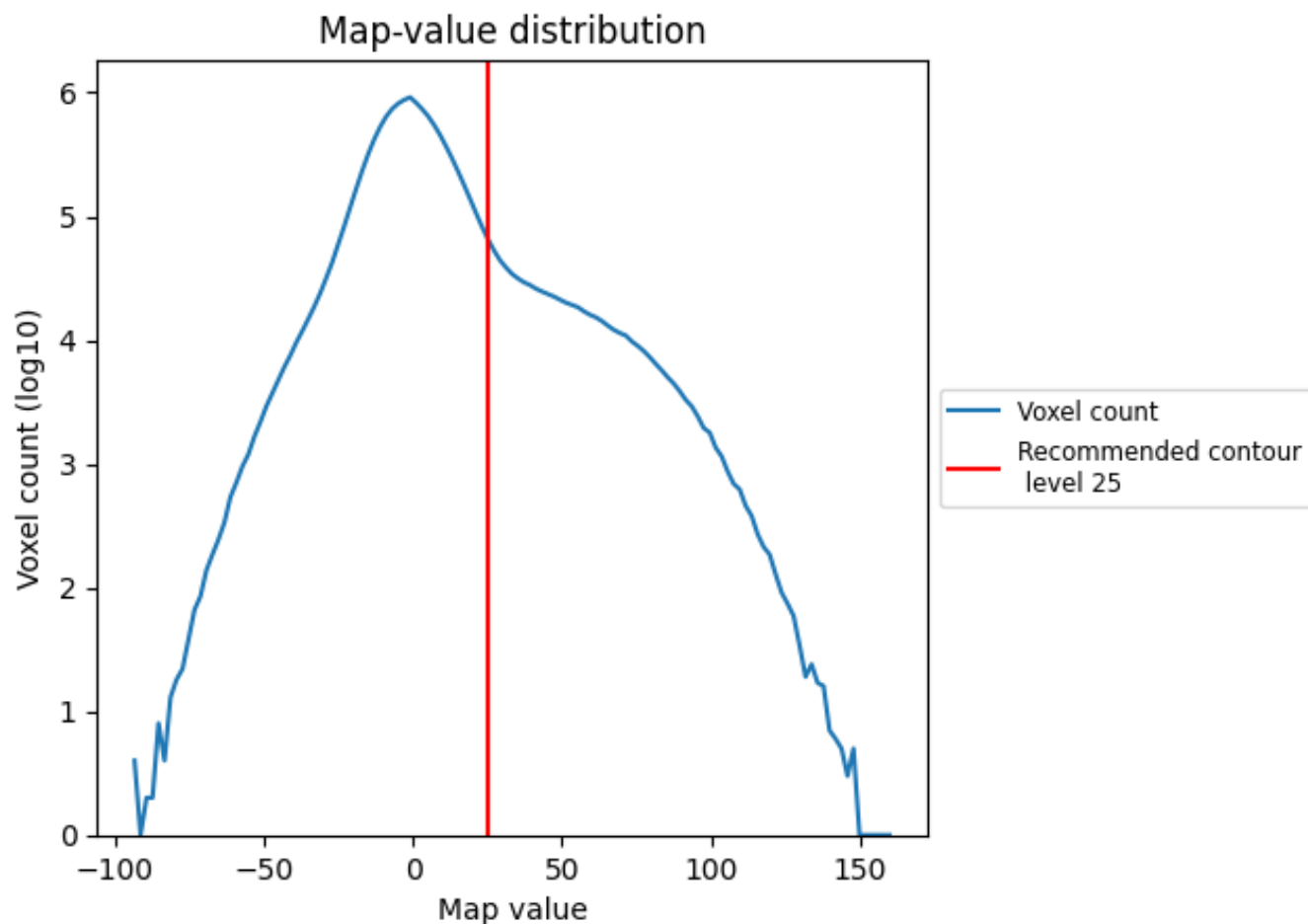
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

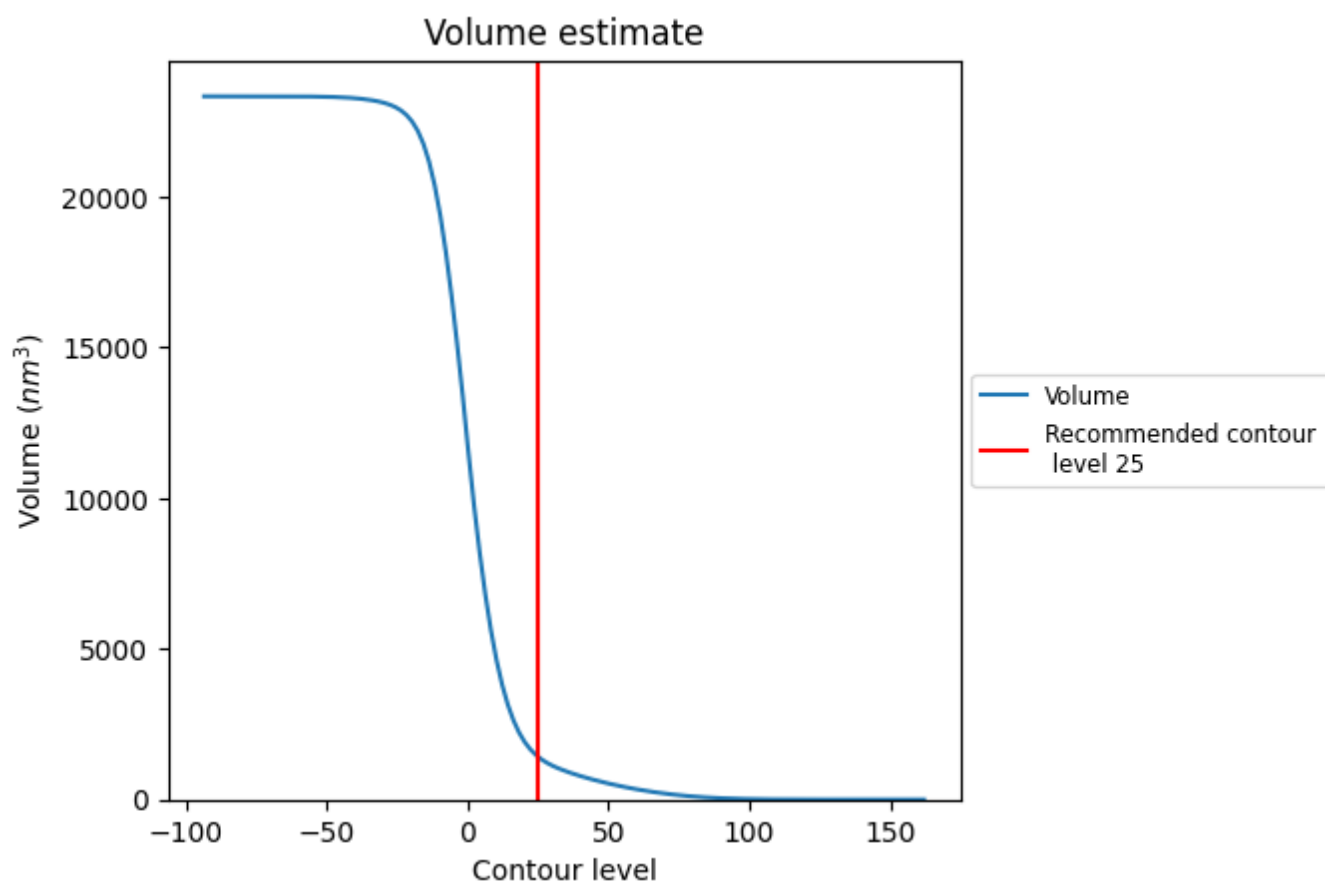
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

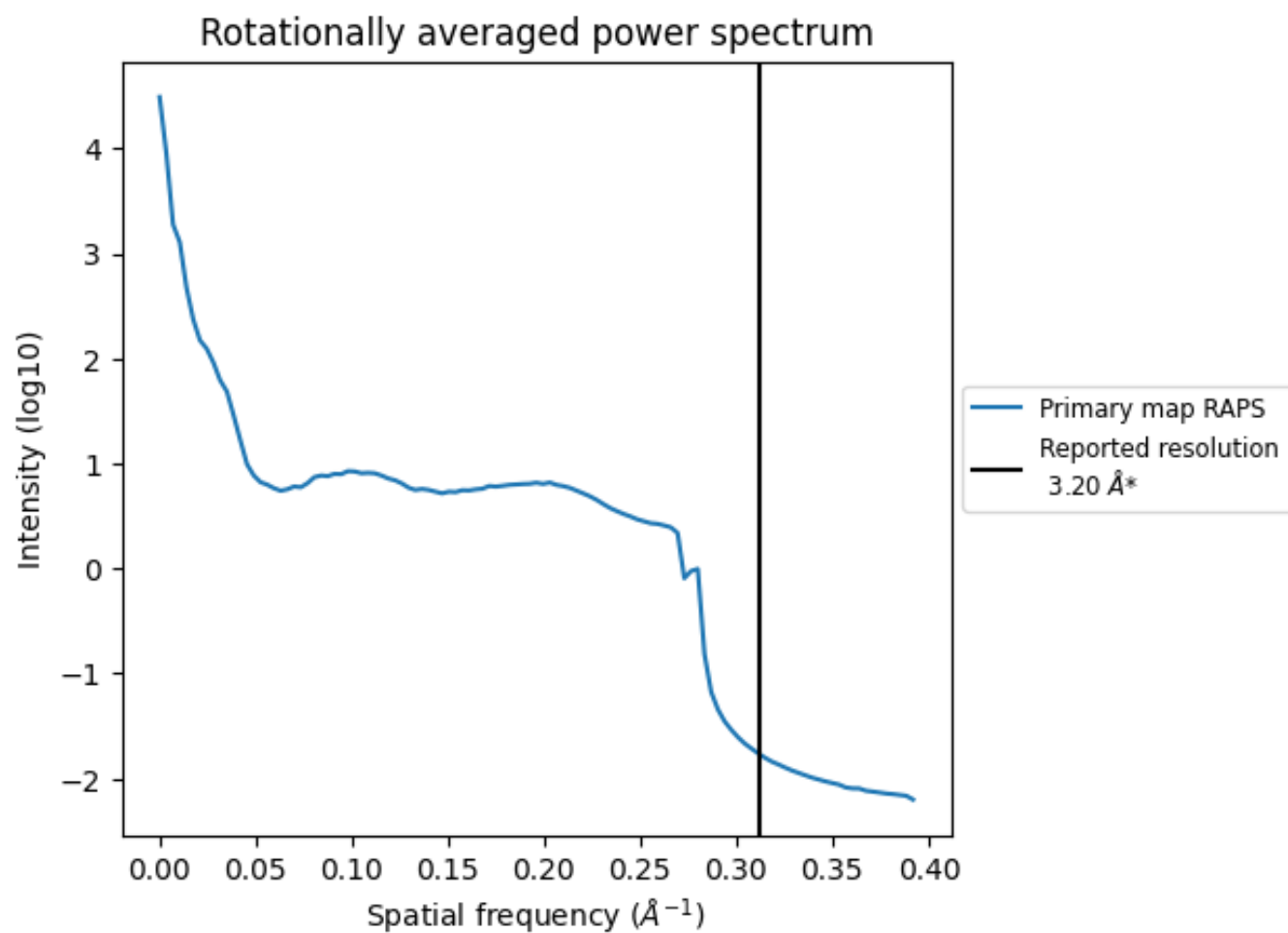
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1428 nm³; this corresponds to an approximate mass of 1290 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.312 Å⁻¹

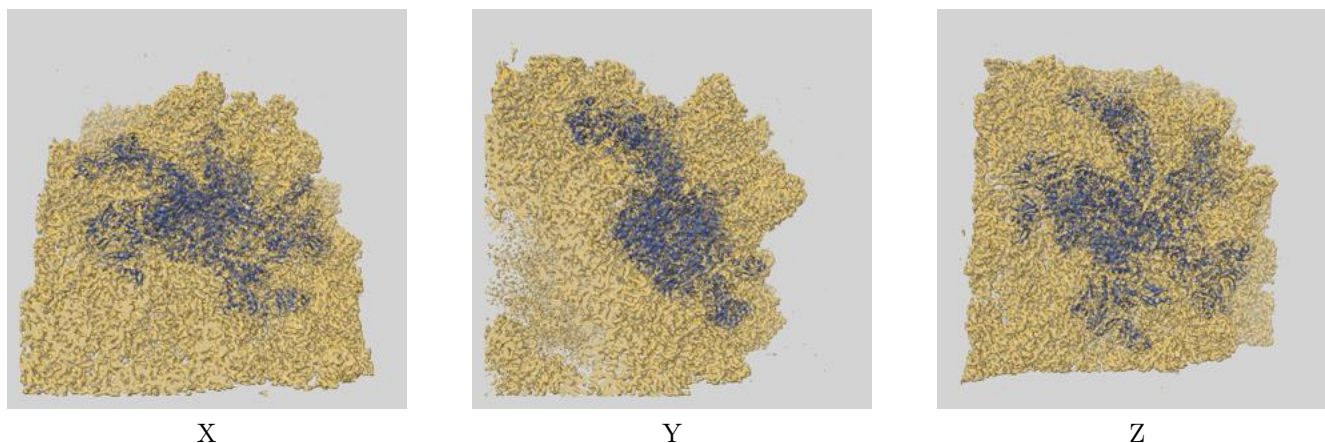
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

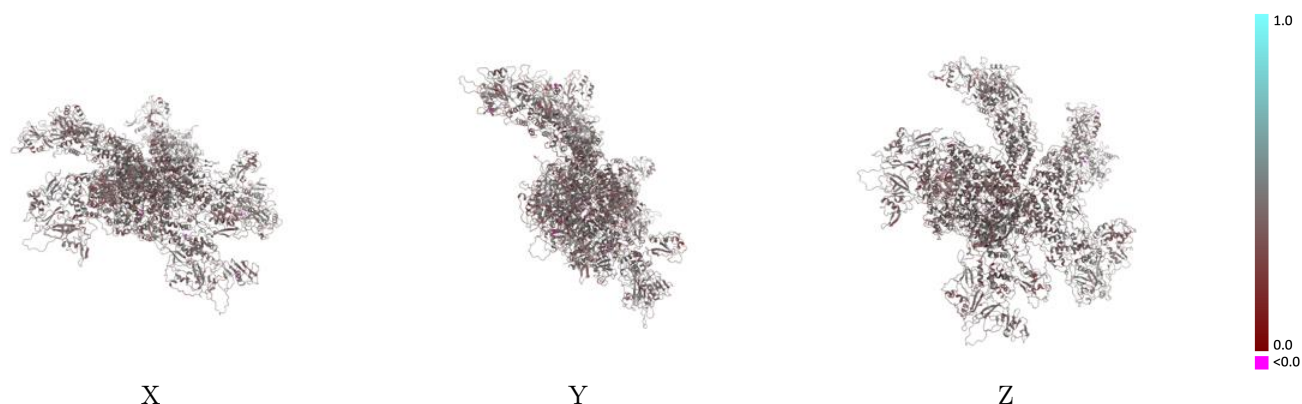
This section contains information regarding the fit between EMDB map EMD-9907 and PDB model 6K32. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

9.1 Map-model overlay [i](#)



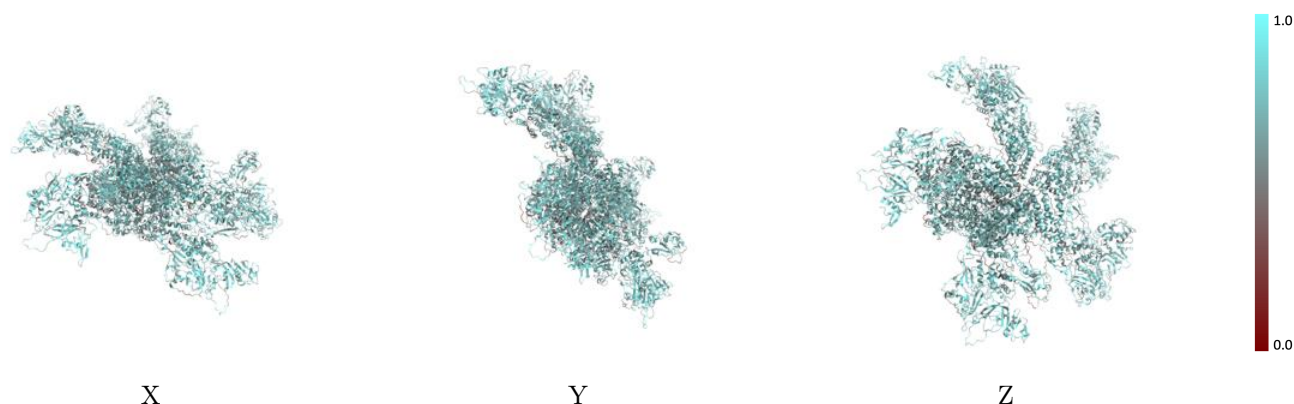
The images above show the 3D surface view of the map at the recommended contour level 25.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



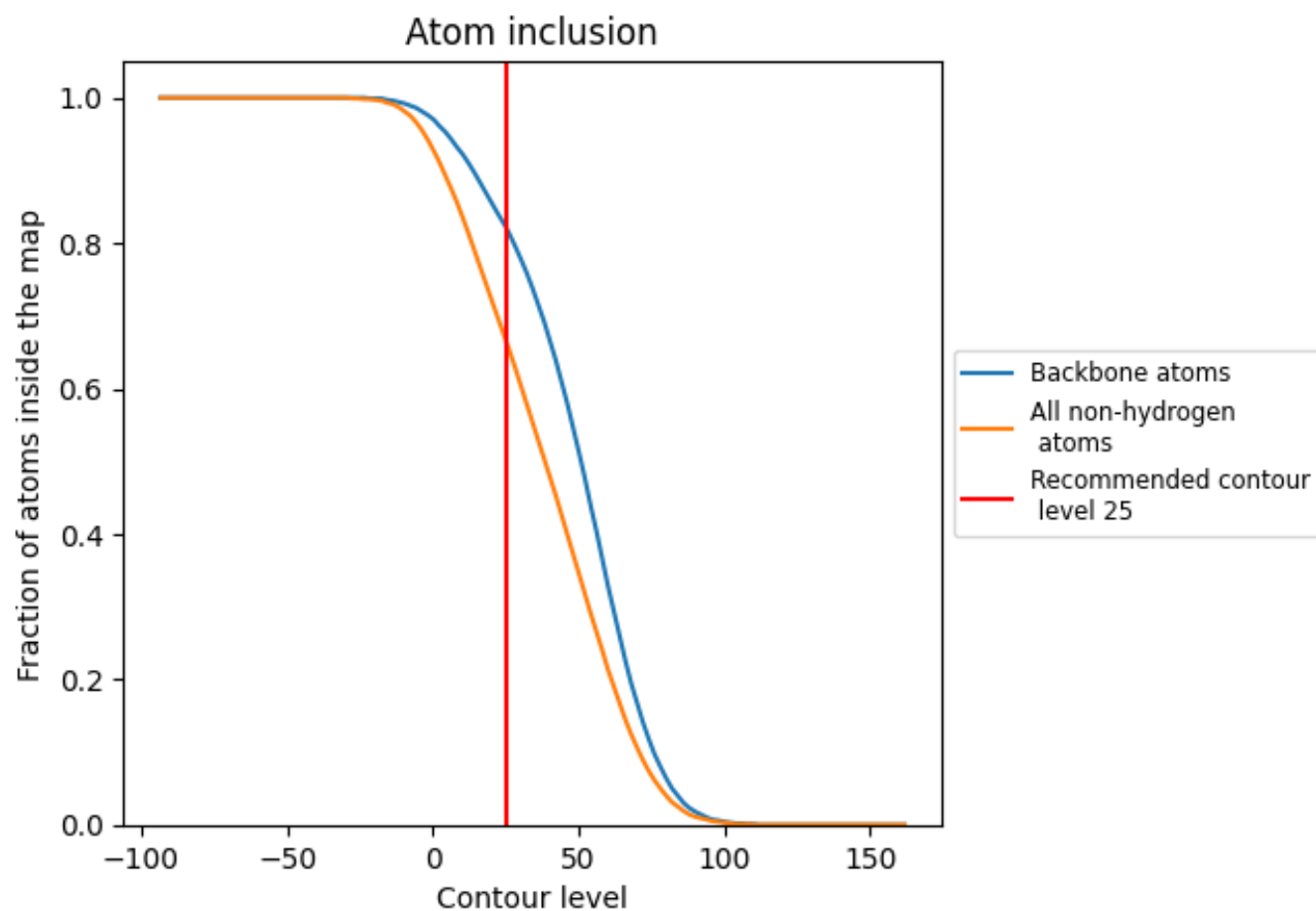
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (25).

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 67% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (25) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6680	0.3950
A	0.6480	0.3890
B	0.6690	0.4030
C	0.6750	0.3950
D	0.6640	0.3900
E	0.6760	0.4010
F	0.6760	0.3940
G	0.6680	0.3980
p	0.5560	0.3690
t	0.5440	0.3410

1.0

0.0

<0.0