



Full wwPDB EM Validation Report ⓘ

Mar 7, 2026 – 02:51 AM UTC

PDB ID : 8DZW / pdb_00008dzw
EMDB ID : EMD-27808
Title : RSV F trimer bound to RSV-199 Fab
Authors : Wen, X.; Jardetzky, T.S.
Deposited on : 2022-08-08
Resolution : 2.46 Å(reported)
Based on initial model : 5U68

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:

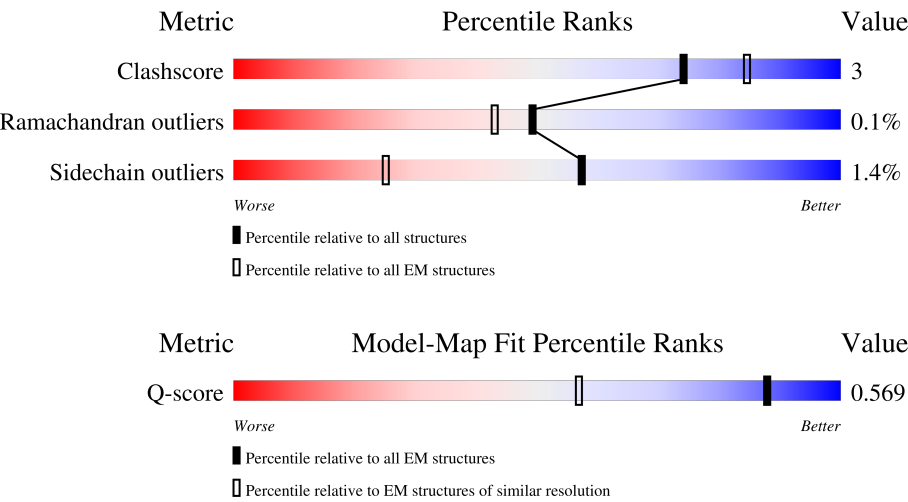
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.46 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	6014 (1.96 - 2.96)

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	484	
1	B	484	
1	C	484	
2	L	216	

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Mol	Chain	Length	Quality of chain
2	M	216	 17%80%19%
2	N	216	 18%82%16%
3	H	225	 19%83%16%
3	I	225	 20%80%18%
3	J	225	 21%83%15%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 40250 atoms, of which 20048 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 RSV fusion protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	444	Total	C	H	N	O	S	0	0
			6917	2173	3478	567	676	23		
1	B	444	Total	C	H	N	O	S	0	0
			6916	2173	3477	567	676	23		
1	C	444	Total	C	H	N	O	S	0	0
			6917	2173	3478	567	676	23		

There are 12 discrepancies between the modelled and reference sequences:

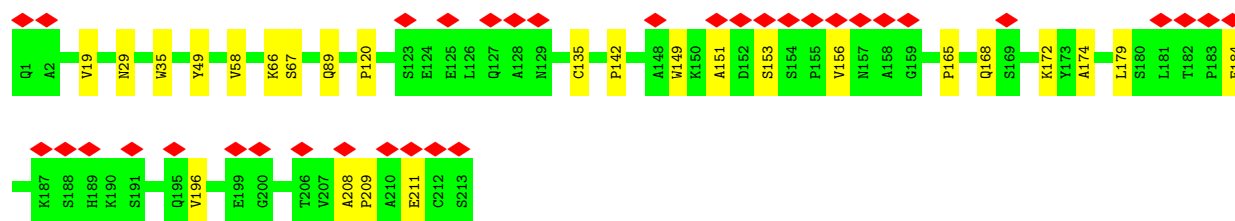
Chain	Residue	Modelled	Actual	Comment	Reference
A	152	VAL	ILE	engineered mutation	UNP A0A2H4WLA4
A	155	CYS	SER	engineered mutation	UNP A0A2H4WLA4
A	190	PHE	SER	engineered mutation	UNP A0A2H4WLA4
A	290	CYS	SER	engineered mutation	UNP A0A2H4WLA4
B	152	VAL	ILE	engineered mutation	UNP A0A2H4WLA4
B	155	CYS	SER	engineered mutation	UNP A0A2H4WLA4
B	190	PHE	SER	engineered mutation	UNP A0A2H4WLA4
B	290	CYS	SER	engineered mutation	UNP A0A2H4WLA4
C	152	VAL	ILE	engineered mutation	UNP A0A2H4WLA4
C	155	CYS	SER	engineered mutation	UNP A0A2H4WLA4
C	190	PHE	SER	engineered mutation	UNP A0A2H4WLA4
C	290	CYS	SER	engineered mutation	UNP A0A2H4WLA4

- Molecule 2 is a protein called RSV-199 Light chain protein.

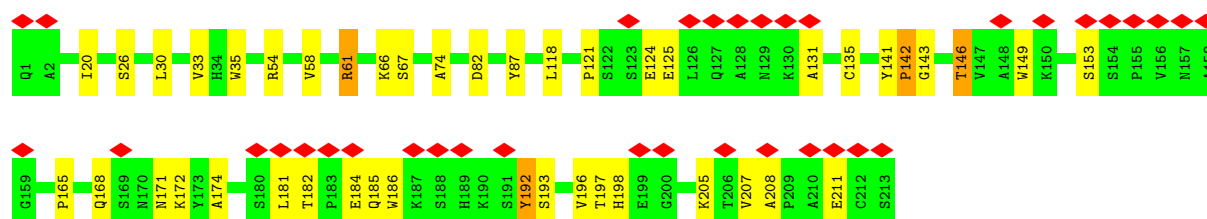
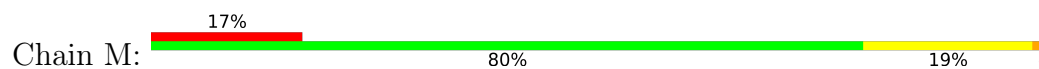
Mol	Chain	Residues	Atoms						AltConf	Trace
2	L	216	Total	C	H	N	O	S	0	0
			3159	1004	1553	272	325	5		
2	M	216	Total	C	H	N	O	S	0	0
			3159	1004	1553	272	325	5		
2	N	216	Total	C	H	N	O	S	0	0
			3159	1004	1553	272	325	5		

- Molecule 3 is a protein called RSV-199 Heavy chain protein.

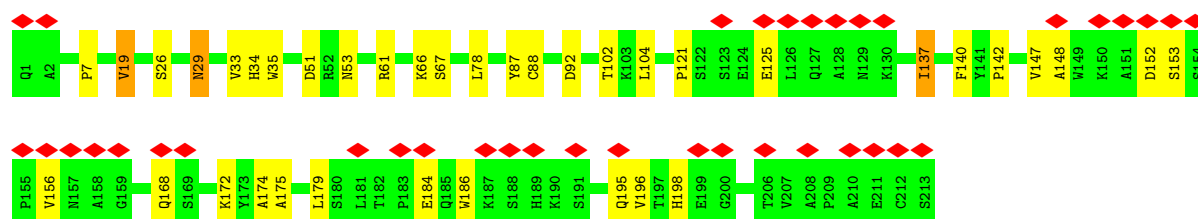
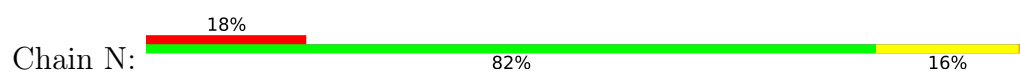
Mol	Chain	Residues	Atoms						AltConf	Trace
3	H	225	Total	C	H	N	O	S	0	0
			3341	1064	1652	287	331	7		
3	I	225	Total	C	H	N	O	S	0	0
			3341	1064	1652	287	331	7		
3	J	225	Total	C	H	N	O	S	0	0
			3341	1064	1652	287	331	7		



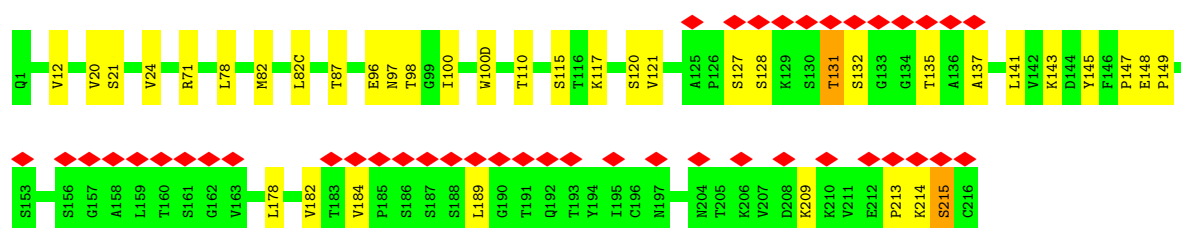
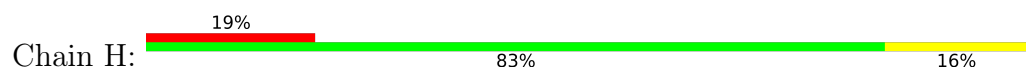
• Molecule 2: RSV-199 Light chain protein



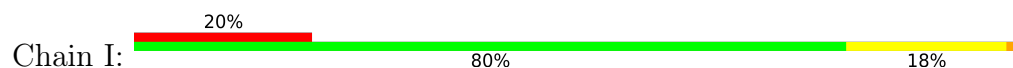
• Molecule 2: RSV-199 Light chain protein



• Molecule 3: RSV-199 Heavy chain protein

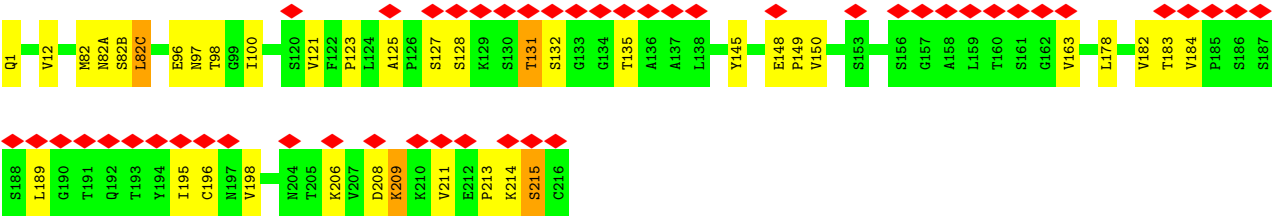
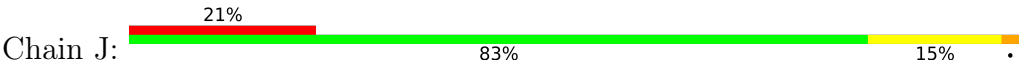


• Molecule 3: RSV-199 Heavy chain protein





• Molecule 3: RSV-199 Heavy chain protein



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	476914	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	4508	Depositor
Maximum defocus (nm)	40000	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	6.454	Depositor
Minimum map value	-4.056	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.111	Depositor
Recommended contour level	0.414	Depositor
Map size ($\text{\AA}$)	355.2, 355.2, 355.2	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.11, 1.11, 1.11	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	0.34	2/3488 (0.1%)	0.53	2/4723 (0.0%)
1	B	0.24	1/3488 (0.0%)	0.41	0/4723
1	C	0.30	1/3488 (0.0%)	0.49	1/4723 (0.0%)
2	L	0.76	4/1646 (0.2%)	1.09	8/2246 (0.4%)
2	M	0.60	2/1646 (0.1%)	0.93	6/2246 (0.3%)
2	N	0.60	3/1646 (0.2%)	0.99	6/2246 (0.3%)
3	H	0.55	1/1731 (0.1%)	0.88	9/2358 (0.4%)
3	I	0.68	6/1731 (0.3%)	0.94	8/2358 (0.3%)
3	J	0.66	4/1731 (0.2%)	0.93	10/2358 (0.4%)
All	All	0.50	24/20595 (0.1%)	0.76	50/27981 (0.2%)

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	49	TYR	C-O	-12.41	1.08	1.23
3	I	184	VAL	C-O	-9.91	1.16	1.24
1	C	242	GLY	C-N	8.49	1.45	1.33
2	L	35	TRP	C-O	-7.98	1.14	1.23
3	I	100	ILE	C-O	-6.97	1.15	1.24
3	J	100	ILE	C-O	-6.82	1.16	1.24
1	A	247	VAL	C-N	6.61	1.42	1.33
3	I	99	GLY	C-O	-6.48	1.15	1.23
2	N	172	LYS	C-O	-6.46	1.16	1.24
2	M	172	LYS	C-O	-6.42	1.15	1.23
2	L	142	PRO	C-O	-6.24	1.15	1.23
3	J	97	ASN	C-O	-6.14	1.16	1.23
3	I	97	ASN	C-O	-6.14	1.16	1.23
3	H	97	ASN	C-O	-6.13	1.16	1.23
2	M	171	ASN	C-O	-6.04	1.15	1.23
3	I	123	PRO	C-O	-6.02	1.16	1.23
3	J	195	ILE	C-O	-5.82	1.18	1.24
2	L	172	LYS	C-O	-5.57	1.16	1.24
3	J	123	PRO	C-O	-5.46	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	31	GLU	C-O	-5.21	1.18	1.24
3	I	127	SER	C-N	5.18	1.40	1.33
2	N	142	PRO	C-O	-5.18	1.16	1.23
2	N	53	ASN	C-O	-5.13	1.17	1.23
1	A	248	SER	CA-CB	-5.07	1.45	1.53

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	120	PRO	CB-CA-C	-8.61	100.42	110.92
1	C	501	GLN	O-C-N	8.25	130.61	122.03
2	M	142	PRO	CA-C-O	-7.88	112.99	121.27
2	N	61	ARG	N-CA-C	-7.85	103.49	113.23
2	L	142	PRO	CA-C-O	-7.66	113.23	121.27
2	M	142	PRO	N-CA-CB	7.50	109.84	103.17
3	J	97	ASN	CA-CB-CG	7.32	119.92	112.60
3	I	97	ASN	CA-CB-CG	7.29	119.89	112.60
3	H	97	ASN	CA-CB-CG	7.28	119.88	112.60
2	L	135	CYS	CA-C-O	-7.12	113.01	120.70
3	I	123	PRO	CA-C-O	-7.11	113.39	122.13
3	J	123	PRO	CA-C-O	-7.05	113.38	122.12
2	L	184	GLU	CA-C-O	-7.04	113.42	120.82
2	N	142	PRO	CA-C-O	-6.86	114.07	121.27
3	H	147	PRO	CA-C-N	6.30	128.71	120.26
3	H	147	PRO	C-N-CA	6.30	128.71	120.26
2	L	58	VAL	CB-CA-C	6.28	122.78	111.36
3	I	207	VAL	CA-C-O	-6.25	116.74	121.68
2	M	61	ARG	N-CA-C	-6.23	105.72	113.38
2	L	151	ALA	CA-C-O	-5.96	113.98	120.36
3	J	208	ASP	CA-CB-CG	5.86	118.46	112.60
3	J	195	ILE	CA-C-O	-5.82	114.19	120.36
3	J	96	GLU	N-CA-C	-5.65	100.57	109.50
3	H	96	GLU	N-CA-C	-5.65	100.58	109.50
3	I	96	GLU	N-CA-C	-5.64	100.59	109.50
3	I	208	ASP	CA-CB-CG	5.58	118.18	112.60
1	A	345	ASN	CA-CB-CG	5.56	118.16	112.60
3	J	127	SER	O-C-N	-5.51	116.46	123.24
3	H	215	SER	CB-CA-C	-5.48	110.23	116.54
2	N	152	ASP	N-CA-C	-5.48	99.98	108.31
3	J	215	SER	CB-CA-C	-5.42	110.31	116.54
2	M	192	TYR	CA-C-O	-5.39	115.32	121.68
2	L	165	PRO	N-CA-CB	5.37	108.09	103.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	215	SER	CB-CA-C	-5.34	110.40	116.54
3	J	209	LYS	CB-CA-C	5.32	118.00	110.24
3	H	131	THR	OG1-CB-CG2	-5.31	98.68	109.30
1	A	26	GLN	CB-CA-C	-5.23	100.17	110.10
3	I	184	VAL	CA-C-O	-5.23	114.70	121.13
2	L	120	PRO	N-CA-CB	5.20	108.13	103.08
2	M	58	VAL	CB-CA-C	5.20	120.82	111.36
3	H	100	ILE	CA-C-O	-5.15	114.34	120.78
2	N	26	SER	CA-C-O	-5.12	115.28	121.32
3	H	147	PRO	N-CA-CB	5.11	107.86	103.31
3	I	131	THR	OG1-CB-CG2	-5.06	99.18	109.30
3	J	121	VAL	CA-C-O	-5.06	115.08	120.85
2	N	19	VAL	CA-C-O	-5.05	116.02	121.92
3	H	127	SER	O-C-N	-5.04	116.59	123.15
3	J	131	THR	OG1-CB-CG2	-5.02	99.25	109.30
2	N	175	ALA	CA-C-O	-5.02	115.67	121.49
2	M	26	SER	CA-C-O	-5.00	115.42	121.32

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	3439	3478	3476	6	0
1	B	3439	3477	3476	16	0
1	C	3439	3478	3476	7	0
2	L	1606	1553	1553	6	0
2	M	1606	1553	1553	21	0
2	N	1606	1553	1553	16	0
3	H	1689	1652	1652	18	0
3	I	1689	1652	1652	25	0
3	J	1689	1652	1652	15	0
All	All	20202	20048	20043	128	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 (128) 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:264:MET:HE1	1:B:274:MET:HE1	1.62	0.81
2:M:146:THR:HG1	2:M:197:THR:HG1	1.26	0.81
2:L:208:ALA:HB3	2:L:211:GLU:HB3	1.63	0.78
3:I:135:THR:HB	3:I:184:VAL:O	1.85	0.75
2:M:208:ALA:HB3	2:M:211:GLU:HB3	1.71	0.73
3:I:82:MET:HE2	3:I:82(C):LEU:HD21	1.72	0.70
3:I:184:VAL:HG11	3:I:194:TYR:CE1	2.28	0.68
2:M:146:THR:OG1	2:M:197:THR:OG1	2.11	0.67
1:A:289:MET:HE1	1:A:297:LEU:CD1	2.26	0.64
2:N:168:GLN:OE1	2:N:174:ALA:HB2	1.98	0.64
1:A:289:MET:HE1	1:A:297:LEU:HD11	1.80	0.63
2:N:19:VAL:HG21	2:N:104:LEU:HD11	1.79	0.62
2:M:168:GLN:OE1	2:M:174:ALA:HB2	1.99	0.62
2:L:168:GLN:OE1	2:L:174:ALA:HB2	2.00	0.62
2:N:29:ASN:OD1	2:N:92:ASP:OD1	2.19	0.60
1:B:426:ASN:OD1	1:B:432:ILE:HD13	2.02	0.60
3:I:145:TYR:CZ	3:I:178:LEU:HD22	2.38	0.58
3:I:128:SER:HA	3:I:131:THR:HG22	1.85	0.58
3:H:145:TYR:OH	3:H:178:LEU:HD22	2.04	0.58
3:J:145:TYR:OH	3:J:178:LEU:HD22	2.05	0.57
1:C:329:GLY:O	1:C:330:SER:OG	2.20	0.57
3:H:128:SER:HA	3:H:131:THR:HG22	1.88	0.56
3:J:145:TYR:CZ	3:J:178:LEU:HD22	2.41	0.55
3:I:145:TYR:OH	3:I:178:LEU:HD22	2.06	0.55
2:M:30:LEU:HD21	2:M:33:VAL:HG22	1.88	0.55
2:N:33:VAL:N	2:N:51:ASP:OD1	2.34	0.55
2:L:89:GLN:NE2	3:H:100(D):TRP:HA	2.22	0.54
3:J:189:LEU:HD21	3:J:213:PRO:HG3	1.88	0.54
1:B:183:ASN:OD1	1:B:185:VAL:HG12	2.07	0.54
3:H:189:LEU:HD21	3:H:213:PRO:HG3	1.89	0.54
3:J:131:THR:HG23	3:J:132:SER:N	2.23	0.54
1:C:289:MET:HE1	1:C:297:LEU:CD1	2.38	0.53
3:J:128:SER:HA	3:J:131:THR:HG22	1.90	0.53
3:J:135:THR:HB	3:J:184:VAL:O	2.08	0.53
3:I:189:LEU:HD21	3:I:213:PRO:HG3	1.89	0.52
3:H:131:THR:HG23	3:H:132:SER:N	2.25	0.52
2:N:156:VAL:HG21	2:N:179:LEU:HD21	1.90	0.52
3:J:196:CYS:SG	3:J:209:LYS:O	2.69	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:145:TYR:CZ	3:H:178:LEU:HD22	2.46	0.51
1:B:392:ASP:OD2	1:B:491:SER:OG	2.26	0.51
2:M:181:LEU:HD11	2:M:192:TYR:CE2	2.46	0.50
1:A:183:ASN:OD1	1:A:185:VAL:HG12	2.12	0.50
2:N:7:PRO:O	2:N:102:THR:OG1	2.24	0.49
3:I:131:THR:HG23	3:I:132:SER:N	2.27	0.49
1:C:354:GLN:O	1:C:357:THR:HG22	2.12	0.49
3:I:40:ALA:HB3	3:I:43:LYS:HB2	1.95	0.49
1:C:316:LEU:HD23	1:C:338:ASP:O	2.13	0.48
1:C:183:ASN:OD1	1:C:185:VAL:HG12	2.14	0.48
3:J:214:LYS:O	3:J:215:SER:C	2.57	0.48
1:B:30:GLU:OE1	1:B:41:SER:OG	2.31	0.48
3:I:214:LYS:O	3:I:215:SER:C	2.57	0.47
3:H:135:THR:HB	3:H:184:VAL:O	2.13	0.47
2:N:35:TRP:HA	2:N:87:TYR:O	2.14	0.47
1:B:188:LEU:HD23	1:B:260:LEU:HD12	1.96	0.47
1:C:188:LEU:HD23	1:C:260:LEU:HD12	1.97	0.47
1:B:398:SER:OG	1:B:399:LYS:N	2.47	0.47
3:H:214:LYS:O	3:H:215:SER:C	2.57	0.46
2:N:137:ILE:HD11	2:N:147:VAL:CG2	2.47	0.45
2:M:35:TRP:HA	2:M:87:TYR:O	2.17	0.45
3:H:87:THR:HG23	3:H:110:THR:HA	1.99	0.45
2:M:61:ARG:NH2	2:M:82:ASP:OD2	2.50	0.45
2:M:184:GLU:OE1	2:M:184:GLU:N	2.49	0.45
3:I:12:VAL:HG11	3:I:82(C):LEU:HD13	1.99	0.45
1:B:432:ILE:HD11	1:B:447:VAL:HG22	1.98	0.44
1:B:206:ILE:HD12	1:B:206:ILE:H	1.82	0.44
1:C:190:PHE:CE2	1:C:192:VAL:HG12	2.52	0.44
3:I:127:SER:O	3:I:131:THR:HB	2.17	0.44
1:B:292:ILE:O	1:B:292:ILE:HG23	2.17	0.44
2:N:184:GLU:OE1	2:N:184:GLU:N	2.51	0.44
2:M:143:GLY:O	2:M:165:PRO:HG3	2.18	0.44
3:I:150:VAL:HG22	3:I:200:HIS:HD2	1.82	0.44
2:L:208:ALA:HB1	2:L:209:PRO:HD2	1.98	0.43
3:H:71:ARG:HA	3:H:78:LEU:HA	2.00	0.43
1:A:68:LYS:O	1:A:69:CYS:C	2.61	0.43
1:B:289:MET:HE1	1:B:297:LEU:CD1	2.49	0.43
3:I:107:THR:HG23	3:I:107:THR:O	2.19	0.43
3:I:148:GLU:N	3:I:149:PRO:HD2	2.34	0.43
3:I:121:VAL:HA	3:I:141:LEU:O	2.19	0.43
3:H:115:SER:O	3:H:117:LYS:HG3	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:66:LYS:HG2	2:N:67:SER:N	2.33	0.42
3:J:196:CYS:SG	3:J:209:LYS:HB3	2.60	0.42
3:J:125:ALA:HA	3:J:211:VAL:CG1	2.49	0.42
3:H:120:SER:O	3:H:143:LYS:N	2.49	0.42
3:I:200:HIS:O	3:I:204:ASN:N	2.52	0.42
2:N:19:VAL:HG21	2:N:104:LEU:CD1	2.47	0.42
2:N:34:HIS:O	2:N:88:CYS:HA	2.19	0.42
2:N:121:PRO:HD2	2:N:186:TRP:CZ2	2.55	0.42
1:B:190:PHE:CE2	1:B:192:VAL:HG12	2.54	0.42
1:B:426:ASN:OD1	1:B:432:ILE:HG21	2.20	0.42
3:J:12:VAL:HG11	3:J:82(C):LEU:HD13	2.01	0.42
2:L:149:TRP:CE3	2:L:179:LEU:HD22	2.55	0.42
1:A:289:MET:HE1	1:A:297:LEU:HD12	2.00	0.42
2:N:148:ALA:HB3	2:N:195:GLN:HE21	1.84	0.42
3:I:184:VAL:HG11	3:I:194:TYR:HE1	1.82	0.42
1:A:401:ASP:OD1	1:A:401:ASP:N	2.53	0.42
3:H:20:VAL:HG12	3:H:21:SER:N	2.35	0.42
3:H:78:LEU:H	3:H:78:LEU:HD23	1.85	0.42
3:H:137:ALA:HA	3:H:182:VAL:O	2.19	0.42
3:I:169:VAL:HG12	3:I:177:SER:O	2.20	0.42
2:M:182:THR:OG1	2:M:185:GLN:HG3	2.20	0.41
3:H:82:MET:HB3	3:H:82(C):LEU:HD21	2.01	0.41
2:M:135:CYS:HB2	2:M:149:TRP:CZ2	2.55	0.41
1:B:426:ASN:ND2	1:B:446:GLY:O	2.33	0.41
2:M:121:PRO:HD2	2:M:186:TRP:CZ2	2.55	0.41
2:M:141:TYR:CD2	2:M:142:PRO:HD3	2.55	0.41
2:M:66:LYS:HG2	2:M:67:SER:N	2.35	0.41
2:L:66:LYS:HG2	2:L:67:SER:N	2.35	0.41
2:M:142:PRO:HD2	2:M:198:HIS:CE1	2.55	0.41
2:M:20:ILE:HD13	2:M:74:ALA:HA	2.02	0.41
2:M:118:LEU:HB3	2:M:205:LYS:HD2	2.02	0.41
3:J:135:THR:OG1	3:J:183:THR:HG23	2.20	0.41
2:M:125:GLU:HG3	2:M:131:ALA:HA	2.03	0.41
2:N:140:PHE:HB2	2:N:198:HIS:CE1	2.56	0.41
3:J:198:VAL:O	3:J:206:LYS:HA	2.21	0.41
3:I:138:LEU:HD21	3:I:194:TYR:CE2	2.56	0.41
3:J:148:GLU:N	3:J:149:PRO:HD2	2.35	0.41
2:M:193:SER:HA	2:M:205:LYS:O	2.21	0.40
3:I:71:ARG:HA	3:I:78:LEU:HA	2.03	0.40
3:I:120:SER:HB3	3:I:143:LYS:HB3	2.03	0.40
3:J:163:VAL:HG22	3:J:182:VAL:HG13	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:121:VAL:HA	3:H:141:LEU:O	2.21	0.40
2:M:124:GLU:HG3	3:I:122:PHE:CD1	2.56	0.40
2:N:78:LEU:HD11	2:N:104:LEU:HD21	2.03	0.40
1:B:64:ILE:O	1:B:87:LYS:NZ	2.54	0.40
3:I:143:LYS:HG2	3:I:144:ASP:N	2.35	0.40
1:B:197:ASN:OD1	1:B:201:LYS:HD2	2.21	0.40
3:H:148:GLU:N	3:H:149:PRO:HD2	2.36	0.40
3:I:1:GLN:O	3:I:1:GLN:HG2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	438/484 (90%)	416 (95%)	22 (5%)	0	100	100
1	B	438/484 (90%)	421 (96%)	16 (4%)	1 (0%)	43	54
1	C	438/484 (90%)	418 (95%)	19 (4%)	1 (0%)	43	54
2	L	214/216 (99%)	201 (94%)	13 (6%)	0	100	100
2	M	214/216 (99%)	199 (93%)	15 (7%)	0	100	100
2	N	214/216 (99%)	200 (94%)	14 (6%)	0	100	100
3	H	223/225 (99%)	208 (93%)	15 (7%)	0	100	100
3	I	223/225 (99%)	206 (92%)	17 (8%)	0	100	100
3	J	223/225 (99%)	209 (94%)	14 (6%)	0	100	100
All	All	2625/2775 (95%)	2478 (94%)	145 (6%)	2 (0%)	49	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	70	ASN
1	B	27	ASN

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	405/441 (92%)	403 (100%)	2 (0%)	81	87
1	B	405/441 (92%)	404 (100%)	1 (0%)	87	92
1	C	405/441 (92%)	403 (100%)	2 (0%)	81	87
2	L	177/177 (100%)	172 (97%)	5 (3%)	38	54
2	M	177/177 (100%)	172 (97%)	5 (3%)	38	54
2	N	177/177 (100%)	172 (97%)	5 (3%)	38	54
3	H	192/192 (100%)	188 (98%)	4 (2%)	47	62
3	I	192/192 (100%)	191 (100%)	1 (0%)	81	87
3	J	192/192 (100%)	185 (96%)	7 (4%)	31	45
All	All	2322/2430 (96%)	2290 (99%)	32 (1%)	57	72

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	172	LEU
1	A	178	VAL
1	B	204	LEU
1	C	99	SER
1	C	295	GLU
2	L	19	VAL
2	L	29	ASN
2	L	153	SER
2	L	156	VAL
2	L	196	VAL
3	H	12	VAL
3	H	24	VAL
3	H	98	THR

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Mol	Chain	Res	Type
3	H	209	LYS
2	M	54	ARG
2	M	146	THR
2	M	153	SER
2	M	196	VAL
2	M	207	VAL
3	I	98	THR
2	N	29	ASN
2	N	125	GLU
2	N	137	ILE
2	N	153	SER
2	N	196	VAL
3	J	1	GLN
3	J	82	MET
3	J	82(A)	ASN
3	J	82(B)	SER
3	J	82(C)	LEU
3	J	98	THR
3	J	150	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	317	HIS
1	B	27	ASN
1	B	383	ASN
1	C	81	GLN
1	C	159	HIS
3	H	105	GLN
3	H	155	ASN
3	H	197	ASN
2	M	37	GLN
2	M	89	GLN
2	N	53	ASN
3	J	155	ASN

5.3.3 RNA ⓘ

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](#)

There are no chain breaks in this entry.

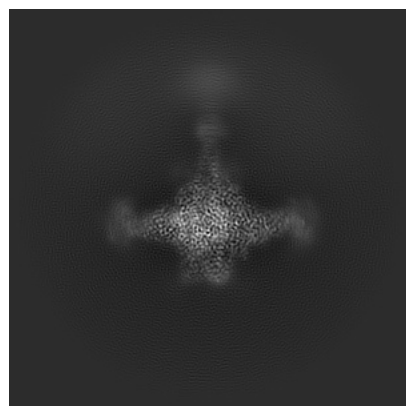
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-27808. These allow visual inspection of the internal detail of the map and identification of artifacts.

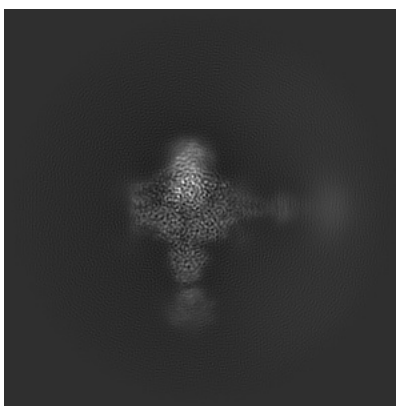
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

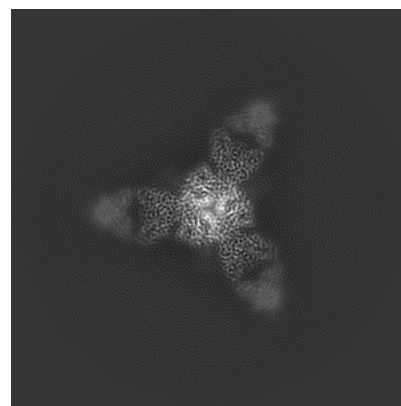
6.1.1 Primary map



X

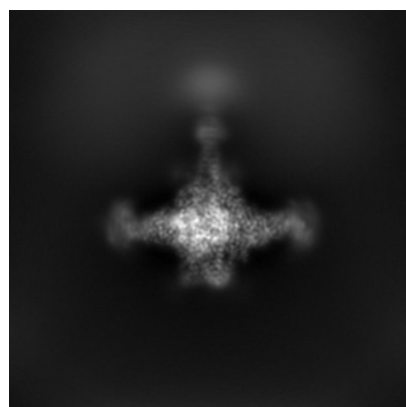


Y

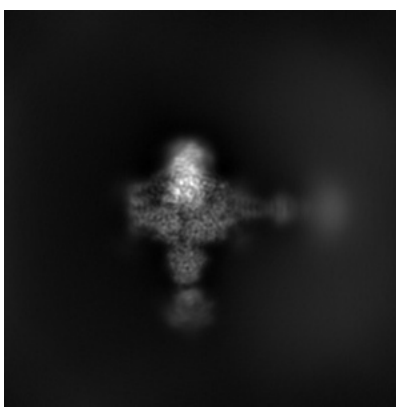


Z

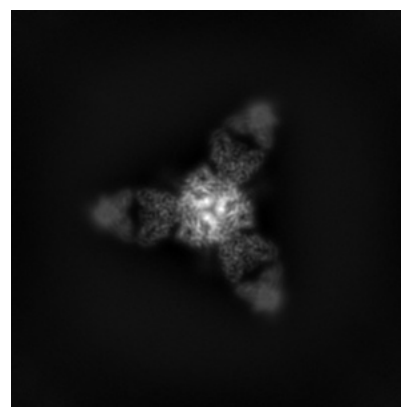
6.1.2 Raw map



X



Y

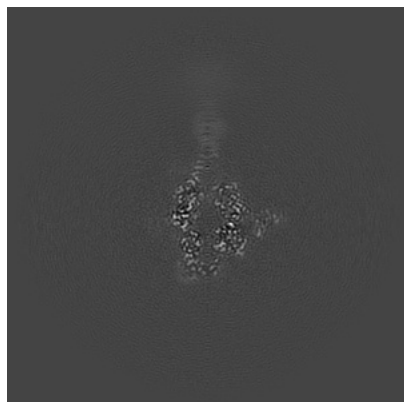


Z

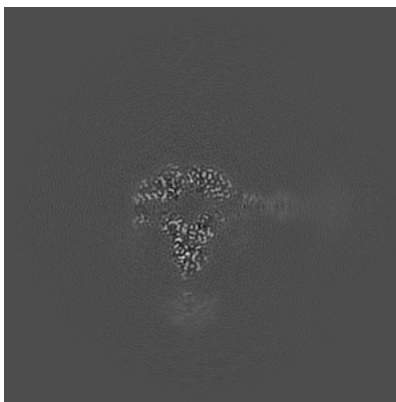
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

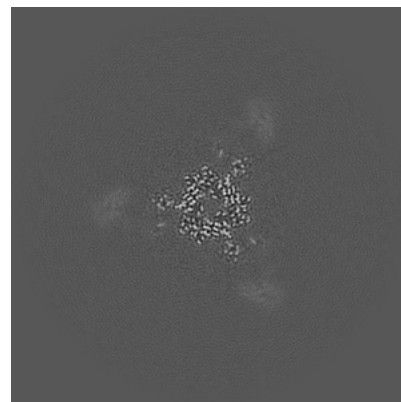
6.2.1 Primary map



X Index: 160



Y Index: 160

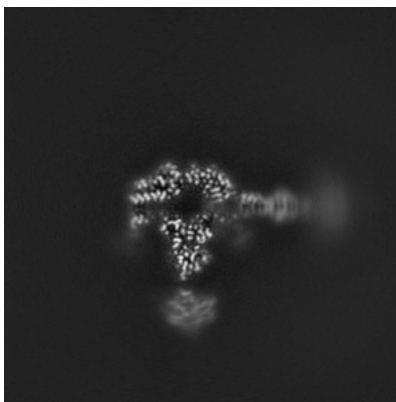


Z Index: 160

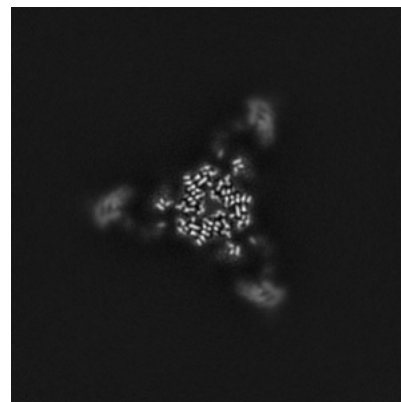
6.2.2 Raw map



X Index: 160



Y Index: 160

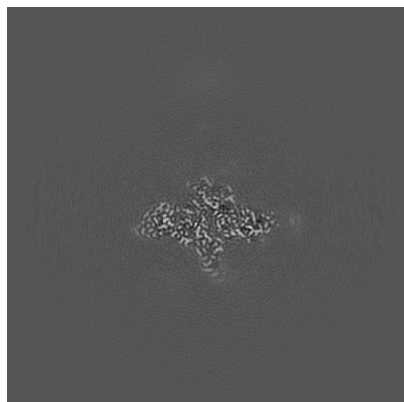


Z Index: 160

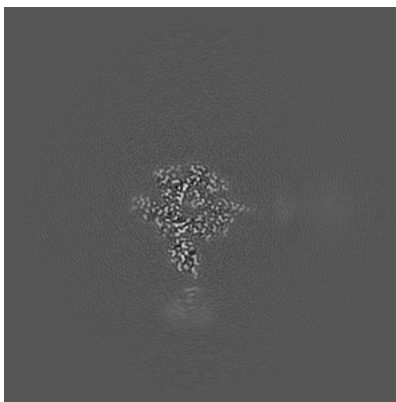
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

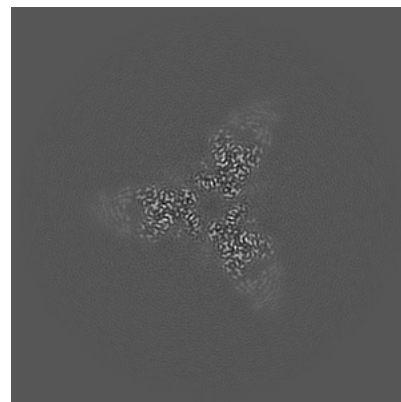
6.3.1 Primary map



X Index: 176



Y Index: 150

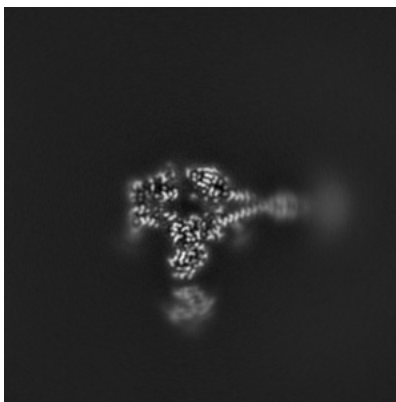


Z Index: 146

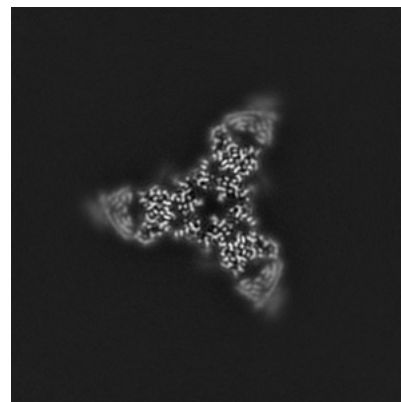
6.3.2 Raw map



X Index: 167



Y Index: 164

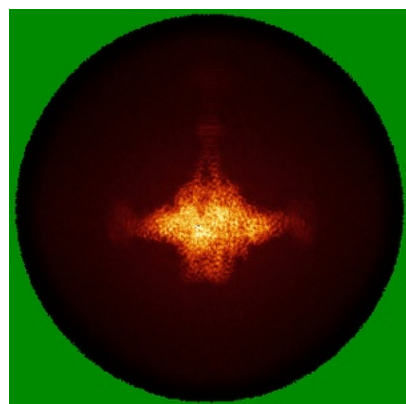


Z Index: 149

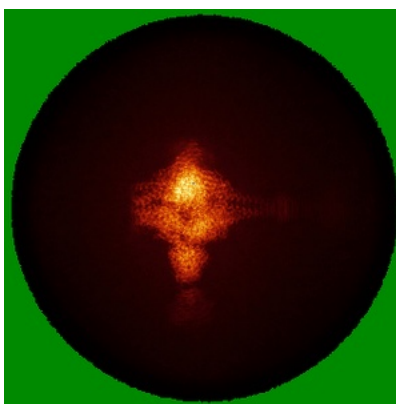
The images above show the largest variance slices of the map in three orthogonal directions.

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

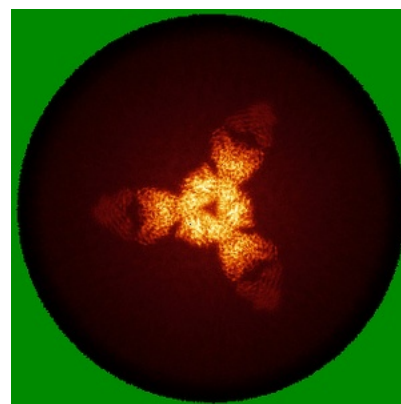
6.4.1 Primary map



X

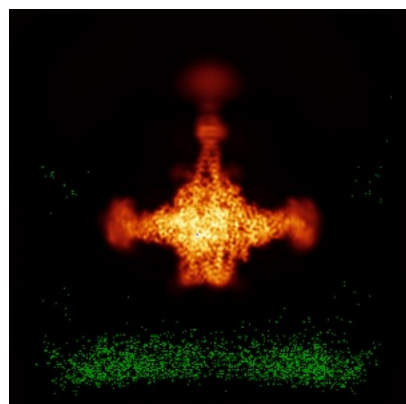


Y

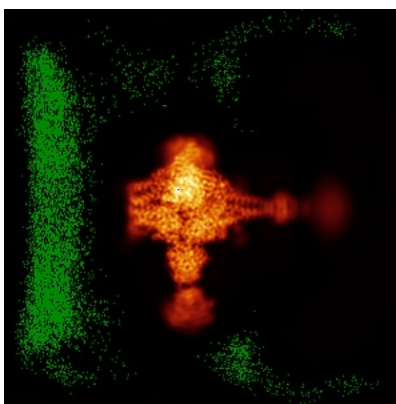


Z

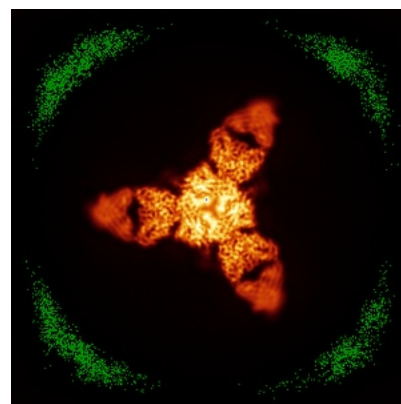
6.4.2 Raw map



X



Y

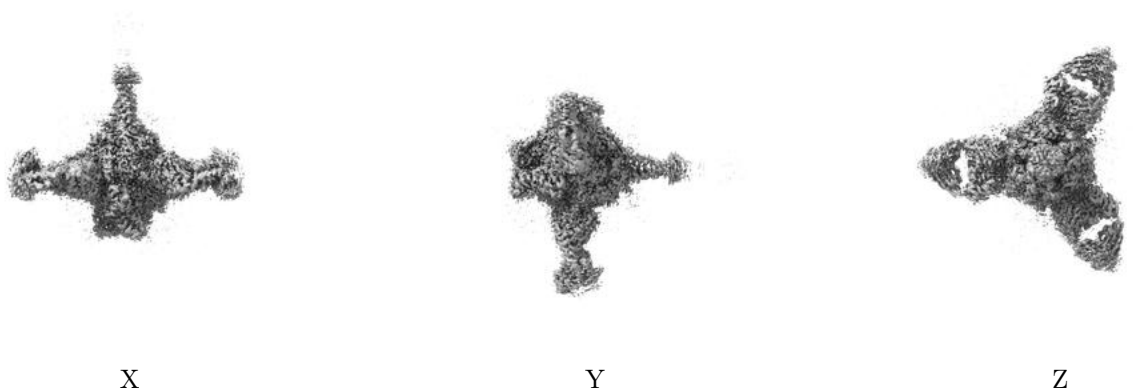


Z

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 0.414. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

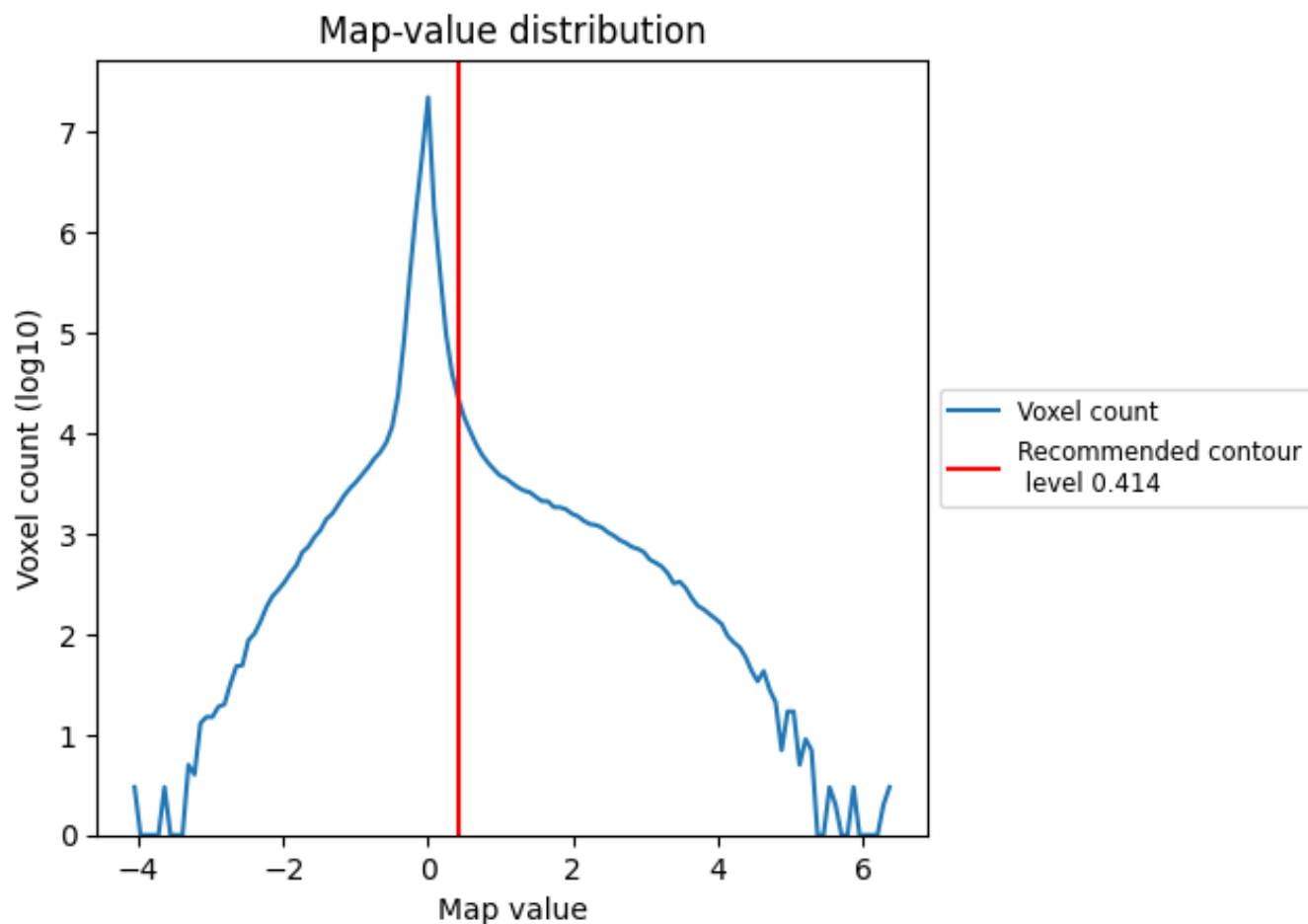
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

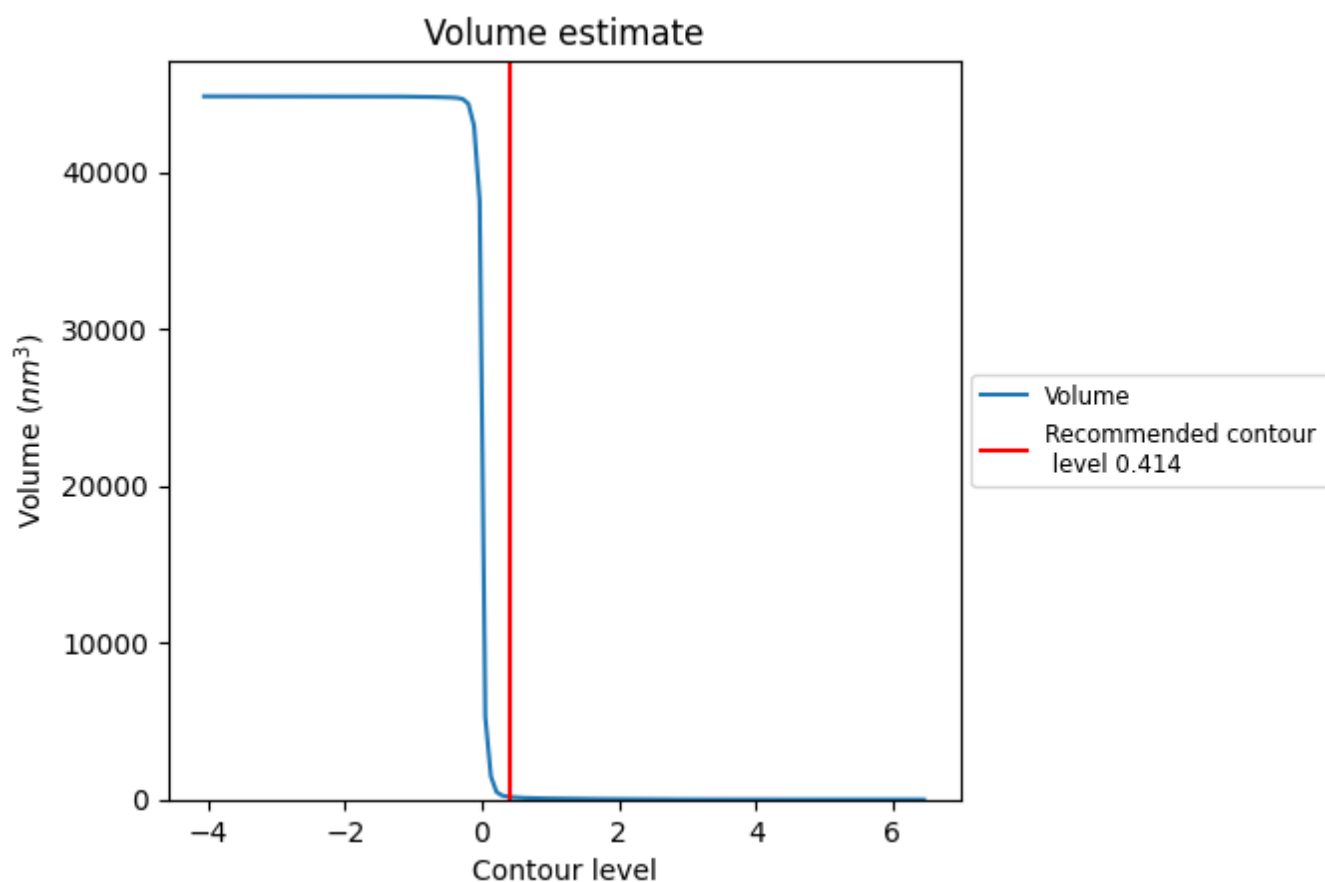
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

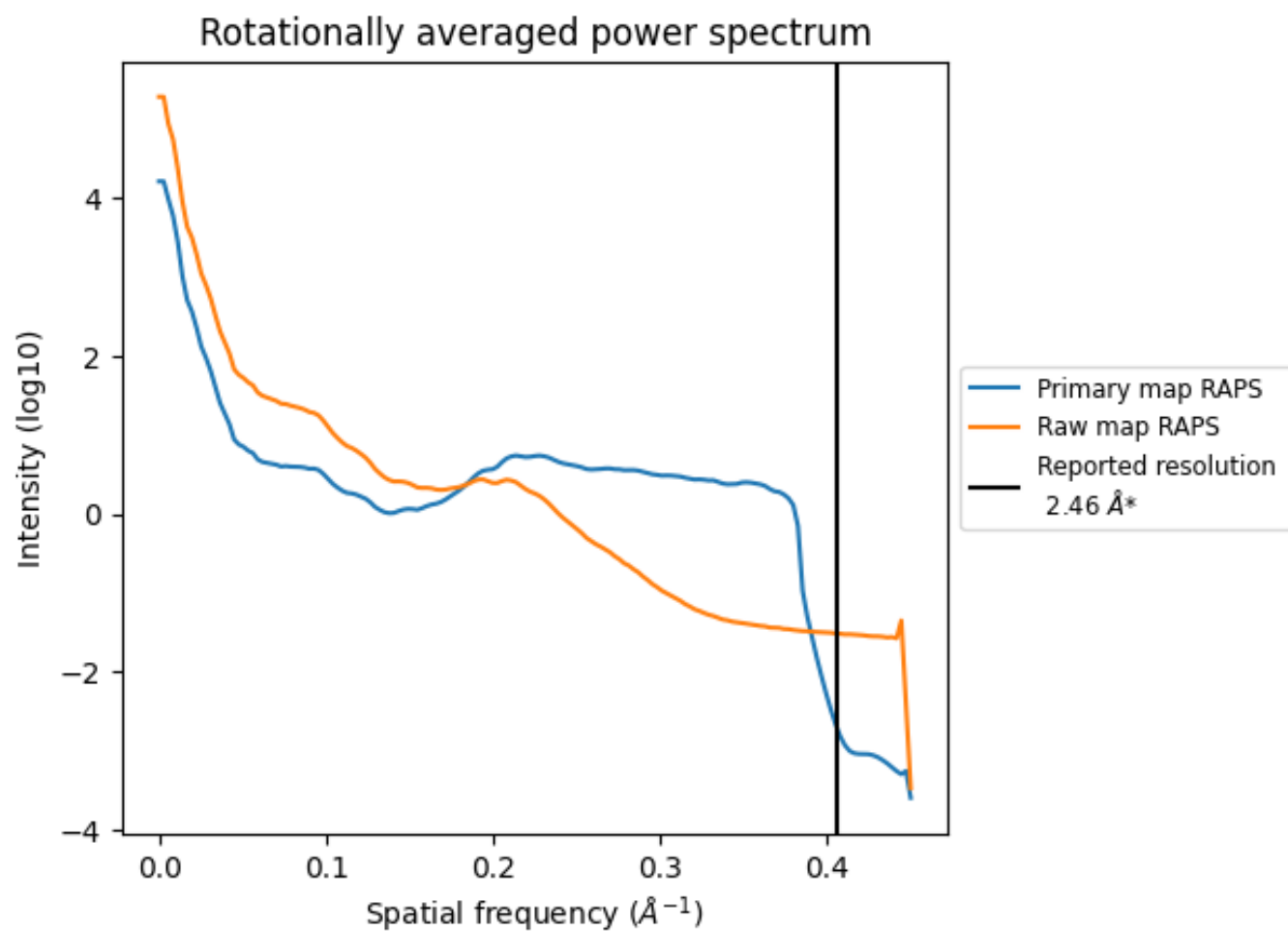
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 166 nm³; this corresponds to an approximate mass of 150 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 ⓘ

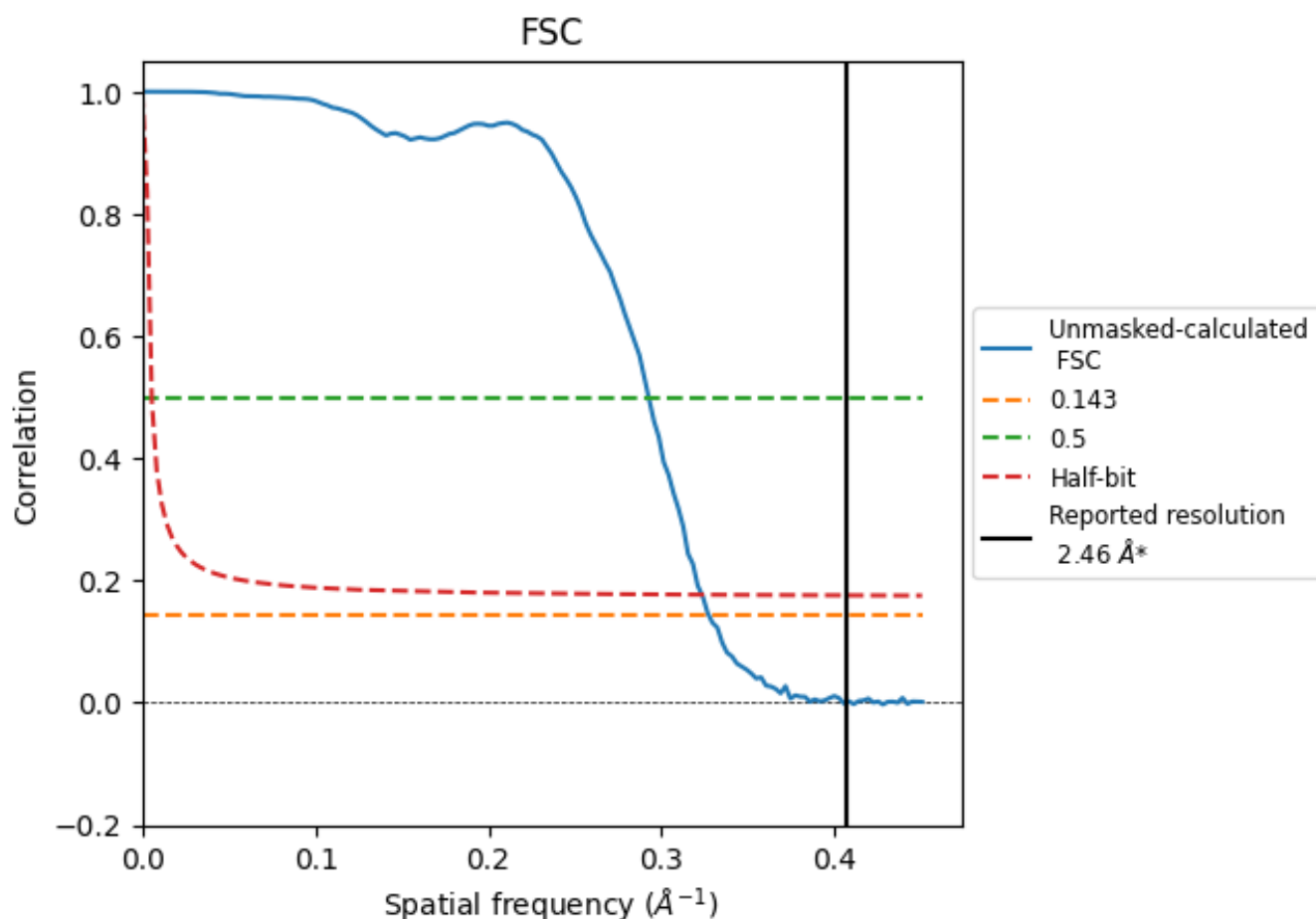


*Reported resolution corresponds to spatial frequency of 0.407 Å⁻¹

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 0.407 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.46	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.06	3.42	3.09

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.06 differs from the reported value 2.46 by more than 10 %

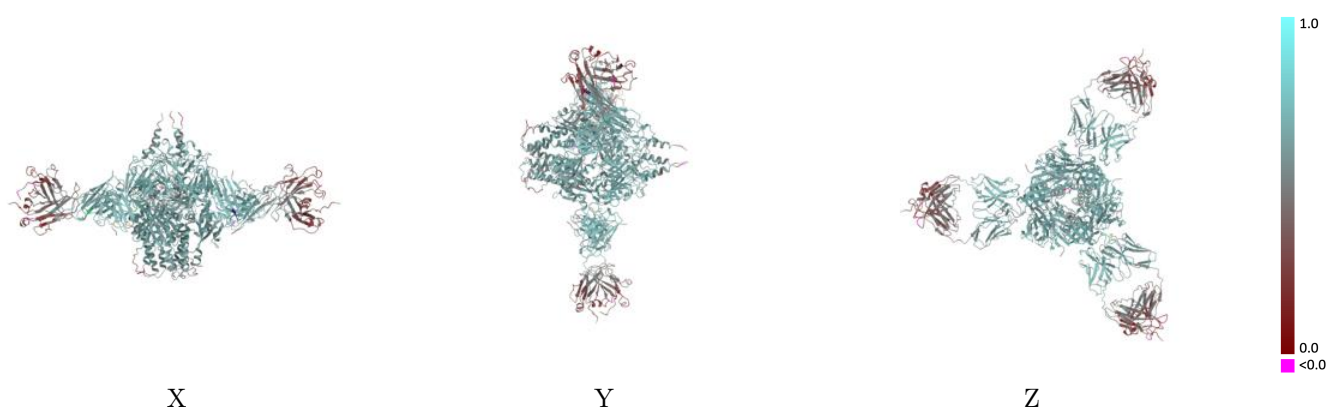
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-27808 and PDB model 8DZW. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

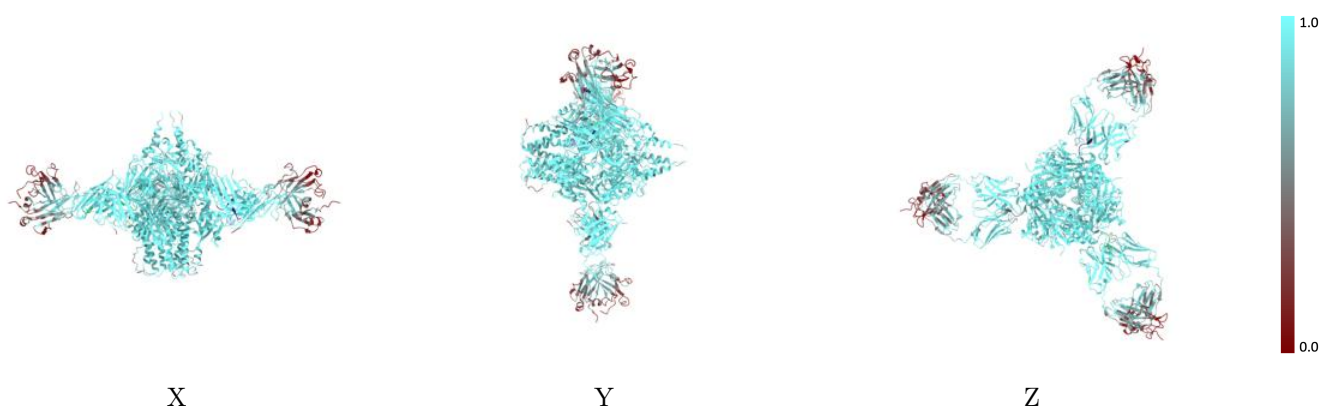
This section was not generated.

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



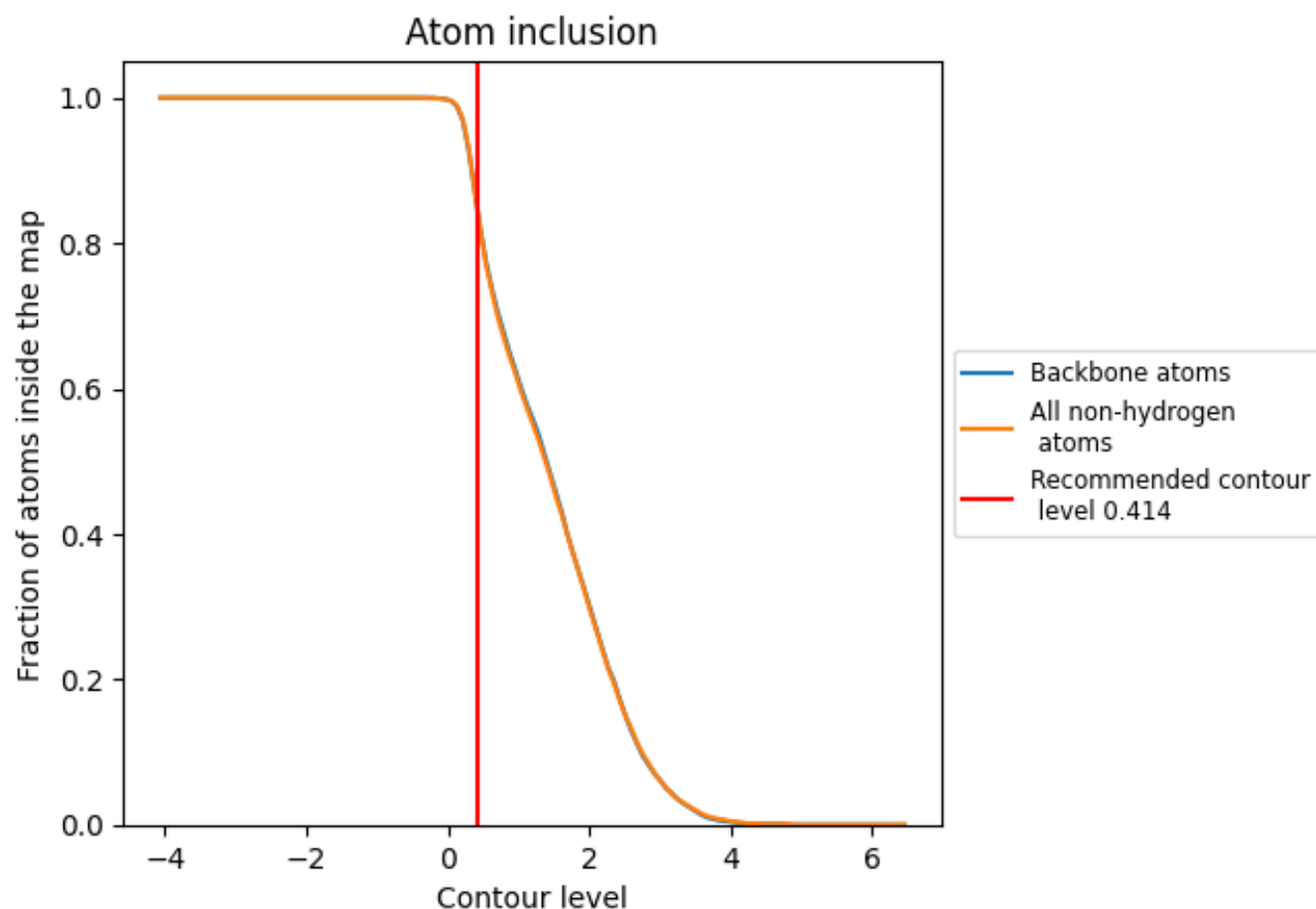
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.414).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 85% 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 (0.414) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8520	0.5690
A	0.9510	0.6240
B	0.9510	0.6250
C	0.9510	0.6260
H	0.7520	0.5300
I	0.7580	0.5320
J	0.7490	0.5290
L	0.7410	0.4850
M	0.7440	0.4940
N	0.7440	0.4920

1.0

0.0

<0.0