Full wwPDB EM Validation Report (i)

Oct 15, 2026 – 05:51 pm BST

PDB ID	:	32LJ / pdb_000032lj
EMDB ID	:	EMD-58942
Title	:	Cryo-EM structure of BnC1 cruciferin homohexamer, an 11S globulin from Brassica napus
Authors	:	Boyens-Thiele, L.; Buell, A.K.; Morth, J.P.
Deposited on	:	2026-07-15
Resolution	:	1.83 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 (i) 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 (i)) were used in the production of this report:

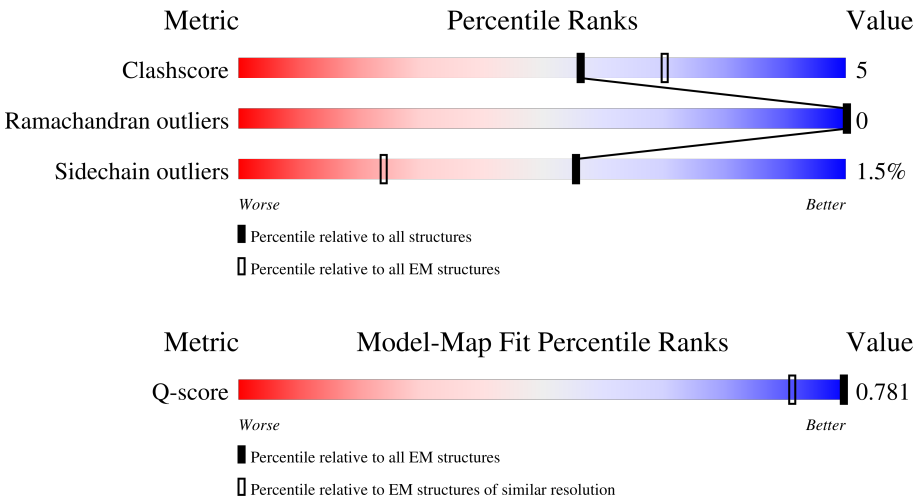
EMDB validation analysis	:	0.0.1.dev133
MolProbity	:	4-5-2 with Phenix2.2
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

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 1.83 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	903 (1.33 - 2.33)

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 <=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	490	69% 9% 22%
1	B	490	67% 11% 22%
1	C	490	68% 10% 22%
1	D	490	67% 11% 22%

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Mol	Chain	Length	Quality of chain
1	E	490	66% 12% 22%
1	F	490	68% 10% 22%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 18616 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 Cruciferin BnC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	384	3023	1901	547	567	8	0	0
1	B	384	3023	1901	547	567	8	0	0
1	C	384	3023	1901	547	567	8	0	0
1	D	384	3023	1901	547	567	8	0	0
1	E	384	3023	1901	547	567	8	0	0
1	F	384	3023	1901	547	567	8	0	0

- Molecule 2 is water.

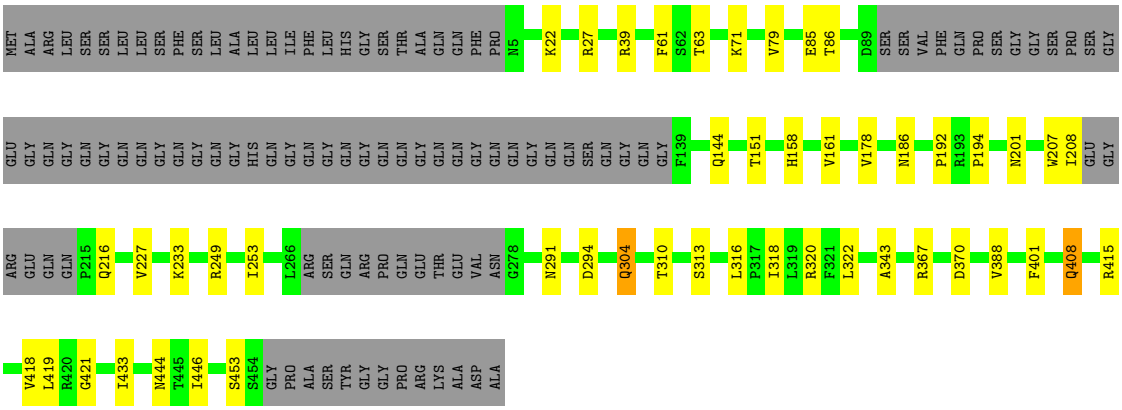
Mol	Chain	Residues	Atoms		AltConf
			Total	O	
2	A	75	75	75	0
2	B	81	81	81	0
2	C	76	76	76	0
2	D	80	80	80	0
2	E	83	83	83	0
2	F	83	83	83	0

3 Residue-property plots (i)

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.

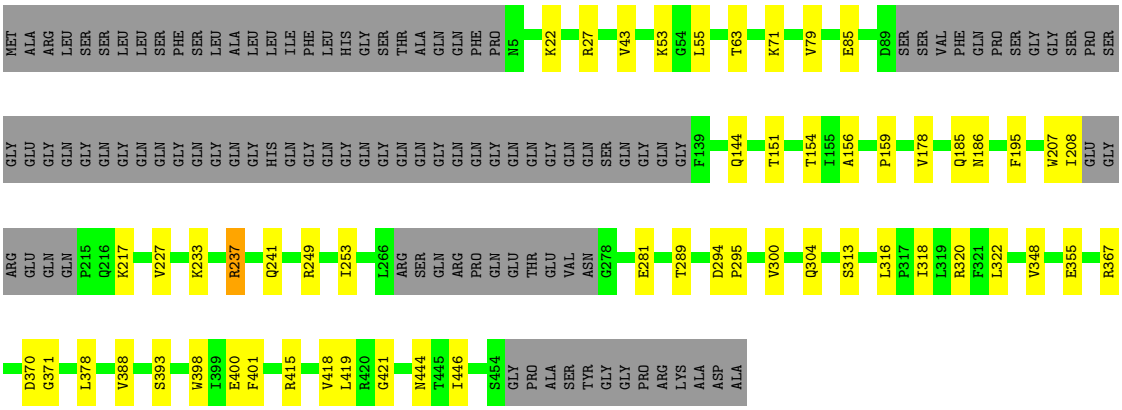
- Molecule 1: Cruciferin BnC1

Chain A:



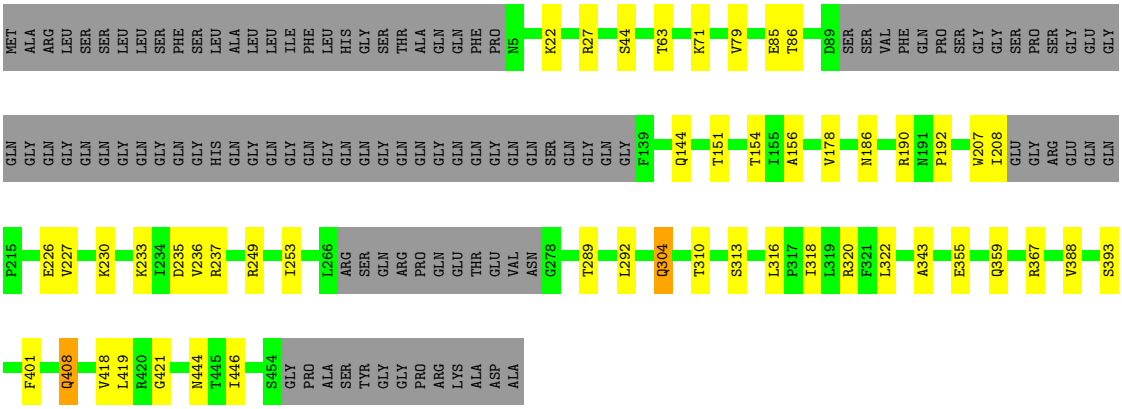
- Molecule 1: Cruciferin BnC1

Chain B:



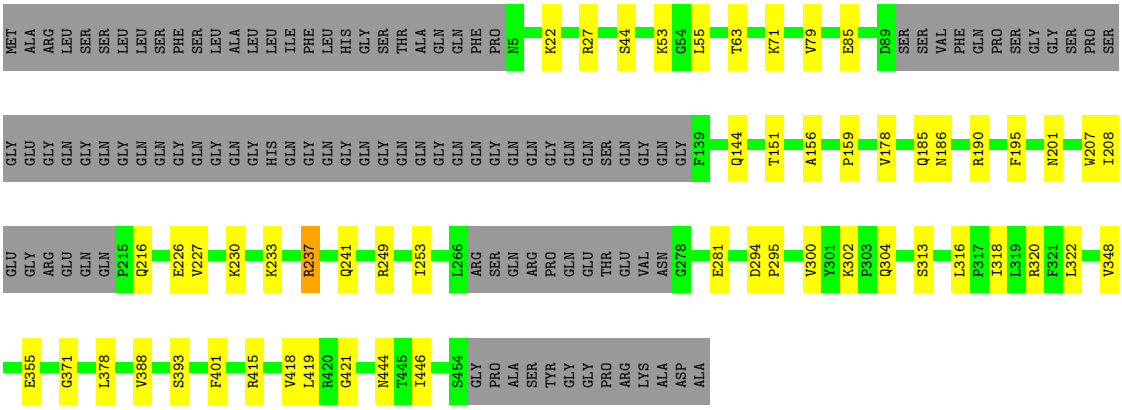
- Molecule 1: Cruciferin BnC1

Chain C:



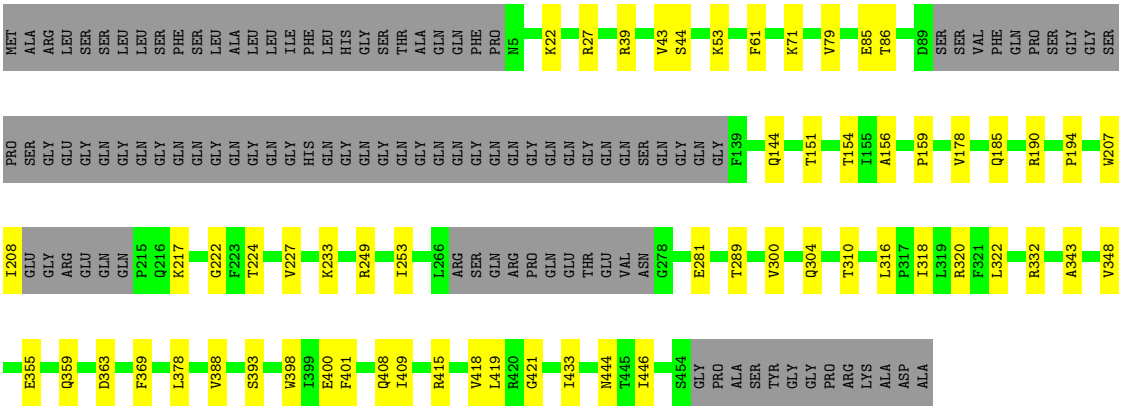
• Molecule 1: Cruciferin BnC1

Chain D: 67% 11% 22%



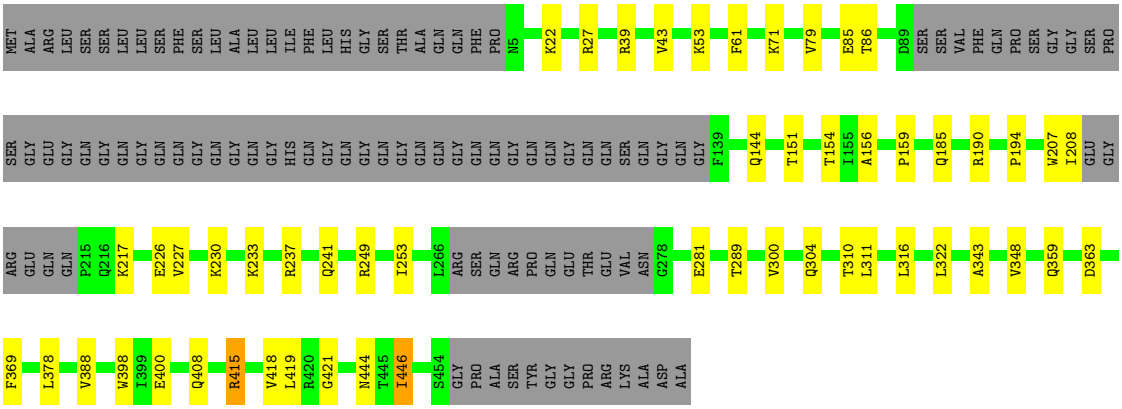
• Molecule 1: Cruciferin BnC1

Chain E: 66% 12% 22%



• Molecule 1: Cruciferin BnC1

Chain F: 68% 10% 22%


MET ALA ARG LEU SER SER LEU LEU SER PHE SER LEU ALA LEU LEU ILE PHE LEU HIS GLY SER THR ALA GLN GLN PHE PRO N5 K22 R27 R39 V43 K53 F61 K71 V79 E85 T86 D89 SER SER VAL PHE GLN PRO SER GLY GLY SER PRO
SER GLY GLU GLY GLN GLY GLN GLY GLN GLN GLY GLN GLY GLN GLY HIS GLN GLY GLN GLY GLN GLY GLN GLN GLY GLN GLN GLY GLN GLN GLN SER GLN GLY GLN GLY F139 Q144 T151 T154 I155 A156 P159 Q185 R190 P194 W207 I208 GLU GLY
ARG GLU GLN GLN P215 Q216 K217 E226 V227 K230 K233 R237 Q241 R249 I253 L266 ARG SER GLN ARG PRO GLN GLU THR GLU VAL ASN G278 E281 T289 V300 Q304 T310 L311 L316 L322 A343 V348 Q359 D363
F369 L378 V388 W398 I399 E400 Q408 R415 V418 L419 R420 G421 N444 T445 I446 S454 GLY PRO ALA SER TYR GLY GLY PRO ARG LYS ALA ASP ALA

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, D3	Depositor
Number of particles used	309000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	59.5	Depositor
Minimum defocus (nm)	300	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.541	Depositor
Minimum map value	-0.261	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.018	Depositor
Recommended contour level	0.05	Depositor
Map size (Å)	232.92, 232.92, 232.92	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.647, 0.647, 0.647	Depositor

5 Model quality (i)

5.1 Standard geometry (i)

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	0.32	0/3084	0.43	0/4187
1	B	0.32	0/3084	0.43	0/4187
1	C	0.32	0/3084	0.43	0/4187
1	D	0.32	0/3084	0.43	0/4187
1	E	0.32	0/3084	0.43	0/4187
1	F	0.32	0/3084	0.43	0/4187
All	All	0.32	0/18504	0.43	0/25122

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (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	3023	0	2977	34	0
1	B	3023	0	2977	36	0
1	C	3023	0	2977	35	0
1	D	3023	0	2977	36	0
1	E	3023	0	2977	38	0
1	F	3023	0	2977	37	0
2	A	75	0	0	0	0
2	B	81	0	0	1	0
2	C	76	0	0	1	0
2	D	80	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	83	0	0	2	0
2	F	83	0	0	1	0
All	All	18616	0	17862	178	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 (178) 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:C:79:VAL:HG11	1:D:418:VAL:HA	1.78	0.64
1:A:79:VAL:HG11	1:B:418:VAL:HA	1.79	0.64
1:D:79:VAL:HG11	1:E:418:VAL:HA	1.80	0.63
1:B:79:VAL:HG11	1:F:418:VAL:HA	1.80	0.62
1:A:158:HIS:CE1	1:A:161:VAL:HG21	2.34	0.62
1:A:418:VAL:HA	1:F:79:VAL:HG11	1.81	0.61
1:C:418:VAL:HA	1:E:79:VAL:HG11	1.82	0.61
1:F:22:LYS:HD2	1:F:27:ARG:HB3	1.83	0.61
1:E:22:LYS:HD2	1:E:27:ARG:HB3	1.83	0.61
1:A:158:HIS:CD2	1:A:318:ILE:HD11	2.38	0.58
1:B:22:LYS:HD2	1:B:27:ARG:HB3	1.85	0.58
1:A:22:LYS:HD2	1:A:27:ARG:HB3	1.84	0.58
1:C:22:LYS:HD2	1:C:27:ARG:HB3	1.86	0.58
1:C:236:VAL:HG12	1:C:237:ARG:HH11	1.70	0.57
1:D:22:LYS:HD2	1:D:27:ARG:HB3	1.86	0.57
1:D:304:GLN:H	1:D:304:GLN:CD	2.13	0.56
1:C:388:VAL:HG11	1:E:227:VAL:HG11	1.88	0.56
1:C:304:GLN:H	1:C:304:GLN:CD	2.14	0.55
1:E:71:LYS:HA	1:E:151:THR:HG22	1.89	0.55
1:B:304:GLN:CD	1:B:304:GLN:H	2.14	0.55
1:F:304:GLN:CD	1:F:304:GLN:H	2.16	0.54
1:D:355:GLU:HG2	1:D:393:SER:HB3	1.89	0.54
1:F:71:LYS:HA	1:F:151:THR:HG22	1.90	0.54
1:E:304:GLN:H	1:E:304:GLN:CD	2.17	0.53
1:C:71:LYS:HA	1:C:151:THR:HG22	1.91	0.52
1:D:71:LYS:HA	1:D:151:THR:HG22	1.91	0.52
1:A:71:LYS:HA	1:A:151:THR:HG22	1.92	0.52
1:A:304:GLN:H	1:A:304:GLN:CD	2.17	0.52
1:A:453:SER:HB2	1:F:230:LYS:HD3	1.92	0.52
1:B:71:LYS:HA	1:B:151:THR:HG22	1.92	0.52
1:C:227:VAL:HG11	1:D:388:VAL:HG11	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:VAL:HG11	1:F:388:VAL:HG11	1.92	0.51
1:D:227:VAL:HG11	1:E:388:VAL:HG11	1.93	0.51
1:E:343:ALA:HB3	1:E:408:GLN:HG3	1.93	0.51
1:A:227:VAL:HG11	1:B:388:VAL:HG11	1.93	0.51
1:B:355:GLU:HG2	1:B:393:SER:HB3	1.92	0.51
1:A:158:HIS:HB3	1:A:322:LEU:HD21	1.93	0.50
1:A:388:VAL:HG11	1:F:227:VAL:HG11	1.93	0.50
1:B:249:ARG:HB3	1:B:253:ILE:HG22	1.94	0.50
1:D:249:ARG:HB3	1:D:253:ILE:HG22	1.94	0.49
1:A:158:HIS:HD2	1:A:318:ILE:HD11	1.76	0.49
1:C:235:ASP:OD1	1:C:237:ARG:HB2	2.12	0.49
1:F:343:ALA:HB3	1:F:408:GLN:HG3	1.94	0.49
1:E:355:GLU:HG2	1:E:393:SER:HB3	1.93	0.49
1:F:226:GLU:HG3	1:F:230:LYS:NZ	2.28	0.49
1:A:313:SER:HB2	1:A:320:ARG:HB2	1.96	0.48
1:C:226:GLU:HG3	1:C:230:LYS:NZ	2.28	0.48
1:C:355:GLU:HG2	1:C:393:SER:HB3	1.94	0.48
1:F:217:LYS:HB2	1:F:217:LYS:HE2	1.66	0.48
1:A:85:GLU:OE1	1:B:421:GLY:HA2	2.13	0.48
1:F:311:LEU:HD21	1:F:316:LEU:HD12	1.97	0.47
1:E:249:ARG:HB3	1:E:253:ILE:HG22	1.96	0.47
1:C:85:GLU:OE1	1:D:421:GLY:HA2	2.14	0.47
1:E:43:VAL:HG21	1:E:348:VAL:HG21	1.96	0.47
1:A:291:ASN:HB3	1:A:294:ASP:OD2	2.15	0.47
1:B:316:LEU:HG	1:B:318:ILE:HG22	1.96	0.47
1:C:421:GLY:HA2	1:E:85:GLU:OE1	2.14	0.47
1:D:316:LEU:HG	1:D:318:ILE:HG22	1.96	0.47
1:E:61:PHE:CE1	1:E:194:PRO:HB3	2.50	0.47
1:E:418:VAL:HG13	1:E:419:LEU:HD12	1.96	0.47
1:A:421:GLY:HA2	1:F:85:GLU:OE1	2.14	0.47
1:A:192:PRO:HG2	1:A:207:TRP:CD1	2.50	0.47
1:D:85:GLU:OE1	1:E:421:GLY:HA2	2.15	0.47
1:B:85:GLU:OE1	1:F:421:GLY:HA2	2.15	0.46
1:C:192:PRO:HG2	1:C:207:TRP:CD1	2.50	0.46
1:F:53:LYS:HG3	2:F:523:HOH:O	2.15	0.46
1:D:313:SER:HB2	1:D:320:ARG:HB2	1.97	0.46
1:E:53:LYS:HG3	2:E:527:HOH:O	2.15	0.46
1:A:61:PHE:CE1	1:A:194:PRO:HB3	2.50	0.46
1:F:418:VAL:HG13	1:F:419:LEU:HD12	1.96	0.46
1:A:207:TRP:CD1	1:A:208:ILE:HG13	2.51	0.46
1:C:316:LEU:HG	1:C:318:ILE:HG22	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:249:ARG:HB3	1:F:253:ILE:HG22	1.97	0.46
1:A:249:ARG:HB3	1:A:253:ILE:HG22	1.98	0.45
1:C:249:ARG:HB3	1:C:253:ILE:HG22	1.97	0.45
1:C:226:GLU:HG3	1:C:230:LYS:HZ2	1.81	0.45
1:D:418:VAL:HG13	1:D:419:LEU:HD12	1.98	0.45
1:E:44:SER:HB2	2:E:508:HOH:O	2.16	0.45
1:F:61:PHE:CE1	1:F:194:PRO:HB3	2.51	0.45
1:D:44:SER:HB2	2:D:506:HOH:O	2.16	0.45
1:B:217:LYS:HB2	1:B:217:LYS:HE2	1.65	0.45
1:E:281:GLU:HB2	1:F:310:THR:HG21	1.98	0.45
1:C:236:VAL:HG12	1:C:237:ARG:NH1	2.31	0.45
1:D:233:LYS:HB2	1:E:444:ASN:HB3	1.99	0.45
1:D:207:TRP:CD1	1:D:208:ILE:HG13	2.53	0.44
1:F:207:TRP:CD1	1:F:208:ILE:HG13	2.52	0.44
1:C:44:SER:HB2	2:C:506:HOH:O	2.17	0.44
1:B:207:TRP:CD1	1:B:208:ILE:HG13	2.53	0.44
1:F:159:PRO:HG3	1:F:185:GLN:HB3	1.99	0.44
1:B:55:LEU:HD23	1:B:253:ILE:HD11	2.00	0.44
1:E:207:TRP:CD1	1:E:208:ILE:HG13	2.52	0.44
1:B:281:GLU:HB2	1:C:310:THR:HG21	2.00	0.44
1:E:310:THR:HG21	1:F:281:GLU:HB2	2.00	0.43
1:A:343:ALA:HB3	1:A:408:GLN:HG3	2.00	0.43
1:B:418:VAL:HG13	1:B:419:LEU:HD12	2.00	0.43
1:D:63:THR:HG22	1:D:186:ASN:HB2	1.99	0.43
1:B:313:SER:HB2	1:B:320:ARG:HB2	2.01	0.43
1:C:85:GLU:OE1	1:C:144:GLN:HG3	2.19	0.43
1:C:367:ARG:HD3	1:E:222:GLY:O	2.18	0.43
1:C:343:ALA:HB3	1:C:408:GLN:HG3	2.00	0.43
1:C:444:ASN:HB3	1:E:233:LYS:HB2	2.00	0.43
1:A:85:GLU:OE1	1:A:144:GLN:HG3	2.18	0.43
1:A:444:ASN:HB3	1:F:233:LYS:HB2	2.01	0.43
1:F:43:VAL:HG21	1:F:348:VAL:HG21	2.00	0.43
1:A:316:LEU:HG	1:A:318:ILE:HG22	2.00	0.43
1:A:433:ILE:HG22	1:F:241:GLN:OE1	2.19	0.43
1:D:55:LEU:HD23	1:D:253:ILE:HD11	2.01	0.43
1:E:159:PRO:HG3	1:E:185:GLN:HB3	2.01	0.43
1:E:178:VAL:HG11	1:E:401:PHE:HB3	2.01	0.43
1:B:233:LYS:HB2	1:F:444:ASN:HB3	2.00	0.43
1:F:348:VAL:CG1	1:F:378:LEU:HD11	2.49	0.43
1:A:178:VAL:HG11	1:A:401:PHE:CB	2.49	0.43
1:A:418:VAL:HG13	1:A:419:LEU:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:ALA:HB1	1:B:322:LEU:HD12	2.01	0.42
1:C:418:VAL:HG13	1:C:419:LEU:HD12	2.01	0.42
1:E:156:ALA:HB1	1:E:322:LEU:HD12	2.01	0.42
1:A:367:ARG:HD3	1:A:370:ASP:HB2	2.01	0.42
1:B:300:VAL:HG22	1:C:86:THR:HB	2.01	0.42
1:E:316:LEU:HG	1:E:318:ILE:HG22	2.01	0.42
1:D:85:GLU:OE1	1:D:144:GLN:HG3	2.20	0.42
1:E:86:THR:HB	1:F:300:VAL:HG22	2.01	0.42
1:B:159:PRO:HG3	1:B:185:GLN:HB3	2.00	0.42
1:B:195:PHE:CE2	1:F:363:ASP:HB3	2.55	0.42
1:D:302:LYS:HB3	1:D:304:GLN:NE2	2.34	0.42
1:F:154:THR:O	1:F:289:THR:HA	2.20	0.42
1:C:233:LYS:HB2	1:D:444:ASN:HB3	2.01	0.42
1:C:178:VAL:HG11	1:C:401:PHE:CB	2.50	0.42
1:C:207:TRP:CD1	1:C:208:ILE:HG13	2.54	0.42
1:B:348:VAL:CG1	1:B:378:LEU:HD11	2.50	0.42
1:C:359:GLN:HB2	1:C:388:VAL:HG12	2.02	0.42
1:D:178:VAL:HG11	1:D:401:PHE:HB3	2.02	0.42
1:E:398:TRP:CD1	1:E:400:GLU:HG3	2.55	0.42
1:D:156:ALA:HB1	1:D:322:LEU:HD12	2.01	0.42
1:B:63:THR:HG22	1:B:186:ASN:HB2	2.01	0.42
1:E:154:THR:O	1:E:289:THR:HA	2.19	0.42
1:E:178:VAL:HG11	1:E:401:PHE:CB	2.50	0.42
1:F:85:GLU:OE1	1:F:144:GLN:HG3	2.20	0.42
1:B:237:ARG:HE	1:B:241:GLN:HG3	1.85	0.42
1:B:367:ARG:HD3	1:B:370:ASP:HB2	2.01	0.42
1:F:156:ALA:HB1	1:F:322:LEU:HD12	2.02	0.42
1:A:233:LYS:HB2	1:B:444:ASN:HB3	2.02	0.41
1:C:313:SER:HB2	1:C:320:ARG:HB2	2.01	0.41
1:D:178:VAL:HG11	1:D:401:PHE:CB	2.50	0.41
1:F:415:ARG:HD3	1:F:446:ILE:O	2.20	0.41
1:A:63:THR:HG22	1:A:186:ASN:HB2	2.02	0.41
1:A:310:THR:HG21	1:D:281:GLU:HB2	2.02	0.41
1:B:178:VAL:HG11	1:B:401:PHE:HB3	2.01	0.41
1:D:159:PRO:HG3	1:D:185:GLN:HB3	2.01	0.41
1:D:237:ARG:HE	1:D:241:GLN:HG3	1.85	0.41
1:A:86:THR:HB	1:D:300:VAL:HG22	2.02	0.41
1:C:63:THR:HG22	1:C:186:ASN:CG	2.46	0.41
1:D:226:GLU:HG3	1:D:230:LYS:NZ	2.35	0.41
1:B:178:VAL:HG11	1:B:401:PHE:CB	2.50	0.41
1:D:348:VAL:CG1	1:D:378:LEU:HD11	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:300:VAL:HG22	1:F:86:THR:HB	2.03	0.41
1:F:359:GLN:HA	1:F:369:PHE:O	2.21	0.41
1:B:154:THR:O	1:B:289:THR:HA	2.21	0.41
1:B:294:ASP:OD1	1:B:295:PRO:HD2	2.21	0.41
1:B:371:GLY:HA3	2:B:581:HOH:O	2.21	0.41
1:D:201:ASN:HB3	1:D:216:GLN:HB3	2.02	0.41
1:C:154:THR:O	1:C:289:THR:HA	2.21	0.41
1:D:195:PHE:CE2	1:E:363:ASP:HB3	2.55	0.41
1:D:371:GLY:HA3	2:D:580:HOH:O	2.21	0.41
1:E:85:GLU:OE1	1:E:144:GLN:HG3	2.20	0.41
1:E:348:VAL:CG1	1:E:378:LEU:HD11	2.50	0.41
1:F:237:ARG:HE	1:F:241:GLN:HG3	1.86	0.41
1:A:201:ASN:HB3	1:A:216:GLN:HB3	2.03	0.41
1:E:359:GLN:HA	1:E:369:PHE:O	2.21	0.41
1:B:43:VAL:HG21	1:B:348:VAL:HG21	2.03	0.40
1:C:292:LEU:HD23	1:C:292:LEU:HA	1.85	0.40
1:D:294:ASP:OD1	1:D:295:PRO:HD2	2.21	0.40
1:E:217:LYS:HE2	1:E:217:LYS:HB2	1.67	0.40
1:F:398:TRP:CD1	1:F:400:GLU:HG3	2.56	0.40
1:A:63:THR:HG22	1:A:186:ASN:CG	2.46	0.40
1:B:85:GLU:OE1	1:B:144:GLN:HG3	2.21	0.40
1:C:156:ALA:HB1	1:C:322:LEU:HD12	2.04	0.40
1:D:241:GLN:OE1	1:E:433:ILE:HG22	2.21	0.40
1:B:398:TRP:CD1	1:B:400:GLU:HG3	2.57	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	376/490 (77%)	370 (98%)	6 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	376/490 (77%)	370 (98%)	6 (2%)	0	100	100
1	C	376/490 (77%)	370 (98%)	6 (2%)	0	100	100
1	D	376/490 (77%)	370 (98%)	6 (2%)	0	100	100
1	E	376/490 (77%)	370 (98%)	6 (2%)	0	100	100
1	F	376/490 (77%)	370 (98%)	6 (2%)	0	100	100
All	All	2256/2940 (77%)	2220 (98%)	36 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	331/409 (81%)	326 (98%)	5 (2%)	57	42
1	B	331/409 (81%)	327 (99%)	4 (1%)	63	51
1	C	331/409 (81%)	327 (99%)	4 (1%)	63	51
1	D	331/409 (81%)	326 (98%)	5 (2%)	57	42
1	E	331/409 (81%)	323 (98%)	8 (2%)	43	25
1	F	331/409 (81%)	327 (99%)	4 (1%)	63	51
All	All	1986/2454 (81%)	1956 (98%)	30 (2%)	55	42

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ARG
1	A	304	GLN
1	A	408	GLN
1	A	415	ARG
1	A	446	ILE
1	B	53	LYS
1	B	237	ARG
1	B	415	ARG

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Mol	Chain	Res	Type
1	B	446	ILE
1	C	190	ARG
1	C	304	GLN
1	C	408	GLN
1	C	446	ILE
1	D	53	LYS
1	D	190	ARG
1	D	237	ARG
1	D	415	ARG
1	D	446	ILE
1	E	39	ARG
1	E	190	ARG
1	E	224	THR
1	E	320	ARG
1	E	332	ARG
1	E	409	ILE
1	E	415	ARG
1	E	446	ILE
1	F	39	ARG
1	F	190	ARG
1	F	415	ARG
1	F	446	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	HIS
1	A	333	GLN
1	A	364	ASN
1	B	34	HIS
1	B	171	GLN
1	B	248	ASN
1	B	304	GLN
1	B	364	ASN
1	B	408	GLN
1	C	171	GLN
1	C	364	ASN
1	D	34	HIS
1	D	171	GLN
1	D	184	HIS
1	D	248	ASN
1	D	304	GLN

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Mol	Chain	Res	Type
1	D	364	ASN
1	D	404	ASN
1	E	34	HIS
1	E	304	GLN
1	E	333	GLN
1	E	364	ASN
1	F	34	HIS
1	F	297	ASN
1	F	304	GLN
1	F	334	ASN
1	F	364	ASN
1	F	406	ASN

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)

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.

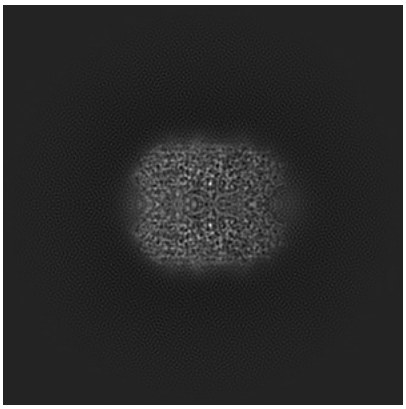
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-58942. These allow visual inspection of the internal detail of the map and identification of artifacts.

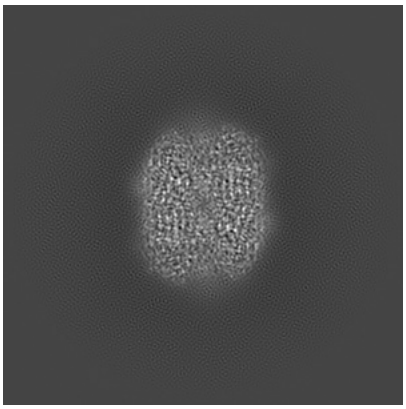
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

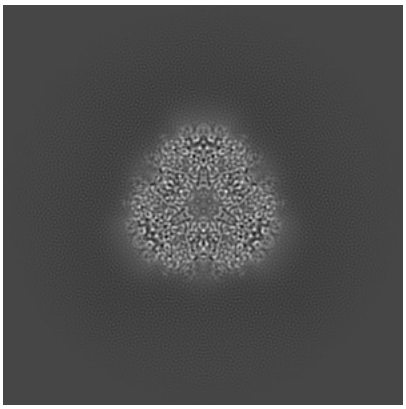
6.1.1 Primary map



X

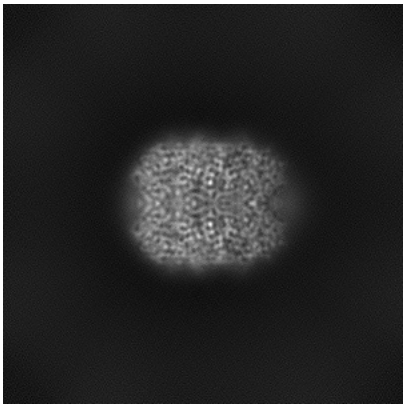


Y

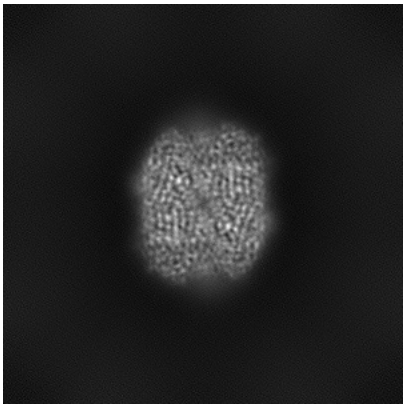


Z

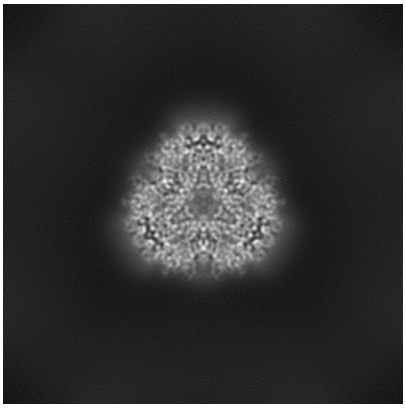
6.1.2 Raw map



X



Y

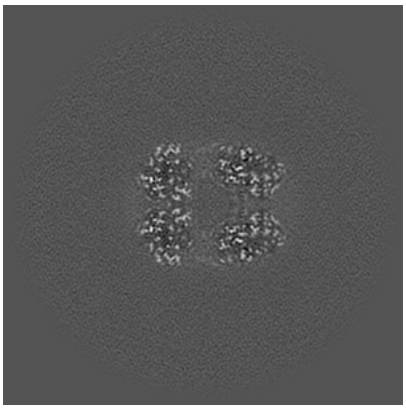


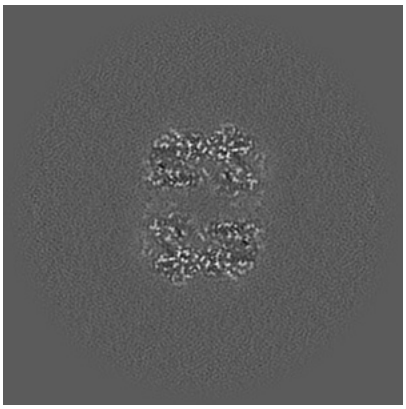
Z

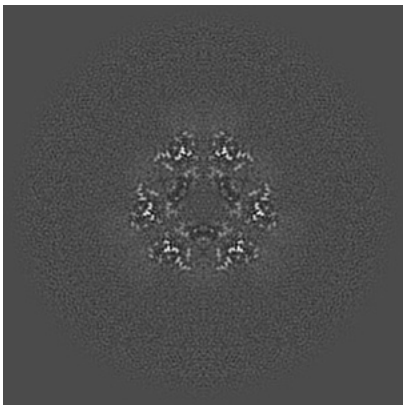
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

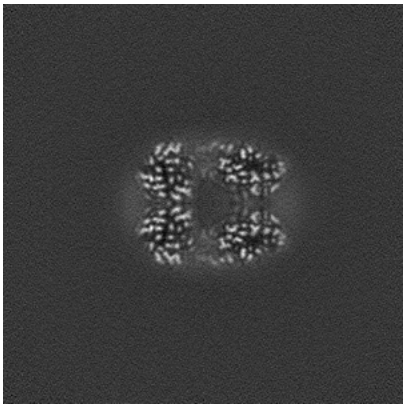
6.2.1 Primary map

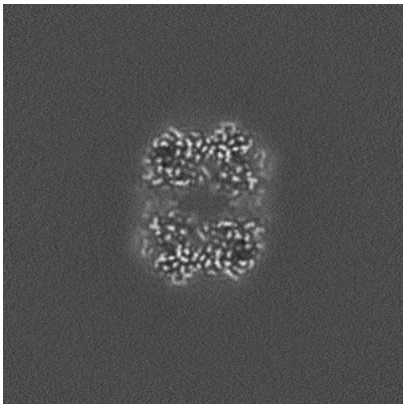

X Index: 180

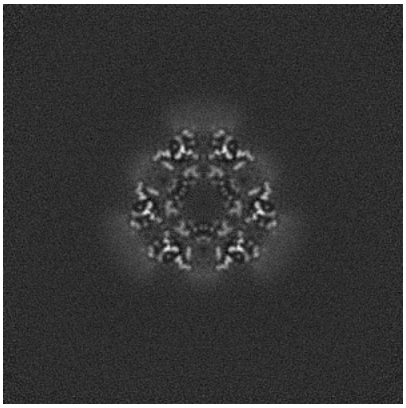

Y Index: 180


Z Index: 180

6.2.2 Raw map


X Index: 180

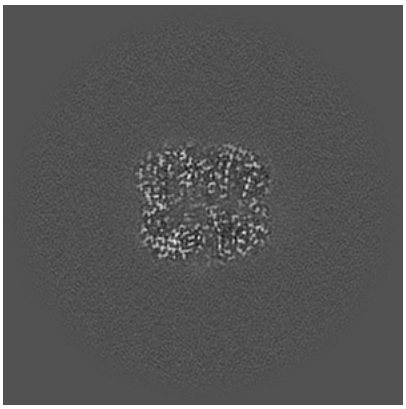

Y Index: 180

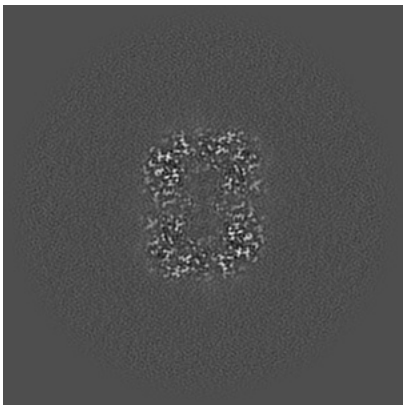

Z Index: 180

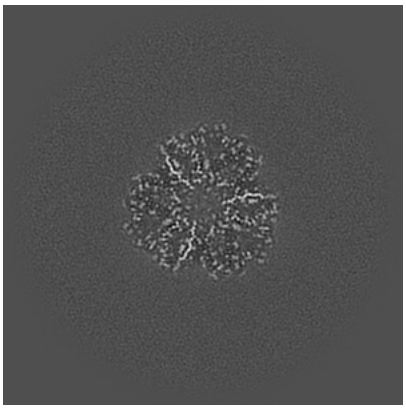
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

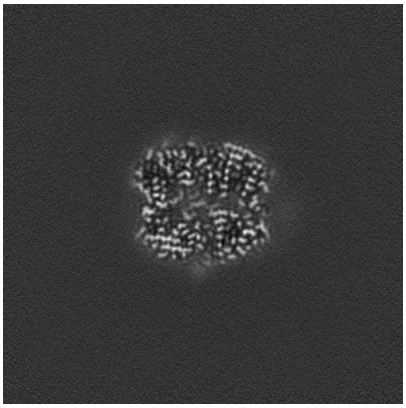
6.3.1 Primary map

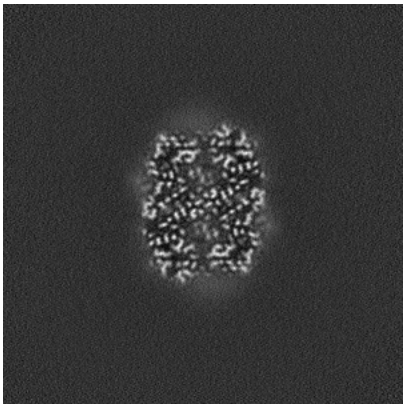

X Index: 149

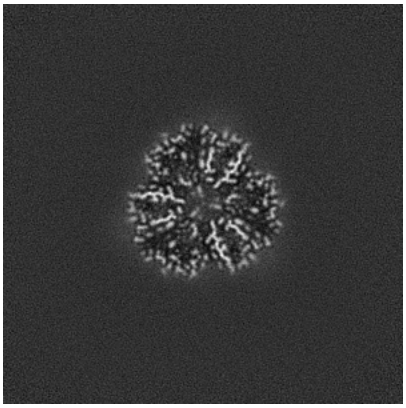

Y Index: 168


Z Index: 155

6.3.2 Raw map


X Index: 148

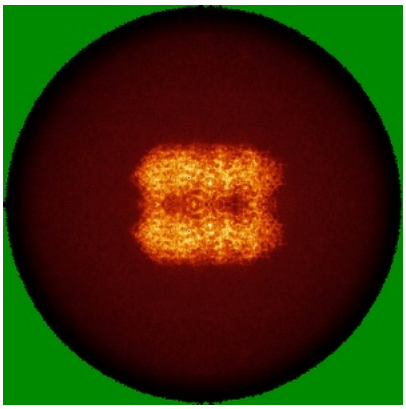

Y Index: 157

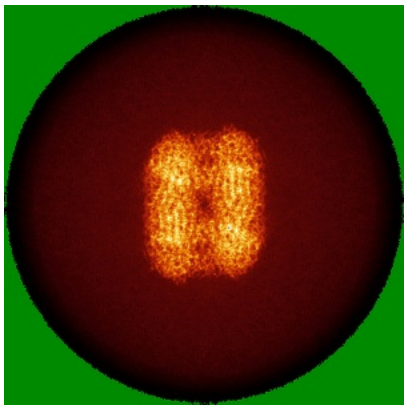

Z Index: 206

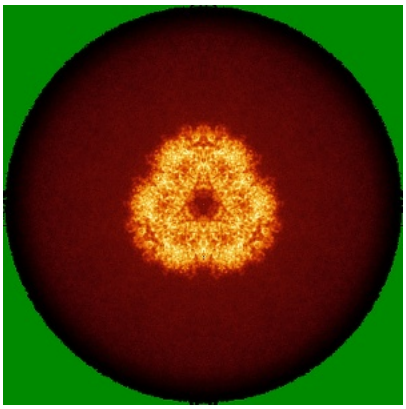
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

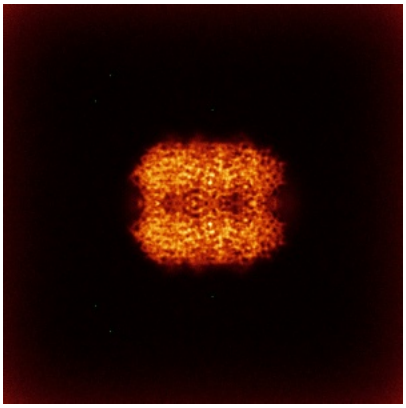
6.4.1 Primary map

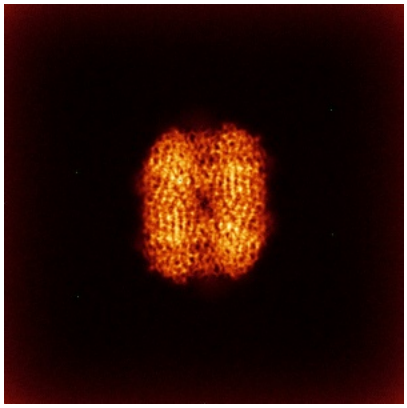

X

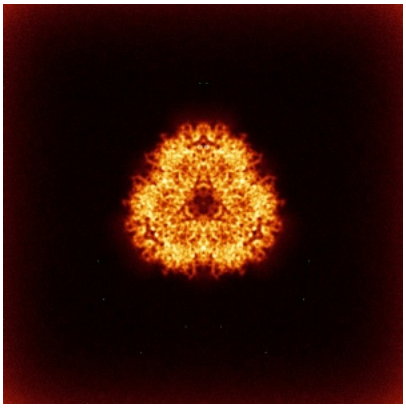

Y


Z

6.4.2 Raw map


X

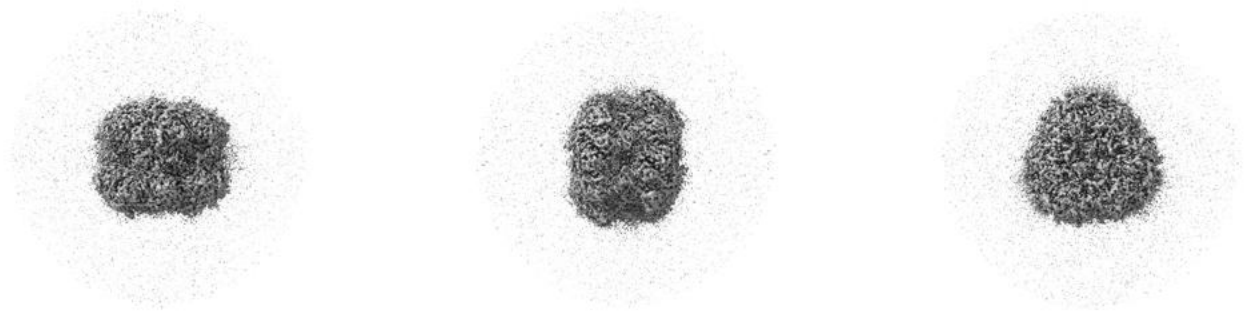

Y


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 0.05. 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.

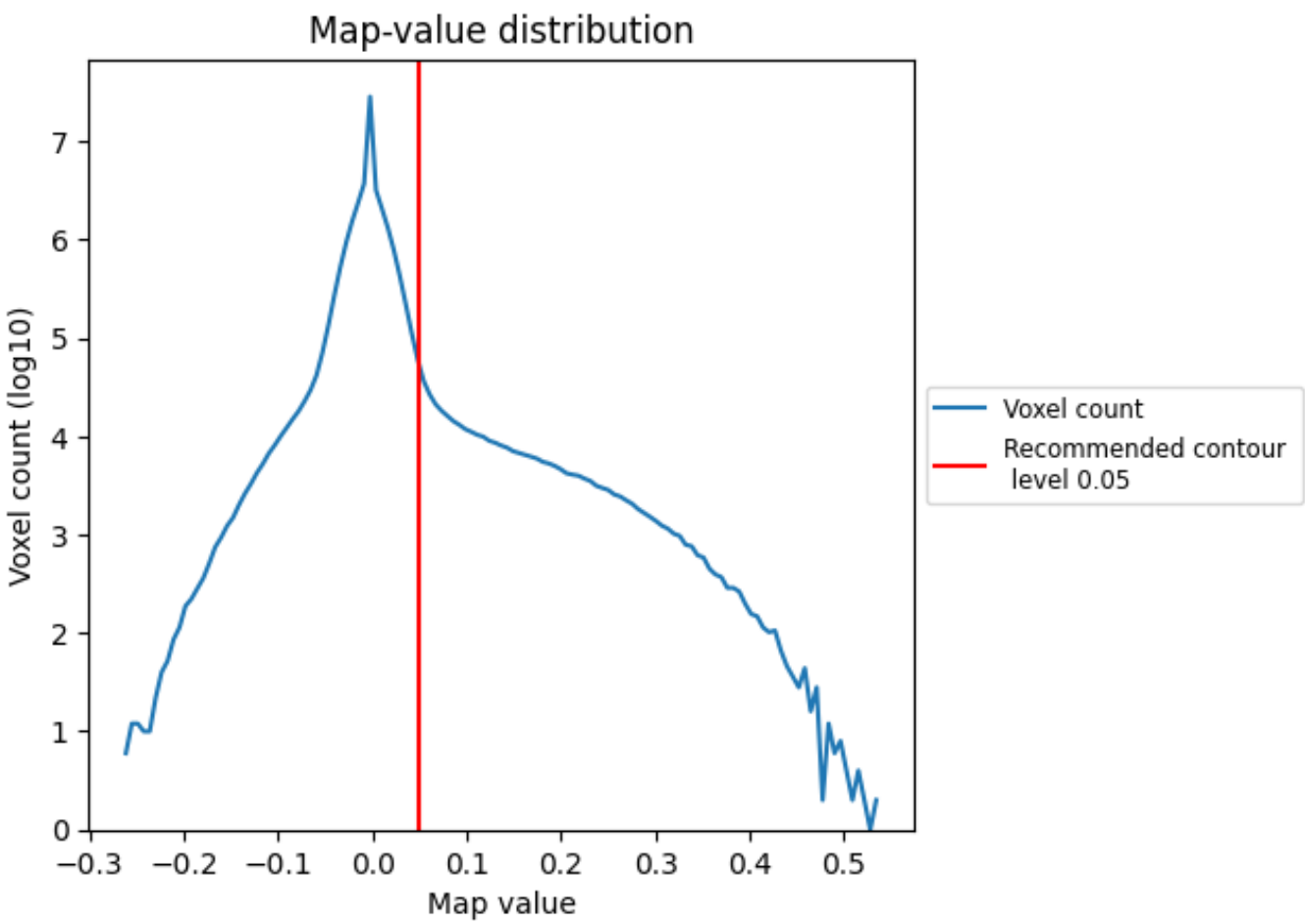
6.6 Mask visualisation (i)

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

7 Map analysis (i)

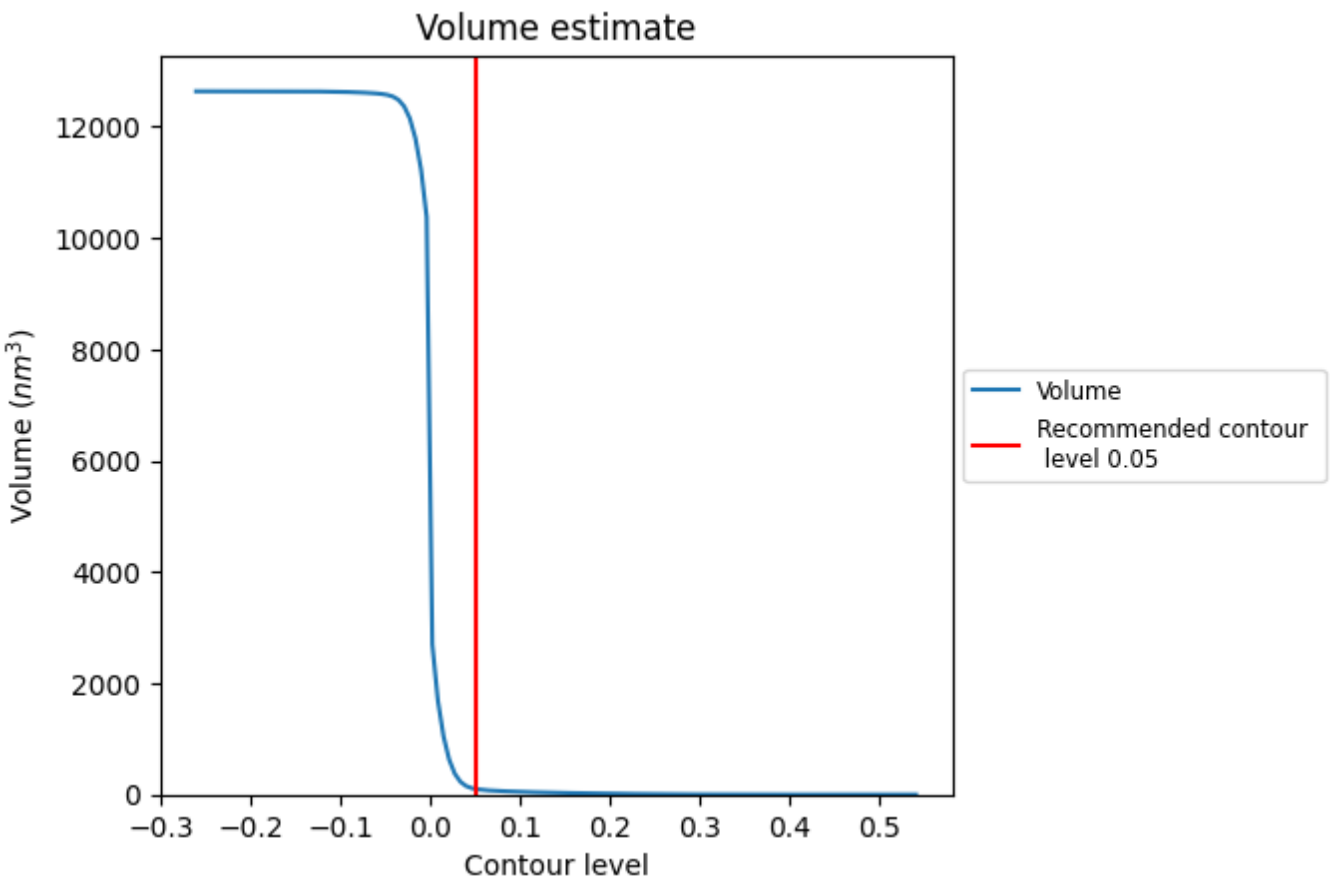
This section contains the results of statistical analysis of the map.

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.

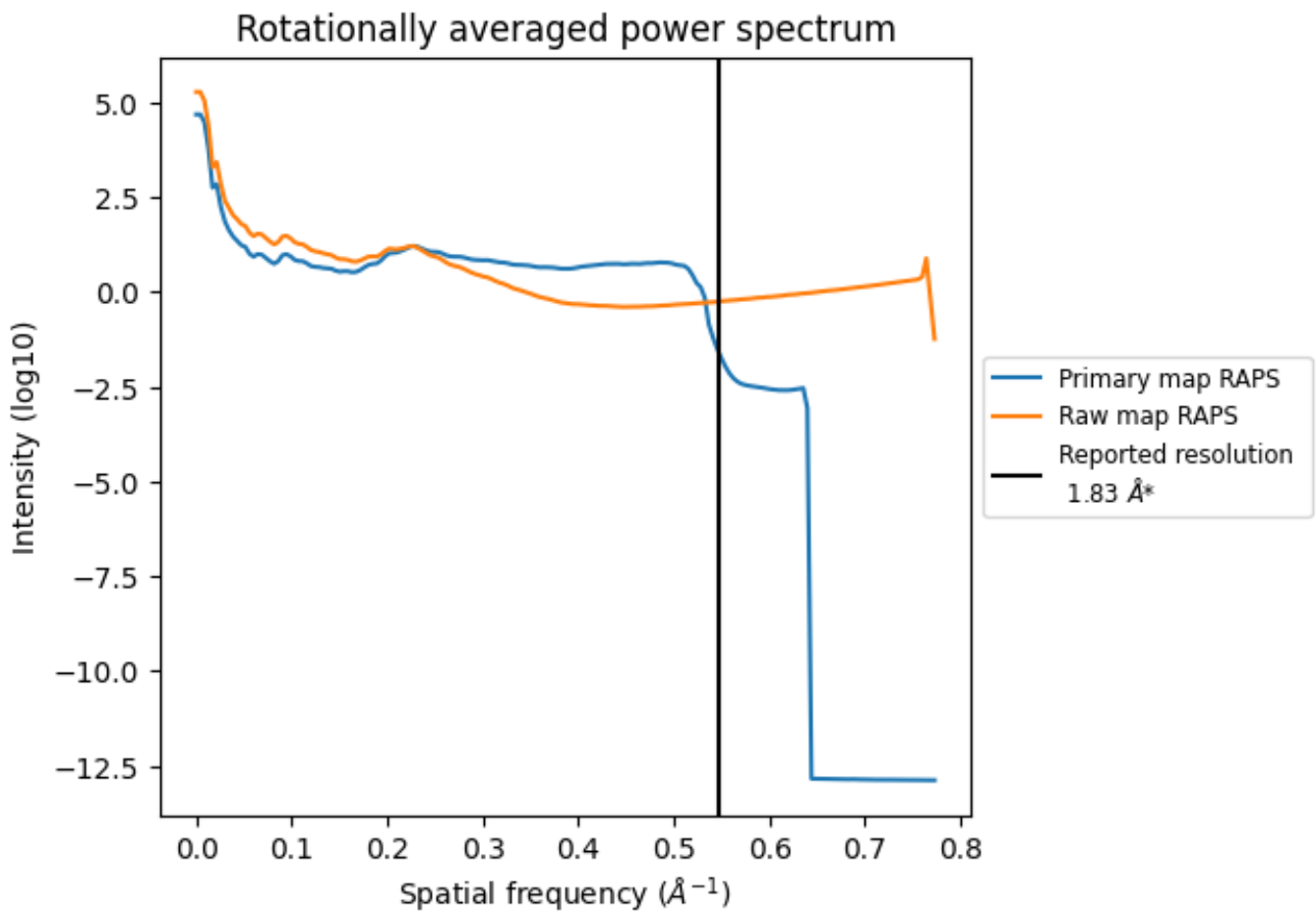
7.2 Volume estimate (i)



The volume at the recommended contour level is 102 nm^3; this corresponds to an approximate mass of 92 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 (i)

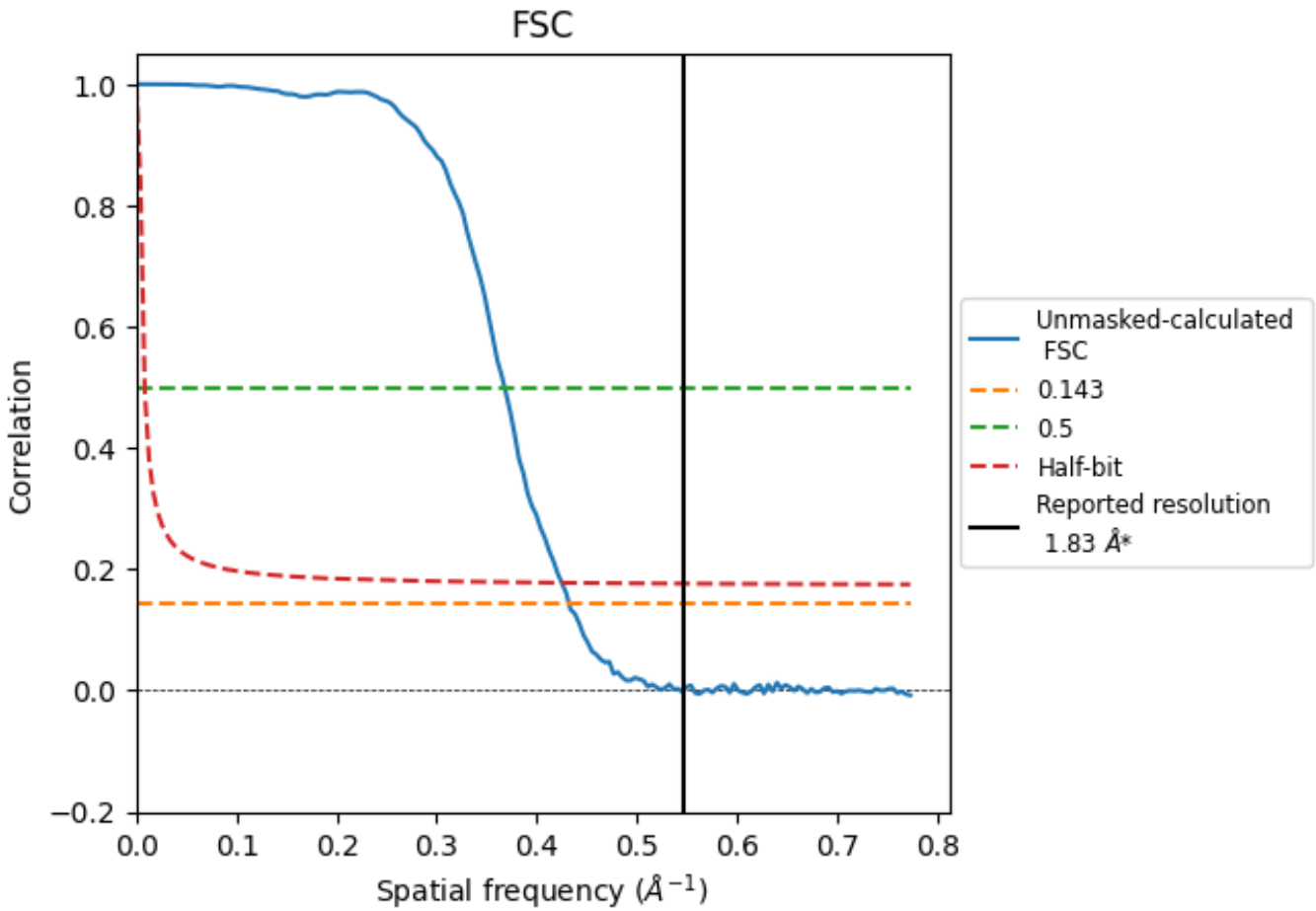


*Reported resolution corresponds to spatial frequency of 0.546 Å^{-1}

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 0.546 Å^{-1}

8.2 Resolution estimates (i)

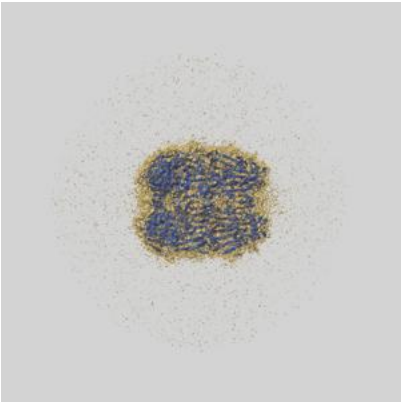
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	1.83	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	2.31	2.72	2.35

*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 2.31 differs from the reported value 1.83 by more than 10 %

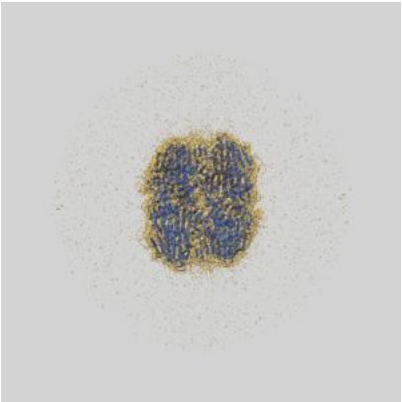
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-58942 and PDB model 32LJ. Per-residue inclusion information can be found in section 3 on page 5.

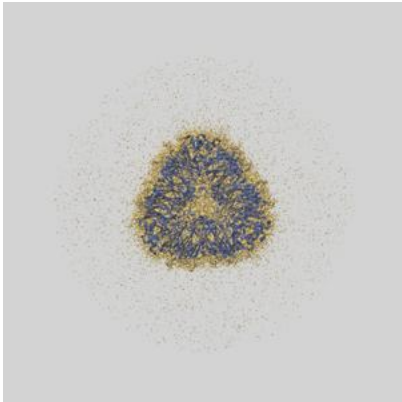
9.1 Map-model overlay (i)



X



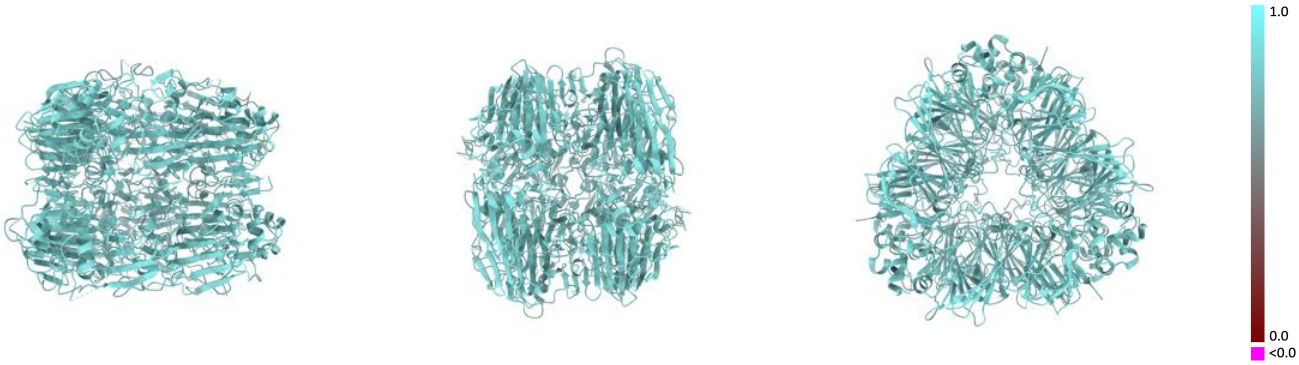
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.05 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.

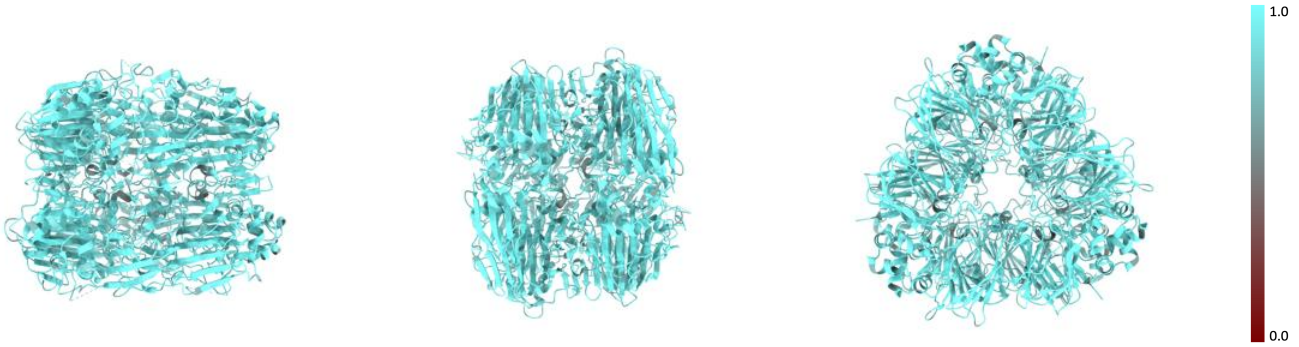
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

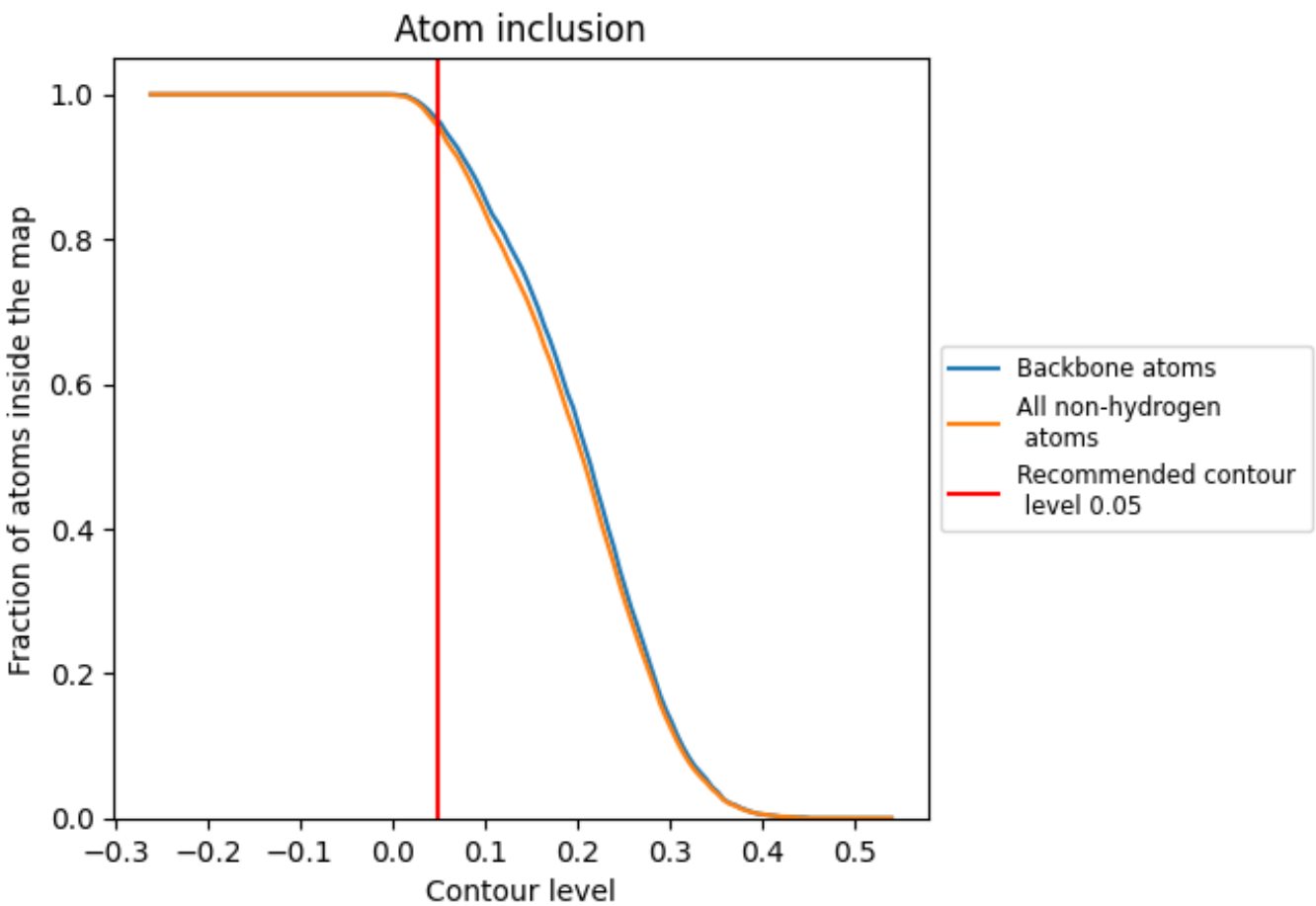
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.05).

9.4 Atom inclusion (i)



At the recommended contour level, 96% of all backbone atoms, 96% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9550	0.7810
A	0.9600	0.7810
B	0.9600	0.7820
C	0.9580	0.7820
D	0.9590	0.7820
E	0.9570	0.7810
F	0.9590	0.7810