



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 07:04 AM UTC

PDB ID : 5I08 / pdb_00005i08
EMDB ID : EMD-8069
Title : Prefusion structure of a human coronavirus spike protein
Authors : Kirchdoerfer, R.N.; Cottrell, C.A.; Wang, N.; Pallesen, J.; Yassine, H.M.;
Turner, H.L.; Corbett, K.S.; Graham, B.S.; McLellan, J.S.; Ward, A.B.
Deposited on : 2016-02-03
Resolution : 4.04 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

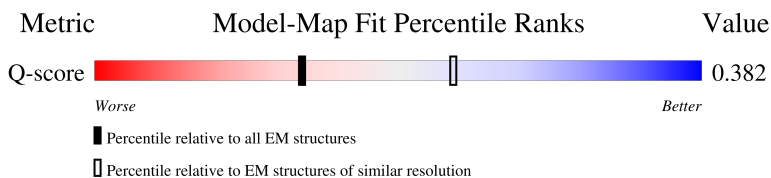
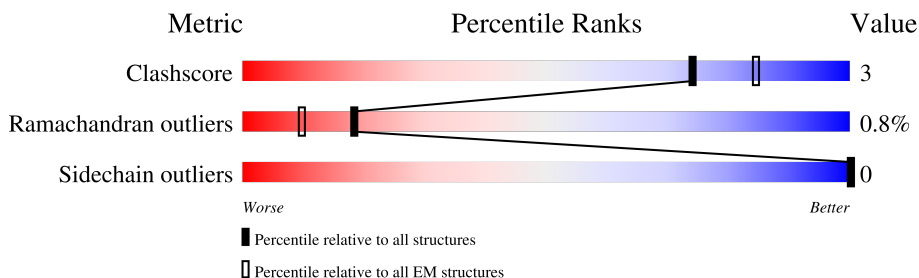
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.04 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	6625 (3.54 - 4.54)

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	1299	
1	B	1299	
1	C	1299	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 22683 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 Spike glycoprotein,Foldon chimera.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	958	Total	C	N	O	S	0	0
			7561	4833	1241	1447	40		
1	B	958	Total	C	N	O	S	0	0
			7561	4833	1241	1447	40		
1	C	958	Total	C	N	O	S	0	0
			7561	4833	1241	1447	40		

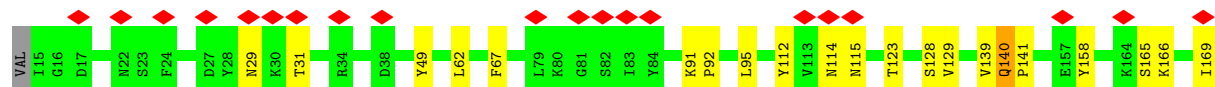
There are 45 discrepancies between the modelled and reference sequences:

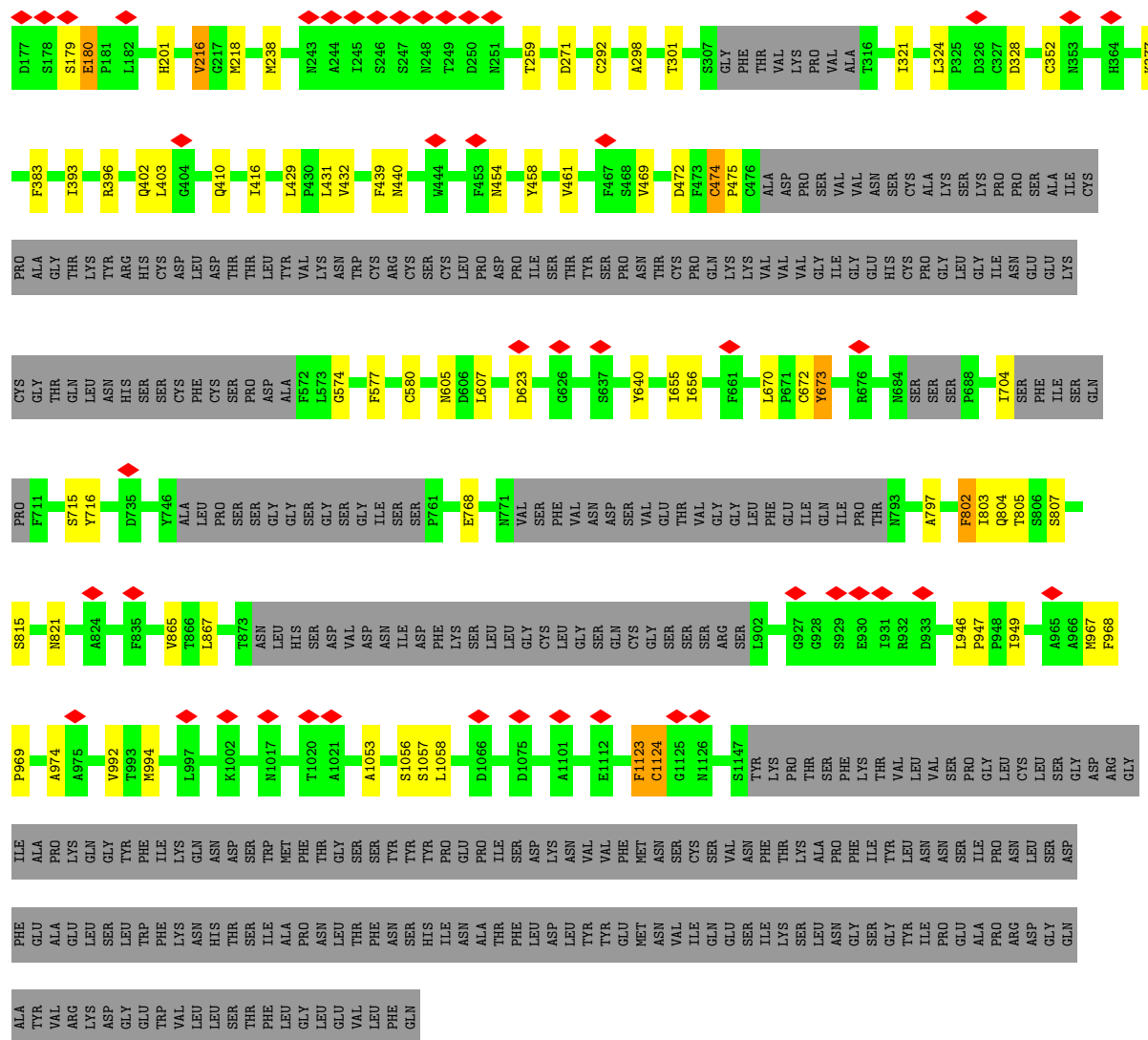
Chain	Residue	Modelled	Actual	Comment	Reference
A	752	GLY	ARG	engineered mutation	UNP Q0ZME7
A	753	GLY	ARG	engineered mutation	UNP Q0ZME7
A	754	SER	LYS	engineered mutation	UNP Q0ZME7
A	755	GLY	ARG	engineered mutation	UNP Q0ZME7
A	756	SER	ARG	engineered mutation	UNP Q0ZME7
A	1277	GLY	-	linker	UNP Q0ZME7
A	1278	SER	-	linker	UNP Q0ZME7
A	1300	LEU	PHE	conflict	UNP P10104
A	1306	GLY	-	expression tag	UNP P10104
A	1307	LEU	-	expression tag	UNP P10104
A	1308	GLU	-	expression tag	UNP P10104
A	1309	VAL	-	expression tag	UNP P10104
A	1310	LEU	-	expression tag	UNP P10104
A	1311	PHE	-	expression tag	UNP P10104
A	1312	GLN	-	expression tag	UNP P10104
B	752	GLY	ARG	engineered mutation	UNP Q0ZME7
B	753	GLY	ARG	engineered mutation	UNP Q0ZME7
B	754	SER	LYS	engineered mutation	UNP Q0ZME7
B	755	GLY	ARG	engineered mutation	UNP Q0ZME7
B	756	SER	ARG	engineered mutation	UNP Q0ZME7
B	1277	GLY	-	linker	UNP Q0ZME7
B	1278	SER	-	linker	UNP Q0ZME7
B	1300	LEU	PHE	conflict	UNP P10104
B	1306	GLY	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1307	LEU	-	expression tag	UNP P10104
B	1308	GLU	-	expression tag	UNP P10104
B	1309	VAL	-	expression tag	UNP P10104
B	1310	LEU	-	expression tag	UNP P10104
B	1311	PHE	-	expression tag	UNP P10104
B	1312	GLN	-	expression tag	UNP P10104
C	752	GLY	ARG	engineered mutation	UNP Q0ZME7
C	753	GLY	ARG	engineered mutation	UNP Q0ZME7
C	754	SER	LYS	engineered mutation	UNP Q0ZME7
C	755	GLY	ARG	engineered mutation	UNP Q0ZME7
C	756	SER	ARG	engineered mutation	UNP Q0ZME7
C	1277	GLY	-	linker	UNP Q0ZME7
C	1278	SER	-	linker	UNP Q0ZME7
C	1300	LEU	PHE	conflict	UNP P10104
C	1306	GLY	-	expression tag	UNP P10104
C	1307	LEU	-	expression tag	UNP P10104
C	1308	GLU	-	expression tag	UNP P10104
C	1309	VAL	-	expression tag	UNP P10104
C	1310	LEU	-	expression tag	UNP P10104
C	1311	PHE	-	expression tag	UNP P10104
C	1312	GLN	-	expression tag	UNP P10104





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	31435	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 ($e^-/\text{\AA}^2$)	57	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	22500	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.150	Depositor
Minimum map value	-0.079	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.0335	Depositor
Map size (Å)	335.36, 335.36, 335.36	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.31, 1.31, 1.31	Depositor

5 Model quality

5.1 Standard geometry

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	0.92	0/7733	1.29	73/10518 (0.7%)
1	B	0.92	0/7733	1.29	73/10518 (0.7%)
1	C	0.92	0/7733	1.29	71/10518 (0.7%)
All	All	0.92	0/23199	1.29	217/31554 (0.7%)

There are no bond length outliers.

All (217) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	946	LEU	CA-C-N	11.19	127.71	119.66
1	A	946	LEU	C-N-CA	11.19	127.71	119.66
1	B	946	LEU	CA-C-N	11.18	127.71	119.66
1	B	946	LEU	C-N-CA	11.18	127.71	119.66
1	C	946	LEU	CA-C-N	11.17	127.70	119.66
1	C	946	LEU	C-N-CA	11.17	127.70	119.66
1	B	947	PRO	CA-C-N	10.50	130.45	119.85
1	B	947	PRO	C-N-CA	10.50	130.45	119.85
1	C	947	PRO	CA-C-N	10.49	130.44	119.85
1	C	947	PRO	C-N-CA	10.49	130.44	119.85
1	A	947	PRO	CA-C-N	10.48	130.44	119.85
1	A	947	PRO	C-N-CA	10.48	130.44	119.85
1	C	454	ASN	N-CA-C	9.52	124.13	111.28
1	A	454	ASN	N-CA-C	9.50	124.11	111.28
1	B	454	ASN	N-CA-C	9.50	124.11	111.28
1	A	968	PHE	CA-C-N	9.09	129.74	120.38
1	A	968	PHE	C-N-CA	9.09	129.74	120.38
1	B	968	PHE	CA-C-N	9.08	129.74	120.38
1	B	968	PHE	C-N-CA	9.08	129.74	120.38
1	C	968	PHE	CA-C-N	9.04	129.69	120.38
1	C	968	PHE	C-N-CA	9.04	129.69	120.38
1	C	768	GLU	CA-C-N	8.84	128.78	119.85
1	C	768	GLU	C-N-CA	8.84	128.78	119.85
1	A	768	GLU	CA-C-N	8.80	128.74	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	768	GLU	C-N-CA	8.80	128.74	119.85
1	B	768	GLU	CA-C-N	8.78	128.72	119.85
1	B	768	GLU	C-N-CA	8.78	128.72	119.85
1	C	802	PHE	N-CA-C	8.50	121.68	111.82
1	A	802	PHE	N-CA-C	8.49	121.67	111.82
1	B	802	PHE	N-CA-C	8.46	121.63	111.82
1	A	324	LEU	CA-C-N	8.07	127.79	119.56
1	A	324	LEU	C-N-CA	8.07	127.79	119.56
1	B	324	LEU	CA-C-N	8.07	127.79	119.56
1	B	324	LEU	C-N-CA	8.07	127.79	119.56
1	C	324	LEU	CA-C-N	8.04	127.77	119.56
1	C	324	LEU	C-N-CA	8.04	127.77	119.56
1	C	969	PRO	N-CA-C	8.00	120.46	110.70
1	B	969	PRO	N-CA-C	7.99	120.44	110.70
1	A	969	PRO	N-CA-C	7.97	120.42	110.70
1	A	321	ILE	CA-C-N	7.83	127.76	119.85
1	A	321	ILE	C-N-CA	7.83	127.76	119.85
1	C	321	ILE	CA-C-N	7.83	127.76	119.85
1	C	321	ILE	C-N-CA	7.83	127.76	119.85
1	B	321	ILE	CA-C-N	7.83	127.75	119.85
1	B	321	ILE	C-N-CA	7.83	127.75	119.85
1	B	474	CYS	CA-C-N	7.78	127.80	119.78
1	B	474	CYS	C-N-CA	7.78	127.80	119.78
1	A	474	CYS	CA-C-N	7.76	127.77	119.78
1	A	474	CYS	C-N-CA	7.76	127.77	119.78
1	C	474	CYS	CA-C-N	7.74	127.75	119.78
1	C	474	CYS	C-N-CA	7.74	127.75	119.78
1	C	259	THR	CA-C-N	7.63	127.56	119.85
1	C	259	THR	C-N-CA	7.63	127.56	119.85
1	B	259	THR	CA-C-N	7.60	127.53	119.85
1	B	259	THR	C-N-CA	7.60	127.53	119.85
1	A	259	THR	CA-C-N	7.59	127.52	119.85
1	A	259	THR	C-N-CA	7.59	127.52	119.85
1	C	292	CYS	CB-CA-C	-7.55	99.02	110.88
1	A	292	CYS	CB-CA-C	-7.55	99.02	110.88
1	B	292	CYS	CB-CA-C	-7.55	99.02	110.88
1	C	91	LYS	CA-C-N	7.24	127.84	120.38
1	C	91	LYS	C-N-CA	7.24	127.84	120.38
1	A	91	LYS	CA-C-N	7.19	127.79	120.38
1	A	91	LYS	C-N-CA	7.19	127.79	120.38
1	B	91	LYS	CA-C-N	7.17	127.76	120.38
1	B	91	LYS	C-N-CA	7.17	127.76	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1057	SER	N-CA-C	6.88	119.64	109.24
1	A	1057	SER	N-CA-C	6.88	119.63	109.24
1	C	1057	SER	N-CA-C	6.88	119.62	109.24
1	B	802	PHE	CA-CB-CG	-6.63	107.17	113.80
1	A	802	PHE	CA-CB-CG	-6.61	107.19	113.80
1	B	238	MET	CA-C-N	6.59	126.51	119.85
1	B	238	MET	C-N-CA	6.59	126.51	119.85
1	A	238	MET	CA-C-N	6.58	126.50	119.85
1	A	238	MET	C-N-CA	6.58	126.50	119.85
1	C	802	PHE	CA-CB-CG	-6.58	107.22	113.80
1	C	238	MET	CA-C-N	6.58	126.49	119.85
1	C	238	MET	C-N-CA	6.58	126.49	119.85
1	A	201	HIS	CA-CB-CG	6.49	120.29	113.80
1	B	201	HIS	CA-CB-CG	6.49	120.28	113.80
1	C	393	ILE	N-CA-C	6.48	115.03	107.77
1	C	201	HIS	CA-CB-CG	6.47	120.27	113.80
1	A	393	ILE	N-CA-C	6.46	115.01	107.77
1	B	393	ILE	N-CA-C	6.45	114.99	107.77
1	B	298	ALA	CA-C-N	6.42	126.39	119.78
1	B	298	ALA	C-N-CA	6.42	126.39	119.78
1	C	298	ALA	CA-C-N	6.42	126.39	119.78
1	C	298	ALA	C-N-CA	6.42	126.39	119.78
1	A	298	ALA	CA-C-N	6.41	126.39	119.78
1	A	298	ALA	C-N-CA	6.41	126.39	119.78
1	C	1123	PHE	N-CA-C	-6.34	100.89	110.28
1	A	1123	PHE	N-CA-C	-6.34	100.90	110.28
1	B	1123	PHE	N-CA-C	-6.32	100.93	110.28
1	C	440	ASN	CA-C-N	6.30	126.21	119.85
1	C	440	ASN	C-N-CA	6.30	126.21	119.85
1	C	429	LEU	CA-C-N	6.29	126.26	119.78
1	C	429	LEU	C-N-CA	6.29	126.26	119.78
1	A	440	ASN	CA-C-N	6.28	126.19	119.85
1	A	440	ASN	C-N-CA	6.28	126.19	119.85
1	B	440	ASN	CA-C-N	6.27	126.18	119.85
1	B	440	ASN	C-N-CA	6.27	126.18	119.85
1	A	429	LEU	CA-C-N	6.27	126.24	119.78
1	A	429	LEU	C-N-CA	6.27	126.24	119.78
1	B	140	GLN	CA-C-N	6.26	126.18	119.85
1	B	140	GLN	C-N-CA	6.26	126.18	119.85
1	B	429	LEU	CA-C-N	6.26	126.23	119.78
1	B	429	LEU	C-N-CA	6.26	126.23	119.78
1	A	49	TYR	N-CA-CB	-6.25	100.78	110.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	GLN	CA-C-N	6.25	126.16	119.85
1	A	140	GLN	C-N-CA	6.25	126.16	119.85
1	B	49	TYR	N-CA-CB	-6.24	100.78	110.65
1	C	49	TYR	N-CA-CB	-6.24	100.79	110.65
1	C	140	GLN	CA-C-N	6.22	126.13	119.85
1	C	140	GLN	C-N-CA	6.22	126.13	119.85
1	B	656	ILE	N-CA-C	6.07	117.03	111.81
1	A	1124	CYS	N-CA-C	-6.07	103.09	111.28
1	C	656	ILE	N-CA-C	6.07	117.03	111.81
1	C	1124	CYS	N-CA-C	-6.07	103.09	111.28
1	A	301	THR	CB-CA-C	-6.06	109.57	116.54
1	A	656	ILE	N-CA-C	6.06	117.02	111.81
1	B	1124	CYS	N-CA-C	-6.05	103.11	111.28
1	B	301	THR	CB-CA-C	-6.03	109.61	116.54
1	C	301	THR	CB-CA-C	-6.03	109.61	116.54
1	B	1058	LEU	N-CA-C	6.00	117.43	108.31
1	A	1058	LEU	N-CA-C	5.98	117.39	108.31
1	B	393	ILE	CA-C-N	5.96	125.92	119.78
1	B	393	ILE	C-N-CA	5.96	125.92	119.78
1	C	1058	LEU	N-CA-C	5.96	117.38	108.31
1	A	393	ILE	CA-C-N	5.91	125.86	119.78
1	A	393	ILE	C-N-CA	5.91	125.86	119.78
1	C	393	ILE	CA-C-N	5.90	125.86	119.78
1	C	393	ILE	C-N-CA	5.90	125.86	119.78
1	B	180	GLU	CA-C-N	5.80	125.76	119.78
1	B	180	GLU	C-N-CA	5.80	125.76	119.78
1	A	180	GLU	CA-C-N	5.79	125.74	119.78
1	A	180	GLU	C-N-CA	5.79	125.74	119.78
1	C	180	GLU	CA-C-N	5.77	125.73	119.78
1	C	180	GLU	C-N-CA	5.77	125.73	119.78
1	B	29	ASN	N-CA-C	5.77	116.70	108.74
1	A	29	ASN	N-CA-C	5.75	116.67	108.74
1	C	29	ASN	N-CA-C	5.73	116.65	108.74
1	C	67	PHE	CA-C-N	5.70	125.61	119.85
1	C	67	PHE	C-N-CA	5.70	125.61	119.85
1	B	67	PHE	CA-C-N	5.69	125.59	119.85
1	B	67	PHE	C-N-CA	5.69	125.59	119.85
1	A	67	PHE	CA-C-N	5.68	125.59	119.85
1	A	67	PHE	C-N-CA	5.68	125.59	119.85
1	B	92	PRO	N-CA-C	5.59	117.52	110.70
1	C	865	VAL	N-CA-C	5.58	115.78	110.42
1	A	865	VAL	N-CA-C	5.58	115.77	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	865	VAL	N-CA-C	5.58	115.77	110.42
1	C	383	PHE	CB-CA-C	-5.57	108.54	115.89
1	A	383	PHE	CB-CA-C	-5.57	108.54	115.89
1	B	383	PHE	CB-CA-C	-5.57	108.54	115.89
1	A	92	PRO	N-CA-C	5.56	117.48	110.70
1	C	92	PRO	N-CA-C	5.53	117.45	110.70
1	C	352	CYS	N-CA-C	5.52	117.19	108.96
1	A	352	CYS	N-CA-C	5.51	117.16	108.96
1	C	158	TYR	CA-C-N	5.51	125.41	119.85
1	C	158	TYR	C-N-CA	5.51	125.41	119.85
1	C	807	SER	CA-C-N	5.50	125.44	119.78
1	C	807	SER	C-N-CA	5.50	125.44	119.78
1	B	352	CYS	N-CA-C	5.49	117.14	108.96
1	A	807	SER	CA-C-N	5.47	125.41	119.78
1	A	807	SER	C-N-CA	5.47	125.41	119.78
1	B	807	SER	CA-C-N	5.46	125.40	119.78
1	B	807	SER	C-N-CA	5.46	125.40	119.78
1	B	1053	ALA	CA-C-N	5.46	127.94	120.35
1	B	1053	ALA	C-N-CA	5.46	127.94	120.35
1	A	158	TYR	CA-C-N	5.46	125.36	119.85
1	A	158	TYR	C-N-CA	5.46	125.36	119.85
1	A	1053	ALA	CA-C-N	5.46	127.94	120.35
1	A	1053	ALA	C-N-CA	5.46	127.94	120.35
1	B	158	TYR	CA-C-N	5.45	125.36	119.85
1	B	158	TYR	C-N-CA	5.45	125.36	119.85
1	C	1053	ALA	CA-C-N	5.45	127.92	120.35
1	C	1053	ALA	C-N-CA	5.45	127.92	120.35
1	A	1056	SER	CA-C-N	-5.42	113.43	122.21
1	A	1056	SER	C-N-CA	-5.42	113.43	122.21
1	B	1056	SER	CA-C-N	-5.42	113.43	122.21
1	B	1056	SER	C-N-CA	-5.42	113.43	122.21
1	C	1056	SER	CA-C-N	-5.41	113.45	122.21
1	C	1056	SER	C-N-CA	-5.41	113.45	122.21
1	B	396	ARG	N-CA-C	5.35	117.11	111.28
1	A	396	ARG	N-CA-C	5.34	117.10	111.28
1	C	396	ARG	N-CA-C	5.32	117.08	111.28
1	C	949	ILE	CB-CA-C	-5.29	105.00	112.14
1	A	949	ILE	CB-CA-C	-5.28	105.01	112.14
1	B	949	ILE	CB-CA-C	-5.27	105.02	112.14
1	C	670	LEU	CA-C-N	5.24	125.14	119.85
1	C	670	LEU	C-N-CA	5.24	125.14	119.85
1	A	670	LEU	CA-C-N	5.22	125.12	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	670	LEU	C-N-CA	5.22	125.12	119.85
1	B	670	LEU	CA-C-N	5.22	125.12	119.85
1	B	670	LEU	C-N-CA	5.22	125.12	119.85
1	A	803	ILE	N-CA-C	5.21	120.17	109.34
1	B	803	ILE	N-CA-C	5.20	120.16	109.34
1	C	803	ILE	N-CA-C	5.20	120.16	109.34
1	C	574	GLY	N-CA-C	5.14	122.44	115.30
1	B	574	GLY	N-CA-C	5.13	122.43	115.30
1	A	574	GLY	N-CA-C	5.13	122.43	115.30
1	B	797	ALA	CA-C-N	5.06	125.14	120.34
1	B	797	ALA	C-N-CA	5.06	125.14	120.34
1	B	95	LEU	CA-C-N	-5.03	113.05	121.75
1	B	95	LEU	C-N-CA	-5.03	113.05	121.75
1	A	95	LEU	CA-C-N	-5.03	113.05	121.75
1	A	95	LEU	C-N-CA	-5.03	113.05	121.75
1	C	95	LEU	CA-C-N	-5.03	113.05	121.75
1	C	95	LEU	C-N-CA	-5.03	113.05	121.75
1	A	797	ALA	CA-C-N	5.03	125.11	120.34
1	A	797	ALA	C-N-CA	5.03	125.11	120.34
1	C	797	ALA	CA-C-N	5.02	125.11	120.34
1	C	797	ALA	C-N-CA	5.02	125.11	120.34
1	B	61	LEU	CA-C-N	-5.02	114.60	122.73
1	B	61	LEU	C-N-CA	-5.02	114.60	122.73
1	A	61	LEU	CA-C-N	-5.01	114.62	122.73
1	A	61	LEU	C-N-CA	-5.01	114.62	122.73

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	7561	0	7292	52	0
1	B	7561	0	7292	45	0
1	C	7561	0	7292	46	0
All	All	22683	0	21876	143	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 3.

All (143) 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:B:802:PHE:O	1:B:802:PHE:CD1	2.17	0.97
1:A:802:PHE:CD1	1:A:802:PHE:O	2.17	0.97
1:C:802:PHE:CD1	1:C:802:PHE:O	2.18	0.96
1:B:804:GLN:HG2	1:B:804:GLN:O	1.80	0.81
1:A:804:GLN:HG2	1:A:804:GLN:O	1.80	0.80
1:C:804:GLN:O	1:C:804:GLN:HG2	1.80	0.79
1:B:994:MET:O	1:B:994:MET:HG2	1.88	0.73
1:A:994:MET:O	1:A:994:MET:HG2	1.88	0.72
1:C:994:MET:O	1:C:994:MET:HG2	1.88	0.71
1:B:607:LEU:HG	1:B:607:LEU:O	1.91	0.70
1:C:607:LEU:O	1:C:607:LEU:HG	1.91	0.70
1:A:607:LEU:HG	1:A:607:LEU:O	1.91	0.69
1:A:216:VAL:HG12	1:A:216:VAL:O	1.92	0.68
1:B:216:VAL:HG12	1:B:216:VAL:O	1.92	0.68
1:C:216:VAL:O	1:C:216:VAL:HG12	1.92	0.67
1:C:458:TYR:CD1	1:C:458:TYR:N	2.64	0.65
1:B:458:TYR:CD1	1:B:458:TYR:N	2.64	0.65
1:A:458:TYR:N	1:A:458:TYR:CD1	2.64	0.64
1:A:580:CYS:SG	1:A:580:CYS:O	2.56	0.64
1:A:62:LEU:C	1:A:62:LEU:HD12	2.24	0.63
1:A:802:PHE:CD1	1:A:802:PHE:C	2.76	0.62
1:C:580:CYS:SG	1:C:580:CYS:O	2.56	0.62
1:C:62:LEU:C	1:C:62:LEU:HD12	2.24	0.62
1:B:580:CYS:SG	1:B:580:CYS:O	2.56	0.62
1:B:62:LEU:C	1:B:62:LEU:HD12	2.24	0.61
1:C:802:PHE:CD1	1:C:802:PHE:C	2.76	0.58
1:B:802:PHE:O	1:B:802:PHE:CG	2.57	0.58
1:A:802:PHE:O	1:A:802:PHE:CG	2.57	0.58
1:C:802:PHE:O	1:C:802:PHE:CG	2.57	0.57
1:A:410:GLN:HA	1:A:410:GLN:OE1	2.07	0.55
1:A:715:SER:OG	1:A:716:TYR:N	2.40	0.55
1:B:62:LEU:HD12	1:B:62:LEU:O	2.07	0.55
1:C:715:SER:OG	1:C:716:TYR:N	2.40	0.55
1:A:62:LEU:HD12	1:A:62:LEU:O	2.07	0.55
1:C:62:LEU:HD12	1:C:62:LEU:O	2.07	0.55
1:B:410:GLN:HA	1:B:410:GLN:OE1	2.07	0.54
1:C:410:GLN:HA	1:C:410:GLN:OE1	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:867:LEU:HD12	1:A:867:LEU:C	2.33	0.54
1:B:867:LEU:HD12	1:B:867:LEU:C	2.33	0.54
1:C:867:LEU:C	1:C:867:LEU:HD12	2.33	0.53
1:C:994:MET:O	1:C:994:MET:CG	2.57	0.53
1:B:605:ASN:OD1	1:B:605:ASN:C	2.52	0.53
1:C:605:ASN:C	1:C:605:ASN:OD1	2.52	0.53
1:B:715:SER:OG	1:B:716:TYR:N	2.40	0.52
1:C:112:TYR:CD2	1:C:112:TYR:O	2.64	0.51
1:A:605:ASN:OD1	1:A:605:ASN:C	2.52	0.51
1:A:123:THR:OG1	1:A:139:VAL:O	2.28	0.51
1:B:112:TYR:O	1:B:112:TYR:CD2	2.64	0.51
1:B:31:THR:HG22	1:B:31:THR:O	2.10	0.50
1:A:112:TYR:O	1:A:112:TYR:CD2	2.64	0.50
1:B:994:MET:O	1:B:994:MET:CG	2.57	0.50
1:A:179:SER:OG	1:A:180:GLU:N	2.45	0.50
1:C:867:LEU:HD12	1:C:867:LEU:O	2.12	0.50
1:A:31:THR:HG22	1:A:31:THR:O	2.10	0.50
1:B:271:ASP:OD1	1:B:271:ASP:C	2.55	0.49
1:C:123:THR:OG1	1:C:139:VAL:O	2.28	0.49
1:A:867:LEU:HD12	1:A:867:LEU:O	2.12	0.49
1:B:867:LEU:HD12	1:B:867:LEU:O	2.12	0.49
1:C:31:THR:O	1:C:31:THR:HG22	2.10	0.49
1:A:128:SER:OG	1:A:129:VAL:N	2.45	0.49
1:A:994:MET:O	1:A:994:MET:CG	2.57	0.49
1:C:128:SER:OG	1:C:129:VAL:N	2.45	0.48
1:C:655:ILE:O	1:C:655:ILE:HG23	2.13	0.48
1:B:655:ILE:O	1:B:655:ILE:HG23	2.13	0.48
1:B:577:PHE:CD1	1:B:577:PHE:C	2.92	0.48
1:C:179:SER:OG	1:C:180:GLU:N	2.45	0.48
1:A:655:ILE:HG23	1:A:655:ILE:O	2.13	0.47
1:B:123:THR:OG1	1:B:139:VAL:O	2.28	0.47
1:C:271:ASP:C	1:C:271:ASP:OD1	2.55	0.47
1:A:704:ILE:HG22	1:A:704:ILE:O	2.15	0.47
1:B:704:ILE:O	1:B:704:ILE:HG22	2.15	0.47
1:C:577:PHE:CD1	1:C:577:PHE:C	2.92	0.46
1:A:577:PHE:CD1	1:A:577:PHE:C	2.92	0.46
1:C:328:ASP:OD1	1:C:328:ASP:C	2.58	0.46
1:A:271:ASP:OD1	1:A:271:ASP:C	2.55	0.46
1:C:704:ILE:O	1:C:704:ILE:HG22	2.15	0.46
1:A:112:TYR:O	1:A:112:TYR:HD2	1.99	0.46
1:B:112:TYR:O	1:B:112:TYR:HD2	1.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:169:ILE:HG13	1:C:169:ILE:O	2.15	0.46
1:A:416:ILE:O	1:A:416:ILE:HG23	2.16	0.46
1:B:169:ILE:HG13	1:B:169:ILE:O	2.15	0.46
1:C:140:GLN:HA	1:C:141:PRO:HD2	1.84	0.46
1:C:805:THR:O	1:C:805:THR:HG22	2.16	0.45
1:B:128:SER:OG	1:B:129:VAL:N	2.45	0.45
1:C:416:ILE:O	1:C:416:ILE:HG23	2.16	0.45
1:A:169:ILE:O	1:A:169:ILE:HG13	2.15	0.45
1:A:805:THR:O	1:A:805:THR:HG22	2.16	0.45
1:B:328:ASP:OD1	1:B:328:ASP:C	2.58	0.45
1:C:112:TYR:O	1:C:112:TYR:HD2	1.99	0.45
1:B:416:ILE:O	1:B:416:ILE:HG23	2.16	0.45
1:A:431:LEU:O	1:A:432:VAL:HB	2.17	0.44
1:C:431:LEU:O	1:C:432:VAL:HB	2.17	0.44
1:B:431:LEU:O	1:B:432:VAL:HB	2.17	0.44
1:B:805:THR:O	1:B:805:THR:HG22	2.16	0.44
1:A:328:ASP:OD1	1:A:328:ASP:C	2.58	0.44
1:B:992:VAL:O	1:B:992:VAL:HG13	2.18	0.44
1:C:992:VAL:HG13	1:C:992:VAL:O	2.18	0.44
1:A:992:VAL:O	1:A:992:VAL:HG13	2.18	0.43
1:B:821:ASN:OD1	1:B:821:ASN:C	2.62	0.43
1:A:377:LYS:HA	1:A:377:LYS:HD3	1.81	0.43
1:A:821:ASN:OD1	1:A:821:ASN:C	2.62	0.43
1:B:114:ASN:O	1:B:115:ASN:HB2	2.19	0.42
1:C:114:ASN:O	1:C:115:ASN:HB2	2.19	0.42
1:A:623:ASP:OD1	1:A:623:ASP:C	2.61	0.42
1:A:114:ASN:O	1:A:115:ASN:HB2	2.19	0.42
1:A:672:CYS:O	1:A:673:TYR:HB2	2.19	0.42
1:C:672:CYS:O	1:C:673:TYR:HB2	2.19	0.42
1:B:377:LYS:HA	1:B:377:LYS:HD3	1.81	0.42
1:C:165:SER:OG	1:C:166:LYS:N	2.52	0.42
1:B:402:GLN:O	1:B:403:LEU:HB2	2.20	0.42
1:B:672:CYS:O	1:B:673:TYR:HB2	2.20	0.42
1:B:623:ASP:OD1	1:B:623:ASP:C	2.61	0.42
1:A:165:SER:OG	1:A:166:LYS:N	2.52	0.41
1:A:996:VAL:HA	1:A:999:LYS:HB2	2.02	0.41
1:B:41:ASP:OD1	1:B:41:ASP:C	2.62	0.41
1:C:472:ASP:OD1	1:C:472:ASP:C	2.63	0.41
1:A:402:GLN:O	1:A:403:LEU:HB2	2.20	0.41
1:B:996:VAL:HA	1:B:999:LYS:HB2	2.02	0.41
1:C:402:GLN:O	1:C:403:LEU:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:821:ASN:OD1	1:C:821:ASN:C	2.62	0.41
1:A:41:ASP:OD1	1:A:41:ASP:C	2.62	0.41
1:A:469:VAL:O	1:A:469:VAL:HG12	2.21	0.41
1:B:429:LEU:HA	1:B:430:PRO:HD3	1.89	0.41
1:B:967:MET:O	1:B:967:MET:HG3	2.20	0.41
1:A:1054:ILE:O	1:A:1054:ILE:HG13	2.21	0.41
1:C:469:VAL:O	1:C:469:VAL:HG12	2.21	0.41
1:C:623:ASP:C	1:C:623:ASP:OD1	2.61	0.41
1:B:469:VAL:O	1:B:469:VAL:HG12	2.21	0.41
1:C:967:MET:O	1:C:967:MET:HG3	2.20	0.41
1:B:290:ILE:HA	1:B:290:ILE:HD12	1.86	0.41
1:C:377:LYS:HA	1:C:377:LYS:HD3	1.81	0.41
1:A:967:MET:O	1:A:967:MET:HG3	2.20	0.40
1:A:1123:PHE:O	1:A:1124:CYS:HB3	2.21	0.40
1:B:576:SER:OG	1:B:577:PHE:N	2.54	0.40
1:C:1123:PHE:O	1:C:1124:CYS:HB3	2.21	0.40
1:A:140:GLN:HA	1:A:141:PRO:HD2	1.84	0.40
1:A:277:THR:OG1	1:A:278:ASN:N	2.54	0.40
1:A:576:SER:OG	1:A:577:PHE:N	2.54	0.40
1:B:43:SER:HA	1:B:279:ALA:O	2.21	0.40
1:A:43:SER:HA	1:A:279:ALA:O	2.21	0.40
1:A:110:LYS:NZ	1:A:196:ASP:OD1	2.55	0.40
1:A:474:CYS:HA	1:A:475:PRO:HD3	1.88	0.40
1:C:474:CYS:HA	1:C:475:PRO:HD3	1.88	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	942/1299 (72%)	868 (92%)	66 (7%)	8 (1%)	16 51

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	942/1299 (72%)	868 (92%)	66 (7%)	8 (1%)	16	51
1	C	942/1299 (72%)	868 (92%)	66 (7%)	8 (1%)	16	51
All	All	2826/3897 (72%)	2604 (92%)	198 (7%)	24 (1%)	18	51

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	218	MET
1	A	673	TYR
1	B	218	MET
1	B	673	TYR
1	C	218	MET
1	C	673	TYR
1	A	461	VAL
1	A	640	TYR
1	B	461	VAL
1	B	640	TYR
1	C	461	VAL
1	C	640	TYR
1	A	439	PHE
1	A	815	SER
1	A	974	ALA
1	B	439	PHE
1	B	815	SER
1	B	974	ALA
1	C	439	PHE
1	C	815	SER
1	C	974	ALA
1	A	216	VAL
1	B	216	VAL
1	C	216	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	860/1161 (74%)	860 (100%)	0	100	100
1	B	860/1161 (74%)	860 (100%)	0	100	100
1	C	860/1161 (74%)	860 (100%)	0	100	100
All	All	2580/3483 (74%)	2580 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	HIS
1	A	188	ASN
1	A	269	ASN
1	A	424	GLN
1	A	465	HIS
1	A	591	ASN
1	A	595	ASN
1	A	621	ASN
1	A	694	ASN
1	B	175	HIS
1	B	188	ASN
1	B	438	ASN
1	B	465	HIS
1	B	595	ASN
1	B	621	ASN
1	B	694	ASN
1	C	175	HIS
1	C	188	ASN
1	C	269	ASN
1	C	278	ASN
1	C	595	ASN
1	C	621	ASN
1	C	694	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

There are no chain breaks in this entry.

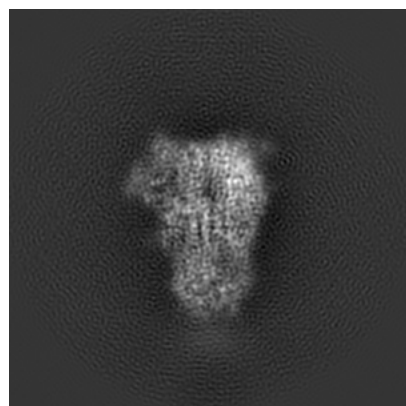
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8069. These allow visual inspection of the internal detail of the map and identification of artifacts.

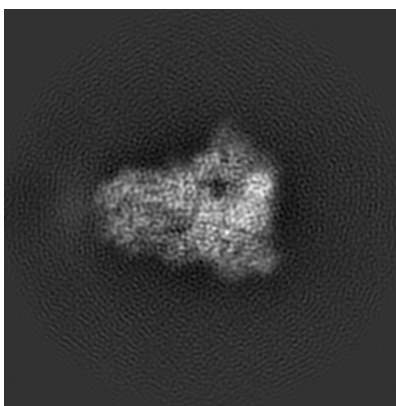
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

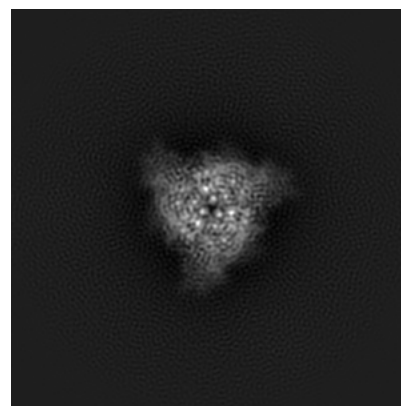
6.1.1 Primary map



X

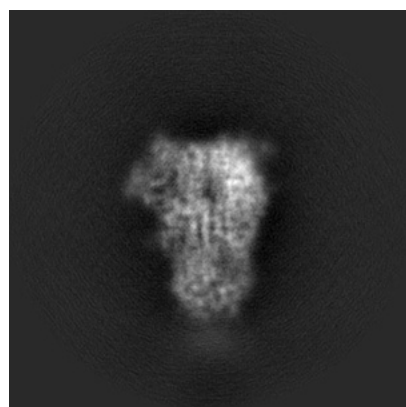


Y

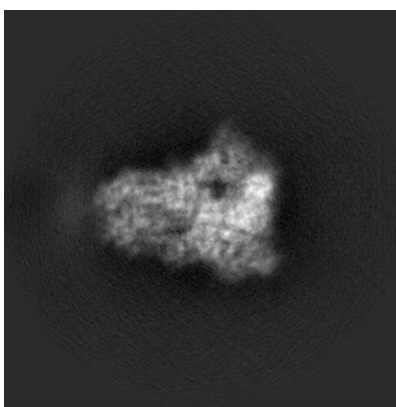


Z

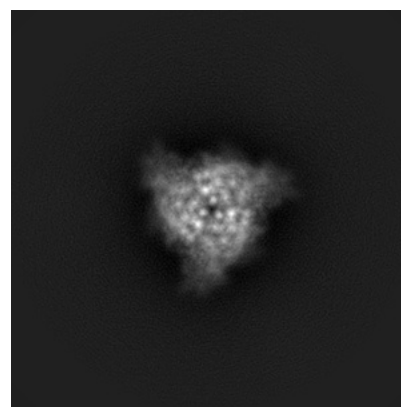
6.1.2 Raw map



X



Y

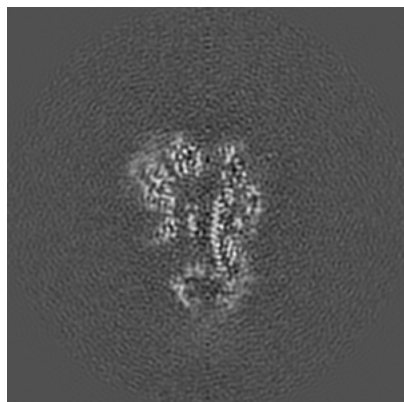


Z

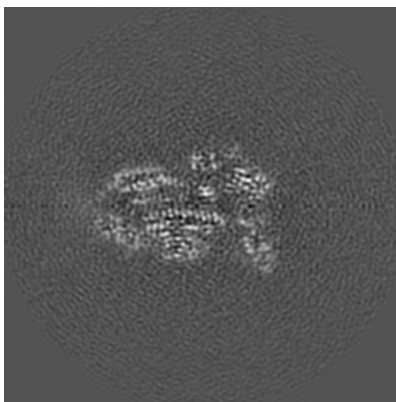
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

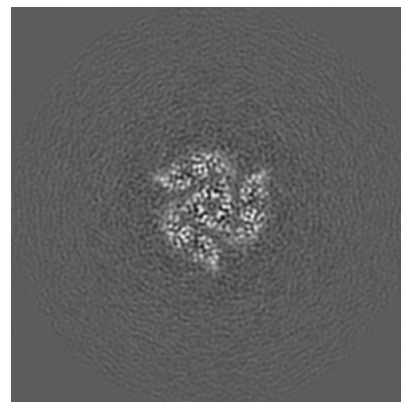
6.2.1 Primary map



X Index: 128

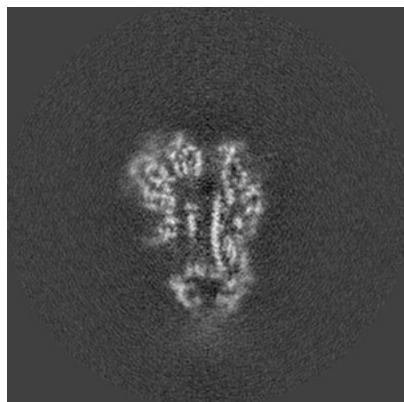


Y Index: 128

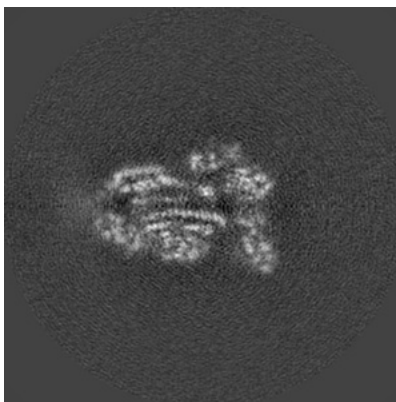


Z Index: 128

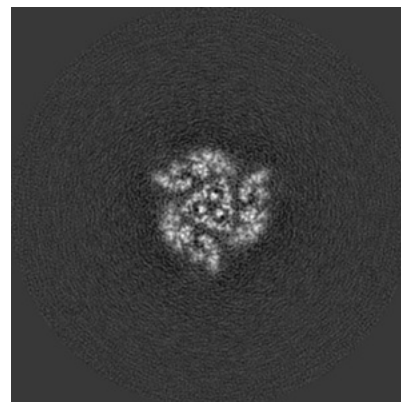
6.2.2 Raw map



X Index: 128



Y Index: 128

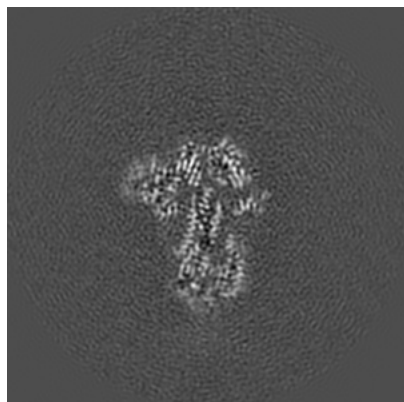


Z Index: 128

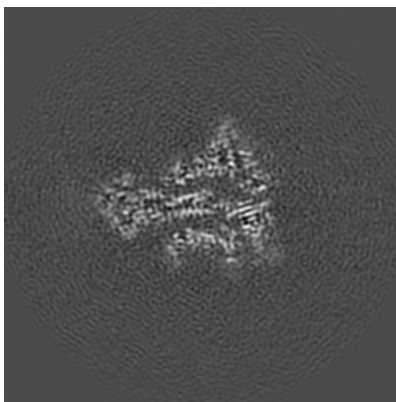
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

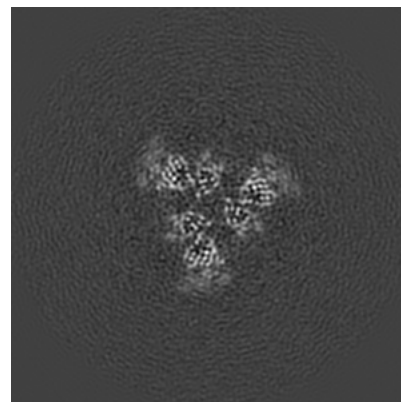
6.3.1 Primary map



X Index: 120

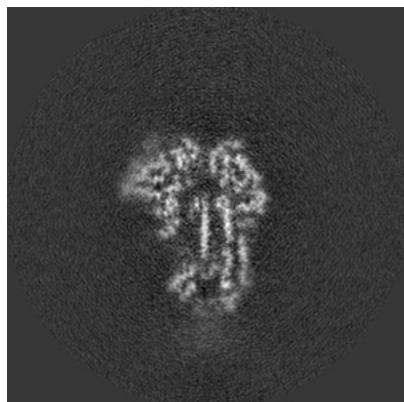


Y Index: 140

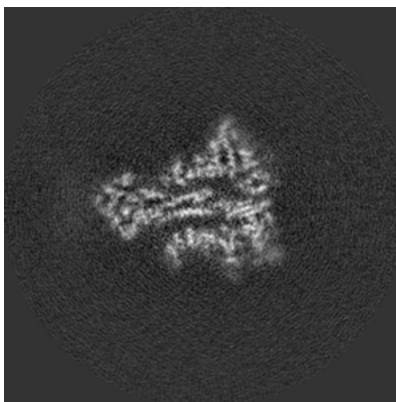


Z Index: 145

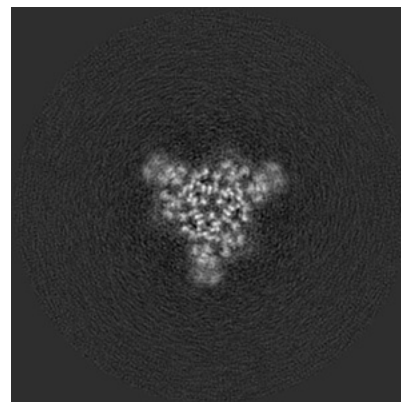
6.3.2 Raw map



X Index: 124



Y Index: 140

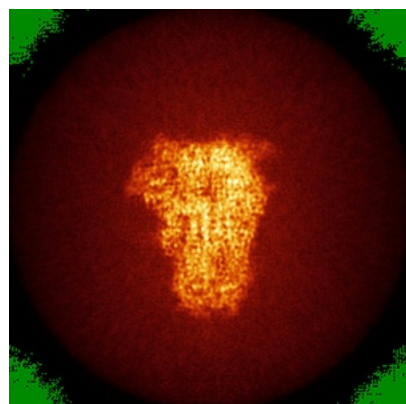


Z Index: 155

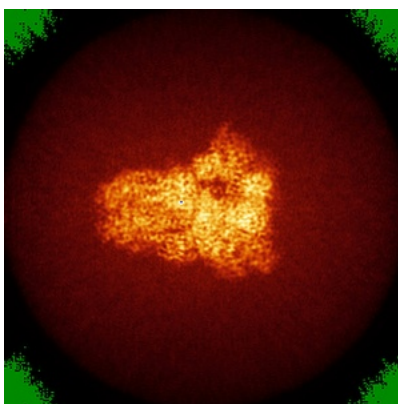
The images above show the largest variance slices of the map in three orthogonal directions.

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

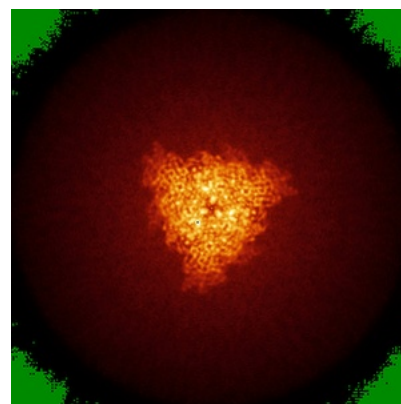
6.4.1 Primary map



X

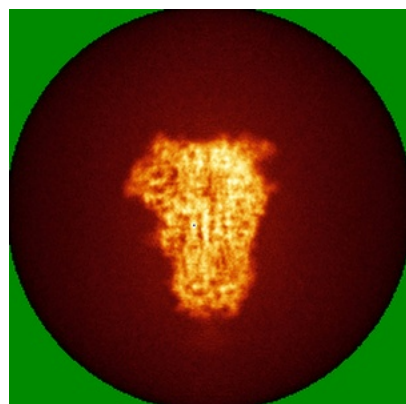


Y

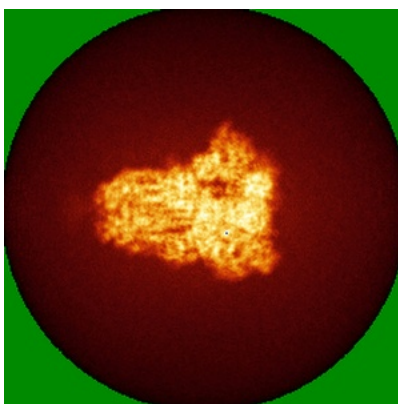


Z

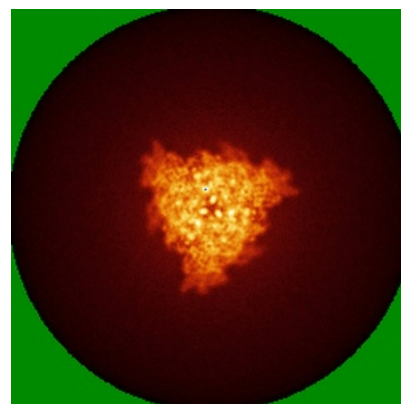
6.4.2 Raw map



X



Y

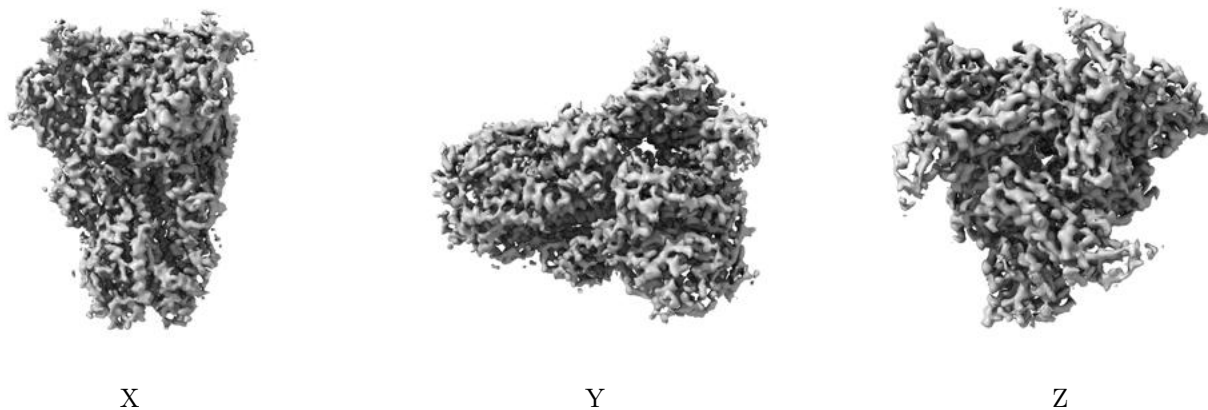


Z

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.

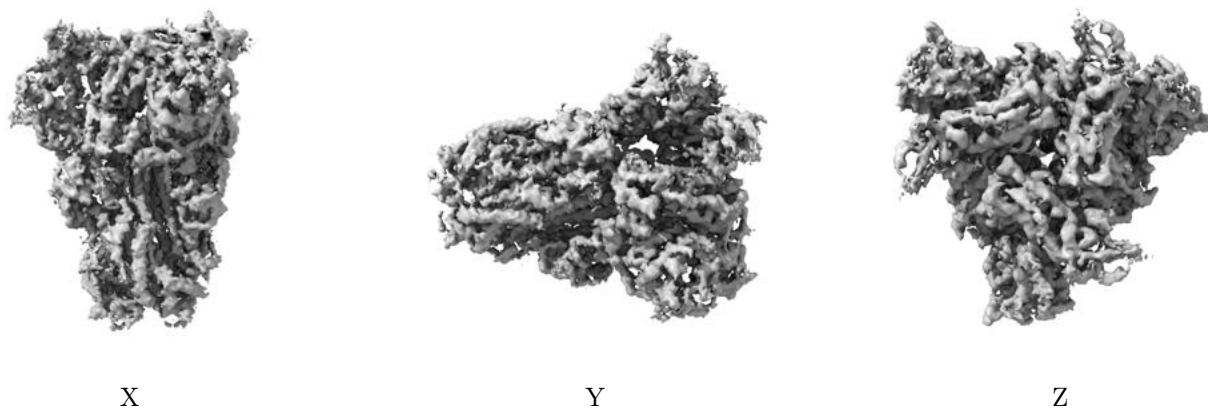
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0335. 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



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.

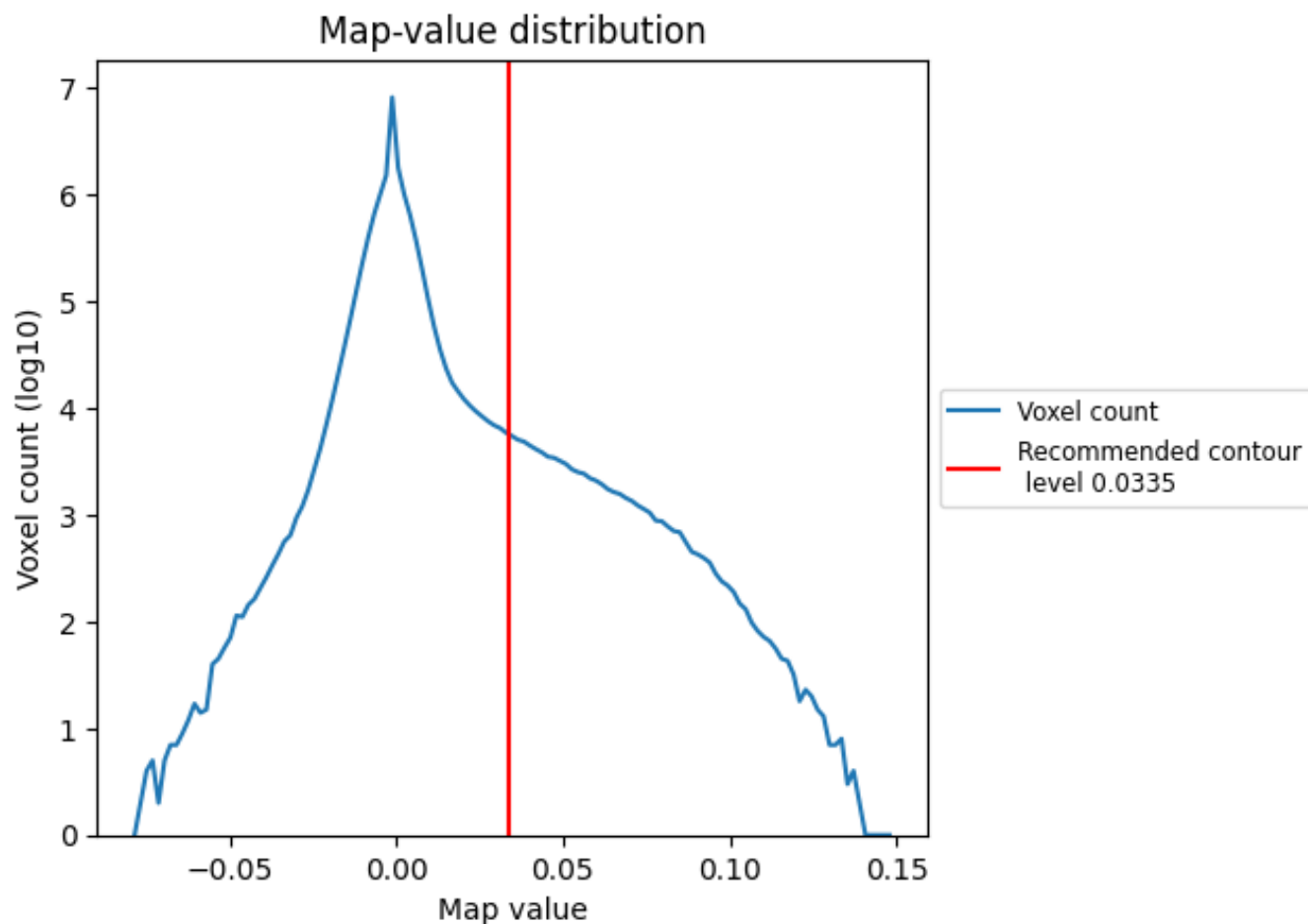
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

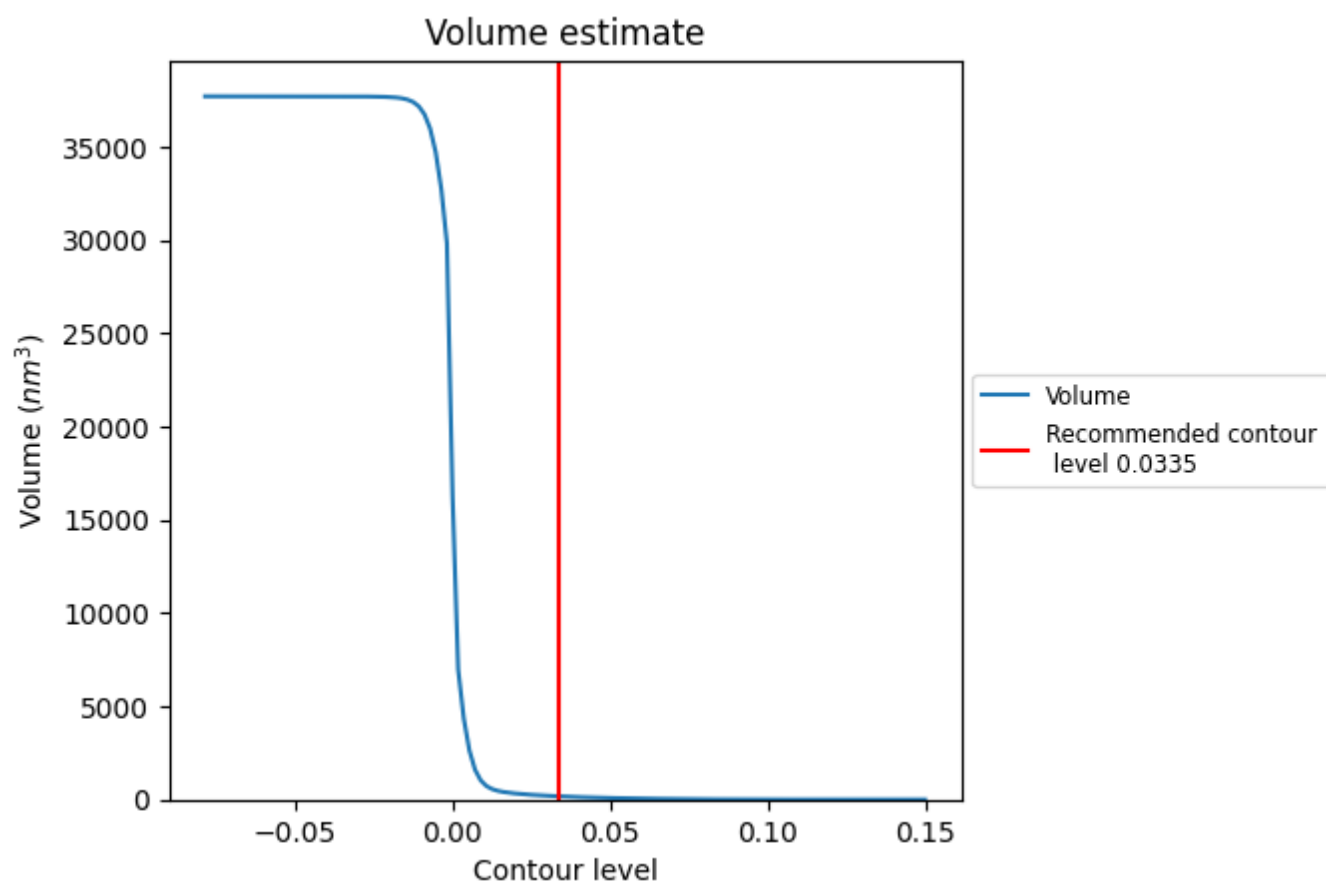
This section contains the results of statistical analysis of the map.

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.

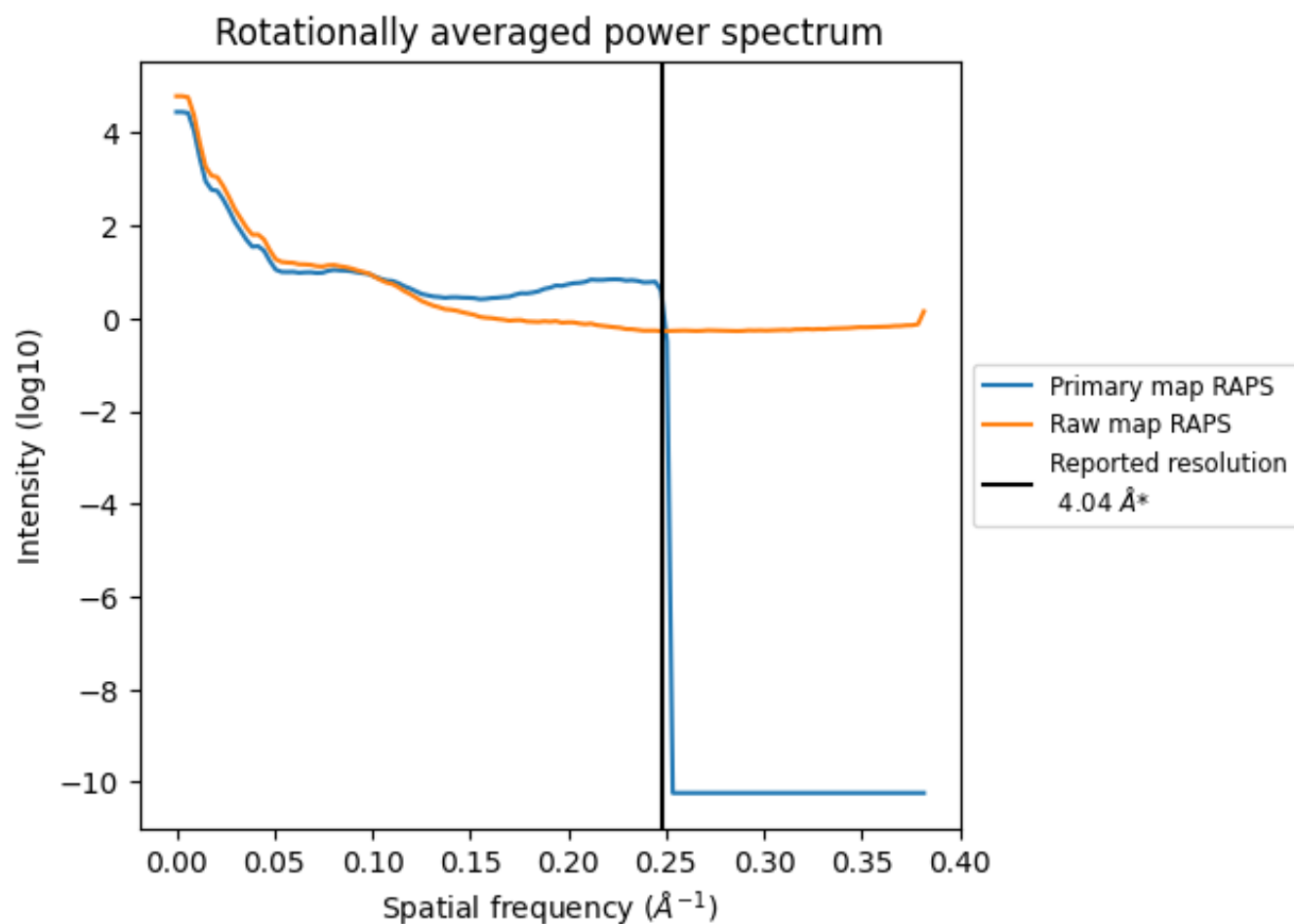
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 176 nm³; this corresponds to an approximate mass of 159 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 ⓘ

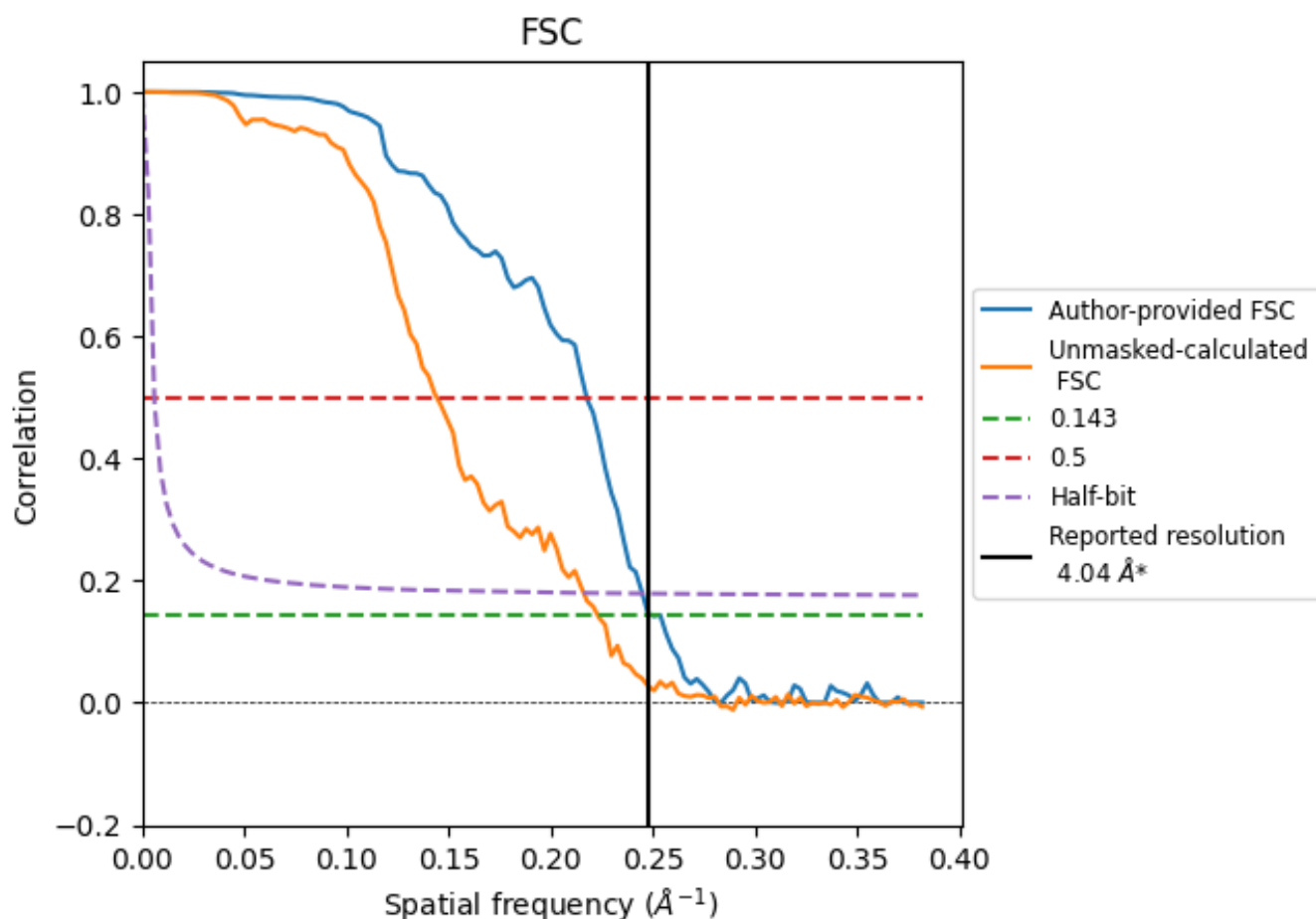


*Reported resolution corresponds to spatial frequency of 0.248 Å⁻¹

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 0.248 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

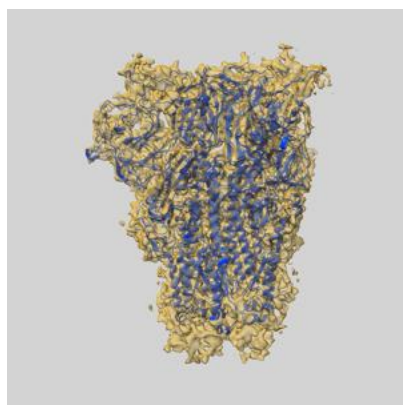
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.04	-	-
Author-provided FSC curve	4.01	4.60	4.08
Unmasked-calculated*	4.48	6.94	4.63

*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 4.48 differs from the reported value 4.04 by more than 10 %

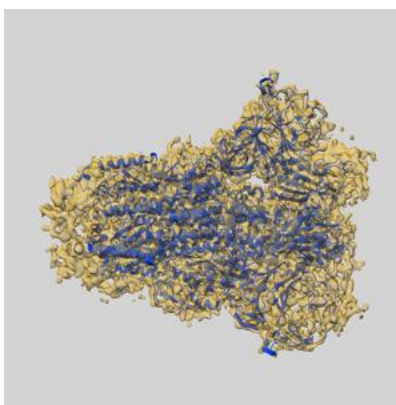
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-8069 and PDB model 5I08. Per-residue inclusion information can be found in section [3](#) on page [5](#).

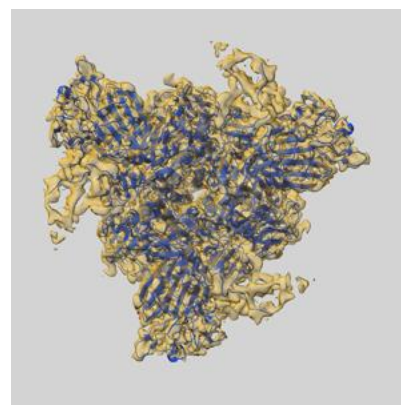
9.1 Map-model overlay [i](#)



X



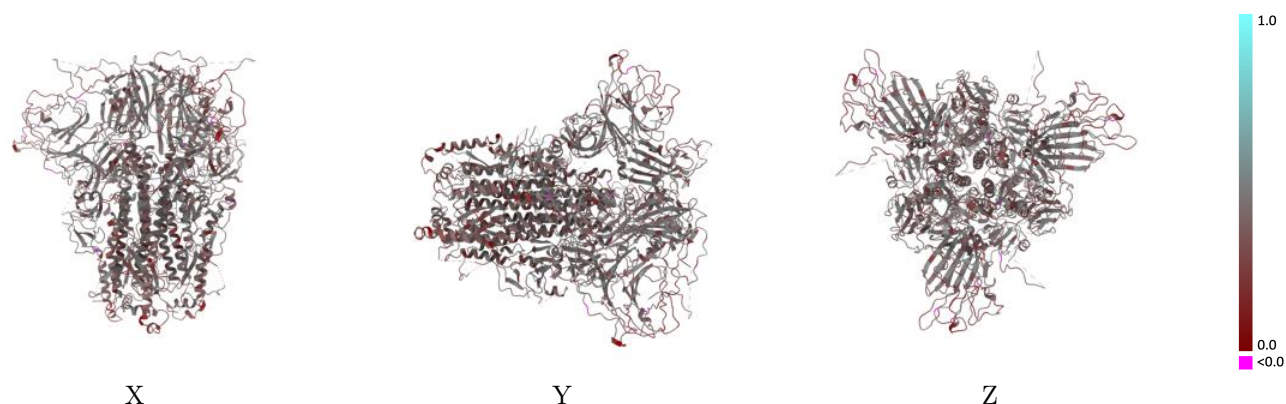
Y



Z

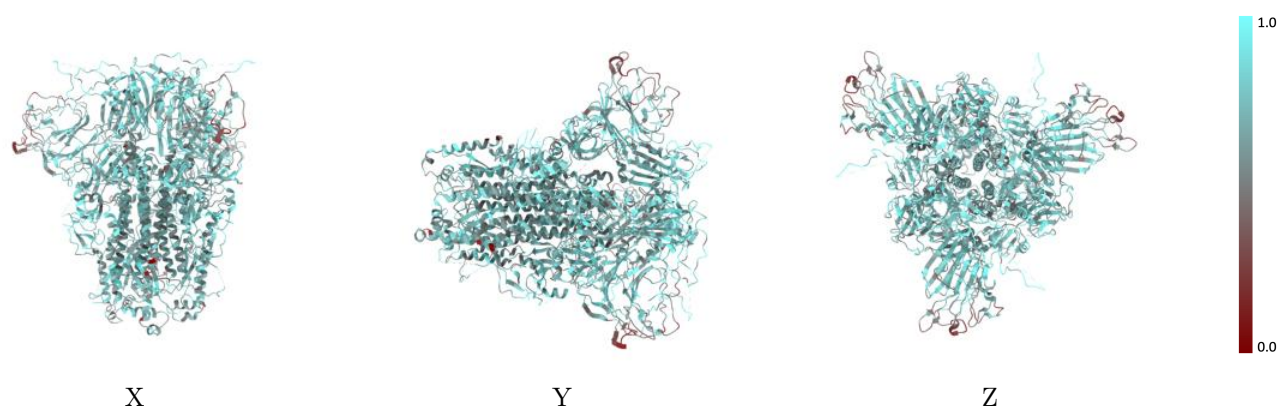
The images above show the 3D surface view of the map at the recommended contour level 0.0335 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



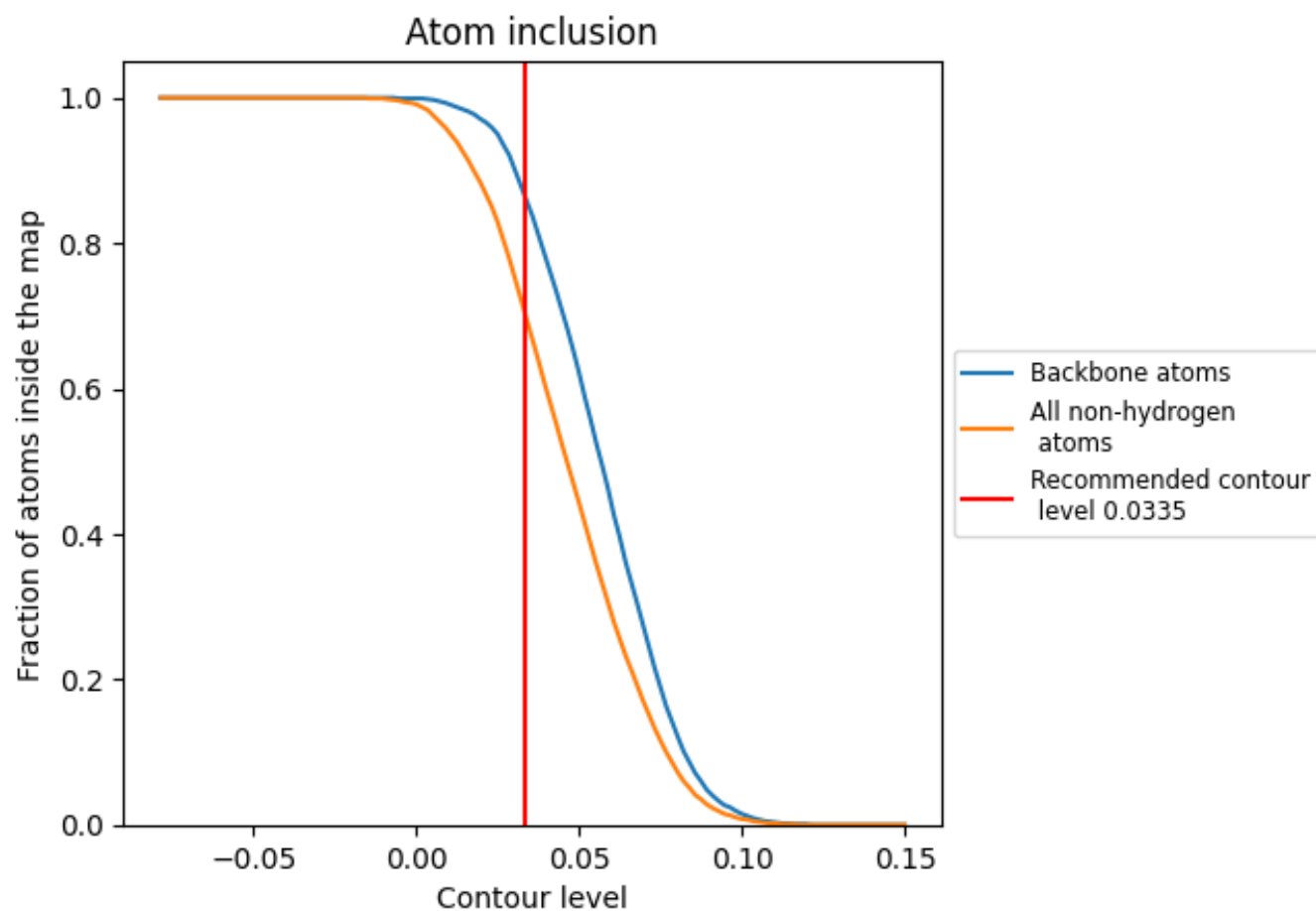
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0335).

9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 70% 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 (0.0335) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7000	0.3820
A	0.6990	0.3830
B	0.7030	0.3830
C	0.6980	0.3820

