



Full wwPDB EM Validation Report ⓘ

Mar 22, 2026 – 05:34 PM UTC

PDB ID : 7B4M / pdb_00007b4m
EMDB ID : EMD-12003
Title : CryoEM structure of the human sodium proton exchanger NHA2 in nanodisc
Authors : Woehlert, D.; Kuhlbrandt, W.; Yildiz, O.
Deposited on : 2020-12-02
Resolution : 7.20 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

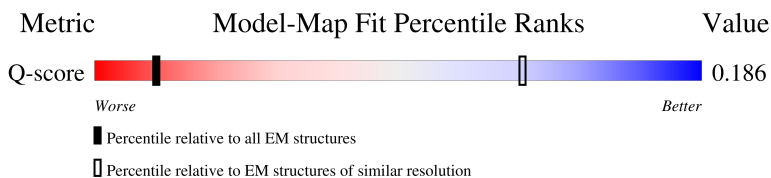
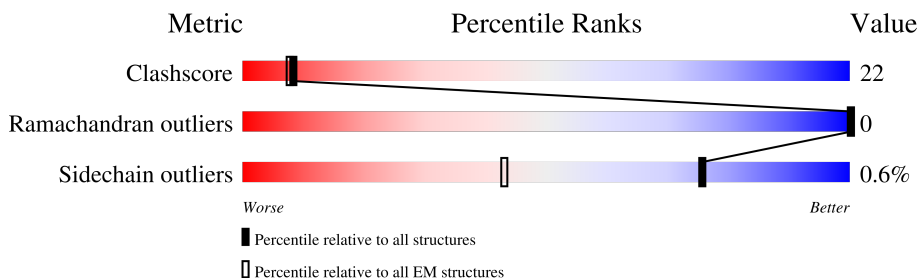
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 7.20 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	444 (6.70 - 7.70)

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	537	 49% 36% 15%
1	B	537	 6% 47% 38% 15%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6760 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.

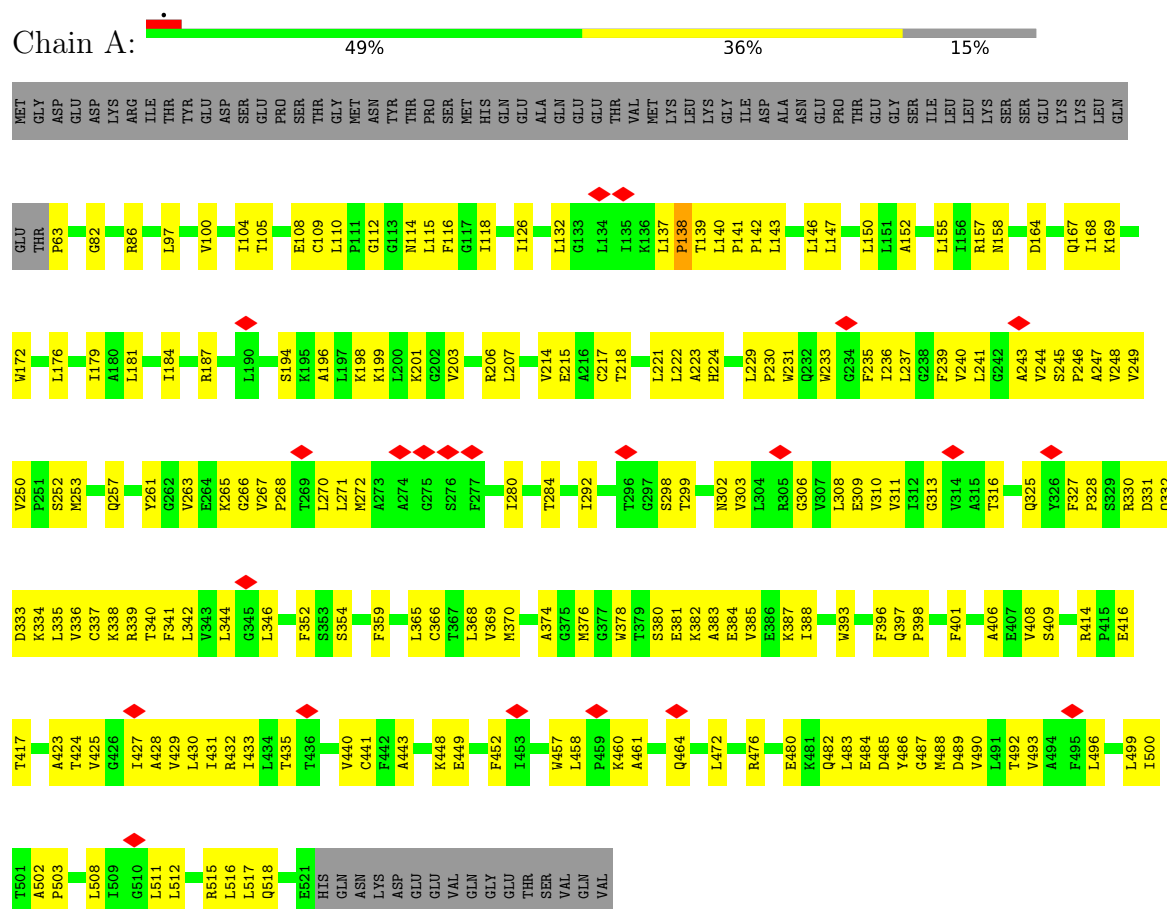
- Molecule 1 is a protein called Sodium/hydrogen exchanger 9B2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	459	Total	C	N	O	S	0	0
			3380	2237	548	577	18		
1	B	459	Total	C	N	O	S	0	0
			3380	2237	548	577	18		

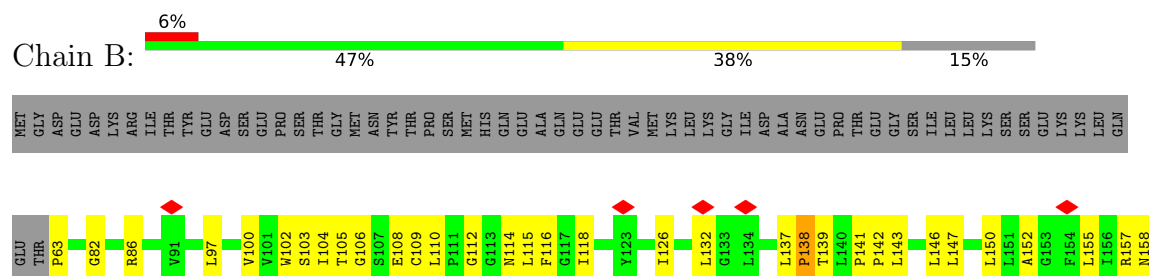
3 Residue-property plots

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: Sodium/hydrogen exchanger 9B2



• Molecule 1: Sodium/hydrogen exchanger 9B2



L486 G487 M488 D489 V490 L491 T492 V493 A494 F495 L496	L400 F401	L404 G405 A406 E407 V408 S409	L414 P415 E416 T417	L423 T424 V425 G426	L427 A428 V429 L430 I431 R432 I433 L434 T435	L438 M439 V440 C441 F442 A443	K448 E449 F452 F455 A456 W457	K460 A461 Q464 L472 R476	E480 K481 Q482 L483 E484 D485	P327 S328 S329 R330 D331 Q332 D333 K334 L335 V336 C337 K338 R339 T340 F341 L342 V343 L344 G345 L346 S347 V348 L349 S353 S354 F359 F360 G361	P246 A247 V248 V249 V250 P251 S252 M253 Q257	Y261 G262 V263 E264 K265 G266 V267 P268 T269 L270 L271 M272 A273 A274 G275 S276 F277	D278 D279 I280 T284 I292 S298 T299 N302 V303 G306 V307 L308 E309 V310 V311 I312 G313 V314 A315 T316 L320 Q325 Y326	D164 Q167 I168 K169 W172 L176 I179 A180 L181 I184 R187 A188 S194 K195 A196 L197 K198 K199 L200 K201 G202 V203 R206 L207 V214 E215 A216 C217 T218 L221 L222 A223 H224 L229 P230 W231 Q232 W233 G234 F235 I236 L237 G238 F239 V240 L241 G242 A243
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	47284	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 ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.810	Depositor
Minimum map value	-0.404	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.057	Depositor
Recommended contour level	0.27	Depositor
Map size (Å)	253.232, 253.232, 253.232	wwPDB
Map dimensions	304, 304, 304	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.833, 0.833, 0.833	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	0.15	0/3450	0.47	3/4700 (0.1%)
1	B	0.15	0/3450	0.47	3/4700 (0.1%)
All	All	0.15	0/6900	0.47	6/9400 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	63	PRO	N-CA-CB	6.71	110.38	103.00
1	B	63	PRO	N-CA-CB	6.70	110.37	103.00
1	A	138	PRO	CA-C-N	5.61	132.25	121.54
1	A	138	PRO	C-N-CA	5.61	132.25	121.54
1	B	138	PRO	CA-C-N	5.57	132.18	121.54
1	B	138	PRO	C-N-CA	5.57	132.18	121.54

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	3380	0	3555	147	0
1	B	3380	0	3555	154	0
All	All	6760	0	7110	299	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 22.

All (299) 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:480:GLU:HB2	1:A:483:LEU:HB2	1.70	0.74
1:B:480:GLU:HB2	1:B:483:LEU:HB2	1.70	0.73
1:A:230:PRO:HG2	1:A:233:TRP:HB3	1.70	0.73
1:B:230:PRO:HG2	1:B:233:TRP:HB3	1.69	0.73
1:A:218:THR:HG21	1:A:431:ILE:HG13	1.70	0.73
1:B:218:THR:HG21	1:B:431:ILE:HG13	1.70	0.72
1:B:378:TRP:O	1:B:382:LYS:N	2.22	0.71
1:B:206:ARG:NH1	1:B:441:CYS:SG	2.66	0.68
1:A:206:ARG:NH1	1:A:441:CYS:SG	2.67	0.68
1:A:214:VAL:HG21	1:A:435:THR:HG21	1.76	0.67
1:B:214:VAL:HG21	1:B:435:THR:HG21	1.76	0.66
1:A:378:TRP:O	1:A:382:LYS:N	2.22	0.66
1:B:229:LEU:HD22	1:B:233:TRP:HZ3	1.61	0.65
1:A:383:ALA:O	1:A:387:LYS:NZ	2.29	0.65
1:A:229:LEU:HD22	1:A:233:TRP:HZ3	1.61	0.65
1:A:147:LEU:HD13	1:A:503:PRO:HG2	1.79	0.65
1:B:147:LEU:HD13	1:B:503:PRO:HG2	1.79	0.64
1:B:157:ARG:HB2	1:B:406:ALA:HA	1.81	0.63
1:A:499:LEU:HD12	1:A:500:ILE:HB	1.81	0.63
1:B:383:ALA:O	1:B:387:LYS:NZ	2.29	0.62
1:A:340:THR:HG21	1:A:388:ILE:HD12	1.81	0.62
1:B:316:THR:OG1	1:B:366:CYS:SG	2.57	0.62
1:A:327:PHE:HA	1:A:330:ARG:HH22	1.65	0.62
1:A:316:THR:OG1	1:A:366:CYS:SG	2.57	0.62
1:B:327:PHE:HA	1:B:330:ARG:HH22	1.65	0.62
1:A:157:ARG:HB2	1:A:406:ALA:HA	1.81	0.62
1:A:280:ILE:HD11	1:A:365:LEU:HD21	1.81	0.61
1:B:340:THR:HG21	1:B:388:ILE:HD12	1.81	0.61
1:B:499:LEU:HD12	1:B:500:ILE:HB	1.81	0.61
1:B:184:ILE:HD11	1:B:250:VAL:HB	1.83	0.61
1:A:146:LEU:HD22	1:A:401:PHE:HE2	1.65	0.61
1:B:146:LEU:HD22	1:B:401:PHE:HE2	1.66	0.61
1:B:280:ILE:HD11	1:B:365:LEU:HD21	1.81	0.60
1:B:229:LEU:HD22	1:B:233:TRP:CZ3	2.36	0.60
1:A:184:ILE:HD11	1:A:250:VAL:HB	1.83	0.60
1:A:229:LEU:HD22	1:A:233:TRP:CZ3	2.36	0.60
1:B:490:VAL:HA	1:B:493:VAL:HG12	1.84	0.60
1:B:215:GLU:HG3	1:B:428:ALA:HB1	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:483:LEU:HD23	1:B:486:TYR:HD2	1.67	0.59
1:A:490:VAL:HA	1:A:493:VAL:HG12	1.84	0.59
1:B:126:ILE:HG22	1:B:398:PRO:HG2	1.85	0.59
1:A:340:THR:O	1:A:344:LEU:HG	2.03	0.58
1:B:340:THR:O	1:B:344:LEU:HG	2.03	0.58
1:A:483:LEU:HD23	1:A:486:TYR:HD2	1.67	0.58
1:A:132:LEU:HD23	1:A:143:LEU:HB3	1.85	0.58
1:B:137:LEU:C	1:B:139:THR:H	2.10	0.58
1:A:308:LEU:HA	1:A:311:VAL:HG22	1.86	0.58
1:B:425:VAL:O	1:B:429:VAL:HG23	2.04	0.58
1:A:126:ILE:HG22	1:A:398:PRO:HG2	1.85	0.58
1:A:425:VAL:O	1:A:429:VAL:HG23	2.04	0.58
1:B:308:LEU:HA	1:B:311:VAL:HG22	1.86	0.58
1:B:334:LYS:HB3	1:B:337:CYS:HB2	1.86	0.57
1:A:215:GLU:HG3	1:A:428:ALA:HB1	1.85	0.57
1:B:172:TRP:O	1:B:176:LEU:HG	2.05	0.57
1:A:137:LEU:C	1:A:139:THR:H	2.10	0.57
1:B:132:LEU:HD23	1:B:143:LEU:HB3	1.86	0.57
1:A:487:GLY:HA2	1:A:490:VAL:HG22	1.87	0.57
1:B:198:LYS:O	1:B:201:LYS:NZ	2.32	0.57
1:A:334:LYS:HB3	1:A:337:CYS:HB2	1.86	0.57
1:B:430:LEU:HD13	1:B:433:ILE:HD12	1.86	0.57
1:A:352:PHE:HB3	1:B:103:SER:HB2	1.87	0.57
1:A:109:CYS:HA	1:A:116:PHE:HB2	1.86	0.56
1:A:241:LEU:HD11	1:A:425:VAL:HG23	1.87	0.56
1:A:284:THR:HG23	1:A:306:GLY:HA3	1.86	0.56
1:B:284:THR:HG23	1:B:306:GLY:HA3	1.86	0.56
1:A:485:ASP:HA	1:A:488:MET:HE3	1.87	0.56
1:B:169:LYS:HG2	1:B:172:TRP:H	1.71	0.56
1:A:172:TRP:O	1:A:176:LEU:HG	2.05	0.56
1:A:432:ARG:NE	1:A:457:TRP:O	2.38	0.56
1:B:109:CYS:HA	1:B:116:PHE:HB2	1.86	0.56
1:B:342:LEU:O	1:B:346:LEU:HG	2.06	0.56
1:B:487:GLY:HA2	1:B:490:VAL:HG22	1.87	0.56
1:A:169:LYS:HG2	1:A:172:TRP:H	1.70	0.56
1:A:430:LEU:HD13	1:A:433:ILE:HD12	1.87	0.56
1:B:432:ARG:NE	1:B:457:TRP:O	2.38	0.55
1:B:485:ASP:HA	1:B:488:MET:HE3	1.87	0.55
1:A:155:LEU:HA	1:A:158:ASN:HD22	1.72	0.55
1:A:427:ILE:O	1:A:431:ILE:HG12	2.07	0.55
1:A:198:LYS:O	1:A:201:LYS:NZ	2.32	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:LEU:HA	1:B:158:ASN:HD22	1.72	0.55
1:B:241:LEU:HD11	1:B:425:VAL:HG23	1.87	0.55
1:A:338:LYS:HA	1:A:341:PHE:CE2	2.43	0.54
1:A:203:VAL:O	1:A:207:LEU:HG	2.07	0.54
1:A:342:LEU:O	1:A:346:LEU:HG	2.06	0.54
1:A:233:TRP:NE1	1:A:236:ILE:HD12	2.21	0.54
1:B:233:TRP:NE1	1:B:236:ILE:HD12	2.21	0.54
1:A:328:PRO:HG2	1:A:339:ARG:NH2	2.23	0.54
1:B:184:ILE:HD12	1:B:187:ARG:HE	1.73	0.54
1:B:203:VAL:O	1:B:207:LEU:HG	2.08	0.53
1:B:280:ILE:HG21	1:B:310:VAL:HG22	1.90	0.53
1:B:338:LYS:HA	1:B:341:PHE:CE2	2.43	0.53
1:A:440:VAL:HG13	1:A:443:ALA:HB3	1.90	0.53
1:A:298:SER:O	1:A:302:ASN:ND2	2.42	0.53
1:B:427:ILE:O	1:B:431:ILE:HG12	2.07	0.53
1:A:184:ILE:HD12	1:A:187:ARG:HE	1.73	0.53
1:A:224:HIS:HA	1:A:229:LEU:H	1.74	0.53
1:B:137:LEU:O	1:B:139:THR:N	2.38	0.53
1:B:328:PRO:HG2	1:B:339:ARG:NH2	2.24	0.52
1:B:440:VAL:HG13	1:B:443:ALA:HB3	1.90	0.52
1:A:164:ASP:O	1:A:167:GLN:NE2	2.43	0.52
1:A:146:LEU:HD12	1:A:248:VAL:HG23	1.91	0.52
1:A:261:TYR:HB3	1:A:517:LEU:HB2	1.92	0.52
1:B:224:HIS:HA	1:B:229:LEU:H	1.74	0.52
1:B:261:TYR:HB3	1:B:517:LEU:HB2	1.92	0.52
1:A:249:VAL:O	1:A:252:SER:OG	2.26	0.52
1:A:280:ILE:HG21	1:A:310:VAL:HG22	1.90	0.52
1:B:298:SER:O	1:B:302:ASN:ND2	2.41	0.51
1:A:414:ARG:NH2	1:A:489:ASP:OD1	2.42	0.51
1:B:146:LEU:HD12	1:B:248:VAL:HG23	1.90	0.51
1:B:414:ARG:NH2	1:B:489:ASP:OD1	2.42	0.51
1:B:164:ASP:O	1:B:167:GLN:NE2	2.43	0.51
1:B:512:LEU:HD12	1:B:516:LEU:HD11	1.92	0.51
1:A:146:LEU:HD22	1:A:401:PHE:CE2	2.46	0.51
1:A:265:LYS:HB3	1:A:267:VAL:HG23	1.93	0.51
1:A:224:HIS:HE2	1:A:231:TRP:CG	2.29	0.50
1:B:146:LEU:HD22	1:B:401:PHE:CE2	2.46	0.50
1:B:184:ILE:HG12	1:B:247:ALA:HA	1.94	0.50
1:B:224:HIS:HE2	1:B:231:TRP:CG	2.30	0.50
1:A:512:LEU:HD12	1:A:516:LEU:HD11	1.92	0.50
1:A:452:PHE:HD1	1:A:516:LEU:HD12	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:LYS:HB3	1:B:267:VAL:HG23	1.93	0.50
1:B:354:SER:HB2	1:B:359:PHE:O	2.11	0.50
1:B:233:TRP:CZ2	1:B:486:TYR:HB3	2.47	0.50
1:B:452:PHE:HD1	1:B:516:LEU:HD12	1.77	0.50
1:B:267:VAL:O	1:B:271:LEU:HG	2.12	0.50
1:A:184:ILE:HG12	1:A:247:ALA:HA	1.94	0.50
1:A:309:GLU:HB2	1:A:359:PHE:HB3	1.93	0.50
1:A:152:ALA:O	1:A:155:LEU:HG	2.12	0.49
1:A:196:ALA:HA	1:A:199:LYS:HE3	1.94	0.49
1:A:354:SER:HB2	1:A:359:PHE:O	2.11	0.49
1:A:233:TRP:CZ2	1:A:486:TYR:HB3	2.47	0.49
1:B:448:LYS:HB3	1:B:516:LEU:HB3	1.94	0.49
1:B:104:ILE:HG22	1:B:105:THR:HG23	1.95	0.49
1:B:152:ALA:O	1:B:155:LEU:HG	2.12	0.49
1:A:480:GLU:HB3	1:A:482:GLN:HE22	1.76	0.49
1:B:309:GLU:HB2	1:B:359:PHE:HB3	1.93	0.49
1:A:313:GLY:HA3	1:A:365:LEU:HD12	1.95	0.49
1:B:196:ALA:HA	1:B:199:LYS:HE3	1.94	0.49
1:A:115:LEU:HD23	1:A:118:ILE:HD12	1.95	0.49
1:B:313:GLY:HA3	1:B:365:LEU:HD12	1.95	0.49
1:A:110:LEU:C	1:A:112:GLY:H	2.21	0.49
1:B:393:TRP:NE1	1:B:397:GLN:HE21	2.11	0.49
1:A:104:ILE:HG22	1:A:105:THR:HG23	1.95	0.48
1:A:397:GLN:HB3	1:A:401:PHE:CE2	2.48	0.48
1:A:448:LYS:HB3	1:A:516:LEU:HB3	1.94	0.48
1:B:115:LEU:HD23	1:B:118:ILE:HD12	1.95	0.48
1:B:333:ASP:OD1	1:B:333:ASP:N	2.46	0.48
1:B:215:GLU:OE2	1:B:460:LYS:NZ	2.46	0.48
1:B:480:GLU:HB3	1:B:482:GLN:HE22	1.76	0.48
1:A:137:LEU:HD22	1:A:142:PRO:HA	1.96	0.48
1:A:267:VAL:O	1:A:271:LEU:HG	2.12	0.48
1:B:334:LYS:O	1:B:338:LYS:N	2.41	0.48
1:A:137:LEU:O	1:A:139:THR:N	2.38	0.48
1:A:460:LYS:HB3	1:A:464:GLN:HE22	1.78	0.48
1:B:115:LEU:HD21	1:B:172:TRP:HZ3	1.79	0.48
1:B:248:VAL:HG11	1:B:502:ALA:HB1	1.95	0.48
1:B:492:THR:O	1:B:496:LEU:HG	2.14	0.48
1:A:248:VAL:HG11	1:A:502:ALA:HB1	1.95	0.48
1:B:397:GLN:HB3	1:B:401:PHE:CE2	2.48	0.48
1:A:333:ASP:OD1	1:A:333:ASP:N	2.46	0.48
1:A:380:SER:OG	1:A:381:GLU:OE1	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:460:LYS:HB3	1:B:464:GLN:HE22	1.78	0.48
1:A:115:LEU:HD21	1:A:172:TRP:HZ3	1.79	0.48
1:A:492:THR:O	1:A:496:LEU:HG	2.14	0.48
1:B:169:LYS:HD3	1:B:172:TRP:CG	2.49	0.47
1:B:449:GLU:OE1	1:B:449:GLU:N	2.46	0.47
1:A:215:GLU:OE2	1:A:460:LYS:NZ	2.46	0.47
1:A:449:GLU:OE1	1:A:449:GLU:N	2.46	0.47
1:B:250:VAL:HG13	1:B:272:MET:SD	2.54	0.47
1:B:137:LEU:HD22	1:B:142:PRO:HA	1.96	0.47
1:B:138:PRO:O	1:B:139:THR:HG22	2.15	0.47
1:B:110:LEU:C	1:B:112:GLY:H	2.21	0.47
1:A:169:LYS:HD3	1:A:172:TRP:CG	2.49	0.47
1:A:250:VAL:HG13	1:A:272:MET:SD	2.54	0.47
1:A:393:TRP:NE1	1:A:397:GLN:HE21	2.11	0.47
1:B:150:LEU:HB3	1:B:401:PHE:HB3	1.96	0.47
1:A:138:PRO:O	1:A:139:THR:HG22	2.15	0.47
1:B:378:TRP:O	1:B:381:GLU:N	2.48	0.47
1:A:378:TRP:O	1:A:381:GLU:N	2.48	0.46
1:A:508:LEU:HB2	1:A:512:LEU:HD23	1.97	0.46
1:B:380:SER:OG	1:B:381:GLU:OE1	2.27	0.46
1:A:214:VAL:HA	1:A:217:CYS:SG	2.55	0.46
1:A:374:ALA:O	1:A:378:TRP:N	2.42	0.46
1:B:214:VAL:HA	1:B:217:CYS:SG	2.55	0.46
1:A:150:LEU:HB3	1:A:401:PHE:HB3	1.97	0.46
1:A:448:LYS:HD3	1:A:516:LEU:HA	1.97	0.46
1:B:292:ILE:HD11	1:B:299:THR:HG21	1.98	0.46
1:B:448:LYS:HD3	1:B:516:LEU:HA	1.97	0.46
1:A:292:ILE:HD11	1:A:299:THR:HG21	1.98	0.46
1:B:472:LEU:HG	1:B:476:ARG:CZ	2.46	0.46
1:B:508:LEU:HB2	1:B:512:LEU:HD23	1.97	0.45
1:A:472:LEU:HG	1:A:476:ARG:CZ	2.46	0.45
1:B:102:TRP:O	1:B:106:GLY:N	2.42	0.45
1:A:339:ARG:HB3	1:A:378:TRP:CH2	2.52	0.45
1:B:339:ARG:HB3	1:B:378:TRP:CH2	2.52	0.45
1:A:215:GLU:CG	1:A:428:ALA:HB1	2.47	0.45
1:A:378:TRP:HB3	1:A:381:GLU:HG2	1.99	0.45
1:A:485:ASP:O	1:A:488:MET:HG2	2.16	0.45
1:A:266:GLY:O	1:A:270:LEU:HG	2.17	0.44
1:B:485:ASP:O	1:B:488:MET:HG2	2.17	0.44
1:A:181:LEU:O	1:A:184:ILE:HG22	2.18	0.44
1:A:235:PHE:HB3	1:A:239:PHE:CZ	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:LEU:HB3	1:B:141:PRO:O	2.18	0.44
1:B:249:VAL:O	1:B:252:SER:OG	2.26	0.44
1:B:266:GLY:O	1:B:270:LEU:HG	2.17	0.44
1:B:344:LEU:O	1:B:347:SER:OG	2.33	0.44
1:B:243:ALA:HB3	1:B:464:GLN:NE2	2.33	0.44
1:B:235:PHE:HB3	1:B:239:PHE:CZ	2.52	0.44
1:B:378:TRP:HB3	1:B:381:GLU:HG2	1.99	0.44
1:A:222:LEU:HD13	1:A:424:THR:HA	2.00	0.44
1:A:370:MET:HE2	1:A:370:MET:HB3	1.94	0.44
1:A:518:GLN:N	1:A:518:GLN:OE1	2.51	0.44
1:B:518:GLN:OE1	1:B:518:GLN:N	2.51	0.44
1:B:223:ALA:C	1:B:229:LEU:HG	2.43	0.43
1:A:257:GLN:HG2	1:A:263:VAL:HG11	2.00	0.43
1:B:181:LEU:O	1:B:184:ILE:HG22	2.18	0.43
1:A:243:ALA:HB3	1:A:464:GLN:NE2	2.33	0.43
1:B:222:LEU:HD13	1:B:424:THR:HA	2.00	0.43
1:A:223:ALA:HB3	1:A:229:LEU:HD11	2.01	0.43
1:B:155:LEU:HA	1:B:158:ASN:ND2	2.33	0.43
1:B:215:GLU:CG	1:B:428:ALA:HB1	2.47	0.43
1:B:336:VAL:O	1:B:340:THR:HG23	2.19	0.43
1:A:336:VAL:O	1:A:340:THR:HG23	2.19	0.43
1:A:137:LEU:HB3	1:A:141:PRO:O	2.18	0.43
1:A:223:ALA:C	1:A:229:LEU:HG	2.43	0.43
1:A:244:VAL:HG13	1:A:246:PRO:HD3	2.01	0.43
1:A:332:GLN:HB2	1:A:335:LEU:HD13	2.00	0.43
1:B:97:LEU:HD12	1:B:100:VAL:HB	2.00	0.43
1:B:244:VAL:HG13	1:B:246:PRO:HD3	2.01	0.43
1:B:257:GLN:HG2	1:B:263:VAL:HG11	2.00	0.43
1:B:370:MET:HE2	1:B:370:MET:HB3	1.94	0.43
1:B:393:TRP:CD1	1:B:397:GLN:HE21	2.37	0.43
1:A:155:LEU:HA	1:A:158:ASN:ND2	2.33	0.43
1:A:334:LYS:O	1:A:338:LYS:N	2.41	0.43
1:A:393:TRP:CD1	1:A:397:GLN:HE21	2.37	0.43
1:B:253:MET:SD	1:B:253:MET:N	2.91	0.43
1:A:416:GLU:O	1:A:417:THR:OG1	2.29	0.42
1:B:299:THR:O	1:B:303:VAL:HG23	2.19	0.42
1:A:253:MET:N	1:A:253:MET:SD	2.91	0.42
1:B:349:LEU:O	1:B:353:SER:OG	2.33	0.42
1:A:97:LEU:HD12	1:A:100:VAL:HB	2.00	0.42
1:A:245:SER:HB2	1:A:461:ALA:H	1.84	0.42
1:A:268:PRO:O	1:A:272:MET:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:GLU:OE1	1:A:417:THR:HG23	2.20	0.42
1:A:299:THR:HA	1:A:302:ASN:HD22	1.85	0.42
1:B:223:ALA:HB3	1:B:229:LEU:HD11	2.01	0.42
1:B:332:GLN:HB2	1:B:335:LEU:HD13	2.00	0.42
1:A:100:VAL:O	1:A:104:ILE:HG13	2.20	0.42
1:A:299:THR:O	1:A:303:VAL:HG23	2.19	0.42
1:B:176:LEU:HA	1:B:179:ILE:HG12	2.02	0.42
1:B:460:LYS:O	1:B:501:THR:OG1	2.36	0.42
1:B:408:VAL:HG22	1:B:409:SER:H	1.85	0.42
1:A:82:GLY:O	1:A:86:ARG:HD3	2.20	0.42
1:B:299:THR:HA	1:B:302:ASN:HD22	1.85	0.42
1:B:382:LYS:HA	1:B:385:VAL:HG22	2.02	0.42
1:A:408:VAL:HG22	1:A:409:SER:H	1.85	0.41
1:B:245:SER:HB2	1:B:461:ALA:H	1.84	0.41
1:B:384:GLU:HA	1:B:387:LYS:HE2	2.02	0.41
1:B:268:PRO:O	1:B:272:MET:HG2	2.19	0.41
1:A:382:LYS:HA	1:A:385:VAL:HG22	2.02	0.41
1:A:176:LEU:HA	1:A:179:ILE:HG12	2.02	0.41
1:B:114:ASN:OD1	1:B:168:ILE:HA	2.20	0.41
1:A:253:MET:HG3	1:A:272:MET:HB3	2.02	0.41
1:B:237:LEU:HA	1:B:240:VAL:HG12	2.02	0.41
1:B:482:GLN:OE1	1:B:482:GLN:N	2.50	0.41
1:A:237:LEU:HA	1:A:240:VAL:HG12	2.02	0.41
1:A:331:ASP:CB	1:B:86:ARG:HH11	2.34	0.41
1:B:221:LEU:HD23	1:B:221:LEU:HA	1.82	0.41
1:A:105:THR:OG1	1:A:108:GLU:HB2	2.21	0.41
1:A:140:LEU:HD12	1:A:140:LEU:HA	1.90	0.41
1:A:393:TRP:CZ3	1:A:396:PHE:HB2	2.56	0.41
1:B:100:VAL:O	1:B:104:ILE:HG13	2.20	0.41
1:B:105:THR:OG1	1:B:108:GLU:HB2	2.21	0.41
1:B:334:LYS:HB2	1:B:338:LYS:HE3	2.02	0.41
1:B:416:GLU:OE1	1:B:417:THR:HG23	2.20	0.41
1:A:325:GLN:NE2	1:A:376:MET:HB2	2.36	0.41
1:A:480:GLU:O	1:A:484:GLU:N	2.54	0.41
1:A:511:LEU:O	1:A:515:ARG:HG2	2.21	0.41
1:B:253:MET:HG3	1:B:272:MET:HB3	2.02	0.41
1:B:423:ALA:O	1:B:427:ILE:HG12	2.21	0.41
1:A:221:LEU:HA	1:A:221:LEU:HD23	1.82	0.40
1:B:393:TRP:CZ3	1:B:396:PHE:HB2	2.56	0.40
1:B:82:GLY:O	1:B:86:ARG:HD3	2.20	0.40
1:B:138:PRO:HG2	1:B:141:PRO:HG2	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:LEU:O	1:A:369:VAL:HG12	2.21	0.40
1:A:384:GLU:HA	1:A:387:LYS:HE2	2.02	0.40
1:A:458:LEU:O	1:A:460:LYS:HG3	2.21	0.40
1:B:262:GLY:C	1:B:268:PRO:HD3	2.47	0.40
1:B:320:LEU:H	1:B:320:LEU:HD12	1.87	0.40
1:B:325:GLN:NE2	1:B:376:MET:HB2	2.36	0.40
1:B:365:LEU:O	1:B:369:VAL:HG12	2.21	0.40
1:B:374:ALA:HB1	1:B:378:TRP:CE2	2.57	0.40
1:B:511:LEU:O	1:B:515:ARG:HG2	2.21	0.40
1:A:114:ASN:OD1	1:A:168:ILE:HA	2.20	0.40
1:A:184:ILE:HD12	1:A:184:ILE:HA	1.87	0.40
1:A:423:ALA:O	1:A:427:ILE:HG12	2.21	0.40
1:B:310:VAL:O	1:B:314:VAL:HG13	2.22	0.40
1:B:399:LEU:HD12	1:B:400:LEU:N	2.37	0.40
1:B:480:GLU:O	1:B:484:GLU:N	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	457/537 (85%)	425 (93%)	32 (7%)	0	100	100
1	B	457/537 (85%)	425 (93%)	32 (7%)	0	100	100
All	All	914/1074 (85%)	850 (93%)	64 (7%)	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	357/444 (80%)	355 (99%)	2 (1%)	78	83
1	B	357/444 (80%)	355 (99%)	2 (1%)	78	83
All	All	714/888 (80%)	710 (99%)	4 (1%)	76	83

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	194	SER
1	A	368	LEU
1	B	194	SER
1	B	368	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	A	165	ASN
1	A	397	GLN
1	A	464	GLN
1	B	81	HIS
1	B	158	ASN
1	B	165	ASN
1	B	397	GLN
1	B	464	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

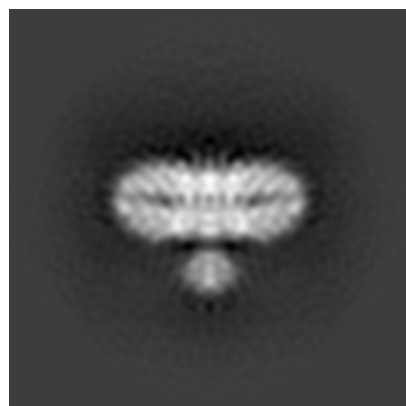
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12003. 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](#)

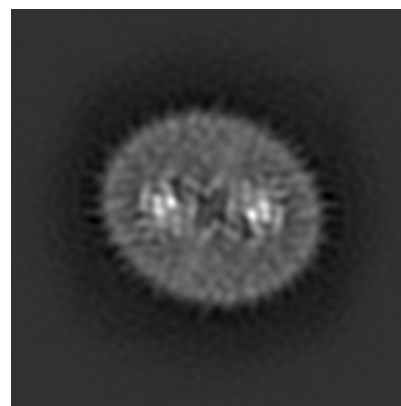
6.1.1 Primary map



X

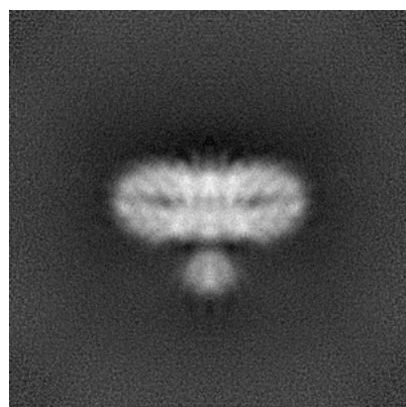


Y

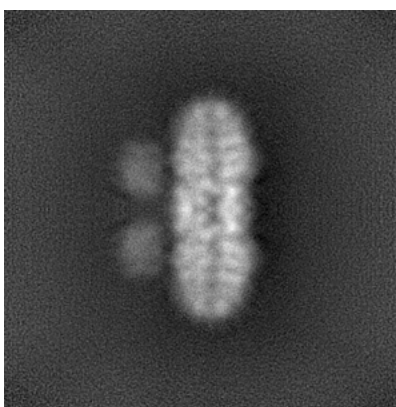


Z

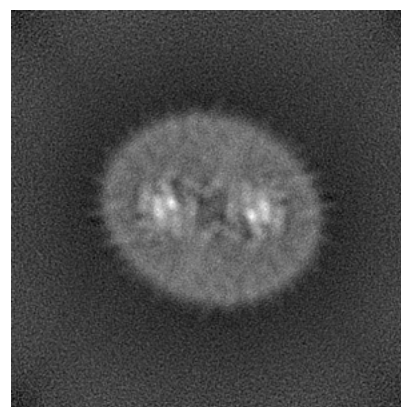
6.1.2 Raw map



X



Y

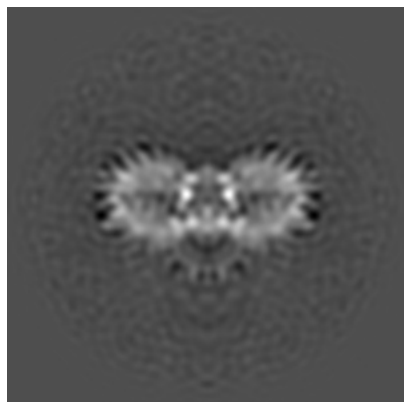


Z

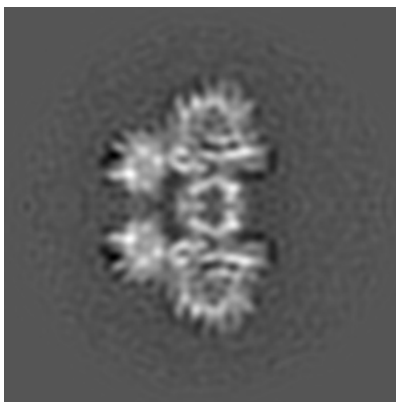
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

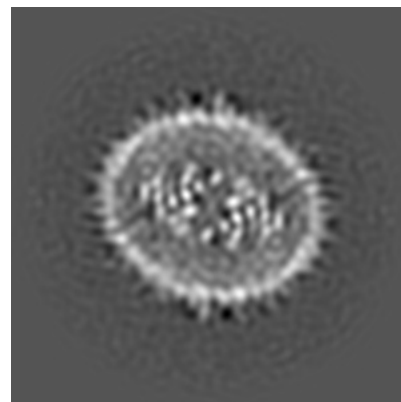
6.2.1 Primary map



X Index: 152

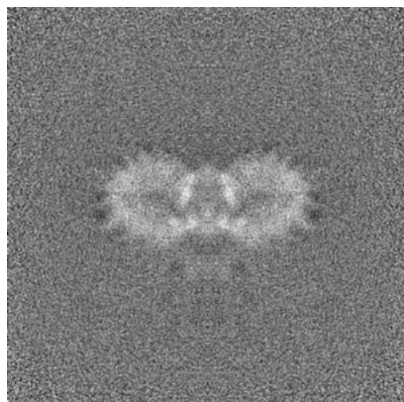


Y Index: 152

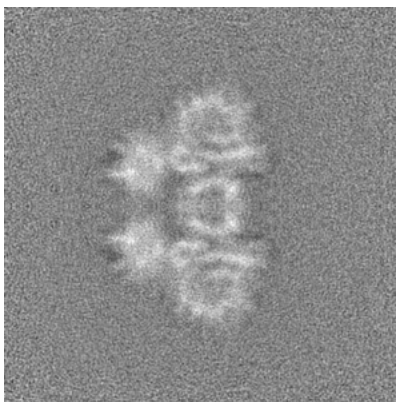


Z Index: 152

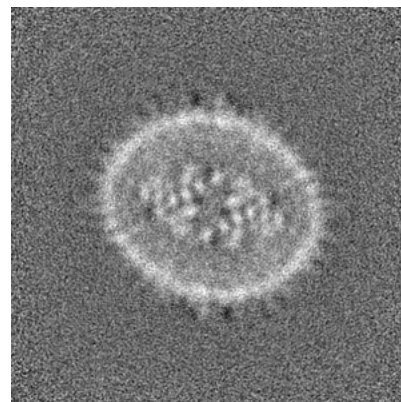
6.2.2 Raw map



X Index: 152



Y Index: 152

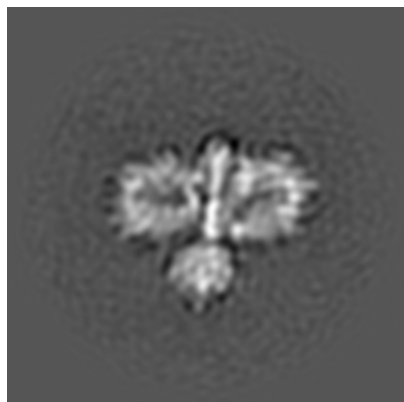


Z Index: 152

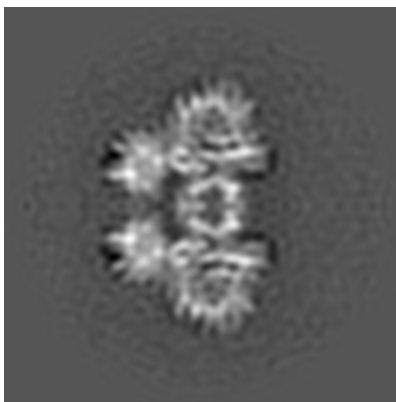
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

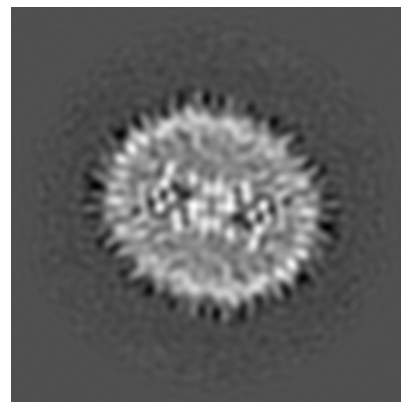
6.3.1 Primary map



X Index: 122

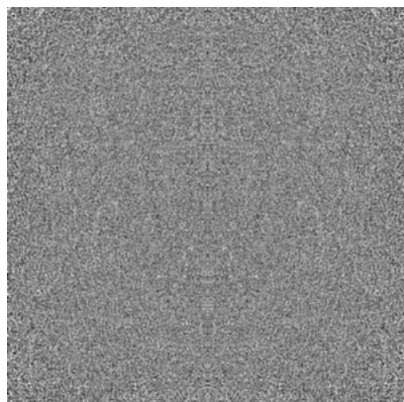


Y Index: 152

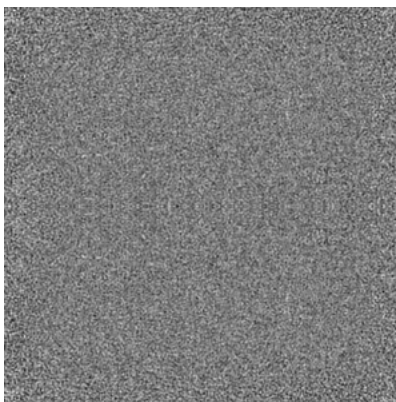


Z Index: 172

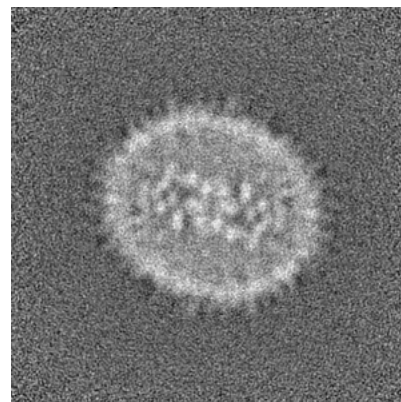
6.3.2 Raw map



X Index: 0



Y Index: 0



Z Index: 170

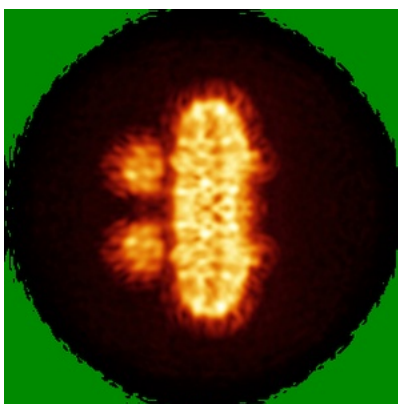
The images above show the largest variance slices of the map in three orthogonal directions.

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

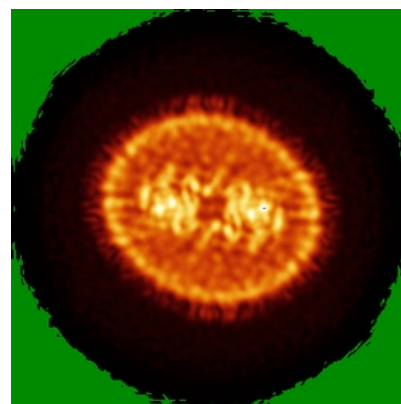
6.4.1 Primary map



X

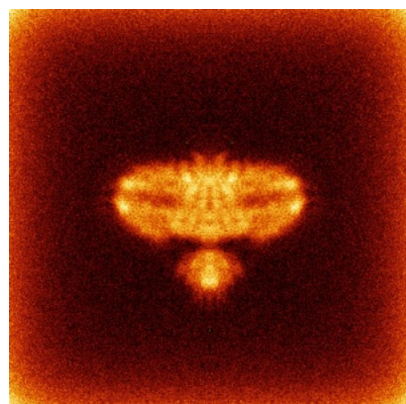


Y

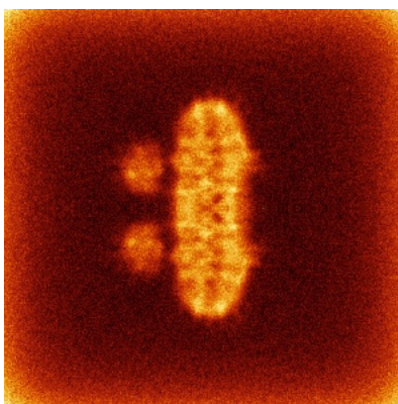


Z

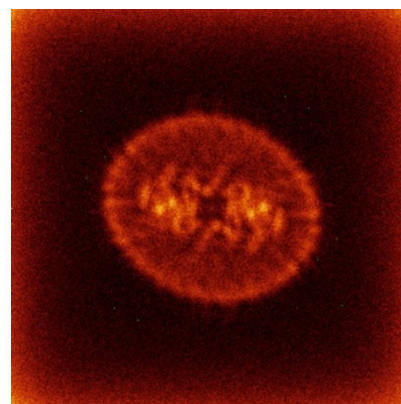
6.4.2 Raw map



X



Y

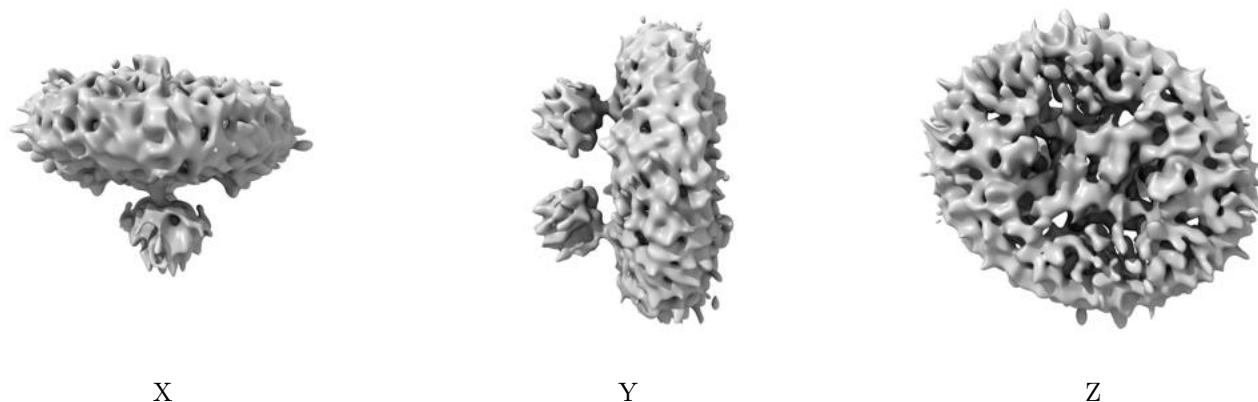


Z

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

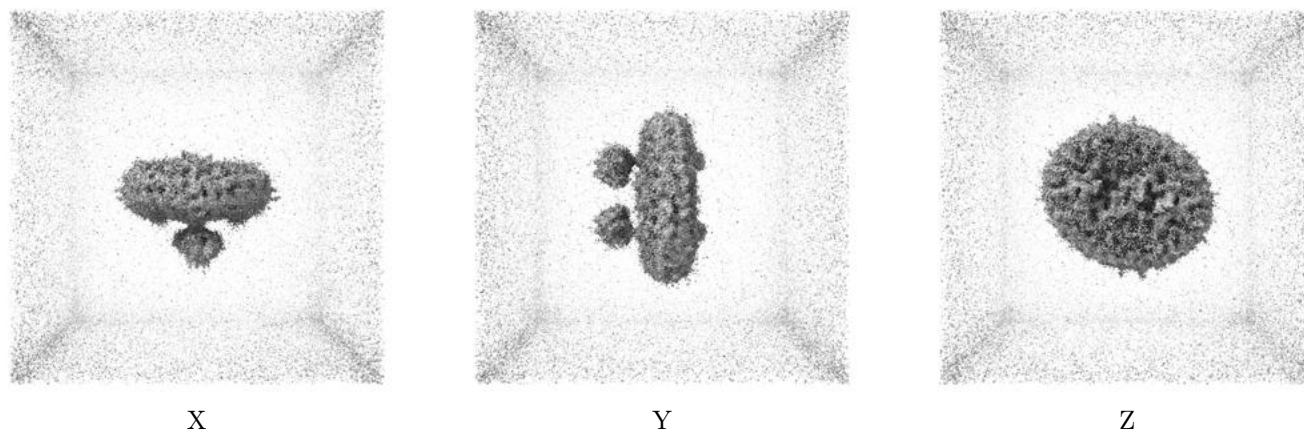
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.27. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

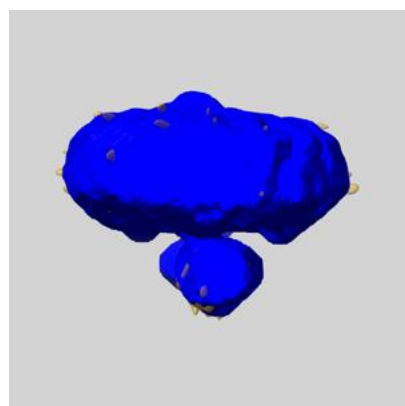
6.6 Mask visualisation [i](#)

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

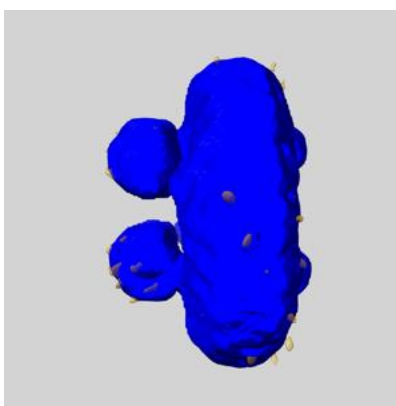
A mask typically either:

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

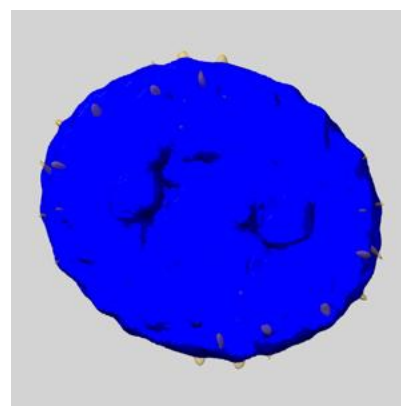
6.6.1 emd_12003_msk_1.map [i](#)



X



Y

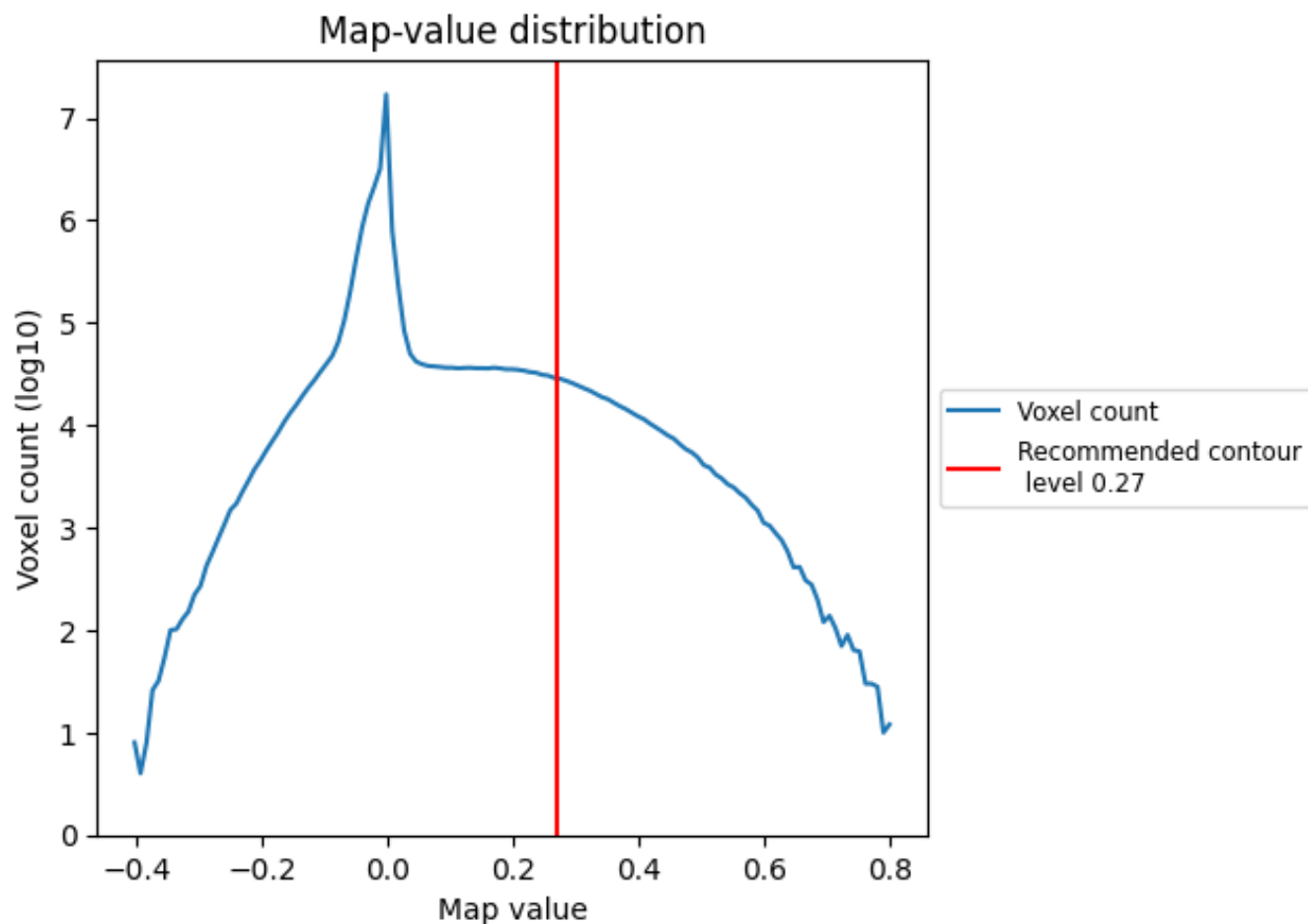


Z

7 Map analysis [i](#)

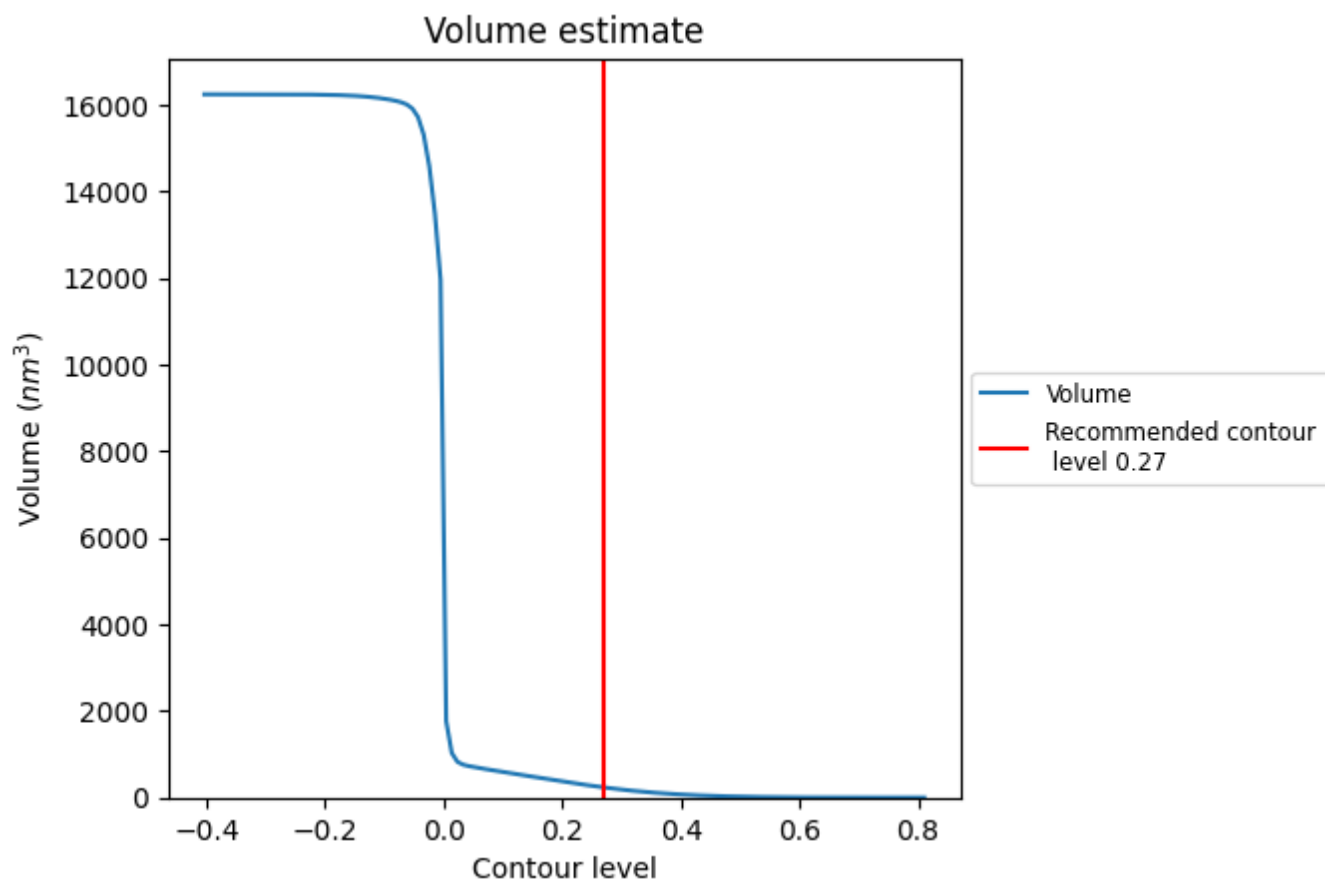
This section contains the results of statistical analysis of the map.

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.

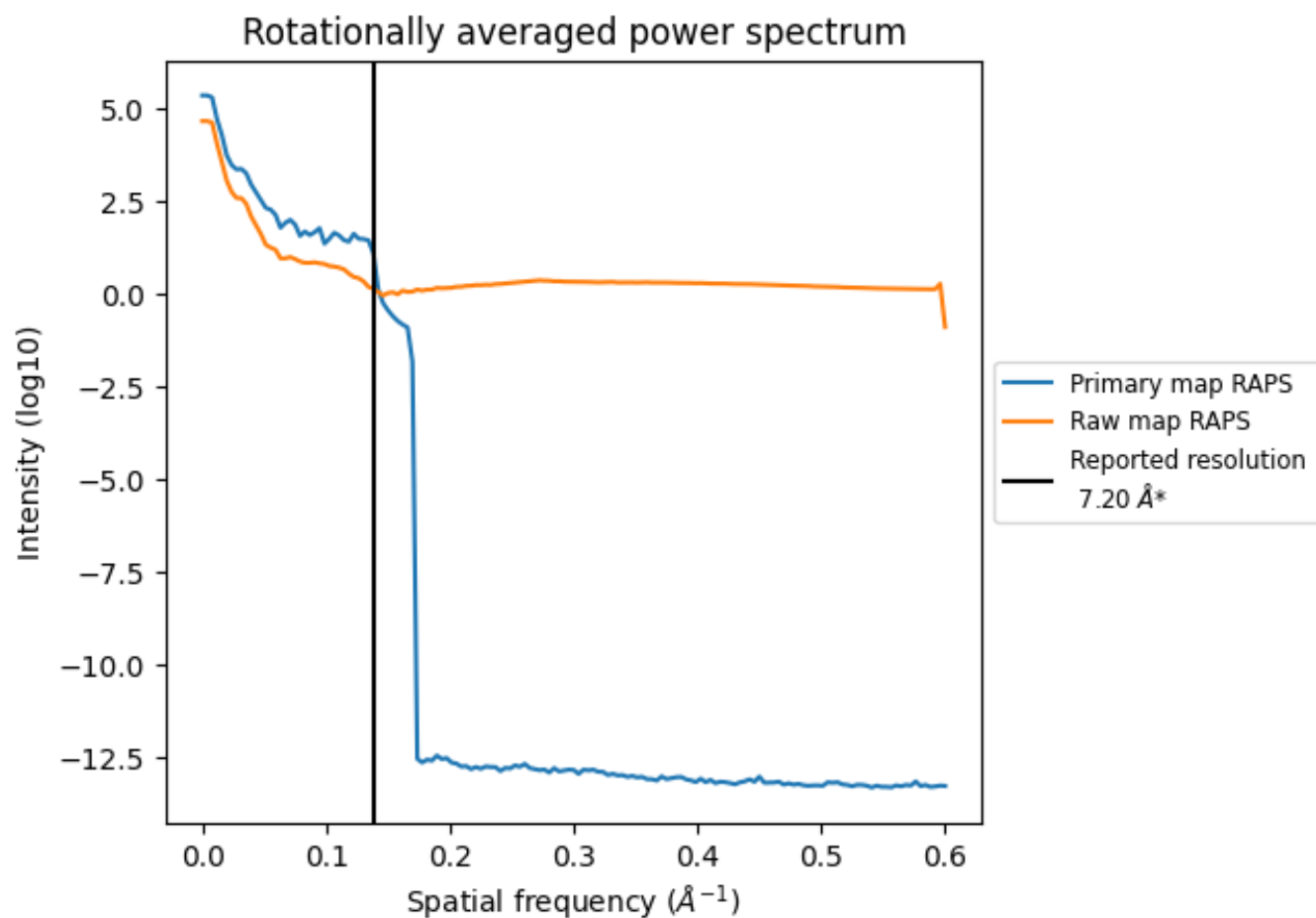
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 235 nm^3 ; this corresponds to an approximate mass of 212 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 ⓘ

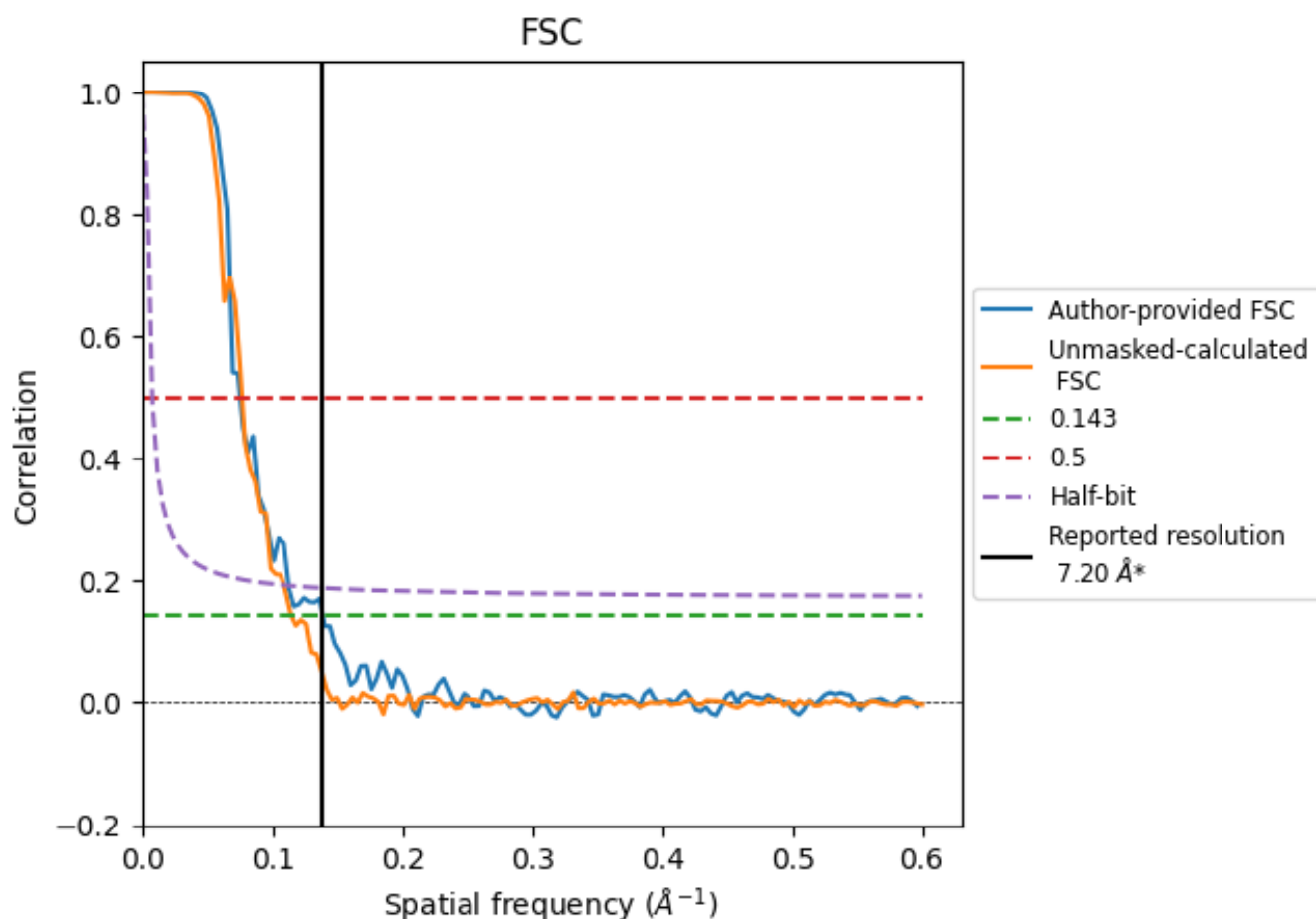


*Reported resolution corresponds to spatial frequency of 0.139 $\text{\AA}^{-1}$

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 0.139 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

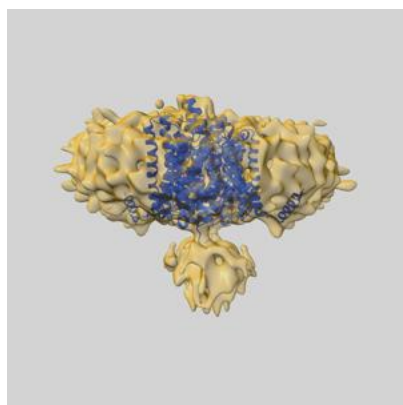
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	7.20	-	-
Author-provided FSC curve	7.18	13.32	8.87
Unmasked-calculated*	8.70	13.04	9.18

*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 8.70 differs from the reported value 7.2 by more than 10 %

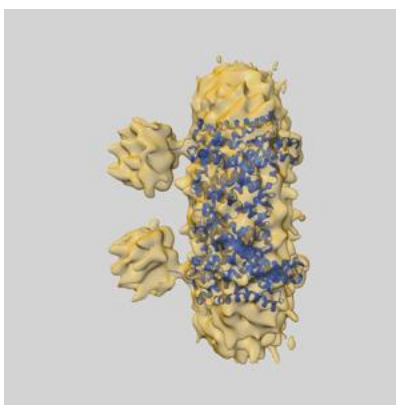
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-12003 and PDB model 7B4M. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

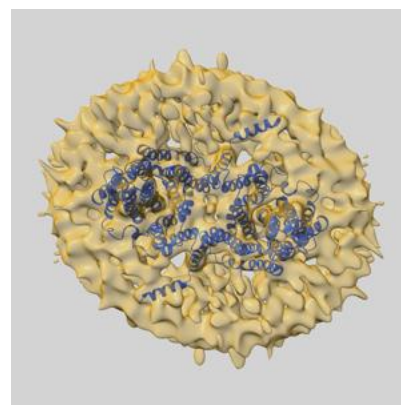
9.1 Map-model overlay [i](#)



X



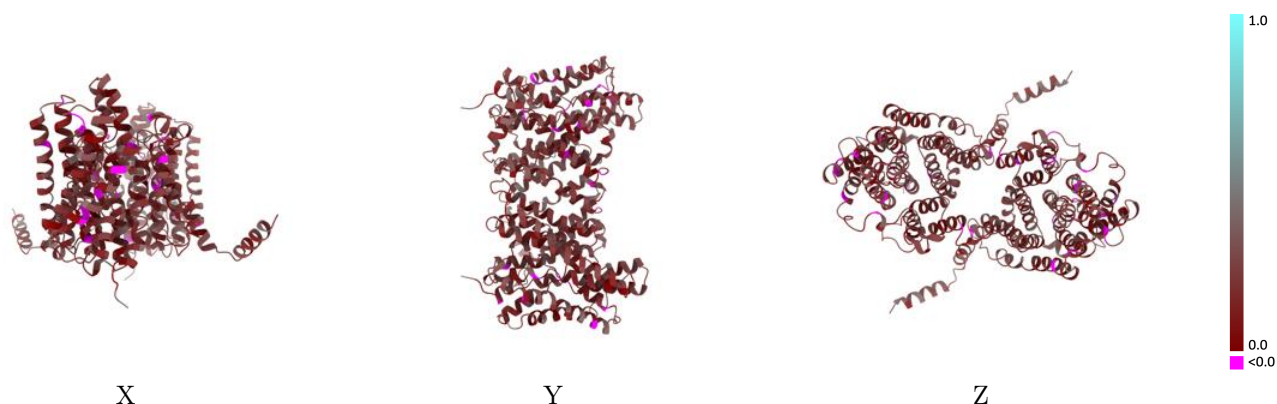
Y



Z

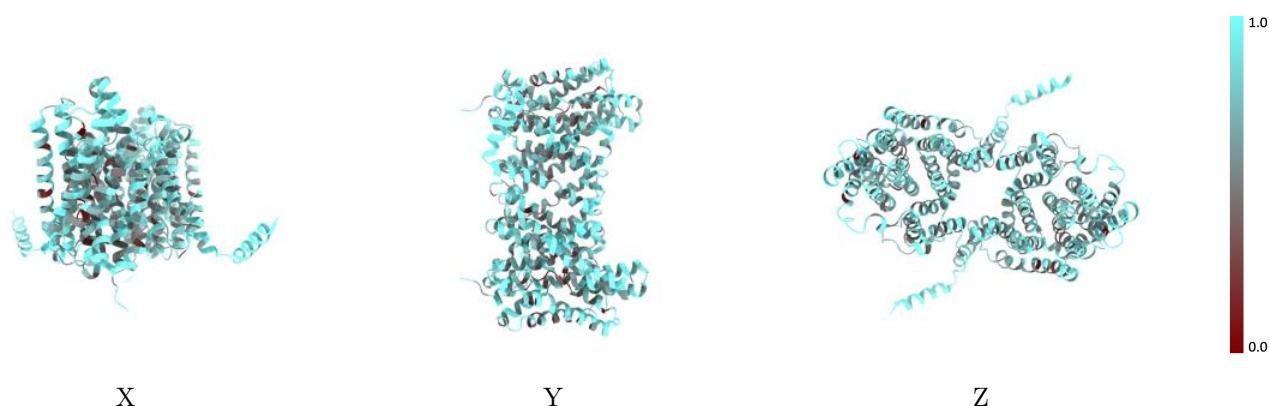
The images above show the 3D surface view of the map at the recommended contour level 0.27 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



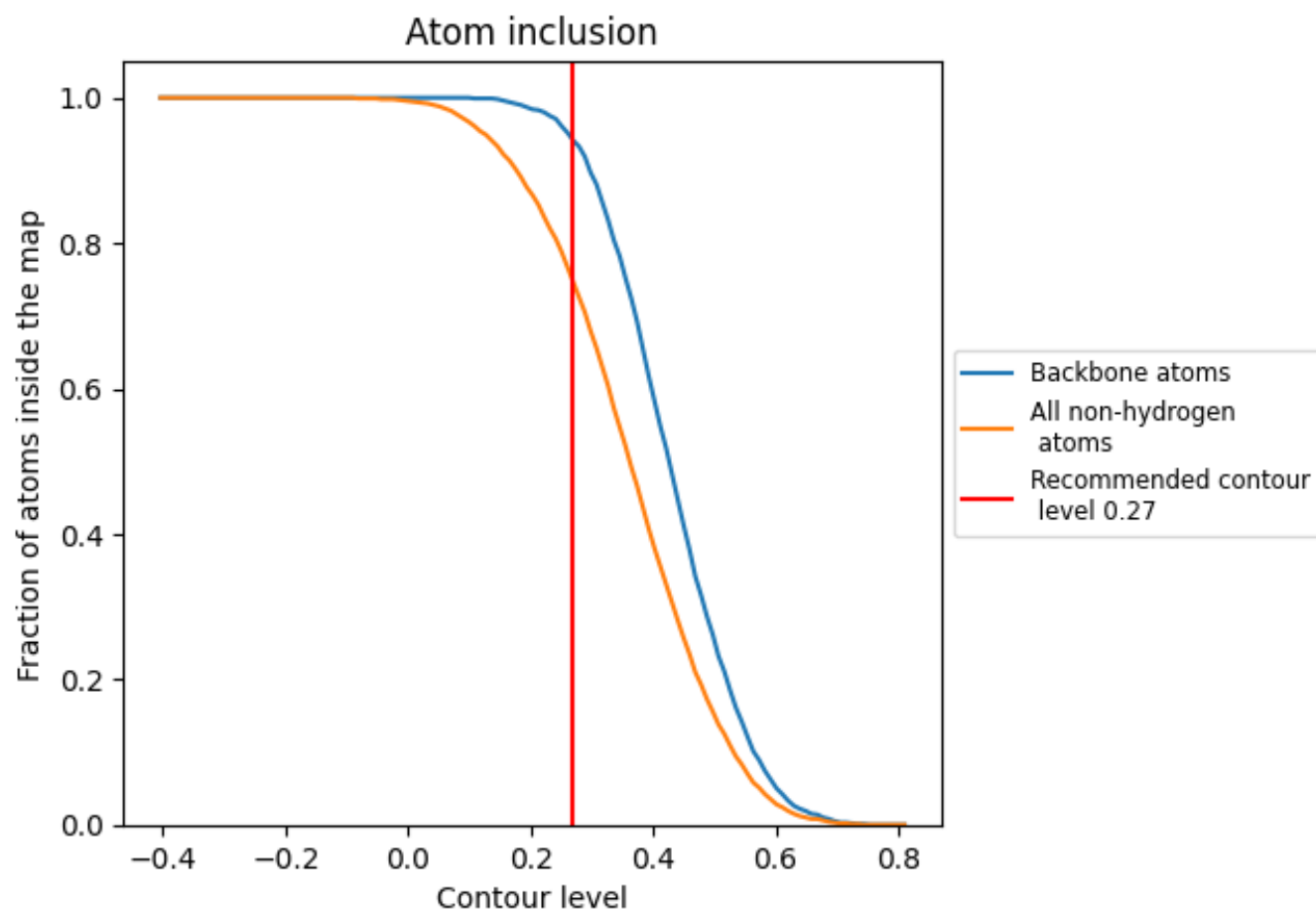
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.27).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 75% 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 (0.27) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7460	0.1860
A	0.7440	0.1870
B	0.7490	0.1860

