



Full wwPDB EM Validation Report ⓘ

Mar 28, 2026 – 07:23 PM UTC

PDB ID : 8VJR / pdb_00008vjr
EMDB ID : EMD-43292
Title : Cryo-EM structure of Tulane virus 9-6-17 variant capsid protein VP1 9-14-18, DTT-treated
Authors : Sun, C.; Jiang, W.
Deposited on : 2024-01-07
Resolution : 2.63 Å(reported)
Based on initial model : 8VG6

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:

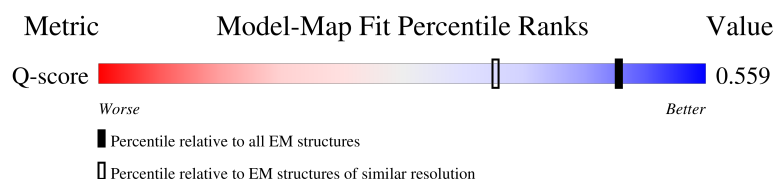
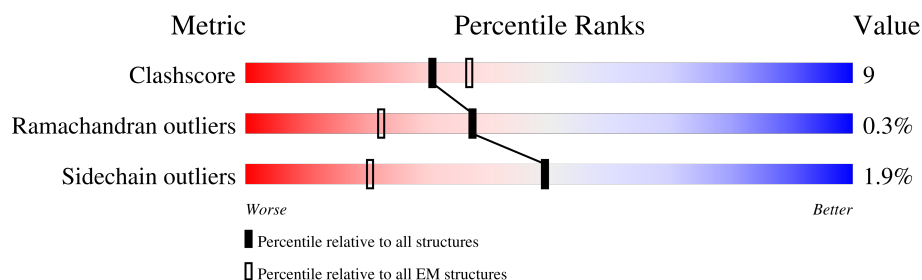
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 2.63 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	-
Q-score	-	25397	8888 (2.13 - 3.13)

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	534	11% 74% 19% • 5%
1	B	534	10% 81% 13% • 5%
1	C	534	57% 81% 16% ••

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 23300 atoms, of which 11495 are hydrogens and 0 are deuteriums.

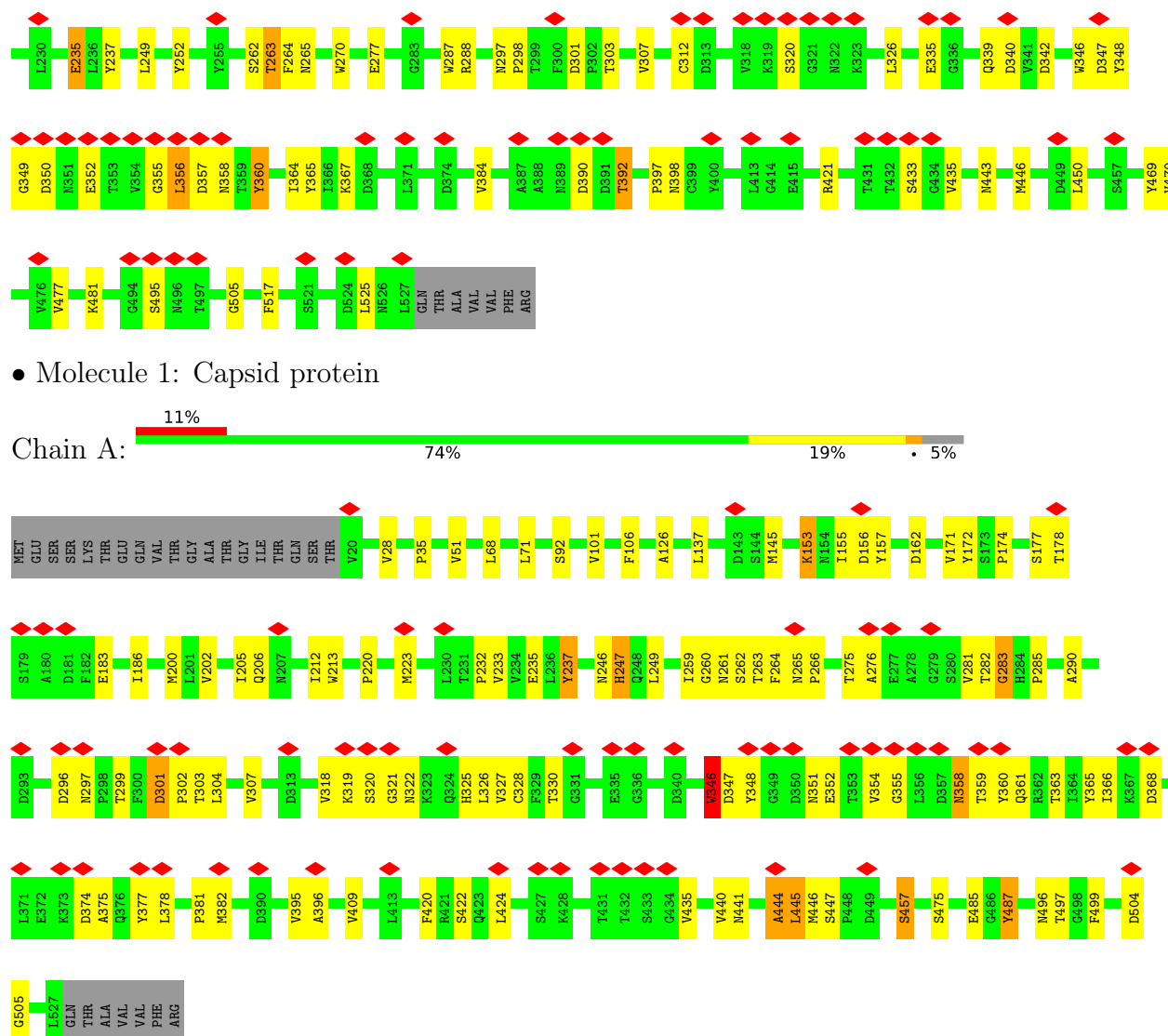
In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Capsid protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	C	526	Total	C	H	N	O	S	0	0
			7928	2583	3909	647	772	17		
1	B	508	Total	C	H	N	O	S	0	0
			7686	2510	3793	626	740	17		
1	A	508	Total	C	H	N	O	S	0	0
			7686	2510	3793	626	740	17		

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	3	SER	ASN	conflict	UNP B2Y6D0
C	284	HIS	ASN	conflict	UNP B2Y6D0
C	334	VAL	PHE	conflict	UNP B2Y6D0
C	335	GLU	ALA	conflict	UNP B2Y6D0
C	343	THR	ALA	conflict	UNP B2Y6D0
C	367	LYS	SER	conflict	UNP B2Y6D0
C	451	MET	ILE	conflict	UNP B2Y6D0
C	452	CYS	ARG	conflict	UNP B2Y6D0
B	3	SER	ASN	conflict	UNP B2Y6D0
B	284	HIS	ASN	conflict	UNP B2Y6D0
B	334	VAL	PHE	conflict	UNP B2Y6D0
B	335	GLU	ALA	conflict	UNP B2Y6D0
B	343	THR	ALA	conflict	UNP B2Y6D0
B	367	LYS	SER	conflict	UNP B2Y6D0
B	451	MET	ILE	conflict	UNP B2Y6D0
B	452	CYS	ARG	conflict	UNP B2Y6D0
A	3	SER	ASN	conflict	UNP B2Y6D0
A	284	HIS	ASN	conflict	UNP B2Y6D0
A	334	VAL	PHE	conflict	UNP B2Y6D0
A	335	GLU	ALA	conflict	UNP B2Y6D0
A	343	THR	ALA	conflict	UNP B2Y6D0
A	367	LYS	SER	conflict	UNP B2Y6D0
A	451	MET	ILE	conflict	UNP B2Y6D0
A	452	CYS	ARG	conflict	UNP B2Y6D0



• Molecule 1: Capsid protein

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	44239	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	23	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	81000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	57.077	Depositor
Minimum map value	-26.310	Depositor
Average map value	0.042	Depositor
Map value standard deviation	2.934	Depositor
Recommended contour level	12.8	Depositor
Map size ($\text{\AA}$)	618.24, 618.24, 618.24	wwPDB
Map dimensions	560, 560, 560	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.104, 1.104, 1.104	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.21	0/4007	0.59	10/5490 (0.2%)
1	B	0.20	0/4011	0.75	15/5502 (0.3%)
1	C	0.19	0/4137	0.54	8/5673 (0.1%)
All	All	0.20	0/12155	0.63	33/16665 (0.2%)

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	7
1	C	0	2
All	All	0	9

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	235	GLU	CA-C-N	23.61	164.19	121.70
1	B	235	GLU	C-N-CA	23.61	164.19	121.70
1	A	264	PHE	CA-C-N	11.48	142.36	121.70
1	A	264	PHE	C-N-CA	11.48	142.36	121.70
1	B	263	THR	CA-C-N	11.31	142.06	121.70
1	B	263	THR	C-N-CA	11.31	142.06	121.70
1	B	177	SER	CA-C-N	10.46	141.97	121.58
1	B	177	SER	C-N-CA	10.46	141.97	121.58
1	C	40	LEU	CA-C-N	10.40	144.19	121.63
1	C	40	LEU	C-N-CA	10.40	144.19	121.63
1	B	104	ASN	CA-C-N	9.49	136.56	120.72
1	B	104	ASN	C-N-CA	9.49	136.56	120.72
1	C	118	PRO	CA-C-N	9.41	136.05	122.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	118	PRO	C-N-CA	9.41	136.05	122.40
1	C	224	VAL	CA-C-N	8.43	136.87	121.70
1	C	224	VAL	C-N-CA	8.43	136.87	121.70
1	A	457	SER	CA-C-N	8.42	134.27	122.06
1	A	457	SER	C-N-CA	8.42	134.27	122.06
1	A	444	ALA	CA-C-N	8.16	136.40	121.70
1	A	444	ALA	C-N-CA	8.16	136.40	121.70
1	A	346	TRP	CA-C-N	6.88	134.08	121.70
1	A	346	TRP	C-N-CA	6.88	134.08	121.70
1	B	92	SER	CA-C-N	6.71	133.78	121.70
1	B	92	SER	C-N-CA	6.71	133.78	121.70
1	B	200	MET	CA-C-N	6.68	133.73	121.70
1	B	200	MET	C-N-CA	6.68	133.73	121.70
1	A	235	GLU	CA-C-N	5.95	130.72	121.26
1	A	235	GLU	C-N-CA	5.95	130.72	121.26
1	B	356	LEU	CA-C-N	5.51	131.62	121.70
1	B	356	LEU	C-N-CA	5.51	131.62	121.70
1	B	392	THR	N-CA-C	-5.34	107.42	114.04
1	C	388	ALA	CA-C-N	5.07	130.82	121.70
1	C	388	ALA	C-N-CA	5.07	130.82	121.70

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	153	LYS	Peptide
1	A	174	PRO	Peptide
1	A	206	GLN	Peptide
1	A	28	VAL	Peptide
1	A	283	GLY	Peptide
1	A	301	ASP	Peptide
1	A	504	ASP	Peptide
1	C	355	GLY	Peptide
1	C	490	ALA	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	3893	3793	3788	88	0
1	B	3893	3793	3791	59	0
1	C	4019	3909	3914	70	0
All	All	11805	11495	11493	208	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 9.

All (208) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:ASP:OD1	1:A:348:TYR:N	2.10	0.83
1:C:359:THR:OG1	1:C:362:ARG:HA	1.78	0.82
1:A:212:ILE:O	1:A:213:TRP:N	2.12	0.82
1:A:232:PRO:O	1:A:422:SER:OG	1.97	0.82
1:A:183:GLU:OE1	1:A:183:GLU:N	2.14	0.79
1:B:86:ARG:NH1	1:B:208:ASN:O	2.17	0.77
1:B:277:GLU:O	1:B:288:ARG:NH2	2.19	0.76
1:B:301:ASP:OD1	1:B:303:THR:OG1	2.04	0.74
1:A:409:VAL:HG11	1:A:487:TYR:OH	1.90	0.71
1:C:316:SER:OG	1:C:395:VAL:O	2.08	0.70
1:A:223:MET:SD	1:A:223:MET:N	2.65	0.70
1:C:227:ARG:NH1	1:C:443:ASN:O	2.24	0.70
1:C:51:VAL:HG22	1:C:78:ASN:OD1	1.92	0.70
1:B:56:PHE:CE2	1:B:69:LEU:HD13	2.28	0.69
1:B:352:GLU:OE1	1:B:352:GLU:N	2.25	0.69
1:A:213:TRP:O	1:A:457:SER:OG	2.10	0.69
1:A:420:PHE:HB2	1:A:444:ALA:HB2	1.75	0.68
1:C:155:ILE:HG23	1:A:155:ILE:HD13	1.75	0.68
1:A:496:ASN:OD1	1:A:497:THR:N	2.29	0.66
1:B:252:TYR:OH	1:B:263:THR:O	2.06	0.65
1:A:351:ASN:OD1	1:A:355:GLY:N	2.29	0.65
1:A:319:LYS:NZ	1:A:352:GLU:OE1	2.20	0.65
1:A:444:ALA:HB3	1:A:445:LEU:C	2.21	0.65
1:A:68:LEU:HD11	1:A:171:VAL:HG23	1.79	0.65
1:C:450:LEU:HD12	1:C:450:LEU:O	1.98	0.64
1:B:235:GLU:O	1:B:421:ARG:N	2.30	0.64
1:C:312:CYS:SG	1:C:314:TRP:NE1	2.71	0.62
1:C:21:THR:HG21	1:C:135:GLU:HG3	1.82	0.62
1:A:276:ALA:HB3	1:A:290:ALA:HA	1.80	0.62
1:C:296:ASP:O	1:C:296:ASP:OD2	2.17	0.62
1:C:312:CYS:HG	1:C:314:TRP:CD1	2.18	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:495:SER:OG	1:B:525:LEU:O	2.18	0.61
1:A:346:TRP:CD1	1:A:347:ASP:HB3	2.35	0.61
1:B:446:MET:SD	1:B:450:LEU:HD23	2.42	0.60
1:C:387:ALA:HB3	1:C:389:ASN:HB3	1.85	0.58
1:C:202:VAL:O	1:C:202:VAL:HG23	2.04	0.58
1:A:275:THR:OG1	1:A:375:ALA:O	2.04	0.57
1:B:200:MET:HB3	1:B:201:LEU:HG	1.86	0.57
1:A:283:GLY:O	1:A:285:PRO:O	2.21	0.57
1:C:78:ASN:ND2	1:C:190:SER:OG	2.37	0.57
1:C:311:PHE:O	1:C:313:ASP:N	2.38	0.56
1:A:233:VAL:HG12	1:A:233:VAL:O	2.05	0.55
1:A:381:PRO:O	1:A:382:MET:HG3	2.06	0.55
1:B:263:THR:OG1	1:B:265:ASN:N	2.39	0.55
1:C:359:THR:CB	1:C:362:ARG:HA	2.36	0.54
1:A:177:SER:O	1:A:177:SER:OG	2.23	0.54
1:C:2:GLU:HA	1:C:20:VAL:O	2.06	0.54
1:C:154:ASN:O	1:A:155:ILE:HG21	2.08	0.54
1:C:56:PHE:CE1	1:C:68:LEU:HD22	2.43	0.54
1:A:223:MET:HB3	1:A:232:PRO:HA	1.90	0.53
1:C:155:ILE:HG23	1:A:155:ILE:CD1	2.36	0.53
1:A:299:THR:HG21	1:A:303:THR:HG22	1.89	0.53
1:A:328:CYS:SG	1:A:330:THR:OG1	2.66	0.53
1:C:5:LYS:NZ	1:C:5:LYS:HB3	2.24	0.53
1:B:349:GLY:O	1:B:357:ASP:N	2.41	0.53
1:A:302:PRO:O	1:A:304:LEU:N	2.42	0.53
1:C:485:GLU:N	1:C:486:GLY:HA2	2.24	0.52
1:B:265:ASN:ND2	1:A:223:MET:C	2.67	0.52
1:A:395:VAL:HB	1:A:396:ALA:HB2	1.90	0.52
1:A:440:VAL:HG22	1:A:441:ASN:H	1.74	0.52
1:A:445:LEU:C	1:A:445:LEU:HD23	2.35	0.52
1:B:355:GLY:O	1:B:358:ASN:N	2.43	0.51
1:C:83:HIS:O	1:C:86:ARG:HG2	2.10	0.51
1:C:295:PRO:O	1:C:296:ASP:OD2	2.28	0.51
1:C:155:ILE:CG2	1:A:155:ILE:HD13	2.40	0.51
1:C:366:ILE:O	1:C:366:ILE:HG22	2.11	0.51
1:B:56:PHE:HE2	1:B:69:LEU:HD13	1.72	0.51
1:B:249:LEU:HD11	1:B:312:CYS:HA	1.92	0.51
1:A:101:VAL:HG13	1:A:186:ILE:CD1	2.40	0.51
1:B:200:MET:SD	1:B:201:LEU:HG	2.51	0.50
1:C:127:TYR:O	1:C:130:THR:OG1	2.28	0.50
1:A:106:PHE:CD1	1:A:106:PHE:N	2.79	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:388:ALA:HB3	1:C:389:ASN:HA	1.93	0.50
1:A:246:ASN:O	1:A:247:HIS:O	2.28	0.50
1:B:433:SER:HB3	1:B:435:VAL:HG12	1.94	0.50
1:C:400:TYR:OH	1:A:301:ASP:OD2	2.07	0.49
1:C:212:ILE:HG22	1:C:509:PHE:HE2	1.77	0.49
1:B:435:VAL:O	1:B:435:VAL:HG13	2.11	0.49
1:B:348:TYR:O	1:B:365:TYR:CE2	2.65	0.49
1:A:325:HIS:H	1:A:348:TYR:HE1	1.61	0.49
1:B:356:LEU:HA	1:B:358:ASN:H	1.78	0.49
1:C:312:CYS:HG	1:C:314:TRP:NE1	2.10	0.49
1:C:318:VAL:HG12	1:C:319:LYS:N	2.28	0.49
1:B:348:TYR:CE1	1:B:364:ILE:HG22	2.48	0.49
1:A:265:ASN:N	1:A:266:PRO:HD3	2.28	0.49
1:A:446:MET:SD	1:A:447:SER:N	2.86	0.48
1:B:160:ARG:NH2	1:B:208:ASN:OD1	2.43	0.48
1:B:390:ASP:O	1:B:392:THR:N	2.39	0.48
1:A:409:VAL:HG21	1:A:487:TYR:CE1	2.47	0.48
1:B:56:PHE:CE2	1:B:69:LEU:HB2	2.49	0.48
1:B:56:PHE:CZ	1:B:69:LEU:HD13	2.48	0.48
1:A:153:LYS:HB2	1:A:157:TYR:CE1	2.49	0.48
1:C:68:LEU:HD11	1:C:171:VAL:HG23	1.96	0.48
1:C:485:GLU:N	1:C:486:GLY:CA	2.76	0.48
1:B:335:GLU:CD	1:B:335:GLU:O	2.57	0.48
1:A:249:LEU:O	1:A:262:SER:OG	2.08	0.47
1:A:327:VAL:HG12	1:A:328:CYS:N	2.28	0.47
1:C:172:TYR:OH	1:B:50:GLU:OE2	2.22	0.47
1:B:41:ASP:OD1	1:B:44:VAL:HG23	2.14	0.47
1:A:354:VAL:O	1:A:358:ASN:N	2.47	0.47
1:C:312:CYS:O	1:C:314:TRP:NE1	2.48	0.47
1:A:301:ASP:HB2	1:A:302:PRO:CD	2.45	0.47
1:A:101:VAL:HG21	1:A:137:LEU:HD21	1.96	0.47
1:A:395:VAL:CB	1:A:396:ALA:HB2	2.45	0.47
1:B:469:TYR:O	1:B:477:VAL:HB	2.15	0.46
1:C:298:PRO:O	1:C:300:PHE:CD2	2.68	0.46
1:B:348:TYR:HE1	1:B:364:ILE:HG22	1.80	0.46
1:A:259:ILE:HG22	1:A:260:GLY:N	2.29	0.46
1:A:281:VAL:HG22	1:A:282:THR:H	1.80	0.46
1:C:483:TYR:O	1:C:486:GLY:HA3	2.16	0.46
1:C:306:PRO:HB2	1:C:360:TYR:O	2.16	0.46
1:C:391:ASP:OD2	1:C:391:ASP:N	2.49	0.46
1:C:491:ASN:C	1:C:491:ASN:HD22	2.24	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:LEU:HB2	1:A:346:TRP:O	2.15	0.46
1:C:483:TYR:O	1:C:486:GLY:CA	2.64	0.46
1:C:195:ASP:OD1	1:C:195:ASP:C	2.58	0.46
1:B:200:MET:HB3	1:B:201:LEU:CG	2.46	0.46
1:A:283:GLY:O	1:A:285:PRO:N	2.49	0.46
1:C:318:VAL:HG12	1:C:319:LYS:H	1.81	0.46
1:A:156:ASP:O	1:A:162:ASP:OD2	2.34	0.45
1:C:358:ASN:HB3	1:C:400:TYR:CE2	2.51	0.45
1:B:307:VAL:HG21	1:B:360:TYR:CD1	2.52	0.45
1:A:321:GLY:O	1:A:325:HIS:NE2	2.50	0.45
1:B:357:ASP:CG	1:B:357:ASP:O	2.59	0.45
1:C:387:ALA:CB	1:C:389:ASN:HB3	2.45	0.45
1:B:339:GLN:O	1:B:340:ASP:C	2.59	0.45
1:C:319:LYS:C	1:C:321:GLY:N	2.75	0.45
1:B:249:LEU:O	1:B:262:SER:OG	2.26	0.45
1:A:359:THR:C	1:A:361:GLN:N	2.74	0.45
1:A:365:TYR:HB3	1:A:366:ILE:N	2.32	0.45
1:C:301:ASP:HB3	1:C:304:LEU:HD23	1.99	0.44
1:C:199:THR:O	1:C:200:MET:C	2.59	0.44
1:C:203:PRO:N	1:C:204:PRO:CD	2.80	0.44
1:C:291:LEU:HD22	1:C:361:GLN:HG2	1.98	0.44
1:B:263:THR:HA	1:B:264:PHE:HB2	1.99	0.44
1:B:443:ASN:OD1	1:B:443:ASN:N	2.51	0.44
1:C:288:ARG:HB3	1:C:288:ARG:NH1	2.33	0.44
1:B:320:SER:OG	1:B:350:ASP:OD1	2.32	0.44
1:A:183:GLU:H	1:A:183:GLU:CD	2.23	0.44
1:A:318:VAL:HG13	1:A:319:LYS:N	2.33	0.44
1:B:41:ASP:OD1	1:B:41:ASP:C	2.61	0.44
1:C:354:VAL:O	1:C:356:LEU:N	2.51	0.44
1:B:470:VAL:O	1:B:505:GLY:HA2	2.18	0.44
1:A:92:SER:HA	1:A:157:TYR:HE2	1.82	0.44
1:A:153:LYS:CB	1:A:157:TYR:CE1	3.01	0.44
1:A:178:THR:HG22	1:A:178:THR:O	2.18	0.44
1:B:88:TYR:CE1	1:B:200:MET:HG3	2.53	0.43
1:B:348:TYR:HB2	1:B:357:ASP:CG	2.43	0.43
1:A:348:TYR:O	1:A:365:TYR:CZ	2.71	0.43
1:C:480:PHE:HD1	1:C:490:ALA:HB2	1.83	0.43
1:C:77:LEU:HA	1:C:78:ASN:OD1	2.18	0.43
1:B:263:THR:CB	1:B:265:ASN:H	2.32	0.43
1:C:131:GLY:O	1:B:80:TYR:OH	2.36	0.43
1:A:330:THR:HA	1:A:378:LEU:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:ASP:O	1:A:297:ASN:OD1	2.35	0.43
1:B:270:TRP:CH2	1:B:312:CYS:HB3	2.53	0.43
1:B:326:LEU:HD23	1:B:384:VAL:HG12	2.01	0.43
1:B:365:TYR:CD1	1:B:365:TYR:C	2.97	0.43
1:C:202:VAL:C	1:C:204:PRO:HD2	2.43	0.43
1:A:200:MET:O	1:A:202:VAL:HG13	2.18	0.43
1:A:237:TYR:HB2	1:A:499:PHE:O	2.18	0.43
1:B:287:TRP:HB3	1:B:367:LYS:HA	2.01	0.43
1:B:346:TRP:O	1:B:347:ASP:C	2.61	0.43
1:C:372:GLU:N	1:C:377:TYR:OH	2.52	0.43
1:C:224:VAL:O	1:C:447:SER:CB	2.67	0.42
1:C:404:ASP:OD2	1:C:404:ASP:N	2.35	0.42
1:A:51:VAL:HG11	1:A:71:LEU:HD21	2.01	0.42
1:A:358:ASN:C	1:A:358:ASN:HD22	2.26	0.42
1:C:400:TYR:CD1	1:C:400:TYR:C	2.98	0.42
1:C:51:VAL:HA	1:C:78:ASN:OD1	2.20	0.42
1:C:352:GLU:OE1	1:C:352:GLU:N	2.48	0.42
1:A:320:SER:O	1:A:322:ASN:OD1	2.37	0.42
1:A:220:PRO:CD	1:A:505:GLY:HA2	2.49	0.42
1:B:40:LEU:HD12	1:B:40:LEU:N	2.35	0.42
1:A:485:GLU:HB2	1:A:487:TYR:CE2	2.55	0.42
1:C:34:ALA:HB2	1:C:157:TYR:CE2	2.55	0.42
1:C:291:LEU:CD2	1:C:363:THR:HG22	2.50	0.42
1:B:297:ASN:N	1:B:298:PRO:CD	2.83	0.41
1:A:445:LEU:HD23	1:A:446:MET:N	2.35	0.41
1:A:263:THR:HB	1:A:265:ASN:O	2.20	0.41
1:C:276:ALA:HB3	1:C:290:ALA:HA	2.02	0.41
1:A:327:VAL:O	1:A:381:PRO:O	2.39	0.41
1:A:374:ASP:CG	1:A:377:TYR:HD1	2.28	0.41
1:C:74:SER:HB3	1:C:164:LEU:HD13	2.02	0.41
1:B:41:ASP:OD1	1:B:41:ASP:O	2.38	0.41
1:B:347:ASP:OD1	1:B:348:TYR:N	2.48	0.41
1:A:366:ILE:O	1:A:368:ASP:N	2.52	0.41
1:A:395:VAL:HA	1:A:396:ALA:HB2	2.03	0.41
1:A:296:ASP:O	1:A:297:ASN:C	2.63	0.41
1:B:339:GLN:HG3	1:B:340:ASP:N	2.35	0.41
1:B:358:ASN:O	1:B:398:ASN:HB2	2.21	0.41
1:B:481:LYS:NZ	1:B:517:PHE:O	2.27	0.41
1:A:395:VAL:CA	1:A:396:ALA:HB2	2.51	0.41
1:A:35:PRO:HD2	1:A:200:MET:HE2	2.03	0.41
1:A:360:TYR:HB2	1:A:363:THR:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:87:MET:O	1:C:202:VAL:N	2.51	0.41
1:B:265:ASN:HD22	1:A:223:MET:HB2	1.86	0.41
1:A:233:VAL:CG1	1:A:420:PHE:HB3	2.51	0.41
1:A:318:VAL:HG22	1:A:319:LYS:H	1.86	0.41
1:A:233:VAL:HG13	1:A:420:PHE:HB3	2.02	0.40
1:A:347:ASP:OD1	1:A:347:ASP:C	2.63	0.40
1:B:360:TYR:HB2	1:B:398:ASN:O	2.21	0.40
1:A:137:LEU:HB2	1:A:145:MET:SD	2.62	0.40
1:C:314:TRP:O	1:C:382:MET:HE1	2.22	0.40
1:A:307:VAL:HG12	1:A:361:GLN:O	2.20	0.40
1:C:288:ARG:O	1:C:365:TYR:HA	2.22	0.40
1:C:480:PHE:CD1	1:C:490:ALA:HB2	2.57	0.40
1:A:126:ALA:HB1	1:A:172:TYR:HD1	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	499/534 (93%)	428 (86%)	70 (14%)	1 (0%)	43	58
1	B	506/534 (95%)	445 (88%)	60 (12%)	1 (0%)	43	58
1	C	524/534 (98%)	469 (90%)	52 (10%)	3 (1%)	21	31
All	All	1529/1602 (95%)	1342 (88%)	182 (12%)	5 (0%)	37	50

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	247	HIS
1	C	211	PRO
1	C	212	ILE
1	C	381	PRO

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Mol	Chain	Res	Type
1	B	397	PRO

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	431/453 (95%)	421 (98%)	10 (2%)	44	65
1	B	431/453 (95%)	424 (98%)	7 (2%)	55	73
1	C	446/453 (98%)	438 (98%)	8 (2%)	51	71
All	All	1308/1359 (96%)	1283 (98%)	25 (2%)	49	70

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	16	THR
1	C	155	ILE
1	C	237	TYR
1	C	244	THR
1	C	312	CYS
1	C	404	ASP
1	C	491	ASN
1	C	504	ASP
1	B	112	ILE
1	B	119	ASN
1	B	156	ASP
1	B	222	SER
1	B	237	TYR
1	B	342	ASP
1	B	360	TYR
1	A	205	ILE
1	A	237	TYR
1	A	261	ASN
1	A	346	TRP
1	A	358	ASN
1	A	424	LEU

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Mol	Chain	Res	Type
1	A	435	VAL
1	A	445	LEU
1	A	475	SER
1	A	487	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	47	ASN
1	C	63	GLN
1	C	104	ASN
1	C	217	GLN
1	C	339	GLN
1	C	394	GLN
1	C	443	ASN
1	C	491	ASN
1	B	47	ASN
1	B	83	HIS
1	B	265	ASN
1	B	441	ASN
1	A	119	ASN
1	A	376	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	4

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	365:TYR	C	366:ILE	N	3.28
1	A	212:ILE	C	213:TRP	N	3.20
1	A	403:TRP	C	404:ASP	N	3.04
1	A	213:TRP	C	214:SER	N	2.95

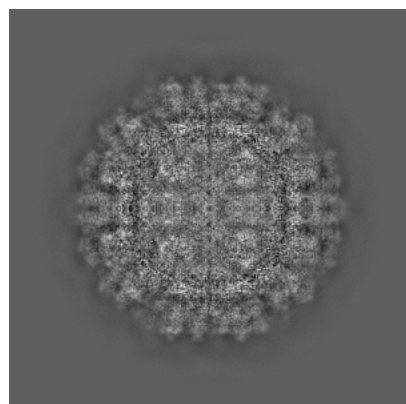
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-43292. These allow visual inspection of the internal detail of the map and identification of artifacts.

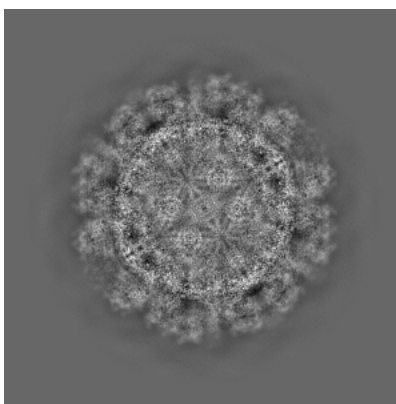
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

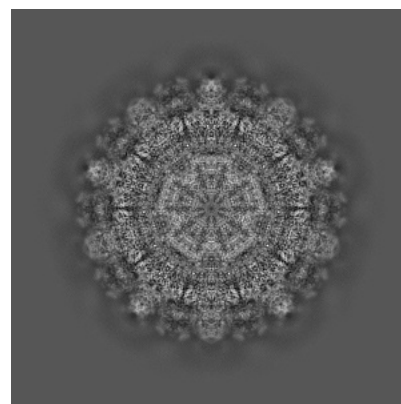
6.1.1 Primary map



X

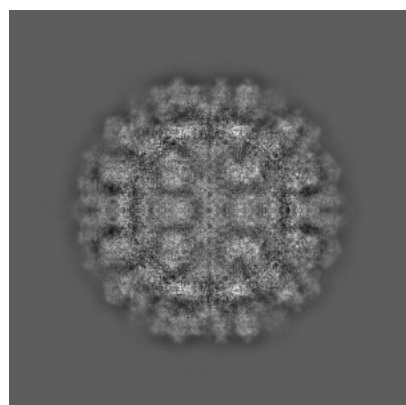


Y

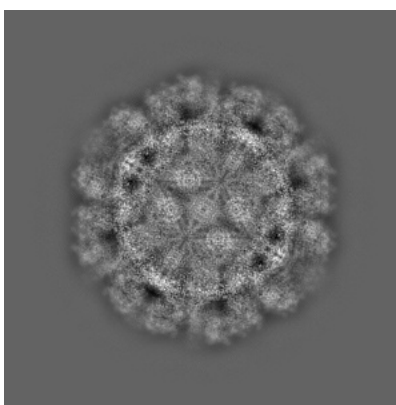


Z

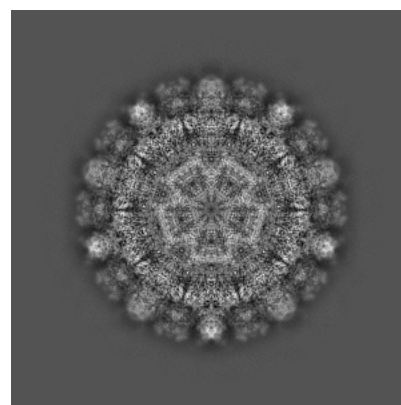
6.1.2 Raw map



X



Y

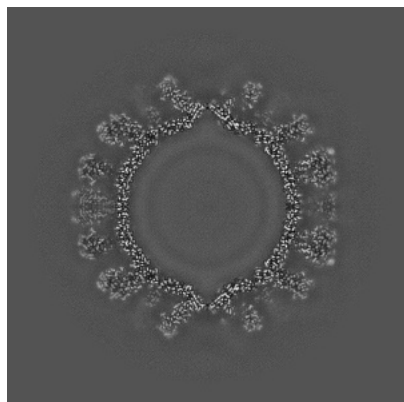


Z

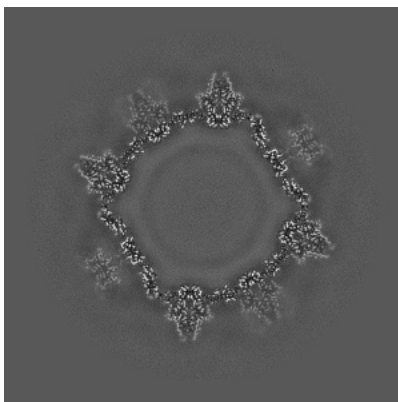
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

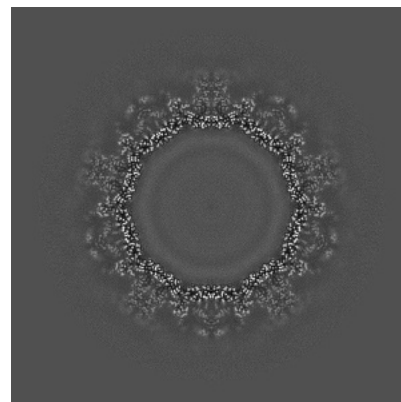
6.2.1 Primary map



X Index: 280

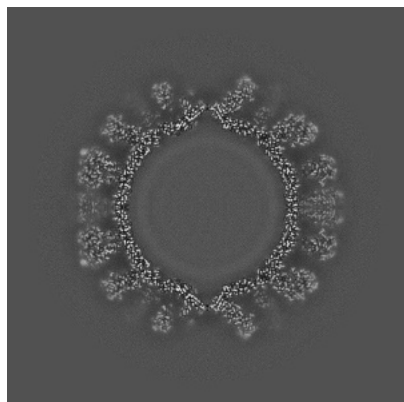


Y Index: 280

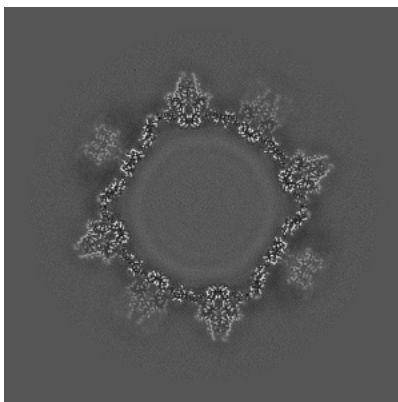


Z Index: 280

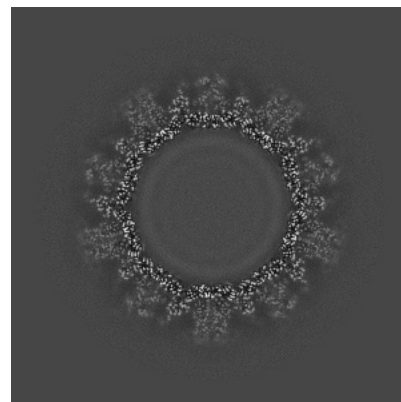
6.2.2 Raw map



X Index: 280



Y Index: 280

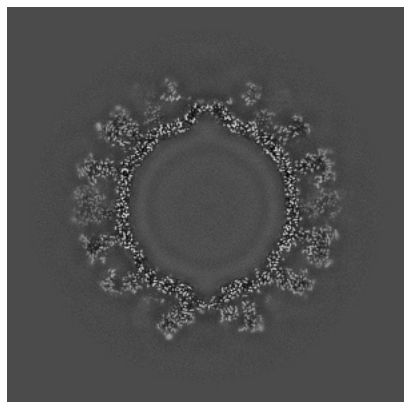


Z Index: 280

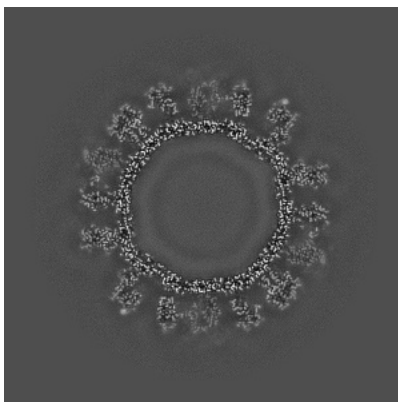
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

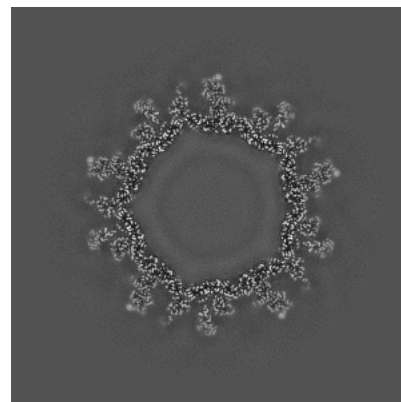
6.3.1 Primary map



X Index: 285

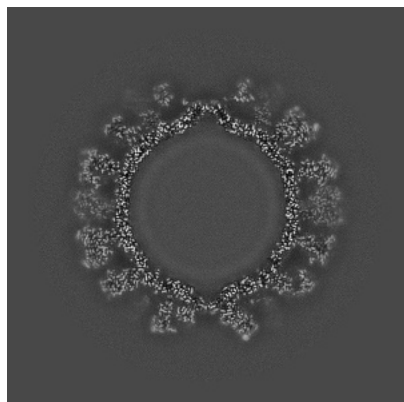


Y Index: 240

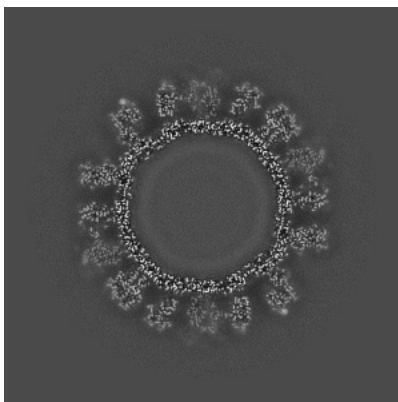


Z Index: 238

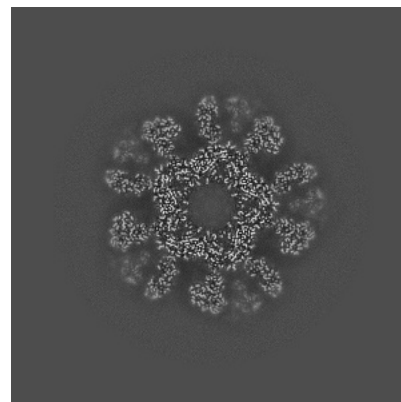
6.3.2 Raw map



X Index: 275



Y Index: 320

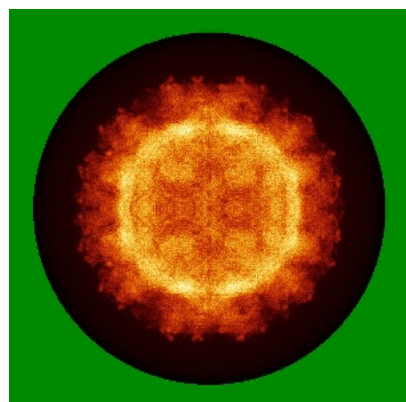


Z Index: 384

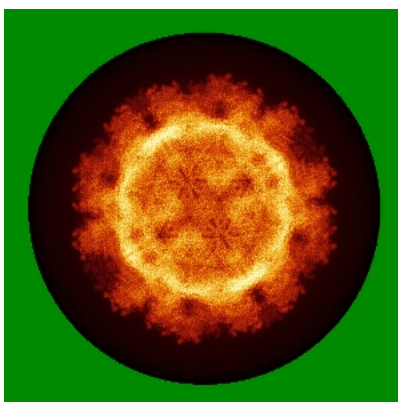
The images above show the largest variance slices of the map in three orthogonal directions.

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

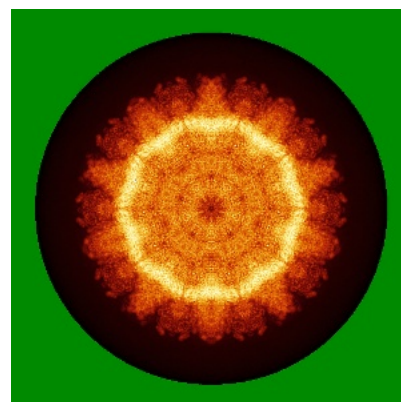
6.4.1 Primary map



X

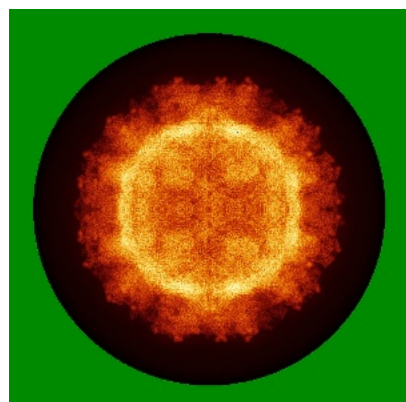


Y

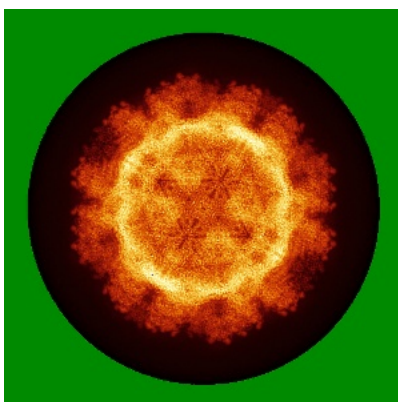


Z

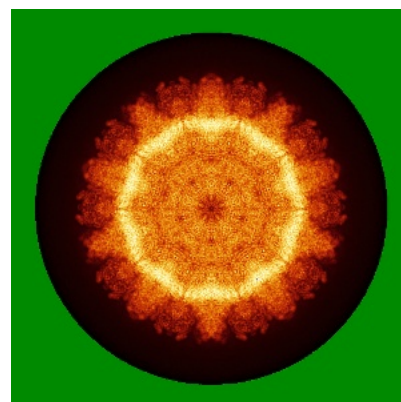
6.4.2 Raw map



X



Y

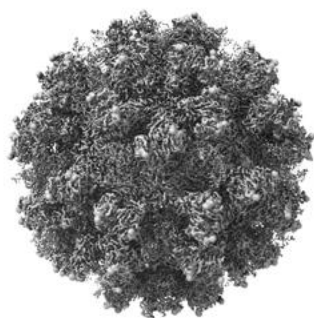


Z

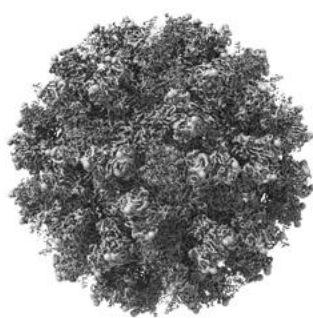
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

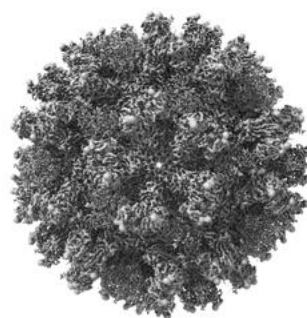
6.5.1 Primary map



X



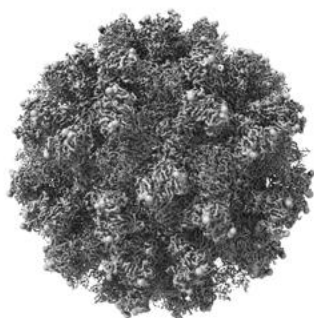
Y



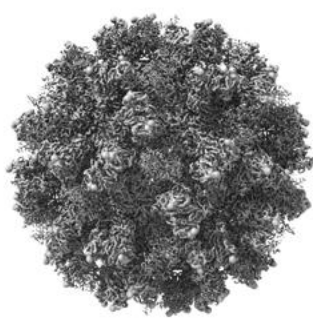
Z

The images above show the 3D surface view of the map at the recommended contour level 12.8. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

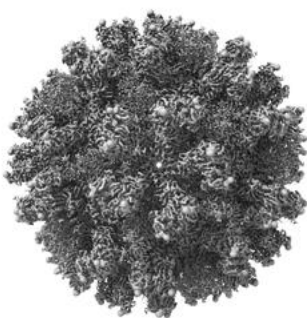
6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

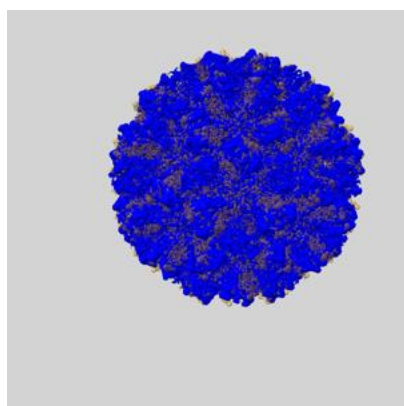
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

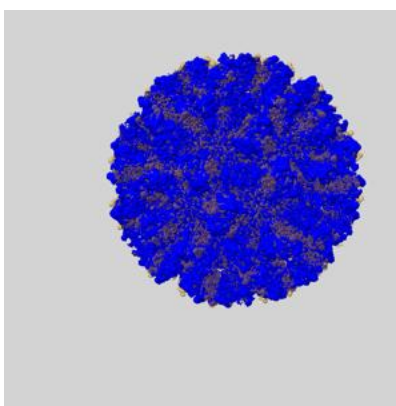
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

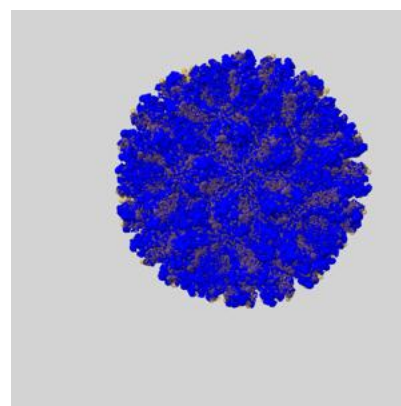
6.6.1 emd_43292_msk_1.map [i](#)



X

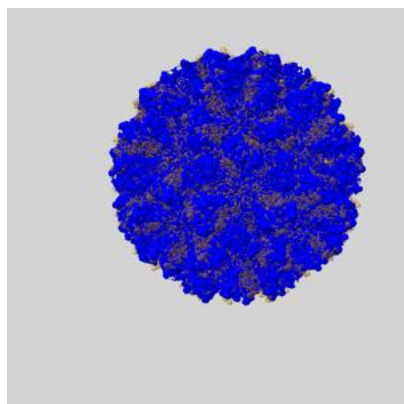


Y

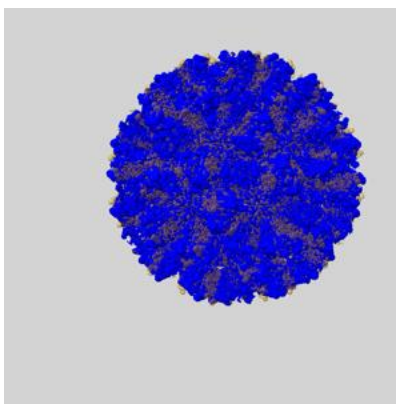


Z

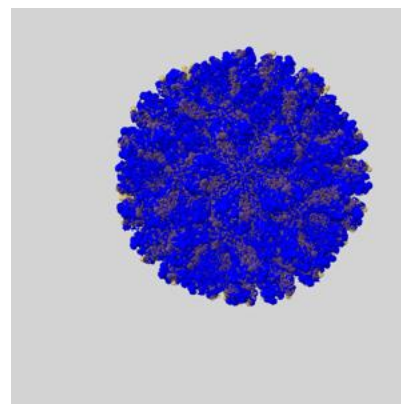
6.6.2 emd_43292_msk_2.map [i](#)



X



Y

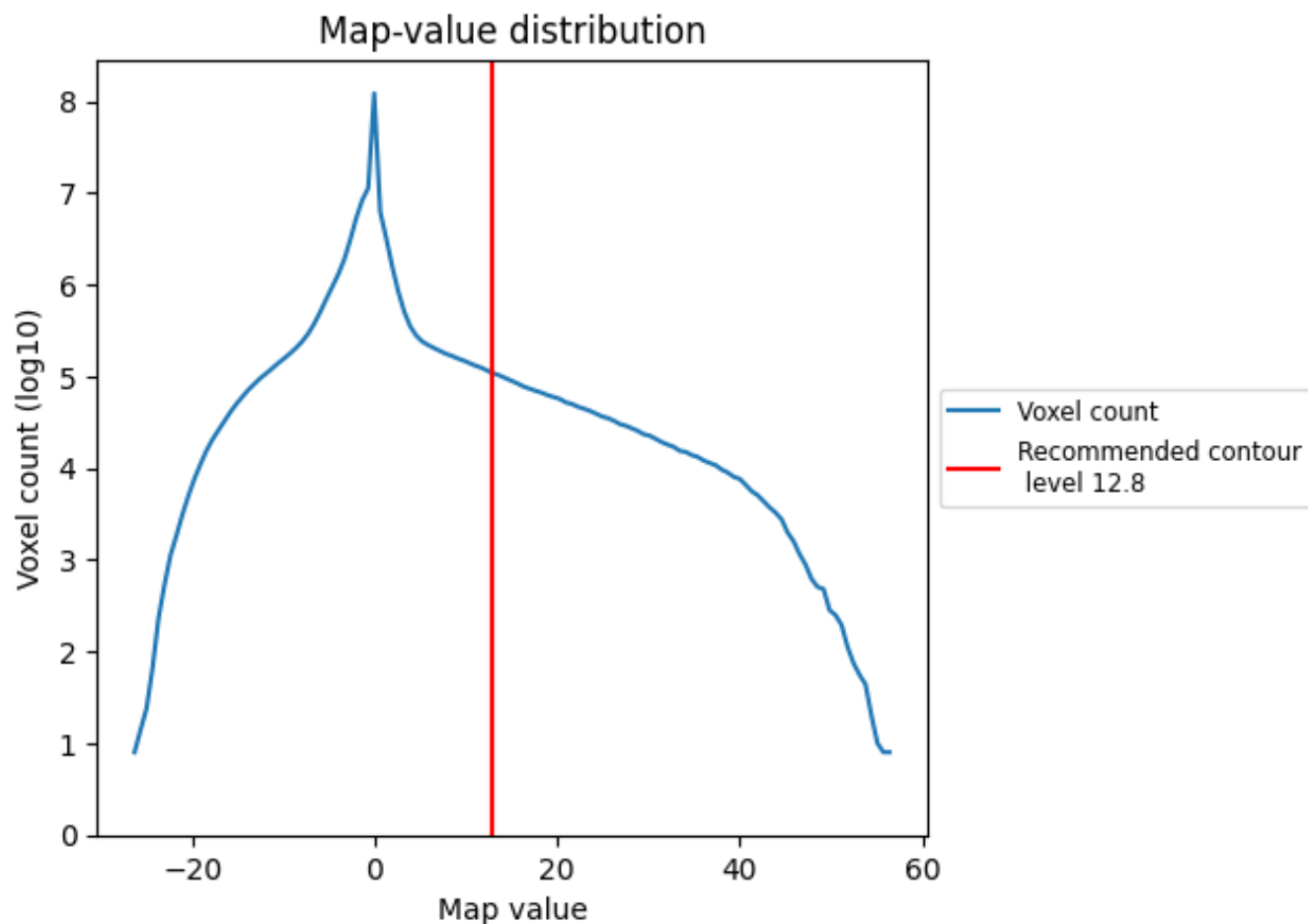


Z

7 Map analysis [i](#)

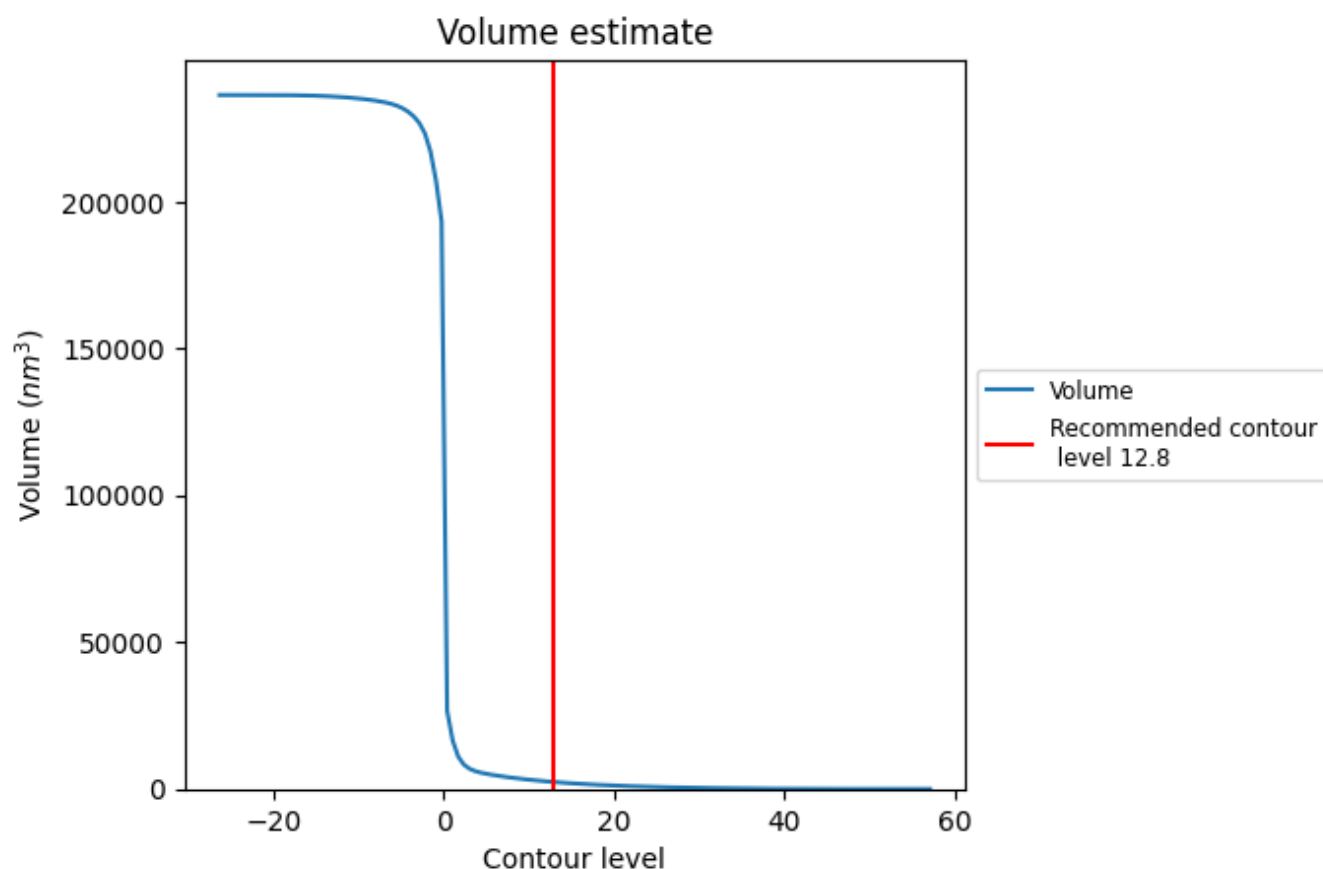
This section contains the results of statistical analysis of the map.

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.

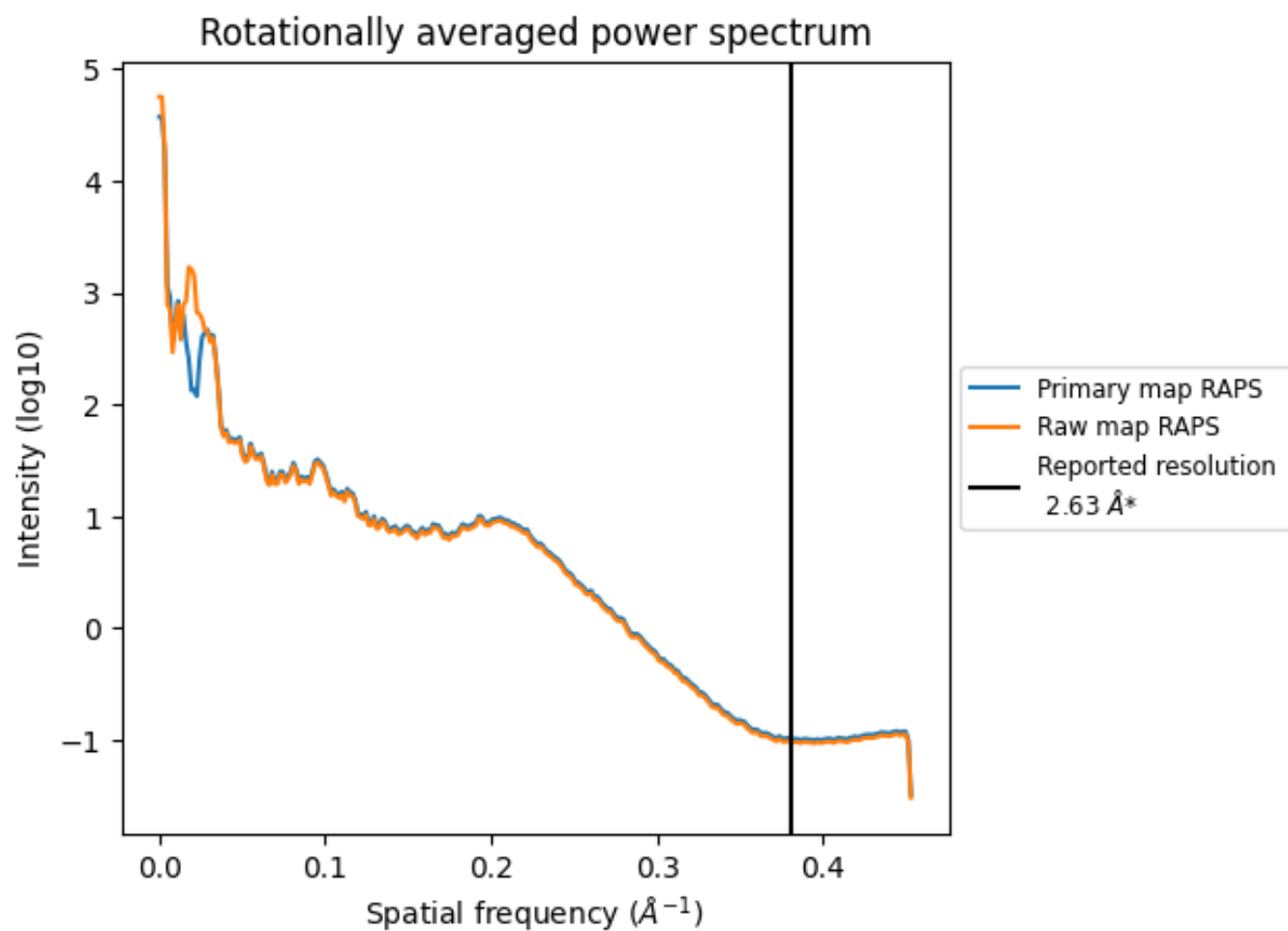
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2390 nm³; this corresponds to an approximate mass of 2159 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 ⓘ

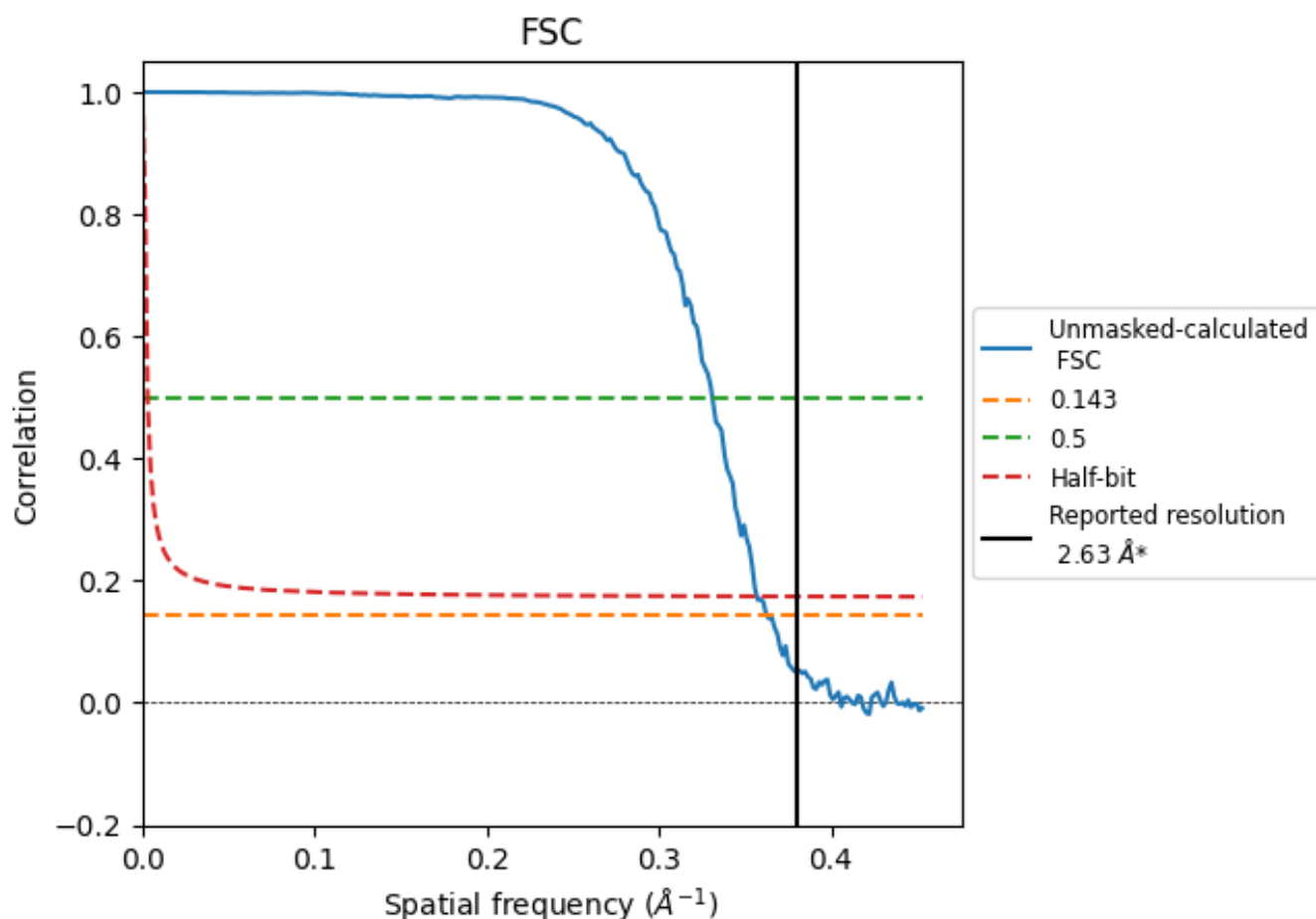


*Reported resolution corresponds to spatial frequency of 0.380 Å⁻¹

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 0.380 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.63	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	2.76	3.02	2.80

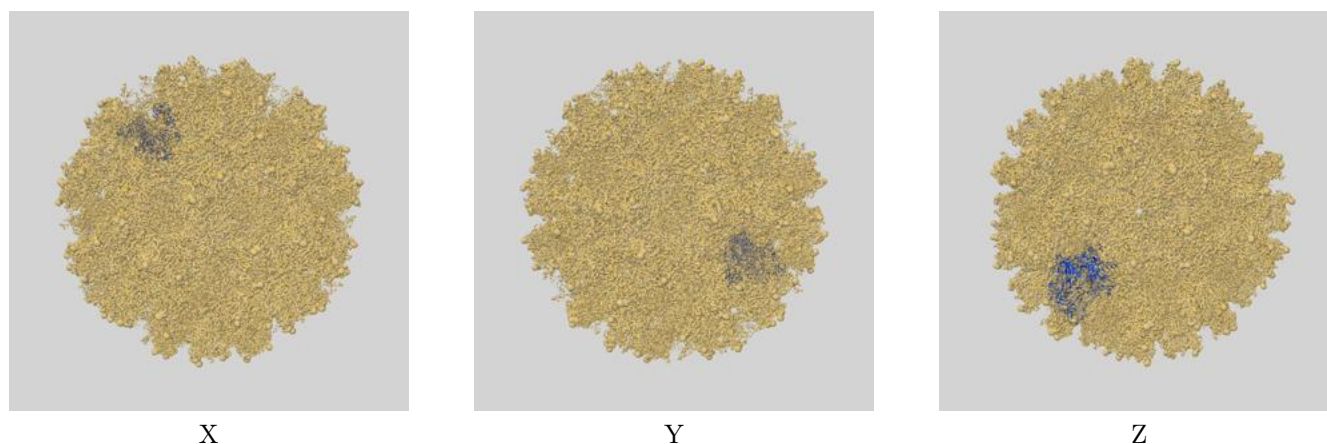
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

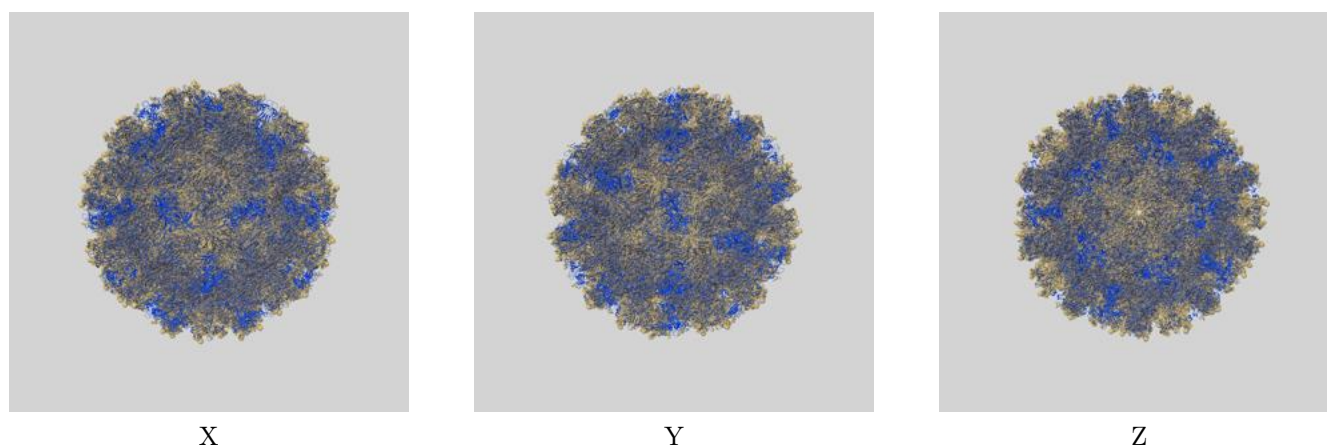
This section contains information regarding the fit between EMDB map EMD-43292 and PDB model 8VJR. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlays

9.1.1 Map-model overlay [i](#)

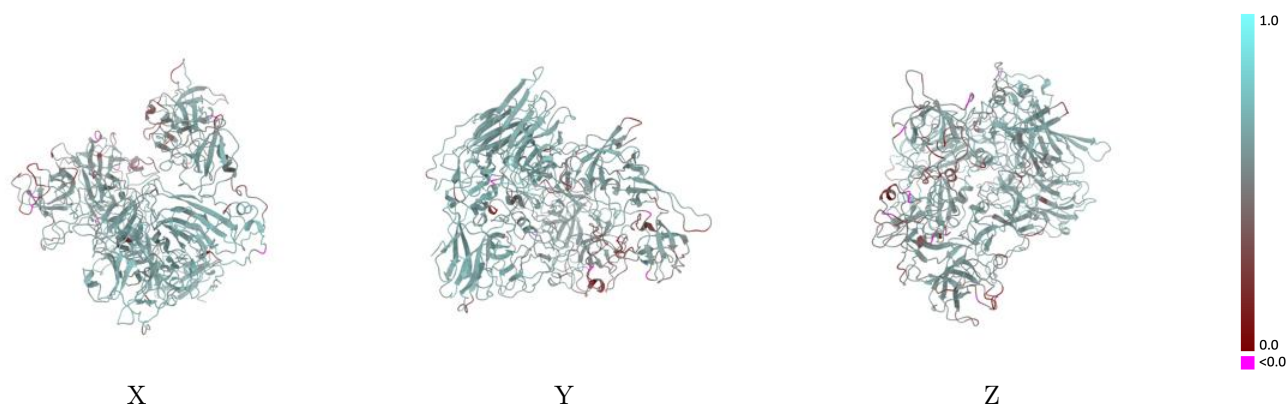


9.1.2 Map-model assembly overlay [i](#)



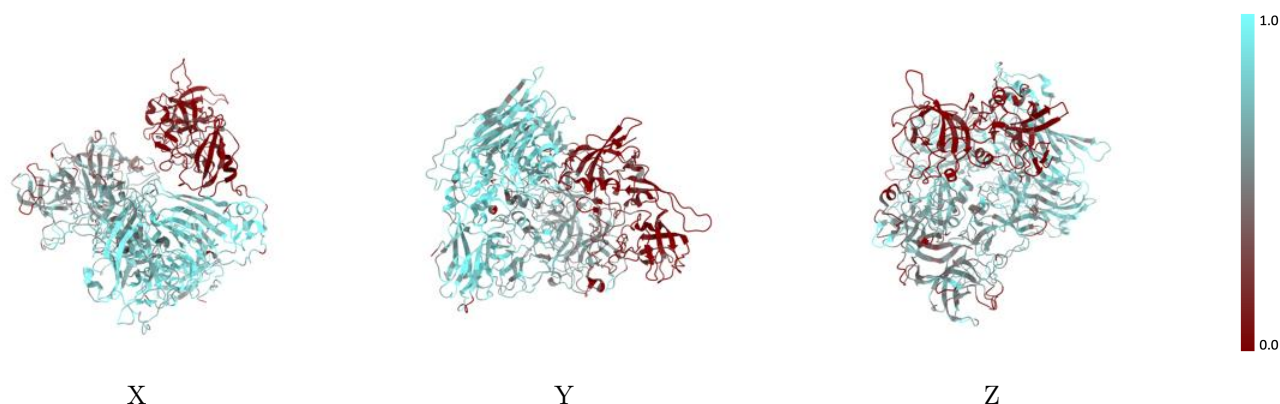
The images above show the 3D surface view of the map at the recommended contour level 12.8 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



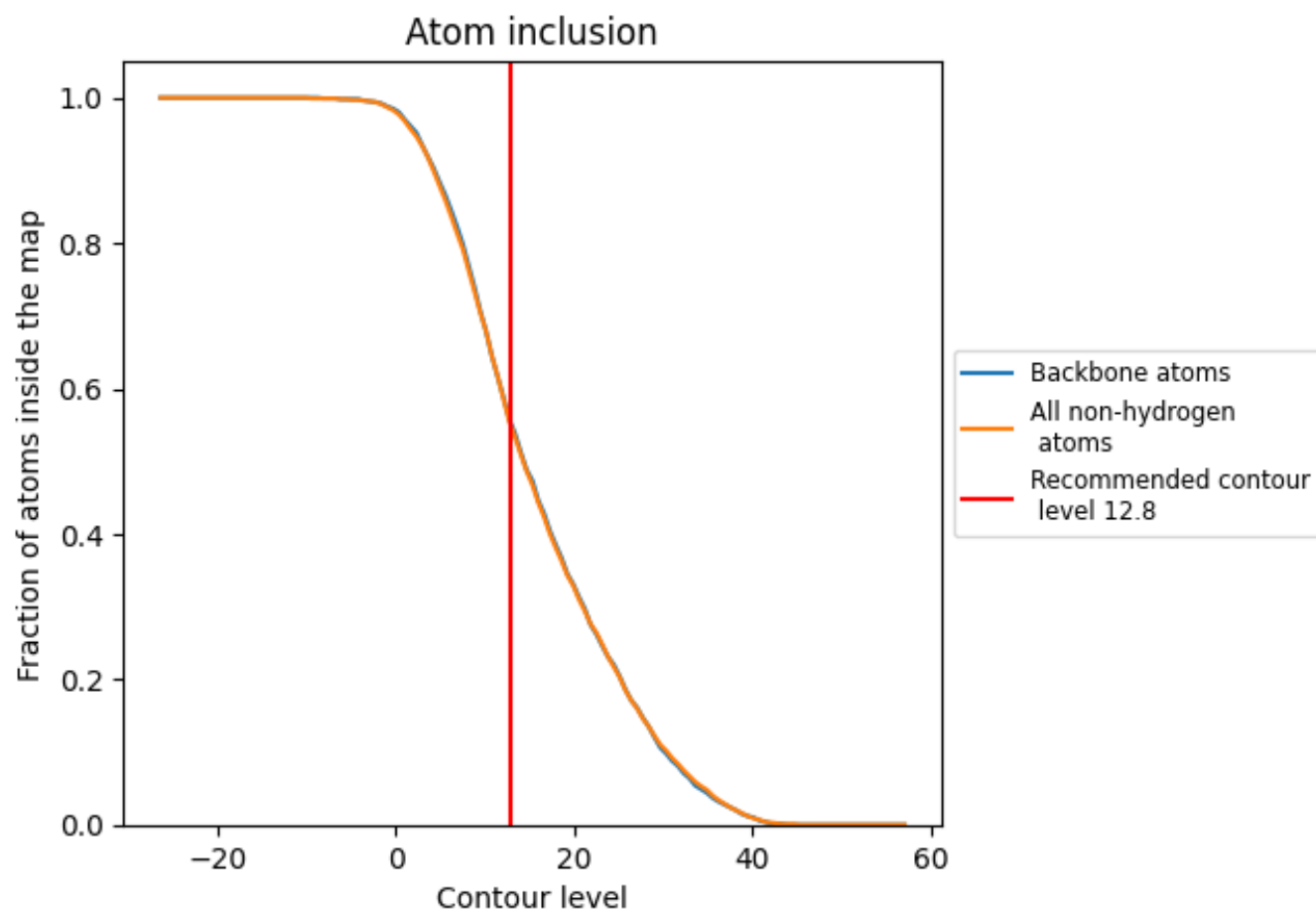
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (12.8).

9.4 Atom inclusion [i](#)



At the recommended contour level, 56% of all backbone atoms, 55% 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 (12.8) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5540	0.5590
A	0.6630	0.5570
B	0.6630	0.5640
C	0.3690	0.5550

