



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 04:07 PM UTC

PDB ID : 7XT4 / pdb_00007xt4
EMDB ID : EMD-33439
Title : Structure of Craspase-NTR
Authors : Feng, Y.; Zhang, L.
Deposited on : 2022-05-16
Resolution : 3.08 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:

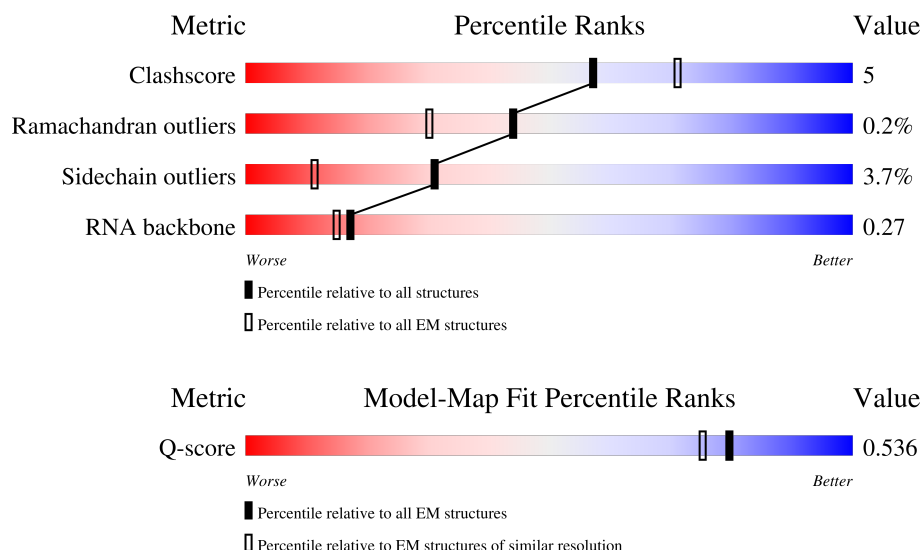
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.08 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	14000 (2.58 - 3.58)

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	C	716	
2	a	46	
3	B	1722	

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Mol	Chain	Length	Quality of chain
4	r	56	16%36%9%39%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 16850 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 CHAT domain protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	676	Total	C	N	O	S	0	0
			5482	3499	935	1023	25		

- Molecule 2 is a RNA chain called RNA (5'-R(P*GP*GP*GP*GP*CP*AP*GP*AP*AP*AP*AP*UP*UP*GP*GP*GP*UP*AP*CP*CP*G)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
2	a	21	Total	C	N	O	P	0	0
			459	204	90	144	21		

- Molecule 3 is a protein called RAMP superfamily protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	1263	Total	C	N	O	S	0	0
			10193	6485	1779	1897	32		

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	conflict	UNP A0A0B0EGF3
B	2	LYS	-	conflict	UNP A0A0B0EGF3
B	3	SER	-	conflict	UNP A0A0B0EGF3
B	4	ASN	-	conflict	UNP A0A0B0EGF3
B	5	ASP	-	conflict	UNP A0A0B0EGF3

- Molecule 4 is a RNA chain called RNA (34-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
4	r	34	Total	C	N	O	P	0	0
			712	319	118	241	34		

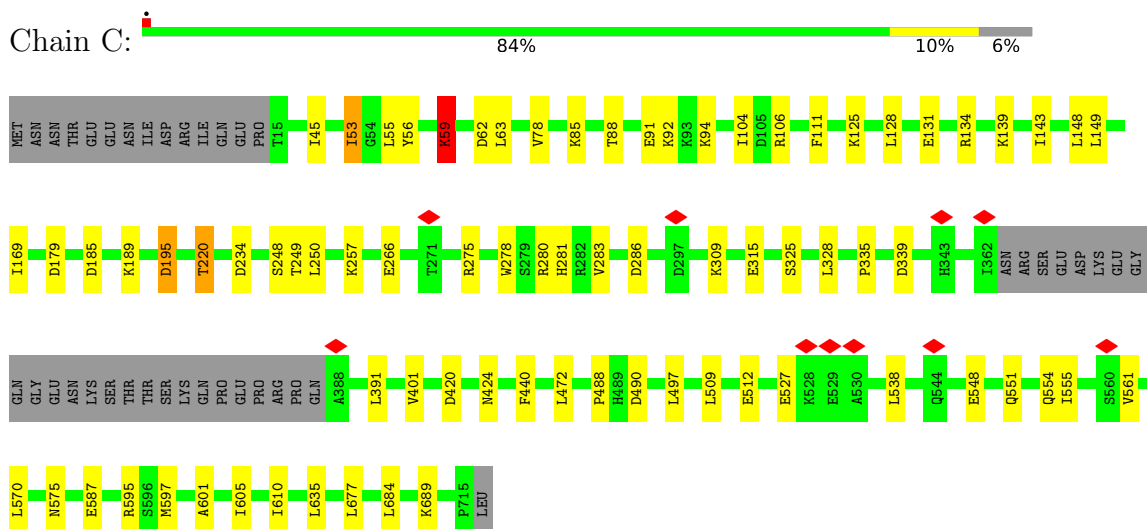
- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
5	B	4	Total	Zn	0
			4	4	

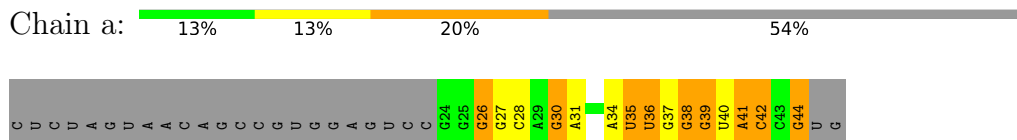
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

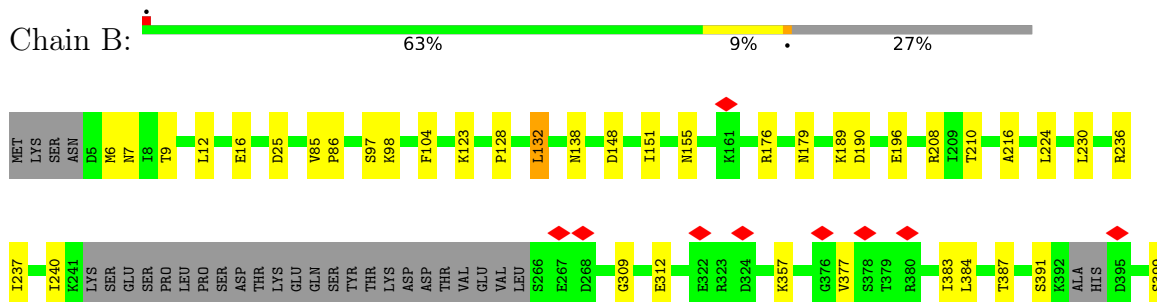
- Molecule 1: CHAT domain protein

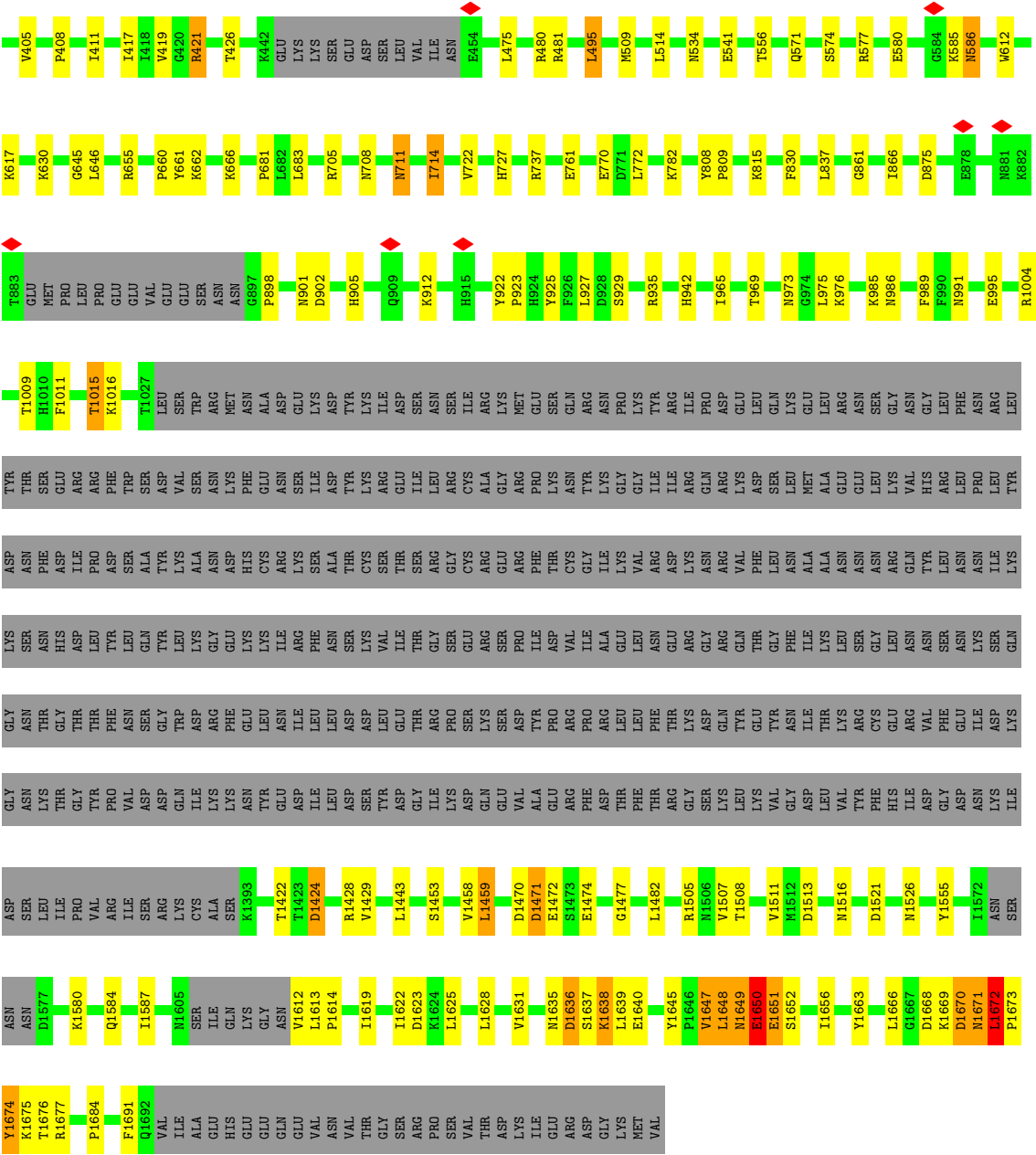


- Molecule 2: RNA (5'-R(P*GP*GP*GP*GP*CP*AP*GP*AP*AP*AP*UP*UP*GP*GP*GP*UP*AP*CP*CP*G)-3')



- Molecule 3: RAMP superfamily protein





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	40262	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)	1700	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.629	Depositor
Minimum map value	-0.069	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.037	Depositor
Recommended contour level	0.056	Depositor
Map size (Å)	271.6, 271.6, 271.6	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.97, 0.97, 0.97	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	C	0.27	0/5599	0.54	8/7548 (0.1%)
2	a	0.72	1/515 (0.2%)	0.80	5/803 (0.6%)
3	B	0.32	3/10431 (0.0%)	0.55	6/14055 (0.0%)
4	r	0.36	0/792	0.58	1/1229 (0.1%)
All	All	0.33	4/17337 (0.0%)	0.56	20/23635 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	B	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	a	37	G	O3'-P	-6.39	1.51	1.61
3	B	1672	LEU	CA-C	-5.82	1.46	1.52
3	B	1650	GLU	CA-C	-5.82	1.45	1.52
3	B	1674	TYR	CA-C	-5.34	1.45	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	a	38	G	C1'-C2'-O2'	-7.33	97.41	108.40
4	r	-13	U	C4'-C3'-O3'	-5.82	104.27	113.00
3	B	1650	GLU	N-CA-C	-5.76	98.52	110.80
3	B	1645	TYR	CA-C-N	-5.75	114.69	120.21
3	B	1645	TYR	C-N-CA	-5.75	114.69	120.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	111	PHE	CA-C-N	5.72	132.73	123.93
1	C	111	PHE	C-N-CA	5.72	132.73	123.93
3	B	1647	VAL	N-CA-C	-5.51	99.70	107.75
2	a	35	U	C2'-C3'-O3'	5.51	117.77	109.50
3	B	1638	LYS	N-CA-C	-5.42	105.82	113.20
2	a	40	U	C4'-C3'-O3'	-5.33	105.00	113.00
1	C	488	PRO	N-CA-C	5.29	120.44	111.03
1	C	325	SER	N-CA-C	5.23	117.03	110.91
1	C	601	ALA	CA-C-N	5.18	129.77	122.46
1	C	601	ALA	C-N-CA	5.18	129.77	122.46
1	C	59	LYS	CA-C-N	5.17	128.57	120.82
1	C	59	LYS	C-N-CA	5.17	128.57	120.82
2	a	41	A	C1'-C2'-O2'	-5.06	100.81	108.40
2	a	39	G	C1'-C2'-O2'	-5.05	100.83	108.40
3	B	1670	ASP	N-CA-C	-5.04	100.07	110.80

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	B	1470	ASP	Peptide
3	B	898	PRO	Peptide

5.2 Too-close contacts [i](#)

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	C	5482	0	5468	40	0
2	a	459	0	229	11	0
3	B	10193	0	10069	105	0
4	r	712	0	365	18	0
5	B	4	0	0	0	0
All	All	16850	0	16131	157	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 5.

All (157) 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)
3:B:1648:LEU:HD13	3:B:1663:TYR:HB2	1.64	0.79
2:a:39:G:H1'	3:B:761:GLU:OE1	1.83	0.77
1:C:59:LYS:HE3	1:C:63:LEU:HD13	1.78	0.66
3:B:985:LYS:HB2	3:B:1507:VAL:HG12	1.81	0.63
3:B:925:TYR:HB2	3:B:1628:LEU:HD21	1.80	0.62
2:a:30:G:H2'	2:a:31:A:H8	1.66	0.60
3:B:973:ASN:O	3:B:986:ASN:ND2	2.33	0.60
3:B:927:LEU:O	3:B:1477:GLY:HA3	2.03	0.59
3:B:965:ILE:HG12	3:B:1511:VAL:HG12	1.86	0.58
3:B:391:SER:HB3	3:B:495:LEU:HB3	1.86	0.57
3:B:711:ASN:N	3:B:711:ASN:OD1	2.38	0.57
1:C:56:TYR:HA	1:C:63:LEU:HD21	1.86	0.57
2:a:34:A:N1	4:r:5:U:H5	2.02	0.57
3:B:1672:LEU:HD23	3:B:1676:THR:HB	1.86	0.57
3:B:25:ASP:N	3:B:25:ASP:OD1	2.38	0.56
3:B:1650:GLU:O	3:B:1652:SER:N	2.38	0.56
2:a:28:C:H5	4:r:11:G:H1	1.53	0.55
4:r:-7:C:H2'	4:r:-4:G:N2	2.21	0.55
1:C:220:THR:HG21	1:C:257:LYS:HB3	1.89	0.55
3:B:1637:SER:C	3:B:1639:LEU:H	2.14	0.55
3:B:737:ARG:NH1	3:B:942:HIS:O	2.40	0.54
3:B:1636:ASP:O	3:B:1637:SER:C	2.50	0.54
1:C:266:GLU:OE2	1:C:275:ARG:NH1	2.38	0.54
2:a:26:G:OP2	3:B:1505:ARG:NH2	2.40	0.54
3:B:309:GLY:HA2	3:B:312:GLU:HB2	1.89	0.54
1:C:59:LYS:HG2	1:C:62:ASP:HB2	1.90	0.54
3:B:16:GLU:OE1	4:r:-5:C:N4	2.41	0.54
3:B:976:LYS:O	3:B:986:ASN:ND2	2.40	0.54
3:B:1424:ASP:OD1	3:B:1424:ASP:N	2.37	0.54
3:B:1637:SER:OG	3:B:1639:LEU:HB2	2.07	0.54
3:B:989:PHE:O	3:B:991:ASN:ND2	2.41	0.54
3:B:705:ARG:HH21	3:B:714:ILE:HD13	1.73	0.54
1:C:234:ASP:OD1	1:C:280:ARG:NH1	2.41	0.53
3:B:1668:ASP:N	3:B:1668:ASP:OD1	2.41	0.53
3:B:708:ASN:ND2	3:B:711:ASN:OD1	2.41	0.53
3:B:236:ARG:HH12	3:B:630:LYS:HE2	1.74	0.52
1:C:59:LYS:O	1:C:63:LEU:HD22	2.10	0.52
3:B:132:LEU:HD12	3:B:216:ALA:HB1	1.92	0.52
3:B:577:ARG:HD2	3:B:655:ARG:HG3	1.92	0.52
3:B:138:ASN:HB2	3:B:148:ASP:HB2	1.92	0.52
3:B:1650:GLU:C	3:B:1652:SER:N	2.65	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:440:PHE:HE2	1:C:635:LEU:HD21	1.75	0.51
3:B:1428:ARG:NH1	3:B:1526:ASN:O	2.44	0.51
3:B:97:SER:OG	3:B:98:LYS:NZ	2.41	0.51
1:C:195:ASP:N	1:C:195:ASP:OD1	2.44	0.51
3:B:875:ASP:N	3:B:875:ASP:OD1	2.44	0.51
3:B:1521:ASP:OD1	3:B:1521:ASP:N	2.42	0.51
2:a:42:C:O2	2:a:42:C:H2'	2.10	0.50
3:B:128:PRO:O	3:B:132:LEU:HB2	2.11	0.50
3:B:782:LYS:NZ	3:B:808:TYR:OH	2.43	0.50
3:B:727:HIS:NE2	3:B:770:GLU:O	2.45	0.50
3:B:190:ASP:N	3:B:190:ASP:OD1	2.45	0.49
3:B:1672:LEU:HG	3:B:1691:PHE:CE2	2.46	0.49
1:C:56:TYR:OH	1:C:106:ARG:NH1	2.46	0.49
1:C:548:GLU:HA	1:C:551:GLN:HE21	1.76	0.49
3:B:577:ARG:NH1	3:B:580:GLU:OE2	2.44	0.49
3:B:1672:LEU:CD2	3:B:1676:THR:HB	2.42	0.49
3:B:986:ASN:HA	3:B:1508:THR:O	2.13	0.49
1:C:179:ASP:OD1	1:C:179:ASP:N	2.42	0.48
3:B:1011:PHE:O	3:B:1015:THR:OG1	2.29	0.48
3:B:1650:GLU:O	3:B:1651:GLU:C	2.54	0.48
1:C:85:LYS:HA	1:C:328:LEU:HB3	1.95	0.48
3:B:1513:ASP:O	3:B:1516:ASN:ND2	2.41	0.48
3:B:1619:ILE:HD12	3:B:1622:ILE:HD12	1.95	0.48
3:B:722:VAL:HB	3:B:772:LEU:HB2	1.94	0.48
4:r:5:U:O2	4:r:5:U:H2'	2.14	0.48
1:C:605:ILE:HG12	1:C:610:ILE:HG13	1.94	0.48
1:C:88:THR:HG22	2:a:44:G:H5'	1.95	0.48
3:B:815:LYS:HE3	3:B:995:GLU:HG2	1.96	0.48
4:r:5:U:H1'	4:r:6:U:C6	2.49	0.48
3:B:176:ARG:NH1	3:B:196:GLU:OE2	2.47	0.47
3:B:681:PRO:O	3:B:861:GLY:N	2.44	0.47
3:B:1471:ASP:N	3:B:1471:ASP:OD1	2.47	0.47
3:B:925:TYR:HE2	3:B:1482:LEU:HD11	1.79	0.47
3:B:1649:ASN:C	3:B:1650:GLU:O	2.55	0.47
1:C:512:GLU:HA	1:C:689:LYS:HD3	1.95	0.47
3:B:1674:TYR:O	3:B:1675:LYS:C	2.56	0.47
2:a:31:A:N1	4:r:8:U:H5	2.13	0.47
3:B:176:ARG:NH2	4:r:-1:A:OP1	2.48	0.47
1:C:283:VAL:HA	1:C:286:ASP:HB2	1.95	0.47
1:C:149:LEU:HD22	1:C:169:ILE:HG23	1.97	0.46
3:B:662:LYS:HA	3:B:666:LYS:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:LYS:HG2	1:C:148:LEU:HD21	1.98	0.46
1:C:131:GLU:OE1	1:C:134:ARG:NH2	2.48	0.46
3:B:1650:GLU:CD	3:B:1651:GLU:H	2.23	0.46
1:C:339:ASP:N	1:C:339:ASP:OD1	2.48	0.46
4:r:5:U:H1'	4:r:6:U:C5	2.51	0.46
3:B:1639:LEU:HD12	3:B:1639:LEU:HA	1.87	0.46
1:C:555:ILE:HG21	1:C:677:LEU:HD13	1.98	0.46
3:B:1453:SER:O	3:B:1453:SER:OG	2.24	0.46
2:a:27:G:H1	4:r:12:C:H5	1.64	0.45
3:B:475:LEU:HB3	3:B:514:LEU:HD13	1.99	0.45
1:C:185:ASP:N	1:C:185:ASP:OD1	2.47	0.45
3:B:617:LYS:HD3	3:B:645:GLY:HA2	1.98	0.45
3:B:1580:LYS:HD3	3:B:1580:LYS:HA	1.79	0.45
3:B:421:ARG:HG2	3:B:556:THR:HG23	1.98	0.45
3:B:1650:GLU:C	3:B:1652:SER:H	2.23	0.45
1:C:587:GLU:HB3	1:C:597:MET:HG2	1.98	0.44
1:C:53:ILE:HD13	1:C:53:ILE:HA	1.87	0.44
1:C:509:LEU:HB3	1:C:512:GLU:HG3	1.97	0.44
1:C:575:ASN:OD1	1:C:575:ASN:N	2.48	0.44
3:B:1650:GLU:OE2	3:B:1651:GLU:N	2.50	0.44
1:C:391:LEU:HD21	1:C:527:GLU:HG2	1.99	0.44
3:B:912:LYS:HB2	3:B:912:LYS:HE2	1.75	0.44
3:B:683:LEU:HB2	3:B:809:PRO:HB3	1.99	0.44
4:r:-7:C:H2'	4:r:-4:G:C2	2.53	0.44
3:B:123:LYS:HB3	3:B:123:LYS:HE2	1.77	0.44
3:B:7:ASN:HD22	3:B:208:ARG:HH21	1.67	0.43
3:B:574:SER:HB3	3:B:661:TYR:HB2	2.01	0.43
3:B:927:LEU:HD11	3:B:1625:LEU:HB2	2.00	0.43
3:B:224:LEU:HD22	3:B:237:ILE:HG21	1.99	0.43
3:B:408:PRO:HG2	3:B:411:ILE:HB	2.00	0.43
3:B:612:TRP:HE3	3:B:646:LEU:HD13	1.83	0.43
1:C:91:GLU:HA	1:C:94:LYS:HD3	2.01	0.43
3:B:1671:ASN:OD1	3:B:1671:ASN:N	2.51	0.43
1:C:139:LYS:HA	1:C:139:LYS:HD3	1.78	0.43
3:B:179:ASN:O	4:r:-3:G:O2'	2.31	0.43
3:B:399:SER:O	3:B:399:SER:OG	2.35	0.43
3:B:905:HIS:NE2	3:B:1623:ASP:OD1	2.45	0.43
3:B:929:SER:HB3	3:B:1016:LYS:HE3	2.01	0.43
2:a:26:G:H1	4:r:13:C:H5	1.66	0.43
3:B:1459:LEU:HD13	3:B:1555:TYR:HE2	1.84	0.43
3:B:1628:LEU:HG	3:B:1684:PRO:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:420:ASP:O	1:C:424:ASN:HA	2.18	0.42
3:B:481:ARG:NH1	4:r:-2:U:O4	2.52	0.42
1:C:570:LEU:HD23	1:C:605:ILE:HG21	2.00	0.42
3:B:9:THR:HB	3:B:240:ILE:HG12	2.01	0.42
3:B:1672:LEU:HG	3:B:1691:PHE:CZ	2.54	0.42
2:a:36:U:H5	4:r:3:A:N1	2.17	0.42
3:B:480:ARG:HB2	3:B:509:MET:HE2	2.02	0.42
3:B:230:LEU:HD12	3:B:230:LEU:HA	1.87	0.42
3:B:1635:ASN:O	3:B:1637:SER:N	2.48	0.42
1:C:538:LEU:H	1:C:561:VAL:HG12	1.84	0.42
3:B:6:MET:HE2	3:B:6:MET:HB2	1.99	0.42
3:B:1422:THR:O	4:r:7:U:O2'	2.33	0.41
3:B:155:ASN:HD22	4:r:-14:U:H6	1.68	0.41
3:B:586:ASN:OD1	3:B:586:ASN:N	2.47	0.41
3:B:1672:LEU:HA	3:B:1673:PRO:HD3	1.84	0.41
3:B:387:THR:O	3:B:387:THR:OG1	2.38	0.41
3:B:901:ASN:O	3:B:902:ASP:C	2.63	0.41
3:B:1674:TYR:O	3:B:1677:ARG:N	2.52	0.41
1:C:335:PRO:HD3	1:C:595:ARG:HH11	1.85	0.41
3:B:1613:LEU:HA	3:B:1614:PRO:HD3	1.93	0.41
3:B:357:LYS:HE2	3:B:357:LYS:HB3	1.73	0.41
3:B:922:TYR:HA	3:B:923:PRO:HD3	1.95	0.41
1:C:189:LYS:HB3	1:C:189:LYS:HE2	1.84	0.41
1:C:78:VAL:HG21	3:B:384:LEU:HD21	2.03	0.41
3:B:571:GLN:HG3	3:B:660:PRO:HG3	2.03	0.41
3:B:585:LYS:HD2	3:B:655:ARG:HH11	1.86	0.41
1:C:59:LYS:HB2	1:C:59:LYS:HE2	1.84	0.40
1:C:278:TRP:CE2	1:C:281:HIS:HB2	2.57	0.40
3:B:104:PHE:O	4:r:-9:G:O2'	2.33	0.40
1:C:248:SER:OG	1:C:249:THR:N	2.54	0.40
1:C:309:LYS:HD3	1:C:309:LYS:HA	1.92	0.40
3:B:85:VAL:HA	3:B:86:PRO:HD3	1.89	0.40
3:B:151:ILE:HA	3:B:210:THR:O	2.21	0.40
3:B:534:ASN:HB2	3:B:541:GLU:HG2	2.04	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	C	672/716 (94%)	645 (96%)	26 (4%)	1 (0%)	48	76
3	B	1247/1722 (72%)	1175 (94%)	69 (6%)	3 (0%)	43	71
All	All	1919/2438 (79%)	1820 (95%)	95 (5%)	4 (0%)	44	71

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	B	1649	ASN
1	C	490	ASP
3	B	1471	ASP
3	B	1472	GLU

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	C	600/638 (94%)	583 (97%)	17 (3%)	38	63
3	B	1103/1525 (72%)	1057 (96%)	46 (4%)	26	55
All	All	1703/2163 (79%)	1640 (96%)	63 (4%)	31	58

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	45	ILE
1	C	53	ILE

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Mol	Chain	Res	Type
1	C	55	LEU
1	C	59	LYS
1	C	92	LYS
1	C	104	ILE
1	C	128	LEU
1	C	143	ILE
1	C	195	ASP
1	C	220	THR
1	C	250	LEU
1	C	315	GLU
1	C	401	VAL
1	C	472	LEU
1	C	497	LEU
1	C	554	GLN
1	C	684	LEU
3	B	12	LEU
3	B	132	LEU
3	B	189	LYS
3	B	377	VAL
3	B	383	ILE
3	B	405	VAL
3	B	417	ILE
3	B	419	VAL
3	B	421	ARG
3	B	426	THR
3	B	495	LEU
3	B	586	ASN
3	B	711	ASN
3	B	714	ILE
3	B	830	PHE
3	B	837	LEU
3	B	866	ILE
3	B	935	ARG
3	B	969	THR
3	B	975	LEU
3	B	1004	ARG
3	B	1009	THR
3	B	1015	THR
3	B	1424	ASP
3	B	1429	VAL
3	B	1443	LEU
3	B	1458	VAL

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Mol	Chain	Res	Type
3	B	1459	LEU
3	B	1474	GLU
3	B	1584	GLN
3	B	1587	ILE
3	B	1612	VAL
3	B	1631	VAL
3	B	1636	ASP
3	B	1638	LYS
3	B	1640	GLU
3	B	1647	VAL
3	B	1648	LEU
3	B	1650	GLU
3	B	1651	GLU
3	B	1656	ILE
3	B	1666	LEU
3	B	1669	LYS
3	B	1670	ASP
3	B	1671	ASN
3	B	1672	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	133	ASN
1	C	183	HIS
1	C	194	ASN
1	C	265	GLN
1	C	395	GLN
1	C	408	ASN
1	C	551	GLN
1	C	552	ASN
1	C	554	GLN
1	C	559	GLN
1	C	576	ASN
3	B	109	HIS
3	B	146	ASN
3	B	152	HIS
3	B	348	GLN
3	B	350	ASN
3	B	365	ASN
3	B	433	GLN
3	B	566	HIS

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Mol	Chain	Res	Type
3	B	712	ASN
3	B	942	HIS
3	B	980	ASN
3	B	991	ASN
3	B	1010	HIS
3	B	1604	ASN
3	B	1649	ASN
3	B	1692	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	a	20/46 (43%)	8 (40%)	0
4	r	33/56 (58%)	13 (39%)	0
All	All	53/102 (51%)	21 (39%)	0

All (21) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	a	26	G
2	a	30	G
2	a	35	U
2	a	36	U
2	a	38	G
2	a	41	A
2	a	42	C
2	a	44	G
4	r	-12	A
4	r	-11	A
4	r	-10	U
4	r	-9	G
4	r	-8	U
4	r	-7	C
4	r	-6	A
4	r	-4	G
4	r	-3	G
4	r	11	G
4	r	14	C
4	r	15	C
4	r	16	G

There are no RNA pucker outliers to report.

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 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-33439. These allow visual inspection of the internal detail of the map and identification of artifacts.

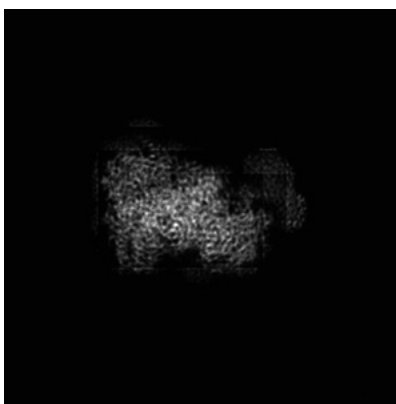
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

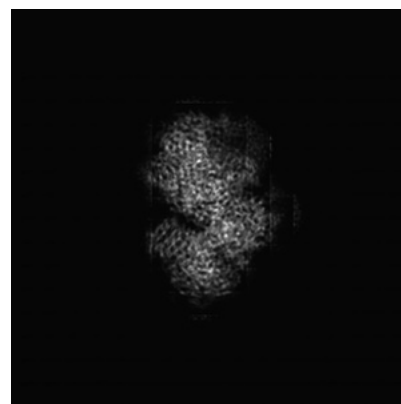
6.1.1 Primary map



X

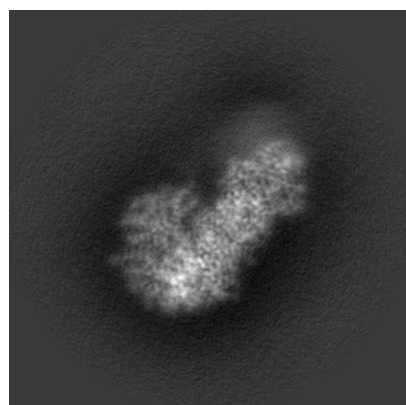


Y

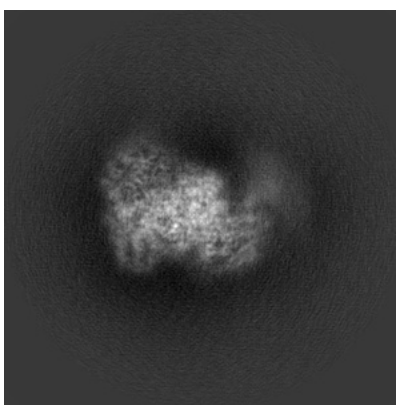


Z

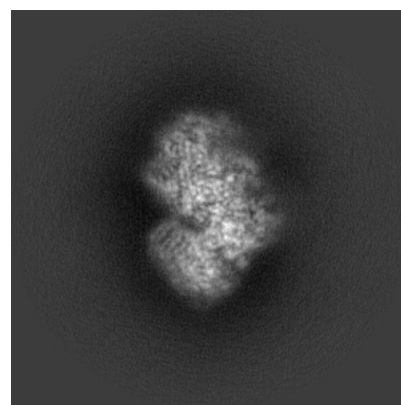
6.1.2 Raw map



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: 140

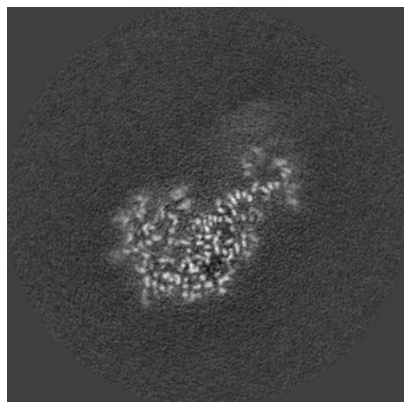


Y Index: 140

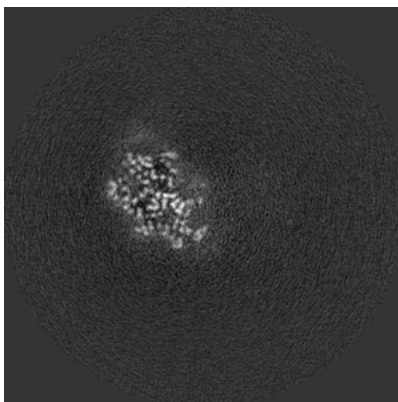


Z Index: 140

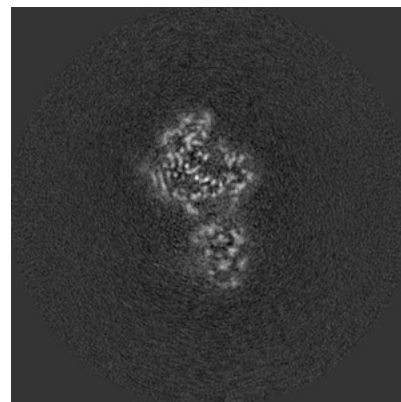
6.2.2 Raw map



X Index: 140



Y Index: 140



Z Index: 140

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: 132

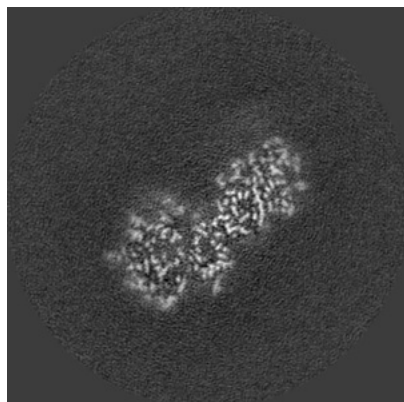


Y Index: 142

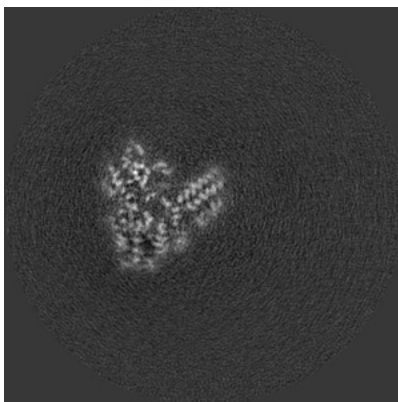


Z Index: 142

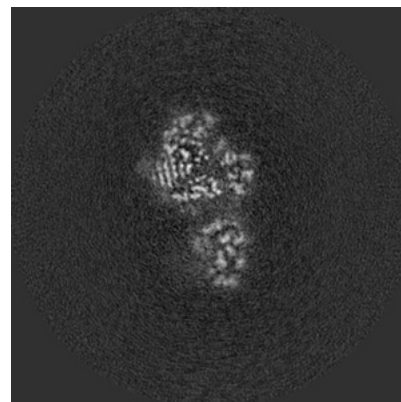
6.3.2 Raw map



X Index: 132



Y Index: 118

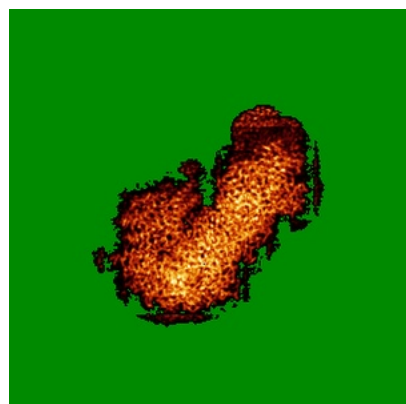


Z Index: 142

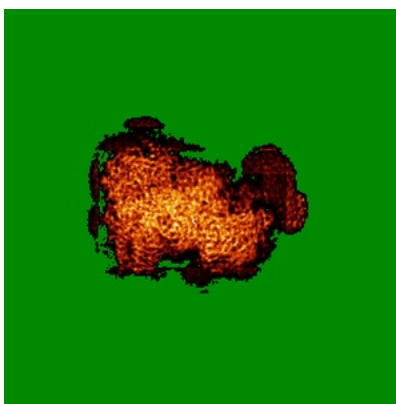
The images above show the largest variance slices of the map in three orthogonal directions.

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

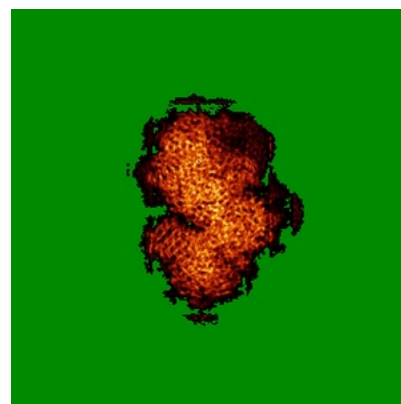
6.4.1 Primary map



X

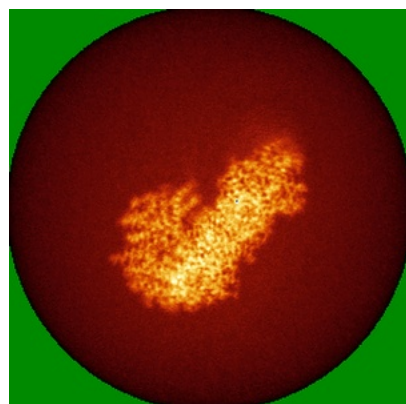


Y

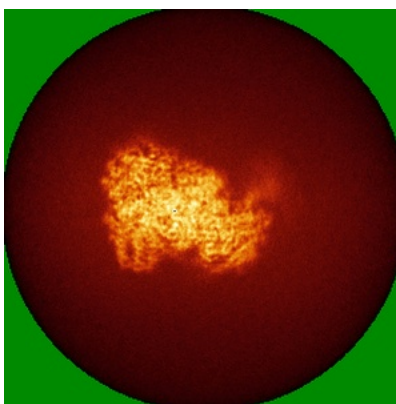


Z

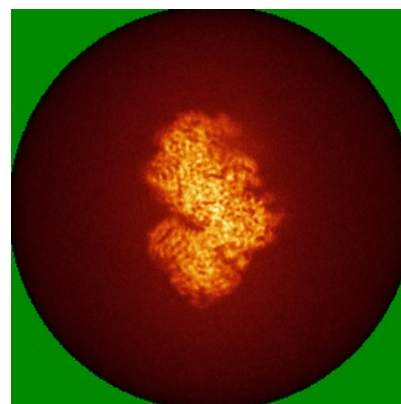
6.4.2 Raw map



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](#)

This section was not generated.

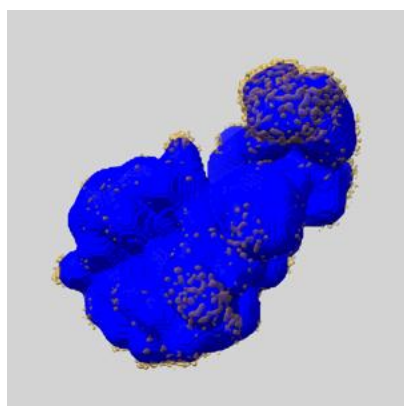
6.6 Mask visualisation [i](#)

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

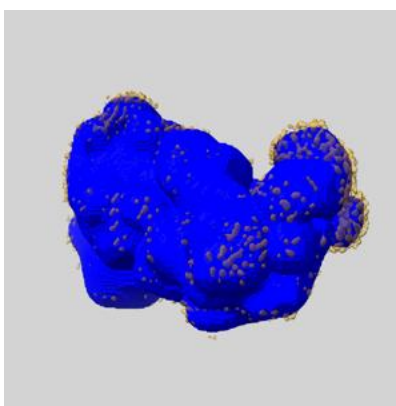
A mask typically either:

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

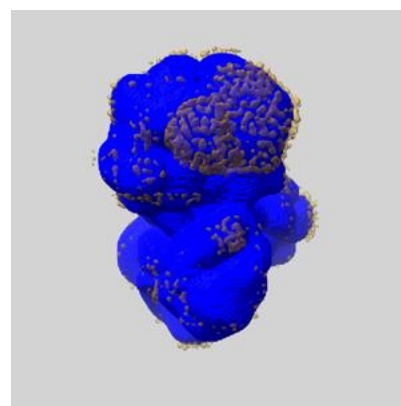
6.6.1 emd_33439_msk_1.map [i](#)



X



Y

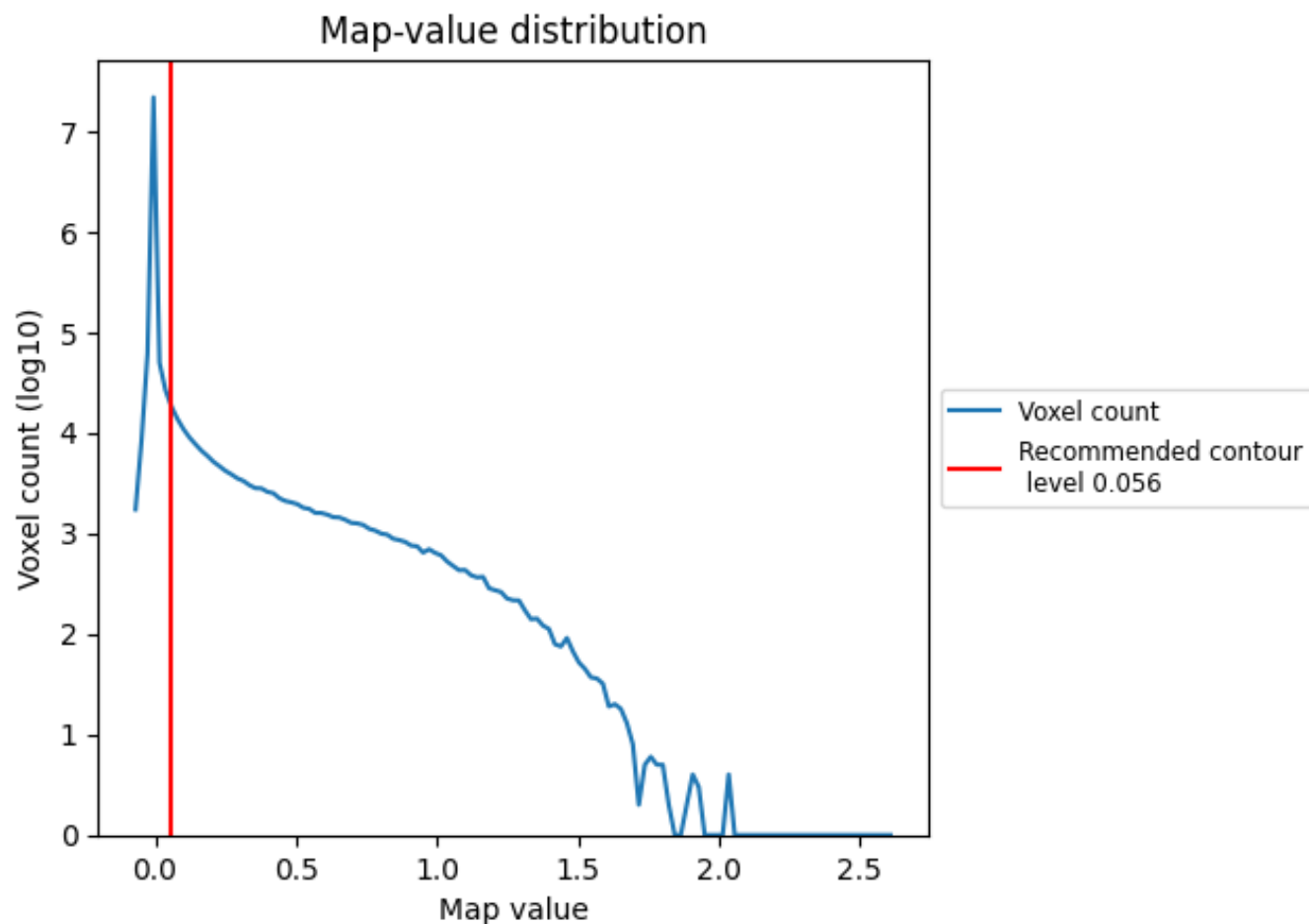


Z

7 Map analysis [i](#)

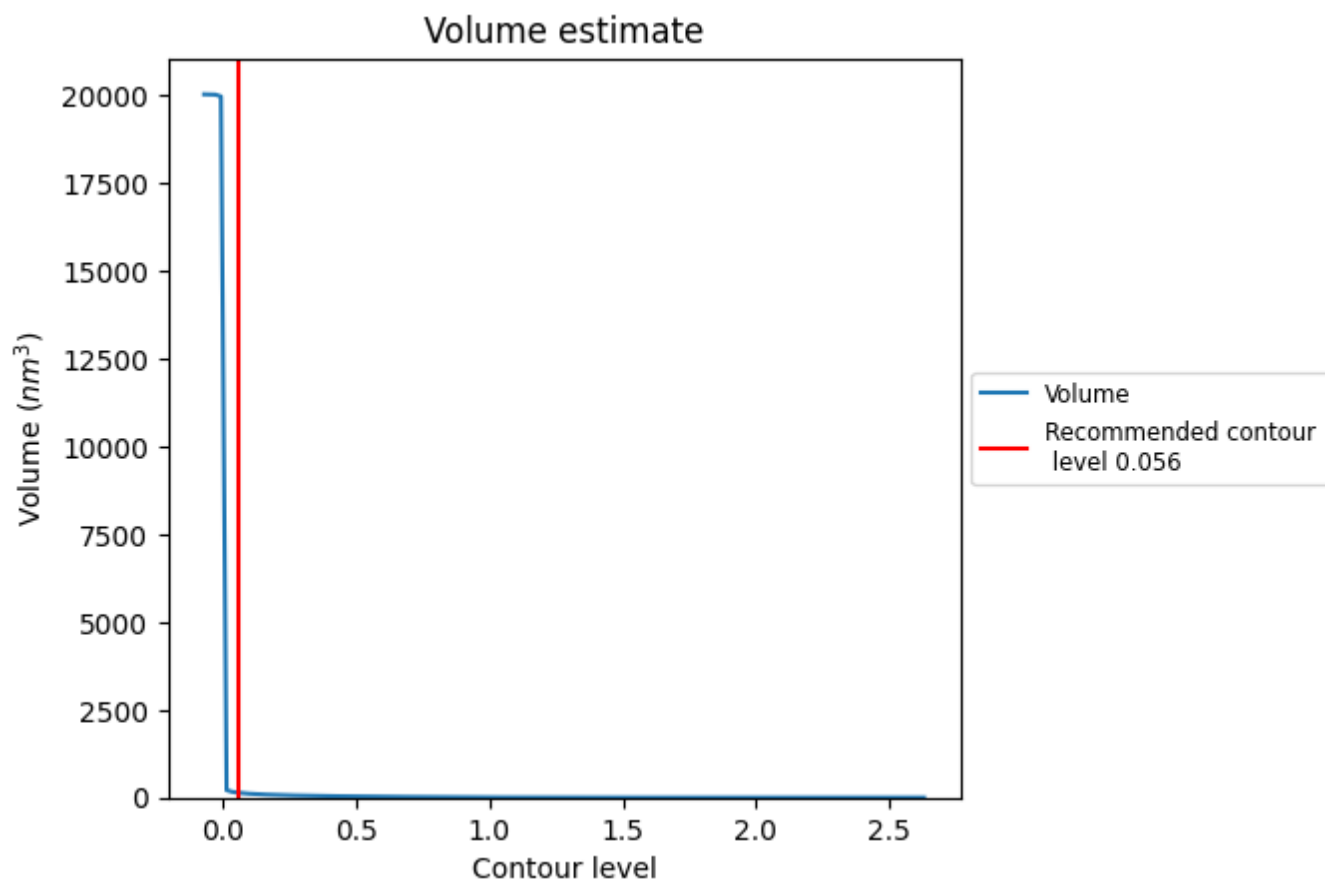
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

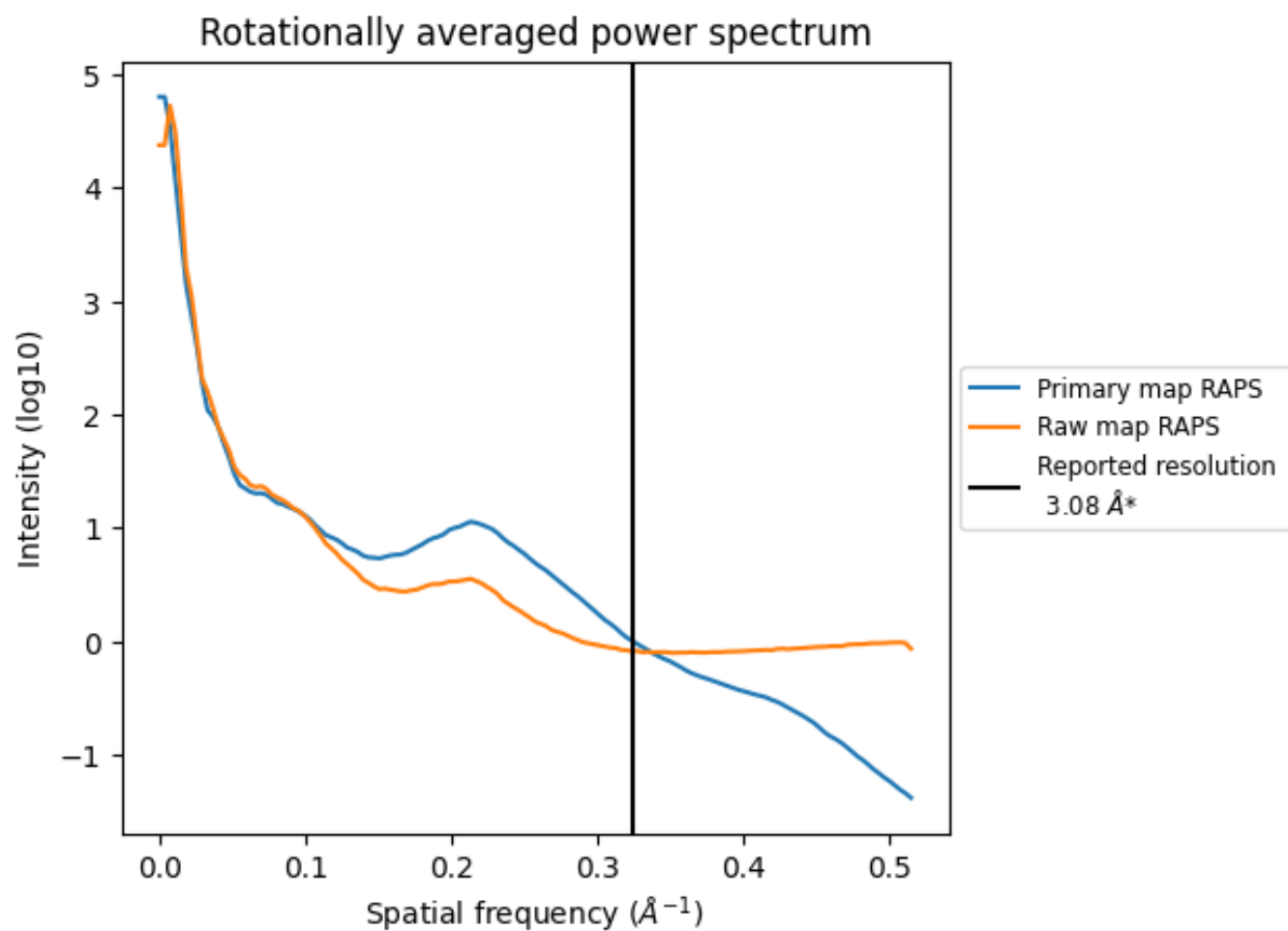
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 140 nm³; this corresponds to an approximate mass of 127 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 ⓘ

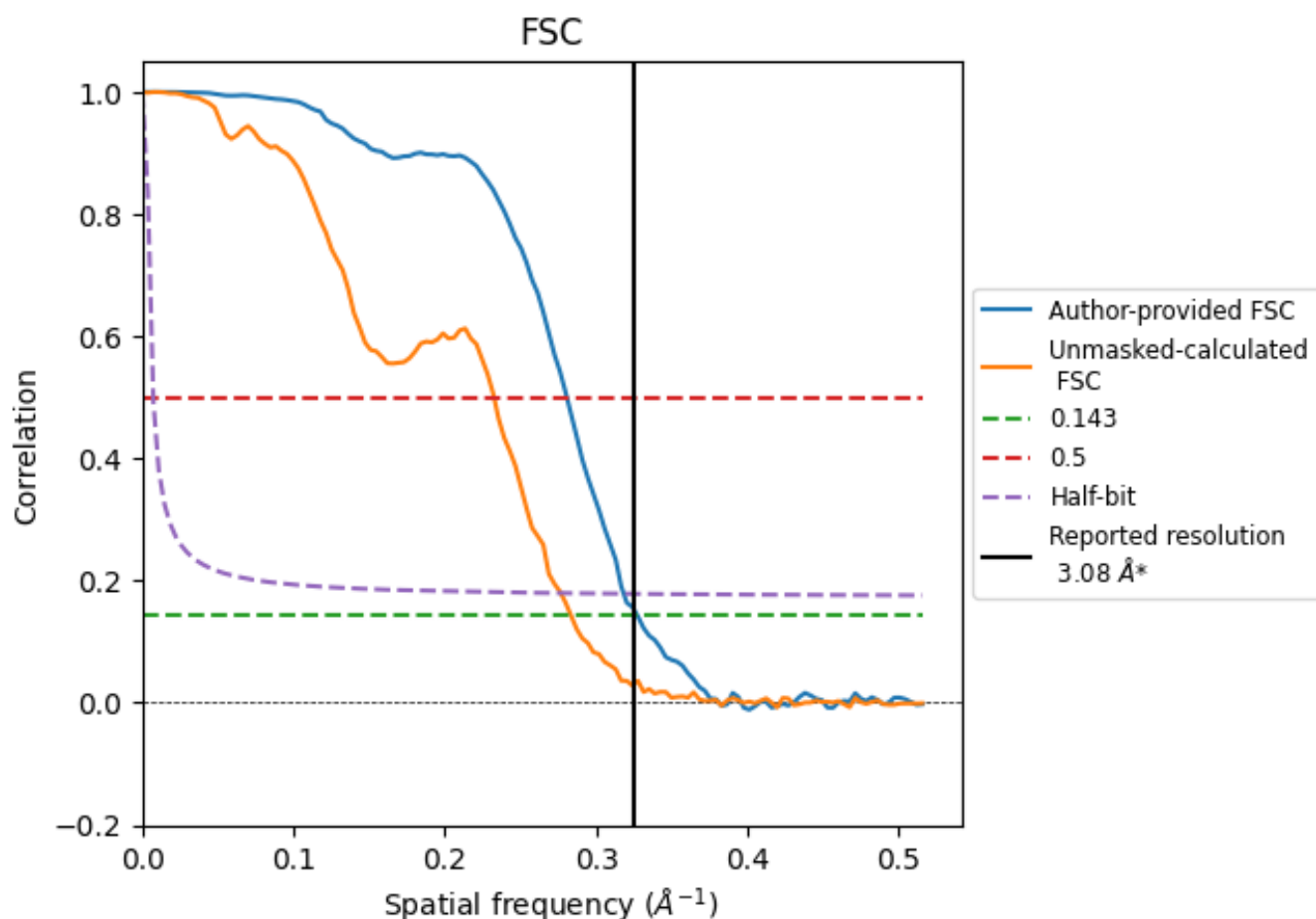


*Reported resolution corresponds to spatial frequency of 0.325 Å⁻¹

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 0.325 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

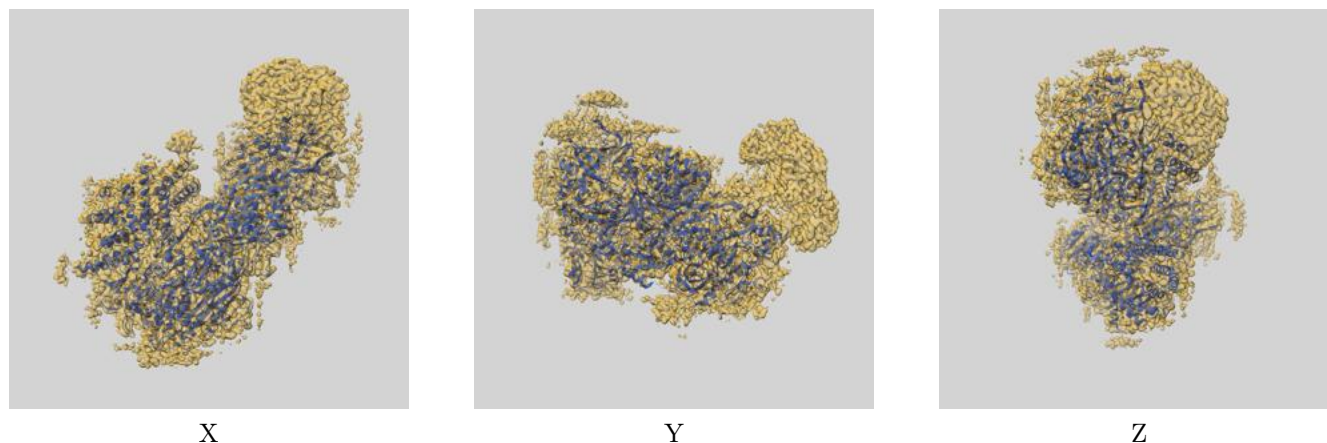
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.08	-	-
Author-provided FSC curve	3.06	3.56	3.14
Unmasked-calculated*	3.53	4.30	3.61

*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 3.53 differs from the reported value 3.08 by more than 10 %

9 Map-model fit [i](#)

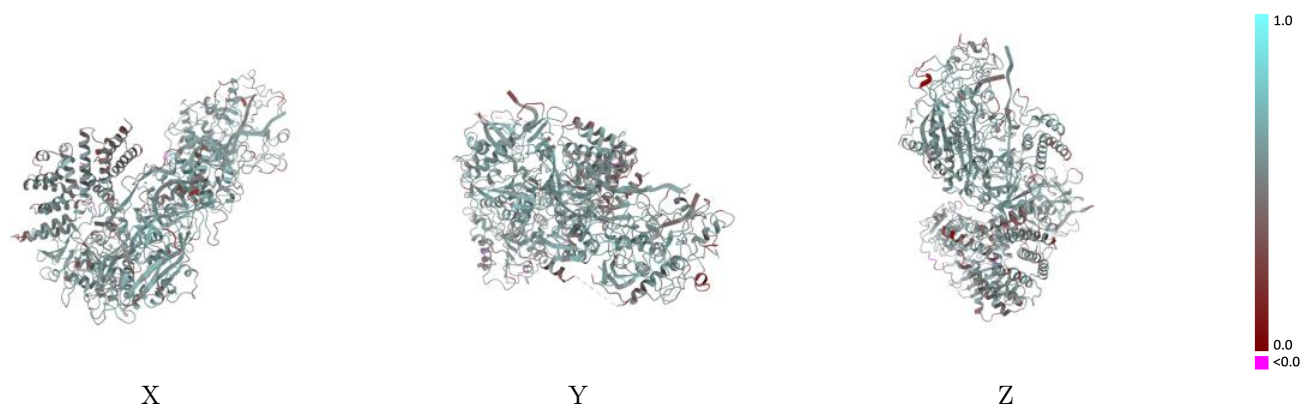
This section contains information regarding the fit between EMDB map EMD-33439 and PDB model 7XT4. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



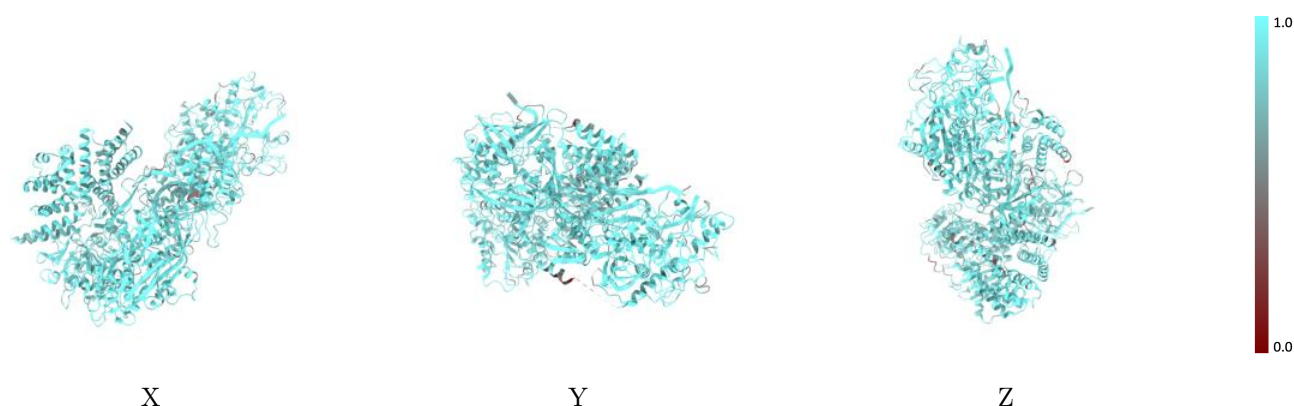
The images above show the 3D surface view of the map at the recommended contour level 0.056 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



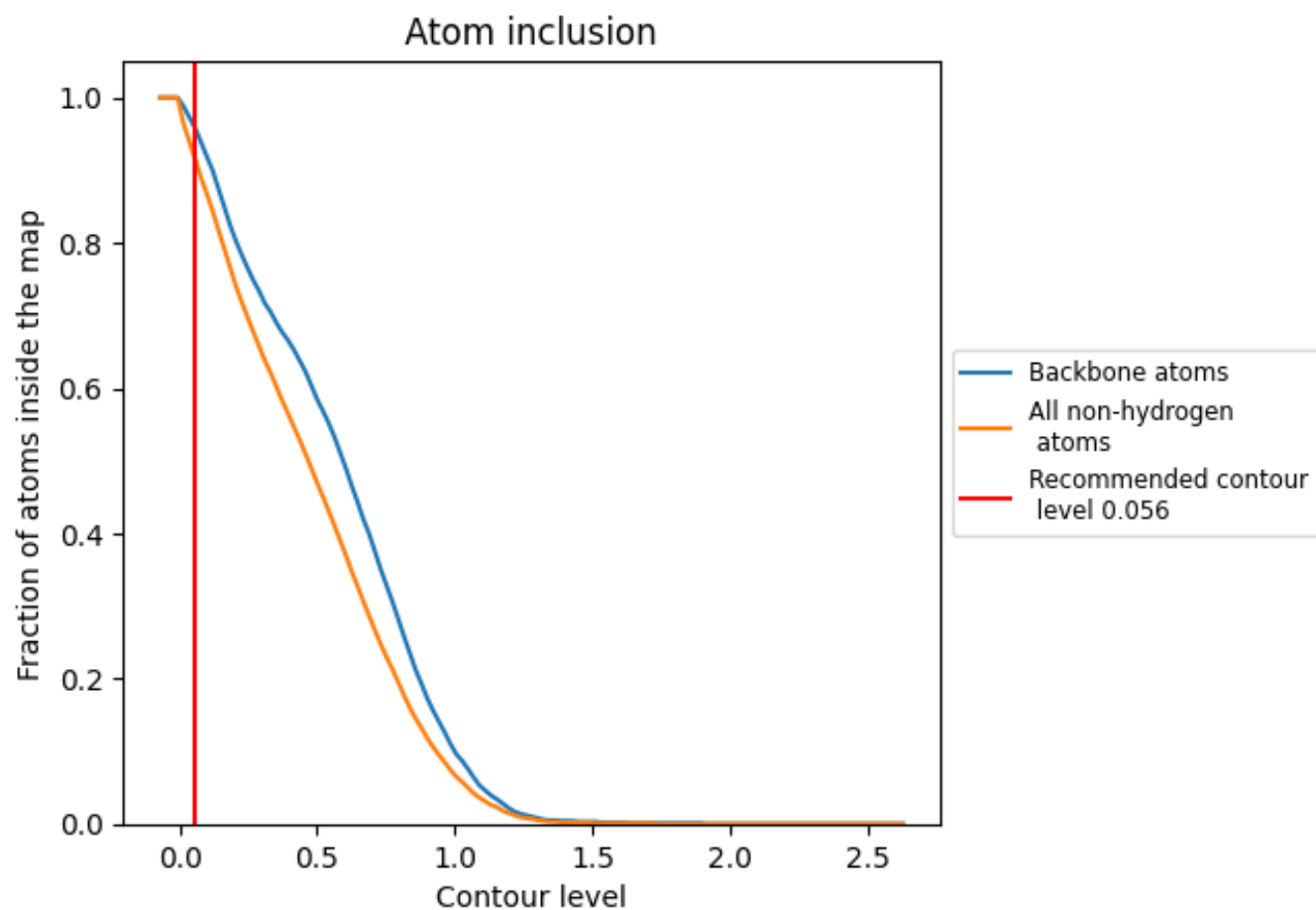
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.056).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 92% 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 (0.056) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9160	0.5360
B	0.9200	0.5500
C	0.8990	0.5050
a	0.9700	0.5830
r	0.9580	0.5580

