



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 11:40 AM UTC

PDB ID : 8I9Q / pdb_00008i9q
EMDB ID : EMD-35280
Title : The focused refinement of CCT4-PhLP2A from TRiC-PhLP2A complex in the open state
Authors : Roh, S.H.; Park, J.; Kim, H.; Lim, S.
Deposited on : 2023-02-07
Resolution : 4.22 Å(reported)
Based on initial model : 6NR8

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

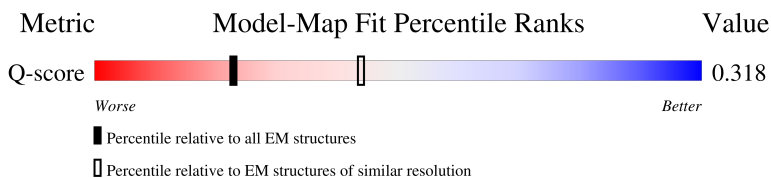
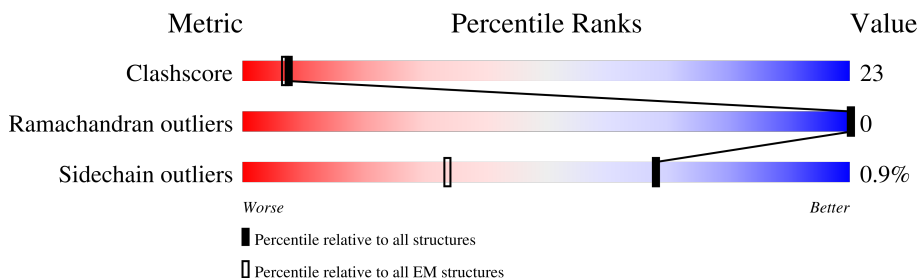
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.22 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	4772 (3.72 - 4.72)

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	D	539	
2	Q	239	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5179 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 T-complex protein 1 subunit delta.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	D	508	Total	C	N	O	S	0	0
			3832	2398	665	746	23		

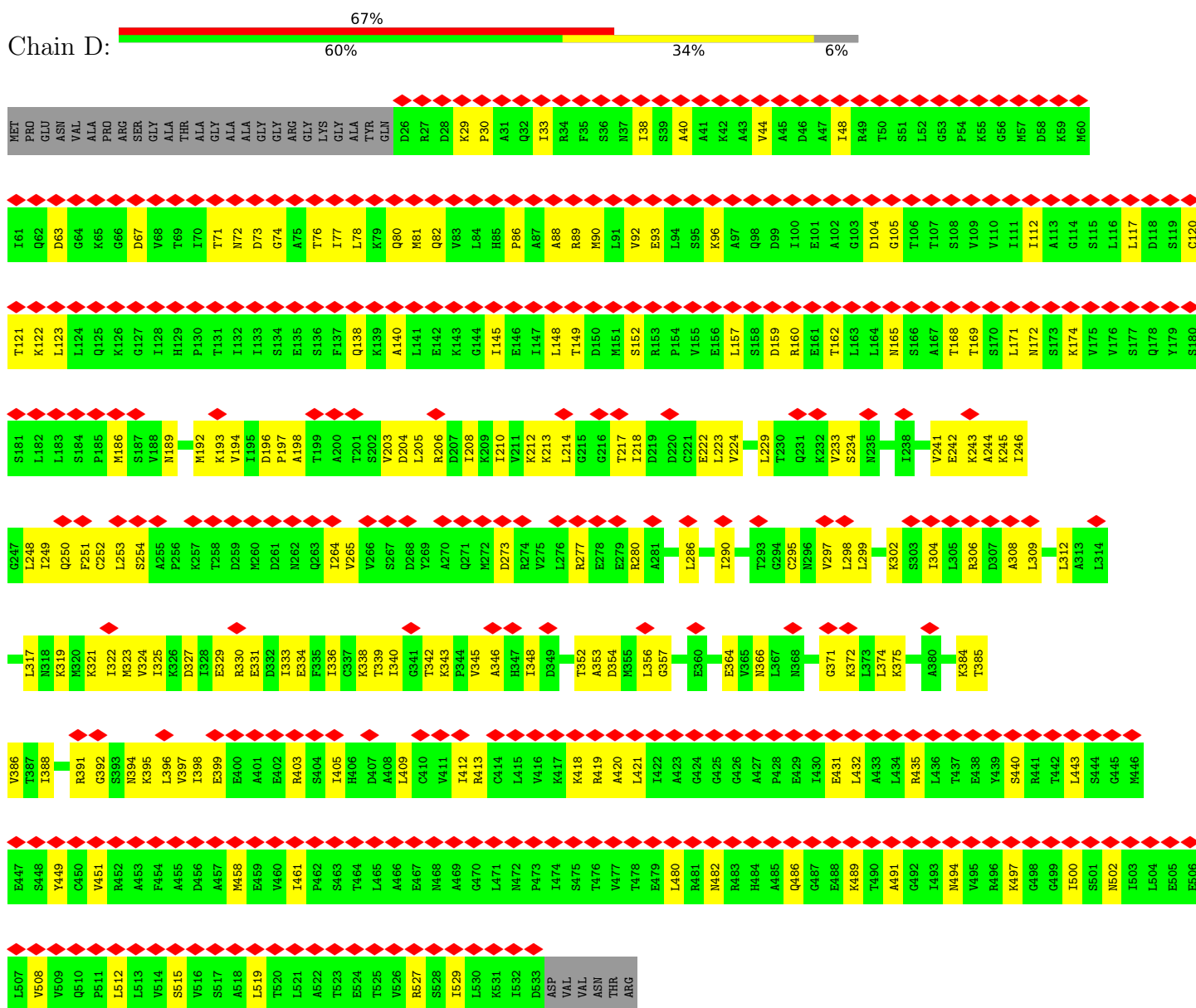
- Molecule 2 is a protein called Phosducin-like protein 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	Q	167	Total	C	N	O	S	0	0
			1347	866	232	243	6		

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: T-complex protein 1 subunit delta



- Molecule 2: Phosducin-like protein 3



MET
GLN
ASP
PRO
ASN
ALA
ASP
THR
GLU
TRP
ASN
ASP
TLE
LEU
ARG
LYS
LYS
GLY
TLE
LEU
PRO
LYS
SER
LEU
LYS
GLU
LEU
GLU
GLU
GLU
ALA
GLU
GLU
GLU
GLN
ARG
TLE
LEU
GLN
GLN
SER
VAL
VAL
LYS
THR
TYR
GLU
ASP
MET
THR
LEU
GLU
LEU
GLU
ASP
HIS
GLU

ASP
GLU
PHE
ASN
GLU
E66
D67
E68
R69
E72
H73
Y74
R75
R76
R77
R78
L79
A80
E81
W82
R83
A84
T85
R86
L87
R88
R89
R90
F91
G92
E93
Y94
L95
G99
K100
D101
Y102
Y103
Q104
E105
V106
T107
K108
A109
G110
E111
G112
L113
W114
V115
T116
L117
H118
L119
Y120
K121
Q122
G123

I130
G136
L137
A138
R139
D143
F146
I147
K148
A149
I150
S151
T152
T153
C154
I155
P156
M157
Y158
P159
D160
R161
M162
L163
P164
T165
I166
F167
V168
Y169
L170
E171
G172
D173
I174
K175
A176
G177
F178
P181
T190
R191
D192
E195
L198
S199
E200
T204
M205
E210

R211
P212
K213
K214
P215
I216
E217
D218
V219
L220
L221
S222
S223
V224
R225
R226
S227
V228
L229
M230
R231
R232
ASP
SER
ASP
SER
GLU
GLY
ASP

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	485964	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE; CTF correction was performed for every micrographs	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	6.843	Depositor
Minimum map value	-4.246	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.059	Depositor
Recommended contour level	0.616	Depositor
Map size (Å)	361.6, 361.6, 361.6	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.13, 1.13, 1.13	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	D	0.23	0/3863	0.47	0/5214
2	Q	0.69	2/1373 (0.1%)	1.10	9/1850 (0.5%)
All	All	0.40	2/5236 (0.0%)	0.69	9/7064 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Q	155	ILE	C-N	11.94	1.50	1.34
2	Q	158	TYR	C-N	11.38	1.50	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	158	TYR	CA-C-N	17.78	140.09	119.47
2	Q	158	TYR	C-N-CA	17.78	140.09	119.47
2	Q	155	ILE	CA-C-N	16.36	137.20	119.05
2	Q	155	ILE	C-N-CA	16.36	137.20	119.05
2	Q	159	PRO	CB-CA-C	-6.13	102.02	112.26
2	Q	159	PRO	N-CD-CG	-5.61	94.78	103.20
2	Q	156	PRO	N-CA-C	5.43	121.08	113.53
2	Q	160	ASP	CA-C-N	-5.31	111.84	121.14
2	Q	160	ASP	C-N-CA	-5.31	111.84	121.14

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	D	3832	0	4042	151	0
2	Q	1347	0	1392	93	0
All	All	5179	0	5434	243	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 23.

All (243) 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)
2:Q:94:VAL:HG11	2:Q:148:LYS:NZ	1.29	1.48
2:Q:121:LYS:CE	2:Q:159:PRO:O	1.70	1.39
2:Q:121:LYS:HE3	2:Q:159:PRO:O	1.19	1.26
2:Q:94:VAL:CG1	2:Q:148:LYS:NZ	1.95	1.26
2:Q:158:TYR:CE2	2:Q:162:ASN:OD1	1.97	1.18
2:Q:94:VAL:CG1	2:Q:148:LYS:HZ2	1.59	1.06
2:Q:121:LYS:HE2	2:Q:159:PRO:O	1.60	0.98
2:Q:119:LEU:HD21	2:Q:151:SER:HB3	1.46	0.98
2:Q:117:LEU:O	2:Q:166:ILE:HG23	1.66	0.95
2:Q:95:LEU:HD23	2:Q:95:LEU:H	1.30	0.95
1:D:245:LYS:HD3	1:D:342:THR:HG21	1.55	0.87
2:Q:161:ARG:CB	2:Q:161:ARG:HH21	1.88	0.86
2:Q:94:VAL:HG11	2:Q:148:LYS:HZ2	0.79	0.86
2:Q:117:LEU:C	2:Q:166:ILE:HG23	2.04	0.83
2:Q:94:VAL:CG1	2:Q:148:LYS:HZ3	1.88	0.82
1:D:213:LYS:HE3	1:D:391:ARG:HB3	1.60	0.82
2:Q:94:VAL:HG11	2:Q:148:LYS:HZ1	1.45	0.80
2:Q:157:ASN:O	2:Q:157:ASN:ND2	2.16	0.79
1:D:160:ARG:HH21	1:D:192:MET:HE1	1.47	0.78
1:D:224:VAL:HG11	1:D:375:LYS:HD2	1.64	0.78
2:Q:158:TYR:N	2:Q:159:PRO:HD2	1.98	0.77
2:Q:161:ARG:HH21	2:Q:161:ARG:HB3	1.50	0.77
2:Q:94:VAL:HG12	2:Q:148:LYS:HZ3	1.49	0.76
2:Q:94:VAL:HG12	2:Q:148:LYS:NZ	1.98	0.75
1:D:343:LYS:HB3	1:D:354:ASP:HB3	1.69	0.74
2:Q:120:TYR:HB2	2:Q:149:ALA:O	1.88	0.74
2:Q:158:TYR:CZ	2:Q:162:ASN:OD1	2.40	0.72
1:D:246:ILE:HA	1:D:297:VAL:HG23	1.72	0.72
2:Q:119:LEU:CD2	2:Q:151:SER:HB3	2.18	0.72
1:D:399:GLU:O	1:D:403:ARG:NH2	2.24	0.70
2:Q:116:ILE:HG23	2:Q:166:ILE:HG21	1.73	0.70
1:D:206:ARG:HH22	1:D:331:GLU:HB3	1.56	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:158:TYR:H	2:Q:159:PRO:HD2	1.57	0.69
2:Q:167:PHE:HB3	2:Q:174:ILE:CG2	2.21	0.69
1:D:405:ILE:O	1:D:409:LEU:HB2	1.93	0.68
1:D:204:ASP:OD1	1:D:205:LEU:N	2.27	0.68
1:D:440:SER:HB3	1:D:451:VAL:HG23	1.76	0.68
2:Q:93:GLU:HG3	2:Q:139:ARG:NH2	2.08	0.68
2:Q:95:LEU:HD23	2:Q:95:LEU:N	2.08	0.68
1:D:364:GLU:HB3	1:D:374:LEU:HA	1.75	0.68
1:D:312:LEU:HD22	2:Q:74:TYR:HE1	1.59	0.68
1:D:245:LYS:HE3	1:D:340:ILE:HG21	1.76	0.67
1:D:277:ARG:HG3	1:D:280:ARG:HE	1.59	0.67
2:Q:93:GLU:H	2:Q:139:ARG:HH21	1.41	0.67
1:D:394:ASN:HB3	1:D:397:VAL:HG23	1.77	0.66
1:D:298:LEU:H	1:D:324:VAL:HG12	1.60	0.65
1:D:157:LEU:HD13	1:D:419:ARG:HB3	1.78	0.65
1:D:290:ILE:HD11	1:D:348:ILE:HG22	1.80	0.64
2:Q:158:TYR:HB3	2:Q:159:PRO:CD	2.27	0.64
2:Q:116:ILE:CG2	2:Q:166:ILE:HG21	2.28	0.64
1:D:117:LEU:O	1:D:121:THR:HG23	1.98	0.64
2:Q:212:PRO:HG2	2:Q:214:LYS:H	1.62	0.64
1:D:210:ILE:HG23	1:D:388:ILE:HB	1.81	0.63
1:D:334:GLU:OE2	1:D:338:LYS:NZ	2.31	0.63
2:Q:79:LEU:HA	2:Q:82:TRP:HB3	1.78	0.63
2:Q:99:GLY:HA3	2:Q:153:THR:HG21	1.81	0.63
2:Q:151:SER:HB2	2:Q:159:PRO:HB3	1.80	0.63
2:Q:167:PHE:HB3	2:Q:174:ILE:HG23	1.80	0.62
1:D:331:GLU:OE1	1:D:331:GLU:N	2.24	0.62
2:Q:94:VAL:HG22	2:Q:146:PHE:HB2	1.81	0.62
1:D:193:LYS:HD3	1:D:223:LEU:HB3	1.82	0.61
1:D:196:ASP:O	1:D:198:ALA:N	2.34	0.61
1:D:122:LYS:NZ	1:D:449:TYR:OH	2.34	0.60
1:D:71:THR:OG1	1:D:73:ASP:OD1	2.19	0.60
1:D:203:VAL:O	1:D:413:ARG:NH1	2.35	0.60
1:D:340:ILE:HG23	1:D:342:THR:HG23	1.84	0.60
2:Q:101:ASP:OD1	2:Q:104:GLN:NE2	2.34	0.60
2:Q:95:LEU:H	2:Q:95:LEU:CD2	2.10	0.59
2:Q:176:ALA:HB1	2:Q:178:PHE:CE1	2.38	0.59
1:D:193:LYS:CE	1:D:223:LEU:HB3	2.32	0.59
2:Q:121:LYS:HE2	2:Q:162:ASN:O	2.02	0.59
1:D:194:VAL:HG22	1:D:386:VAL:HG21	1.85	0.59
1:D:193:LYS:CD	1:D:223:LEU:HB3	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:431:GLU:OE2	1:D:435:ARG:NH1	2.36	0.59
1:D:193:LYS:HE2	1:D:223:LEU:HB3	1.84	0.58
1:D:93:GLU:HA	1:D:96:LYS:HG2	1.85	0.58
2:Q:73:MET:HA	2:Q:76:ARG:HB2	1.85	0.58
1:D:81:MET:HE3	1:D:82:GLN:H	1.68	0.58
2:Q:192:ASP:OD1	2:Q:192:ASP:N	2.36	0.58
2:Q:204:ILE:HG23	2:Q:205:MET:H	1.69	0.58
1:D:38:ILE:HD13	1:D:121:THR:HG22	1.86	0.58
2:Q:119:LEU:HD21	2:Q:151:SER:CB	2.29	0.58
2:Q:117:LEU:H	2:Q:166:ILE:CG2	2.16	0.58
2:Q:93:GLU:N	2:Q:139:ARG:HH21	2.02	0.57
1:D:395:LYS:O	1:D:398:ILE:HG22	2.05	0.56
1:D:112:ILE:HD11	1:D:461:ILE:HD11	1.88	0.56
1:D:330:ARG:O	1:D:333:ILE:HG22	2.05	0.56
1:D:197:PRO:HG3	1:D:384:LYS:HB2	1.89	0.55
1:D:159:ASP:OD1	1:D:162:THR:HG23	2.07	0.55
1:D:160:ARG:NH2	1:D:192:MET:HE1	2.20	0.55
1:D:418:LYS:HG3	1:D:420:ALA:H	1.70	0.55
1:D:432:LEU:HB3	1:D:458:MET:HE3	1.89	0.55
2:Q:158:TYR:HB3	2:Q:159:PRO:HD2	1.87	0.55
1:D:336:ILE:HA	1:D:339:THR:HG22	1.88	0.55
1:D:88:ALA:O	1:D:92:VAL:HG23	2.07	0.55
2:Q:116:ILE:HG23	2:Q:166:ILE:CG2	2.38	0.54
1:D:171:LEU:HA	1:D:174:LYS:HB2	1.89	0.54
2:Q:167:PHE:HB3	2:Q:174:ILE:HG21	1.89	0.54
1:D:295:CYS:O	1:D:297:VAL:HG22	2.07	0.54
2:Q:117:LEU:O	2:Q:166:ILE:CG2	2.47	0.54
1:D:233:VAL:HG23	1:D:234:SER:H	1.73	0.54
1:D:273:ASP:O	1:D:277:ARG:NH1	2.41	0.53
2:Q:74:TYR:O	2:Q:78:ARG:N	2.41	0.53
1:D:233:VAL:HG12	1:D:325:ILE:HG13	1.89	0.53
2:Q:114:TRP:CD1	2:Q:170:LEU:HD13	2.44	0.53
1:D:117:LEU:HA	1:D:120:CYS:SG	2.49	0.53
2:Q:190:THR:OG1	2:Q:192:ASP:OD1	2.15	0.53
1:D:194:VAL:HG11	1:D:208:ILE:HD11	1.91	0.53
1:D:251:PHE:CE2	1:D:253:LEU:HB3	2.43	0.53
1:D:208:ILE:HD13	1:D:409:LEU:HD11	1.90	0.52
1:D:515:SER:O	1:D:519:LEU:HG	2.09	0.52
1:D:193:LYS:HE2	1:D:223:LEU:HD23	1.90	0.52
2:Q:109:ALA:HB3	2:Q:171:GLU:HG3	1.92	0.52
2:Q:69:ARG:O	2:Q:73:MET:HG2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:248:LEU:HD21	1:D:333:ILE:HD11	1.92	0.52
2:Q:158:TYR:HE2	2:Q:162:ASN:OD1	1.84	0.52
2:Q:161:ARG:HG2	2:Q:161:ARG:O	2.09	0.52
1:D:120:CYS:HA	1:D:123:LEU:HB3	1.92	0.51
1:D:145:ILE:O	1:D:149:THR:HG23	2.11	0.51
1:D:214:LEU:HD23	1:D:214:LEU:H	1.73	0.51
2:Q:221:LEU:HD22	2:Q:225:ARG:HH22	1.75	0.51
1:D:264:ILE:HG22	1:D:265:VAL:HG23	1.91	0.51
1:D:329:GLU:CD	1:D:330:ARG:H	2.19	0.51
1:D:243:LYS:HD3	1:D:244:ALA:N	2.26	0.50
1:D:491:ALA:HA	1:D:502:ASN:HA	1.92	0.50
2:Q:95:LEU:N	2:Q:95:LEU:CD2	2.73	0.50
2:Q:158:TYR:CB	2:Q:159:PRO:HD2	2.41	0.50
1:D:243:LYS:HD3	1:D:244:ALA:H	1.76	0.49
1:D:152:SER:OG	1:D:421:LEU:O	2.30	0.49
2:Q:117:LEU:HD12	2:Q:167:PHE:HD1	1.78	0.49
2:Q:176:ALA:HB1	2:Q:178:PHE:CZ	2.48	0.49
2:Q:154:CYS:SG	2:Q:155:ILE:N	2.85	0.49
2:Q:166:ILE:HD12	2:Q:166:ILE:H	1.78	0.49
1:D:251:PHE:HZ	1:D:286:LEU:HD23	1.77	0.49
1:D:366:ASN:HD21	1:D:372:LYS:HB2	1.77	0.49
1:D:336:ILE:O	1:D:339:THR:HG22	2.13	0.49
1:D:241:VAL:HG12	1:D:242:GLU:H	1.78	0.48
1:D:443:LEU:HD21	1:D:451:VAL:HG21	1.95	0.48
1:D:193:LYS:HZ3	1:D:222:GLU:C	2.22	0.48
1:D:162:THR:HA	1:D:165:ASN:HD21	1.78	0.48
1:D:394:ASN:CG	1:D:396:LEU:HB2	2.38	0.48
1:D:169:THR:HA	1:D:172:ASN:HD21	1.77	0.48
1:D:233:VAL:HG11	1:D:323:MET:HE1	1.95	0.48
1:D:138:GLN:HB2	1:D:527:ARG:HH12	1.77	0.48
1:D:189:ASN:O	1:D:193:LYS:HB2	2.14	0.48
1:D:394:ASN:OD1	1:D:396:LEU:N	2.40	0.48
2:Q:192:ASP:HA	2:Q:195:GLU:HG2	1.96	0.47
1:D:213:LYS:HD2	1:D:371:GLY:HA3	1.95	0.47
1:D:317:LEU:HB3	1:D:322:ILE:HB	1.95	0.47
1:D:40:ALA:O	1:D:44:VAL:HG23	2.14	0.47
2:Q:67:ASP:OD1	2:Q:68:GLU:N	2.48	0.47
2:Q:115:VAL:HG22	2:Q:169:TYR:O	2.15	0.47
1:D:78:LEU:HD21	1:D:92:VAL:HA	1.97	0.47
1:D:217:THR:HG22	1:D:391:ARG:NH2	2.30	0.47
1:D:189:ASN:OD1	1:D:388:ILE:HD12	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:250:GLN:HG3	1:D:346:ALA:HA	1.97	0.47
2:Q:77:ARG:HD3	2:Q:77:ARG:HA	1.68	0.47
2:Q:137:LEU:HD21	2:Q:191:ARG:HD2	1.97	0.47
2:Q:116:ILE:HD13	2:Q:146:PHE:CE1	2.51	0.46
1:D:208:ILE:O	1:D:210:ILE:HG13	2.15	0.46
1:D:248:LEU:O	1:D:345:VAL:HG12	2.15	0.46
1:D:76:THR:O	1:D:80:GLN:HG2	2.15	0.46
2:Q:104:GLN:OE1	2:Q:108:LYS:HB2	2.14	0.46
1:D:319:LYS:O	1:D:321:LYS:NZ	2.40	0.46
1:D:245:LYS:N	1:D:357:GLY:O	2.45	0.46
2:Q:214:LYS:O	2:Q:215:PRO:C	2.58	0.46
1:D:251:PHE:CE1	1:D:348:ILE:HG23	2.51	0.46
1:D:71:THR:HG23	1:D:77:ILE:HD11	1.98	0.45
1:D:73:ASP:OD1	1:D:73:ASP:N	2.46	0.45
1:D:418:LYS:HD2	1:D:512:LEU:HB3	1.97	0.45
1:D:205:LEU:N	1:D:413:ARG:HH22	2.15	0.45
1:D:304:ILE:HD12	1:D:304:ILE:HA	1.85	0.45
2:Q:130:ILE:HD12	2:Q:130:ILE:HA	1.81	0.45
1:D:86:PRO:HA	1:D:89:ARG:HG3	1.99	0.45
2:Q:220:LEU:O	2:Q:223:SER:OG	2.28	0.45
1:D:48:ILE:HG21	1:D:74:GLY:HA2	1.99	0.45
1:D:218:ILE:HG12	1:D:397:VAL:CG1	2.47	0.44
1:D:229:LEU:HB2	1:D:374:LEU:HD11	1.99	0.44
1:D:168:THR:O	1:D:172:ASN:ND2	2.50	0.44
2:Q:105:GLU:HA	2:Q:108:LYS:HB3	1.99	0.44
1:D:241:VAL:HG12	1:D:242:GLU:N	2.33	0.44
1:D:245:LYS:CD	1:D:342:THR:HG21	2.38	0.44
2:Q:174:ILE:H	2:Q:174:ILE:HG13	1.52	0.44
1:D:193:LYS:HD3	1:D:223:LEU:CB	2.48	0.44
1:D:212:LYS:HB2	1:D:398:ILE:HD11	2.00	0.44
1:D:366:ASN:ND2	1:D:372:LYS:HB2	2.31	0.43
2:Q:114:TRP:CZ2	2:Q:170:LEU:HD22	2.52	0.43
1:D:306:ARG:NH2	1:D:309:LEU:HB2	2.33	0.43
1:D:352:THR:OG1	1:D:353:ALA:N	2.51	0.43
2:Q:155:ILE:HG23	2:Q:155:ILE:O	2.18	0.43
1:D:194:VAL:HG21	1:D:208:ILE:HD13	1.99	0.43
1:D:203:VAL:O	1:D:413:ARG:CZ	2.67	0.43
1:D:329:GLU:O	1:D:333:ILE:N	2.51	0.43
1:D:384:LYS:O	1:D:385:THR:OG1	2.32	0.43
1:D:508:VAL:HG23	1:D:508:VAL:O	2.19	0.43
1:D:251:PHE:CG	1:D:252:CYS:N	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:117:LEU:H	2:Q:166:ILE:HG22	1.80	0.43
1:D:117:LEU:HD21	1:D:529:ILE:HD11	2.01	0.43
1:D:418:LYS:O	1:D:419:ARG:HG3	2.19	0.43
1:D:297:VAL:HG12	1:D:323:MET:H	1.83	0.42
1:D:302:LYS:HA	1:D:327:ASP:OD1	2.18	0.42
1:D:317:LEU:HB2	1:D:322:ILE:HD12	2.00	0.42
2:Q:99:GLY:HA3	2:Q:153:THR:CG2	2.49	0.42
1:D:140:ALA:HB2	1:D:451:VAL:HG12	2.01	0.42
2:Q:100:LYS:HA	2:Q:100:LYS:HD3	1.92	0.42
2:Q:161:ARG:HH21	2:Q:161:ARG:CG	2.32	0.42
1:D:138:GLN:HB2	1:D:527:ARG:NH1	2.34	0.42
1:D:249:ILE:HD11	1:D:298:LEU:HD13	2.01	0.42
1:D:251:PHE:HE2	1:D:253:LEU:HB3	1.85	0.42
1:D:277:ARG:HA	1:D:280:ARG:NE	2.34	0.42
1:D:29:LYS:HB3	1:D:30:PRO:HD3	2.02	0.42
1:D:302:LYS:NZ	1:D:327:ASP:O	2.52	0.42
1:D:480:LEU:HD12	1:D:480:LEU:HA	1.87	0.42
1:D:489:LYS:HA	1:D:489:LYS:HD2	1.87	0.42
2:Q:163:LEU:HG	2:Q:181:PRO:HD3	2.01	0.42
2:Q:198:LEU:O	2:Q:200:GLU:N	2.53	0.42
1:D:491:ALA:HB3	1:D:500:ILE:HG22	2.02	0.42
2:Q:117:LEU:HA	2:Q:117:LEU:HD23	1.74	0.42
1:D:63:ASP:HB3	1:D:67:ASP:OD1	2.20	0.41
1:D:254:SER:OG	1:D:308:ALA:O	2.38	0.41
1:D:342:THR:HG22	1:D:356:LEU:O	2.20	0.41
1:D:148:LEU:HG	1:D:512:LEU:HD11	2.01	0.41
2:Q:93:GLU:CG	2:Q:139:ARG:NH2	2.80	0.41
2:Q:163:LEU:HD12	2:Q:181:PRO:HA	2.02	0.41
1:D:72:ASN:OD1	1:D:72:ASN:N	2.53	0.41
1:D:409:LEU:HA	1:D:412:ILE:HG12	2.03	0.41
1:D:169:THR:HA	1:D:172:ASN:ND2	2.36	0.41
1:D:196:ASP:HB2	1:D:197:PRO:HD3	2.03	0.41
1:D:297:VAL:HG12	1:D:323:MET:N	2.35	0.41
1:D:443:LEU:HD12	1:D:443:LEU:O	2.20	0.41
1:D:299:LEU:HD22	1:D:325:ILE:HB	2.02	0.41
1:D:494:ASN:HD22	1:D:497:LYS:HG2	1.86	0.41
1:D:30:PRO:HA	1:D:33:ILE:HG22	2.02	0.41
1:D:245:LYS:HD2	1:D:340:ILE:HG12	2.02	0.40
2:Q:99:GLY:CA	2:Q:153:THR:HG21	2.50	0.40
2:Q:162:ASN:HD22	2:Q:162:ASN:N	2.18	0.40
1:D:104:ASP:OD1	1:D:105:GLY:N	2.47	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:217:THR:C	1:D:392:GLY:HA2	2.46	0.40
1:D:482:ASN:O	1:D:486:GLN:HG2	2.21	0.40
2:Q:161:ARG:CB	2:Q:161:ARG:NH2	2.70	0.40
1:D:90:MET:O	1:D:93:GLU:HG3	2.22	0.40
1:D:186:MET:HE3	1:D:186:MET:HB3	1.97	0.40
2:Q:109:ALA:O	2:Q:113:LEU:HB3	2.22	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	D	506/539 (94%)	451 (89%)	55 (11%)	0	100	100
2	Q	165/239 (69%)	143 (87%)	22 (13%)	0	100	100
All	All	671/778 (86%)	594 (88%)	77 (12%)	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	D	433/452 (96%)	433 (100%)	0	100	100
2	Q	147/215 (68%)	142 (97%)	5 (3%)	32	54

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	580/667 (87%)	575 (99%)	5 (1%)	68 76

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	Q	160	ASP
2	Q	161	ARG
2	Q	163	LEU
2	Q	165	THR
2	Q	167	PHE

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

Mol	Chain	Res	Type
1	D	85	HIS
1	D	98	GLN
1	D	172	ASN
1	D	347	HIS
1	D	350	GLN
1	D	366	ASN
1	D	472	ASN
1	D	484	HIS
1	D	494	ASN
2	Q	131	ASN
2	Q	157	ASN
2	Q	162	ASN

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 ⓘ

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.

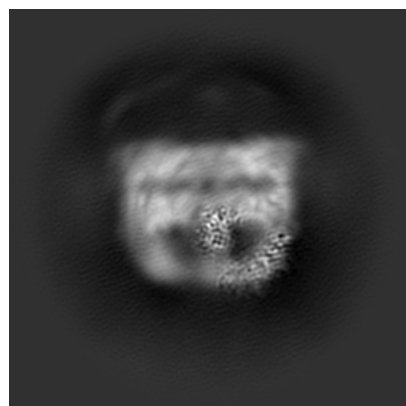
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-35280. These allow visual inspection of the internal detail of the map and identification of artifacts.

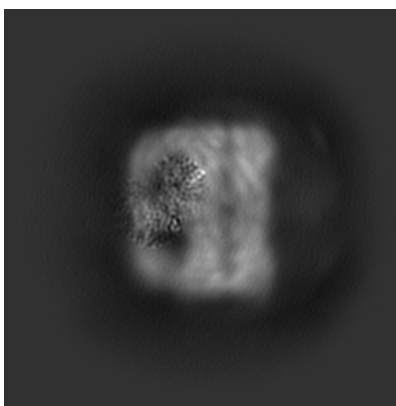
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

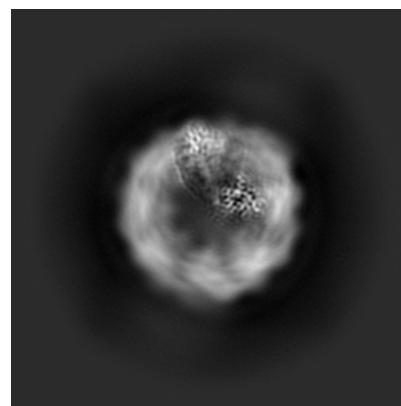
6.1.1 Primary map



X

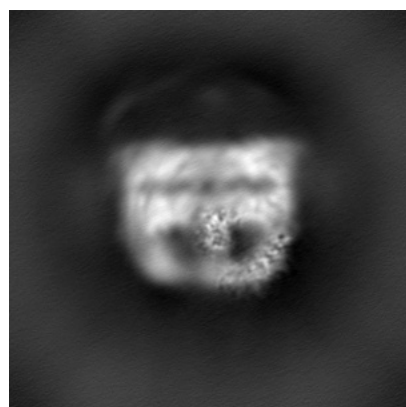


Y

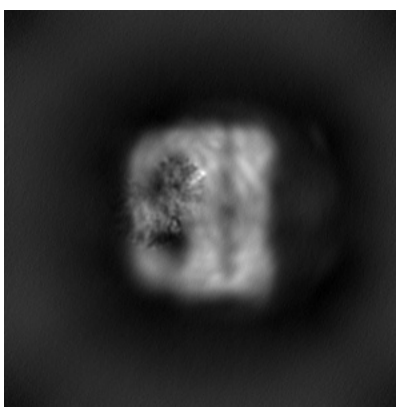


Z

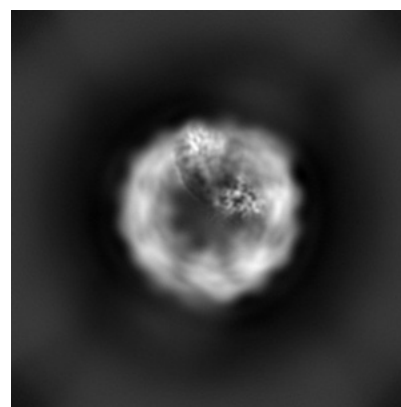
6.1.2 Raw map



X



Y

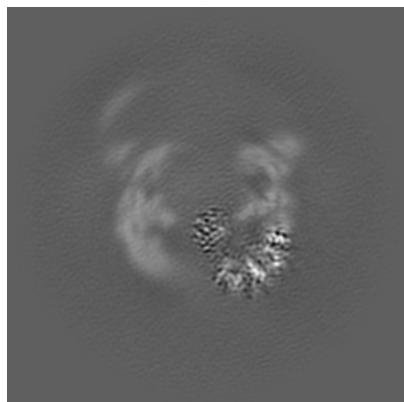


Z

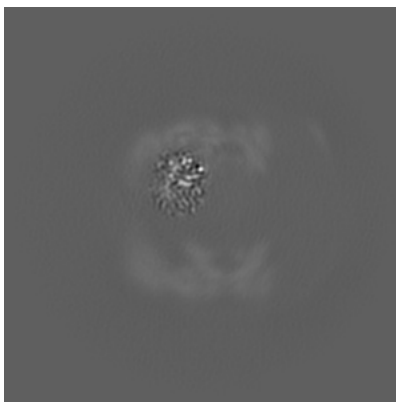
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

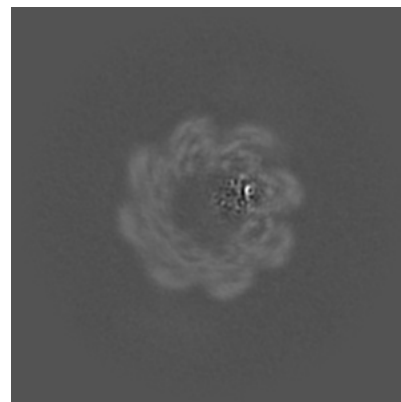
6.2.1 Primary map



X Index: 160

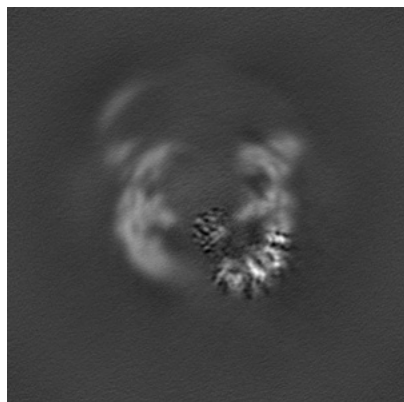


Y Index: 160

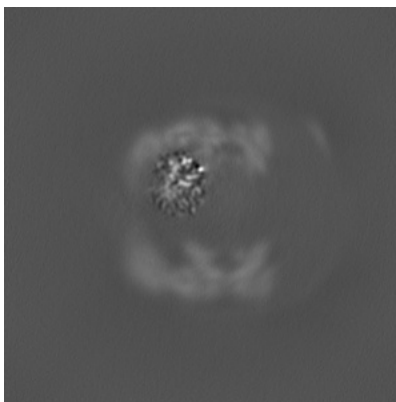


Z Index: 160

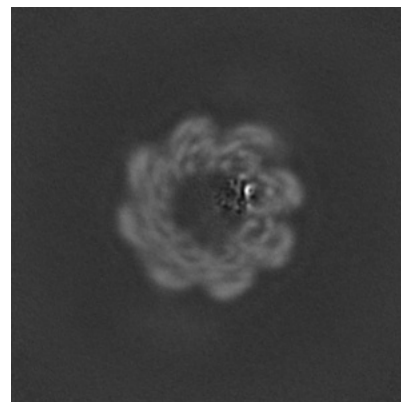
6.2.2 Raw map



X Index: 160



Y Index: 160

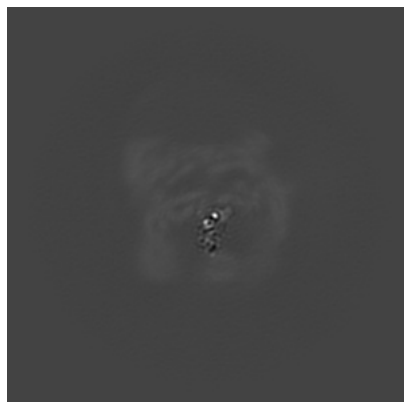


Z Index: 160

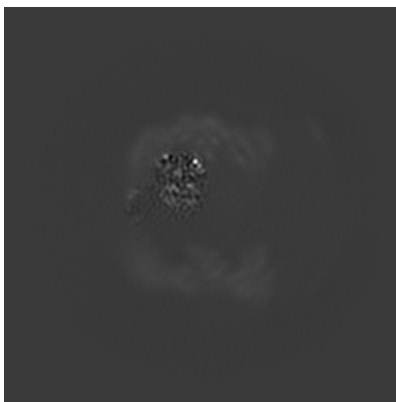
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

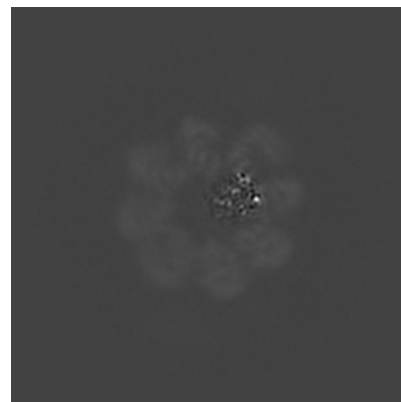
6.3.1 Primary map



X Index: 196

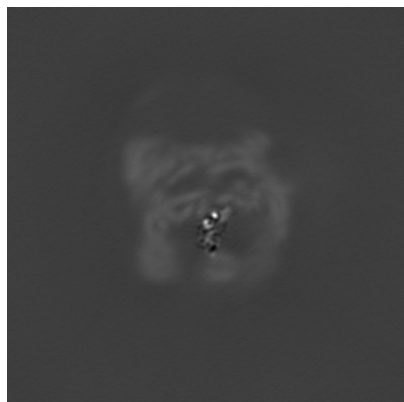


Y Index: 166

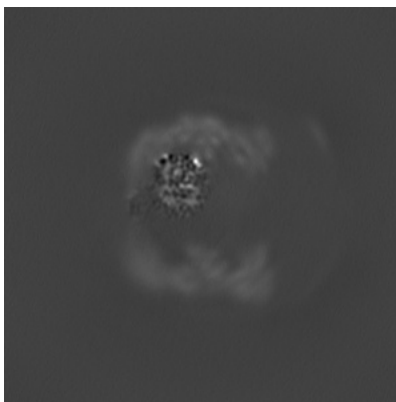


Z Index: 153

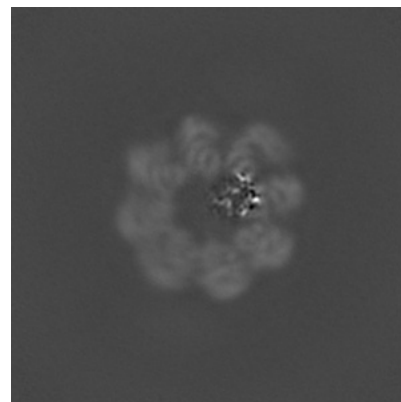
6.3.2 Raw map



X Index: 196



Y Index: 165

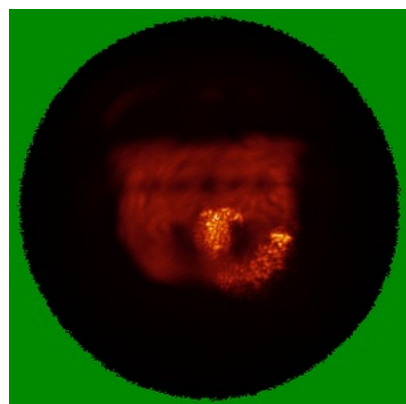


Z Index: 154

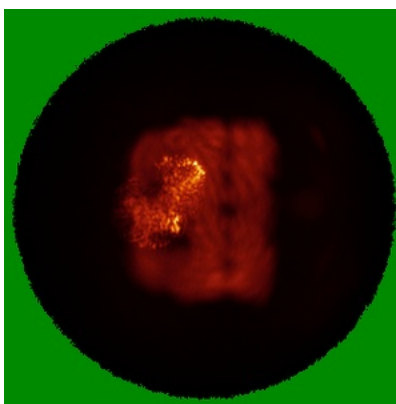
The images above show the largest variance slices of the map in three orthogonal directions.

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

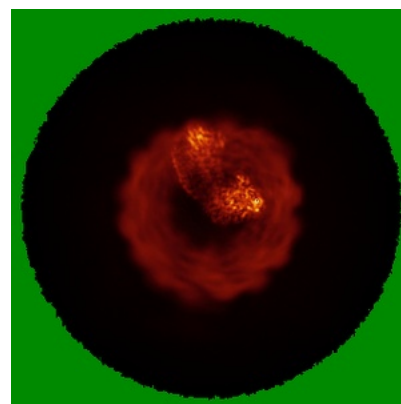
6.4.1 Primary map



X

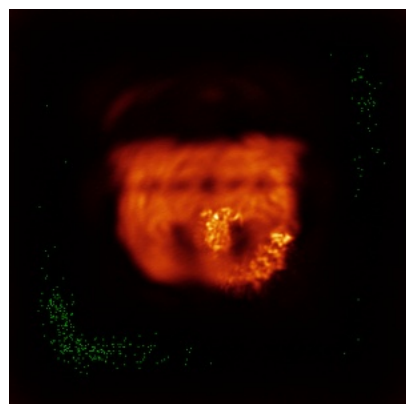


Y

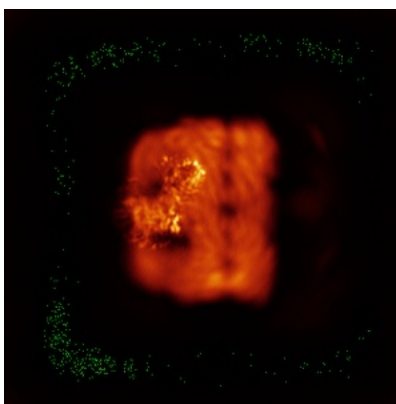


Z

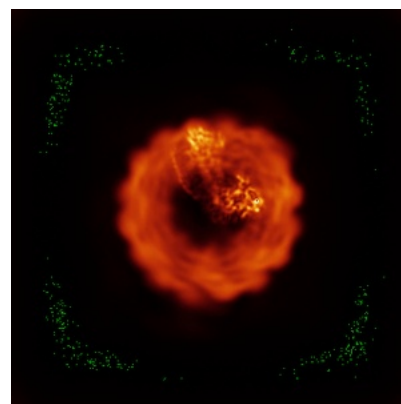
6.4.2 Raw map



X



Y

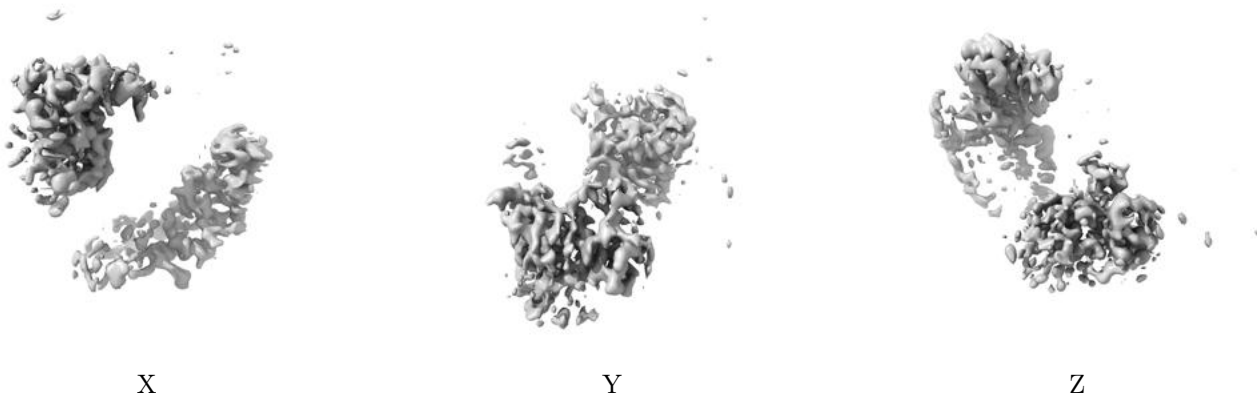


Z

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 0.616. 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.

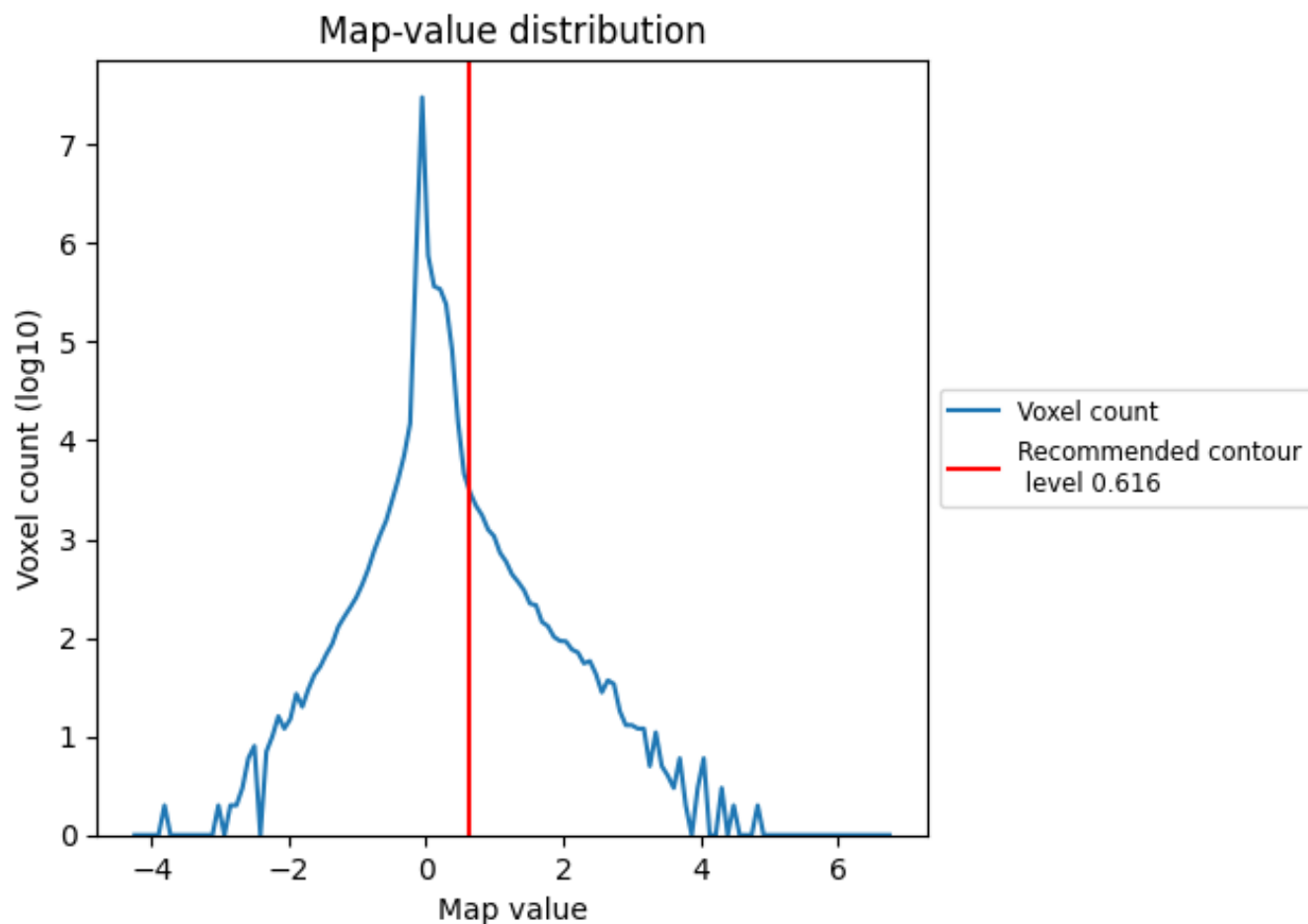
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

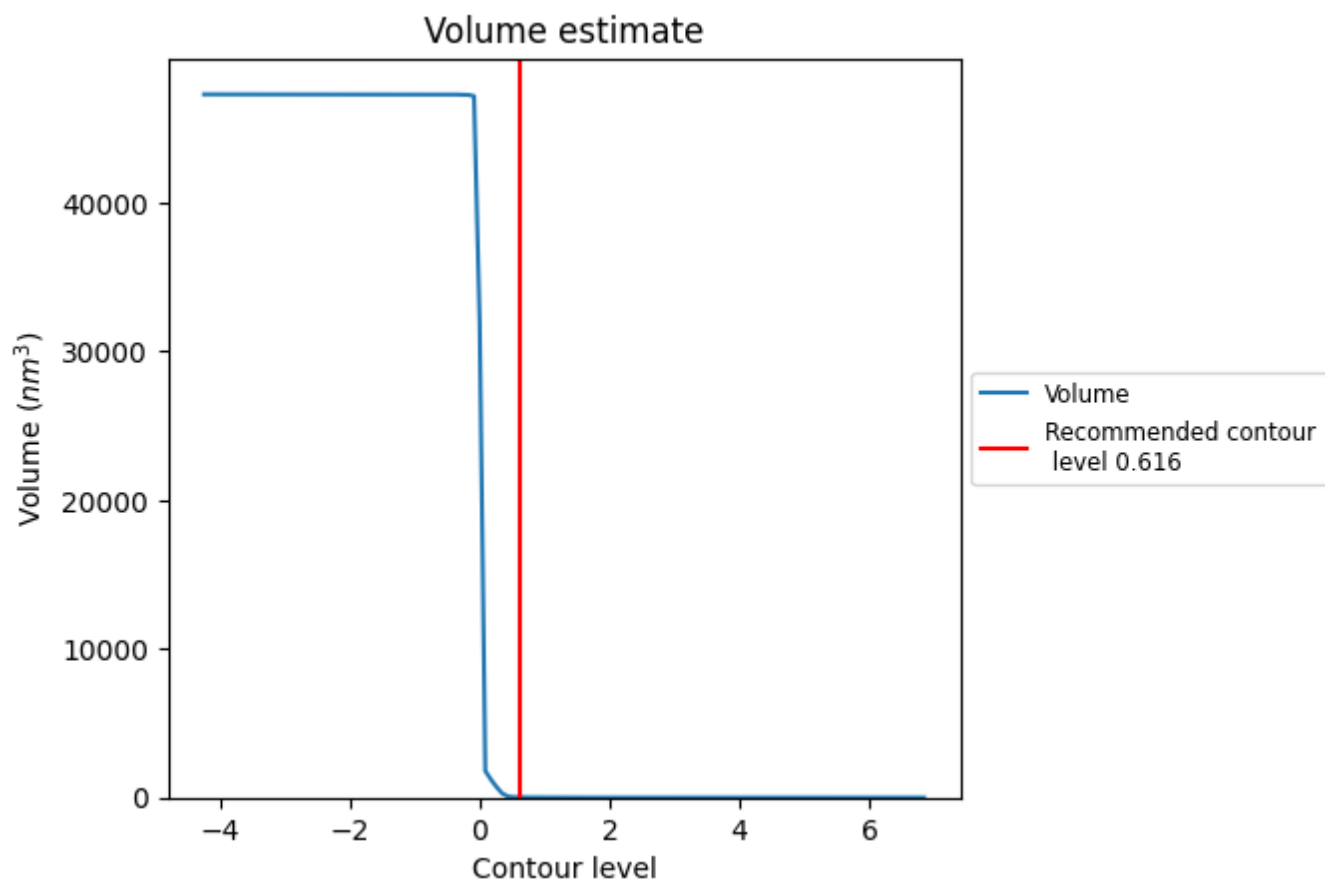
This section contains the results of statistical analysis of the map.

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.

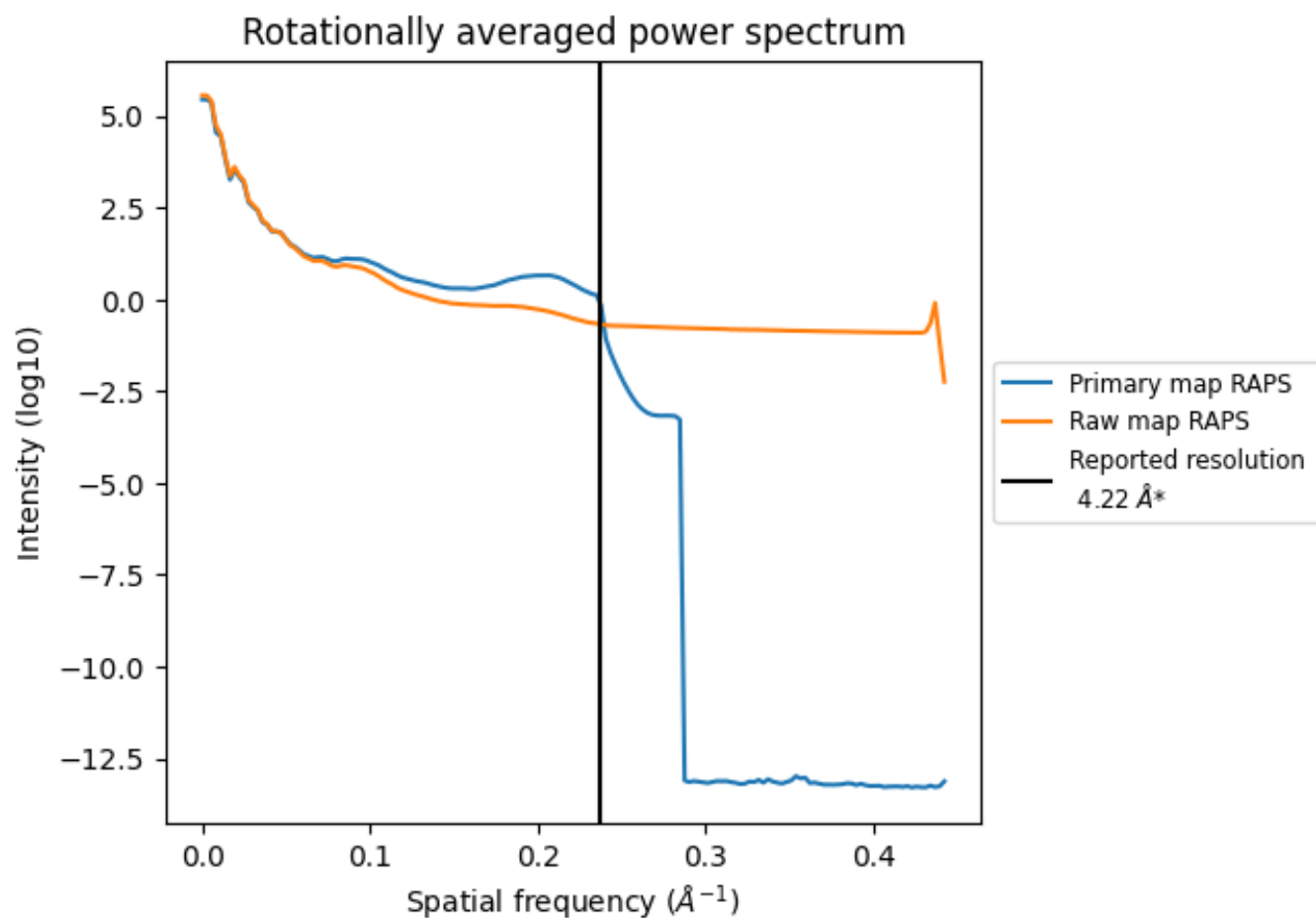
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 21 nm^3 ; this corresponds to an approximate mass of 19 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 ⓘ

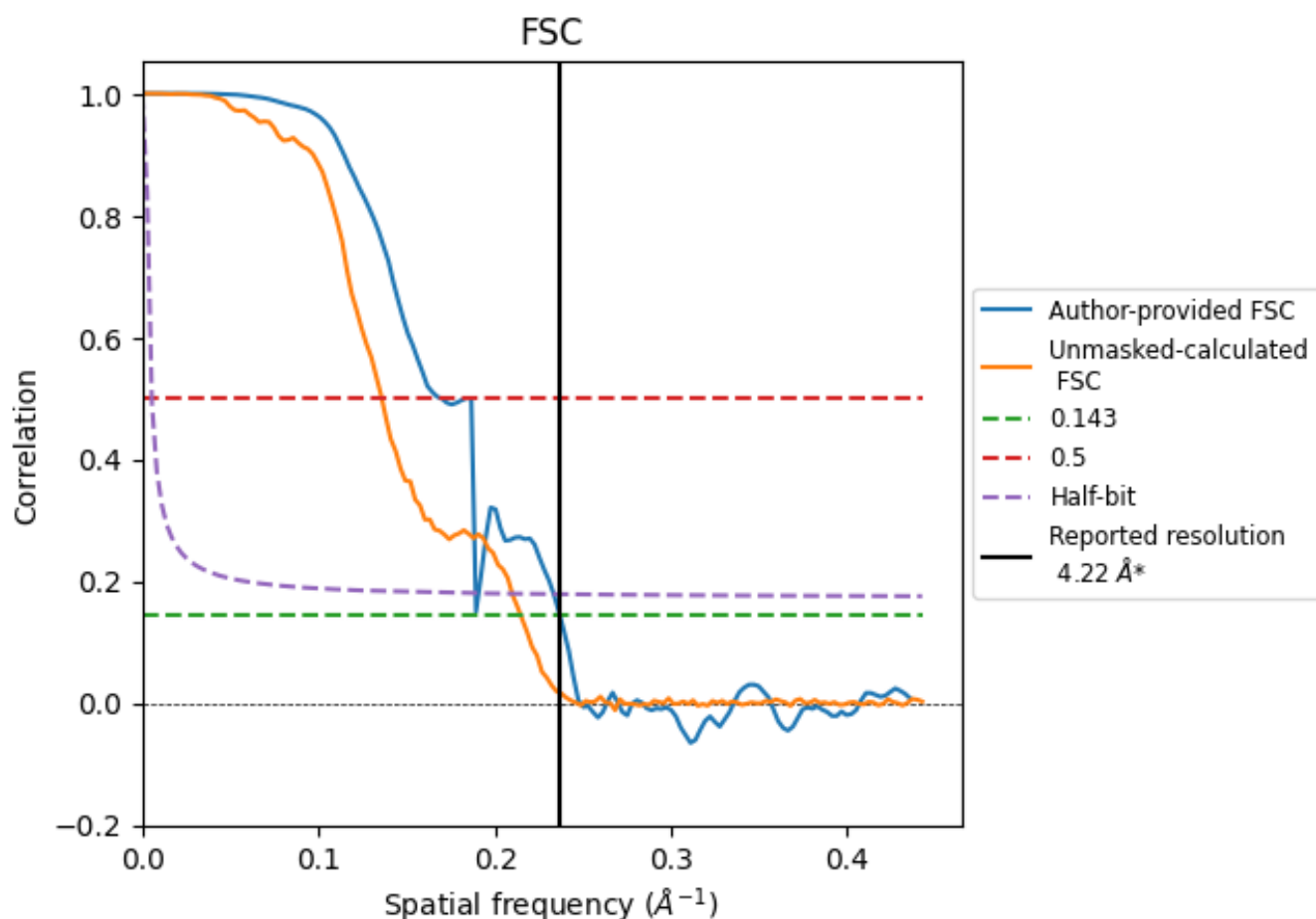


*Reported resolution corresponds to spatial frequency of 0.237 Å⁻¹

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 0.237 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

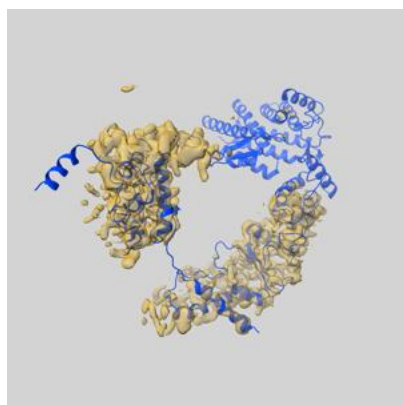
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.22	-	-
Author-provided FSC curve	4.22	5.90	5.29
Unmasked-calculated*	4.65	7.36	4.76

*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 4.65 differs from the reported value 4.22 by more than 10 %

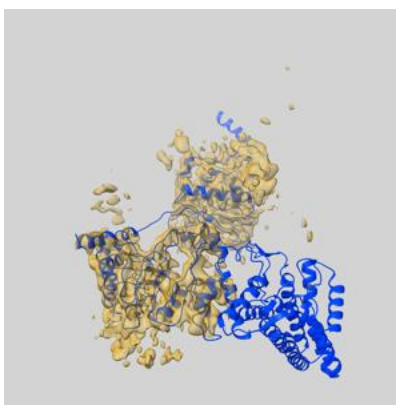
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-35280 and PDB model 8I9Q. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

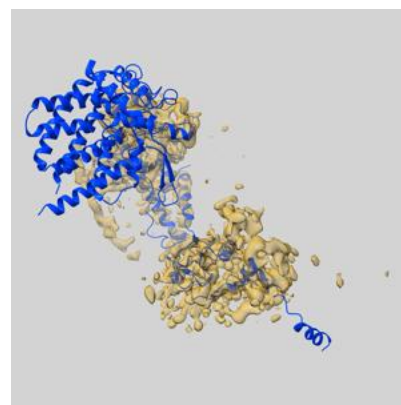
9.1 Map-model overlay [i](#)



X



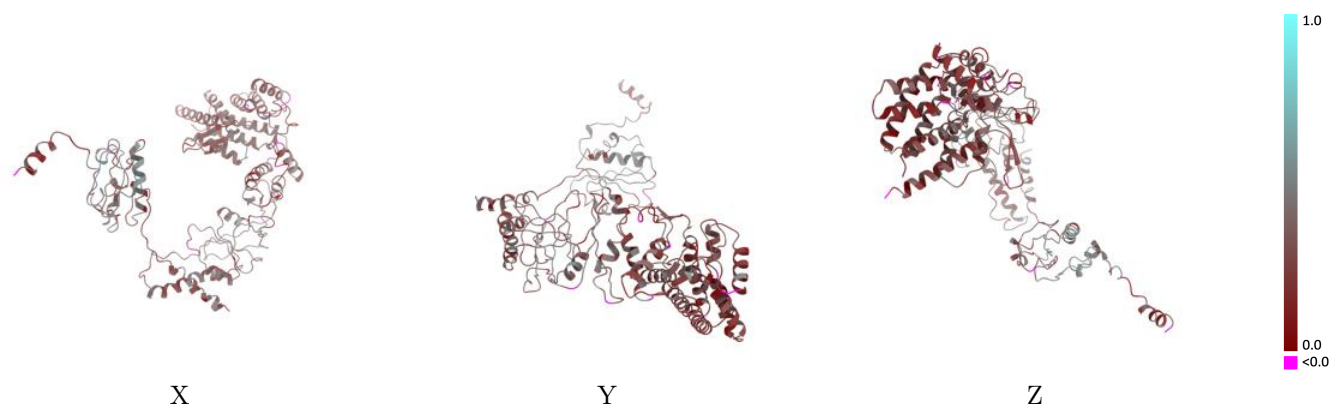
Y



Z

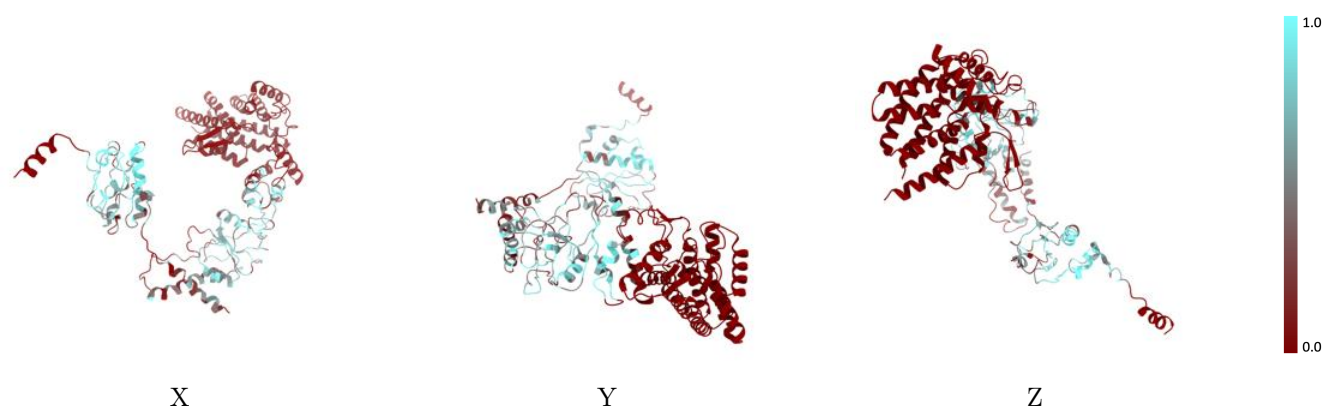
The images above show the 3D surface view of the map at the recommended contour level 0.616 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



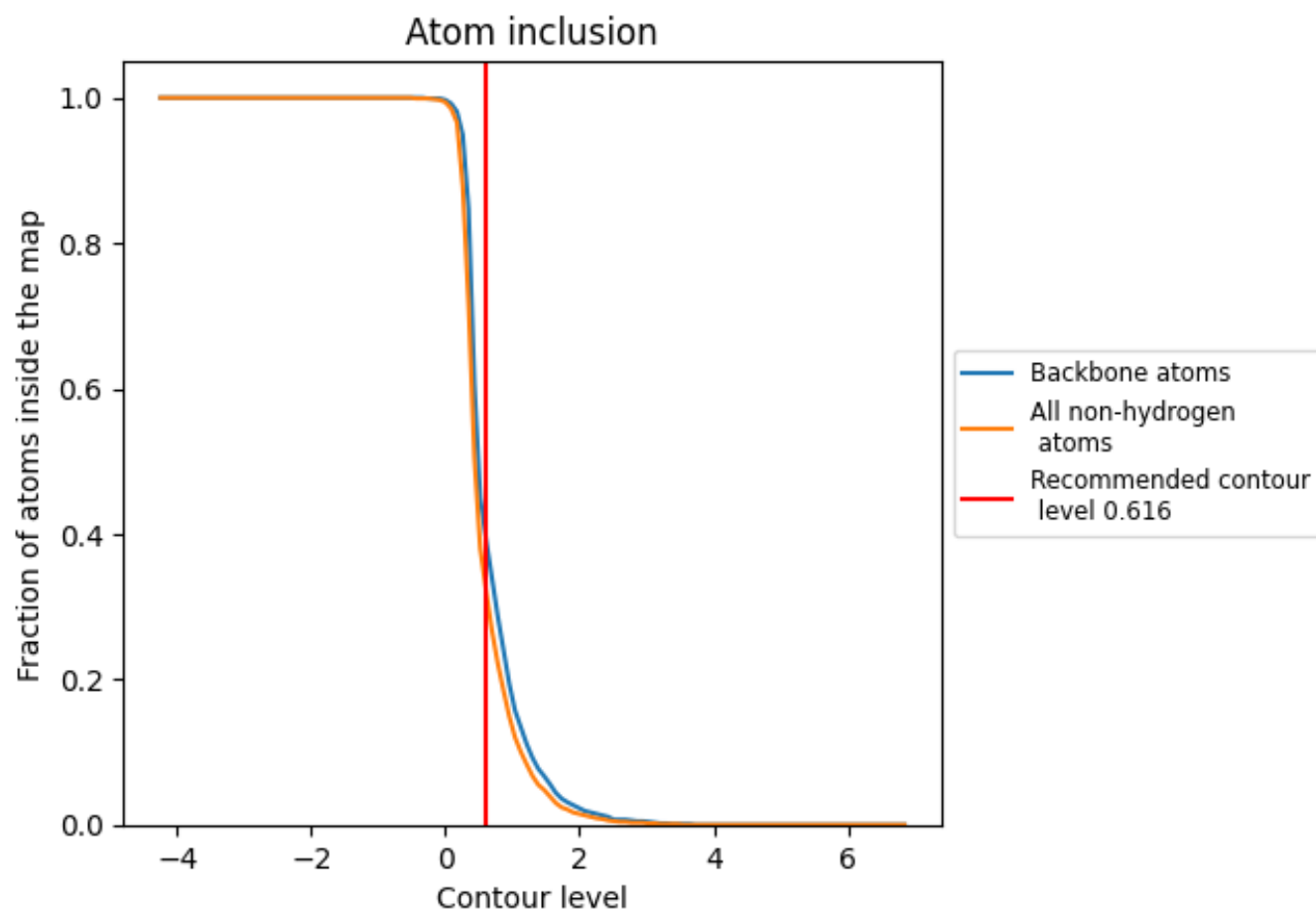
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.616).

9.4 Atom inclusion [i](#)



At the recommended contour level, 39% of all backbone atoms, 32% 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 (0.616) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.3160	0.3180
D	0.2380	0.2980
Q	0.5420	0.3770

