



Full wwPDB EM Validation Report ⓘ

Mar 10, 2025 – 08:09 PM JST

PDB ID : 9J82
EMDB ID : EMD-61213
Title : Cryo-EM structure of wild type Aquifex aeolicus RseP in complex with Fab
Authors : Asahi, K.; Hirose, M.; Aruga, R.; Kato, T.; Nogi, T.
Deposited on : 2024-08-20
Resolution : 3.95 Å(reported)
Based on initial models : ., 8ZAY, 6AKQ, 3WKM

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.dev117
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.41.2

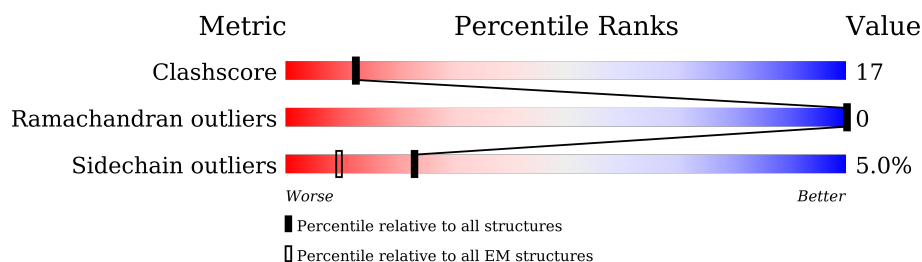
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.95 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	441	
2	H	239	
3	L	214	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6724 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 Putative zinc metalloprotease aq_1964.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	429	Total	C	N	O	S	0	0
			3386	2250	556	575	5		

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	430	GLY	-	expression tag	UNP O67776
A	431	ARG	-	expression tag	UNP O67776
A	432	GLY	-	expression tag	UNP O67776
A	433	SER	-	expression tag	UNP O67776
A	434	HIS	-	expression tag	UNP O67776
A	435	HIS	-	expression tag	UNP O67776
A	436	HIS	-	expression tag	UNP O67776
A	437	HIS	-	expression tag	UNP O67776
A	438	HIS	-	expression tag	UNP O67776
A	439	HIS	-	expression tag	UNP O67776
A	440	HIS	-	expression tag	UNP O67776
A	441	HIS	-	expression tag	UNP O67776

- Molecule 2 is a protein called VH-CH1 region of mouse monoclonal antibody IgG 4A9.

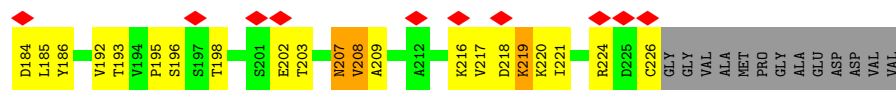
Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	219	Total	C	N	O	S	0	0
			1670	1059	279	325	7		

- Molecule 3 is a protein called L chain of mouse monoclonal antibody IgG 4A9.

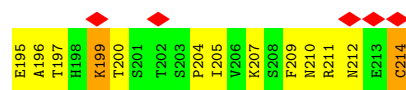
Mol	Chain	Residues	Atoms					AltConf	Trace
3	L	214	Total	C	N	O	S	0	0
			1667	1037	279	343	8		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	Zn	0
			1	1	



• Molecule 3: L chain of mouse monoclonal antibody IgG 4A9



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	88124	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	162	Depositor
Maximum defocus (nm)	3647	Depositor
Magnification	103703	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.489	Depositor
Minimum map value	-0.199	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.018	Depositor
Recommended contour level	0.16	Depositor
Map size (Å)	243.0, 243.0, 243.0	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.35, 1.35, 1.35	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, FME, SNN

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	0.25	0/3445	0.52	1/4668 (0.0%)
2	H	0.27	0/1713	0.59	0/2339
3	L	0.27	0/1705	0.53	0/2315
All	All	0.26	0/6863	0.54	1/9322 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	359	MET	CA-CB-CG	5.03	121.84	113.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	149	ILE	Mainchain

5.2 Too-close contacts

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	3386	0	3593	111	0
2	H	1670	0	1633	61	0
3	L	1667	0	1581	59	0
4	A	1	0	0	0	0
All	All	6724	0	6807	225	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 17.

All (225) 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:H:145:SER:N	2:H:196:SER:HG	1.69	0.90
1:A:148:LYS:HB3	1:A:178:PHE:HB2	1.57	0.86
1:A:377:ILE:HD11	1:A:405:GLN:HA	1.63	0.80
1:A:219:PRO:HB2	1:A:272:LEU:HD11	1.64	0.80
2:H:68:VAL:HB	2:H:72:PHE:HD2	1.49	0.78
1:A:226:LYS:HB2	1:A:263:ARG:HH12	1.49	0.77
2:H:135:LEU:HB3	3:L:118:PHE:HD1	1.50	0.76
2:H:8:GLU:H	2:H:116:GLN:HE22	1.32	0.75
1:A:113:GLY:HA2	1:A:296:GLU:O	1.86	0.75
1:A:233:GLU:HB2	1:A:260:LYS:HD3	1.67	0.74
1:A:349:ALA:HA	1:A:356:TYR:HE1	1.52	0.74
1:A:279:LYS:HG3	1:A:280:THR:HG23	1.72	0.71
1:A:412:ILE:O	1:A:416:THR:HG23	1.91	0.71
3:L:155:ARG:NH1	3:L:185:GLU:OE2	2.24	0.70
3:L:189:HIS:O	3:L:211:ARG:NH1	2.25	0.69
2:H:54:ARG:NH2	2:H:63:GLU:OE2	2.26	0.68
2:H:33:ASP:O	2:H:55:ASN:ND2	2.26	0.68
1:A:149:ILE:HD13	1:A:177:LEU:HD12	1.74	0.67
1:A:34:ILE:HB	1:A:65:LYS:HB2	1.77	0.67
3:L:110:ASP:OD2	3:L:199:LYS:NZ	2.28	0.67
2:H:220:LYS:NZ	2:H:221:ILE:O	2.27	0.66
1:A:263:ARG:O	1:A:266:LYS:NZ	2.26	0.65
2:H:71:ARG:NH2	2:H:89:THR:O	2.29	0.65
1:A:410:ALA:O	1:A:414:THR:HG23	1.97	0.65
3:L:113:PRO:HB3	3:L:139:PHE:HB3	1.79	0.64
1:A:210:VAL:HG13	1:A:290:LYS:HB3	1.80	0.63
1:A:42:PRO:HA	1:A:56:ALA:HA	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:GLN:OE1	1:A:422:ASN:ND2	2.32	0.62
2:H:135:LEU:HB3	3:L:118:PHE:CD1	2.32	0.62
3:L:54:ARG:NH1	3:L:59:PRO:O	2.32	0.62
3:L:117:ILE:HD12	3:L:209:PHE:HD2	1.64	0.62
2:H:130:PRO:HB3	2:H:156:TYR:HB3	1.81	0.62
2:H:25:VAL:HG22	2:H:82:ILE:HD12	1.80	0.62
3:L:117:ILE:C	3:L:118:PHE:HD2	2.03	0.62
1:A:69:GLU:HA	1:A:79:PHE:HB3	1.80	0.62
1:A:322:ILE:HD13	1:A:325:LEU:HD21	1.82	0.61
1:A:28:PHE:O	1:A:82:LYS:NZ	2.34	0.61
1:A:148:LYS:NZ	1:A:150:SNM:O	2.32	0.61
3:L:16:GLY:N	3:L:78:VAL:O	2.31	0.60
1:A:339:ILE:HG21	1:A:427:LEU:HD11	1.82	0.60
2:H:174:VAL:HG22	2:H:192:VAL:HG12	1.81	0.60
3:L:87:PHE:HB3	3:L:101:GLY:HA2	1.84	0.60
3:L:150:ILE:HD12	3:L:192:TYR:HA	1.84	0.60
3:L:33:VAL:HG21	3:L:71:PHE:CE2	2.36	0.60
1:A:38:GLY:H	1:A:62:GLY:HA2	1.66	0.60
1:A:104:LEU:O	1:A:108:LEU:HG	2.01	0.59
1:A:413:ILE:O	1:A:417:ILE:HG12	2.02	0.59
3:L:159:VAL:HG22	3:L:179:LEU:HD11	1.84	0.59
1:A:267:MET:HE2	1:A:267:MET:HA	1.83	0.59
1:A:392:ARG:HG3	1:A:394:ARG:HG2	1.83	0.59
3:L:181:LEU:HD13	3:L:185:GLU:HG3	1.84	0.59
1:A:224:GLY:O	1:A:263:ARG:NH2	2.37	0.58
3:L:61:ARG:NH1	3:L:82:ASP:OD2	2.33	0.58
1:A:277:ASP:HB2	1:A:284:PHE:HB3	1.86	0.58
2:H:224:ARG:HH22	3:L:119:PRO:HB2	1.68	0.58
2:H:72:PHE:HD1	2:H:87:VAL:HG22	1.67	0.58
2:H:126:LYS:HD3	2:H:127:THR:N	2.19	0.58
3:L:115:VAL:HG12	3:L:136:LEU:HD13	1.84	0.58
1:A:156:THR:HG23	1:A:158:GLU:H	1.69	0.57
1:A:47:GLN:NE2	1:A:49:GLY:O	2.38	0.57
1:A:250:ARG:O	1:A:253:GLN:NE2	2.36	0.57
2:H:41:GLN:NE2	2:H:45:LYS:O	2.28	0.56
1:A:154:VAL:HG23	1:A:159:ASP:HB2	1.87	0.56
2:H:135:LEU:HD21	2:H:152:LEU:HB2	1.88	0.56
3:L:33:VAL:HG21	3:L:71:PHE:HE2	1.70	0.56
1:A:19:PHE:CE1	1:A:23:LEU:HD21	2.40	0.56
1:A:349:ALA:HA	1:A:356:TYR:CE1	2.38	0.56
1:A:353:PHE:HA	1:A:356:TYR:CD2	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:MET:O	1:A:363:ILE:HG12	2.05	0.56
1:A:382:LEU:H	1:A:382:LEU:HD23	1.70	0.56
1:A:103:ILE:HG22	1:A:311:THR:HG21	1.88	0.56
1:A:200:GLU:OE1	1:A:202:GLY:N	2.39	0.55
1:A:241:THR:HG22	1:A:243:TYR:H	1.71	0.55
3:L:19:VAL:HG12	3:L:75:ILE:HB	1.89	0.55
2:H:181:LEU:HB2	2:H:186:TYR:CE1	2.42	0.55
2:H:72:PHE:CD1	2:H:87:VAL:HG22	2.43	0.54
1:A:86:GLN:O	1:A:90:ILE:HG12	2.07	0.54
2:H:35:TYR:HB2	2:H:103:ASP:CG	2.28	0.54
2:H:22:LEU:HD13	2:H:85:LEU:HD23	1.90	0.53
2:H:68:VAL:HB	2:H:72:PHE:CD2	2.36	0.53
2:H:8:GLU:H	2:H:116:GLN:NE2	2.04	0.53
3:L:144:ILE:HD12	3:L:145:ASN:H	1.74	0.53
1:A:39:PHE:N	1:A:58:LEU:O	2.35	0.53
2:H:46:ALA:HB2	3:L:100:ALA:HA	1.91	0.53
1:A:115:GLU:HG3	1:A:295:VAL:HG22	1.91	0.53
3:L:26:SER:OG	3:L:27:GLN:OE1	2.27	0.52
3:L:190:ASN:HD21	3:L:212:ASN:HD21	1.55	0.52
3:L:210:ASN:HB3	3:L:212:ASN:OD1	2.09	0.52
1:A:83:LYS:N	1:A:86:GLN:OE1	2.40	0.52
1:A:239:ILE:HG12	1:A:244:GLU:HB3	1.91	0.52
1:A:374:PRO:O	1:A:375:LEU:HD23	2.09	0.52
2:H:74:ILE:HG22	2:H:85:LEU:HD13	1.91	0.52
3:L:24:LYS:HA	3:L:69:THR:O	2.10	0.52
1:A:84:PRO:HA	1:A:87:LYS:HD3	1.91	0.51
1:A:355:PRO:HA	1:A:358:VAL:HG12	1.91	0.51
1:A:302:GLU:OE1	1:A:302:GLU:N	2.41	0.51
1:A:19:PHE:O	1:A:23:LEU:HG	2.11	0.51
3:L:21:ILE:HG23	3:L:73:LEU:HB3	1.93	0.51
3:L:112:ALA:HA	3:L:200:THR:HG21	1.93	0.51
2:H:3:GLU:OE2	2:H:3:GLU:N	2.30	0.50
1:A:240:ASN:N	1:A:244:GLU:OE1	2.45	0.50
2:H:145:SER:OG	2:H:146:MET:N	2.44	0.50
2:H:203:THR:HB	2:H:220:LYS:HE2	1.94	0.50
2:H:221:ILE:O	2:H:221:ILE:HG13	2.12	0.50
3:L:168:SER:OG	3:L:169:LYS:NZ	2.42	0.50
2:H:81:SER:C	2:H:82:ILE:HD13	2.32	0.50
1:A:31:LYS:NZ	1:A:50:GLU:OE2	2.42	0.50
3:L:18:ARG:HA	3:L:75:ILE:O	2.11	0.50
3:L:85:ASP:N	3:L:85:ASP:OD1	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:LYS:O	1:A:218:SER:OG	2.25	0.49
2:H:198:THR:O	2:H:202:GLU:HG3	2.13	0.49
1:A:213:GLY:HA2	1:A:227:PRO:HB3	1.95	0.49
2:H:164:THR:HG23	2:H:207:ASN:HB2	1.94	0.49
2:H:42:PRO:HB2	2:H:45:LYS:HB2	1.94	0.49
3:L:62:PHE:CD1	3:L:75:ILE:HD12	2.48	0.49
3:L:149:LYS:HD2	3:L:152:GLY:HA2	1.94	0.49
2:H:183:SER:O	2:H:185:LEU:HD22	2.13	0.49
1:A:426:ARG:HD3	1:A:429:ARG:HH21	1.76	0.48
2:H:162:THR:OG1	2:H:209:ALA:HB3	2.13	0.48
2:H:195:PRO:HB2	2:H:198:THR:HG23	1.94	0.48
1:A:100:LEU:O	1:A:103:ILE:HG13	2.14	0.48
1:A:217:GLY:O	1:A:222:GLN:NE2	2.47	0.48
2:H:37:SER:HG	2:H:111:PHE:HE2	1.59	0.48
3:L:22:THR:HG22	3:L:72:THR:HG22	1.96	0.48
1:A:377:ILE:HG12	1:A:408:GLY:HA3	1.96	0.47
1:A:128:TYR:HB3	1:A:204:ALA:HB3	1.96	0.47
3:L:124:GLN:HE22	3:L:131:SER:HG	1.61	0.47
1:A:415:LEU:O	1:A:419:VAL:HG13	2.14	0.47
1:A:180:GLU:OE2	1:A:183:GLY:N	2.46	0.47
2:H:146:MET:SD	2:H:193:THR:HG22	2.55	0.47
3:L:11:MET:N	3:L:105:GLU:OE1	2.47	0.47
3:L:197:THR:HG22	3:L:204:PRO:HB3	1.96	0.47
1:A:214:VAL:HG12	1:A:227:PRO:HG3	1.96	0.47
1:A:234:VAL:HG11	1:A:248:GLU:HG2	1.97	0.47
1:A:412:ILE:O	1:A:415:LEU:HG	2.15	0.47
1:A:219:PRO:O	1:A:223:VAL:HG23	2.14	0.46
2:H:224:ARG:HG3	2:H:226:CYS:O	2.15	0.46
1:A:340:ALA:O	1:A:343:GLN:HG2	2.15	0.46
2:H:136:ALA:O	2:H:224:ARG:HD3	2.14	0.46
1:A:211:VAL:HG21	1:A:287:LEU:HD11	1.97	0.46
2:H:29:PHE:CE2	2:H:102:ARG:HD2	2.50	0.46
3:L:24:LYS:HD2	3:L:69:THR:O	2.16	0.46
1:A:15:TRP:O	1:A:19:PHE:HB2	2.16	0.46
2:H:126:LYS:HE2	2:H:126:LYS:HA	1.98	0.46
1:A:124:VAL:HA	1:A:157:TRP:NE1	2.31	0.46
2:H:126:LYS:HD3	2:H:127:THR:H	1.80	0.45
1:A:211:VAL:HG22	1:A:287:LEU:HD21	1.97	0.45
1:A:403:TYR:O	1:A:407:VAL:HG13	2.16	0.45
2:H:81:SER:H	2:H:81:SER:HG	1.50	0.45
1:A:6:PHE:HB2	1:A:342:ALA:HA	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:116:SER:HB2	3:L:118:PHE:HE2	1.82	0.45
2:H:31:PHE:O	2:H:76:ARG:NH2	2.50	0.45
3:L:195:GLU:OE1	3:L:195:GLU:N	2.38	0.45
3:L:184:ASP:O	3:L:188:ARG:HG2	2.17	0.45
1:A:362:PHE:O	1:A:365:LEU:HG	2.16	0.45
3:L:196:ALA:HB3	3:L:205:ILE:HG23	2.00	0.44
2:H:15:GLN:OE1	2:H:15:GLN:N	2.47	0.44
2:H:220:LYS:HA	2:H:220:LYS:HD2	1.79	0.44
1:A:226:LYS:HB2	1:A:263:ARG:NH1	2.26	0.44
1:A:295:VAL:O	1:A:297:LYS:NZ	2.51	0.44
1:A:250:ARG:HA	1:A:283:TYR:CZ	2.53	0.44
2:H:40:ARG:HG3	2:H:48:GLU:HB3	1.99	0.44
3:L:194:CYS:O	3:L:194:CYS:SG	2.76	0.44
1:A:232:LEU:CD2	1:A:262:LEU:HB2	2.49	0.43
3:L:212:ASN:OD1	3:L:212:ASN:N	2.48	0.43
2:H:51:ALA:HB3	2:H:74:ILE:HG21	2.00	0.43
3:L:188:ARG:O	3:L:189:HIS:ND1	2.52	0.43
1:A:252:SER:O	1:A:255:LYS:HG2	2.19	0.43
2:H:226:CYS:HB3	3:L:214:CYS:HB2	1.62	0.43
3:L:48:ILE:HG13	3:L:53:ILE:O	2.18	0.43
3:L:68:GLY:N	3:L:71:PHE:HE1	2.17	0.43
3:L:71:PHE:N	3:L:71:PHE:CD1	2.85	0.43
1:A:246:VAL:O	1:A:250:ARG:HG3	2.19	0.43
1:A:262:LEU:HG	1:A:265:GLY:H	1.83	0.43
1:A:99:PHE:O	1:A:103:ILE:HG23	2.19	0.43
1:A:152:TYR:CZ	1:A:166:ARG:HD2	2.53	0.43
1:A:418:PHE:O	1:A:421:ILE:HG13	2.19	0.43
2:H:184:ASP:O	2:H:185:LEU:HD13	2.18	0.43
2:H:95:SER:HG	2:H:122:VAL:H	1.62	0.42
1:A:311:THR:HA	1:A:314:LEU:CD2	2.49	0.42
1:A:30:VAL:HG22	1:A:66:LEU:HD12	2.01	0.42
1:A:89:LEU:HD12	1:A:92:LEU:HD21	2.00	0.42
2:H:89:THR:O	2:H:91:ARG:HD3	2.19	0.42
1:A:43:ILE:HG22	1:A:55:ILE:O	2.20	0.42
1:A:272:LEU:HD12	1:A:273:ILE:H	1.85	0.42
1:A:43:ILE:N	1:A:55:ILE:O	2.51	0.42
1:A:106:PHE:HA	1:A:109:VAL:HG12	2.01	0.42
1:A:134:ILE:O	1:A:138:ILE:HG22	2.19	0.42
2:H:208:VAL:HG23	2:H:217:VAL:HB	2.01	0.42
3:L:160:LEU:HD23	3:L:160:LEU:HA	1.85	0.42
3:L:36:TYR:HB2	3:L:87:PHE:CE1	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:142:LYS:O	3:L:144:ILE:HG22	2.19	0.42
1:A:277:ASP:O	1:A:281:GLY:N	2.46	0.42
1:A:21:HIS:CE1	1:A:94:GLY:H	2.38	0.41
1:A:290:LYS:HE2	1:A:290:LYS:HB2	1.71	0.41
1:A:235:ASN:HD21	1:A:257:ILE:HA	1.85	0.41
2:H:130:PRO:HB2	2:H:153:VAL:HG23	2.01	0.41
1:A:124:VAL:HG23	1:A:157:TRP:CE2	2.56	0.41
1:A:164:LEU:HD12	1:A:165:ILE:N	2.36	0.41
3:L:17:ASP:O	3:L:78:VAL:N	2.48	0.41
1:A:39:PHE:O	1:A:58:LEU:N	2.53	0.41
3:L:149:LYS:O	3:L:150:ILE:HD13	2.21	0.41
2:H:81:SER:O	2:H:82:ILE:HD13	2.21	0.41
3:L:21:ILE:HD13	3:L:102:THR:HG21	2.03	0.41
1:A:326:ILE:HD12	1:A:326:ILE:HA	1.93	0.41
2:H:114:TRP:HB2	3:L:43:SER:HB2	2.03	0.41
3:L:170:ASP:OD2	3:L:172:THR:OG1	2.24	0.41
1:A:51:THR:O	1:A:51:THR:OG1	2.39	0.41
1:A:248:GLU:O	1:A:251:LYS:HG3	2.20	0.41
1:A:252:SER:HA	1:A:255:LYS:HE3	2.02	0.41
3:L:6:GLN:NE2	3:L:102:THR:OG1	2.54	0.41
3:L:83:LEU:HD13	3:L:83:LEU:HA	1.94	0.41
1:A:254:GLY:HA2	1:A:275:ALA:HA	2.01	0.41
2:H:219:LYS:HA	2:H:219:LYS:HD3	1.83	0.41
1:A:28:PHE:CE2	1:A:89:LEU:HD23	2.56	0.40
1:A:241:THR:N	1:A:244:GLU:OE1	2.54	0.40
1:A:385:LEU:HD11	1:A:404:TRP:NE1	2.36	0.40
1:A:181:ARG:HG2	1:A:182:ASN:OD1	2.20	0.40
1:A:27:LEU:HD12	1:A:27:LEU:HA	1.83	0.40
1:A:374:PRO:HB3	1:A:384:LEU:HD11	2.03	0.40
2:H:36:MET:HB3	2:H:83:LEU:HD22	2.02	0.40
1:A:28:PHE:HE2	1:A:89:LEU:HD23	1.87	0.40
1:A:102:THR:HG21	1:A:365:LEU:HB3	2.03	0.40
2:H:112:ASP:OD2	2:H:113:TYR:CD2	2.74	0.40
2:H:146:MET:HE2	2:H:146:MET:HA	2.03	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	425/441 (96%)	409 (96%)	16 (4%)	0	100	100
2	H	215/239 (90%)	213 (99%)	2 (1%)	0	100	100
3	L	212/214 (99%)	202 (95%)	10 (5%)	0	100	100
All	All	852/894 (95%)	824 (97%)	28 (3%)	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	359/369 (97%)	341 (95%)	18 (5%)	20	43
2	H	189/201 (94%)	179 (95%)	10 (5%)	19	42
3	L	191/191 (100%)	182 (95%)	9 (5%)	22	45
All	All	739/761 (97%)	702 (95%)	37 (5%)	23	43

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	PHE
1	A	19	PHE
1	A	29	ARG
1	A	35	PHE
1	A	89	LEU

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Mol	Chain	Res	Type
1	A	106	PHE
1	A	152	TYR
1	A	182	ASN
1	A	226	LYS
1	A	251	LYS
1	A	264	ASN
1	A	267	MET
1	A	314	LEU
1	A	359	MET
1	A	360	MET
1	A	394	ARG
1	A	404	TRP
1	A	426	ARG
2	H	40	ARG
2	H	86	GLN
2	H	112	ASP
2	H	123	SER
2	H	151	CYS
2	H	207	ASN
2	H	208	VAL
2	H	216	LYS
2	H	218	ASP
2	H	219	LYS
3	L	18	ARG
3	L	26	SER
3	L	48	ILE
3	L	85	ASP
3	L	118	PHE
3	L	178	THR
3	L	199	LYS
3	L	207	LYS
3	L	214	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	333	GLN
1	A	422	ASN
2	H	116	GLN
3	L	190	ASN
3	L	210	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
1	SNN	A	150	1	7,8,8	5.28	5 (71%)	7,11,11	1.67	1 (14%)
1	FME	A	1	1	8,9,10	0.37	0	7,9,11	0.87	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SNN	A	150	1	-	-	0/1/1/1
1	FME	A	1	1	-	1/7/9/11	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	150	SNN	C-N1	9.45	1.49	1.37
1	A	150	SNN	C4-C5	8.88	1.63	1.51
1	A	150	SNN	C5-N1	4.10	1.43	1.37
1	A	150	SNN	O-C	-2.29	1.18	1.23
1	A	150	SNN	O5-C5	-2.15	1.18	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	150	SNN	O5-C5-C4	3.67	131.24	126.39

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	FME	CB-CG-SD-CE

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	150	SNN	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

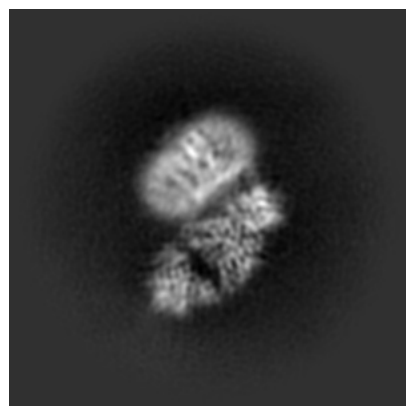
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-61213. These allow visual inspection of the internal detail of the map and identification of artifacts.

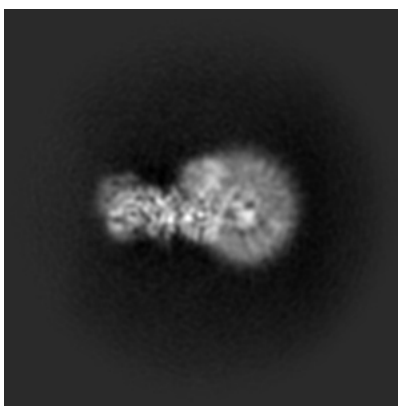
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

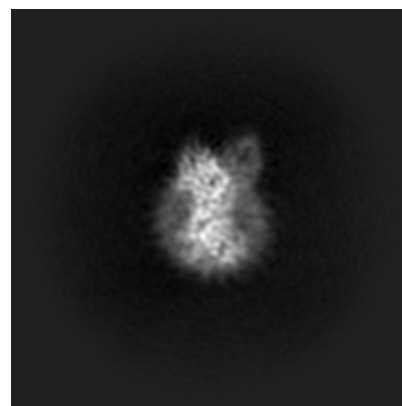
6.1.1 Primary map



X

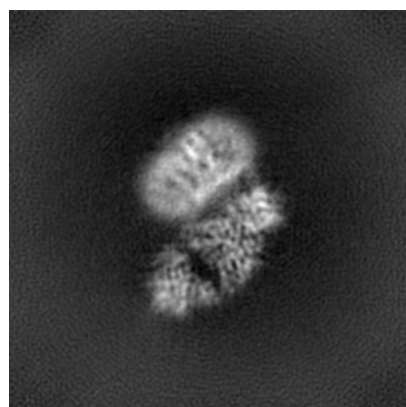


Y

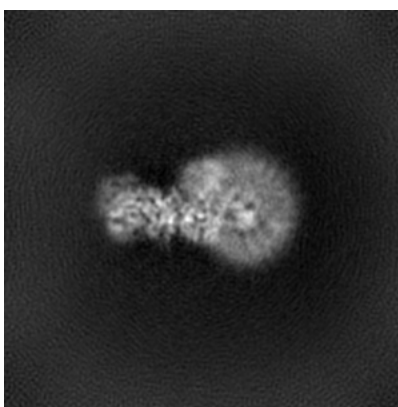


Z

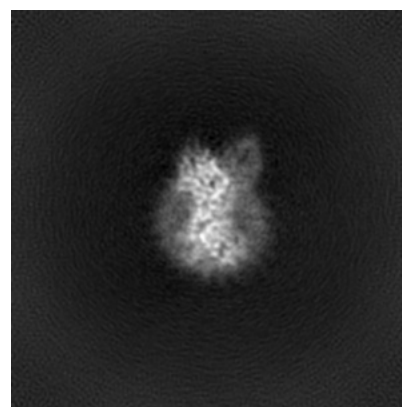
6.1.2 Raw map



X



Y

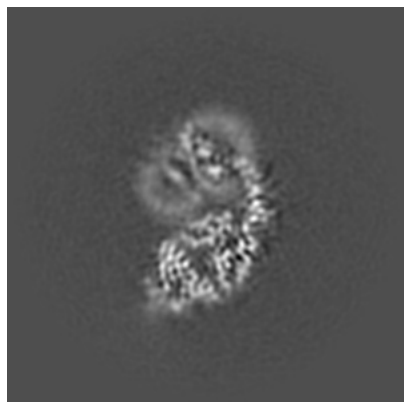


Z

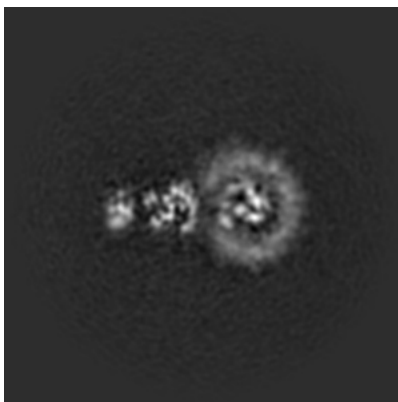
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

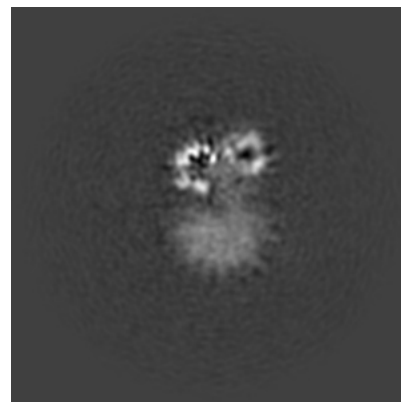
6.2.1 Primary map



X Index: 90

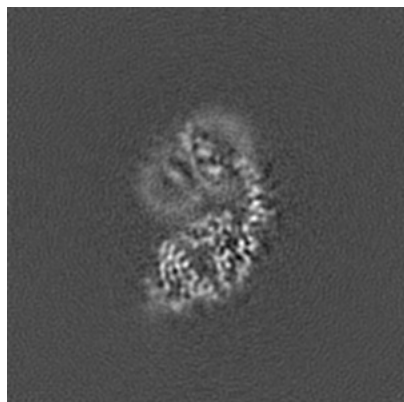


Y Index: 90

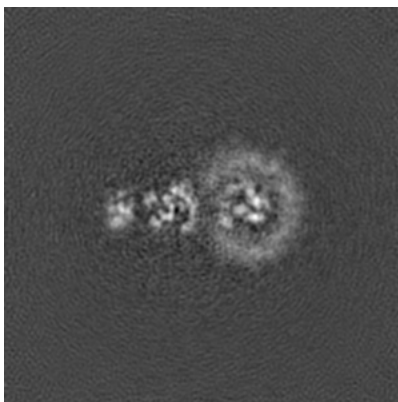


Z Index: 90

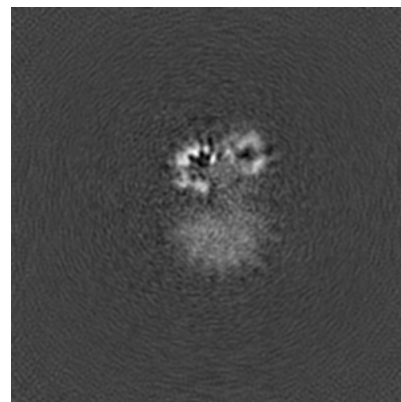
6.2.2 Raw map



X Index: 90



Y Index: 90

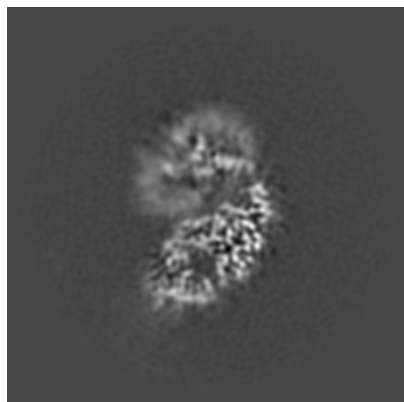


Z Index: 90

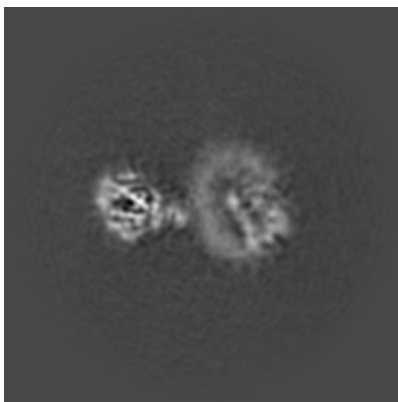
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

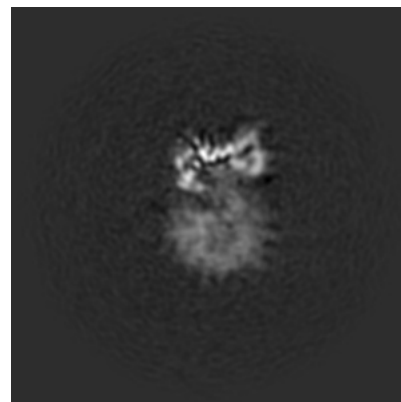
6.3.1 Primary map



X Index: 87

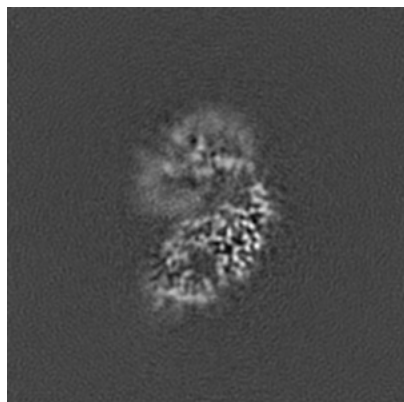


Y Index: 77

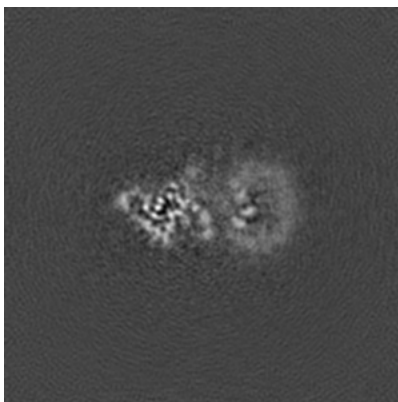


Z Index: 92

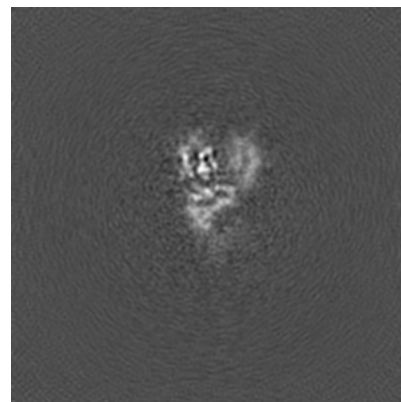
6.3.2 Raw map



X Index: 87



Y Index: 99

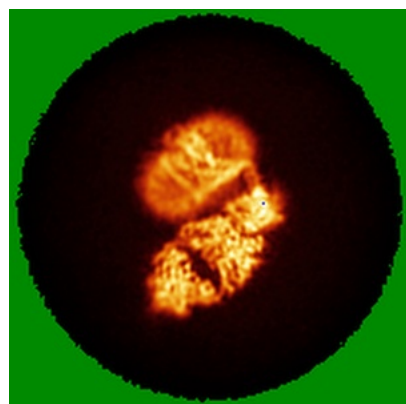


Z Index: 83

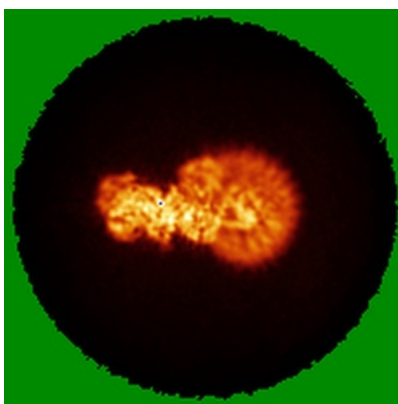
The images above show the largest variance slices of the map in three orthogonal directions.

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

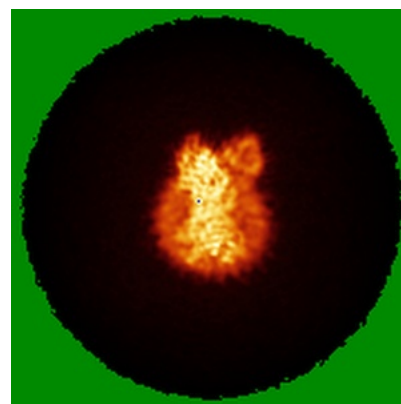
6.4.1 Primary map



X

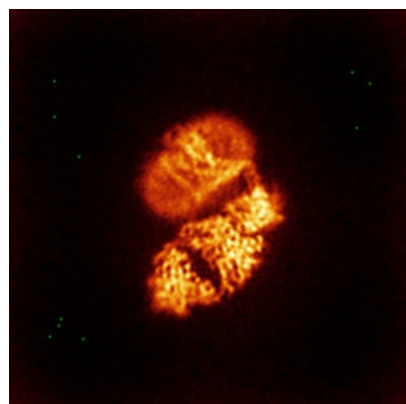


Y

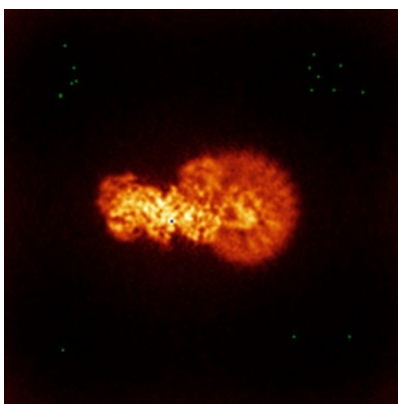


Z

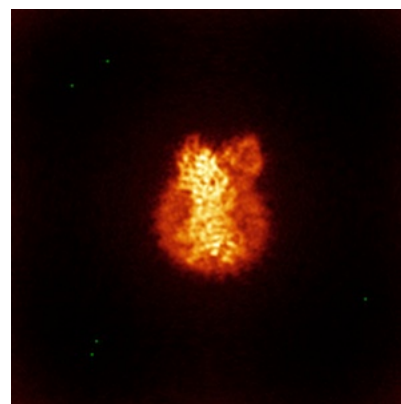
6.4.2 Raw map



X



Y

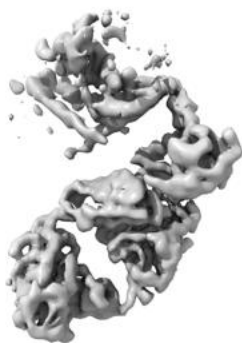


Z

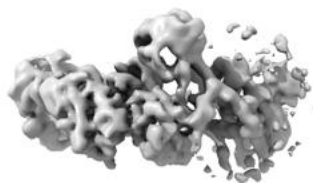
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

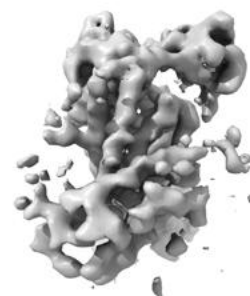
6.5.1 Primary map



X



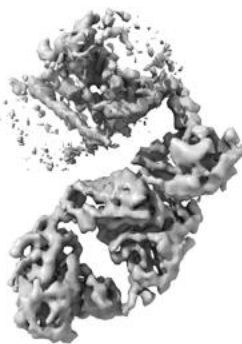
Y



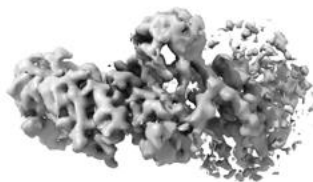
Z

The images above show the 3D surface view of the map at the recommended contour level 0.16. 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.

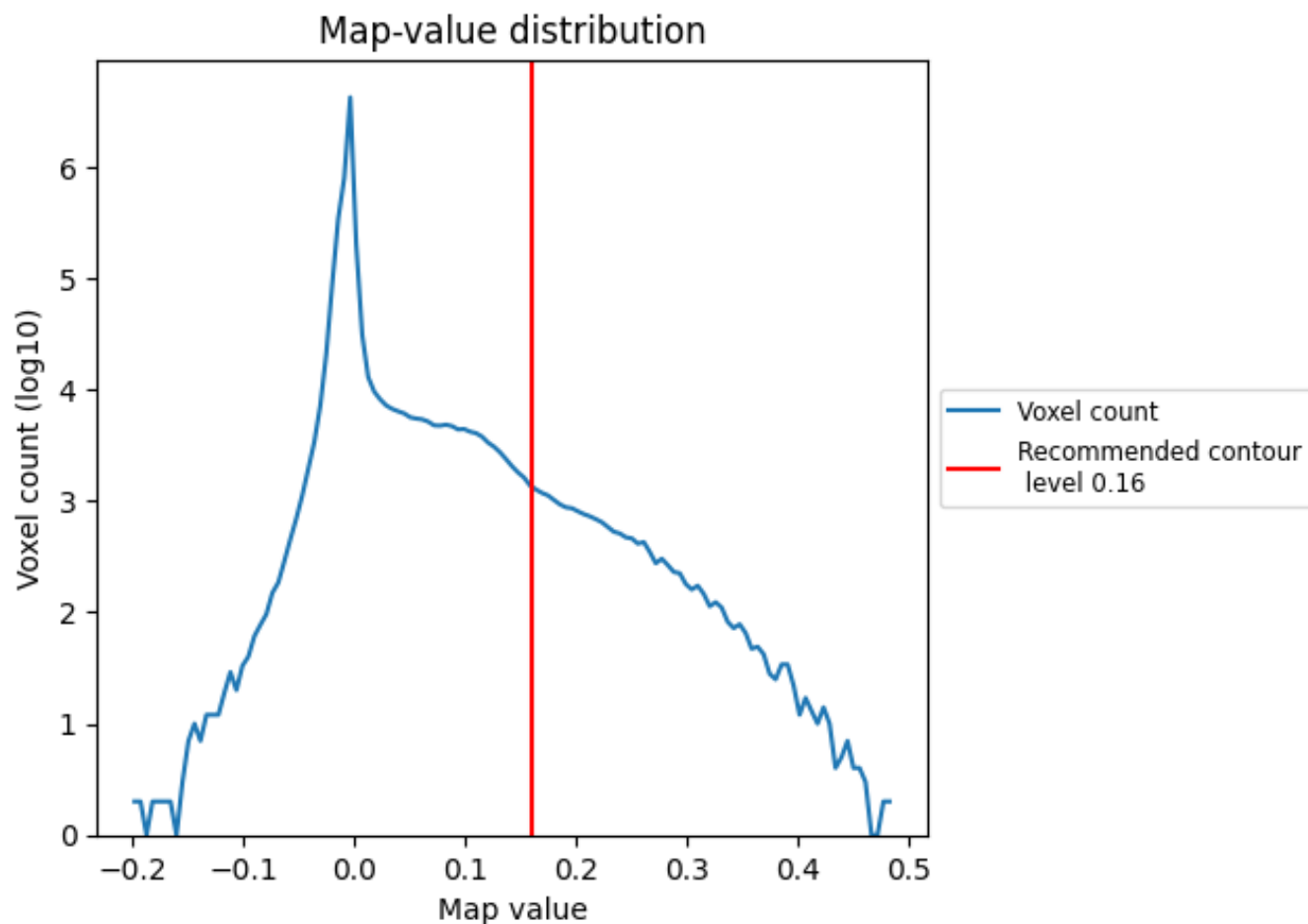
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

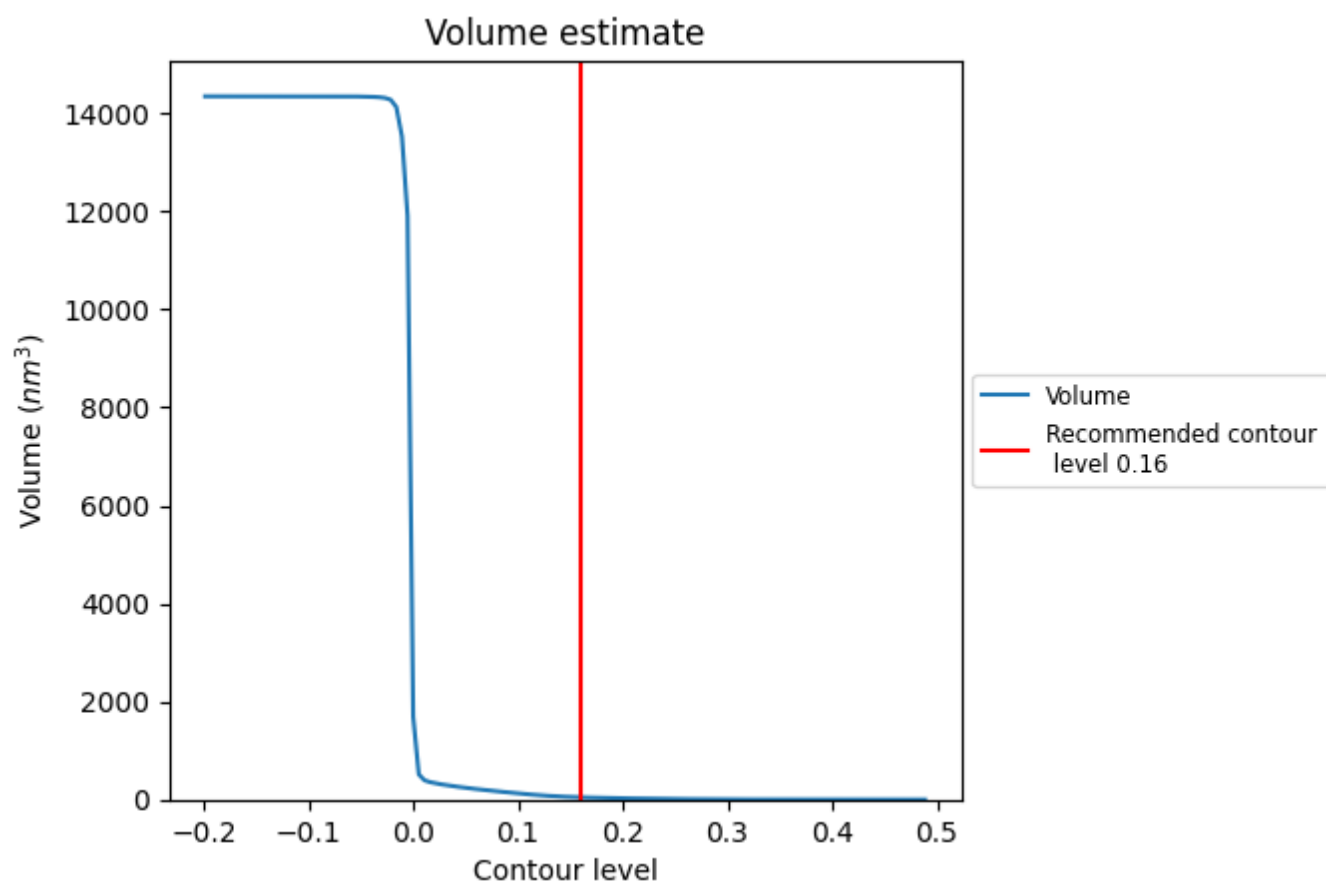
This section contains the results of statistical analysis of the map.

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.

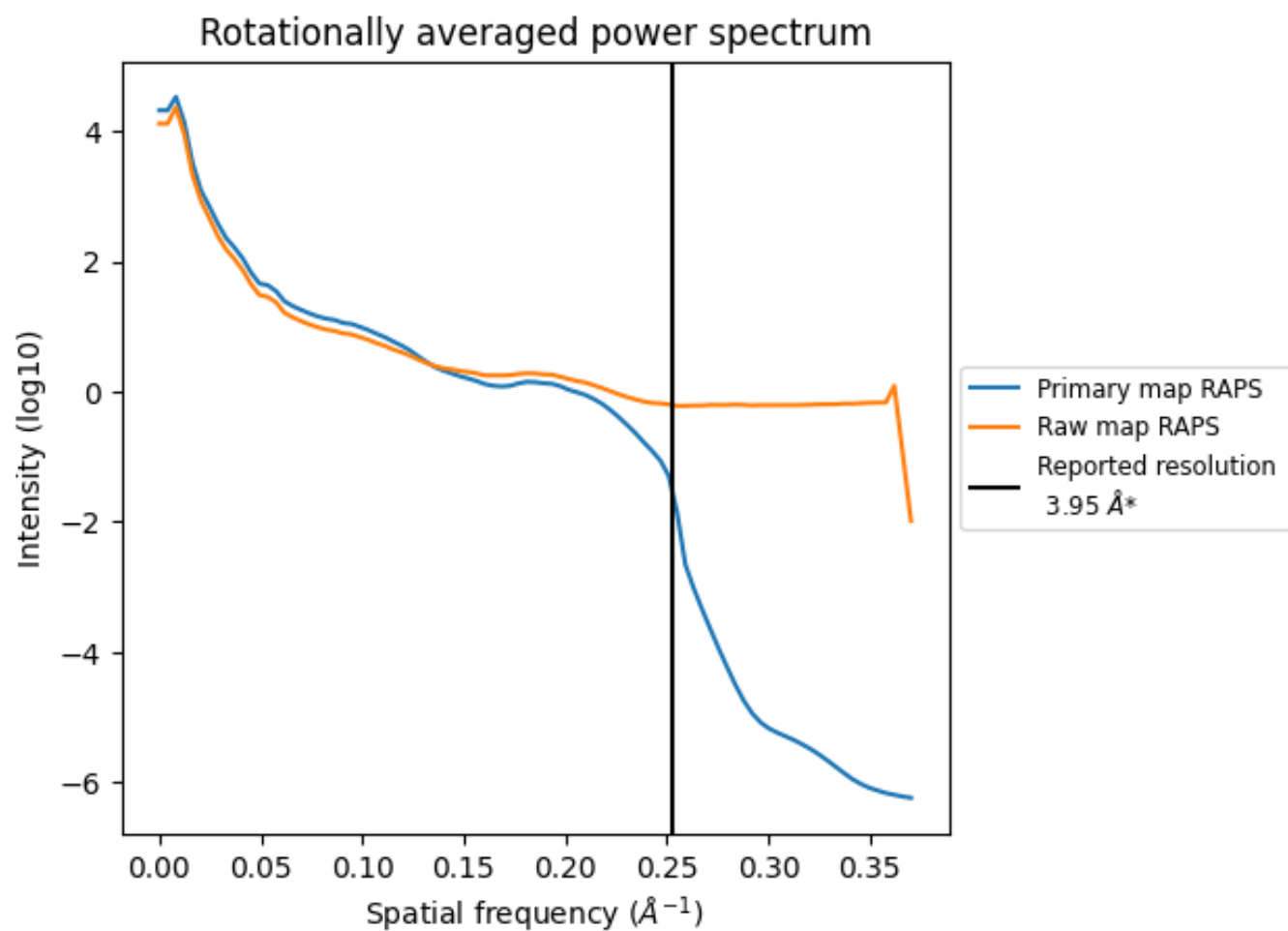
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 46 nm³; this corresponds to an approximate mass of 41 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 ⓘ

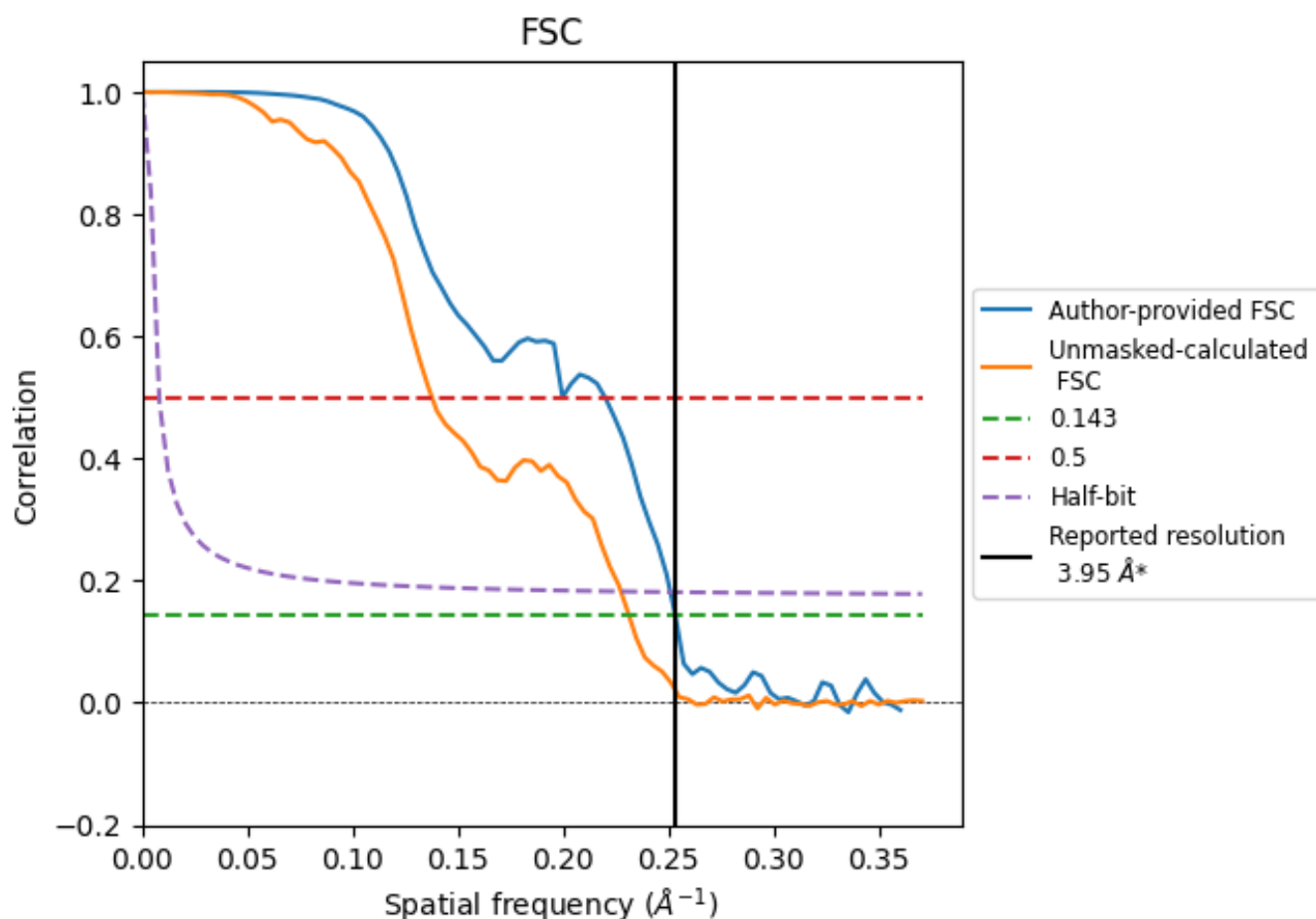


*Reported resolution corresponds to spatial frequency of 0.253 Å⁻¹

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 0.253 \AA^{-1}

8.2 Resolution estimates [i](#)

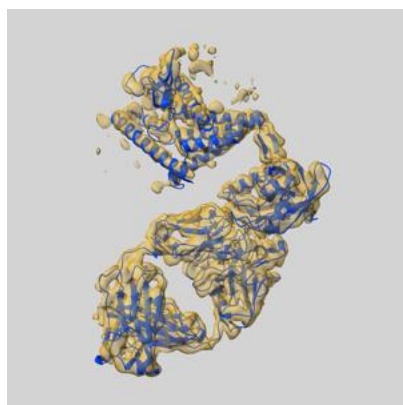
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.95	-	-
Author-provided FSC curve	3.95	4.55	3.99
Unmasked-calculated*	4.33	7.27	4.40

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

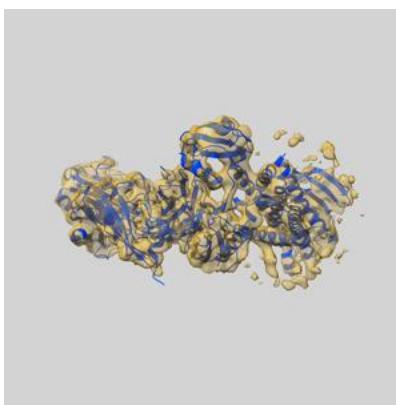
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-61213 and PDB model 9J82. Per-residue inclusion information can be found in section [3](#) on page [5](#).

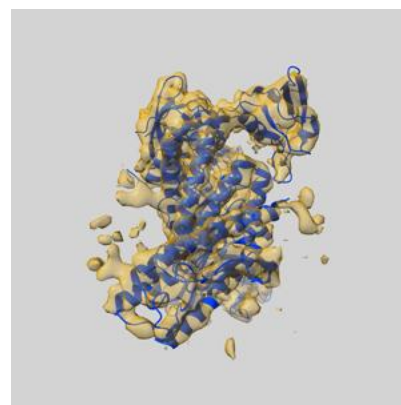
9.1 Map-model overlay [i](#)



X



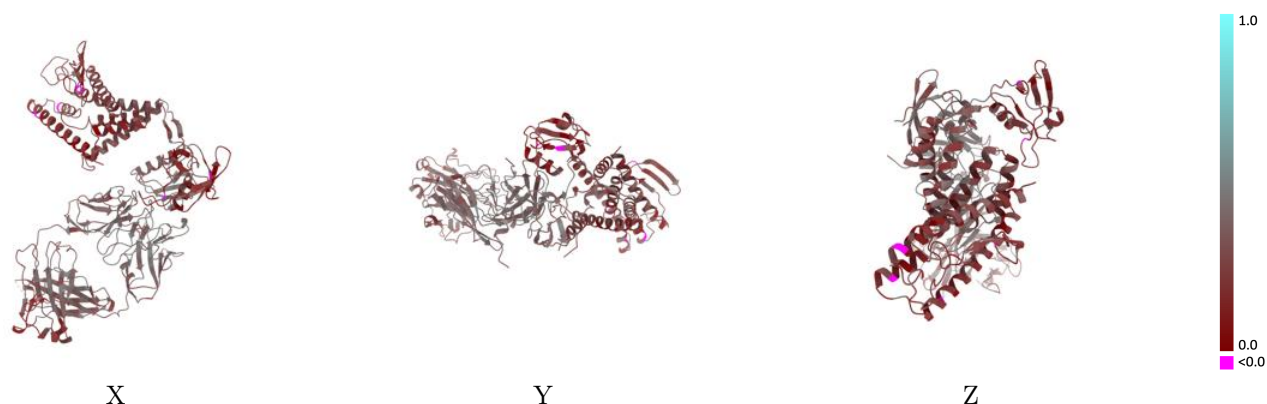
Y



Z

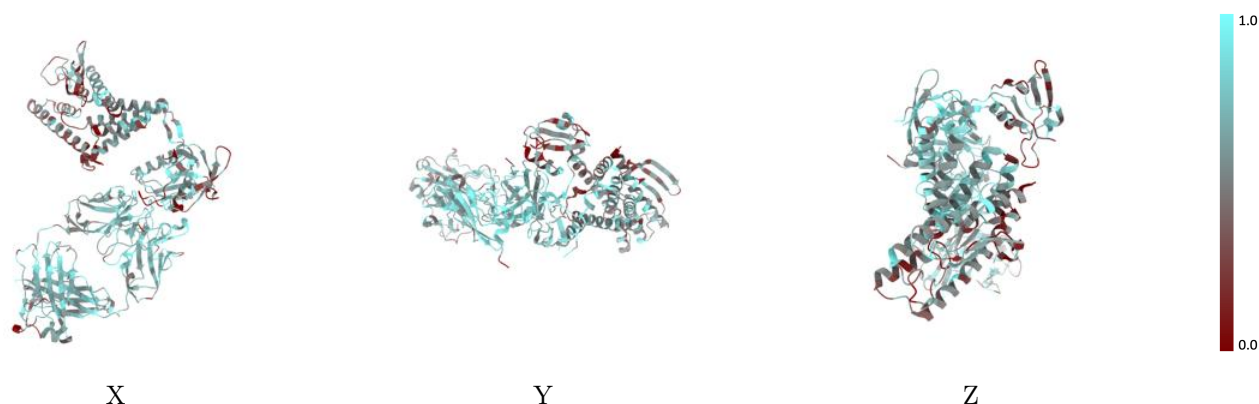
The images above show the 3D surface view of the map at the recommended contour level 0.16 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



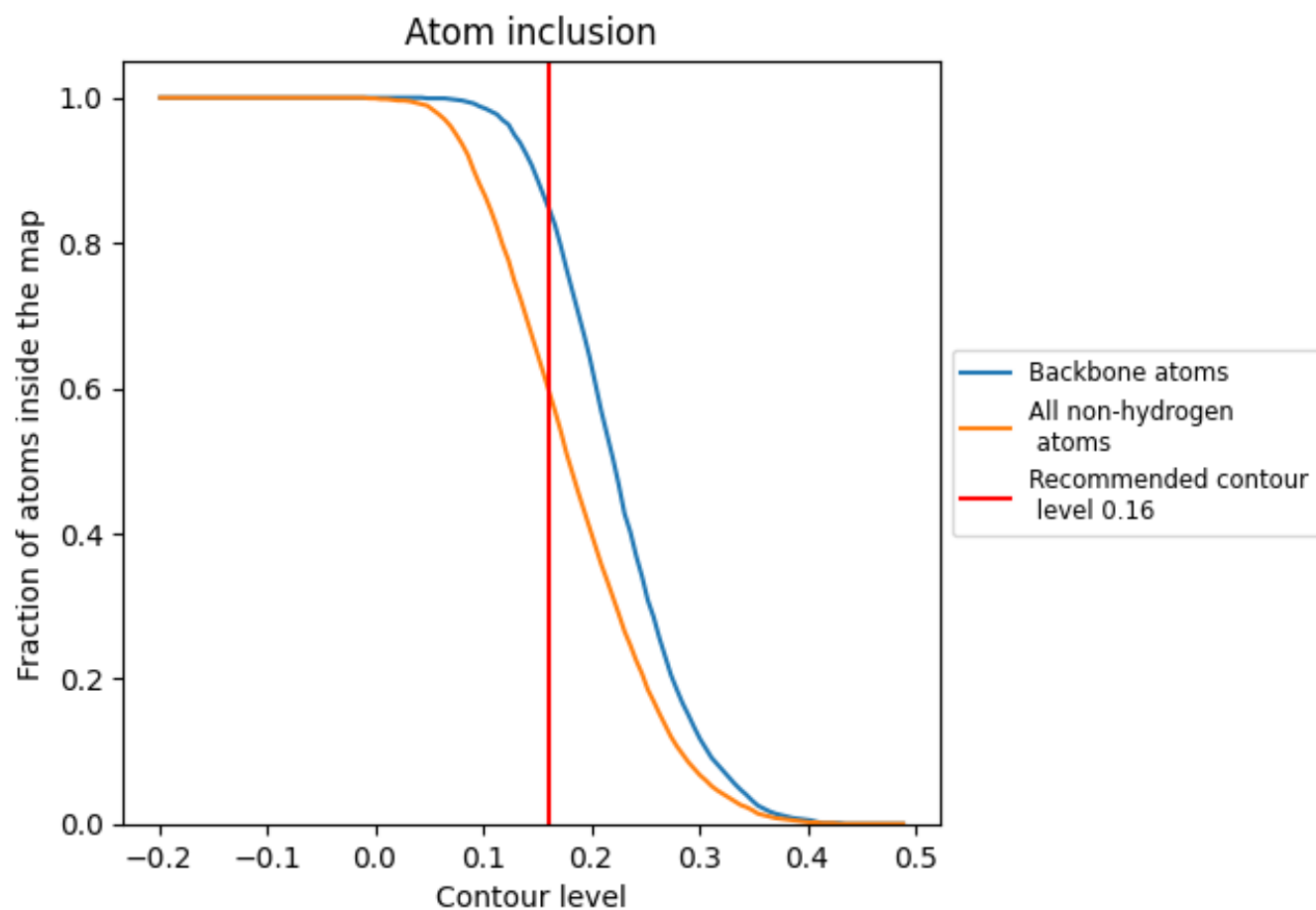
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.16).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 60% 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 (0.16) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.6000	<div></div> 0.3060
A	<div></div> 0.5220	<div></div> 0.2540
H	<div></div> 0.7010	<div></div> 0.3670
L	<div></div> 0.6580	<div></div> 0.3510

