



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 09:30 PM UTC

PDB ID : 6BIH / pdb_00006bih
EMDB ID : EMD-7100
Title : The Structure of the Actin-Smooth Muscle Myosin Motor Domain Complex
in the Rigor State
Authors : Taylor, K.A.; Banerjee, C.; Hu, Z.
Deposited on : 2017-11-02
Resolution : 6.00 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
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

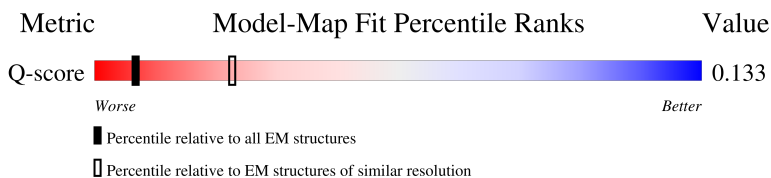
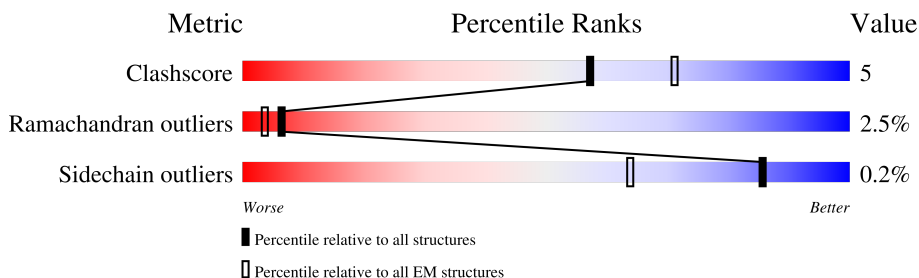
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 6.00 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	525 (5.50 - 6.50)

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	H	798	41% 82% 7% 9%
2	C	377	49% 91% 8% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17605 atoms, of which 8773 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 Myosin-11.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	H	729	Total	C	H	N	O	S	0	0
			11758	3739	5887	1005	1098	29		

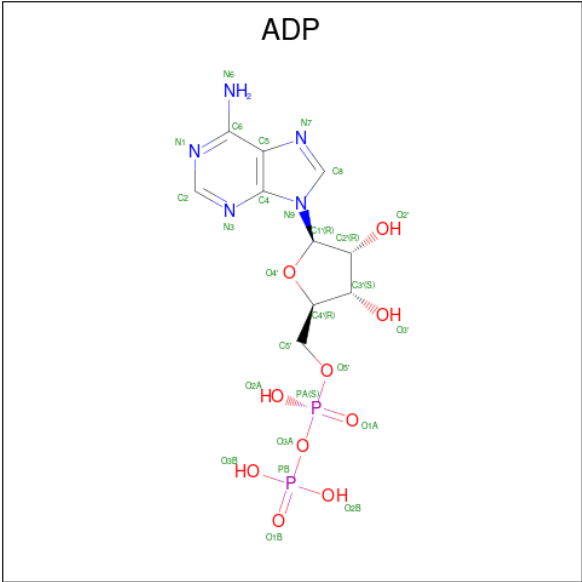
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	791	ASP	-	expression tag	UNP P10587
H	792	TYR	-	expression tag	UNP P10587
H	793	LYS	-	expression tag	UNP P10587
H	794	ASP	-	expression tag	UNP P10587
H	795	ASP	-	expression tag	UNP P10587
H	796	ASP	-	expression tag	UNP P10587
H	797	ASP	-	expression tag	UNP P10587
H	798	LYS	-	expression tag	UNP P10587

- Molecule 2 is a protein called Actin, alpha skeletal muscle.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	C	375	Total	C	H	N	O	S	0	0
			5819	1854	2886	493	565	21		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					AltConf
3	C	1	Total	C	N	O	P	0
			27	10	5	10	2	

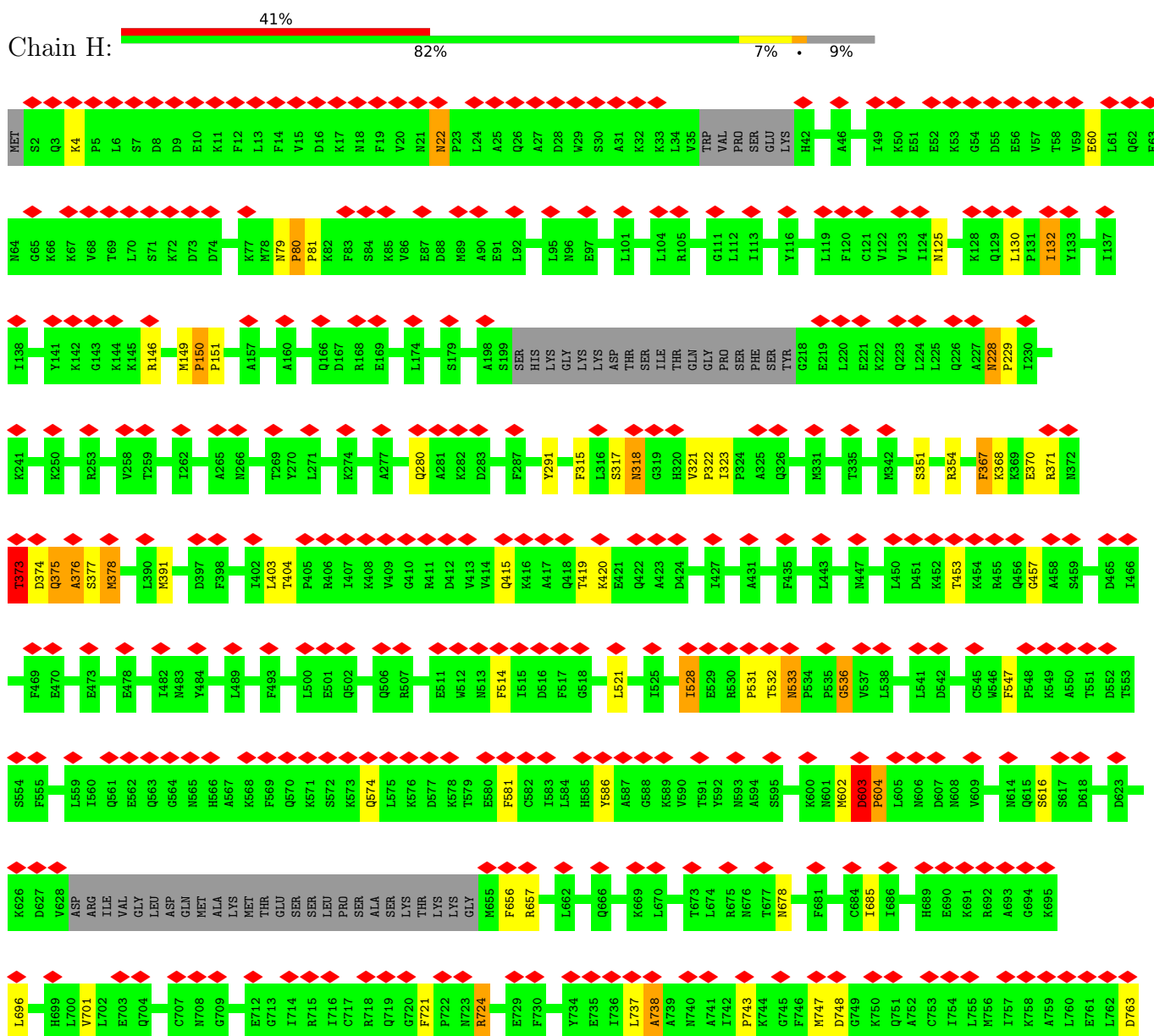
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

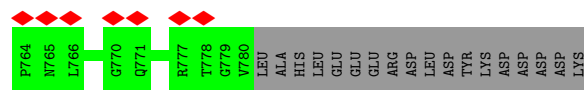
Mol	Chain	Residues	Atoms		AltConf
4	C	1	Total	Mg	0
			1	1	

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

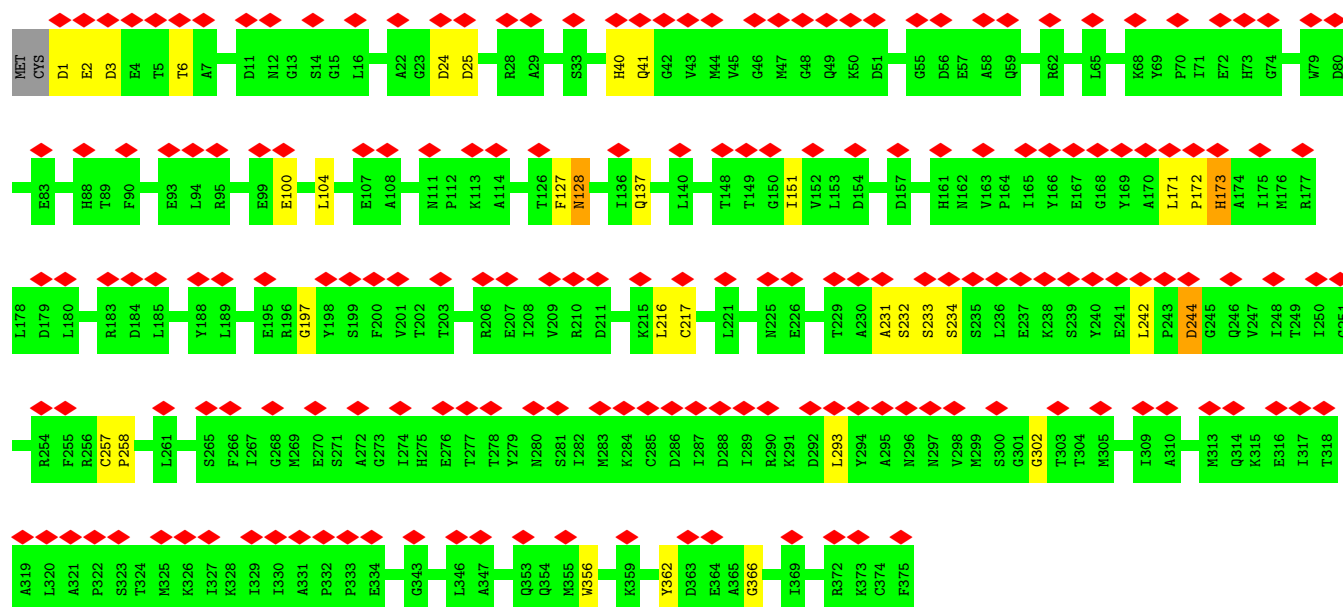
• Molecule 1: Myosin-11





- Molecule 2: Actin, alpha skeletal muscle

Chain C: 49% 91% 8% ..



4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=-166.779310°, rise=28.186850 Å, axial sym=C1	Depositor
Number of segments used	85000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{Å}^2$)	60	Depositor
Minimum defocus (nm)	2200	Depositor
Maximum defocus (nm)	5000	Depositor
Magnification	Not provided	
Image detector	DIRECT ELECTRON DE-20 (5k x 3k)	Depositor
Maximum map value	0.050	Depositor
Minimum map value	-0.029	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.016	Depositor
Map size (Å)	252.4416, 252.4416, 252.4416	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.9861, 0.9861, 0.9861	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ADP

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	H	0.94	3/5979 (0.1%)	1.13	42/8058 (0.5%)
2	C	0.55	0/2995	0.65	0/4055
All	All	0.83	3/8974 (0.0%)	0.99	42/12113 (0.3%)

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	H	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	528	ILE	C-N	-22.76	1.01	1.33
1	H	536	GLY	C-N	-16.58	1.12	1.33
1	H	60	GLU	CA-C	-5.36	1.46	1.52

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	375	GLN	N-CA-C	-18.71	90.12	111.82
1	H	373	THR	N-CA-C	-11.97	91.99	109.59
1	H	367	PHE	N-CA-C	9.72	123.68	110.55
1	H	80	PRO	CA-C-N	8.64	129.49	119.47
1	H	80	PRO	C-N-CA	8.64	129.49	119.47
1	H	547	PHE	CA-C-N	8.49	129.00	119.32
1	H	547	PHE	C-N-CA	8.49	129.00	119.32
1	H	125	ASN	CA-C-N	7.84	127.56	119.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	125	ASN	C-N-CA	7.84	127.56	119.64
1	H	130	LEU	CA-C-N	7.77	129.56	119.84
1	H	130	LEU	C-N-CA	7.77	129.56	119.84
1	H	4	LYS	CA-C-N	7.58	128.27	119.47
1	H	4	LYS	C-N-CA	7.58	128.27	119.47
1	H	404	THR	CA-C-N	7.58	128.04	119.93
1	H	404	THR	C-N-CA	7.58	128.04	119.93
1	H	149	MET	CA-C-N	7.38	127.98	120.38
1	H	149	MET	C-N-CA	7.38	127.98	120.38
1	H	22	ASN	CA-C-N	7.17	128.80	119.84
1	H	22	ASN	C-N-CA	7.17	128.80	119.84
1	H	685	ILE	N-CA-C	7.10	118.05	108.11
1	H	763	ASP	CA-C-N	7.02	128.62	119.84
1	H	763	ASP	C-N-CA	7.02	128.62	119.84
1	H	721	PHE	CA-C-N	6.68	127.22	119.47
1	H	721	PHE	C-N-CA	6.68	127.22	119.47
1	H	150	PRO	CA-C-N	6.38	126.76	119.93
1	H	150	PRO	C-N-CA	6.38	126.76	119.93
1	H	228	ASN	CA-C-N	6.35	126.27	119.28
1	H	228	ASN	C-N-CA	6.35	126.27	119.28
1	H	378	MET	CA-C-N	6.24	126.71	119.47
1	H	378	MET	C-N-CA	6.24	126.71	119.47
1	H	678	ASN	CA-C-N	6.08	127.44	119.84
1	H	678	ASN	C-N-CA	6.08	127.44	119.84
1	H	323	ILE	CA-C-N	6.05	126.41	119.93
1	H	323	ILE	C-N-CA	6.05	126.41	119.93
1	H	701	VAL	N-CA-C	6.02	116.19	110.53
1	H	315	PHE	N-CA-C	-5.95	102.11	110.50
1	H	514	PHE	CB-CA-C	-5.75	109.41	117.23
1	H	603	ASP	CA-C-N	5.69	126.95	119.84
1	H	603	ASP	C-N-CA	5.69	126.95	119.84
1	H	747	MET	N-CA-C	5.14	117.30	111.02
1	H	79	ASN	CA-C-N	5.08	125.61	120.38
1	H	79	ASN	C-N-CA	5.08	125.61	120.38

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	528	ILE	Mainchain
1	H	536	GLY	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	H	5871	5887	5876	66	0
2	C	2933	2886	2893	18	0
3	C	27	0	12	2	0
4	C	1	0	0	0	0
All	All	8832	8773	8781	84	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (84) 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:H:367:PHE:CD2	1:H:376:ALA:HB1	1.60	1.35
1:H:368:LYS:HE2	1:H:378:MET:CA	1.51	1.17
1:H:368:LYS:CE	1:H:378:MET:N	2.04	1.13
1:H:374:ASP:CG	1:H:376:ALA:HA	1.74	1.10
1:H:367:PHE:CD2	1:H:376:ALA:CB	2.39	1.06
1:H:368:LYS:HE2	1:H:378:MET:N	1.19	1.05
1:H:374:ASP:CG	1:H:376:ALA:CA	2.30	1.04
1:H:374:ASP:CB	1:H:420:LYS:HB3	1.87	1.03
1:H:374:ASP:HB3	1:H:420:LYS:CB	1.90	1.00
1:H:368:LYS:HE3	1:H:377:SER:OG	1.57	0.99
1:H:374:ASP:OD1	1:H:376:ALA:N	2.03	0.91
1:H:373:THR:OG1	1:H:419:THR:HG21	1.51	0.87
1:H:374:ASP:OD2	1:H:376:ALA:HA	1.72	0.87
1:H:374:ASP:HB3	1:H:420:LYS:HB3	0.94	0.86
1:H:373:THR:HG1	1:H:419:THR:HG21	1.38	0.85
1:H:374:ASP:CG	1:H:376:ALA:N	2.35	0.84
1:H:374:ASP:O	1:H:419:THR:HA	1.80	0.82
1:H:368:LYS:CE	1:H:377:SER:OG	2.19	0.81
1:H:367:PHE:HD2	1:H:376:ALA:HB1	1.00	0.78
1:H:374:ASP:CG	1:H:376:ALA:H	1.92	0.77
1:H:367:PHE:HE2	1:H:403:LEU:CD2	2.00	0.74
1:H:374:ASP:O	1:H:419:THR:C	2.38	0.67
1:H:374:ASP:O	1:H:419:THR:CA	2.43	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:375:GLN:O	1:H:376:ALA:HB3	1.96	0.65
1:H:367:PHE:HE2	1:H:403:LEU:HD23	1.62	0.64
1:H:375:GLN:O	1:H:376:ALA:CB	2.46	0.64
1:H:374:ASP:C	1:H:376:ALA:N	2.51	0.64
1:H:374:ASP:O	1:H:420:LYS:N	2.33	0.62
1:H:374:ASP:OD2	1:H:376:ALA:CA	2.38	0.62
1:H:367:PHE:CE2	1:H:376:ALA:CB	2.83	0.61
1:H:367:PHE:CE2	1:H:403:LEU:HD23	2.36	0.60
1:H:603:ASP:N	1:H:604:PRO:HD3	2.17	0.59
1:H:132:ILE:HG13	1:H:132:ILE:O	2.01	0.58
1:H:373:THR:O	1:H:373:THR:HG22	2.02	0.58
1:H:368:LYS:HB2	1:H:374:ASP:OD2	2.04	0.57
1:H:367:PHE:HE2	1:H:403:LEU:HD21	1.69	0.55
1:H:374:ASP:CB	1:H:376:ALA:HA	2.35	0.54
2:C:356:TRP:O	2:C:356:TRP:HD1	1.91	0.54
1:H:370:GLU:HB2	1:H:374:ASP:CG	2.33	0.54
1:H:370:GLU:CB	1:H:374:ASP:CG	2.80	0.54
1:H:367:PHE:HD2	1:H:376:ALA:CB	1.92	0.54
1:H:367:PHE:CE2	1:H:376:ALA:HB3	2.44	0.52
2:C:104:LEU:C	2:C:104:LEU:HD23	2.35	0.52
2:C:127:PHE:O	2:C:128:ASN:C	2.52	0.52
1:H:391:MET:O	1:H:616:SER:OG	2.27	0.52
1:H:368:LYS:HE3	1:H:377:SER:HG	1.70	0.50
2:C:231:ALA:O	2:C:233:SER:N	2.45	0.49
1:H:374:ASP:OD2	1:H:376:ALA:C	2.55	0.48
2:C:40:HIS:O	2:C:41:GLN:HB2	2.14	0.48
1:H:368:LYS:N	1:H:377:SER:O	2.47	0.48
1:H:532:THR:O	1:H:533:ASN:HB2	2.14	0.47
2:C:171:LEU:O	2:C:173:HIS:N	2.45	0.47
2:C:302:GLY:CA	3:C:401:ADP:H4'	2.45	0.47
1:H:420:LYS:HD2	1:H:420:LYS:C	2.40	0.47
1:H:581:PHE:CD2	1:H:581:PHE:C	2.92	0.47
1:H:370:GLU:HB3	1:H:374:ASP:CG	2.40	0.47
1:H:737:LEU:O	1:H:738:ALA:CB	2.63	0.46
1:H:737:LEU:O	1:H:738:ALA:HB3	2.15	0.46
2:C:362:TYR:O	2:C:366:GLY:N	2.48	0.46
1:H:370:GLU:HB2	1:H:374:ASP:OD2	2.16	0.45
2:C:172:PRO:O	2:C:173:HIS:HB3	2.16	0.45
1:H:351:SER:OG	1:H:354:ARG:NH2	2.50	0.45
2:C:356:TRP:CD1	2:C:356:TRP:C	2.92	0.45
1:H:317:SER:O	1:H:318:ASN:HB2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:24:ASP:OD1	2:C:25:ASP:N	2.50	0.44
1:H:724:ARG:O	1:H:724:ARG:HG3	2.18	0.43
1:H:521:LEU:HD21	1:H:586:TYR:CD1	2.54	0.43
1:H:280:GLN:OE1	1:H:291:TYR:OH	2.37	0.43
2:C:242:LEU:O	2:C:244:ASP:N	2.52	0.43
1:H:150:PRO:HA	1:H:151:PRO:HD3	1.87	0.43
1:H:228:ASN:HB2	1:H:229:PRO:CD	2.49	0.43
2:C:257:CYS:HB2	2:C:258:PRO:HD3	2.00	0.43
1:H:80:PRO:HA	1:H:81:PRO:HD3	1.86	0.42
1:H:696:LEU:C	1:H:696:LEU:HD23	2.44	0.42
2:C:151:ILE:HG13	2:C:293:LEU:HB3	2.00	0.42
2:C:302:GLY:HA3	3:C:401:ADP:H4'	2.01	0.42
1:H:368:LYS:HB2	1:H:376:ALA:C	2.38	0.42
2:C:356:TRP:O	2:C:356:TRP:CD1	2.73	0.42
1:H:602:MET:O	1:H:603:ASP:CB	2.67	0.42
1:H:748:ASP:OD1	1:H:748:ASP:C	2.62	0.41
1:H:22:ASN:OD1	1:H:22:ASN:C	2.63	0.41
2:C:216:LEU:O	2:C:217:CYS:C	2.64	0.41
1:H:321:VAL:HB	1:H:322:PRO:HD3	2.04	0.40
2:C:1:ASP:O	2:C:3:ASP:N	2.54	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	H	721/798 (90%)	658 (91%)	46 (6%)	17 (2%)	4	27
2	C	371/377 (98%)	339 (91%)	22 (6%)	10 (3%)	4	25
All	All	1092/1175 (93%)	997 (91%)	68 (6%)	27 (2%)	6	26

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	146	ARG
1	H	318	ASN
1	H	376	ALA
1	H	531	PRO
1	H	738	ALA
2	C	128	ASN
2	C	173	HIS
2	C	232	SER
2	C	2	GLU
2	C	6	THR
2	C	137	GLN
2	C	234	SER
2	C	244	ASP
1	H	415	GLN
1	H	453	THR
1	H	657	ARG
1	H	724	ARG
2	C	100	GLU
1	H	603	ASP
1	H	574	GLN
1	H	457	GLY
1	H	604	PRO
1	H	656	PHE
1	H	533	ASN
1	H	743	PRO
2	C	197	GLY
1	H	132	ILE

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	H	641/704 (91%)	639 (100%)	2 (0%)	86	85
2	C	318/320 (99%)	318 (100%)	0	100	100
All	All	959/1024 (94%)	957 (100%)	2 (0%)	85	86

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	371	ARG
1	H	373	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	226	GLN
1	H	491	GLN
1	H	522	GLN
1	H	615	GLN
2	C	128	ASN
2	C	137	GLN
2	C	161	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ADP	C	401	-	28,29,29	1.10	2 (7%)	43,45,45	1.23	6 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ADP	C	401	-	-	3/16/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	401	ADP	C2-N1	3.09	1.39	1.33
3	C	401	ADP	C5-N7	-3.06	1.33	1.39

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	401	ADP	C5-C4-N3	2.51	130.17	126.72
3	C	401	ADP	C4-C5-N7	2.43	113.36	110.58
3	C	401	ADP	C4-N9-C8	-2.28	103.35	105.74
3	C	401	ADP	C2-N1-C6	2.27	122.46	118.73
3	C	401	ADP	N3-C2-N1	-2.10	125.40	128.58
3	C	401	ADP	O3B-PB-O2B	2.07	115.58	107.80

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	401	ADP	C3'-C4'-C5'-O5'
3	C	401	ADP	O4'-C4'-C5'-O5'
3	C	401	ADP	PA-O3A-PB-O3B

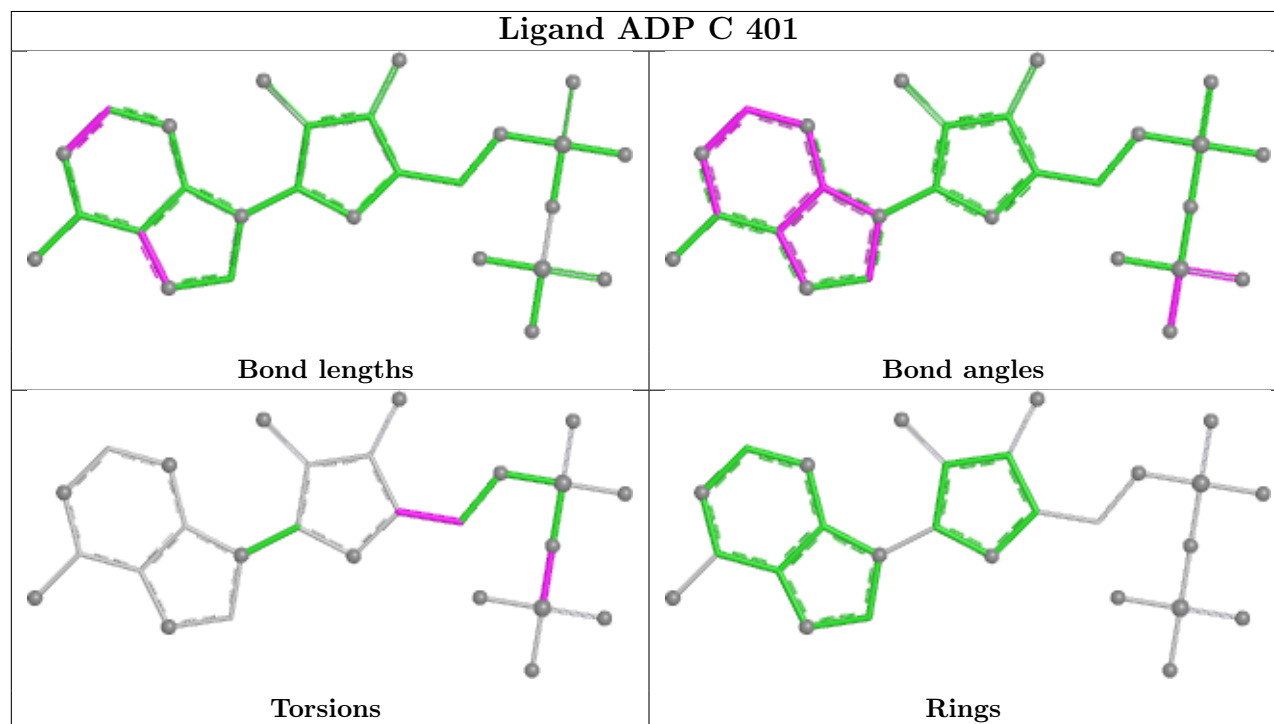
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	401	ADP	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	H	2
2	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	72:GLU	C	73:HIS	N	3.67
1	H	536:GLY	C	537:VAL	N	1.12

Continued on next page...

Continued from previous page...

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	528:ILE	C	529:GLU	N	1.01

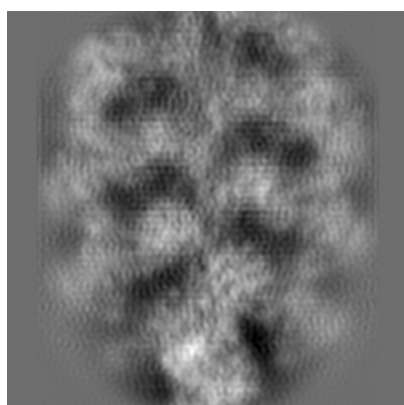
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-7100. These allow visual inspection of the internal detail of the map and identification of artifacts.

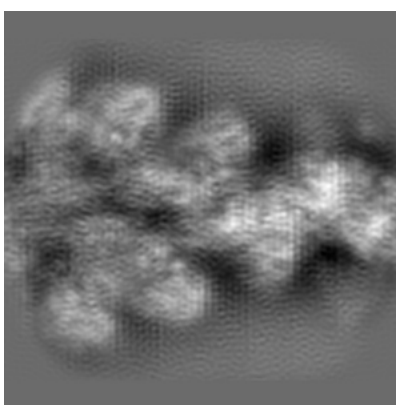
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

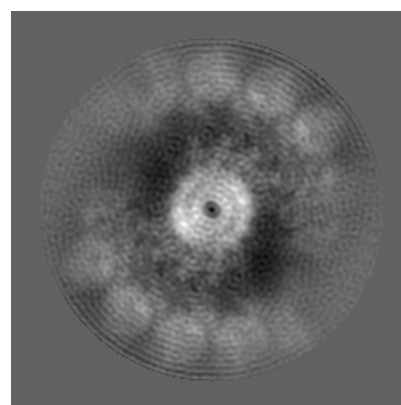
6.1.1 Primary map



X



Y

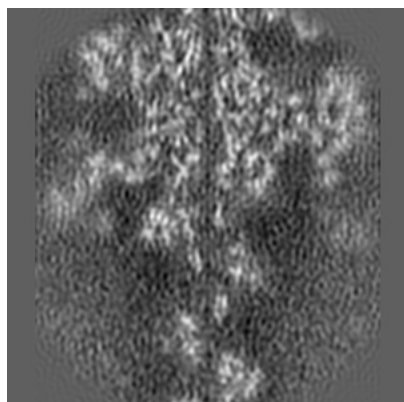


Z

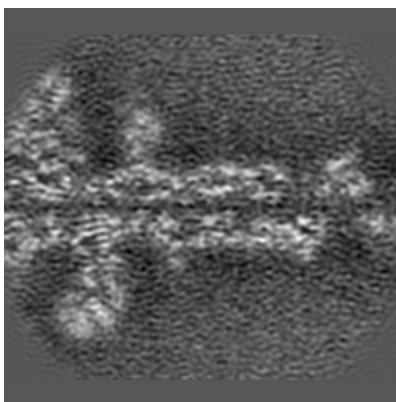
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

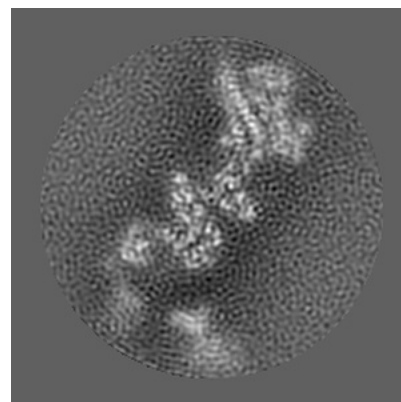
6.2.1 Primary map



X Index: 128



Y Index: 128

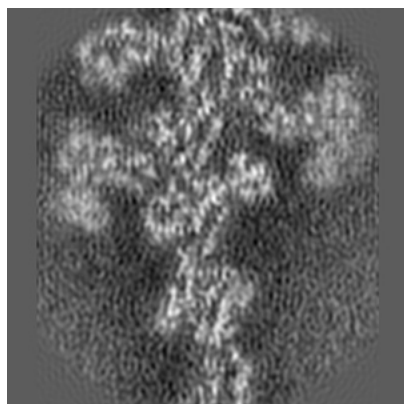


Z Index: 128

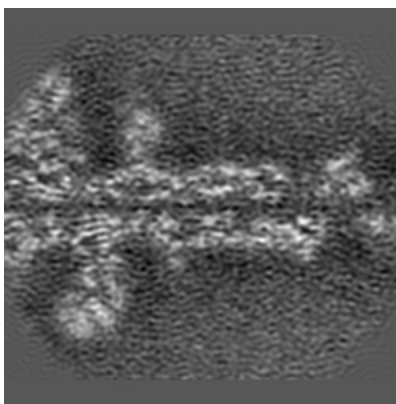
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

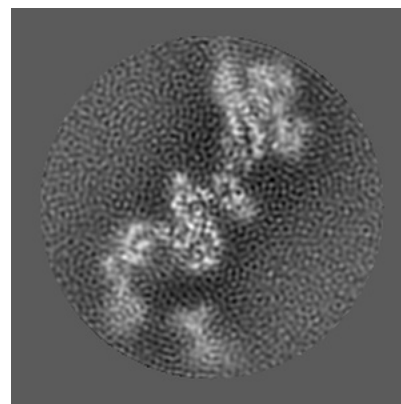
6.3.1 Primary map



X Index: 120



Y Index: 128

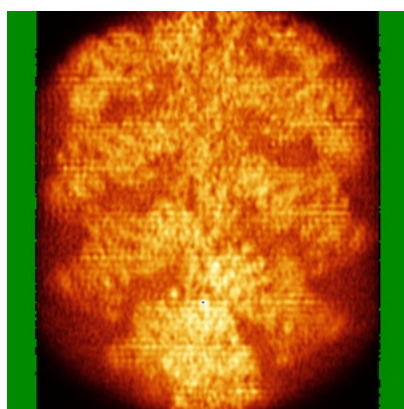


Z Index: 126

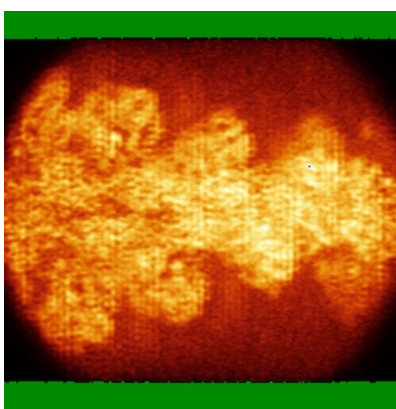
The images above show the largest variance slices of the map in three orthogonal directions.

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

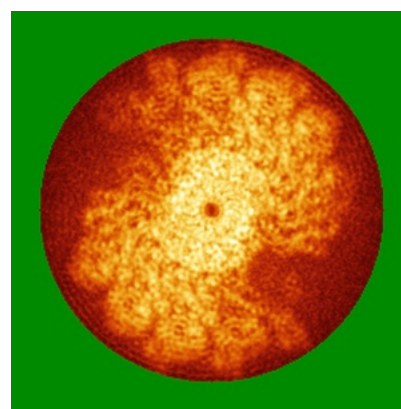
6.4.1 Primary map



X



Y

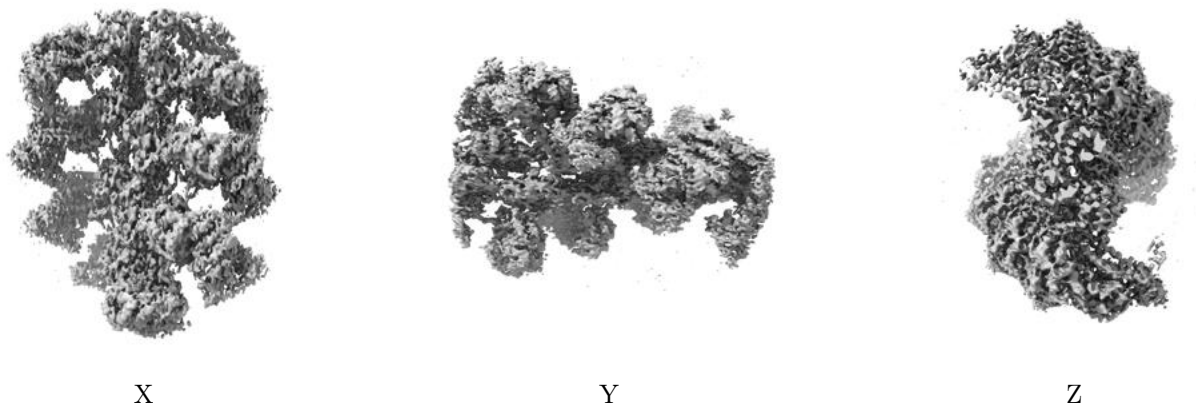


Z

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 0.016. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

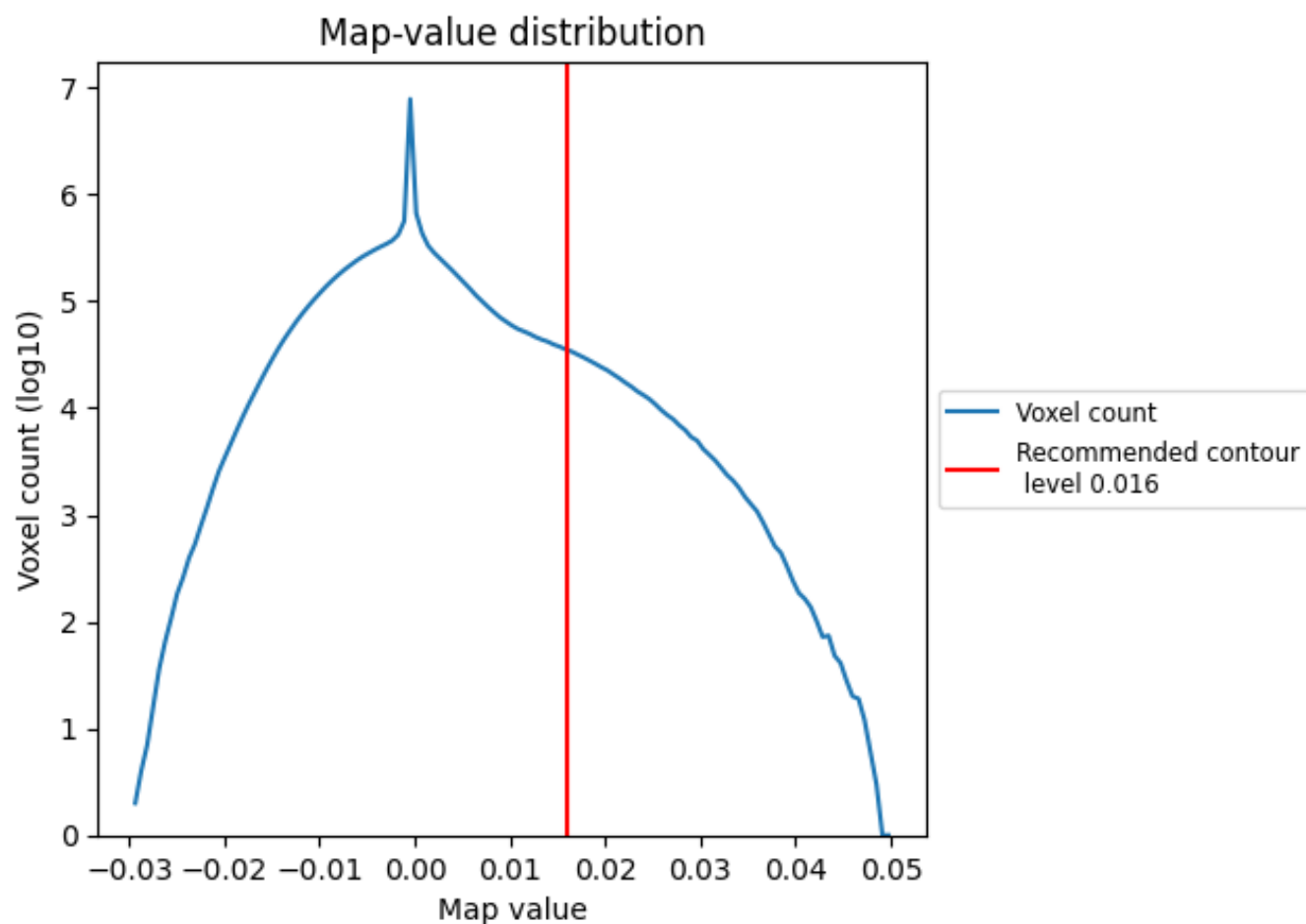
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

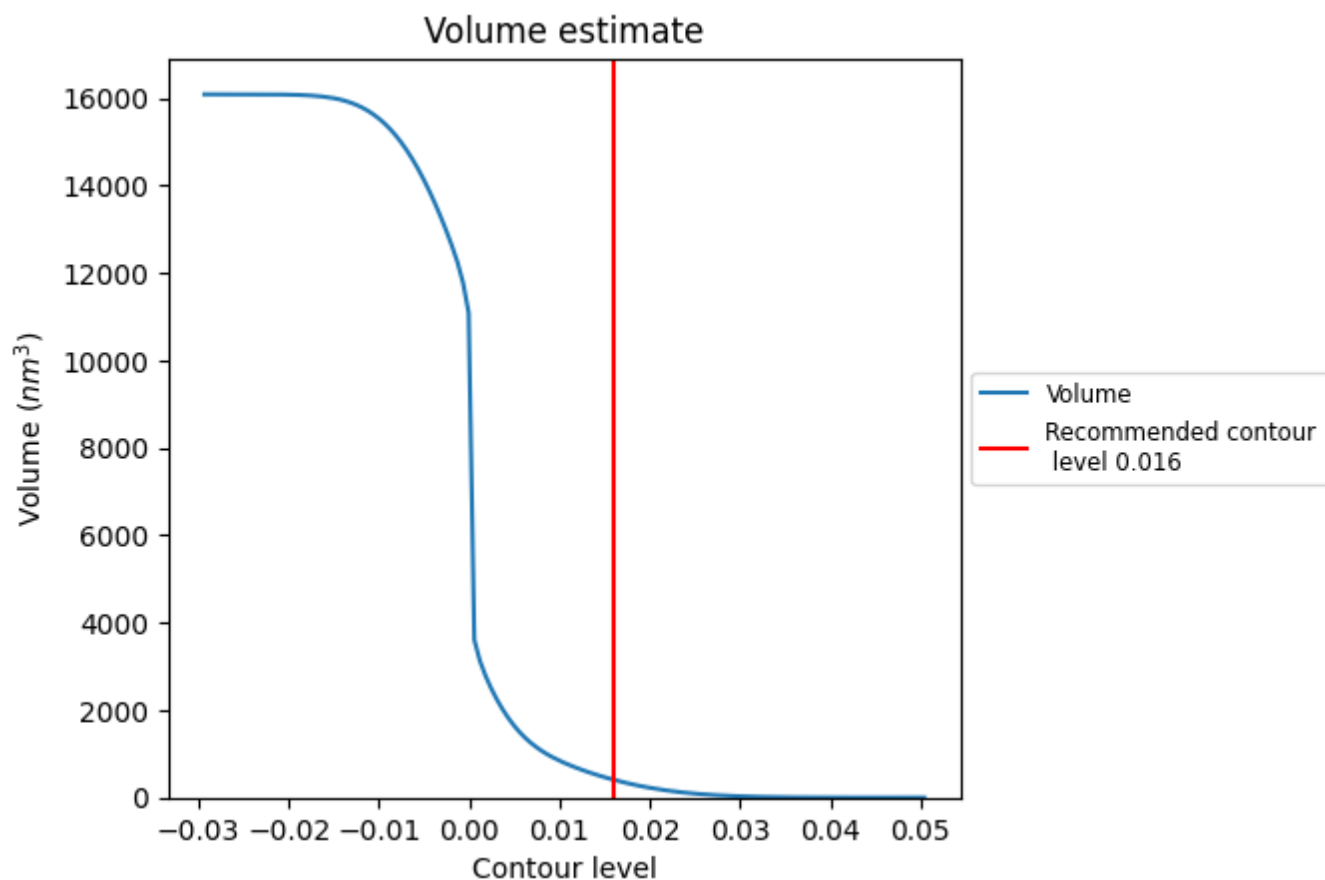
This section contains the results of statistical analysis of the map.

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.

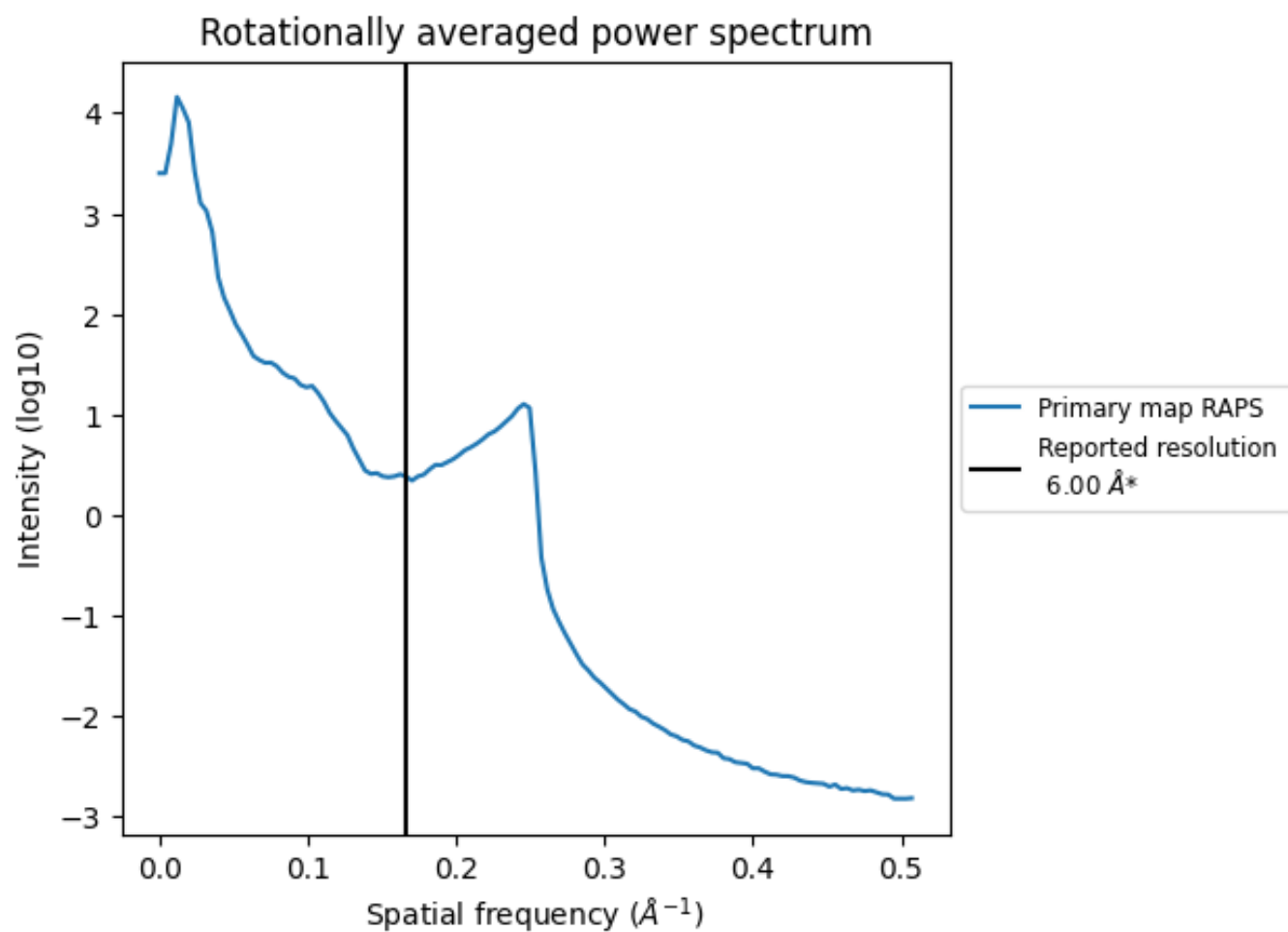
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 407 nm³; this corresponds to an approximate mass of 367 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 ⓘ

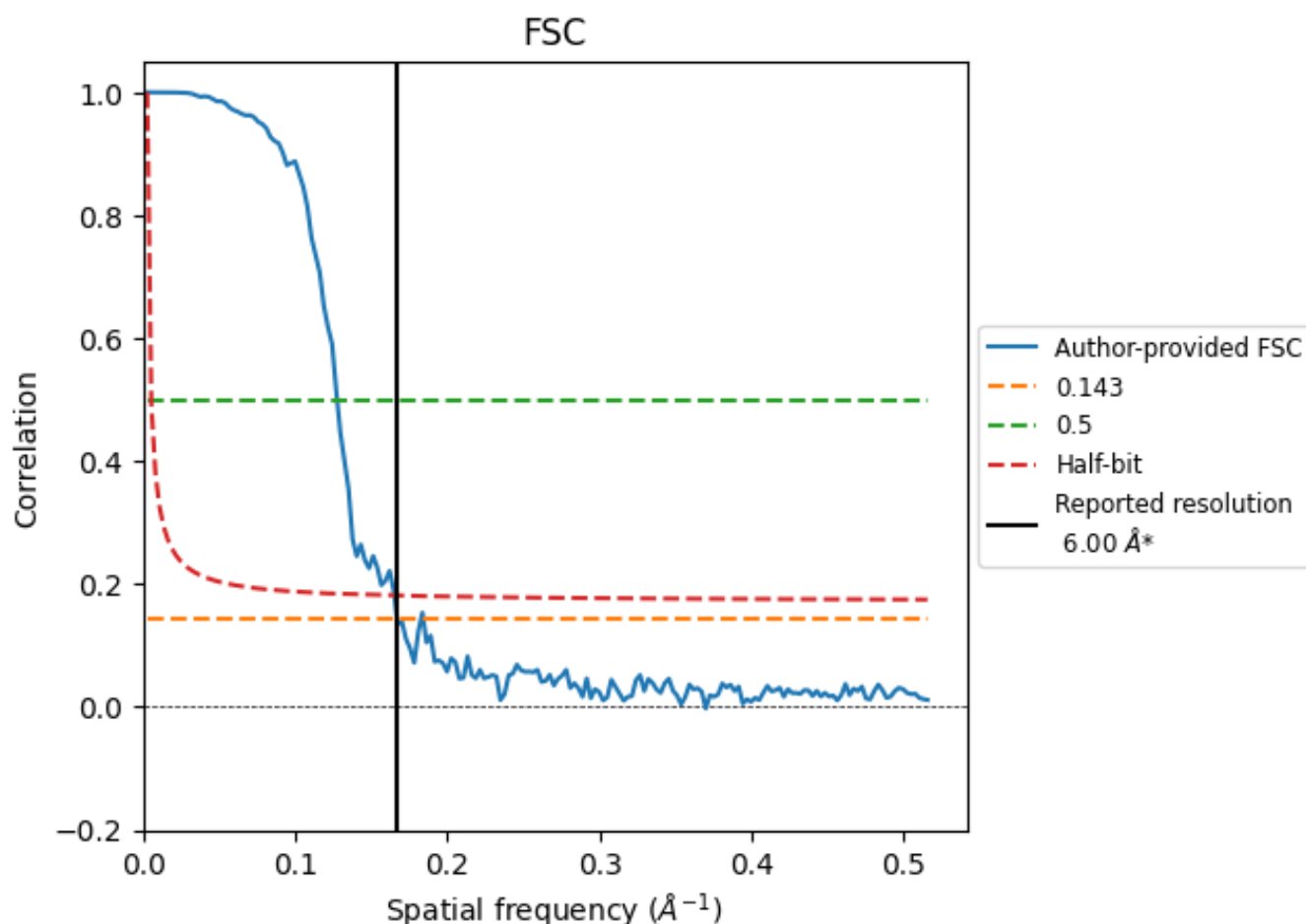


*Reported resolution corresponds to spatial frequency of 0.167 Å⁻¹

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 0.167 Å⁻¹

8.2 Resolution estimates [i](#)

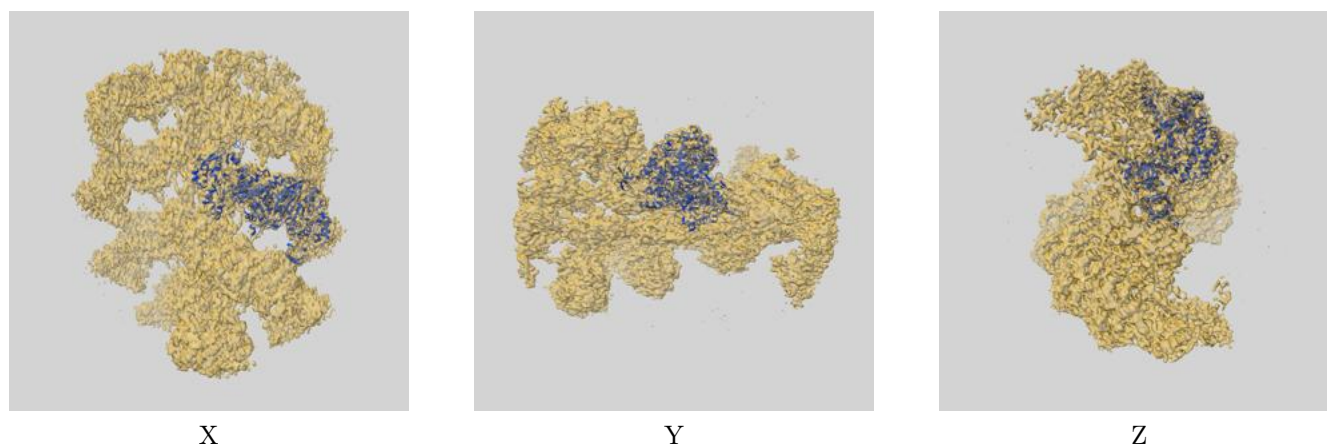
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	6.00	-	-
Author-provided FSC curve	5.99	7.84	6.06
Unmasked-calculated*	-	-	-

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

9 Map-model fit [i](#)

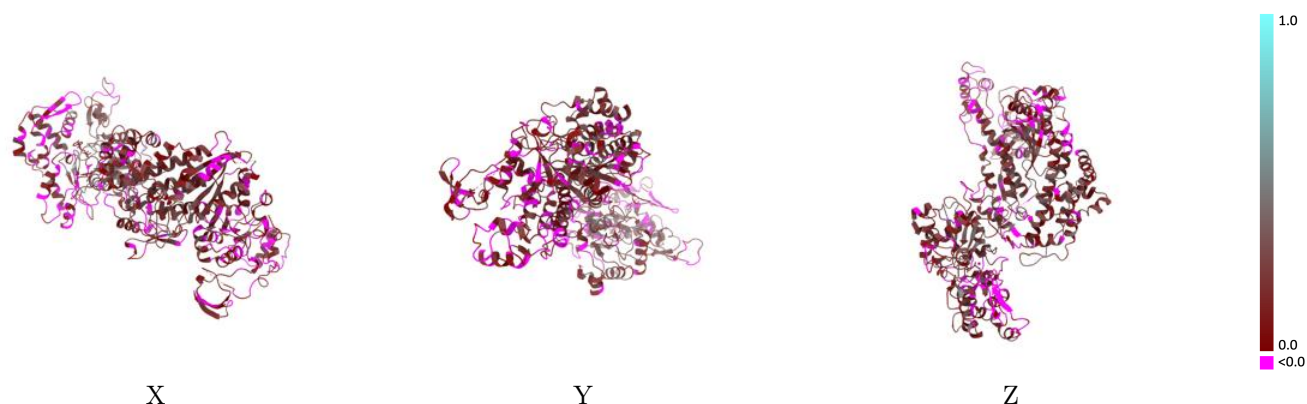
This section contains information regarding the fit between EMDB map EMD-7100 and PDB model 6BIH. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



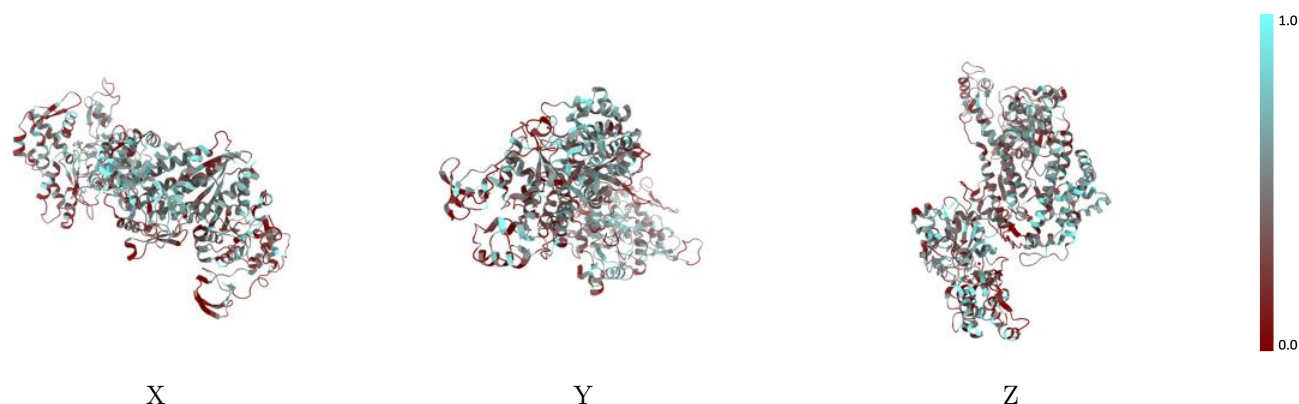
The images above show the 3D surface view of the map at the recommended contour level 0.016 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



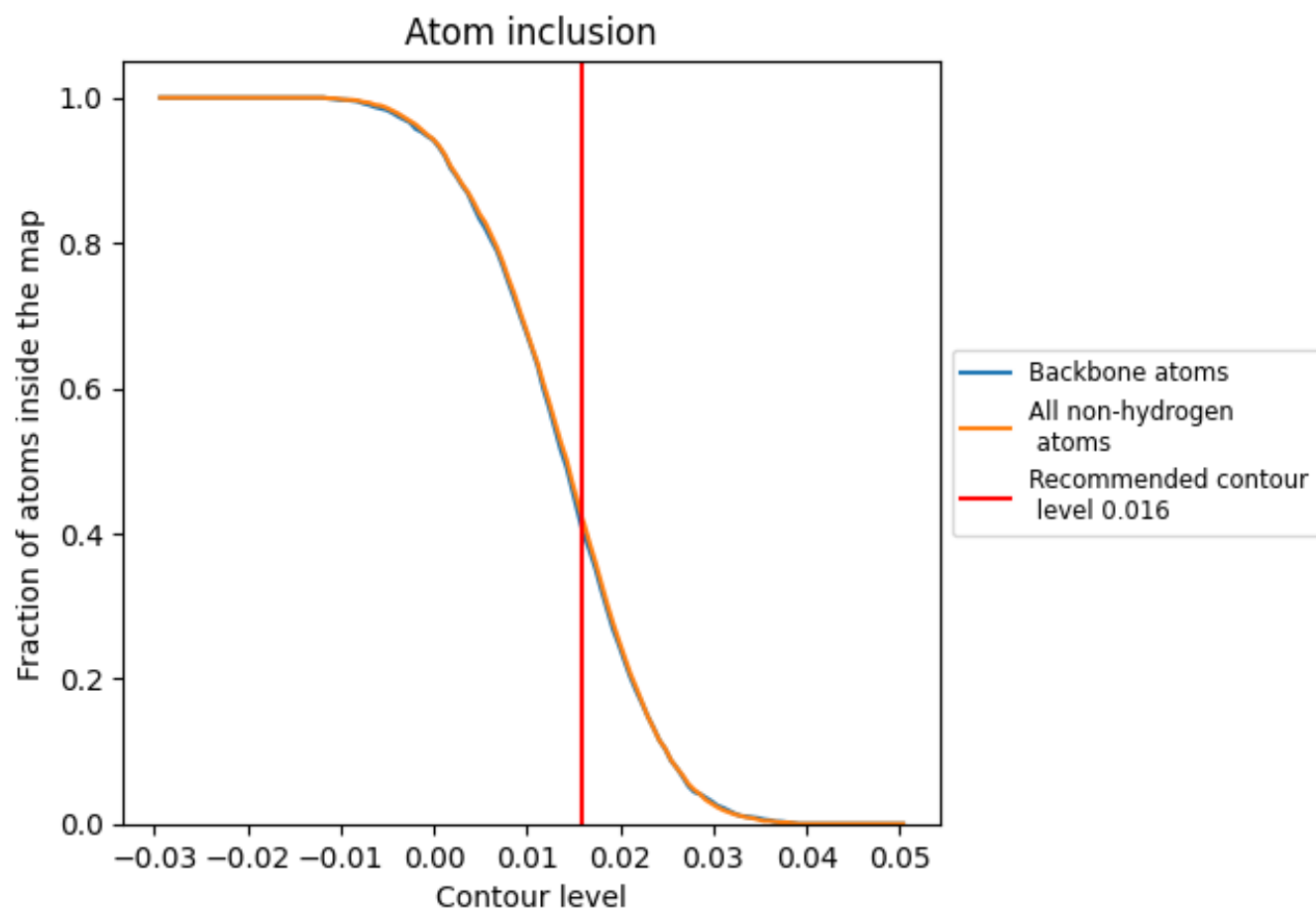
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.016).

9.4 Atom inclusion [i](#)



At the recommended contour level, 40% of all backbone atoms, 42% 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 (0.016) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.4180	0.1330
C	0.4100	0.1260
H	0.4260	0.1370

