



Full wwPDB EM Validation Report ⓘ

Mar 24, 2026 – 12:11 PM UTC

PDB ID : 8GHE / pdb_00008ghe
EMDB ID : EMD-40042
Title : The structure of h12-LOX in tetrameric form bound to endogenous inhibitor oleoyl-CoA
Authors : Black, K.A.; Mobbs, J.I.; Venugopal, H.; Thal, D.M.; Glukhova, A.
Deposited on : 2023-03-09
Resolution : 2.05 Å(reported)
Based on initial model : .

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
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

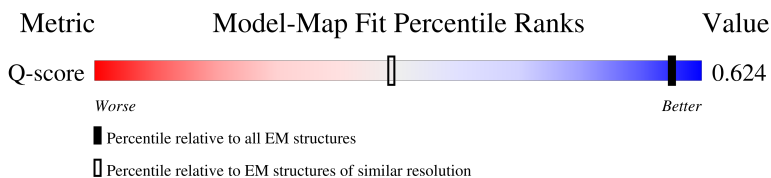
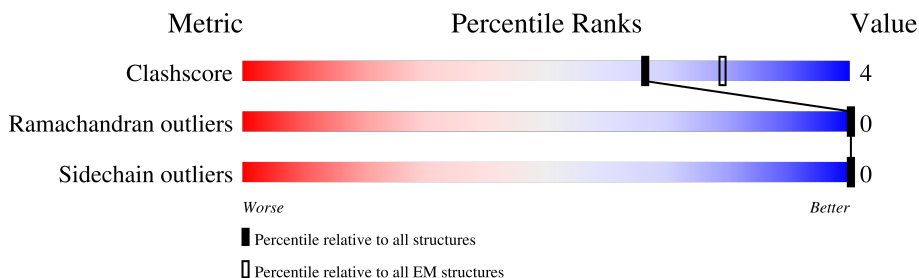
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.05 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	1849 (1.55 - 2.55)

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	669	 88% 11% .
1	B	669	 88% 11% .
1	C	669	 87% 12% .
1	D	669	 88% 11% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 22193 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 Polyunsaturated fatty acid lipoygenase ALOX12.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	662	Total	C	N	O	S	2	0
			5327	3402	930	961	34		
1	C	662	Total	C	N	O	S	1	0
			5319	3397	927	961	34		
1	D	662	Total	C	N	O	S	1	0
			5319	3397	927	961	34		
1	A	662	Total	C	N	O	S	2	0
			5327	3402	930	961	34		

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	MET	-	initiating methionine	UNP P18054
B	-4	HIS	-	expression tag	UNP P18054
B	-3	HIS	-	expression tag	UNP P18054
B	-2	HIS	-	expression tag	UNP P18054
B	-1	HIS	-	expression tag	UNP P18054
B	0	HIS	-	expression tag	UNP P18054
B	1	HIS	-	expression tag	UNP P18054
B	322	SER	ASN	variant	UNP P18054
C	-5	MET	-	initiating methionine	UNP P18054
C	-4	HIS	-	expression tag	UNP P18054
C	-3	HIS	-	expression tag	UNP P18054
C	-2	HIS	-	expression tag	UNP P18054
C	-1	HIS	-	expression tag	UNP P18054
C	0	HIS	-	expression tag	UNP P18054
C	1	HIS	-	expression tag	UNP P18054
C	322	SER	ASN	variant	UNP P18054
D	-5	MET	-	initiating methionine	UNP P18054
D	-4	HIS	-	expression tag	UNP P18054
D	-3	HIS	-	expression tag	UNP P18054
D	-2	HIS	-	expression tag	UNP P18054
D	-1	HIS	-	expression tag	UNP P18054
D	0	HIS	-	expression tag	UNP P18054

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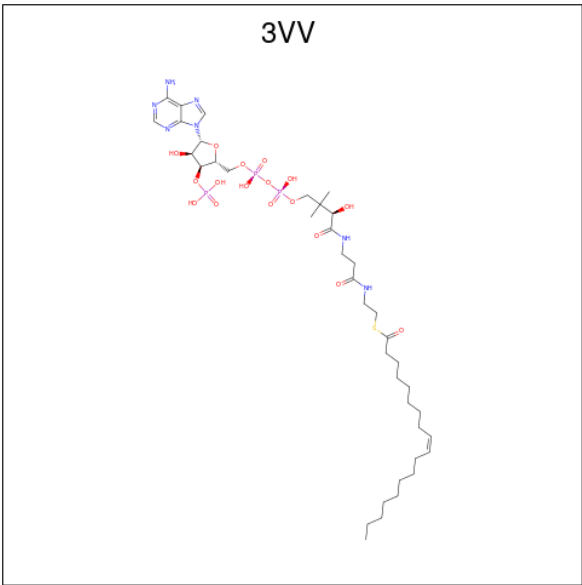
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Chain	Residue	Modelled	Actual	Comment	Reference
D	1	HIS	-	expression tag	UNP P18054
D	322	SER	ASN	variant	UNP P18054
A	-5	MET	-	initiating methionine	UNP P18054
A	-4	HIS	-	expression tag	UNP P18054
A	-3	HIS	-	expression tag	UNP P18054
A	-2	HIS	-	expression tag	UNP P18054
A	-1	HIS	-	expression tag	UNP P18054
A	0	HIS	-	expression tag	UNP P18054
A	1	HIS	-	expression tag	UNP P18054
A	322	SER	ASN	variant	UNP P18054

- Molecule 2 is FE (II) ION (CCD ID: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms	AltConf
2	B	1	Total Fe 1 1	0
2	C	1	Total Fe 1 1	0
2	D	1	Total Fe 1 1	0
2	A	1	Total Fe 1 1	0

- Molecule 3 is S-{(3R,5R,9R)-1-[(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-4-hydroxy-3-(phosphonooxy)tetrahydrofuran-2-yl]-3,5,9-trihydroxy-8,8-dimethyl-3,5-dioxido-10,14-dioxo-2,4,6-trioxa-11,15-diaza-3lambda 5 ,5lambda 5 -diphosphaheptadecan-17-yl} (9Z)-octadec-9-enethioate (non-preferred name) (CCD ID: 3VV) (formula: C₃₉H₆₈N₇O₁₇P₃S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						AltConf
			Total	C	N	O	P	S	
3	A	1	67	39	7	17	3	1	0

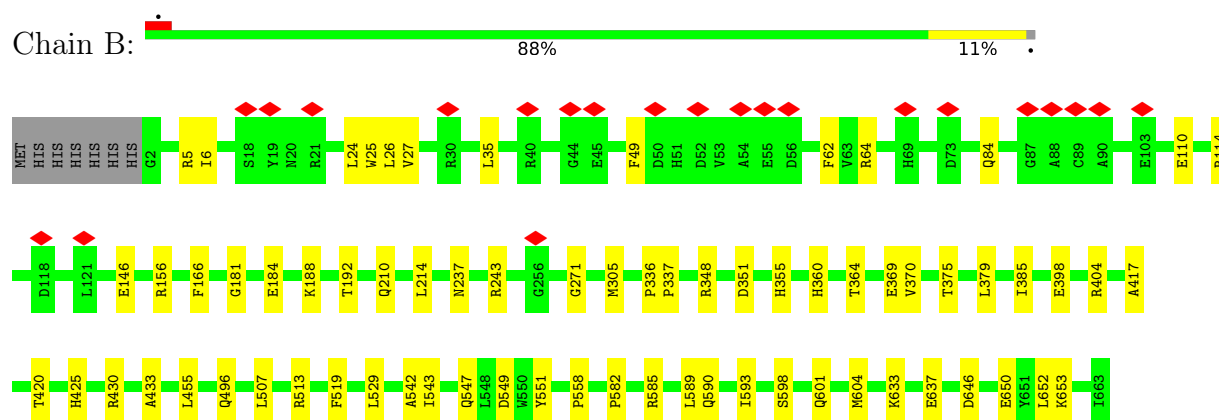
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		AltConf
4	B	197	Total	O	0
			197	197	
4	C	212	Total	O	0
			212	212	
4	D	198	Total	O	0
			198	198	
4	A	223	Total	O	0
			223	223	

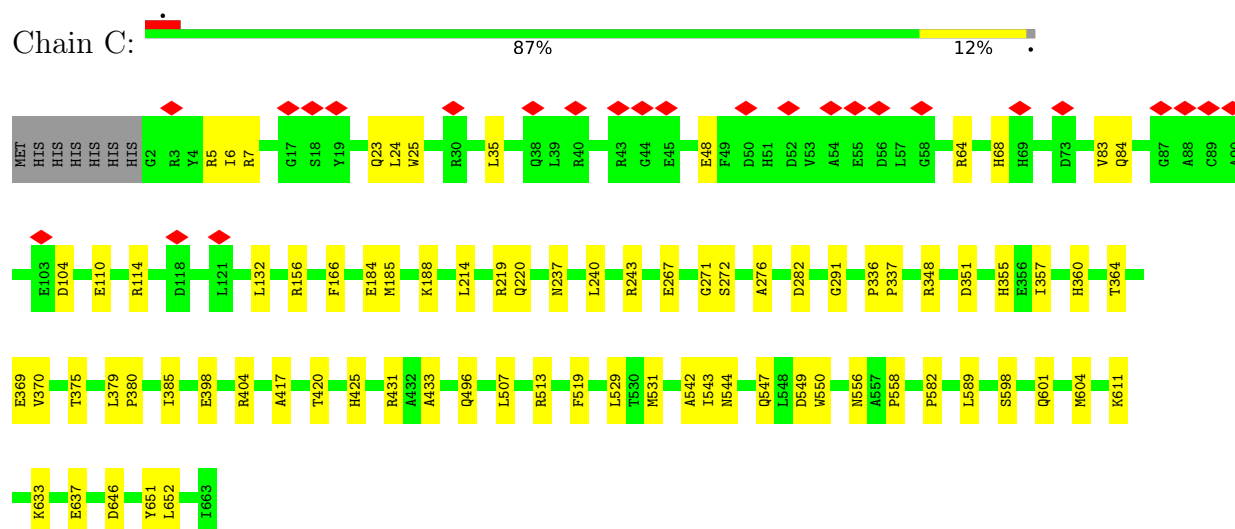
3 Residue-property plots

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.

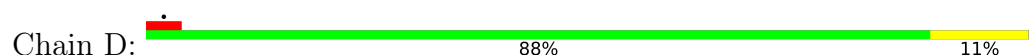
• Molecule 1: Polyunsaturated fatty acid lipoxygenase ALOX12

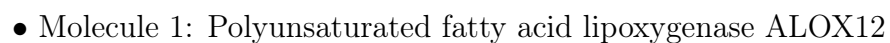


• Molecule 1: Polyunsaturated fatty acid lipoxygenase ALOX12



• Molecule 1: Polyunsaturated fatty acid lipoxygenase ALOX12





P443	P450	D463	W468	I471	E504	Q511	T530	A542	Q547	L548	W549	Y551	N556	A557	P558	P562	R565	L569	H596	Q601	M604	K633	E637	R641	M642	P648	L652	L663														
L148	L187	R188	R189	K211	L214	A215	E216	N237	R242	R243	Q271	N287	G291	P299	K304	M305	E306	Q317	R348	D351	F352	H355	H360	L361	T364	E369	L379	F386	R404	D415	K416	A417	T420	H425								
MET	HIS	HIS	HIS	HIS	HIS	G2	R3	Y4	H5	I6	R7	A8	Y9	L14	S18	Y19	W25	R30	R40	G44	E45	E48	F49	D50	H51	D52	V53	A54	F55	R64	D73	D74	R80	Q84	G87	A88	C89	A90	E103	D118	L121	K145

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	311579	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.009	Depositor
Minimum map value	-0.935	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.070	Depositor
Recommended contour level	0.35	Depositor
Map size (Å)	295.2, 295.2, 295.2	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.82000005, 0.82000005, 0.82000005	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: FE2, 3VV

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	0.23	0/5474	0.46	0/7438
1	B	0.18	0/5474	0.36	0/7438
1	C	0.17	0/5463	0.35	0/7424
1	D	0.17	0/5463	0.35	0/7424
All	All	0.19	0/21874	0.38	0/29724

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

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	585[A]	ARG	Sidechain
1	B	585[B]	ARG	Sidechain

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	A	5327	0	5266	42	0
1	B	5327	0	5266	38	0
1	C	5319	0	5253	46	0
1	D	5319	0	5253	42	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	67	0	64	7	0
4	A	223	0	0	1	0
4	B	197	0	0	0	0
4	C	212	0	0	0	0
4	D	198	0	0	0	0
All	All	22193	0	21102	165	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 4.

All (165) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:547:GLN:O	1:A:551:TYR:HB2	1.90	0.72
1:C:5:ARG:HB3	1:C:84:GLN:HB3	1.81	0.63
1:D:5:ARG:HB3	1:D:84:GLN:HB3	1.84	0.60
1:A:361:LEU:HB2	3:A:702:3VV:H1	1.83	0.59
1:A:7:ARG:HG2	1:A:48:GLU:HG2	1.84	0.59
1:B:181:GLY:HA2	1:B:210:GLN:HE22	1.71	0.56
1:B:5:ARG:HB3	1:B:84:GLN:HB3	1.87	0.56
1:D:184:GLU:HG2	1:D:188:LYS:HE3	1.89	0.55
1:B:156:ARG:NE	1:A:291:GLY:O	2.40	0.54
1:A:504:GLU:OE1	1:A:511:GLN:NE2	2.39	0.54
1:A:299:PRO:HA	1:A:317:GLN:O	2.08	0.53
1:B:355:HIS:NE2	1:B:543:ILE:O	2.38	0.53
1:C:369:GLU:OE2	1:C:404:ARG:NH2	2.42	0.53
1:B:6:ILE:HB	1:B:49:PHE:HB2	1.91	0.52
1:B:417:ALA:HB2	1:B:589:LEU:HD23	1.92	0.52
1:C:184:GLU:HG2	1:C:188:LYS:HE3	1.92	0.51
1:A:585:ARG:NH2	3:A:702:3VV:O55	2.43	0.51
1:D:360:HIS:O	1:D:364:THR:OG1	2.29	0.51
1:A:189[B]:ARG:NH1	4:A:836:HOH:O	2.44	0.50
1:D:348:ARG:NH1	1:D:351:ASP:OD2	2.44	0.50
1:D:355:HIS:NE2	1:D:543:ILE:O	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:HIS:O	1:A:364:THR:OG1	2.26	0.50
1:A:417:ALA:HB2	1:A:589:LEU:HD23	1.92	0.50
1:B:369:GLU:OE2	1:B:404:ARG:NH2	2.43	0.50
1:B:375:THR:HG23	1:B:379:LEU:HD12	1.93	0.50
1:C:348:ARG:NH1	1:C:351:ASP:OD2	2.45	0.50
1:C:360:HIS:O	1:C:364:THR:OG1	2.29	0.50
1:C:433:ALA:O	1:C:513:ARG:NH1	2.45	0.49
1:B:25:TRP:HB2	1:B:64:ARG:HB2	1.93	0.49
1:A:633:LYS:NZ	1:A:637:GLU:OE2	2.45	0.49
1:A:352:PHE:HZ	3:A:702:3VV:H22	1.77	0.49
1:D:243:ARG:NH1	1:D:271:GLY:O	2.45	0.49
1:B:184:GLU:HG2	1:B:188:LYS:HE3	1.94	0.48
1:C:375:THR:HG23	1:C:379:LEU:HD12	1.95	0.48
1:D:433:ALA:O	1:D:513:ARG:NH1	2.47	0.48
1:A:369:GLU:OE2	1:A:404:ARG:NH2	2.47	0.48
1:B:6:ILE:HD11	1:B:26:LEU:HD11	1.95	0.48
1:D:219:ARG:NH1	1:D:220:GLN:OE1	2.47	0.48
1:B:214:LEU:HD23	1:B:549:ASP:HB3	1.96	0.47
1:B:348:ARG:NH1	1:B:351:ASP:OD2	2.47	0.47
1:C:25:TRP:HB2	1:C:64:ARG:HB2	1.94	0.47
1:B:633:LYS:NZ	1:B:637:GLU:OE2	2.47	0.47
1:D:496:GLN:HE22	1:D:519:PHE:HB2	1.78	0.47
1:C:104:ASP:N	1:C:104:ASP:OD1	2.48	0.47
1:B:542:ALA:HB2	1:B:652:LEU:HD21	1.96	0.47
1:B:601:GLN:HB2	1:B:604:MET:HB3	1.97	0.47
1:D:633:LYS:NZ	1:D:637:GLU:OE2	2.47	0.47
1:D:369:GLU:OE2	1:D:404:ARG:NH2	2.45	0.47
1:B:420:THR:O	1:B:425:HIS:N	2.48	0.47
1:C:237:ASN:HB2	1:C:355:HIS:HB2	1.97	0.47
1:C:633:LYS:NZ	1:C:637:GLU:OE2	2.48	0.47
1:D:110:GLU:OE1	1:D:114:ARG:NH2	2.41	0.47
1:B:370:VAL:HG11	1:B:507:LEU:HG	1.98	0.46
1:C:110:GLU:OE1	1:C:114:ARG:NH2	2.42	0.46
1:A:416:LYS:HG3	3:A:702:3VV:H59	1.96	0.46
1:B:146:GLU:O	1:B:430:ARG:NE	2.43	0.46
1:C:156:ARG:NE	1:D:291:GLY:O	2.46	0.46
1:C:355:HIS:NE2	1:C:543:ILE:O	2.42	0.46
1:D:25:TRP:HB2	1:D:64:ARG:HB2	1.97	0.46
1:D:375:THR:HG23	1:D:379:LEU:HD12	1.97	0.46
1:B:646:ASP:N	1:B:646:ASP:OD1	2.49	0.46
1:C:6:ILE:HG13	1:C:83:VAL:HG13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:237:ASN:HB2	1:D:355:HIS:HB2	1.98	0.46
1:C:420:THR:O	1:C:425:HIS:N	2.49	0.46
1:D:544:ASN:O	1:D:547:GLN:NE2	2.48	0.46
1:C:542:ALA:HB2	1:C:652:LEU:HD21	1.98	0.46
1:D:542:ALA:HB2	1:D:652:LEU:HD21	1.98	0.46
1:A:287:ASN:ND2	1:A:415:ASP:O	2.43	0.46
1:A:642:ASN:HB3	1:A:648:PRO:HB3	1.98	0.46
1:B:385:ILE:HD12	1:B:529:LEU:HD11	1.98	0.45
1:B:433:ALA:O	1:B:513:ARG:NH1	2.49	0.45
1:B:305:MET:HE3	1:B:455:LEU:HD22	1.99	0.45
1:B:360:HIS:O	1:B:364:THR:OG1	2.31	0.45
1:D:214:LEU:HD23	1:D:549:ASP:HB3	1.99	0.45
1:D:385:ILE:HD12	1:D:529:LEU:HD11	1.97	0.45
1:A:304:LYS:NZ	1:A:306:GLU:OE2	2.47	0.45
1:B:110:GLU:OE1	1:B:114:ARG:NH2	2.42	0.45
1:C:214:LEU:HD23	1:C:549:ASP:HB3	1.98	0.45
1:A:379:LEU:HB3	1:A:386:PHE:HB2	1.98	0.45
1:C:385:ILE:HD12	1:C:529:LEU:HD11	1.97	0.45
1:A:14:LEU:O	1:A:74:ASP:HA	2.16	0.45
1:D:305:MET:HE3	1:D:455:LEU:HD22	1.99	0.44
1:D:550:TRP:NE1	1:D:651:TYR:OH	2.49	0.44
1:B:237:ASN:HB2	1:B:355:HIS:HB2	1.99	0.44
1:C:24:LEU:HD22	1:C:35:LEU:HD22	2.00	0.44
1:A:558:PRO:HB3	1:A:582:PRO:HD3	1.99	0.44
1:A:596:HIS:CG	3:A:702:3VV:H13	2.52	0.44
1:B:166:PHE:N	1:B:398:GLU:OE1	2.50	0.44
1:D:420:THR:O	1:D:425:HIS:N	2.51	0.44
1:A:25:TRP:HB2	1:A:64:ARG:HB2	1.99	0.44
1:D:24:LEU:HD22	1:D:35:LEU:HD22	1.99	0.44
1:A:420:THR:O	1:A:425:HIS:N	2.51	0.44
3:A:702:3VV:H56	3:A:702:3VV:H65	1.99	0.44
1:C:336:PRO:HA	1:C:337:PRO:HD3	1.93	0.44
1:D:7:ARG:HG2	1:D:48:GLU:HG2	2.00	0.44
1:B:496:GLN:HE22	1:B:519:PHE:HB2	1.83	0.44
1:C:166:PHE:N	1:C:398:GLU:OE1	2.51	0.44
1:A:463:ASP:OD1	1:A:641:ARG:NH1	2.50	0.44
1:A:5:ARG:HB3	1:A:84:GLN:HB3	1.99	0.44
1:A:237:ASN:HB2	1:A:355:HIS:HB2	1.99	0.44
1:C:240:LEU:HD23	1:C:276:ALA:HB1	2.00	0.43
1:B:27:VAL:HB	1:B:62:PHE:HB2	2.00	0.43
1:B:547:GLN:HB2	1:B:598:SER:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:LYS:HB2	1:A:148:LEU:HD12	2.00	0.43
1:C:370:VAL:HG11	1:C:507:LEU:HG	2.00	0.43
1:D:483:TYR:OH	1:D:492:ASP:OD2	2.33	0.43
1:D:187:LEU:HD22	1:A:187:LEU:HD22	2.01	0.43
1:A:243:ARG:NH1	1:A:271:GLY:O	2.48	0.43
1:C:611:LYS:HA	1:C:611:LYS:HD3	1.93	0.43
1:A:348:ARG:NH1	1:A:351:ASP:OD2	2.51	0.43
1:D:267:GLU:OE1	1:D:272:SER:OG	2.36	0.43
1:D:166:PHE:N	1:D:398:GLU:OE1	2.47	0.43
1:C:417:ALA:HB2	1:C:589:LEU:HD23	2.00	0.42
1:D:132:LEU:HD11	1:D:380:PRO:HD3	2.01	0.42
1:C:7:ARG:HG2	1:C:48:GLU:HG2	2.00	0.42
1:C:219:ARG:NH1	1:C:220:GLN:OE1	2.51	0.42
1:A:443:PRO:HD2	1:A:468:TRP:CE3	2.54	0.42
1:D:370:VAL:HG11	1:D:507:LEU:HG	2.02	0.42
1:A:84:GLN:NE2	1:A:89:CYS:O	2.43	0.42
1:C:267:GLU:OE1	1:C:272:SER:OG	2.33	0.42
1:C:282:ASP:O	1:C:431:ARG:NH1	2.35	0.42
1:C:291:GLY:O	1:D:156:ARG:NE	2.52	0.42
1:D:357:ILE:HD13	1:D:357:ILE:HA	1.93	0.42
1:C:24:LEU:O	1:C:35:LEU:N	2.45	0.42
1:D:652:LEU:HD23	1:D:652:LEU:HA	1.88	0.42
1:A:242:ARG:HG2	1:A:450:ARG:HG2	2.01	0.42
1:B:24:LEU:HD22	1:B:35:LEU:HD22	2.02	0.42
1:C:357:ILE:HD13	1:C:357:ILE:HA	1.93	0.42
1:D:601:GLN:HB2	1:D:604:MET:HB3	2.01	0.42
1:C:544:ASN:O	1:C:547:GLN:NE2	2.52	0.42
1:C:550:TRP:NE1	1:C:651:TYR:OH	2.52	0.42
1:C:547:GLN:HB2	1:C:598:SER:HB3	2.02	0.42
1:A:211:LYS:NZ	1:A:216:GLU:OE2	2.45	0.42
1:D:531:MET:HE3	1:D:531:MET:HB3	1.87	0.41
1:A:596:HIS:CD2	3:A:702:3VV:H13	2.55	0.41
1:B:243:ARG:NH1	1:B:271:GLY:O	2.45	0.41
1:A:542:ALA:HB2	1:A:652:LEU:HD21	2.01	0.41
1:A:601:GLN:HB2	1:A:604:MET:HB3	2.02	0.41
1:D:6:ILE:HG13	1:D:83:VAL:HG13	2.01	0.41
1:D:646:ASP:N	1:D:646:ASP:OD1	2.49	0.41
1:C:23:GLN:HG3	1:C:68:HIS:HB2	2.02	0.41
1:C:185:MET:HE2	1:C:185:MET:HB3	1.93	0.41
1:B:593:ILE:HD13	1:B:593:ILE:HA	1.91	0.41
1:C:496:GLN:HE22	1:C:519:PHE:HB2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:LEU:HD23	1:A:549:ASP:HB3	2.02	0.41
1:B:650:GLU:HB3	1:B:653:LYS:HE2	2.02	0.41
1:C:243:ARG:NH1	1:C:271:GLY:O	2.47	0.41
1:A:471:ILE:HG22	1:A:530:THR:HG23	2.02	0.41
1:B:551:TYR:HB3	1:B:590:GLN:HE21	1.86	0.41
1:C:531:MET:HE3	1:C:531:MET:HB3	1.91	0.41
1:D:547:GLN:HB2	1:D:598:SER:HB3	2.03	0.41
1:B:336:PRO:HA	1:B:337:PRO:HD3	1.88	0.40
1:C:558:PRO:HB3	1:C:582:PRO:HD3	2.02	0.40
1:C:646:ASP:N	1:C:646:ASP:OD1	2.49	0.40
1:A:550:TRP:O	1:A:556:ASN:ND2	2.46	0.40
1:B:188:LYS:O	1:B:192:THR:OG1	2.31	0.40
1:B:558:PRO:HB3	1:B:582:PRO:HD3	2.03	0.40
1:C:550:TRP:O	1:C:556:ASN:ND2	2.47	0.40
1:C:601:GLN:HB2	1:C:604:MET:HB3	2.03	0.40
1:D:241:LEU:HD11	1:D:274:PHE:HB3	2.03	0.40
1:D:417:ALA:HB2	1:D:589:LEU:HD23	2.02	0.40
1:A:9:ALA:HB3	1:A:80:ARG:HG2	2.02	0.40
1:C:132:LEU:HD11	1:C:380:PRO:HD3	2.04	0.40
1:D:24:LEU:O	1:D:35:LEU:N	2.46	0.40
1:A:652:LEU:HD23	1:A:652:LEU:HA	1.87	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	662/669 (99%)	653 (99%)	9 (1%)	0	100	100
1	B	662/669 (99%)	650 (98%)	12 (2%)	0	100	100
1	C	661/669 (99%)	648 (98%)	13 (2%)	0	100	100
1	D	661/669 (99%)	650 (98%)	11 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2646/2676 (99%)	2601 (98%)	45 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	572/578 (99%)	572 (100%)	0	100	100
1	B	572/578 (99%)	572 (100%)	0	100	100
1	C	571/578 (99%)	571 (100%)	0	100	100
1	D	571/578 (99%)	571 (100%)	0	100	100
All	All	2286/2312 (99%)	2286 (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 (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	20	ASN
1	B	68	HIS
1	B	237	ASN
1	B	308	ASN
1	B	358	GLN
1	B	400	ASN
1	C	125	GLN
1	C	626	GLN
1	D	51	HIS
1	D	84	GLN
1	D	119	ASN
1	D	237	ASN
1	D	358	GLN
1	D	406	GLN
1	A	167	HIS

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Mol	Chain	Res	Type
1	A	358	GLN
1	A	520	GLN
1	A	547	GLN
1	A	596	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	3VV	A	702	-	67,69,69	0.44	0	89,95,95	0.67	2 (2%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	3VV	A	702	-	-	7/68/84/84	0/3/3/3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	A	702	3VV	C16-C17-C18	2.58	117.96	112.27
3	A	702	3VV	P54-O53-C52	-2.56	116.59	123.43

There are no chirality outliers.

All (7) torsion outliers are listed below:

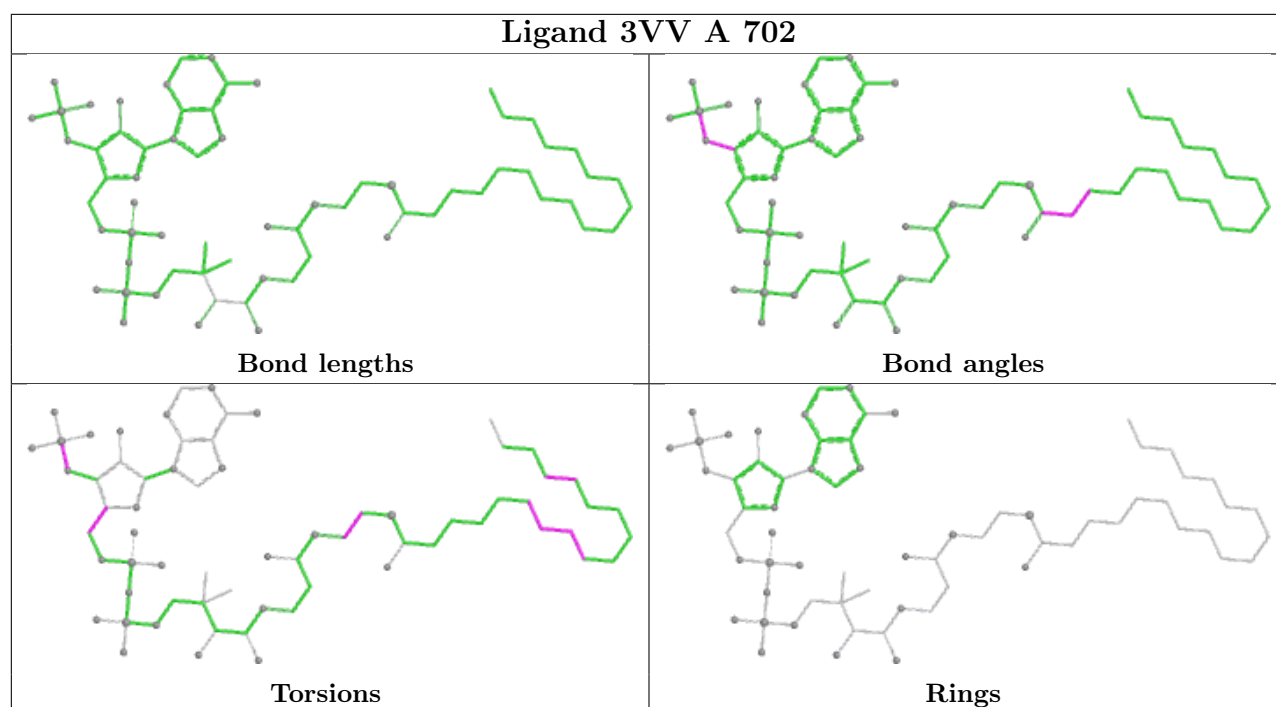
Mol	Chain	Res	Type	Atoms
3	A	702	3VV	S20-C21-C22-N23
3	A	702	3VV	C03-C04-C05-C06
3	A	702	3VV	O45-C46-C47-O48
3	A	702	3VV	C11-C12-C13-C14
3	A	702	3VV	C09-C10-C11-C12
3	A	702	3VV	C10-C11-C12-C13
3	A	702	3VV	C52-O53-P54-O55

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	3VV	7	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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.

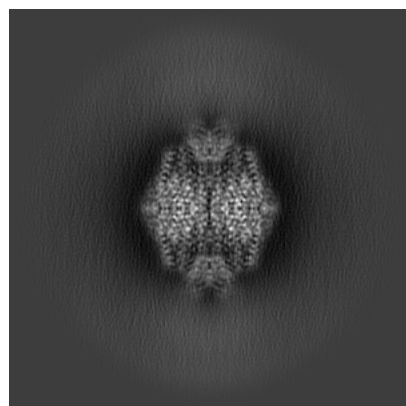
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40042. These allow visual inspection of the internal detail of the map and identification of artifacts.

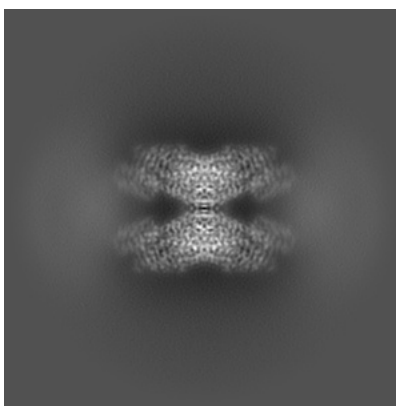
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

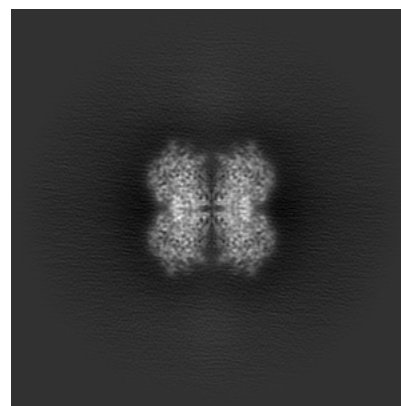
6.1.1 Primary map



X

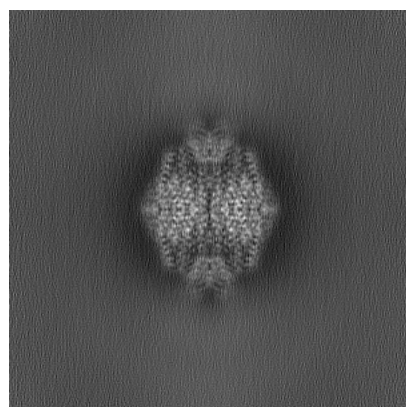


Y

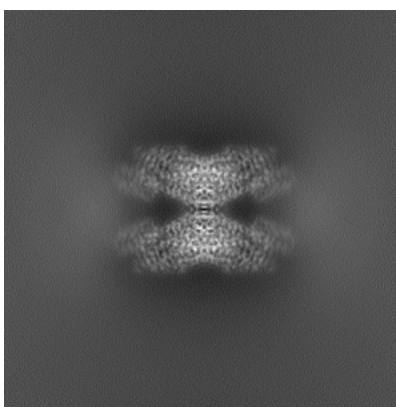


Z

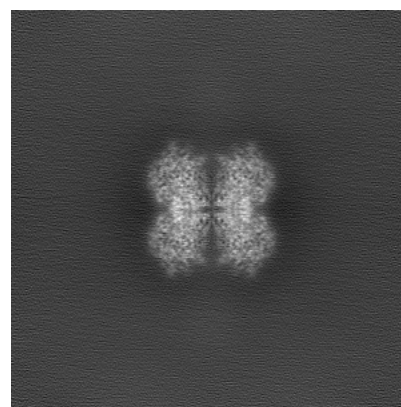
6.1.2 Raw map



X



Y

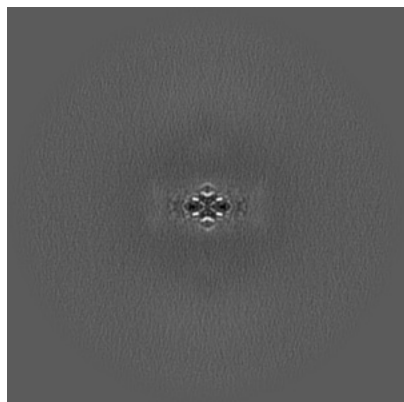


Z

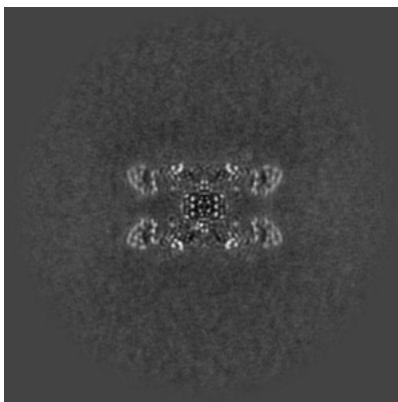
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

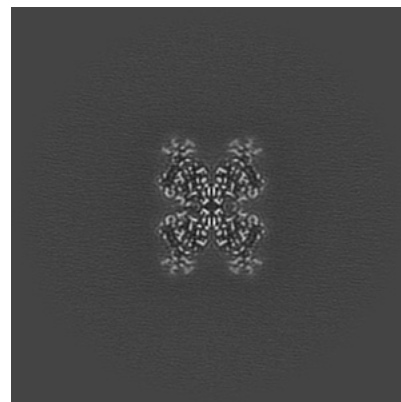
6.2.1 Primary map



X Index: 180

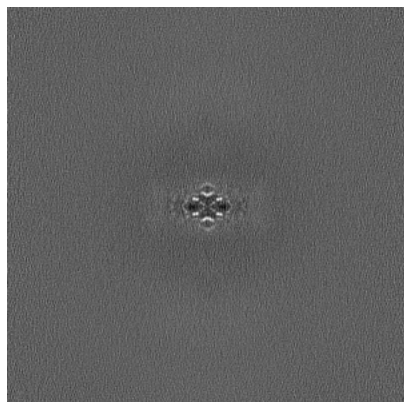


Y Index: 180

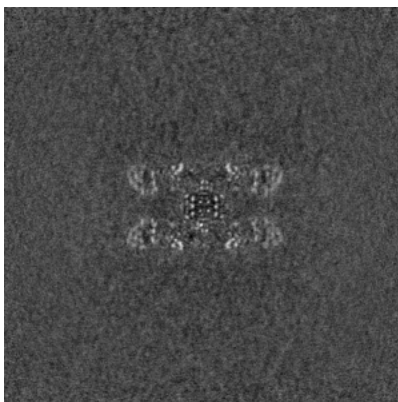


Z Index: 180

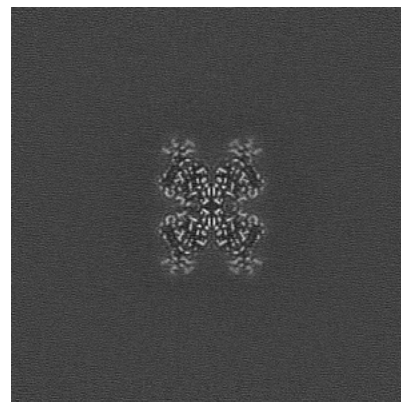
6.2.2 Raw map



X Index: 180



Y Index: 180

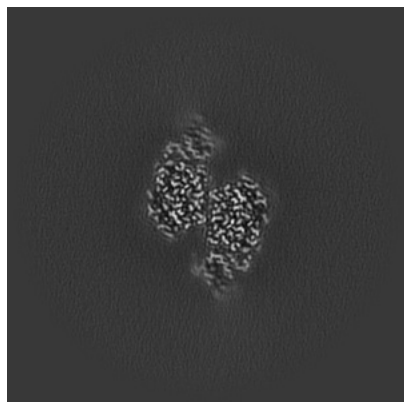


Z Index: 180

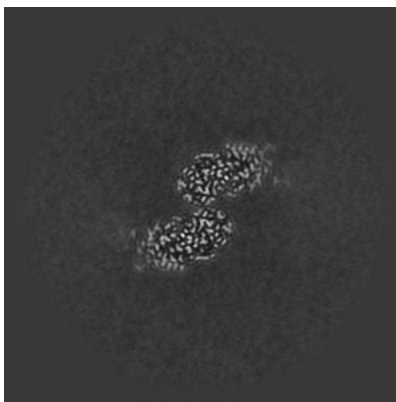
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

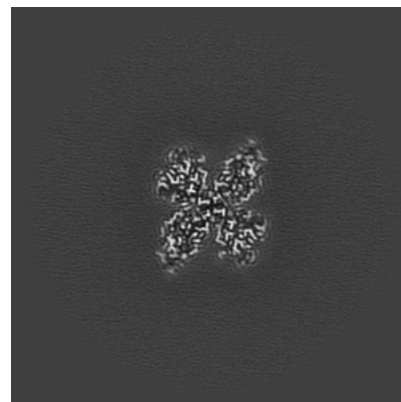
6.3.1 Primary map



X Index: 152

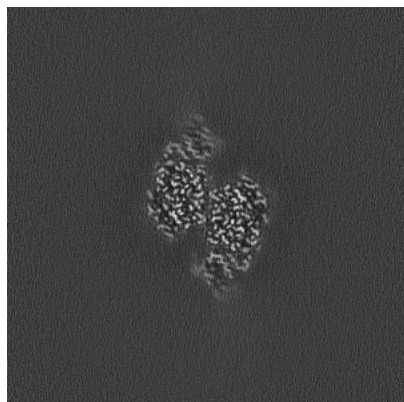


Y Index: 203

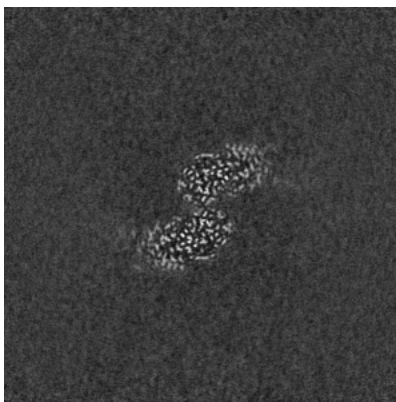


Z Index: 172

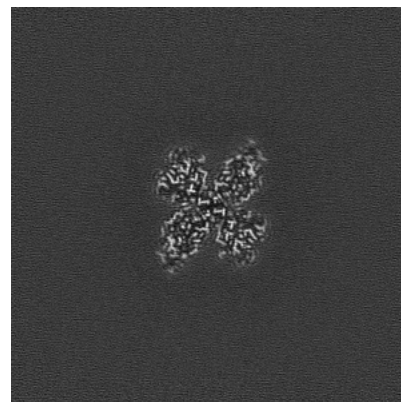
6.3.2 Raw map



X Index: 152



Y Index: 203

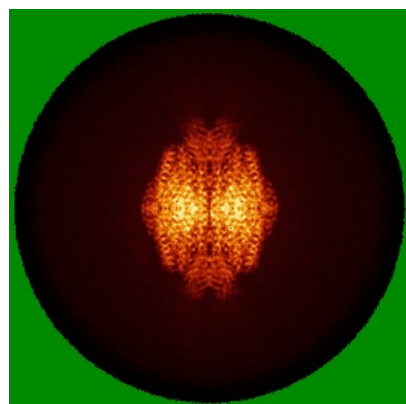


Z Index: 172

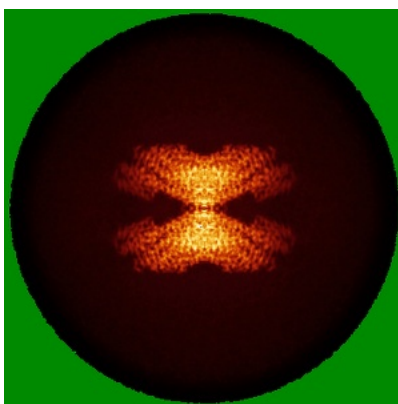
The images above show the largest variance slices of the map in three orthogonal directions.

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

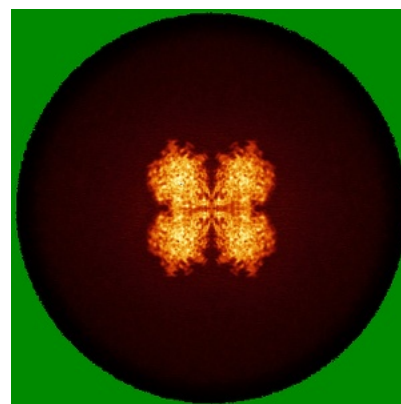
6.4.1 Primary map



X

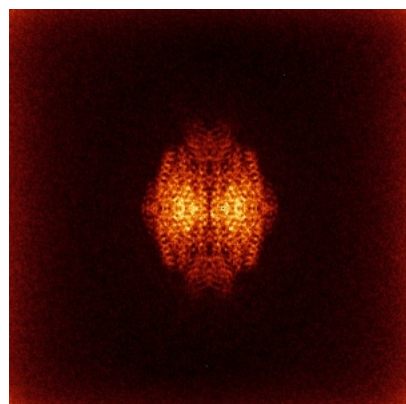


Y

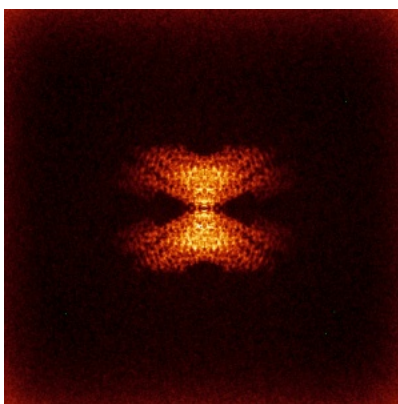


Z

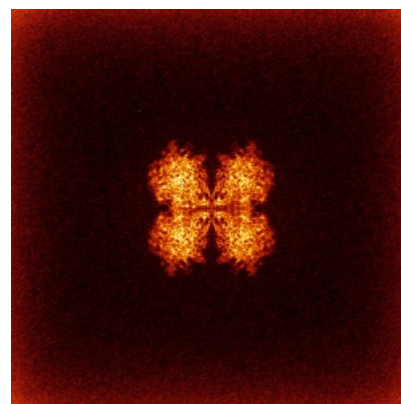
6.4.2 Raw map



X



Y

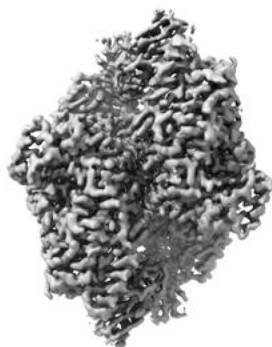


Z

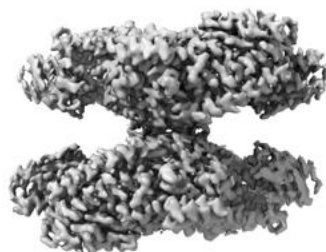
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

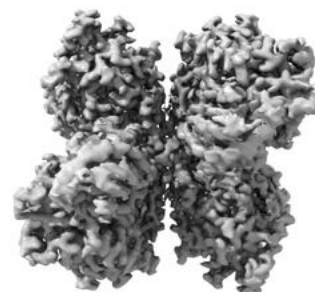
6.5.1 Primary map



X



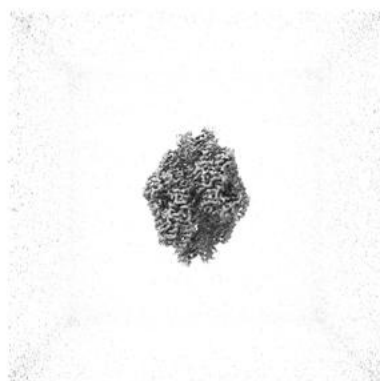
Y



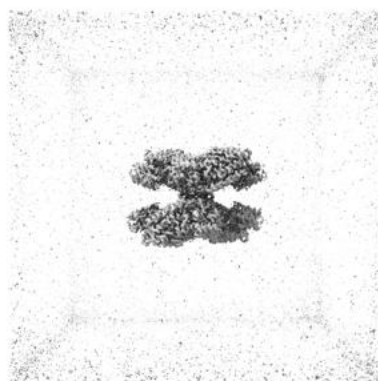
Z

The images above show the 3D surface view of the map at the recommended contour level 0.35. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

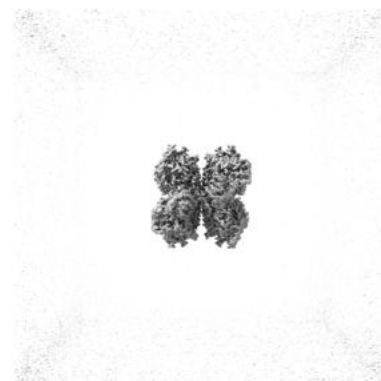
6.5.2 Raw map



X



Y



Z

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

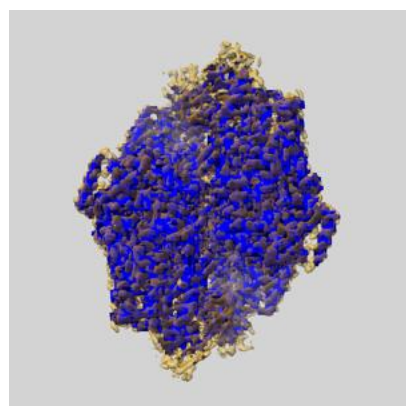
6.6 Mask visualisation [i](#)

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

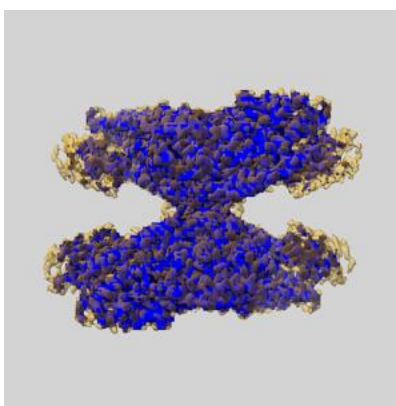
A mask typically either:

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

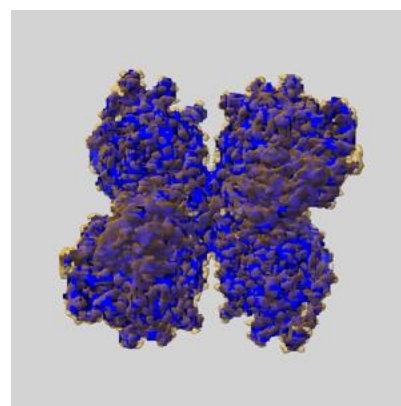
6.6.1 emd_40042_msk_1.map [i](#)



X



Y

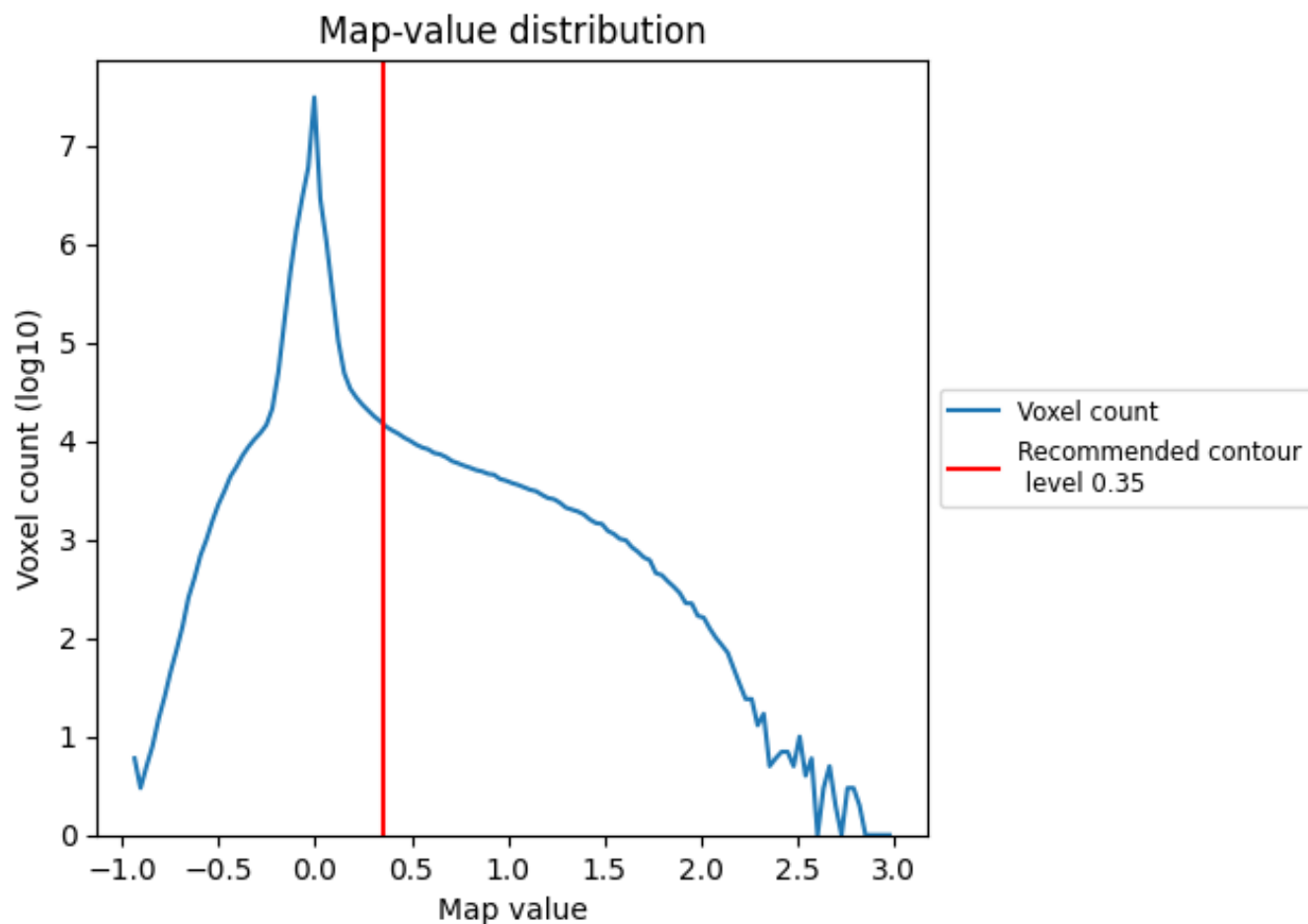


Z

7 Map analysis [i](#)

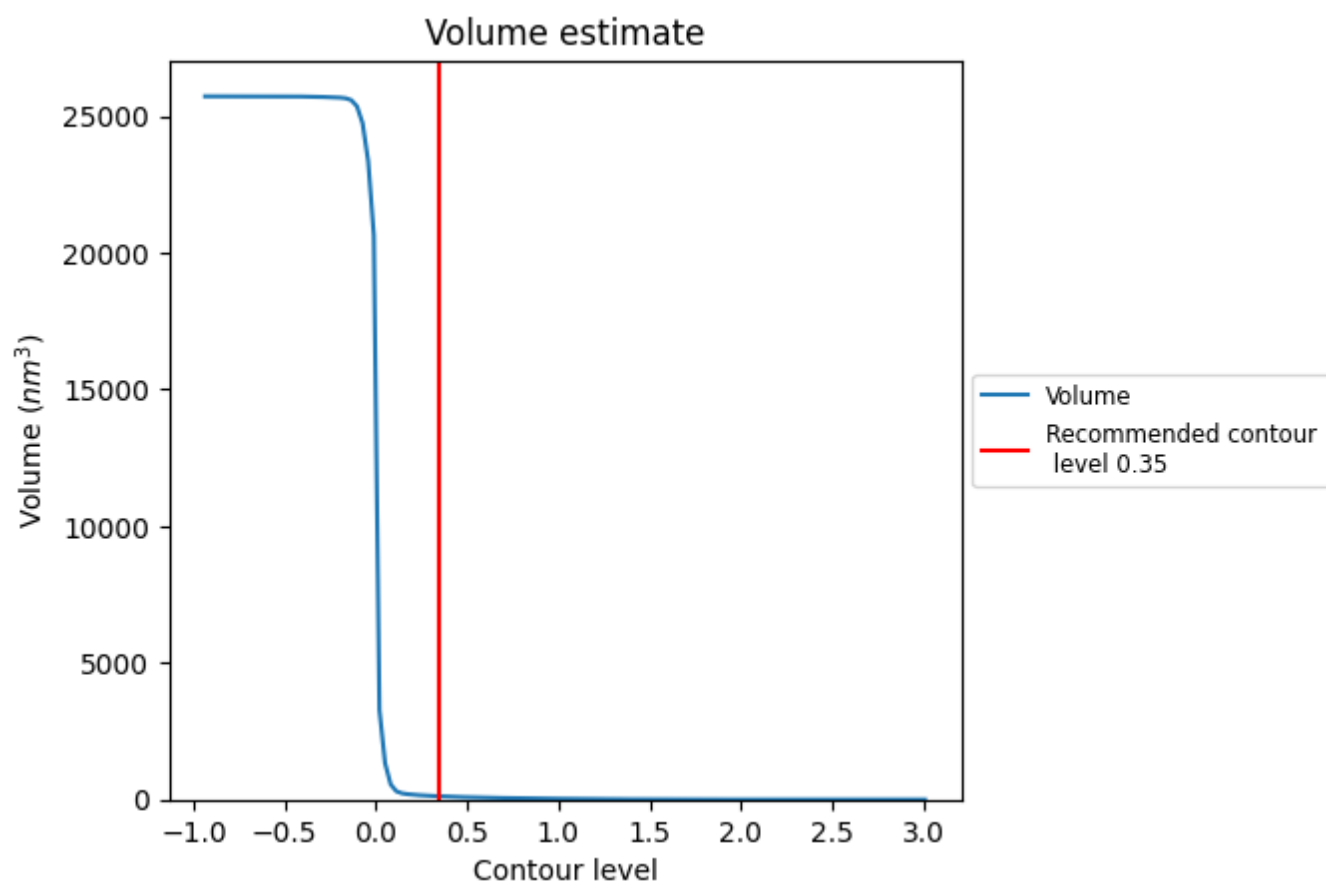
This section contains the results of statistical analysis of the map.

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.

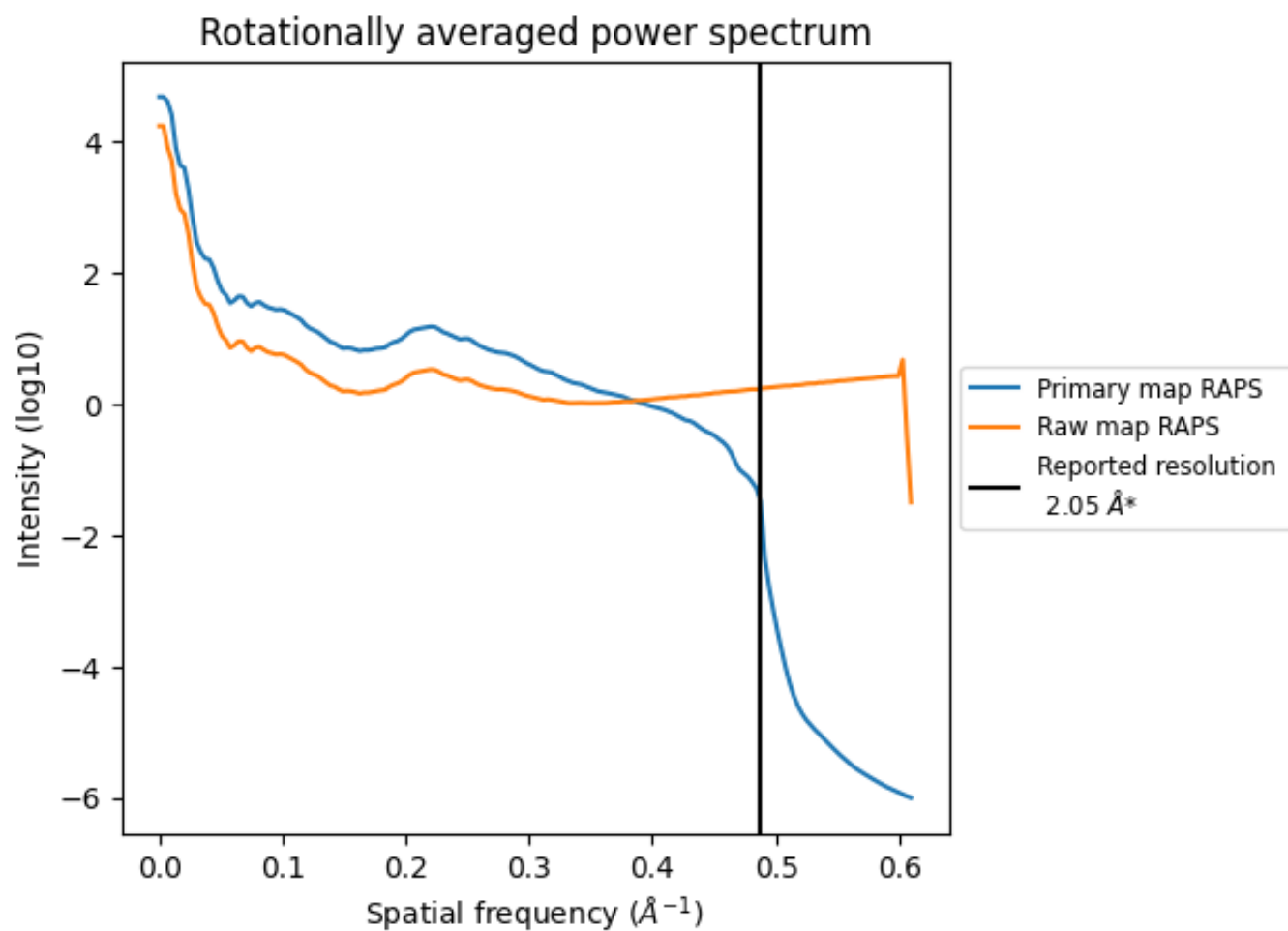
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 120 nm^3 ; this corresponds to an approximate mass of 108 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 ⓘ

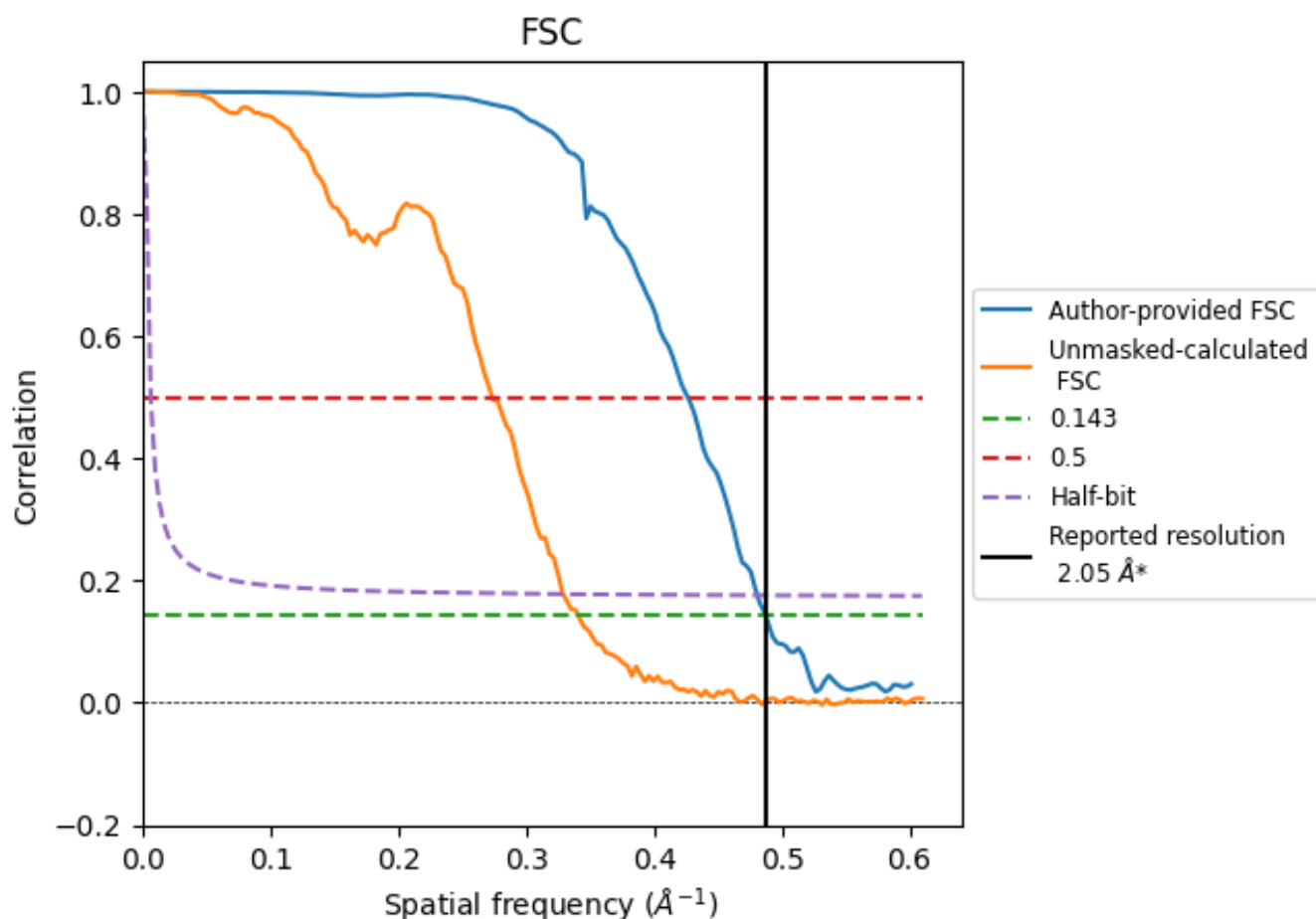


*Reported resolution corresponds to spatial frequency of 0.488 Å⁻¹

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 $0.488 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

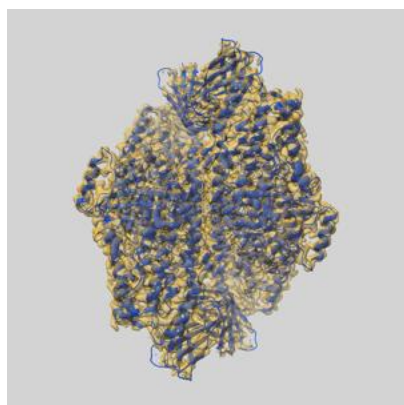
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.05	-	-
Author-provided FSC curve	2.05	2.34	2.08
Unmasked-calculated*	2.93	3.65	3.04

*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 2.93 differs from the reported value 2.05 by more than 10 %

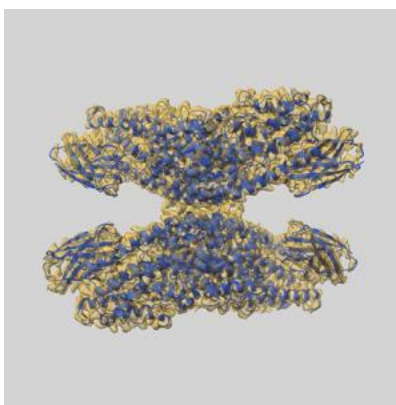
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-40042 and PDB model 8GHE. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

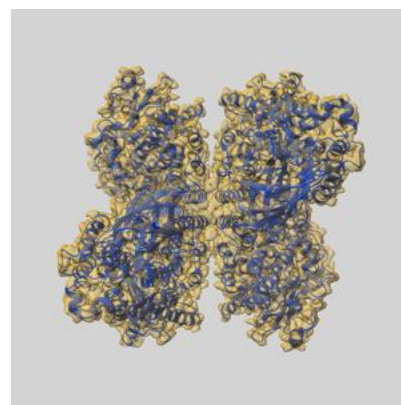
9.1 Map-model overlay [i](#)



X



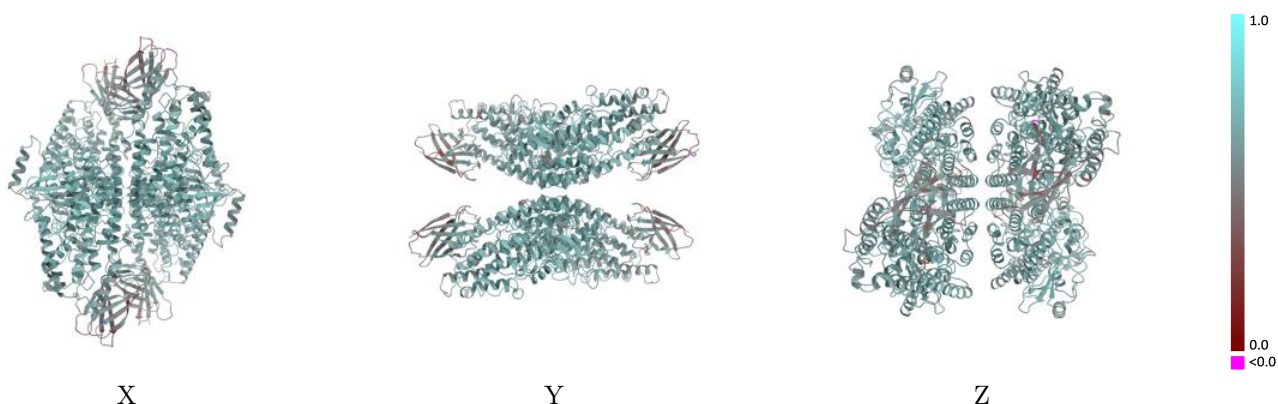
Y



Z

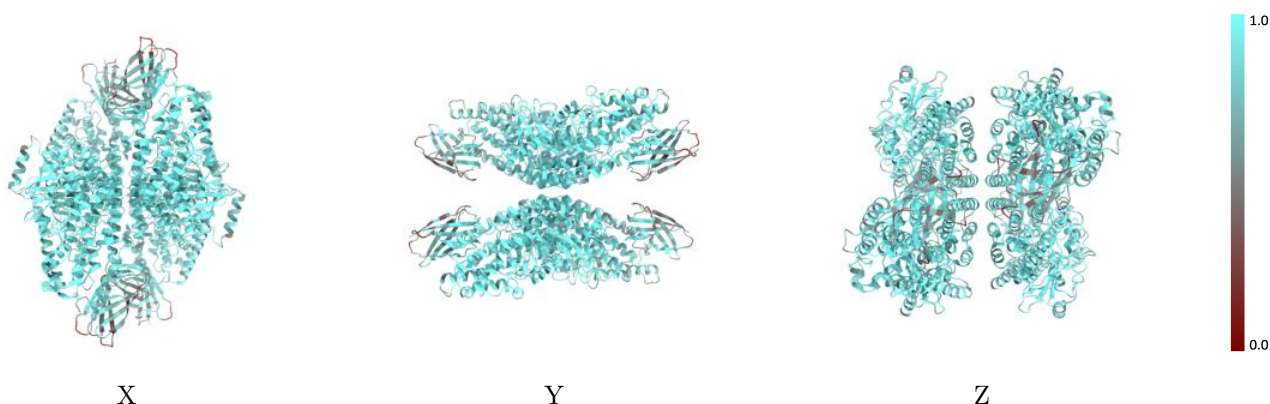
The images above show the 3D surface view of the map at the recommended contour level 0.35 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



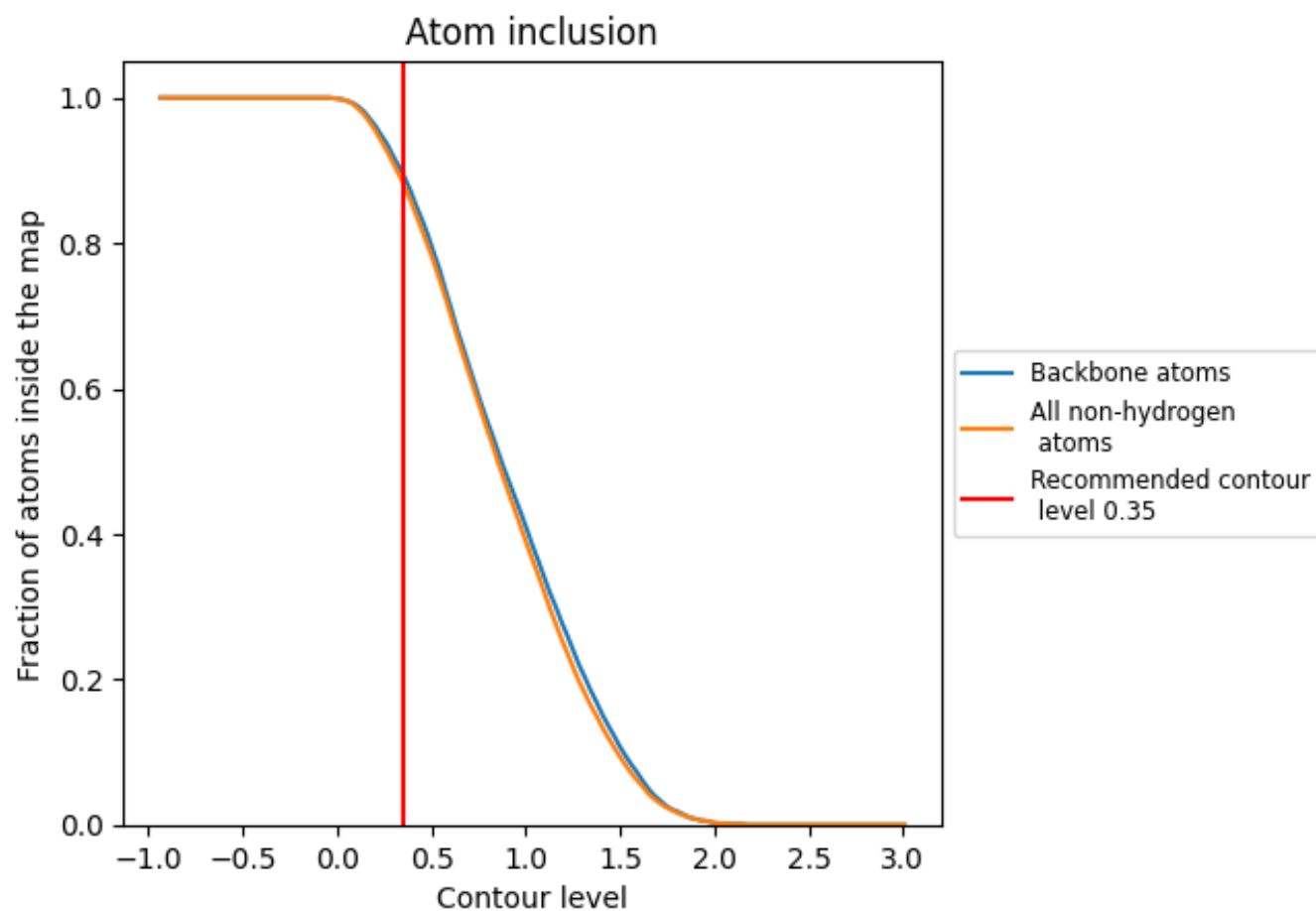
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.35).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 88% 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 (0.35) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8830	0.6240
A	0.9000	0.6310
B	0.8850	0.6230
C	0.8800	0.6220
D	0.8750	0.6180

