



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 08:30 AM UTC

PDB ID : 8T13 / pdb_00008t13
EMDB ID : EMD-40953
Title : Cryo-EM structure of DENV2 NS5 in complex with human STAT2 with the N-terminal domain of STAT2 disordered
Authors : Biswal, M.; Lu, J.; Song, J.
Deposited on : 2023-06-01
Resolution : 3.45 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
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

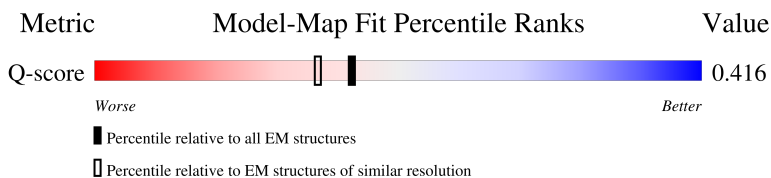
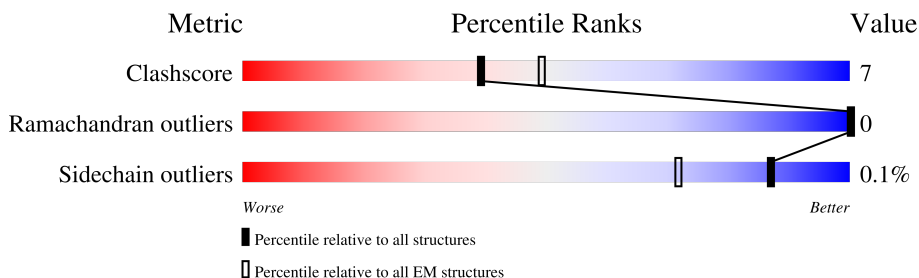
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.45 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	13836 (2.95 - 3.95)

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	851	
2	B	900	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9304 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 Signal transducer and activator of transcription 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	290	Total	C	N	O	S	0	0
			2311	1475	403	427	6		

- Molecule 2 is a protein called Non-structural protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	877	Total	C	N	O	S	0	0
			6991	4395	1249	1305	42		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
3	B	2	Total	Zn	0
			2	2	

[illegible]

- Molecule 2: Non-structural protein 5

Chain B: 82% 15% .

I867	I868	R871	S872	L873	N876	Y879	M883	PRO	SER	MET	LYS	ARG	PHE	ARG	ARG	GLU	GLU	GLU	GLY	ALA	GLY	VAL	LEU	TRP	D683	M684	K689	Q692	E695	M718	L724	V725	M730	E733	R737	S747	L748	T751	L754	Y758	R769	L772	V783	P784	S785	H786	W787	M809	N814	E825	D826	V830	Q845	S849	L850	I851	G852	L853	T854	W859	O864	E192	K193	M194	R207	L210	T215	M218	S222	N223	N227	S231	I235	T240	N241	R242	F243	T244	K249	R263	E268	K283	W292	H306	V321	N322	V325	R326	E357	K358	E365	P366	T370	M374	K375	R392	T395	E398	V403	L409	G410	A411	I412	M417	K418	W419	V426	L441	T448	W453	R458	K467	Y476	G516	K524	A536	E550	H561	I592	T613	V637	T638	W646	R654	M658	V667	D671	D672	R673	L678	N682	GLY	THR	ASN	THR	GLY	E7	W13	K14	L17	L20	E24	R38	K42	E43	K46	V55	S56	R57	K61	L62	V66	R84	K95	K101	K105	E111	M116	R125	P136	P137	T142	S151	E157	Q177	N184	GLY
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	86928	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.149	Depositor
Minimum map value	-1.329	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.025	Depositor
Recommended contour level	0.175	Depositor
Map size (Å)	444.0, 444.0, 444.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.11, 1.11, 1.11	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.09	0/2344	0.25	0/3161
2	B	0.10	0/7150	0.29	0/9669
All	All	0.10	0/9494	0.28	0/12830

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2311	0	2370	41	0
2	B	6991	0	6824	92	0
3	B	2	0	0	0	0
All	All	9304	0	9194	133	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 7.

All (133) 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:235:LEU:HB3	1:A:314:ARG:HH21	1.58	0.67
2:B:868:ASN:HA	2:B:871:ARG:HD3	1.75	0.67
2:B:868:ASN:OD1	2:B:871:ARG:NH1	2.30	0.65
2:B:403:VAL:HG11	2:B:426:VAL:HG21	1.79	0.65
2:B:207:ARG:NH2	2:B:215:THR:O	2.29	0.64
2:B:754:LEU:HG	2:B:758:TYR:CE2	2.33	0.64
2:B:783:VAL:HG12	2:B:785:SER:H	1.64	0.63
2:B:157:GLU:OE1	2:B:184:ASN:ND2	2.34	0.61
1:A:376:PHE:HB3	1:A:405:LEU:HD22	1.83	0.60
1:A:170:GLN:HE22	1:A:285:LEU:HD23	1.67	0.59
2:B:43:GLU:HA	2:B:46:LYS:HD2	1.83	0.59
1:A:376:PHE:HD2	1:A:405:LEU:HB3	1.68	0.59
1:A:153:ARG:HD3	1:A:268:LEU:HD11	1.84	0.59
1:A:235:LEU:HD23	1:A:314:ARG:HE	1.67	0.59
2:B:737:ARG:O	2:B:737:ARG:HD3	2.03	0.58
2:B:784:PRO:HA	2:B:879:TYR:HA	1.84	0.58
2:B:241:ASN:O	2:B:244:THR:OG1	2.20	0.58
1:A:236:GLU:OE2	1:A:314:ARG:NH1	2.37	0.57
2:B:785:SER:OG	2:B:786:HIS:N	2.38	0.57
2:B:658:MET:HG2	2:B:667:VAL:HG22	1.87	0.56
2:B:322:ASN:OD1	2:B:751:THR:OG1	2.24	0.56
2:B:678:LEU:O	2:B:682:ASN:ND2	2.39	0.56
1:A:155:MET:HE1	1:A:220:LEU:HB2	1.88	0.56
2:B:730:ASN:HB3	2:B:733:GLU:HG3	1.89	0.55
1:A:156:MET:HE1	1:A:271:HIS:HB3	1.88	0.55
1:A:255:LEU:HD12	1:A:258:LEU:HD13	1.88	0.55
2:B:101:LYS:HE3	2:B:125:ARG:HH12	1.72	0.54
1:A:140:SER:OG	1:A:141:GLN:NE2	2.38	0.54
1:A:324:MET:SD	1:A:324:MET:N	2.78	0.53
2:B:142:THR:HG23	2:B:177:GLN:HB2	1.89	0.53
2:B:671:ASP:OD2	2:B:673:ARG:NE	2.36	0.52
2:B:809:MET:HE3	2:B:809:MET:HA	1.92	0.52
1:A:142:GLN:O	1:A:146:GLU:HG2	2.09	0.52
2:B:453:ASN:HB2	2:B:476:TYR:HB2	1.90	0.51
2:B:864:GLN:HA	2:B:867:ILE:HG22	1.91	0.51
2:B:222:SER:OG	2:B:223:ASN:OD1	2.27	0.51
2:B:637:VAL:O	2:B:638:THR:OG1	2.25	0.51
2:B:849:SER:OG	2:B:850:LEU:N	2.43	0.51
1:A:424:VAL:HG13	1:A:425:THR:HG23	1.92	0.51
2:B:56:SER:OG	2:B:57:ARG:N	2.44	0.51
2:B:646:TRP:CH2	2:B:654:ARG:HG3	2.46	0.51
2:B:357:GLU:HG2	2:B:358:LYS:HG3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:VAL:HG21	1:A:450:THR:HG22	1.93	0.50
2:B:283:LYS:HE2	2:B:448:GLU:HG2	1.93	0.50
2:B:306:HIS:HB2	2:B:592:ILE:HG13	1.93	0.50
2:B:718:MET:HE3	2:B:724:LEU:HD11	1.94	0.50
2:B:550:GLU:OE2	2:B:613:THR:OG1	2.30	0.49
1:A:151:ASP:OD1	1:A:152:LEU:N	2.45	0.49
2:B:868:ASN:HA	2:B:871:ARG:HH11	1.77	0.49
2:B:227:ASN:O	2:B:231:SER:OG	2.25	0.49
2:B:42:LYS:O	2:B:46:LYS:HG3	2.14	0.48
1:A:335:LYS:H	1:A:340:PHE:HZ	1.62	0.47
1:A:178:LYS:O	1:A:182:LYS:HG2	2.14	0.47
1:A:321:GLN:HG2	1:A:451:LEU:HD12	1.97	0.47
2:B:218:MET:HE2	2:B:235:ILE:HG13	1.97	0.47
2:B:240:ILE:HA	2:B:243:PHE:CE2	2.50	0.47
1:A:210:ARG:HG2	1:A:293:LEU:HD11	1.97	0.46
2:B:516:GLY:HA3	2:B:725:VAL:HG11	1.96	0.46
2:B:671:ASP:OD1	2:B:672:ASP:N	2.49	0.46
1:A:235:LEU:HD11	1:A:317:VAL:HG21	1.97	0.46
1:A:273:ARG:HH12	1:A:442:LEU:HB2	1.80	0.46
2:B:13:TRP:HB2	2:B:240:ILE:HD11	1.97	0.46
2:B:95:LYS:O	2:B:263:ARG:NH2	2.48	0.46
1:A:232:LEU:O	1:A:314:ARG:NH2	2.49	0.46
2:B:825:GLU:HG2	2:B:826:ASP:H	1.80	0.46
2:B:374:MET:HE2	2:B:374:MET:HB3	1.86	0.45
2:B:392:ARG:O	2:B:561:HIS:NE2	2.48	0.45
1:A:149:ILE:O	1:A:153:ARG:HG2	2.15	0.45
2:B:637:VAL:HG12	2:B:638:THR:H	1.80	0.45
2:B:249:LYS:HE3	2:B:249:LYS:HB3	1.64	0.45
2:B:851:ILE:HB	2:B:859:TRP:HZ3	1.80	0.45
2:B:747:SER:OG	2:B:748:LEU:N	2.49	0.45
2:B:283:LYS:HE3	2:B:449:THR:HG22	1.97	0.45
2:B:692:GLN:HB3	2:B:695:GLU:HB2	1.98	0.45
2:B:20:LEU:HD23	2:B:20:LEU:HA	1.88	0.45
2:B:151:SER:N	2:B:157:GLU:OE2	2.48	0.44
2:B:365:GLU:HG2	2:B:366:PRO:HD2	1.99	0.44
2:B:370:THR:O	2:B:374:MET:HG2	2.18	0.44
2:B:825:GLU:H	2:B:825:GLU:CD	2.26	0.44
2:B:814:ASN:ND2	2:B:830:VAL:O	2.44	0.44
1:A:209:ARG:HA	1:A:209:ARG:HD2	1.71	0.44
2:B:737:ARG:HD3	2:B:737:ARG:C	2.43	0.44
2:B:441:LEU:HD23	2:B:441:LEU:HA	1.77	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:ASP:O	1:A:155:MET:HG3	2.16	0.44
1:A:238:TRP:HH2	1:A:253:HIS:HA	1.83	0.44
2:B:55:VAL:HG22	2:B:116:MET:HE3	1.99	0.43
2:B:194:MET:HA	2:B:194:MET:HE2	2.00	0.43
1:A:283:SER:O	1:A:287:SER:OG	2.31	0.43
2:B:375:LYS:HB2	2:B:375:LYS:HE2	1.79	0.43
1:A:346:LEU:HD11	1:A:348:VAL:HG12	1.99	0.43
2:B:374:MET:HE1	2:B:684:MET:HG2	1.98	0.43
2:B:769:ARG:HD2	2:B:772:LEU:HD12	2.00	0.43
1:A:390:LYS:HA	1:A:390:LYS:HD3	1.84	0.43
2:B:395:THR:N	2:B:398:GLU:OE2	2.44	0.43
2:B:20:LEU:HD22	2:B:24:GLU:HG2	2.01	0.43
2:B:62:LEU:O	2:B:66:VAL:HG23	2.19	0.43
2:B:411:ALA:HB1	2:B:476:TYR:CE1	2.53	0.43
2:B:14:LYS:O	2:B:17:LEU:HG	2.18	0.42
2:B:105:LYS:H	2:B:105:LYS:HG2	1.67	0.42
1:A:258:LEU:HD12	1:A:258:LEU:H	1.85	0.42
2:B:417:ASN:CG	2:B:418:LYS:HG3	2.44	0.42
1:A:196:GLN:NE2	1:A:197:LYS:HG2	2.34	0.42
2:B:358:LYS:HB3	2:B:358:LYS:HE3	1.66	0.42
2:B:418:LYS:HB2	2:B:419:TRP:CE3	2.55	0.42
2:B:825:GLU:OE1	2:B:825:GLU:N	2.53	0.42
1:A:199:LEU:O	1:A:202:THR:OG1	2.34	0.41
2:B:292:TRP:HH2	2:B:592:ILE:HD11	1.85	0.41
2:B:845:GLN:NE2	2:B:851:ILE:O	2.53	0.41
1:A:140:SER:HG	1:A:141:GLN:HE21	1.60	0.41
1:A:345:ARG:HE	1:A:345:ARG:HB3	1.63	0.41
1:A:185:THR:O	1:A:192:GLN:NE2	2.53	0.41
2:B:136:PRO:HA	2:B:137:PRO:HD3	1.95	0.41
1:A:214:LEU:HD11	1:A:293:LEU:HD23	2.03	0.41
2:B:192:GLU:N	2:B:192:GLU:OE1	2.53	0.41
2:B:268:GLU:N	2:B:268:GLU:OE1	2.54	0.41
2:B:84:ARG:HG3	2:B:111:GLU:OE2	2.21	0.41
1:A:225:THR:O	1:A:229:GLU:HG2	2.20	0.41
2:B:61:LYS:HG2	2:B:210:LEU:HD13	2.02	0.41
2:B:467:LYS:NZ	2:B:737:ARG:HH21	2.17	0.41
2:B:524:LYS:HD3	2:B:524:LYS:HA	1.88	0.41
2:B:536:ALA:HB2	2:B:689:LYS:HB2	2.02	0.41
2:B:785:SER:C	2:B:787:TRP:H	2.29	0.41
2:B:859:TRP:CG	2:B:859:TRP:O	2.73	0.41
2:B:38:ARG:HE	2:B:42:LYS:HE2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:361:VAL:HG12	1:A:382:ASN:HB2	2.02	0.40
1:A:153:ARG:HA	1:A:153:ARG:HD2	1.84	0.40
2:B:322:ASN:HB3	2:B:325:VAL:HG12	2.03	0.40
2:B:873:LEU:O	2:B:873:LEU:HD23	2.21	0.40
2:B:410:GLY:O	2:B:412:ILE:HG22	2.21	0.40
1:A:153:ARG:O	1:A:157:GLU:HG2	2.21	0.40
2:B:321:VAL:HG11	2:B:326:ARG:HD3	2.03	0.40
2:B:409:LEU:HD12	2:B:409:LEU:HA	1.91	0.40
2:B:769:ARG:HG3	2:B:772:LEU:HB2	2.04	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	282/851 (33%)	277 (98%)	5 (2%)	0	100	100
2	B	875/900 (97%)	824 (94%)	51 (6%)	0	100	100
All	All	1157/1751 (66%)	1101 (95%)	56 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	254/784 (32%)	253 (100%)	1 (0%)	84	81
2	B	745/780 (96%)	745 (100%)	0	100	100
All	All	999/1564 (64%)	998 (100%)	1 (0%)	87	87

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	228	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	170	GLN
1	A	241	GLN
1	A	303	GLN
2	B	166	ASN
2	B	177	GLN
2	B	293	HIS
2	B	575	ASN
2	B	581	GLN
2	B	845	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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

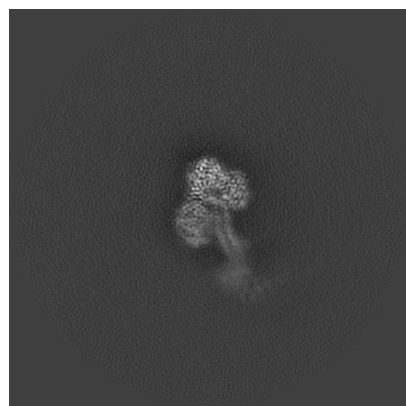
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40953. These allow visual inspection of the internal detail of the map and identification of artifacts.

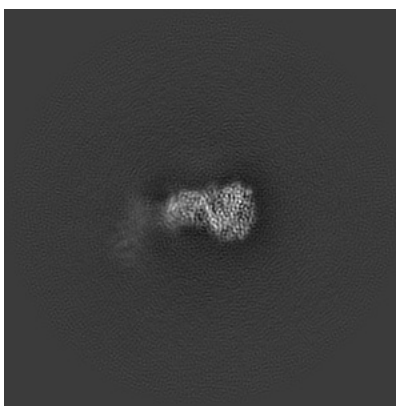
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

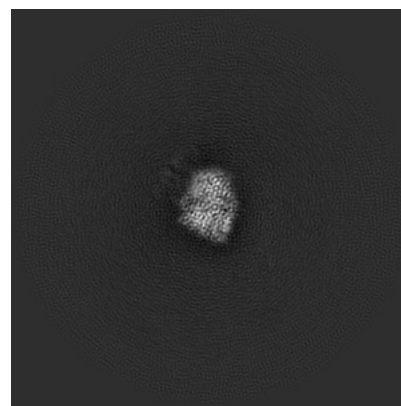
6.1.1 Primary map



X

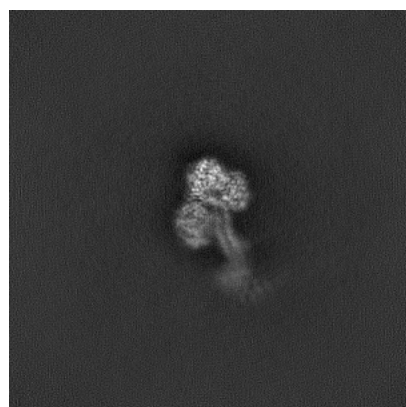


Y

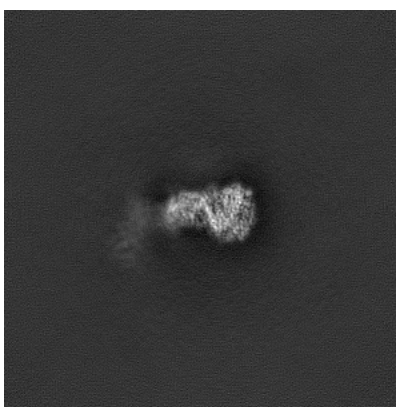


Z

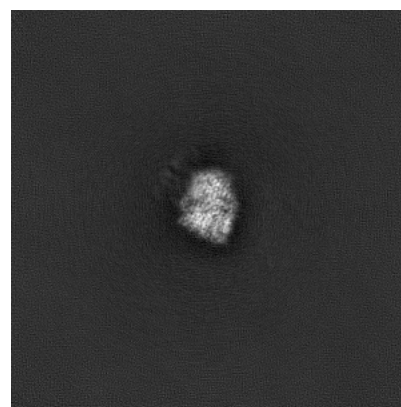
6.1.2 Raw map



X



Y

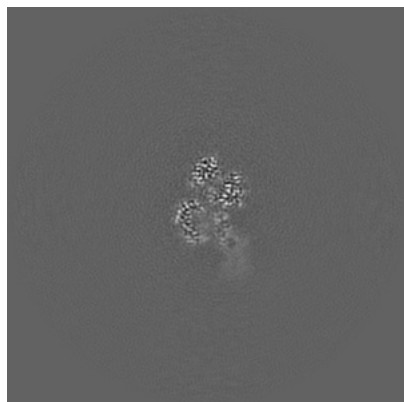


Z

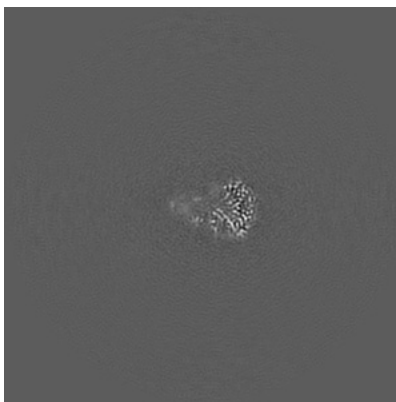
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

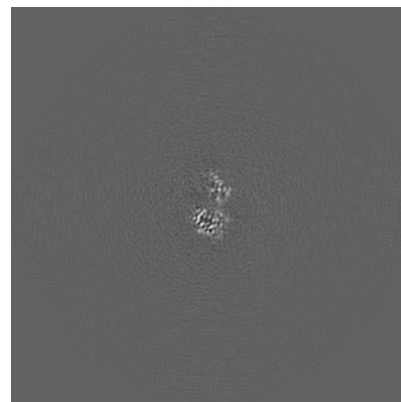
6.2.1 Primary map



X Index: 200

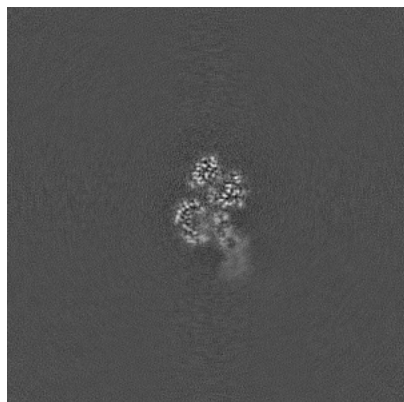


Y Index: 200

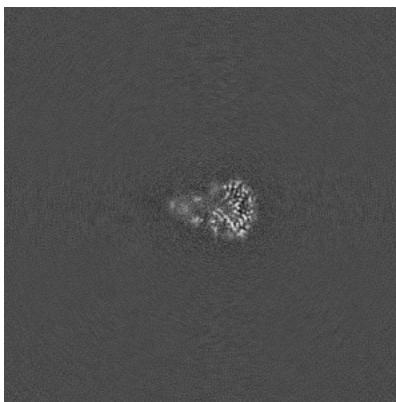


Z Index: 200

6.2.2 Raw map



X Index: 200



Y Index: 200

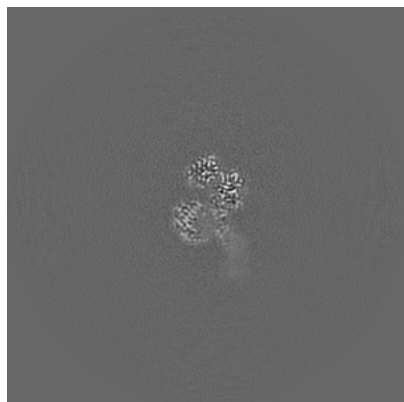


Z Index: 200

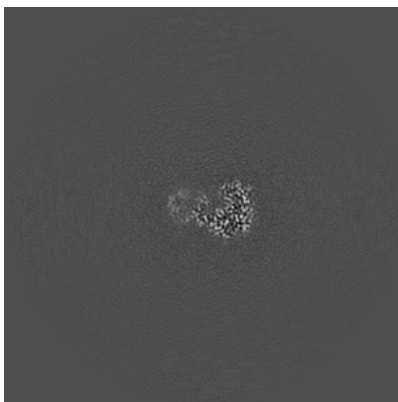
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

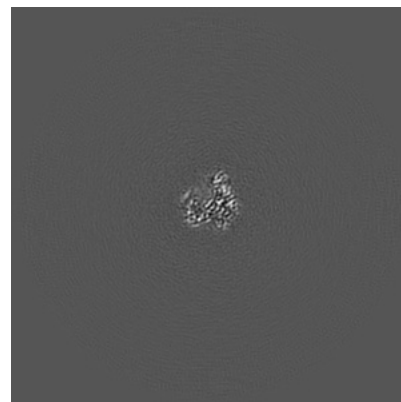
6.3.1 Primary map



X Index: 204

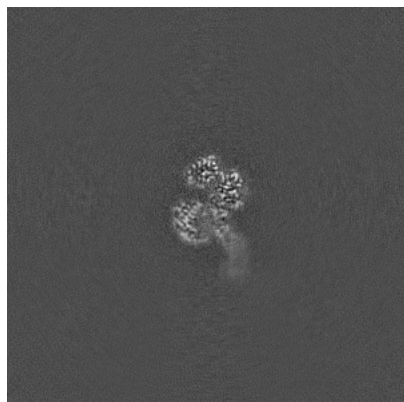


Y Index: 192

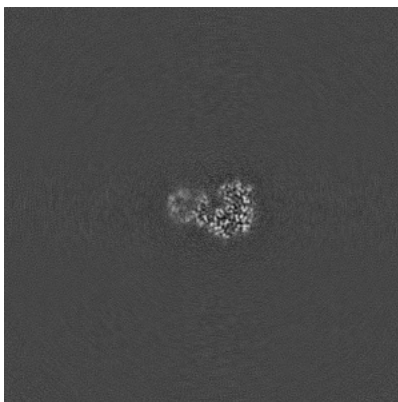


Z Index: 231

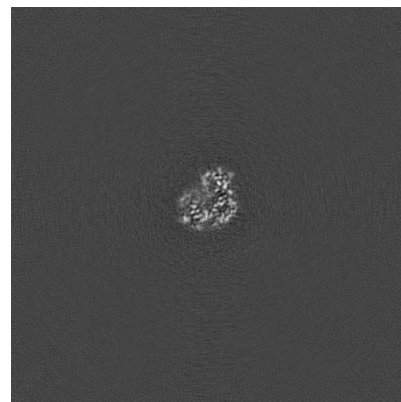
6.3.2 Raw map



X Index: 204



Y Index: 192

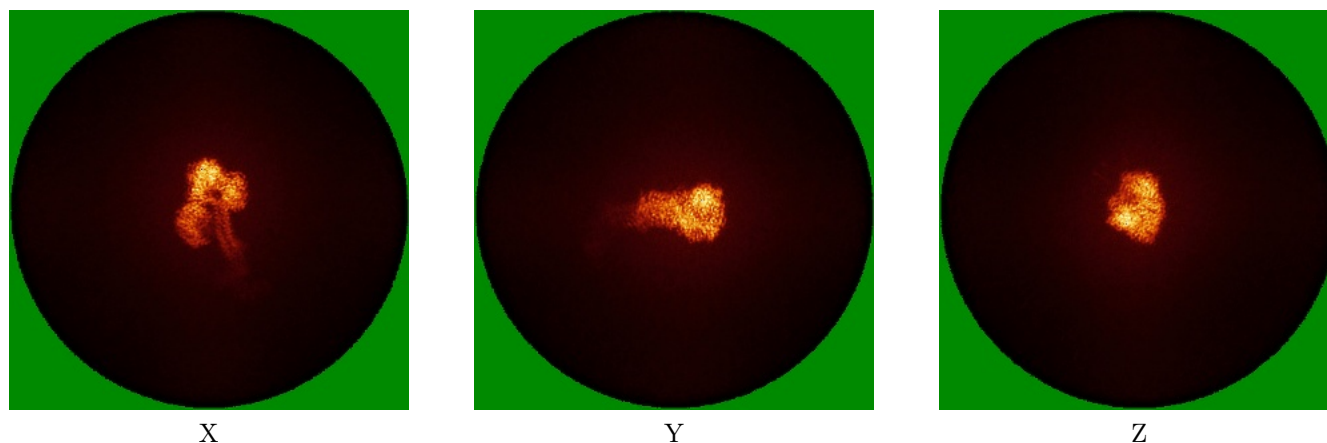


Z Index: 227

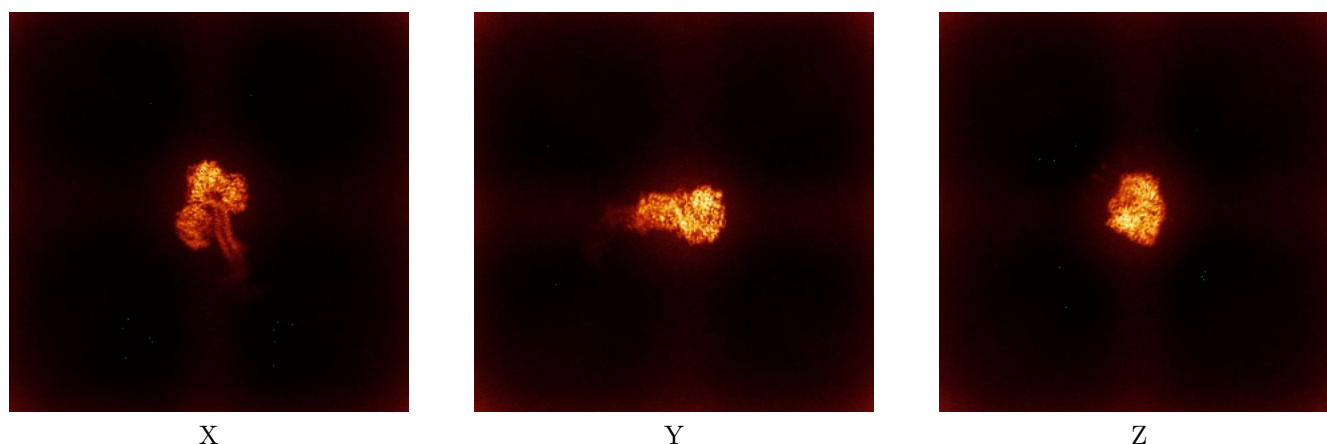
The images above show the largest variance slices of the map in three orthogonal directions.

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.

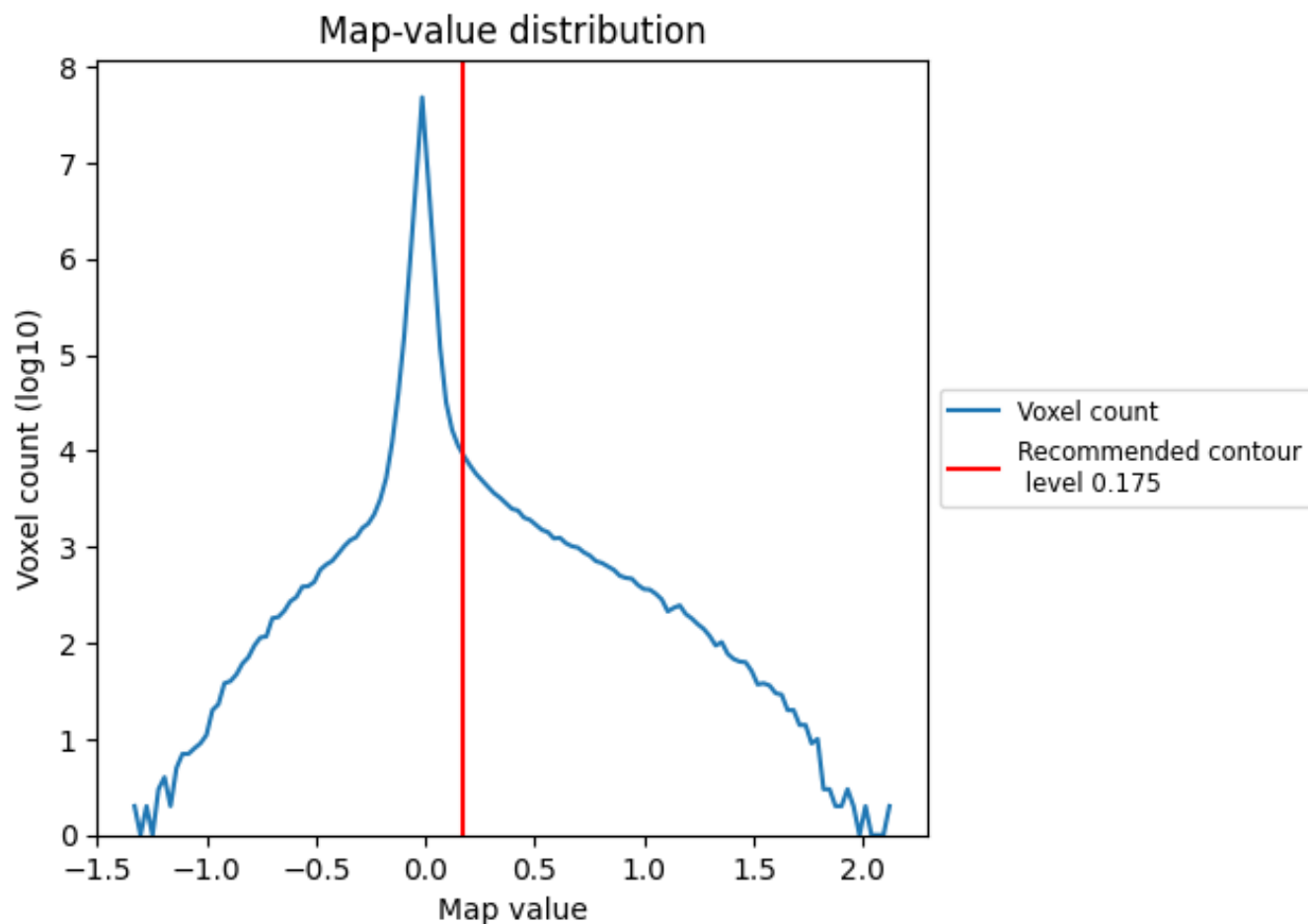
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

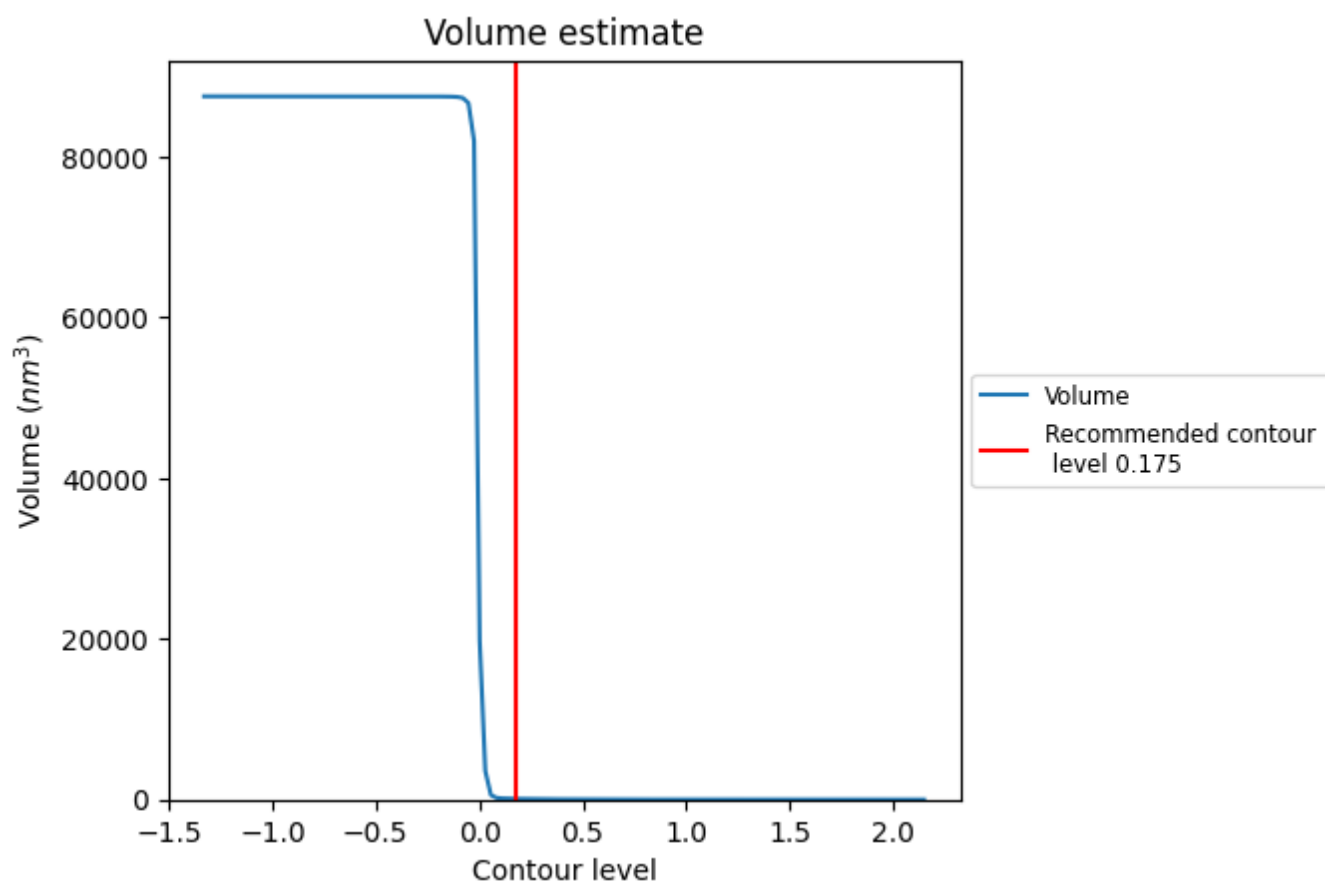
This section contains the results of statistical analysis of the map.

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.

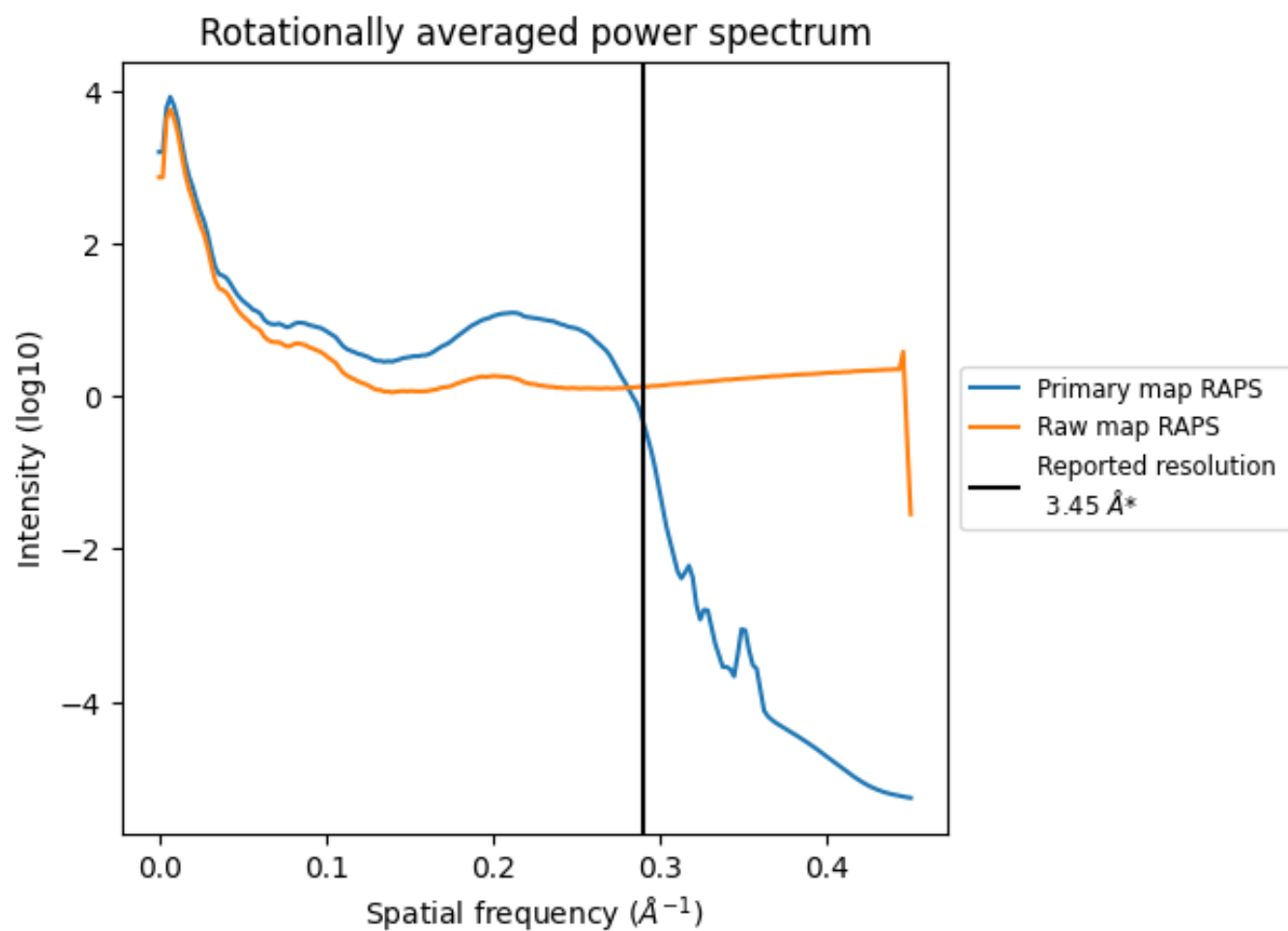
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 96 nm^3 ; this corresponds to an approximate mass of 87 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 ⓘ

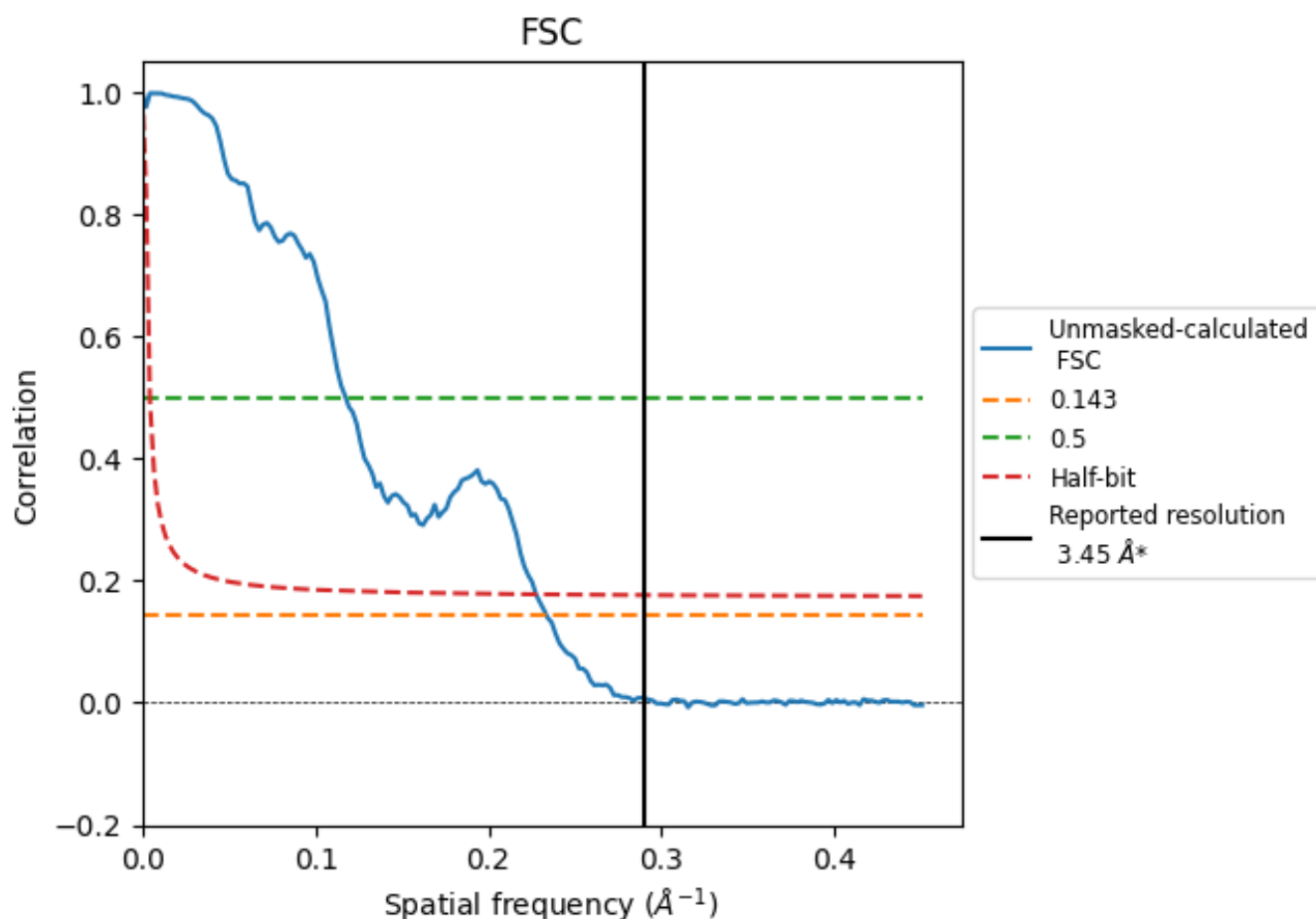


*Reported resolution corresponds to spatial frequency of 0.290 Å⁻¹

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 0.290 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.45	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.28	8.51	4.40

*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 4.28 differs from the reported value 3.45 by more than 10 %

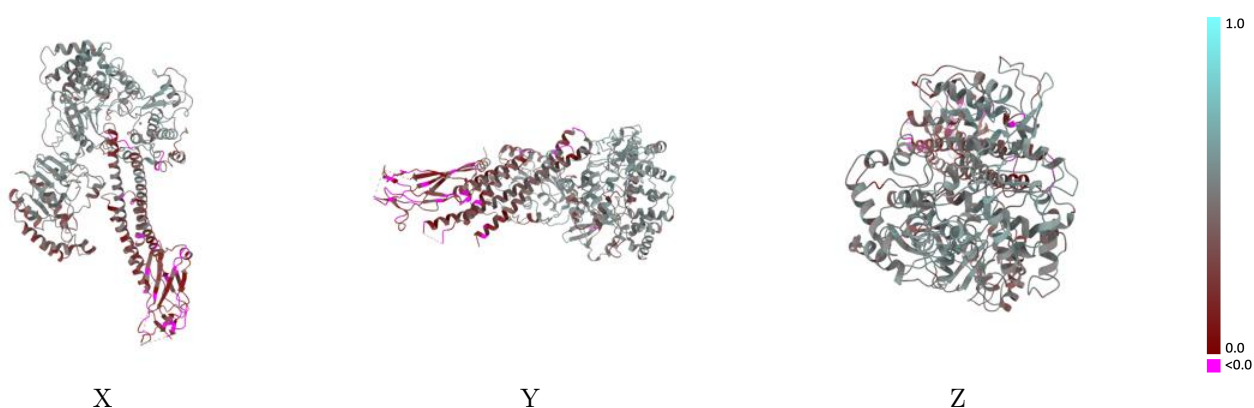
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-40953 and PDB model 8T13. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlay [i](#)

This section was not generated.

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



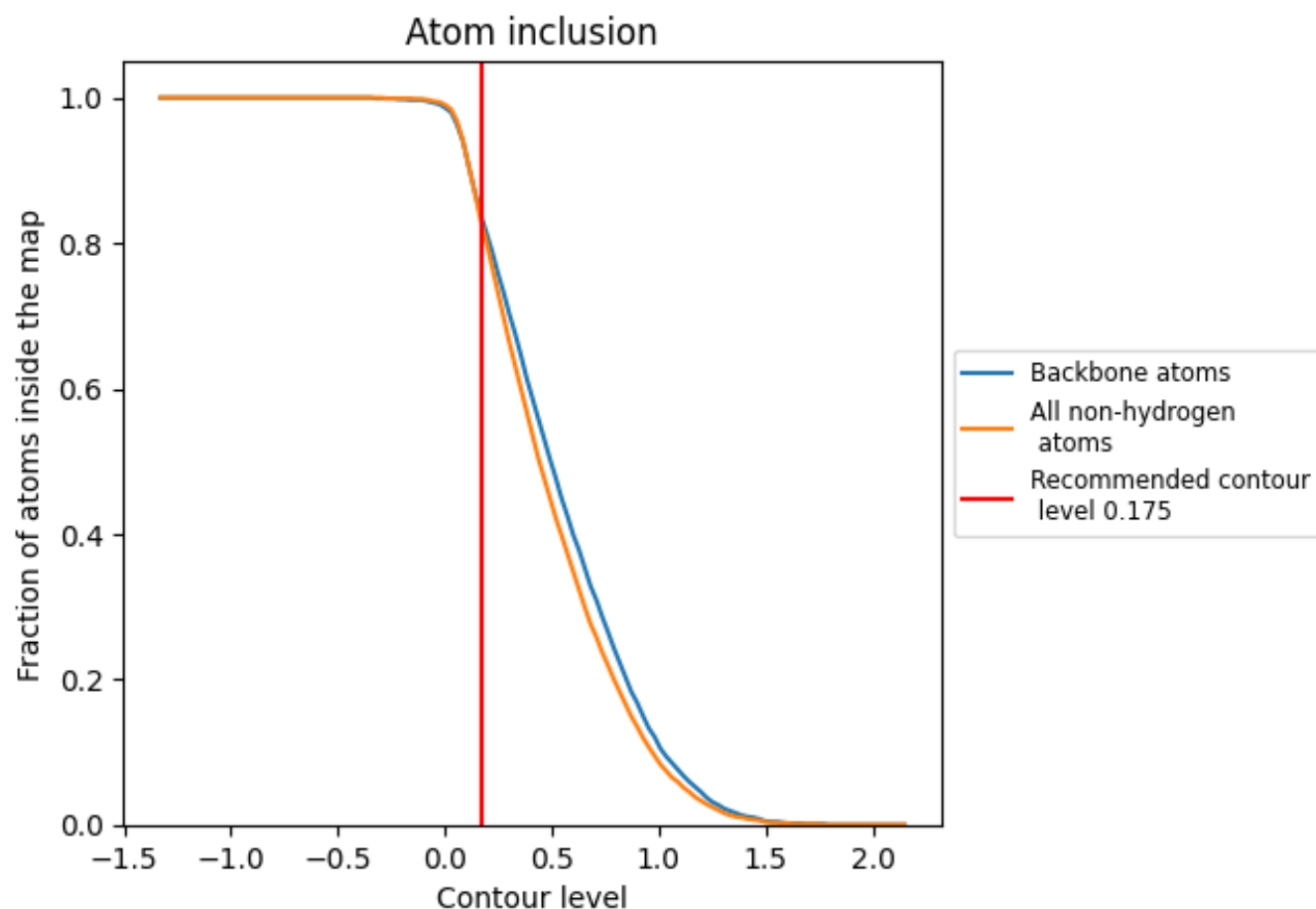
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.175).

9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 82% 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 (0.175) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8220	0.4160
A	0.4990	0.2290
B	0.9290	0.4780

