



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 03:55 AM UTC

PDB ID : 8SQK / pdb_00008sqk
EMDB ID : EMD-40708
Title : SARS-CoV-2 replication-transcription complex bound to RNA-nsp9 and GDP-betaS, as a pre-catalytic deRNAylation/mRNA capping intermediate
Authors : Small, G.I.; Darst, S.A.; Campbell, E.A.
Deposited on : 2023-05-04
Resolution : 3.01 Å(reported)

This is a Full wwPDB EM 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/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:

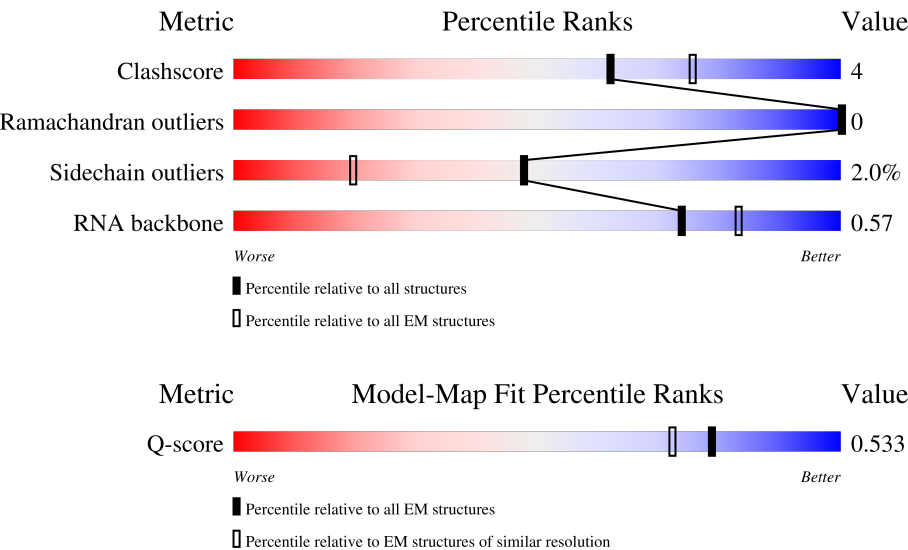
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 3.01 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	13882 (2.51 - 3.51)

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	929	91% 9%
2	B	198	14% 85% 8% 6%
2	D	198	9% 76% 17% 7%

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Mol	Chain	Length	Quality of chain
3	C	83	75% 12% • 12%
4	G	113	7% 85% 14% •
5	P	32	16% 84% 16%
6	T	35	17% 91% 9%
7	O	20	5% 95%

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 25574 atoms, of which 12229 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-directed RNA polymerase nsp12.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	929	Total	C	H	N	O	S	1	0
			14710	4781	7221	1258	1396	54		

- Molecule 2 is a protein called Non-structural protein 8.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	187	Total	C	H	N	O	S	0	0
			2825	886	1410	244	274	11		
2	D	185	Total	C	H	N	O	S	0	0
			2842	889	1428	242	272	11		

- Molecule 3 is a protein called Non-structural protein 7.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	C	73	Total	C	H	N	O	S	0	0
			1153	354	591	92	110	6		

- Molecule 4 is a protein called Non-structural protein 9.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	G	113	Total	C	H	N	O	S	0	0
			1651	528	822	143	153	5		

- Molecule 5 is a RNA chain called Primer RNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	P	32	Total	C	H	N	O	P	0	0
			1023	303	346	119	223	32		

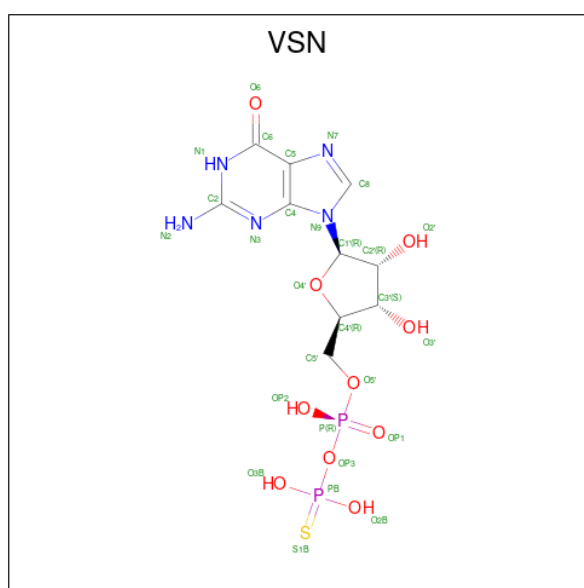
- Molecule 6 is a RNA chain called Template RNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
6	T	35	Total	C	H	N	O	P	0	0
			1123	333	376	131	248	35		

- Molecule 7 is a RNA chain called SARS-CoV-2 5' UTR.

Mol	Chain	Residues	Atoms						AltConf	Trace
7	O	1	Total	C	H	N	O	P	0	0
			33	10	11	5	6	1		

- Molecule 8 is 5'-O-[(R)-hydroxy(thiophosphonoxy)phosphoryl]guanosine (CCD ID: VSN) (formula: C₁₀H₁₅N₅O₁₀P₂S).



Mol	Chain	Residues	Atoms							AltConf
8	A	1	Total	C	H	N	O	P	S	0
			42	10	14	5	10	2	1	
8	A	1	Total	C	H	N	O	P	S	0
			38	10	10	5	10	2	1	

- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
9	A	2	Total	Zn	0
			2	2	

- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
10	A	1	Total 1	Mg 1	0

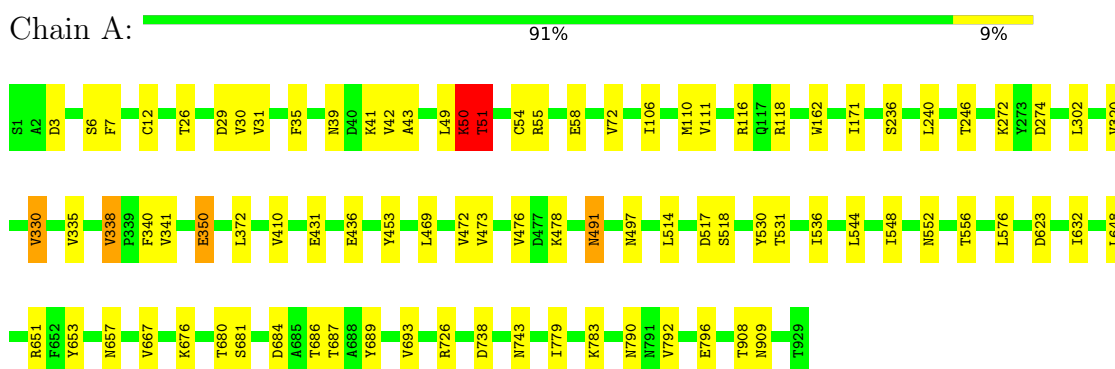
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		AltConf
11	A	110	Total 110	O 110	0
11	B	3	Total 3	O 3	0
11	C	5	Total 5	O 5	0
11	P	5	Total 5	O 5	0
11	T	7	Total 7	O 7	0
11	O	1	Total 1	O 1	0

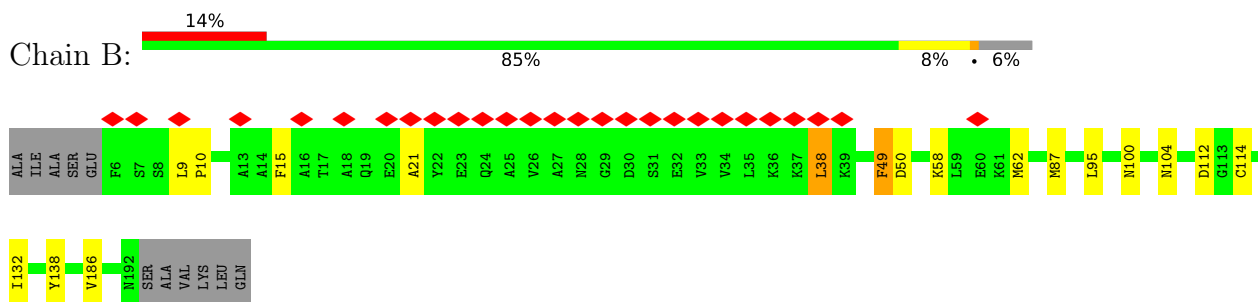
3 Residue-property plots

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.

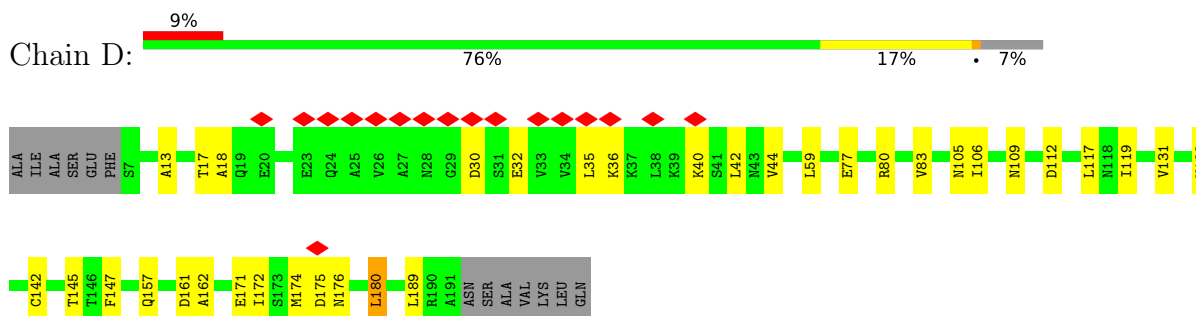
- Molecule 1: RNA-directed RNA polymerase nsp12



- Molecule 2: Non-structural protein 8

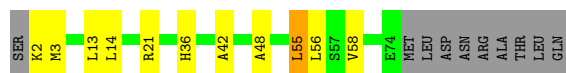


- Molecule 2: Non-structural protein 8



- Molecule 3: Non-structural protein 7

Chain C:  75% 12% • 12%



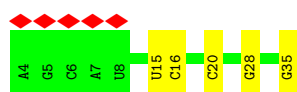
- Molecule 4: Non-structural protein 9

Chain G:  7% 85% 14% •

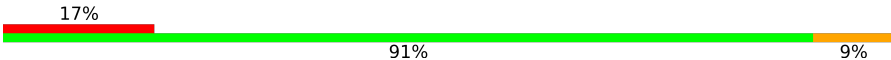


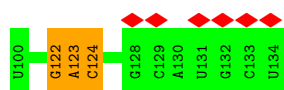
- Molecule 5: Primer RNA

Chain P:  16% 84% 16%



- Molecule 6: Template RNA

Chain T:  17% 91% 9%



- Molecule 7: SARS-CoV-2 5' UTR

Chain O:  5% 95%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	113552	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	51.51	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	43.190	Depositor
Minimum map value	-27.514	Depositor
Average map value	0.001	Depositor
Map value standard deviation	1.029	Depositor
Recommended contour level	2.83	Depositor
Map size (Å)	277.12, 277.12, 277.12	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.0825, 1.0825, 1.0825	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: VSN, ZN, MG

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	0.19	1/7681 (0.0%)	0.39	4/10423 (0.0%)
2	B	0.13	0/1434	0.28	0/1945
2	D	0.14	0/1433	0.33	0/1942
3	C	0.16	0/565	0.29	0/761
4	G	0.15	0/845	0.34	0/1154
5	P	0.12	0/755	0.21	0/1173
6	T	0.13	0/834	0.23	0/1298
7	O	0.12	0/24	0.32	0/35
All	All	0.17	1/13571 (0.0%)	0.35	4/18731 (0.0%)

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	50	LYS	N-CA	-5.89	1.39	1.45

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	51	THR	CA-C-N	11.80	136.55	120.38
1	A	51	THR	C-N-CA	11.80	136.55	120.38
1	A	51	THR	CB-CA-C	11.64	133.59	110.42
1	A	50	LYS	N-CA-CB	-5.47	101.98	110.46

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	51	THR	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	7489	7221	7232	53	0
2	B	1415	1410	1408	10	0
2	D	1414	1428	1428	26	0
3	C	562	591	591	7	0
4	G	829	822	802	10	0
5	P	677	346	346	1	0
6	T	747	376	376	3	0
7	O	22	11	12	1	0
8	A	56	24	0	7	0
9	A	2	0	0	0	0
10	A	1	0	0	0	0
11	A	110	0	0	8	0
11	B	3	0	0	0	0
11	C	5	0	0	0	0
11	O	1	0	0	1	0
11	P	5	0	0	0	0
11	T	7	0	0	0	0
All	All	13345	12229	12195	110	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 4.

All (110) 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)
8:A:1002:VSN:PB	8:A:1002:VSN:S1B	1.51	1.50
6:T:123:A:O2'	6:T:124:C:OP1	1.87	0.91
1:A:436:GLU:N	1:A:436:GLU:OE2	2.08	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:796:GLU:OE2	11:A:1101:HOH:O	1.94	0.85
1:A:497:ASN:OD1	11:A:1102:HOH:O	1.98	0.81
2:B:100:ASN:O	2:B:104:ASN:ND2	2.14	0.79
2:D:40:LYS:NZ	6:T:122:G:OP1	2.18	0.74
8:A:1002:VSN:S1B	8:A:1002:VSN:O3B	2.45	0.74
1:A:431:GLU:OE2	11:A:1103:HOH:O	2.06	0.73
7:O:1:A:N1	11:O:101:HOH:O	2.21	0.73
2:D:161:ASP:OD1	2:D:162:ALA:N	2.22	0.72
4:G:15:ALA:HB1	4:G:24:THR:O	1.90	0.71
1:A:576:LEU:HD11	1:A:686:THR:HG22	1.72	0.69
8:A:1002:VSN:S1B	8:A:1002:VSN:O2B	2.47	0.69
1:A:272:LYS:NZ	1:A:274:ASP:O	2.26	0.69
2:D:77:GLU:OE2	2:D:80:ARG:NH2	2.26	0.69
2:D:105:ASN:OD1	2:D:109:ASN:ND2	2.29	0.66
1:A:684:ASP:OD1	1:A:687:THR:OG1	2.13	0.66
4:G:15:ALA:HB2	4:G:26:ASP:N	2.11	0.65
1:A:29:ASP:O	1:A:51:THR:O	2.14	0.65
1:A:453:TYR:CE1	1:A:667:VAL:HG21	2.32	0.64
1:A:469:LEU:O	1:A:473:VAL:HG23	1.98	0.63
8:A:1002:VSN:S1B	8:A:1002:VSN:OP3	2.54	0.62
1:A:651:ARG:NH1	11:A:1107:HOH:O	2.31	0.61
1:A:30:VAL:HG12	1:A:51:THR:HA	1.82	0.60
4:G:45:LEU:HD22	4:G:86:LYS:HE3	1.82	0.59
2:D:142:CYS:HG	2:D:147:PHE:HD1	1.51	0.58
4:G:70:GLU:OE1	4:G:99:ARG:NH2	2.37	0.58
6:T:123:A:HO2'	6:T:124:C:P	2.22	0.58
8:A:1001:VSN:O4'	5:P:35:G:O2'	2.21	0.58
1:A:330:VAL:HG13	1:A:341:VAL:HG13	1.87	0.56
1:A:478:LYS:NZ	1:A:743:ASN:OD1	2.24	0.56
2:D:138:TYR:CE2	2:D:172:ILE:HG21	2.41	0.56
1:A:29:ASP:OD1	1:A:30:VAL:N	2.39	0.56
2:D:145:THR:HG22	2:D:145:THR:O	2.07	0.55
1:A:7:PHE:HZ	1:A:72:VAL:HG11	1.71	0.55
1:A:472:VAL:O	1:A:476:VAL:HG23	2.07	0.54
1:A:110:MET:HA	1:A:110:MET:HE2	1.90	0.54
1:A:58:GLU:OE1	1:A:118:ARG:NH1	2.41	0.53
4:G:25:ASP:OD1	4:G:26:ASP:N	2.42	0.53
4:G:9:LEU:HD23	4:G:31:TYR:HB3	1.90	0.53
2:D:30:ASP:C	2:D:35:LEU:HD23	2.34	0.53
3:C:3:MET:HE1	3:C:48:ALA:HB2	1.92	0.51
1:A:552:ASN:OD1	1:A:552:ASN:O	2.27	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:15:ALA:HB2	4:G:26:ASP:H	1.76	0.51
1:A:726:ARG:NH2	1:A:738:ASP:OD2	2.44	0.50
11:A:1105:HOH:O	3:C:2:LYS:N	2.44	0.50
1:A:491:ASN:N	1:A:491:ASN:OD1	2.44	0.50
1:A:372:LEU:HD12	2:B:87:MET:HE1	1.93	0.50
3:C:58:VAL:HG22	2:D:119:ILE:HG12	1.94	0.49
2:D:13:ALA:O	2:D:17:THR:HG23	2.13	0.49
1:A:3:ASP:O	1:A:6:SER:OG	2.23	0.49
1:A:42:VAL:HG22	1:A:43:ALA:N	2.28	0.48
2:B:58:LYS:O	2:B:62:MET:HG2	2.13	0.48
1:A:116:ARG:NH2	8:A:1002:VSN:OP2	2.46	0.48
2:D:171:GLU:OE1	2:D:180:LEU:HD21	2.14	0.48
2:B:132:ILE:HG21	2:B:138:TYR:HB2	1.96	0.47
1:A:517:ASP:C	1:A:517:ASP:OD1	2.57	0.47
1:A:544:LEU:HD12	1:A:556:THR:HG22	1.95	0.47
1:A:676:LYS:NZ	1:A:681:SER:OG	2.46	0.47
3:C:56:LEU:HD23	2:D:106:ILE:HD11	1.97	0.47
2:D:35:LEU:HD12	2:D:36:LYS:N	2.29	0.47
2:D:32:GLU:N	2:D:32:GLU:OE1	2.47	0.46
1:A:779:ILE:HG22	11:A:1101:HOH:O	2.16	0.46
2:D:175:ASP:OD1	2:D:176:ASN:N	2.48	0.46
3:C:14:LEU:HD22	3:C:36:HIS:CG	2.50	0.46
2:D:157:GLN:HE21	2:D:189:LEU:HD22	1.81	0.45
1:A:514:LEU:O	1:A:518:SER:OG	2.30	0.45
1:A:530:TYR:CD2	1:A:536:ILE:HD11	2.52	0.45
1:A:517:ASP:OD1	1:A:518:SER:N	2.50	0.45
2:D:117:LEU:HD21	2:D:131:VAL:HG21	1.97	0.45
1:A:302:LEU:O	1:A:648:LEU:HD21	2.17	0.45
1:A:653:TYR:O	1:A:657:ASN:ND2	2.43	0.45
1:A:162:TRP:O	1:A:171:ILE:HD13	2.17	0.44
1:A:335:VAL:O	1:A:338:VAL:HG12	2.17	0.44
2:D:112:ASP:OD1	2:D:112:ASP:C	2.60	0.44
1:A:531:THR:HG22	1:A:536:ILE:HD12	1.99	0.44
1:A:623:ASP:N	1:A:623:ASP:OD1	2.49	0.44
1:A:689:TYR:O	1:A:693:VAL:HG23	2.18	0.43
4:G:80:PRO:O	4:G:81:LYS:C	2.61	0.43
1:A:31:VAL:O	1:A:49:LEU:HD12	2.19	0.43
1:A:39:ASN:OD1	1:A:41:LYS:N	2.48	0.43
2:D:174:MET:O	2:D:174:MET:HE3	2.19	0.43
1:A:783:LYS:HB2	11:A:1101:HOH:O	2.17	0.43
1:A:35:PHE:CZ	1:A:50:LYS:HG3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:9:LEU:HB2	2:B:10:PRO:HD3	1.99	0.43
2:B:112:ASP:OD1	2:B:112:ASP:N	2.49	0.43
1:A:106:ILE:HD12	1:A:111:VAL:HG21	2.00	0.43
4:G:5:SER:HB3	4:G:101:MET:HG2	2.01	0.42
2:B:21:ALA:HB1	2:B:38:LEU:HB2	2.01	0.42
2:D:109:ASN:O	2:D:112:ASP:OD1	2.38	0.42
2:B:49:PHE:O	2:B:50:ASP:C	2.63	0.42
2:D:138:TYR:HE2	2:D:172:ILE:HG21	1.85	0.42
3:C:13:LEU:HG	3:C:55:LEU:HD22	2.01	0.42
2:B:95:LEU:C	2:B:95:LEU:HD23	2.45	0.41
2:D:138:TYR:CD1	2:D:142:CYS:HB2	2.55	0.41
2:D:138:TYR:CE1	2:D:142:CYS:HB2	2.55	0.41
1:A:320:VAL:O	1:A:320:VAL:HG12	2.20	0.41
1:A:576:LEU:HD12	11:A:1187:HOH:O	2.19	0.41
1:A:55:ARG:HD3	8:A:1002:VSN:N7	2.36	0.41
2:D:83:VAL:HG22	2:D:83:VAL:O	2.20	0.41
1:A:340:PHE:CE1	2:B:95:LEU:HD12	2.56	0.41
3:C:3:MET:HE2	3:C:42:ALA:O	2.21	0.41
1:A:350:GLU:HG3	1:A:632:ILE:HD11	2.03	0.41
2:D:59:LEU:C	2:D:59:LEU:HD23	2.46	0.40
4:G:68:GLU:O	4:G:69:LEU:HD12	2.21	0.40
2:D:18:ALA:HB3	2:D:42:LEU:HD13	2.03	0.40
1:A:236:SER:O	1:A:240:LEU:HD12	2.21	0.40
1:A:410:VAL:HG23	1:A:544:LEU:HD23	2.04	0.40
1:A:530:TYR:CE2	1:A:536:ILE:HD11	2.57	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 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	928/929 (100%)	892 (96%)	36 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	185/198 (93%)	177 (96%)	8 (4%)	0	100	100
2	D	183/198 (92%)	174 (95%)	9 (5%)	0	100	100
3	C	71/83 (86%)	70 (99%)	1 (1%)	0	100	100
4	G	111/113 (98%)	104 (94%)	7 (6%)	0	100	100
All	All	1478/1521 (97%)	1417 (96%)	61 (4%)	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 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	821/820 (100%)	806 (98%)	15 (2%)	51	76
2	B	149/167 (89%)	144 (97%)	5 (3%)	32	64
2	D	151/167 (90%)	149 (99%)	2 (1%)	61	80
3	C	68/77 (88%)	66 (97%)	2 (3%)	37	68
4	G	82/94 (87%)	80 (98%)	2 (2%)	43	72
All	All	1271/1325 (96%)	1245 (98%)	26 (2%)	48	75

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	CYS
1	A	26	THR
1	A	50	LYS
1	A	54	CYS
1	A	246	THR
1	A	330	VAL
1	A	338	VAL
1	A	350	GLU
1	A	491	ASN
1	A	548	ILE

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Mol	Chain	Res	Type
1	A	680	THR
1	A	790	ASN
1	A	792	VAL
1	A	908	THR
1	A	909	ASN
2	B	15	PHE
2	B	38	LEU
2	B	49	PHE
2	B	114	CYS
2	B	186	VAL
3	C	21	ARG
3	C	55	LEU
2	D	44	VAL
2	D	180	LEU
4	G	45	LEU
4	G	73	CYS

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

Mol	Chain	Res	Type
1	A	64	ASN
1	A	79	ASN
1	A	81	GLN
1	A	209	ASN
1	A	297	ASN
1	A	355	HIS
1	A	599	HIS
3	C	63	GLN
3	C	69	ASN
4	G	1	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
5	P	31/32 (96%)	4 (12%)	0
6	T	34/35 (97%)	3 (8%)	1 (2%)
7	O	0/20	-	-
All	All	65/87 (74%)	7 (10%)	1 (1%)

All (7) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	P	15	U
5	P	16	C
5	P	20	C
5	P	28	G
6	T	122	G
6	T	123	A
6	T	124	C

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
6	T	123	A

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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	$\# Z > 2$
8	VSN	A	1001	-	29,30,30	3.26	14 (48%)	42,47,47	1.82	12 (28%)
8	VSN	A	1002	10	29,30,30	3.57	4 (13%)	42,47,47	1.95	8 (19%)

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
8	VSN	A	1001	-	-	4/15/32/32	0/3/3/3
8	VSN	A	1002	10	-	7/15/32/32	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	1002	VSN	PB-S1B	-18.16	1.51	1.90
8	A	1001	VSN	O4'-C4'	7.45	1.61	1.45
8	A	1001	VSN	C2'-C3'	-7.17	1.33	1.53
8	A	1001	VSN	P-OP3	6.04	1.66	1.59
8	A	1001	VSN	C2-N2	5.40	1.46	1.34
8	A	1001	VSN	O4'-C1'	-5.30	1.29	1.42
8	A	1001	VSN	C2'-C1'	4.57	1.67	1.53
8	A	1001	VSN	C5'-C4'	-4.40	1.38	1.51
8	A	1001	VSN	PB-O2B	3.60	1.66	1.54
8	A	1001	VSN	PB-O3B	3.19	1.65	1.54
8	A	1002	VSN	C5-C4	3.05	1.47	1.38
8	A	1001	VSN	C6-N1	-2.99	1.33	1.38
8	A	1002	VSN	C6-N1	-2.59	1.34	1.38
8	A	1001	VSN	O3'-C3'	2.56	1.49	1.43
8	A	1001	VSN	C5-N7	-2.50	1.34	1.39
8	A	1001	VSN	P-O5'	2.50	1.69	1.59
8	A	1001	VSN	C8-N9	-2.35	1.32	1.37
8	A	1002	VSN	C5-N7	-2.32	1.34	1.39

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1002	VSN	C5-C4-N3	-6.47	118.10	128.39
8	A	1001	VSN	C5-C4-N3	-5.54	119.56	128.39
8	A	1002	VSN	C2-N3-C4	5.20	121.26	112.30
8	A	1002	VSN	N9-C4-N3	4.81	135.56	125.95
8	A	1001	VSN	C2-N3-C4	4.62	120.25	112.30
8	A	1001	VSN	N9-C4-N3	3.90	133.75	125.95
8	A	1002	VSN	P-OP3-PB	-3.43	120.60	133.17
8	A	1001	VSN	C3'-C2'-C1'	3.25	107.61	101.46
8	A	1002	VSN	C6-C5-N7	3.07	135.88	130.29
8	A	1002	VSN	C4-C5-N7	-2.64	106.49	110.67
8	A	1001	VSN	C2-N1-C6	-2.55	120.48	125.11
8	A	1001	VSN	C5-C6-N1	2.50	119.62	113.25
8	A	1001	VSN	P-OP3-PB	-2.44	124.25	133.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1001	VSN	O6-C6-C5	-2.42	120.14	126.53
8	A	1001	VSN	C6-C5-N7	2.31	134.50	130.29
8	A	1002	VSN	C3'-C2'-C1'	2.12	105.48	101.46
8	A	1002	VSN	O6-C6-C5	-2.10	120.99	126.53
8	A	1001	VSN	C8-N7-C5	2.08	107.96	104.26
8	A	1001	VSN	C4-C5-N7	-2.06	107.40	110.67
8	A	1001	VSN	N9-C8-N7	-2.01	109.67	113.40

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	1001	VSN	C5'-O5'-P-OP2
8	A	1001	VSN	C5'-O5'-P-OP3
8	A	1001	VSN	C5'-O5'-P-OP1
8	A	1002	VSN	C5'-O5'-P-OP2
8	A	1002	VSN	C5'-O5'-P-OP3
8	A	1002	VSN	C3'-C4'-C5'-O5'
8	A	1002	VSN	O4'-C4'-C5'-O5'
8	A	1001	VSN	PB-OP3-P-O5'
8	A	1002	VSN	PB-OP3-P-O5'
8	A	1002	VSN	C2'-C1'-N9-C8
8	A	1002	VSN	C2'-C1'-N9-C4

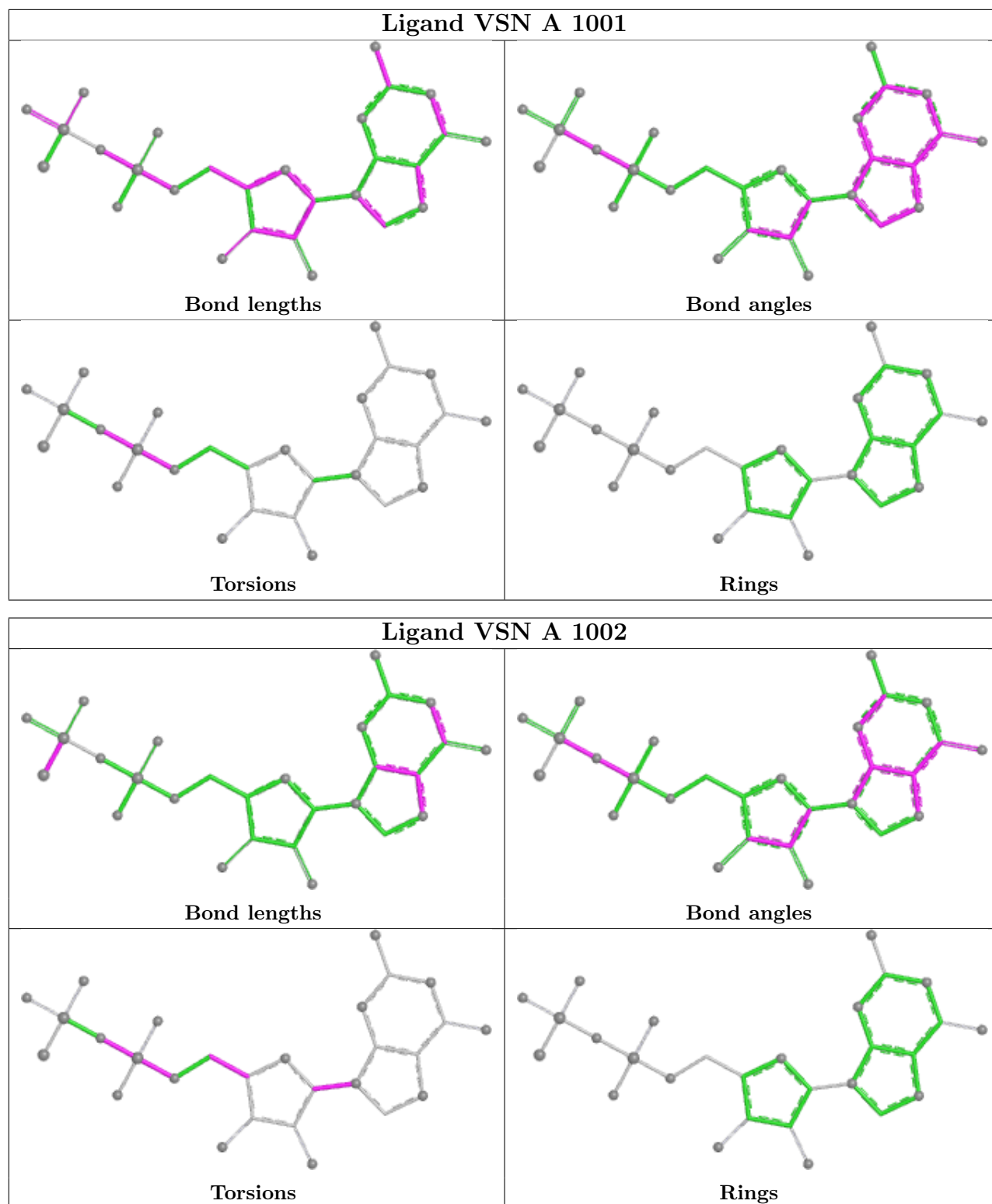
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	1001	VSN	1	0
8	A	1002	VSN	6	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.

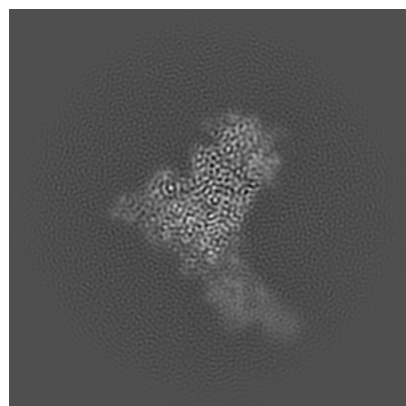
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40708. These allow visual inspection of the internal detail of the map and identification of artifacts.

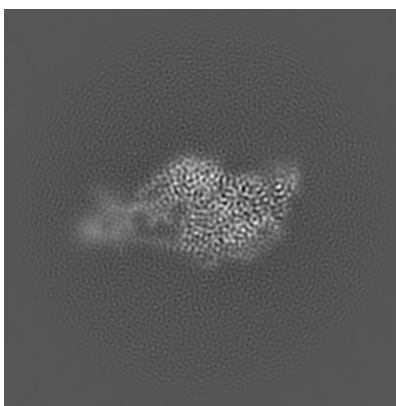
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

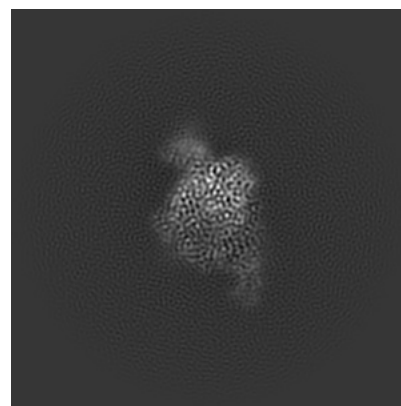
6.1.1 Primary map



X

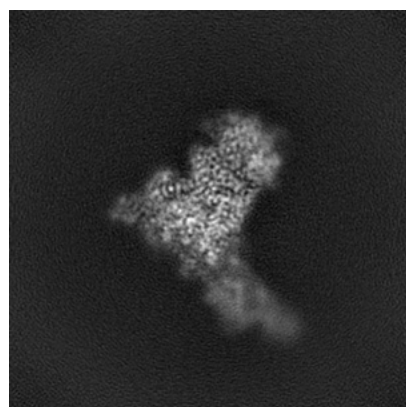


Y

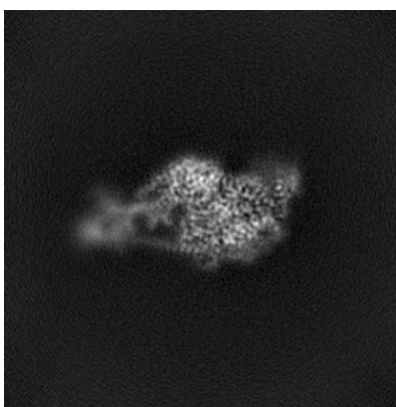


Z

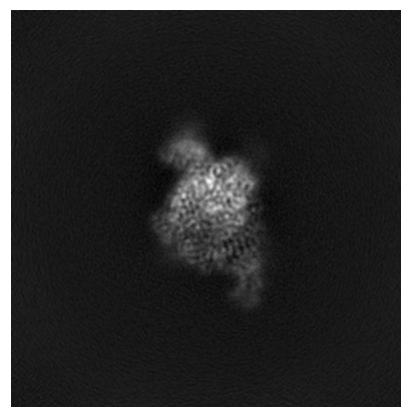
6.1.2 Raw map



X



Y

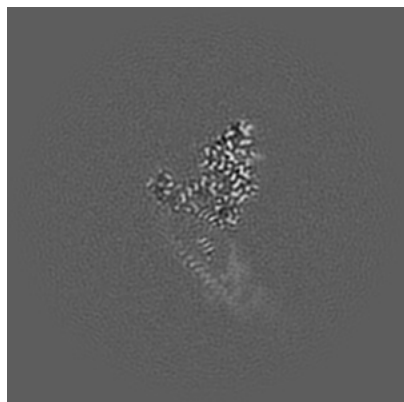


Z

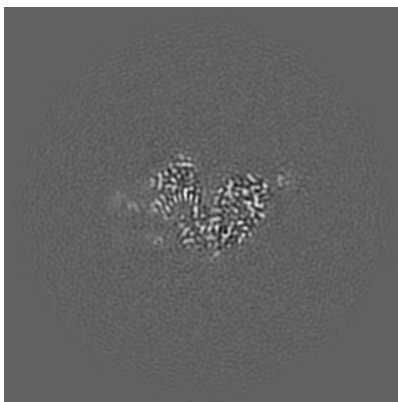
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

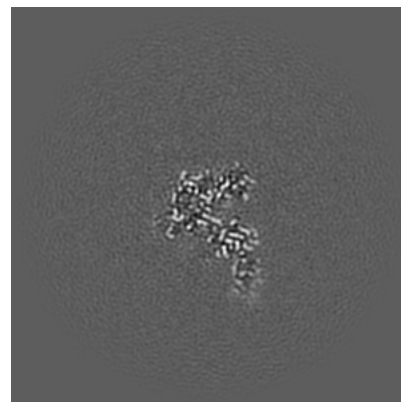
6.2.1 Primary map



X Index: 128

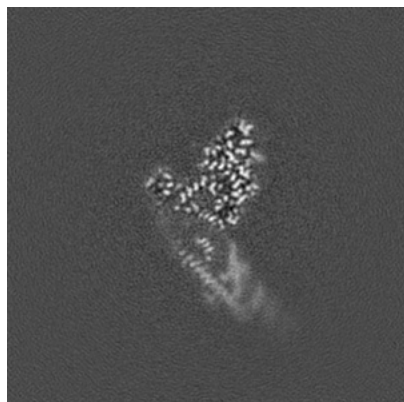


Y Index: 128

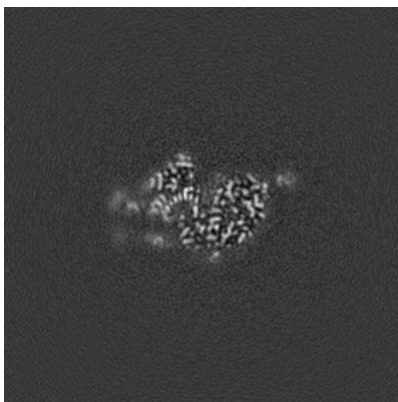


Z Index: 128

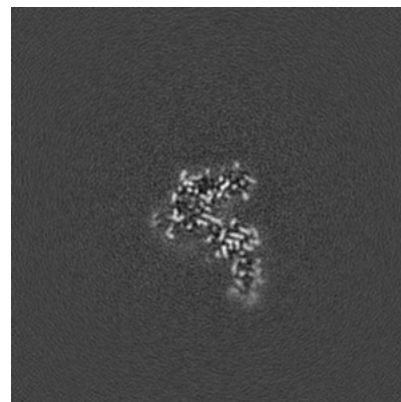
6.2.2 Raw map



X Index: 128



Y Index: 128

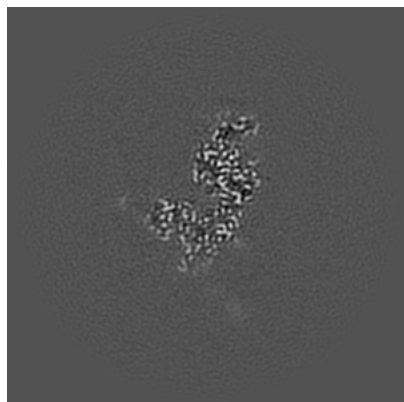


Z Index: 128

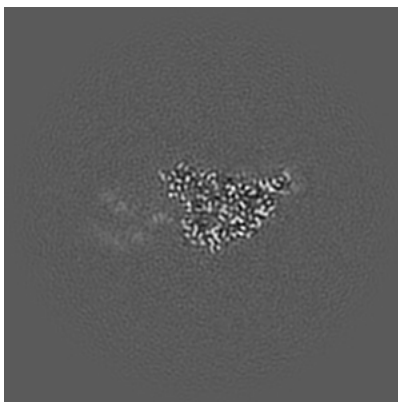
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

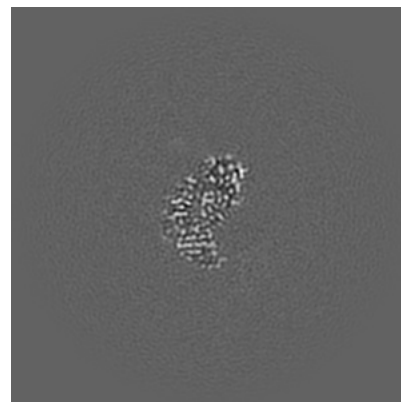
6.3.1 Primary map



X Index: 140

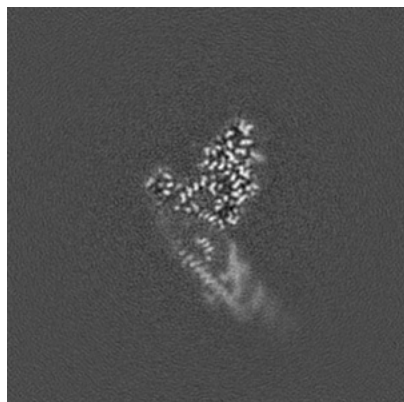


Y Index: 136

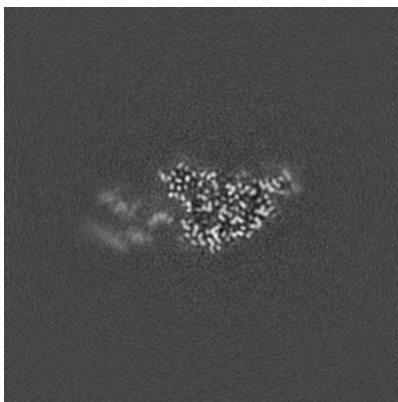


Z Index: 141

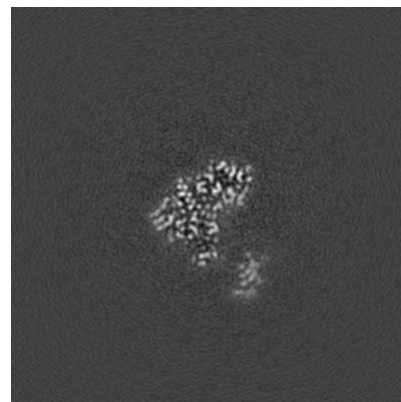
6.3.2 Raw map



X Index: 128



Y Index: 136

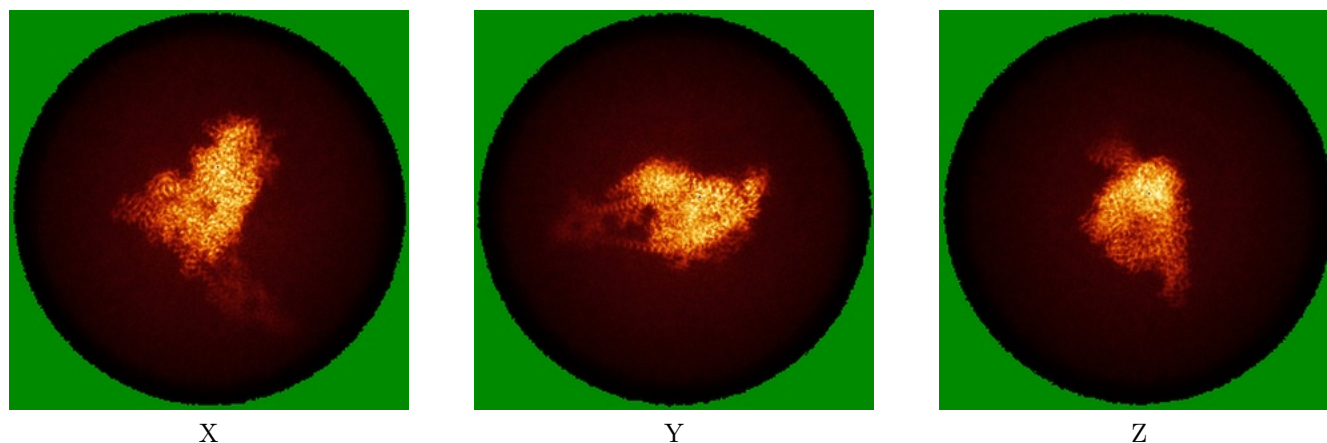


Z Index: 135

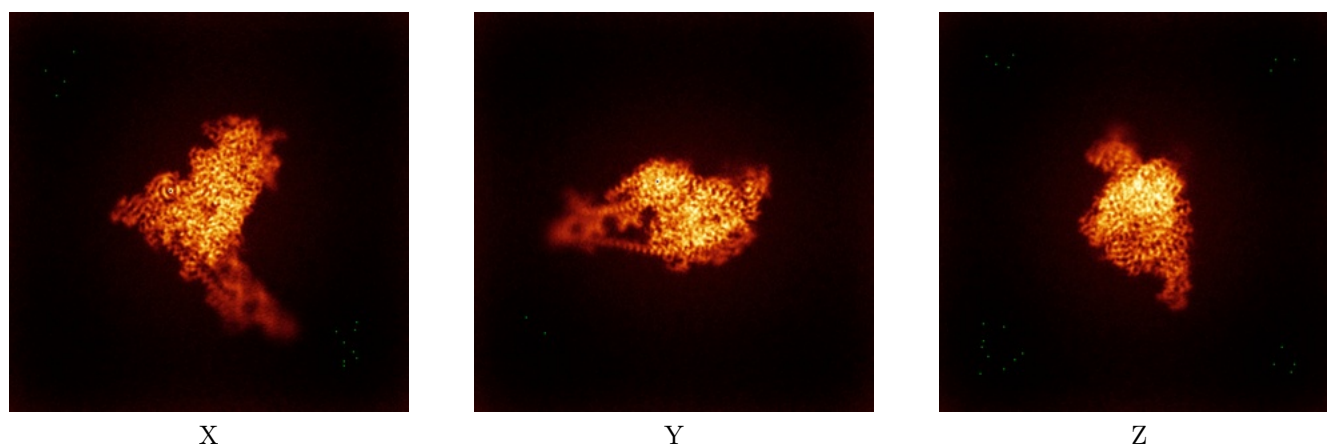
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

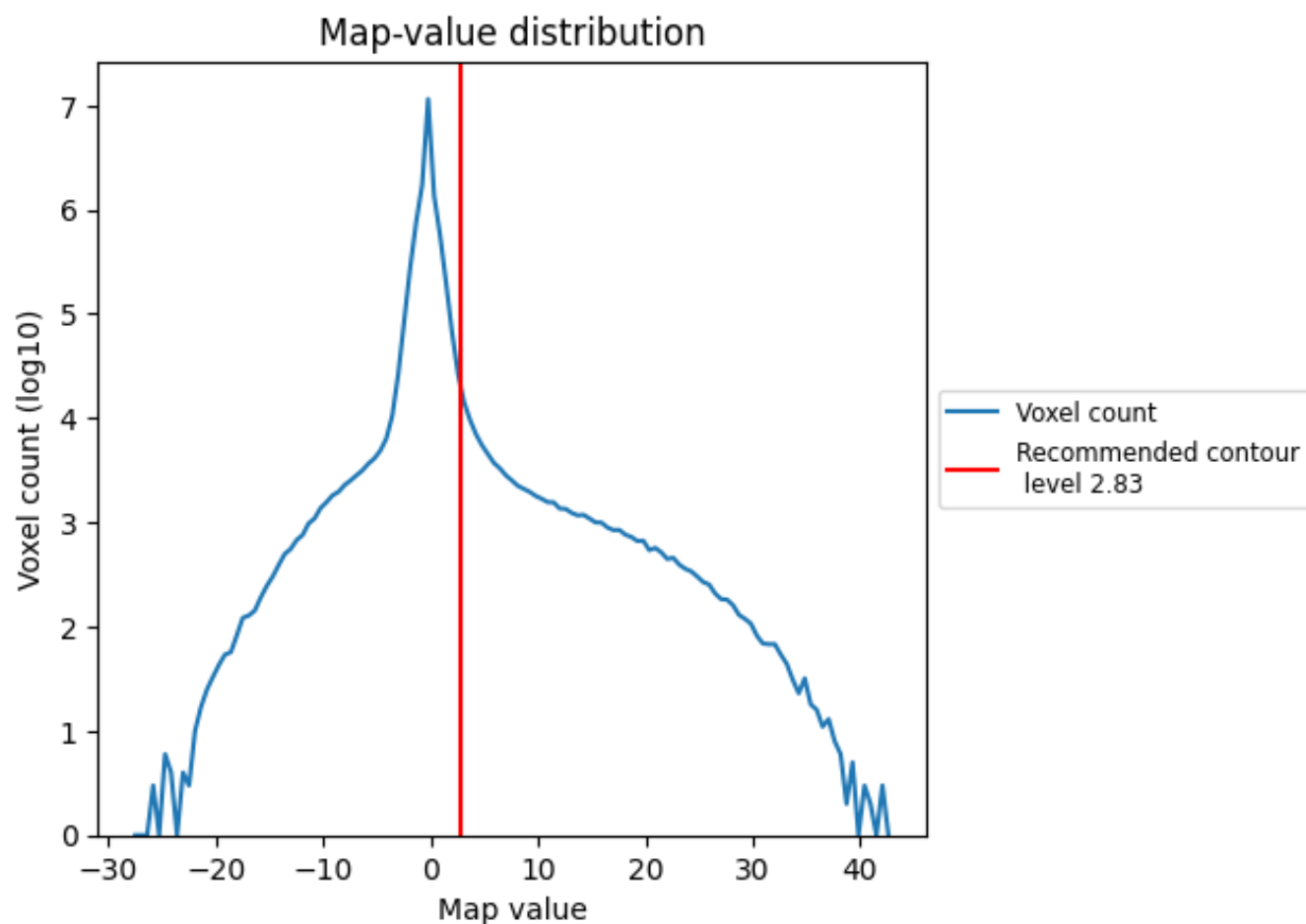
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

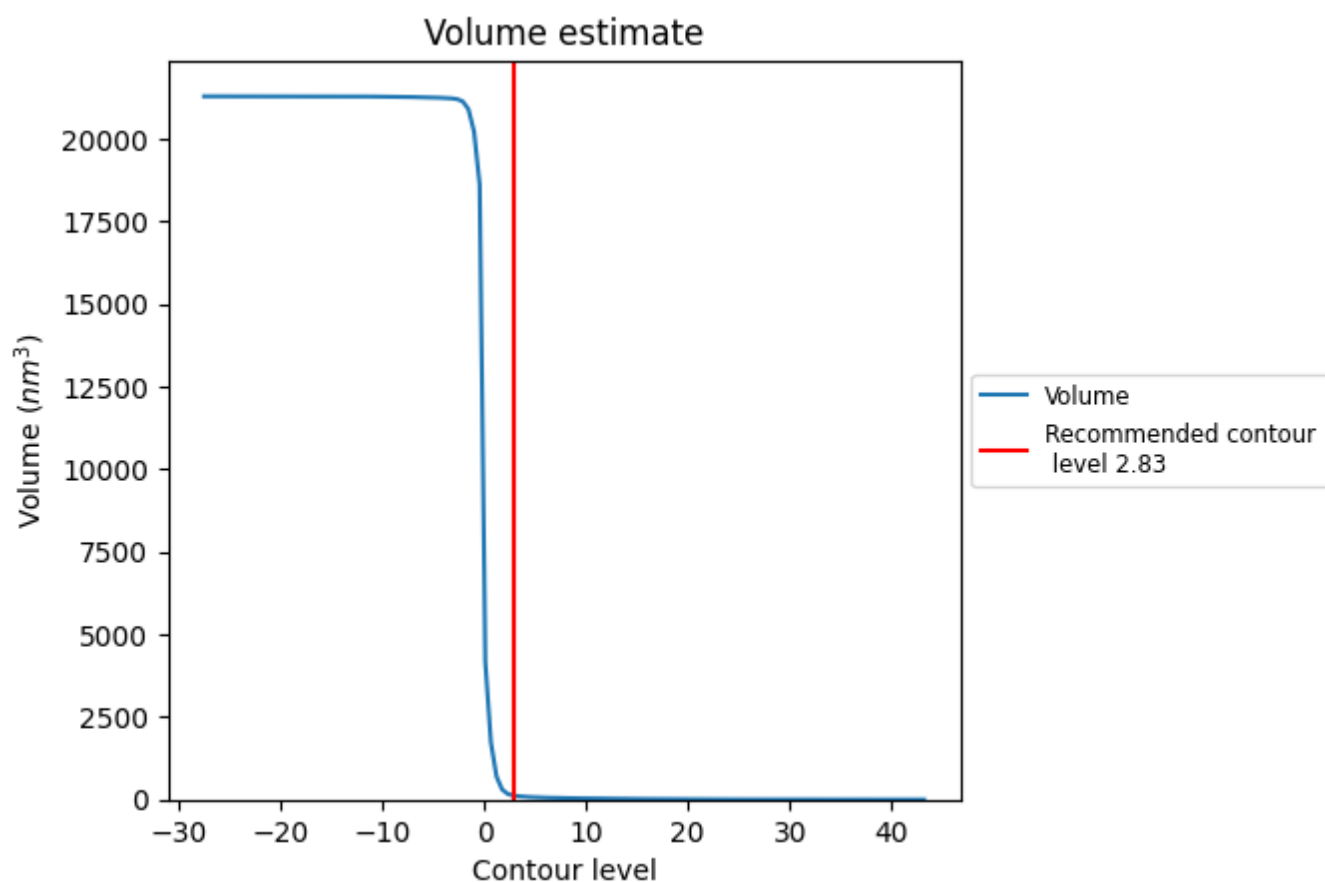
This section contains the results of statistical analysis of the map.

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.

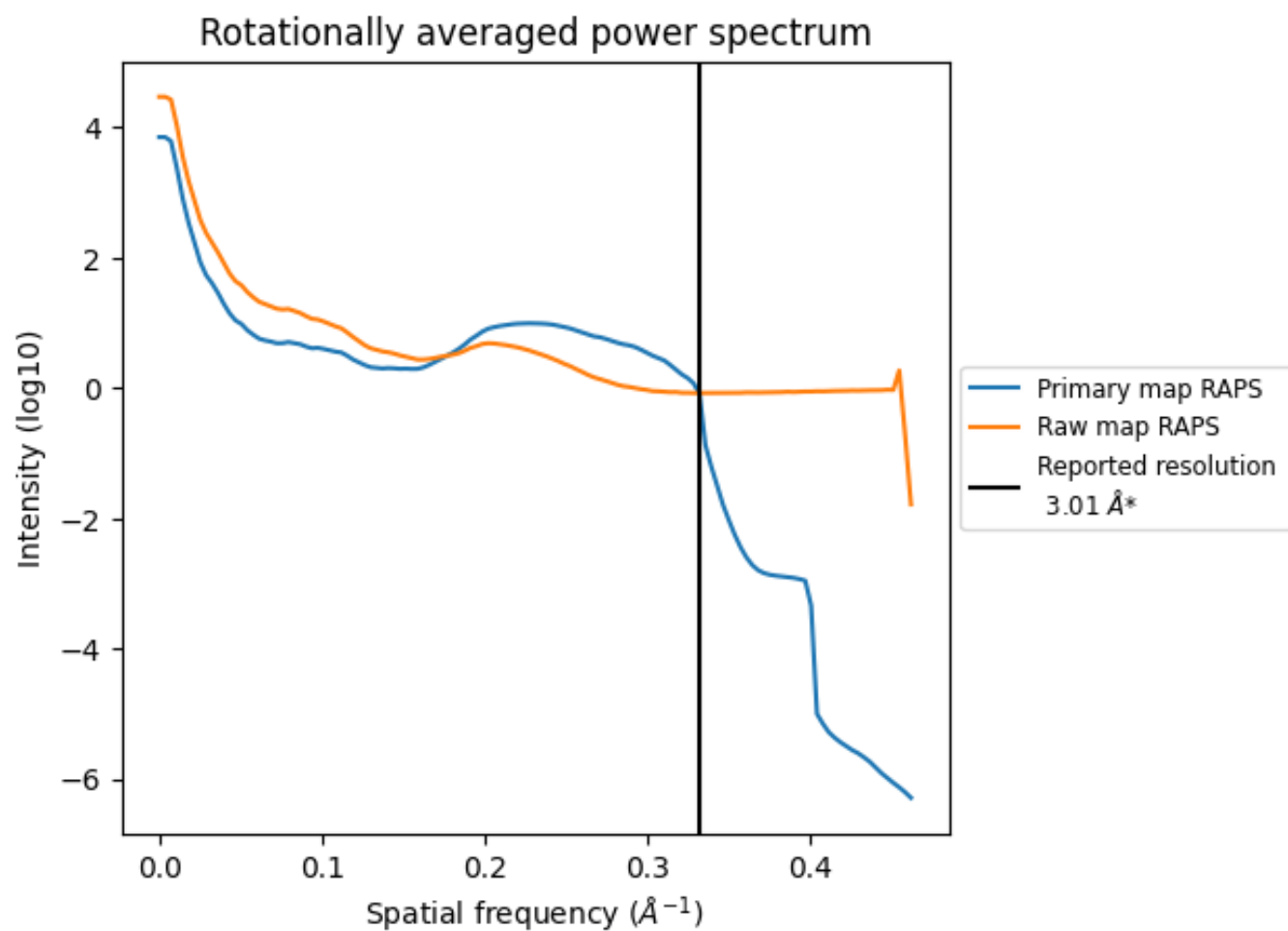
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 125 nm^3 ; this corresponds to an approximate mass of 113 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 ⓘ

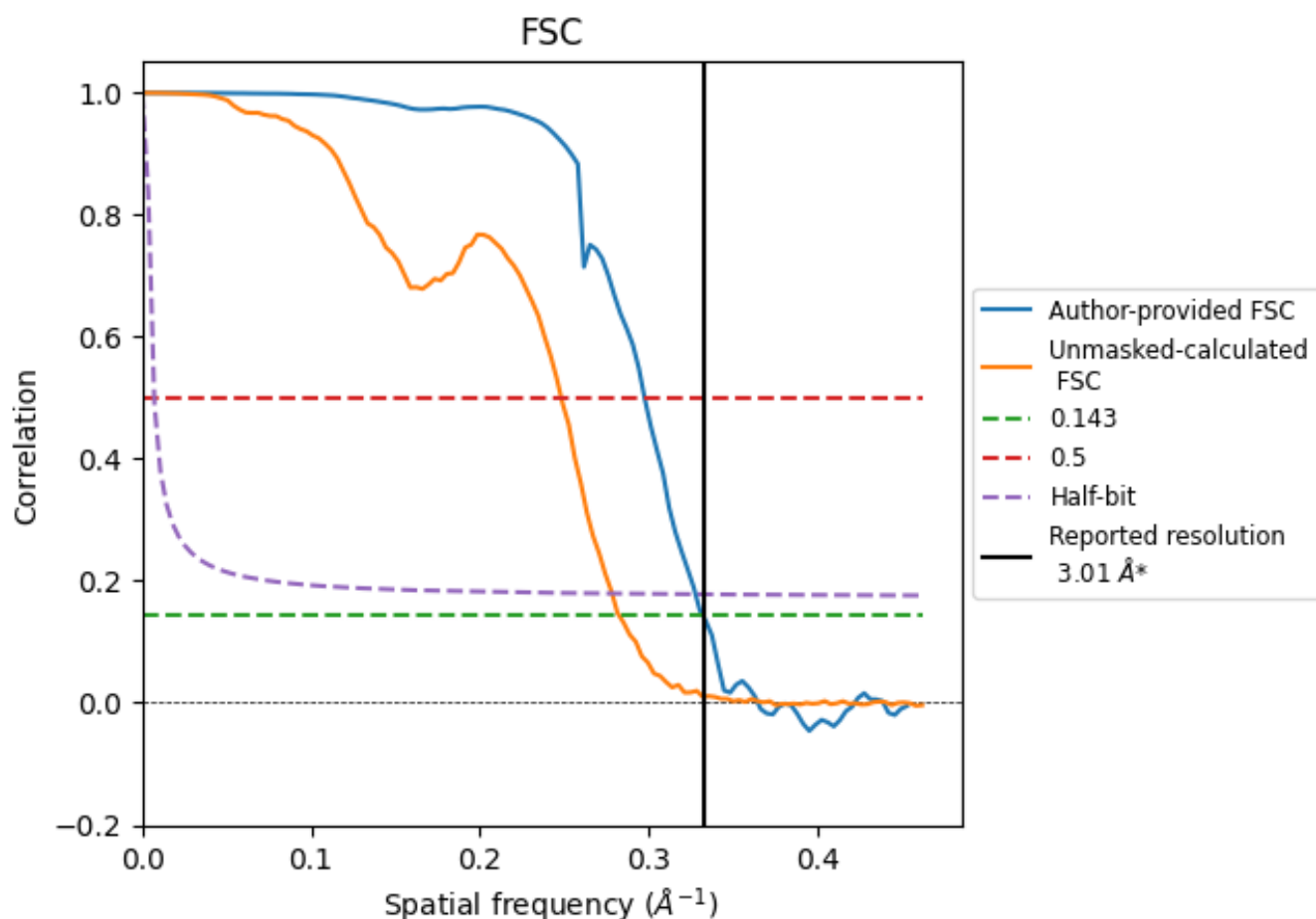


*Reported resolution corresponds to spatial frequency of 0.332 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.332 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

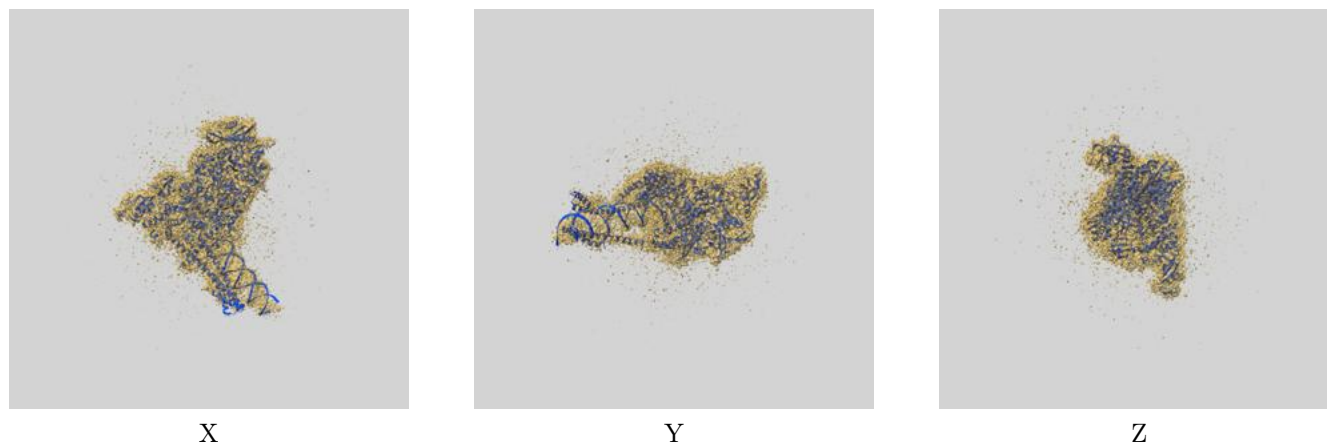
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.01	-	-
Author-provided FSC curve	3.01	3.36	3.05
Unmasked-calculated*	3.54	4.03	3.59

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.54 differs from the reported value 3.01 by more than 10 %

9 Map-model fit [i](#)

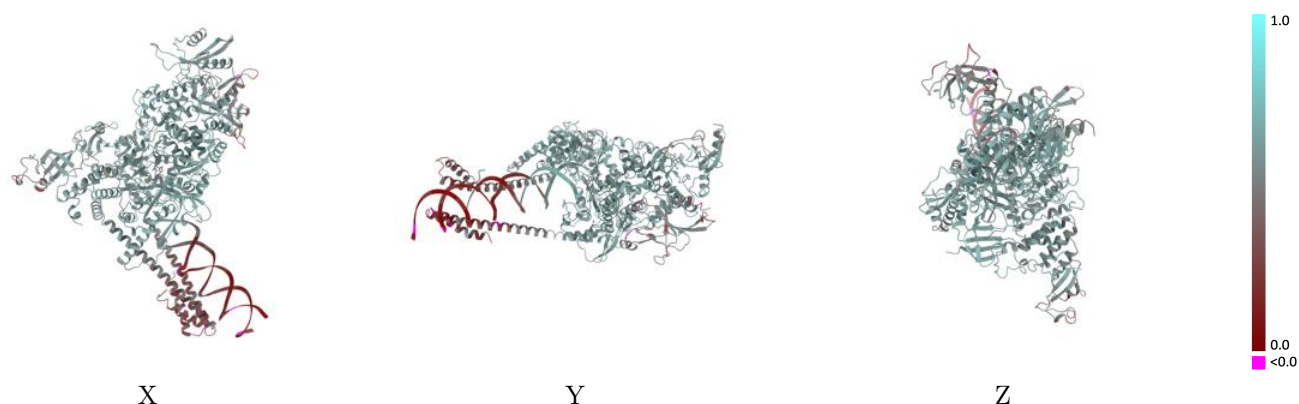
This section contains information regarding the fit between EMDB map EMD-40708 and PDB model 8SQK. Per-residue inclusion information can be found in section [3](#) on page [7](#).

9.1 Map-model overlay [i](#)



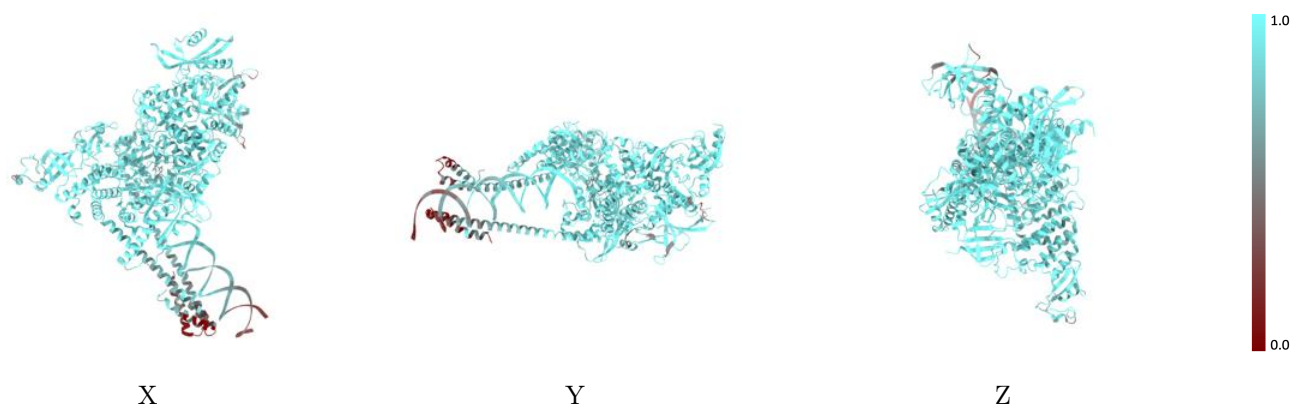
The images above show the 3D surface view of the map at the recommended contour level 2.83 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



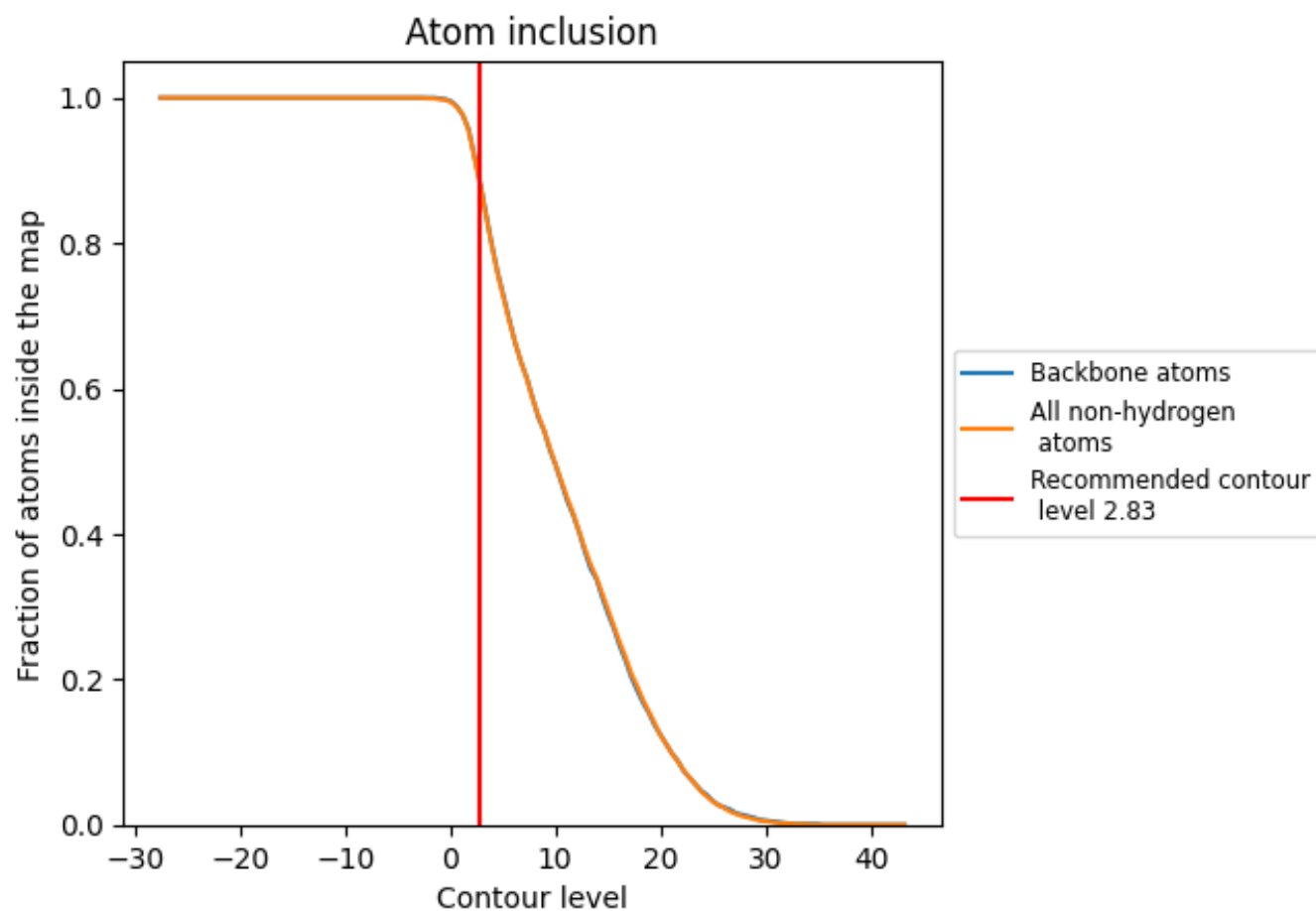
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (2.83).

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 88% 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 (2.83) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8830	0.5330
A	0.9590	0.5980
B	0.7850	0.4850
C	0.9280	0.5800
D	0.7800	0.4730
G	0.8510	0.4750
O	0.8640	0.4820
P	0.7390	0.2870
T	0.7480	0.3340

1.0

0.0

<0.0