Full wwPDB EM Validation Report (i)

Dec 13, 2021 – 12:10 pm GMT

PDB ID	:	7PM0
EMDB ID	:	EMD-13508
Title	:	Cryo-EM structure of the actomyosin-V complex in the rigor state (central 1er, young JASP-stabilized F-actin, class 1)
Authors	:	Pospich, S.; Sweeney, H.L.; Houdusse, A.; Raunser, S.
Deposited on	:	2021-09-01
Resolution	:	3.60 Å(reported)
Based on initial model	:	7PLY

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (i)) were used in the production of this report:

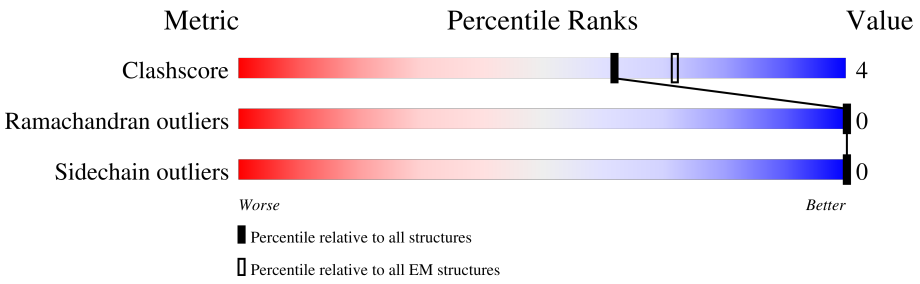
EMDB validation analysis	:	0.0.0.dev97
Mogul	:	1.8.4, CSD as541be (2020)
MolProbity	:	4.02b-467
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.24

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.60 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion < 40%). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	B	151	18% 73% 19% 8%
2	A	792	85% 10% 5%
3	C	377	86% 12%

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 10169 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 Myosin light chain 6B.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	B	139	1059	673	174	206	6	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	initiating methionine	UNP P14649

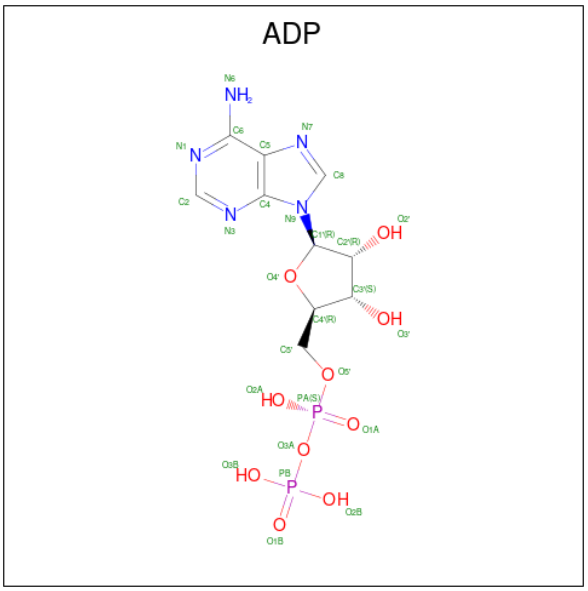
- Molecule 2 is a protein called Unconventional myosin-Va.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	A	751	6130	3926	1051	1118	35	0	0

- Molecule 3 is a protein called Actin, alpha skeletal muscle.

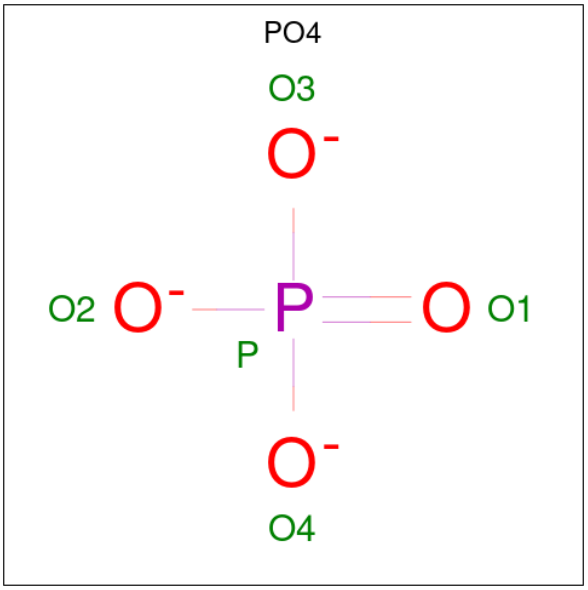
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	371	2900	1837	489	553	21	0	0

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).

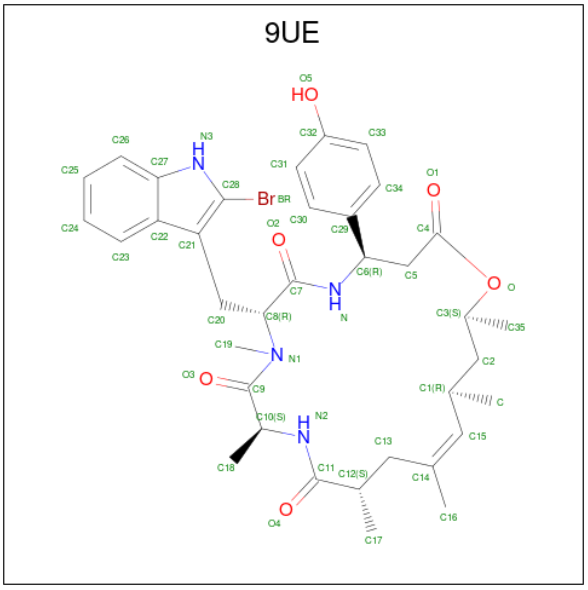


Mol	Chain	Residues	Atoms			AltConf
			Total	O	P	
5	C	1	5	4	1	0

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 7 is Jasplakinolide (three-letter code: 9UE) (formula: $C_{36}H_{45}BrN_4O_6$).



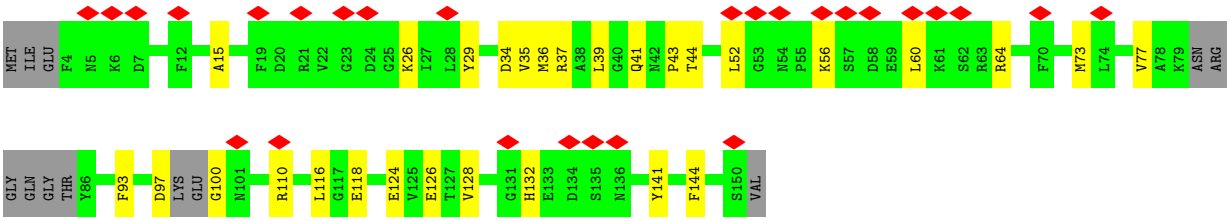
Mol	Chain	Residues	Atoms					AltConf
			Total	Br	C	N	O	
7	C	1	47	1	36	4	6	0

3 Residue-property plots (i)

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

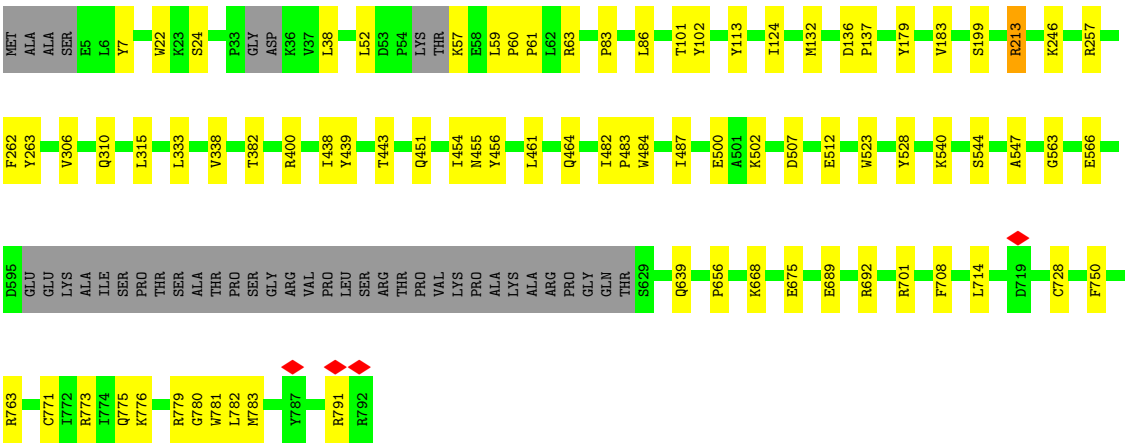
- Molecule 1: Myosin light chain 6B

Chain B:



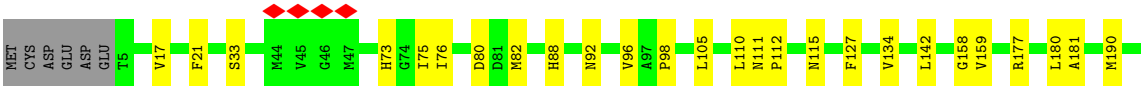
- Molecule 2: Unconventional myosin-Va

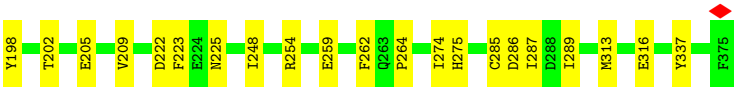
Chain A:



- Molecule 3: Actin, alpha skeletal muscle

Chain C:




Y198
T202
E205
V209
D222
F223
E224
N225
I248
R254
E259
F262
Q263
P264
I274
H275
C285
D286
I287
D288
I289
M313
E316
Y337
F375

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	POINT, Not provided	
Number of segments used	110797	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 ($e^-/Å^2$)	80	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.036	Depositor
Minimum map value	-0.015	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.006	Depositor
Map size (Å)	352.0, 352.0, 352.0	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	Depositor

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: HIC, PO4, ADP, MG, 9UE

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	B	0.26	0/1075	0.48	0/1448
2	A	0.32	0/6265	0.51	0/8442
3	C	0.35	0/2950	0.52	0/3994
All	All	0.32	0/10290	0.51	0/13884

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
2	A	0	1
3	C	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	213	ARG	Sidechain
3	C	254	ARG	Sidechain

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	B	1059	0	998	22	0
2	A	6130	0	6123	50	0
3	C	2900	0	2872	28	0
4	C	27	0	12	0	0
5	C	5	0	0	0	0
6	C	1	0	0	0	0
7	C	47	0	0	1	0
All	All	10169	0	10005	90	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 4.

All (90) 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:B:34:ASP:HA	2:A:783:MET:SD	2.28	0.74
3:C:76:ILE:HD13	3:C:82:MET:HE1	1.69	0.74
1:B:37:ARG:NH2	1:B:43:PRO:O	2.21	0.73
1:B:124:GLU:HG2	2:A:782:LEU:HD22	1.70	0.72
3:C:285:CYS:HB3	3:C:289:ILE:HD11	1.73	0.70
3:C:112:PRO:O	3:C:115:ASN:ND2	2.27	0.68
2:A:456:TYR:OH	2:A:500:GLU:OE1	2.09	0.66
1:B:118:GLU:O	2:A:779:ARG:NH2	2.21	0.63
3:C:110:LEU:O	3:C:111:ASN:ND2	2.33	0.61
2:A:438:ILE:HD13	2:A:461:LEU:HD11	1.82	0.61
1:B:37:ARG:HD2	2:A:783:MET:SD	2.41	0.61
2:A:512:GLU:HG2	2:A:523:TRP:HB2	1.84	0.60
2:A:701:ARG:HG2	2:A:750:PHE:CD1	2.38	0.59
2:A:540:LYS:NZ	2:A:544:SER:O	2.37	0.57
3:C:88:HIS:CG	3:C:92:ASN:HD21	2.23	0.56
2:A:771:CYS:SG	2:A:775:GLN:NE2	2.74	0.56
2:A:213:ARG:NH1	2:A:443:THR:O	2.39	0.55
1:B:93:PHE:CD2	1:B:141:TYR:HB2	2.43	0.54
3:C:275:HIS:NE2	3:C:316:GLU:OE1	2.41	0.53
2:A:102:TYR:CE2	2:A:137:PRO:HB3	2.44	0.53
3:C:21:PHE:HZ	3:C:96:VAL:HG21	1.73	0.52
3:C:190:MET:HG2	3:C:209:VAL:HG11	1.90	0.52
3:C:17:VAL:HG23	3:C:33:SER:HB2	1.92	0.52
2:A:528:TYR:OH	2:A:547:ALA:O	2.25	0.52
2:A:246:LYS:NZ	2:A:639:GLN:OE1	2.37	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:306:VAL:O	2:A:310:GLN:HG2	2.10	0.51
1:B:26:LYS:O	1:B:64:ARG:NH2	2.44	0.51
2:A:484:TRP:O	2:A:487:ILE:HG22	2.11	0.51
1:B:29:TYR:HB2	1:B:60:LEU:HA	1.94	0.50
2:A:102:TYR:OH	2:A:132:MET:SD	2.63	0.50
1:B:36:MET:SD	1:B:77:VAL:HG21	2.52	0.50
2:A:199:SER:HA	2:A:315:LEU:HD11	1.92	0.49
2:A:113:TYR:CE2	2:A:668:LYS:HA	2.47	0.49
1:B:37:ARG:HA	1:B:41:GLN:O	2.13	0.49
3:C:262:PHE:HD1	3:C:275:HIS:NE2	2.12	0.48
1:B:110:ARG:HH12	1:B:126:GLU:HA	1.79	0.47
2:A:689:GLU:CD	2:A:692:ARG:HE	2.18	0.47
1:B:37:ARG:HD3	2:A:780:GLY:HA2	1.95	0.47
2:A:482:ILE:HG23	2:A:483:PRO:HD2	1.95	0.46
2:A:38:LEU:HD11	2:A:52:LEU:HD21	1.97	0.46
2:A:179:TYR:CZ	2:A:183:VAL:HG11	2.50	0.46
3:C:286:ASP:OD1	3:C:287:ILE:N	2.47	0.46
2:A:464:GLN:HE22	2:A:656:PRO:HG2	1.80	0.46
2:A:22:TRP:CD2	2:A:83:PRO:HG3	2.51	0.45
3:C:202:THR:HG22	3:C:205:GLU:HG3	1.98	0.45
1:B:128:VAL:HG23	2:A:781:TRP:CD1	2.51	0.45
1:B:35:VAL:O	1:B:39:LEU:HD13	2.16	0.45
2:A:24:SER:OG	2:A:63:ARG:NH1	2.50	0.45
2:A:86:LEU:HD21	2:A:675:GLU:HB3	1.98	0.45
2:A:124:ILE:HG23	2:A:179:TYR:CD1	2.51	0.45
2:A:439:TYR:CD2	2:A:454:ILE:HG23	2.51	0.44
1:B:29:TYR:HD1	1:B:52:LEU:HD22	1.83	0.44
1:B:116:LEU:O	2:A:775:GLN:NE2	2.38	0.44
3:C:158:GLY:O	3:C:181:ALA:HB2	2.18	0.44
1:B:97:ASP:OD1	1:B:100:GLY:HA2	2.18	0.43
2:A:773:ARG:O	2:A:776:LYS:HG2	2.18	0.43
2:A:7:TYR:HE2	2:A:59:LEU:HD13	1.83	0.43
3:C:80:ASP:OD2	3:C:80:ASP:N	2.49	0.43
2:A:338:VAL:O	2:A:400:ARG:NH1	2.52	0.43
2:A:257:ARG:HB3	2:A:263:TYR:CE2	2.55	0.42
3:C:222:ASP:OD2	3:C:225:ASN:ND2	2.47	0.42
2:A:714:LEU:HD21	2:A:763:ARG:HA	2.02	0.42
3:C:198:TYR:CE2	3:C:248:ILE:HB	2.54	0.42
2:A:52:LEU:HD12	2:A:57:LYS:HE2	2.01	0.42
2:A:382:THR:HG21	3:C:337:TYR:CE2	2.54	0.42
2:A:563:GLY:HA2	2:A:566:GLU:OE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:159:VAL:HG22	3:C:177:ARG:NH2	2.34	0.42
3:C:223:PHE:CE1	3:C:259:GLU:HG2	2.55	0.42
2:A:136:ASP:OD2	2:A:136:ASP:N	2.53	0.42
3:C:105:LEU:O	3:C:134:VAL:HA	2.20	0.42
1:B:56:LYS:HA	1:B:56:LYS:HD3	1.91	0.41
3:C:202:THR:CG2	3:C:205:GLU:HG3	2.49	0.41
2:A:502:LYS:HA	2:A:507:ASP:OD2	2.21	0.41
1:B:52:LEU:HD23	1:B:73:MET:HE1	2.02	0.41
2:A:262:PHE:CD1	2:A:333:LEU:HD21	2.55	0.41
1:B:15:ALA:HB2	2:A:791:ARG:NH2	2.35	0.41
3:C:180:LEU:HD11	3:C:264:PRO:HG3	2.02	0.41
3:C:142:LEU:HD23	3:C:142:LEU:HA	1.87	0.41
3:C:75:ILE:HD11	7:C:404:9UE:C22	2.50	0.41
2:A:701:ARG:HG2	2:A:750:PHE:HD1	1.86	0.41
3:C:262:PHE:HD1	3:C:275:HIS:CE1	2.39	0.41
1:B:44:THR:HG22	2:A:776:LYS:HD3	2.03	0.41
1:B:132:HIS:CD2	1:B:144:PHE:HD1	2.39	0.41
2:A:451:GLN:O	2:A:455:ASN:ND2	2.54	0.41
2:A:708:PHE:HB2	2:A:728:CYS:SG	2.60	0.41
3:C:76:ILE:HG21	3:C:82:MET:HE2	2.03	0.40
2:A:101:THR:HG22	2:A:102:TYR:N	2.36	0.40
2:A:60:PRO:HA	2:A:61:PRO:HD3	1.97	0.40
3:C:98:PRO:HG3	3:C:127:PHE:CD1	2.57	0.40
3:C:274:ILE:HD12	3:C:313:MET:CE	2.52	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	B	133/151 (88%)	131 (98%)	2 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	743/792 (94%)	743 (100%)	0	0	100	100
3	C	368/377 (98%)	368 (100%)	0	0	100	100
All	All	1244/1320 (94%)	1242 (100%)	2 (0%)	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	B	108/129 (84%)	108 (100%)	0	100	100
2	A	663/695 (95%)	663 (100%)	0	100	100
3	C	313/319 (98%)	313 (100%)	0	100	100
All	All	1084/1143 (95%)	1084 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	75	GLN
2	A	200	ASN
2	A	264	GLN
2	A	497	ASN
2	A	632	HIS
2	A	638	HIS
2	A	642	ASN
2	A	726	GLN
3	C	41	GLN
3	C	111	ASN
3	C	162	ASN
3	C	296	ASN
3	C	371	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)

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
3	HIC	C	73	3	8,11,12	1.65	2 (25%)	6,14,16	0.71	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
3	HIC	C	73	3	-	1/5/6/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	73	HIC	CD2-CG	3.51	1.41	1.36
3	C	73	HIC	CZ-NE2	-2.06	1.42	1.48

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	73	HIC	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	$\#\|Z\| > 2$
5	PO4	C	402	6	4,4,4	0.96	0	6,6,6	0.44	0
7	9UE	C	404	-	47,50,50	0.62	1 (2%)	60,71,71	1.66	14 (23%)
4	ADP	C	401	6	24,29,29	0.95	1 (4%)	29,45,45	1.38	4 (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
7	9UE	C	404	-	-	15/54/56/56	0/3/4/4
4	ADP	C	401	6	-	0/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	401	ADP	C5-C4	2.21	1.46	1.40
7	C	404	9UE	BR-C28	-2.01	1.87	1.90

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	C	404	9UE	C5-C6-N	4.64	117.66	108.69
7	C	404	9UE	C6-C5-C4	3.95	121.21	112.80
4	C	401	ADP	PA-O3A-PB	-3.76	119.91	132.83

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	C	404	9UE	C6-N-C7	-3.61	117.78	123.19
7	C	404	9UE	C20-C21-C28	3.47	131.29	127.15
7	C	404	9UE	C1-C15-C14	3.29	137.53	127.66
7	C	404	9UE	C12-C11-N2	3.18	119.83	116.42
4	C	401	ADP	N3-C2-N1	-3.15	123.75	128.68
7	C	404	9UE	C3-O-C4	-2.92	114.12	117.88
7	C	404	9UE	C20-C8-C7	-2.90	103.84	111.05
4	C	401	ADP	C4-C5-N7	-2.84	106.44	109.40
7	C	404	9UE	C7-C8-N1	2.73	117.50	110.01
7	C	404	9UE	O4-C11-C12	-2.72	116.14	121.33
7	C	404	9UE	C17-C12-C13	2.59	115.59	111.39
7	C	404	9UE	C10-C9-N1	2.47	123.38	118.45
7	C	404	9UE	C10-N2-C11	2.46	126.61	121.29
7	C	404	9UE	C-C1-C15	2.22	115.43	109.99
4	C	401	ADP	O4'-C1'-C2'	-2.19	103.72	106.93

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	C	404	9UE	C-C1-C15-C14
7	C	404	9UE	O1-C4-O-C3
7	C	404	9UE	C5-C4-O-C3
7	C	404	9UE	N2-C11-C12-C13
7	C	404	9UE	O4-C11-C12-C13
7	C	404	9UE	C34-C29-C6-N
7	C	404	9UE	C30-C29-C6-N
7	C	404	9UE	N2-C11-C12-C17
7	C	404	9UE	O4-C11-C12-C17
7	C	404	9UE	C17-C12-C13-C14
7	C	404	9UE	C11-C12-C13-C14
7	C	404	9UE	C5-C6-N-C7
7	C	404	9UE	C9-C10-N2-C11
7	C	404	9UE	C15-C1-C2-C3
7	C	404	9UE	O1-C4-C5-C6

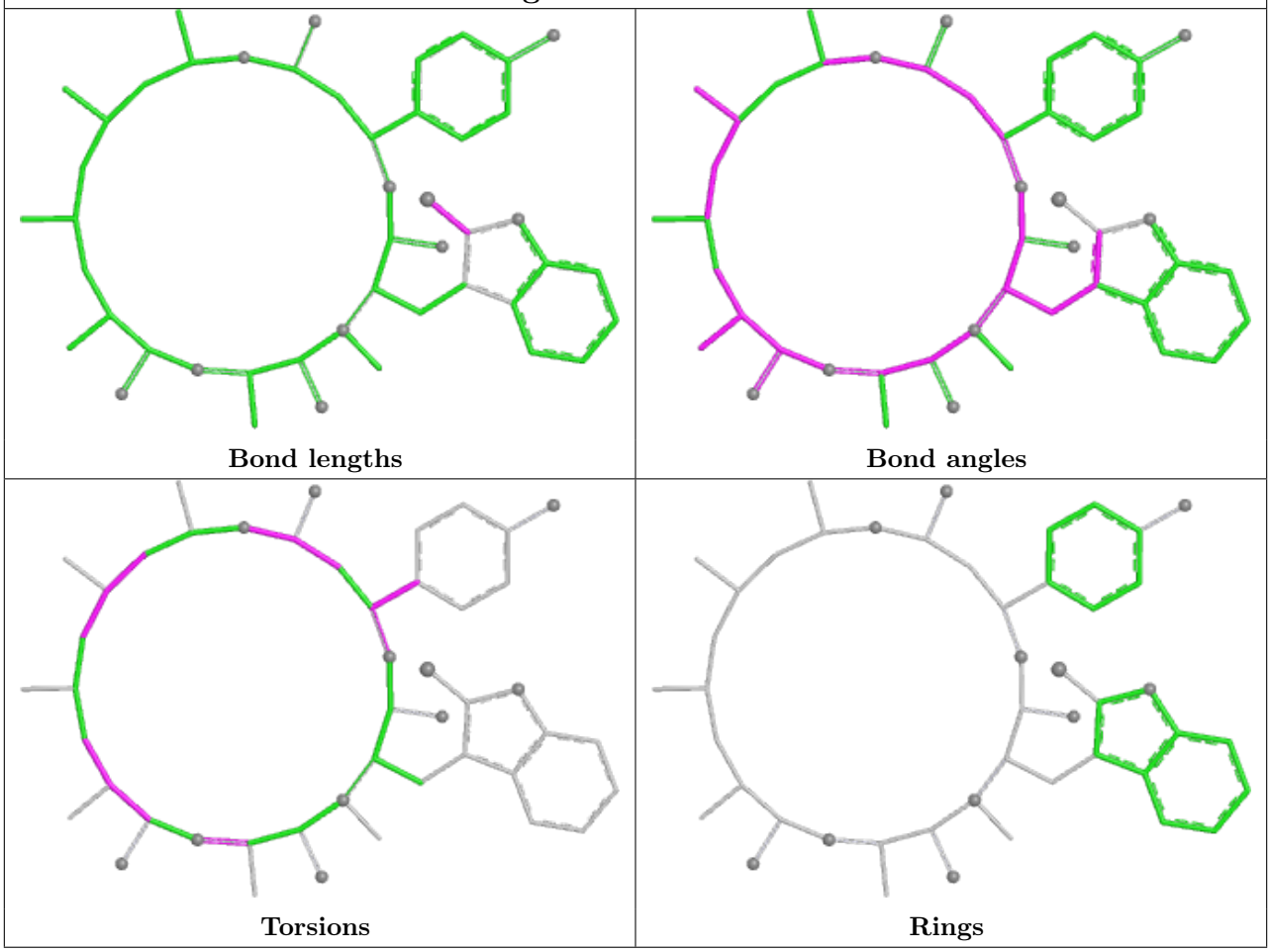
There are no ring outliers.

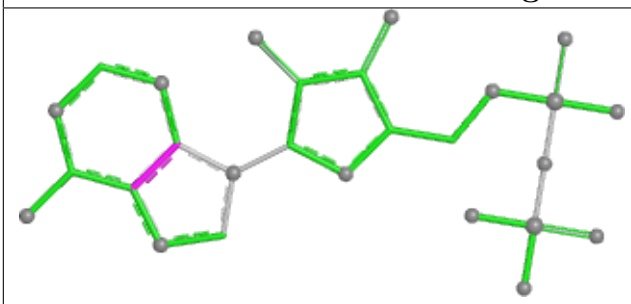
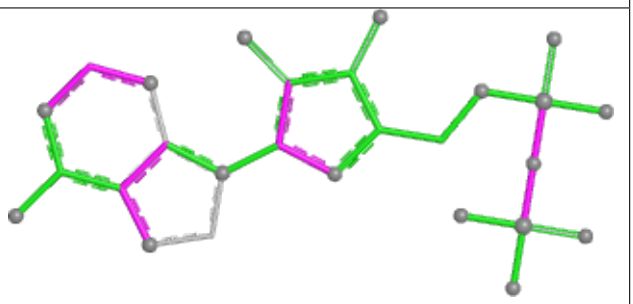
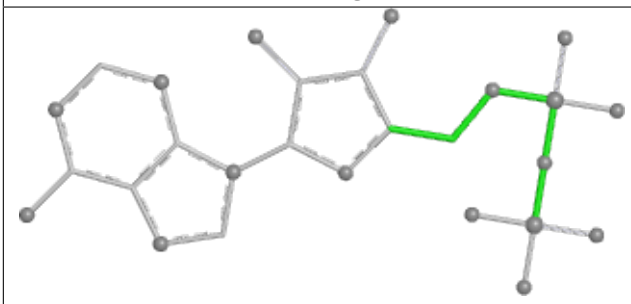
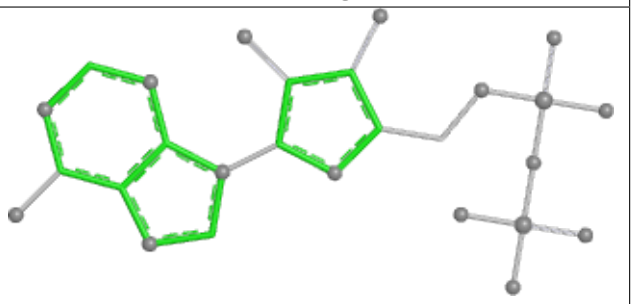
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	404	9UE	1	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 then 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.

Ligand 9UE C 404



Ligand ADP C 401	
 Bond lengths	 Bond angles
 Torsions	 Rings

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.

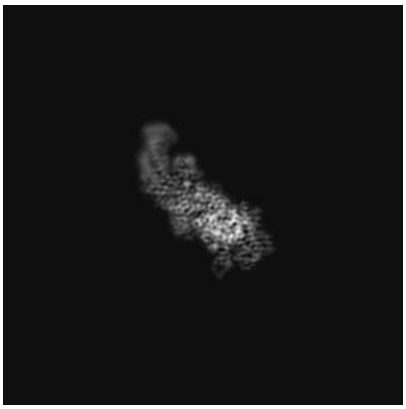
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-13508. These allow visual inspection of the internal detail of the map and identification of artifacts.

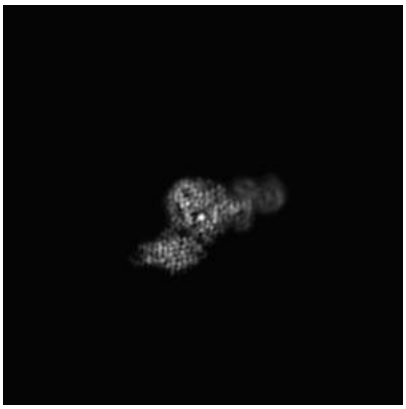
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

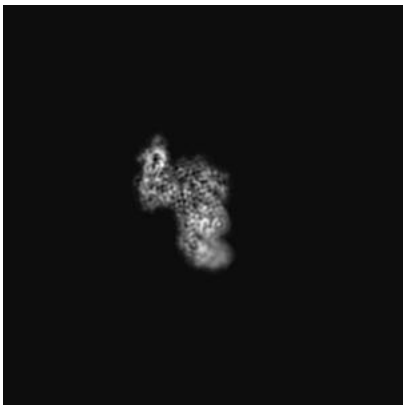
6.1.1 Primary map



X

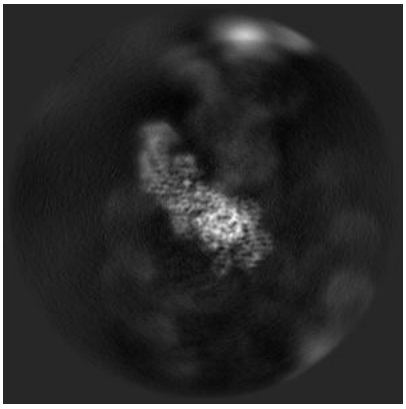


Y

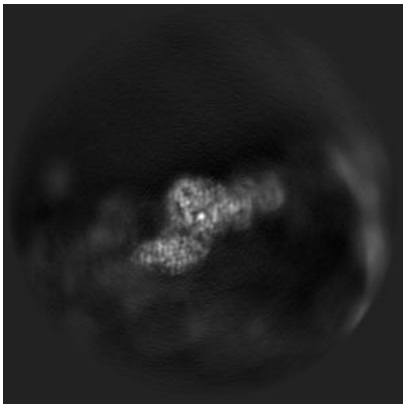


Z

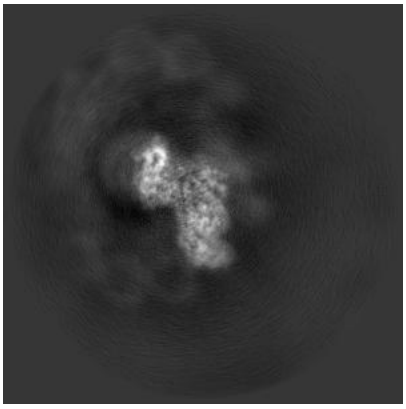
6.1.2 Raw map



X



Y

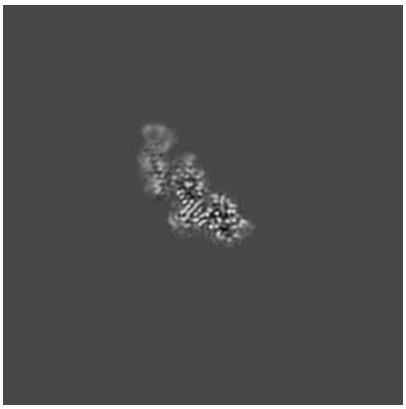


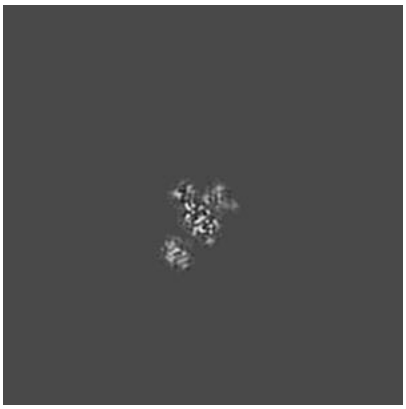
Z

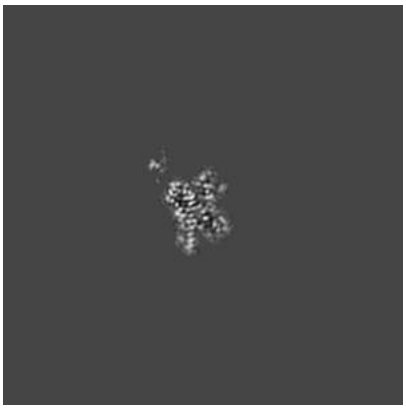
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

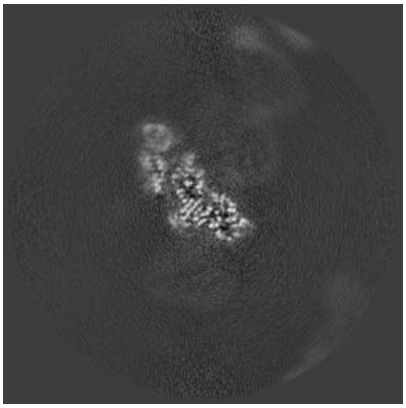
6.2.1 Primary map

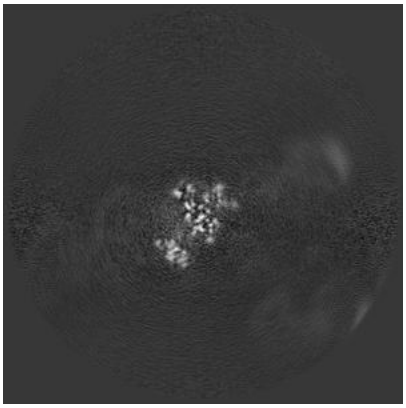

X Index: 160

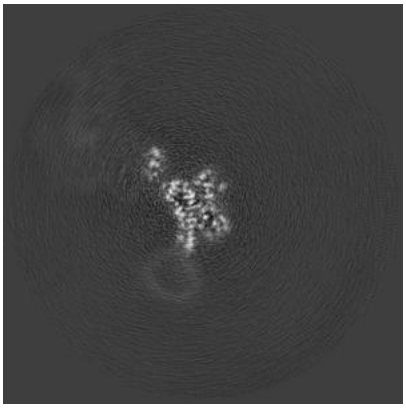

Y Index: 160


Z Index: 160

6.2.2 Raw map


X Index: 160

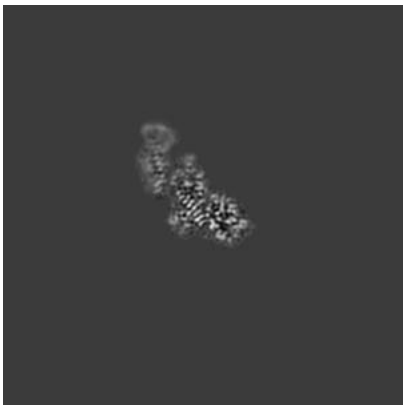

Y Index: 160

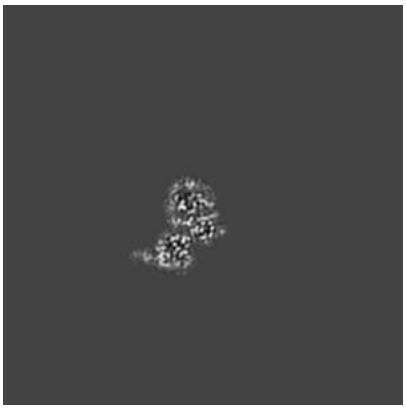

Z Index: 160

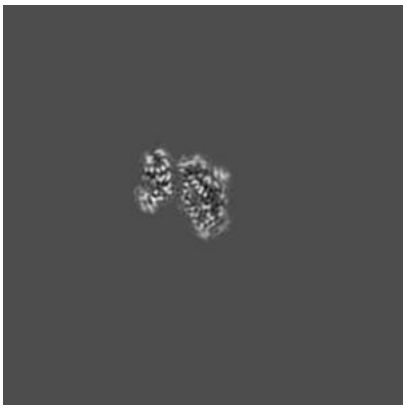
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

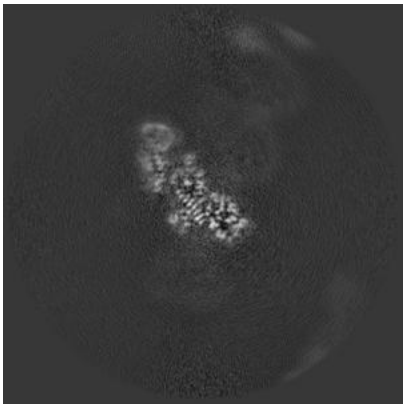
6.3.1 Primary map

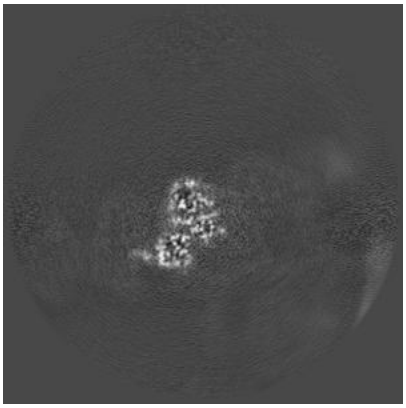

X Index: 161

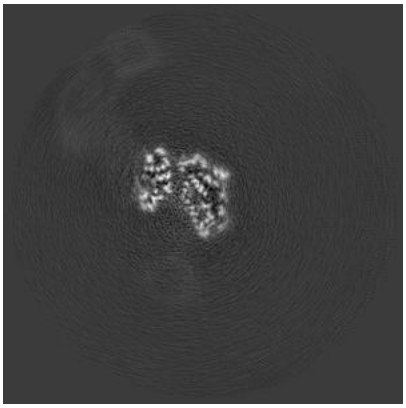

Y Index: 170


Z Index: 145

6.3.2 Raw map


X Index: 161


Y Index: 170


Z Index: 145

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

6.4 Orthogonal surface views (i)

6.4.1 Primary map

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.006. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 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.

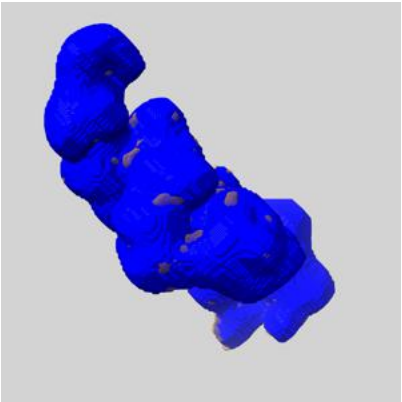
6.5 Mask visualisation (i)

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

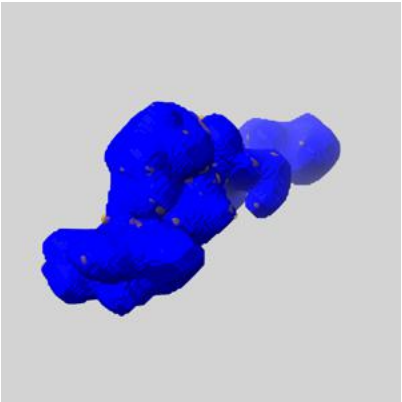
A mask typically either:

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

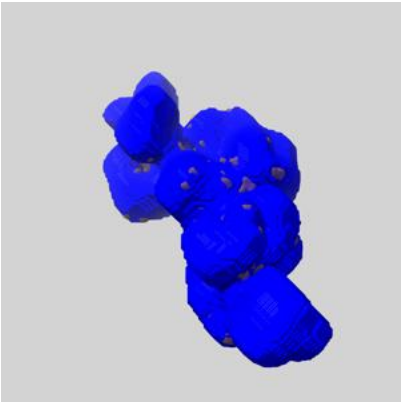
6.5.1 emd_13508_msk_1.map (i)



X



Y

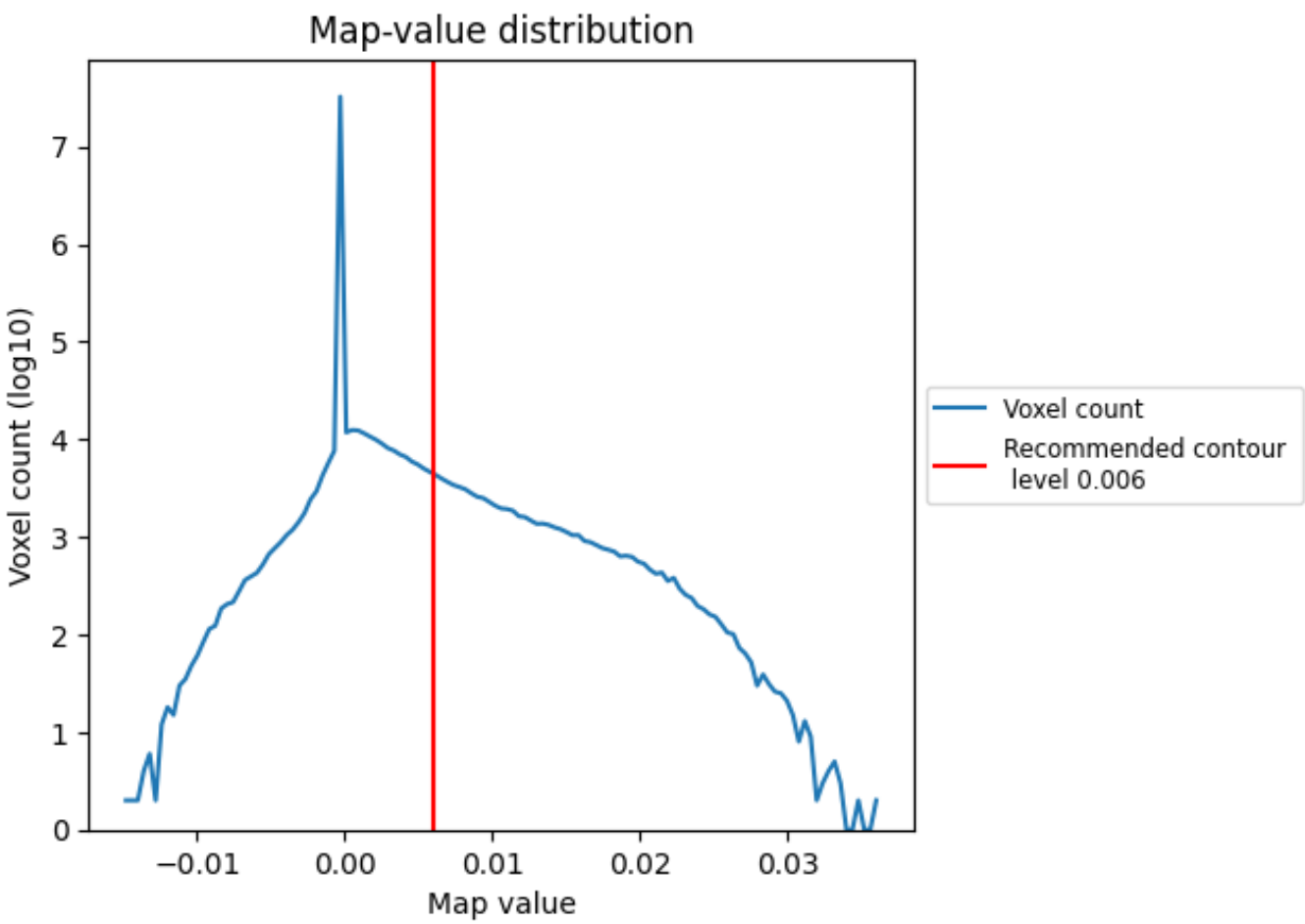


Z

7 Map analysis (i)

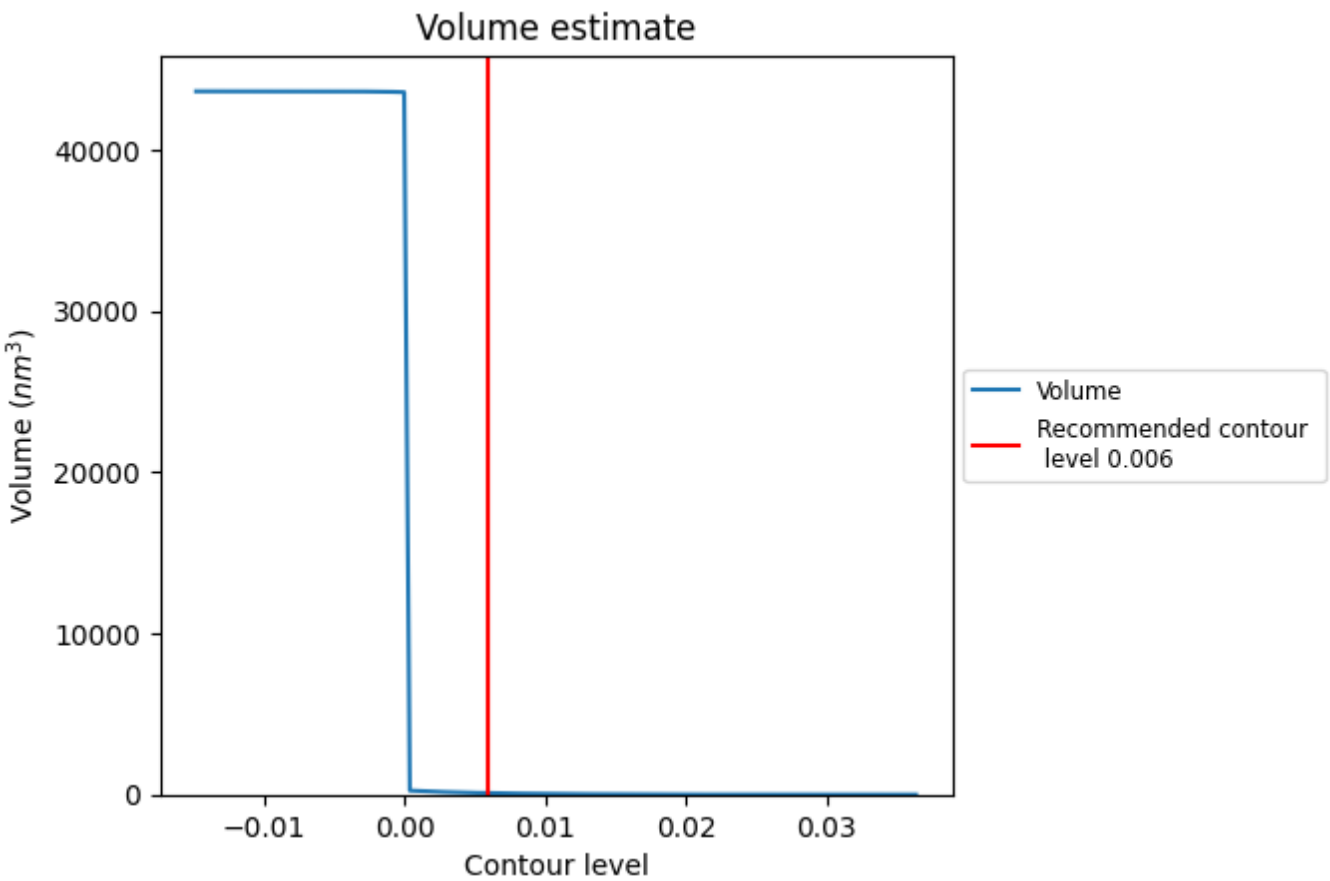
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution (i)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

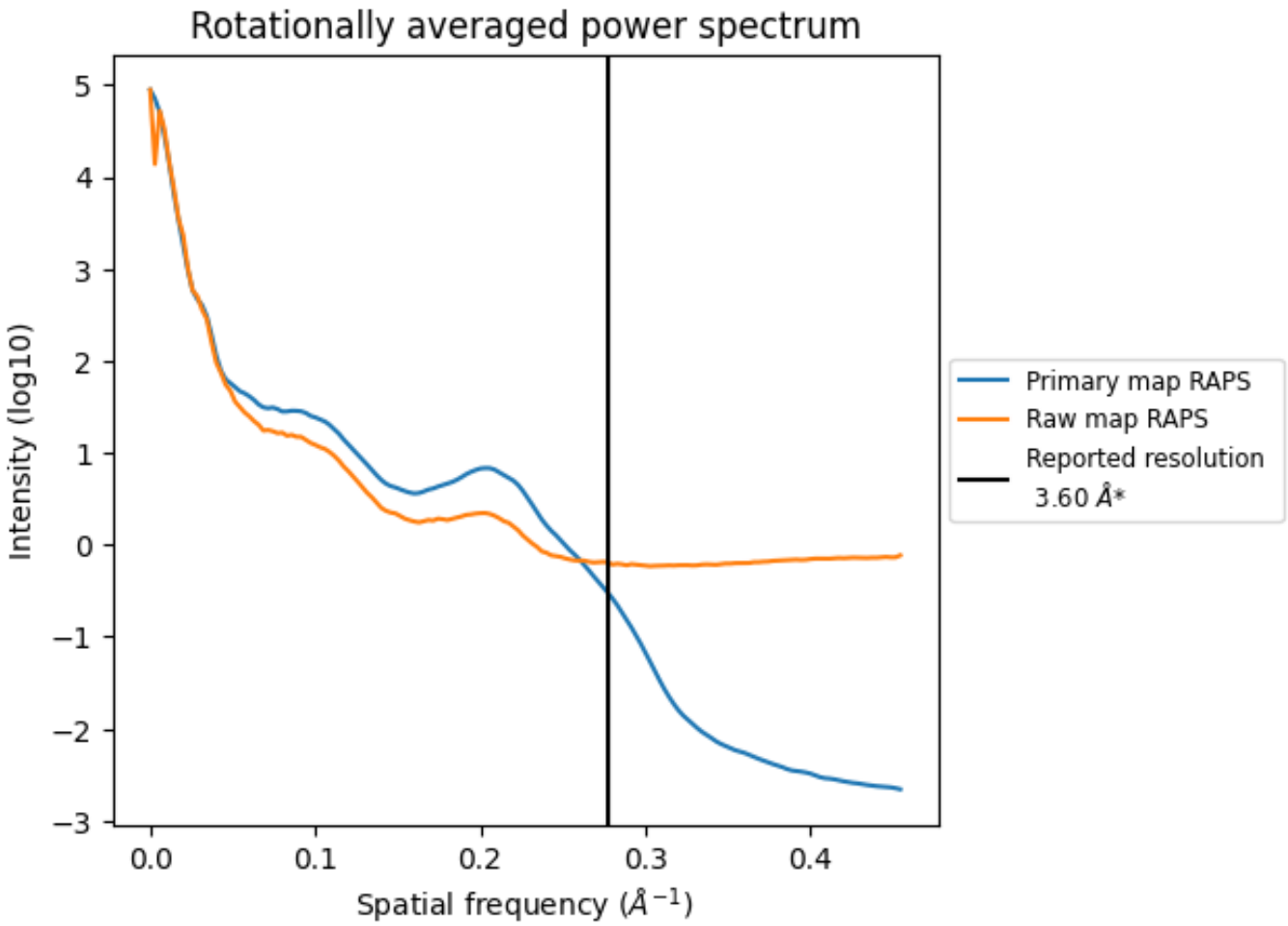
7.2 Volume estimate (i)



The volume at the recommended contour level is 91 nm^3; this corresponds to an approximate mass of 82 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum (i)

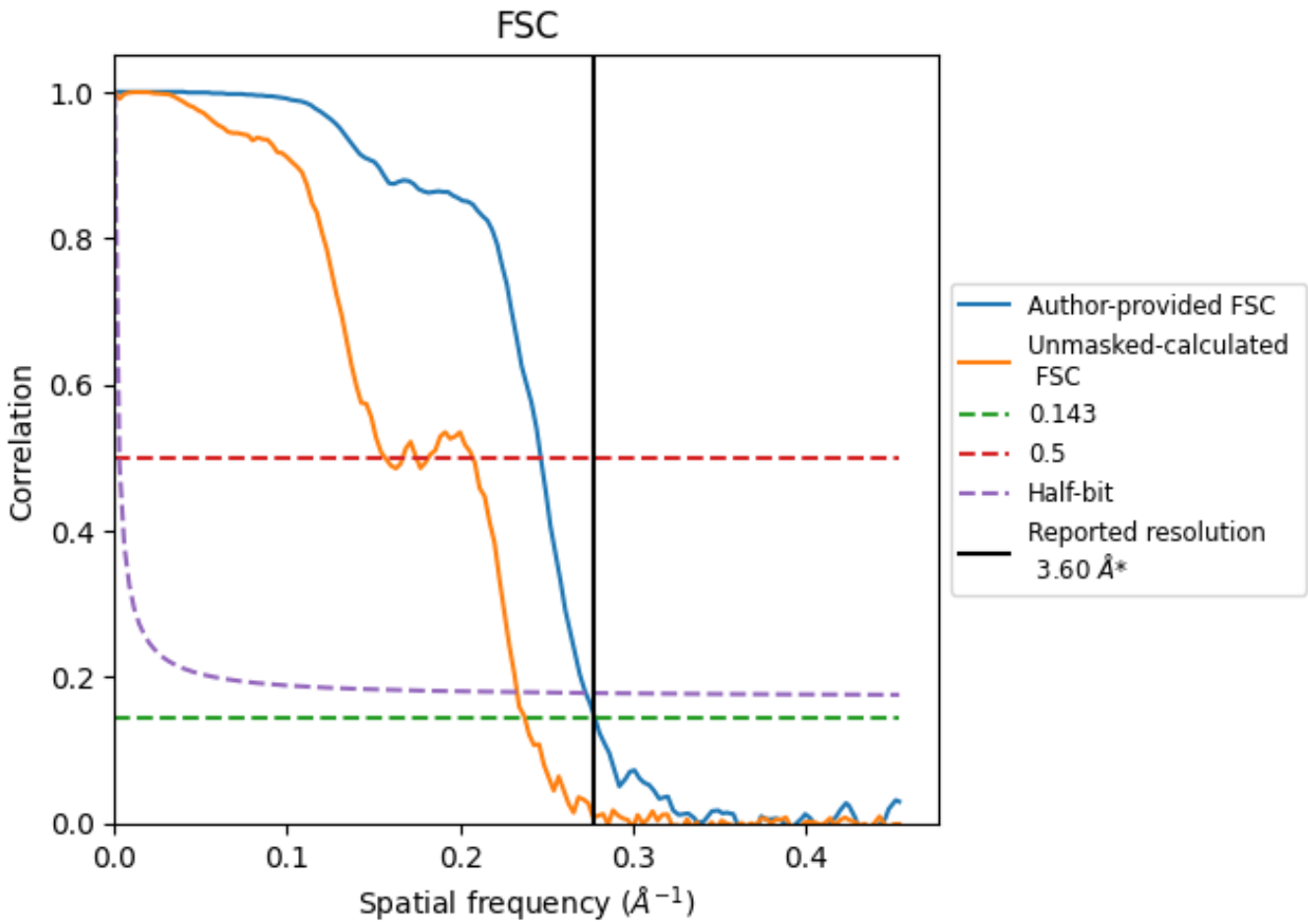


*Reported resolution corresponds to spatial frequency of 0.278 Å^{-1}

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 0.278 Å^{-1}

8.2 Resolution estimates (i)

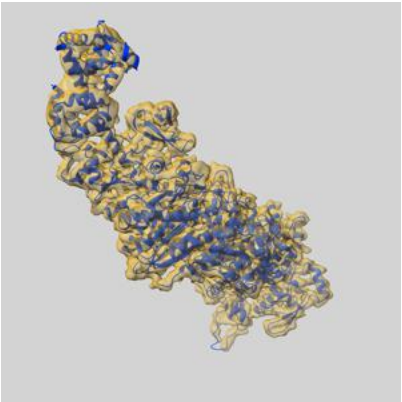
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.60	-	-
Author-provided FSC curve	3.59	4.05	3.66
Unmasked-calculated*	4.21	6.39	4.29

*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 4.21 differs from the reported value 3.6 by more than 10 %

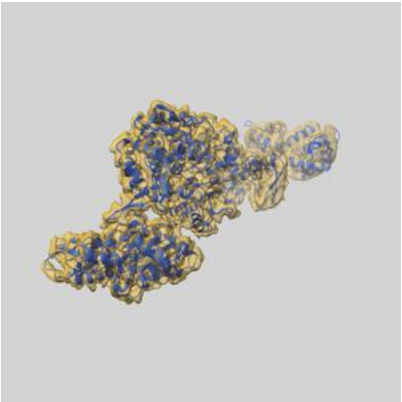
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-13508 and PDB model 7PM0. Per-residue inclusion information can be found in section 3 on page 6.

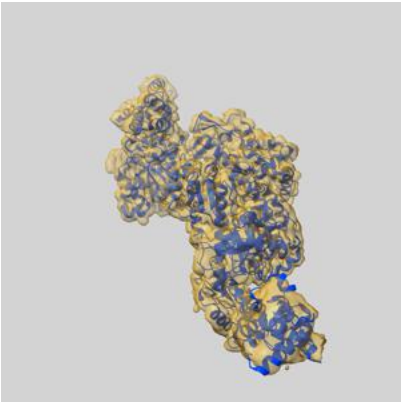
9.1 Map-model overlay (i)



X



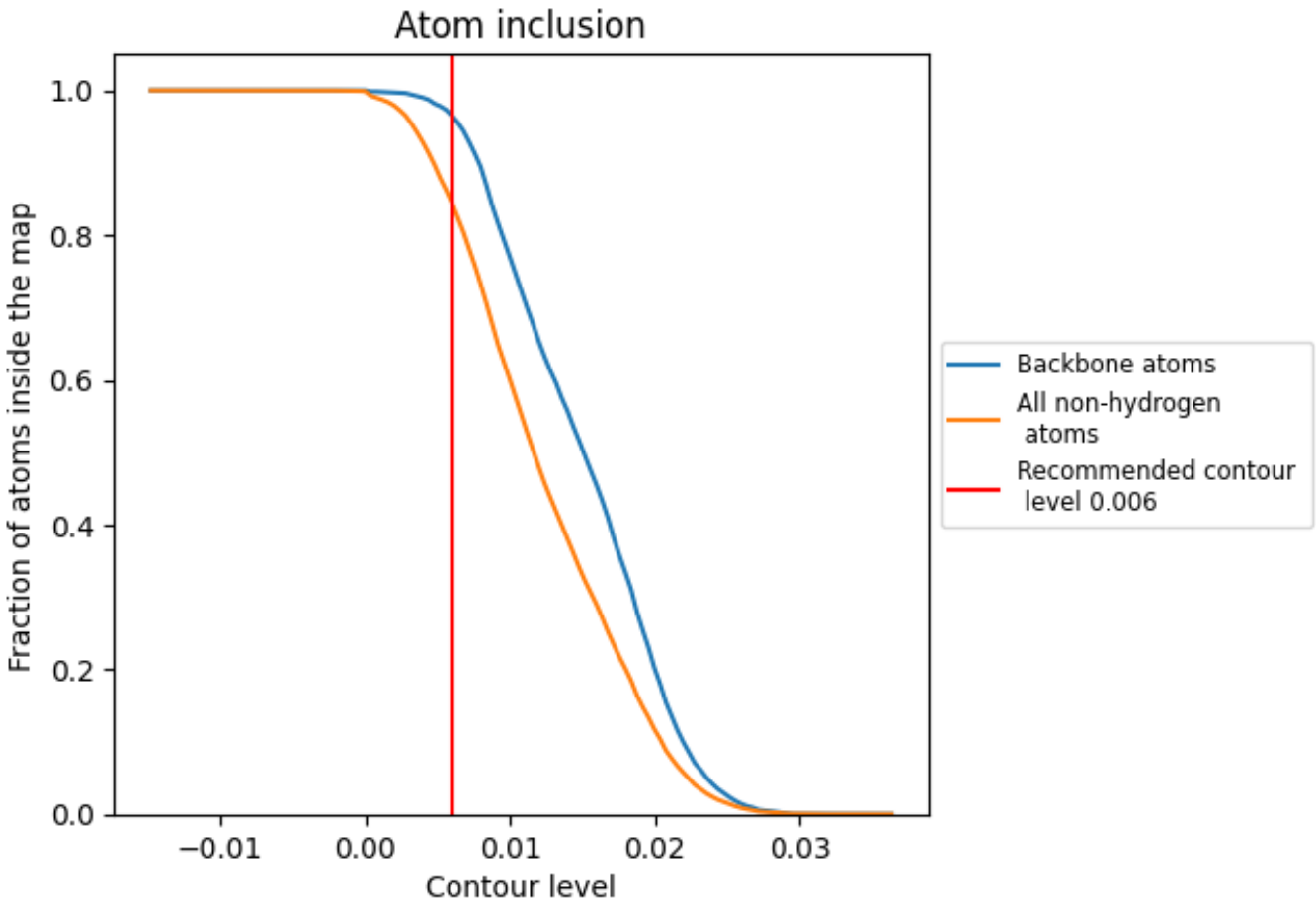
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.006 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 Atom inclusion (i)



At the recommended contour level, 96% of all backbone atoms, 84% of all non-hydrogen atoms, are inside the map.