



## Full wwPDB EM Validation Report ⓘ

Nov 13, 2022 – 10:05 AM EST

PDB ID : 6VSA  
EMDB ID : EMD-21373  
Title : Single particle reconstruction of HemQ from *Geobacillus* based on data acquired in the presence of substantial aberrations  
Authors : Bromberg, R.; Guo, Y.; Borek, D.; Otwinowski, Z.  
Deposited on : 2020-02-10  
Resolution : 2.32 Å(reported)  
Based on initial model : 1T0T

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

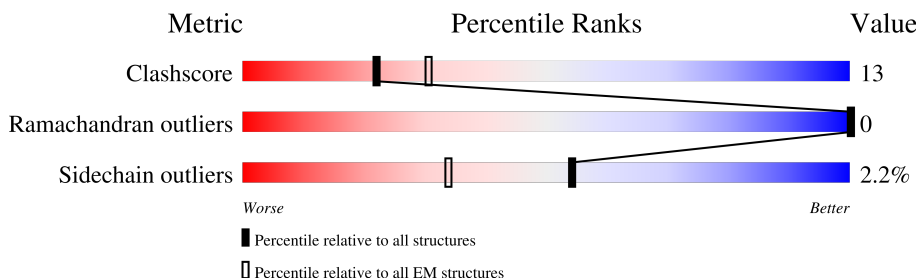
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 2.32 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  $\geq 3$ , 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	V	248	<div> <div>13%</div> <div>62%</div> <div>32%</div> <div>5%</div> </div>
1	W	248	<div> <div>12%</div> <div>65%</div> <div>29%</div> <div>• 5%</div> </div>
1	X	248	<div> <div>12%</div> <div>62%</div> <div>31%</div> <div>• 5%</div> </div>
1	Y	248	<div> <div>12%</div> <div>62%</div> <div>31%</div> <div>• 5%</div> </div>
1	Z	248	<div> <div>12%</div> <div>66%</div> <div>26%</div> <div>• 5%</div> </div>

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 9712 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 HemQ.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	V	235	Total	C	N	O	S	0	0
			1934	1246	322	356	10		
1	W	235	Total	C	N	O	S	0	0
			1934	1246	322	356	10		
1	X	235	Total	C	N	O	S	0	0
			1934	1246	322	356	10		
1	Y	235	Total	C	N	O	S	0	0
			1934	1246	322	356	10		
1	Z	235	Total	C	N	O	S	0	0
			1934	1246	322	356	10		

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
V	235	THR	ILE	conflict	UNP A0A0E0TGF4
W	235	THR	ILE	conflict	UNP A0A0E0TGF4
X	235	THR	ILE	conflict	UNP A0A0E0TGF4
Y	235	THR	ILE	conflict	UNP A0A0E0TGF4
Z	235	THR	ILE	conflict	UNP A0A0E0TGF4

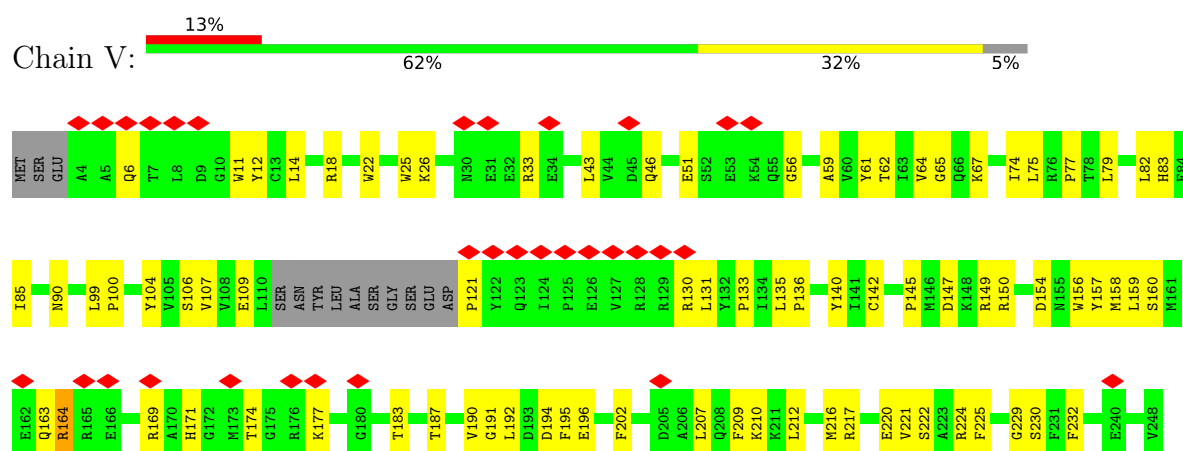
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		AltConf
2	V	8	Total	O	0
			8	8	
2	W	9	Total	O	0
			9	9	
2	X	8	Total	O	0
			8	8	
2	Y	8	Total	O	0
			8	8	
2	Z	9	Total	O	0
			9	9	

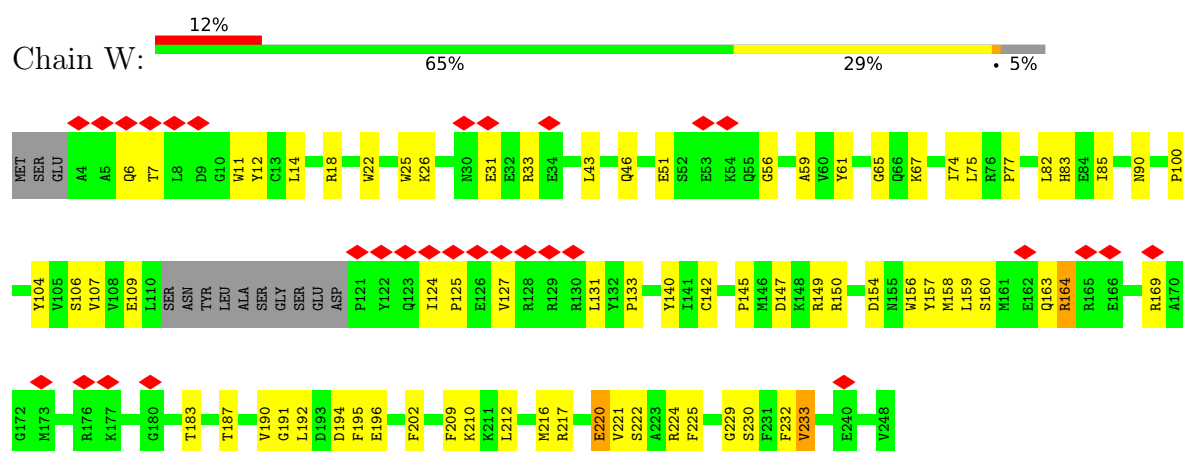
### 3 Residue-property plots

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

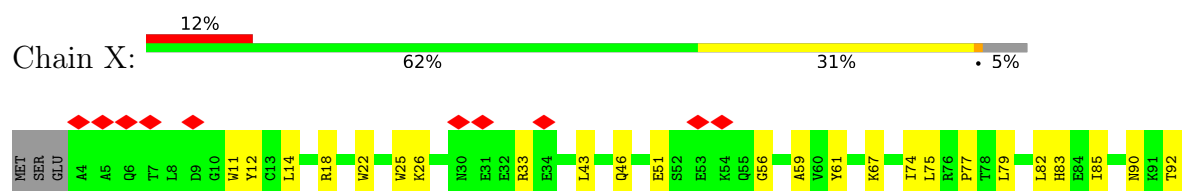
#### • Molecule 1: HemQ

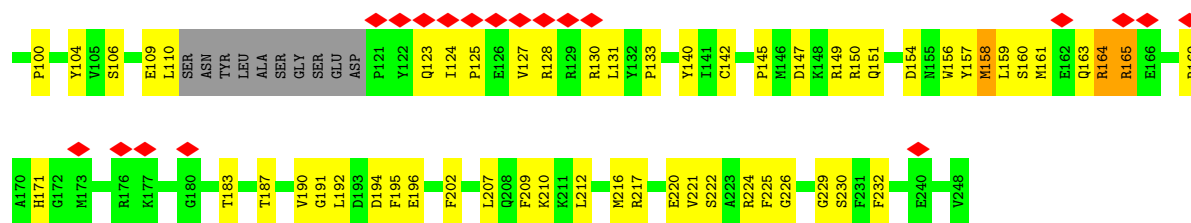


#### • Molecule 1: HemQ

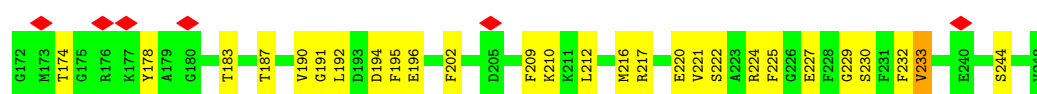


#### • Molecule 1: HemQ

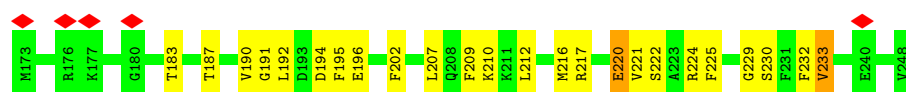
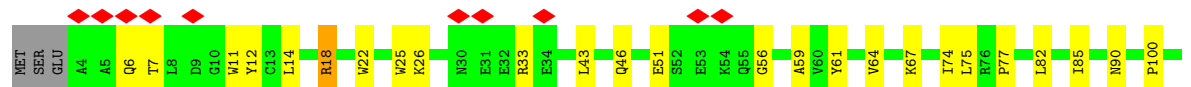




• Molecule 1: HemQ



• Molecule 1: HemQ



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C5	Depositor
Number of particles used	81302	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS TALOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	120	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.499	Depositor
Minimum map value	-0.238	Depositor
Average map value	0.020	Depositor
Map value standard deviation	0.129	Depositor
Recommended contour level	0.9	Depositor
Map size (Å)	184.32, 184.32, 184.32	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.72, 0.72, 0.72	Depositor

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	V	0.63	1/1984 (0.1%)	1.10	6/2685 (0.2%)
1	W	0.64	1/1984 (0.1%)	1.09	7/2685 (0.3%)
1	X	0.64	1/1984 (0.1%)	1.10	6/2685 (0.2%)
1	Y	0.64	1/1984 (0.1%)	1.09	6/2685 (0.2%)
1	Z	0.63	1/1984 (0.1%)	1.10	9/2685 (0.3%)
All	All	0.64	5/9920 (0.1%)	1.10	34/13425 (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	Y	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	W	230	SER	CA-CB	-6.02	1.44	1.52
1	X	230	SER	CA-CB	-5.91	1.44	1.52
1	Z	230	SER	CA-CB	-5.90	1.44	1.52
1	Y	230	SER	CA-CB	-5.83	1.44	1.52
1	V	230	SER	CA-CB	-5.73	1.44	1.52

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	V	121	PRO	CA-N-CD	-8.84	99.12	111.50
1	X	128	ARG	NE-CZ-NH1	-8.56	116.02	120.30
1	Z	169	ARG	CB-CG-CD	7.23	130.39	111.60
1	Y	127	VAL	CA-CB-CG1	6.71	120.97	110.90
1	X	164	ARG	CG-CD-NE	-6.67	97.78	111.80
1	Y	164	ARG	CG-CD-NE	-6.66	97.81	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	W	164	ARG	CG-CD-NE	-6.62	97.89	111.80
1	V	164	ARG	CG-CD-NE	-6.60	97.95	111.80
1	Z	164	ARG	CG-CD-NE	-6.59	97.96	111.80
1	Y	127	VAL	CA-CB-CG2	6.14	120.11	110.90
1	Z	18	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	V	164	ARG	NE-CZ-NH1	-5.78	117.41	120.30
1	W	169	ARG	CG-CD-NE	5.78	123.93	111.80
1	W	164	ARG	NE-CZ-NH1	-5.76	117.42	120.30
1	X	164	ARG	NE-CZ-NH1	-5.75	117.42	120.30
1	Z	164	ARG	NE-CZ-NH1	-5.67	117.46	120.30
1	Z	146	MET	CB-CA-C	-5.66	99.07	110.40
1	X	128	ARG	NE-CZ-NH2	5.57	123.09	120.30
1	V	100	PRO	N-CD-CG	-5.52	94.91	103.20
1	X	100	PRO	N-CD-CG	-5.44	95.04	103.20
1	W	100	PRO	N-CD-CG	-5.33	95.20	103.20
1	Z	100	PRO	N-CD-CG	-5.33	95.21	103.20
1	Y	164	ARG	NE-CZ-NH1	-5.25	117.67	120.30
1	W	140	TYR	CB-CA-C	-5.19	100.03	110.40
1	V	140	TYR	CB-CA-C	-5.13	100.14	110.40
1	Z	140	TYR	CB-CA-C	-5.11	100.18	110.40
1	X	140	TYR	CB-CA-C	-5.11	100.19	110.40
1	Y	140	TYR	CB-CA-C	-5.11	100.19	110.40
1	Z	146	MET	N-CA-CB	-5.08	101.45	110.60
1	Y	19	THR	CA-CB-OG1	-5.08	98.34	109.00
1	V	130	ARG	CG-CD-NE	5.06	122.43	111.80
1	W	169	ARG	NE-CZ-NH2	-5.06	117.77	120.30
1	W	220	GLU	N-CA-CB	5.00	119.61	110.60
1	Z	220	GLU	N-CA-CB	5.00	119.60	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	Y	178	TYR	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	V	1934	0	1882	60	0
1	W	1934	0	1882	58	0
1	X	1934	0	1882	55	0
1	Y	1934	0	1882	59	0
1	Z	1934	0	1882	55	0
2	V	8	0	0	1	0
2	W	9	0	0	1	0
2	X	8	0	0	0	0
2	Y	8	0	0	1	0
2	Z	9	0	0	1	0
All	All	9712	0	9410	245	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 13.

All (245) 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:Y:191:GLY:O	1:Y:192:LEU:HD23	1.72	0.90
1:V:191:GLY:O	1:V:192:LEU:HD23	1.72	0.89
1:X:191:GLY:O	1:X:192:LEU:HD23	1.72	0.89
1:Z:191:GLY:O	1:Z:192:LEU:HD23	1.72	0.89
1:W:191:GLY:O	1:W:192:LEU:HD23	1.72	0.88
1:W:67:LYS:NZ	1:Y:194:ASP:OD1	2.05	0.88
1:X:67:LYS:NZ	1:Z:194:ASP:OD1	2.05	0.88
1:X:194:ASP:OD1	1:Y:67:LYS:NZ	2.07	0.88
1:Y:126:GLU:O	1:Y:127:VAL:HB	1.74	0.87
1:V:194:ASP:OD1	1:Z:67:LYS:NZ	2.07	0.86
1:V:67:LYS:NZ	1:W:194:ASP:OD1	2.07	0.86
1:V:160:SER:HB3	1:V:163:GLN:HG3	1.69	0.75
1:W:160:SER:HB3	1:W:163:GLN:HG3	1.69	0.75
1:Z:22:TRP:O	1:Z:26:LYS:HG3	1.87	0.75
1:Y:160:SER:HB3	1:Y:163:GLN:HG3	1.69	0.75
1:W:22:TRP:O	1:W:26:LYS:HG3	1.87	0.74
1:X:22:TRP:O	1:X:26:LYS:HG3	1.87	0.74
1:X:160:SER:HB3	1:X:163:GLN:HG3	1.69	0.73
1:V:22:TRP:O	1:V:26:LYS:HG3	1.87	0.73
1:Y:22:TRP:O	1:Y:26:LYS:HG3	1.87	0.72
1:V:107:VAL:CG1	1:Z:210:LYS:HE3	2.20	0.72
1:Z:107:VAL:HG23	1:Z:192:LEU:CD1	2.21	0.71
1:V:65:GLY:O	2:V:301:HOH:O	2.09	0.70
1:Y:65:GLY:O	2:Y:301:HOH:O	2.07	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:65:GLY:O	2:W:301:HOH:O	2.09	0.70
1:Y:107:VAL:HG23	1:Y:192:LEU:CD1	2.22	0.70
1:W:107:VAL:HG23	1:W:192:LEU:CD1	2.22	0.69
1:Z:157:TYR:O	2:Z:301:HOH:O	2.11	0.67
1:X:149:ARG:HG2	1:X:151:GLN:HB2	1.80	0.63
1:Y:149:ARG:HG2	1:Y:151:GLN:HB2	1.82	0.62
1:V:171:HIS:HB2	1:V:221:VAL:HG11	1.82	0.61
1:X:190:VAL:O	1:Y:210:LYS:HA	2.01	0.61
1:W:171:HIS:HB2	1:W:221:VAL:HG11	1.83	0.61
1:X:171:HIS:HB2	1:X:221:VAL:HG11	1.82	0.61
1:W:106:SER:HB2	1:W:187:THR:HB	1.83	0.60
1:V:106:SER:HB2	1:V:187:THR:HB	1.83	0.60
1:Y:14:LEU:HB2	1:Y:82:LEU:HD13	1.84	0.60
1:W:210:LYS:HA	1:Y:190:VAL:O	2.01	0.60
1:X:106:SER:HB2	1:X:187:THR:HB	1.83	0.60
1:Y:106:SER:HB2	1:Y:187:THR:HB	1.84	0.60
1:X:14:LEU:HB2	1:X:82:LEU:HD13	1.83	0.60
1:Z:14:LEU:HB2	1:Z:82:LEU:HD13	1.84	0.60
1:Z:171:HIS:HB2	1:Z:221:VAL:HG11	1.82	0.59
1:W:124:ILE:HD11	1:W:127:VAL:HG23	1.84	0.59
1:Y:171:HIS:HB2	1:Y:221:VAL:HG11	1.83	0.59
1:W:14:LEU:HB2	1:W:82:LEU:HD13	1.84	0.59
1:Y:145:PRO:HG2	1:Y:229:GLY:HA3	1.84	0.59
1:Z:106:SER:HB2	1:Z:187:THR:HB	1.84	0.59
1:X:210:LYS:HA	1:Z:190:VAL:O	2.01	0.59
1:V:145:PRO:HG2	1:V:229:GLY:HA3	1.84	0.59
1:Y:174:THR:CG2	1:Y:216:MET:HG2	2.33	0.59
1:Z:145:PRO:HG2	1:Z:229:GLY:HA3	1.84	0.59
1:V:212:LEU:O	1:V:216:MET:HG3	2.03	0.59
1:X:124:ILE:HD11	1:X:127:VAL:HG23	1.84	0.58
1:X:212:LEU:O	1:X:216:MET:HG3	2.03	0.58
1:W:145:PRO:HG2	1:W:229:GLY:HA3	1.84	0.58
1:V:190:VAL:O	1:Z:210:LYS:HA	2.03	0.58
1:X:145:PRO:HG2	1:X:229:GLY:HA3	1.84	0.58
1:V:14:LEU:HB2	1:V:82:LEU:HD13	1.84	0.58
1:Z:212:LEU:O	1:Z:216:MET:HG3	2.03	0.58
1:Y:212:LEU:O	1:Y:216:MET:HG3	2.03	0.57
1:V:210:LYS:HA	1:W:190:VAL:O	2.03	0.57
1:W:212:LEU:O	1:W:216:MET:HG3	2.05	0.57
1:Y:174:THR:HG23	1:Y:216:MET:HG2	1.87	0.56
1:W:158:MET:HA	1:W:158:MET:CE	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:107:VAL:HG13	1:Z:210:LYS:HE3	1.89	0.55
1:X:158:MET:HA	1:X:158:MET:CE	2.36	0.55
1:V:158:MET:HA	1:V:158:MET:CE	2.36	0.55
1:V:104:TYR:OH	1:V:195:PHE:HB2	2.07	0.55
1:Z:104:TYR:OH	1:Z:195:PHE:HB2	2.07	0.55
1:Z:18:ARG:NH2	1:Z:90:ASN:OD1	2.35	0.54
1:Z:158:MET:HA	1:Z:158:MET:CE	2.36	0.54
1:X:104:TYR:OH	1:X:195:PHE:HB2	2.08	0.54
1:Y:104:TYR:OH	1:Y:195:PHE:HB2	2.07	0.54
1:W:104:TYR:OH	1:W:195:PHE:HB2	2.07	0.54
1:X:110:LEU:HD11	1:X:130:ARG:HE	1.74	0.53
1:X:147:ASP:OD1	1:X:196:GLU:N	2.42	0.53
1:W:147:ASP:OD1	1:W:196:GLU:N	2.42	0.53
1:V:147:ASP:OD1	1:V:196:GLU:N	2.42	0.53
1:Z:147:ASP:OD1	1:Z:196:GLU:N	2.42	0.53
1:Y:149:ARG:CG	1:Y:151:GLN:HB2	2.39	0.53
1:X:149:ARG:CG	1:X:151:GLN:HB2	2.39	0.53
1:Y:147:ASP:OD1	1:Y:196:GLU:N	2.42	0.53
1:W:212:LEU:HD11	1:W:216:MET:HE3	1.92	0.52
1:X:192:LEU:HA	1:Y:233:VAL:HG11	1.91	0.52
1:V:192:LEU:HA	1:Z:233:VAL:HG11	1.91	0.51
1:V:150:ARG:CG	1:V:157:TYR:HD2	2.24	0.51
1:W:150:ARG:CG	1:W:157:TYR:HD2	2.24	0.50
1:V:43:LEU:O	1:V:46:GLN:HB2	2.11	0.50
1:V:109:GLU:O	1:V:109:GLU:HG3	2.11	0.50
1:X:150:ARG:CG	1:X:157:TYR:HD2	2.25	0.50
1:Y:150:ARG:CG	1:Y:157:TYR:HD2	2.24	0.50
1:Z:150:ARG:CG	1:Z:157:TYR:HD2	2.25	0.50
1:Y:43:LEU:O	1:Y:46:GLN:HB2	2.11	0.50
1:W:43:LEU:O	1:W:46:GLN:HB2	2.11	0.50
1:W:233:VAL:HG11	1:Y:192:LEU:HA	1.92	0.50
1:Z:109:GLU:O	1:Z:109:GLU:HG3	2.12	0.50
1:X:109:GLU:HG3	1:X:109:GLU:O	2.12	0.50
1:X:161:MET:SD	1:X:165:ARG:NH1	2.84	0.49
1:X:43:LEU:O	1:X:46:GLN:HB2	2.11	0.49
1:W:221:VAL:O	1:W:225:PHE:HB2	2.12	0.49
1:V:221:VAL:O	1:V:225:PHE:HB2	2.12	0.49
1:Z:43:LEU:O	1:Z:46:GLN:HB2	2.11	0.49
1:V:174:THR:HA	1:V:177:LYS:HE3	1.93	0.49
1:W:109:GLU:O	1:W:109:GLU:HG3	2.12	0.49
1:Y:109:GLU:O	1:Y:109:GLU:HG3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:221:VAL:O	1:X:225:PHE:HB2	2.12	0.49
1:Y:221:VAL:O	1:Y:225:PHE:HB2	2.12	0.48
1:V:18:ARG:NH2	1:V:90:ASN:OD1	2.42	0.48
1:Z:221:VAL:O	1:Z:225:PHE:HB2	2.12	0.48
1:Z:159:LEU:O	1:Z:164:ARG:NH1	2.48	0.47
1:X:18:ARG:NH2	1:X:90:ASN:OD1	2.42	0.47
1:V:150:ARG:NH1	1:Z:149:ARG:HH22	2.12	0.47
1:X:159:LEU:O	1:X:164:ARG:NH1	2.48	0.47
1:X:150:ARG:NH1	1:Y:149:ARG:HH22	2.13	0.47
1:V:159:LEU:O	1:V:164:ARG:NH1	2.48	0.47
1:W:149:ARG:HH22	1:Y:150:ARG:NH1	2.14	0.46
1:X:224:ARG:HG3	1:Z:158:MET:HG3	1.98	0.46
1:X:192:LEU:HA	1:Y:233:VAL:CG1	2.46	0.46
1:X:217:ARG:HA	1:X:222:SER:HB2	1.98	0.46
1:X:149:ARG:HH22	1:Z:150:ARG:NH1	2.13	0.46
1:W:159:LEU:O	1:W:164:ARG:NH1	2.48	0.46
1:W:233:VAL:CG1	1:Y:192:LEU:HA	2.46	0.46
1:Z:158:MET:HA	1:Z:158:MET:HE2	1.97	0.46
1:V:192:LEU:HA	1:Z:233:VAL:CG1	2.45	0.46
1:Y:159:LEU:O	1:Y:164:ARG:NH1	2.48	0.46
1:V:158:MET:HG3	1:Z:224:ARG:HG3	1.98	0.45
1:V:62:THR:HB	1:W:83:HIS:CE1	2.52	0.45
1:W:183:THR:HB	1:W:202:PHE:HB2	1.98	0.45
1:Y:154:ASP:HA	1:Y:224:ARG:HH21	1.82	0.45
1:Y:217:ARG:HA	1:Y:222:SER:HB2	1.97	0.45
1:W:233:VAL:HG21	1:Y:79:LEU:HD21	1.99	0.45
1:V:79:LEU:HD21	1:Z:233:VAL:HG21	1.98	0.45
1:X:83:HIS:CE1	1:Y:62:THR:HB	2.52	0.45
1:Y:107:VAL:HG23	1:Y:192:LEU:HD12	1.98	0.45
1:W:217:ARG:HA	1:W:222:SER:HB2	1.98	0.45
1:X:183:THR:HB	1:X:202:PHE:HB2	1.98	0.45
1:Y:183:THR:HB	1:Y:202:PHE:HB2	1.98	0.45
1:W:59:ALA:HB1	1:W:61:TYR:HE1	1.82	0.45
1:Z:59:ALA:HB1	1:Z:61:TYR:HE1	1.82	0.45
1:V:154:ASP:HA	1:V:224:ARG:HH21	1.82	0.45
1:X:149:ARG:NH2	1:Z:150:ARG:NH1	2.64	0.45
1:Z:154:ASP:HA	1:Z:224:ARG:HH21	1.82	0.45
1:V:183:THR:HB	1:V:202:PHE:HB2	1.98	0.45
1:V:224:ARG:HG3	1:W:158:MET:HG3	1.98	0.45
1:W:107:VAL:HG23	1:W:192:LEU:HD12	1.98	0.45
1:W:154:ASP:HA	1:W:224:ARG:HH21	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Z:107:VAL:HG23	1:Z:192:LEU:HD11	1.98	0.45
1:Z:217:ARG:HA	1:Z:222:SER:HB2	1.98	0.45
1:Z:183:THR:HB	1:Z:202:PHE:HB2	1.98	0.45
1:V:217:ARG:HA	1:V:222:SER:HB2	1.98	0.44
1:W:220:GLU:O	1:W:224:ARG:HB2	2.18	0.44
1:V:149:ARG:HH22	1:W:150:ARG:NH1	2.15	0.44
1:X:59:ALA:HB1	1:X:61:TYR:HE1	1.82	0.44
1:W:149:ARG:NH2	1:Y:150:ARG:NH1	2.66	0.44
1:V:59:ALA:HB1	1:V:61:TYR:HE1	1.82	0.44
1:V:150:ARG:NH1	1:Z:149:ARG:NH2	2.65	0.44
1:Y:59:ALA:HB1	1:Y:61:TYR:HE1	1.83	0.44
1:V:220:GLU:O	1:V:224:ARG:HB2	2.17	0.44
1:X:154:ASP:HA	1:X:224:ARG:HH21	1.82	0.44
1:Z:220:GLU:O	1:Z:224:ARG:HB2	2.18	0.44
1:Y:220:GLU:O	1:Y:224:ARG:HB2	2.17	0.44
1:X:11:TRP:CD1	1:X:77:PRO:HA	2.54	0.43
1:X:79:LEU:HD21	1:Y:233:VAL:HG21	1.99	0.43
1:V:149:ARG:NH2	1:W:150:ARG:NH1	2.66	0.43
1:Z:11:TRP:CD1	1:Z:77:PRO:HA	2.54	0.43
1:Z:131:LEU:C	1:Z:133:PRO:HD3	2.39	0.43
1:V:131:LEU:C	1:V:133:PRO:HD3	2.39	0.43
1:X:207:LEU:HD12	1:X:207:LEU:HA	1.88	0.43
1:X:220:GLU:O	1:X:224:ARG:HB2	2.18	0.43
1:Z:142:CYS:HB2	1:Z:209:PHE:CD2	2.54	0.43
1:V:191:GLY:C	1:V:192:LEU:HD23	2.38	0.43
1:X:131:LEU:C	1:X:133:PRO:HD3	2.39	0.43
1:X:150:ARG:NH1	1:Y:149:ARG:NH2	2.66	0.43
1:Y:11:TRP:CD1	1:Y:77:PRO:HA	2.54	0.43
1:Y:131:LEU:C	1:Y:133:PRO:HD3	2.39	0.43
1:Y:142:CYS:HB2	1:Y:209:PHE:CD2	2.54	0.43
1:X:142:CYS:HB2	1:X:209:PHE:CD2	2.54	0.43
1:Z:160:SER:OG	1:Z:163:GLN:HG3	2.18	0.43
1:W:124:ILE:HA	1:W:125:PRO:HD3	1.92	0.43
1:W:11:TRP:CD1	1:W:77:PRO:HA	2.54	0.42
1:V:11:TRP:CD1	1:V:77:PRO:HA	2.54	0.42
1:W:131:LEU:C	1:W:133:PRO:HD3	2.39	0.42
1:W:224:ARG:HG3	1:Y:158:MET:HG3	2.01	0.42
1:Y:191:GLY:C	1:Y:192:LEU:HD23	2.38	0.42
1:V:158:MET:HA	1:V:158:MET:HE2	2.02	0.42
1:W:12:TYR:O	1:W:75:LEU:HA	2.20	0.42
1:V:135:LEU:HA	1:V:136:PRO:HD3	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:142:CYS:HB2	1:V:209:PHE:CD2	2.54	0.42
1:V:232:PHE:CD1	1:V:232:PHE:N	2.88	0.42
1:W:142:CYS:HB2	1:W:209:PHE:CD2	2.54	0.42
1:W:195:PHE:HB3	1:W:196:GLU:H	1.76	0.42
1:W:18:ARG:NH2	1:W:90:ASN:OD1	2.42	0.42
1:W:232:PHE:CD1	1:W:232:PHE:N	2.88	0.42
1:Y:129:ARG:CZ	1:Y:129:ARG:HB3	2.50	0.42
1:X:195:PHE:HB3	1:X:196:GLU:H	1.76	0.42
1:Z:25:TRP:O	1:Z:33:ARG:HD3	2.20	0.42
1:Y:174:THR:HG21	1:Y:216:MET:HG2	2.01	0.42
1:W:74:ILE:HG21	1:W:85:ILE:HD12	2.01	0.42
1:W:82:LEU:HA	1:W:82:LEU:HD23	1.74	0.42
1:W:107:VAL:HG23	1:W:192:LEU:HD11	1.99	0.42
1:X:232:PHE:CD1	1:X:232:PHE:N	2.88	0.42
1:Y:124:ILE:O	1:Y:128:ARG:HB2	2.20	0.42
1:V:12:TYR:O	1:V:75:LEU:HA	2.20	0.41
1:V:25:TRP:O	1:V:33:ARG:HD3	2.20	0.41
1:V:82:LEU:HD23	1:V:82:LEU:HA	1.74	0.41
1:V:74:ILE:HG21	1:V:85:ILE:HD12	2.02	0.41
1:Y:25:TRP:O	1:Y:33:ARG:HD3	2.20	0.41
1:Z:74:ILE:HG21	1:Z:85:ILE:HD12	2.02	0.41
1:X:25:TRP:O	1:X:33:ARG:HD3	2.20	0.41
1:X:12:TYR:O	1:X:75:LEU:HA	2.20	0.41
1:Z:232:PHE:CD1	1:Z:232:PHE:N	2.88	0.41
1:V:64:VAL:HG22	1:W:83:HIS:HB2	2.02	0.41
1:W:142:CYS:HA	1:W:232:PHE:O	2.21	0.41
1:Y:12:TYR:O	1:Y:75:LEU:HA	2.20	0.41
1:Y:74:ILE:HG21	1:Y:85:ILE:HD12	2.02	0.41
1:V:156:TRP:CE2	1:V:164:ARG:HD2	2.56	0.41
1:V:83:HIS:HB2	1:Z:64:VAL:HG22	2.02	0.41
1:W:191:GLY:C	1:W:192:LEU:HD23	2.38	0.41
1:Y:232:PHE:CD1	1:Y:232:PHE:N	2.88	0.41
1:Y:51:GLU:HA	1:Y:56:GLY:O	2.21	0.41
1:Z:51:GLU:HA	1:Z:56:GLY:O	2.21	0.41
1:Z:107:VAL:HG23	1:Z:192:LEU:HD12	1.98	0.41
1:Z:156:TRP:O	1:Z:164:ARG:HD3	2.21	0.41
1:W:25:TRP:O	1:W:33:ARG:HD3	2.20	0.41
1:X:74:ILE:HG21	1:X:85:ILE:HD12	2.02	0.41
1:X:156:TRP:O	1:X:164:ARG:HD3	2.21	0.41
1:Z:142:CYS:HA	1:Z:232:PHE:O	2.20	0.41
1:V:142:CYS:HA	1:V:232:PHE:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:207:LEU:HD12	1:V:207:LEU:HA	1.87	0.40
1:W:51:GLU:HA	1:W:56:GLY:O	2.21	0.40
1:V:51:GLU:HA	1:V:56:GLY:O	2.21	0.40
1:X:92:THR:O	1:Y:26:LYS:NZ	2.55	0.40
1:Z:12:TYR:O	1:Z:75:LEU:HA	2.20	0.40
1:Z:156:TRP:CE2	1:Z:164:ARG:HD2	2.56	0.40
1:V:99:LEU:N	1:V:99:LEU:HD23	2.36	0.40
1:X:33:ARG:HD3	1:X:33:ARG:HH11	1.76	0.40
1:W:26:LYS:NZ	1:Y:92:THR:O	2.54	0.40
1:W:156:TRP:CE2	1:W:164:ARG:HD2	2.56	0.40
1:X:51:GLU:HA	1:X:56:GLY:O	2.21	0.40
1:X:124:ILE:HA	1:X:125:PRO:HD3	1.93	0.40
1:V:145:PRO:HB2	1:V:195:PHE:CG	2.57	0.40
1:V:156:TRP:O	1:V:164:ARG:HD3	2.21	0.40
1:X:226:GLY:O	1:Z:150:ARG:NH1	2.49	0.40
1:Y:107:VAL:HG23	1:Y:192:LEU:HD11	1.98	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	V	231/248 (93%)	226 (98%)	5 (2%)	0	100	100
1	W	231/248 (93%)	226 (98%)	5 (2%)	0	100	100
1	X	231/248 (93%)	226 (98%)	5 (2%)	0	100	100
1	Y	231/248 (93%)	225 (97%)	6 (3%)	0	100	100
1	Z	231/248 (93%)	226 (98%)	5 (2%)	0	100	100
All	All	1155/1240 (93%)	1129 (98%)	26 (2%)	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	V	207/218 (95%)	205 (99%)	2 (1%)	76	87
1	W	207/218 (95%)	203 (98%)	4 (2%)	57	73
1	X	207/218 (95%)	203 (98%)	4 (2%)	57	73
1	Y	207/218 (95%)	201 (97%)	6 (3%)	42	57
1	Z	207/218 (95%)	200 (97%)	7 (3%)	37	51
All	All	1035/1090 (95%)	1012 (98%)	23 (2%)	54	68

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	V	6	GLN
1	V	169	ARG
1	W	6	GLN
1	W	7	THR
1	W	31	GLU
1	W	233	VAL
1	X	123	GLN
1	X	158	MET
1	X	165	ARG
1	X	169	ARG
1	Y	7	THR
1	Y	161	MET
1	Y	169	ARG
1	Y	227	GLU
1	Y	233	VAL
1	Y	244	SER
1	Z	6	GLN
1	Z	7	THR
1	Z	146	MET
1	Z	160	SER
1	Z	169	ARG
1	Z	207	LEU
1	Z	233	VAL



Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	W	83	HIS
1	W	171	HIS
1	X	83	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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

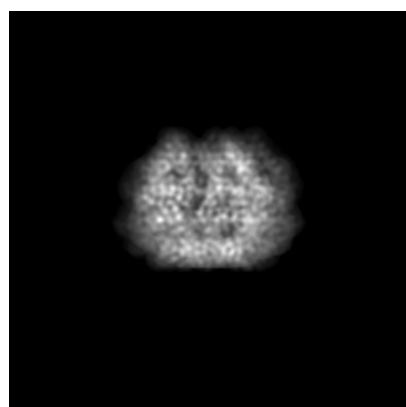
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21373. These allow visual inspection of the internal detail of the map and identification of artifacts.

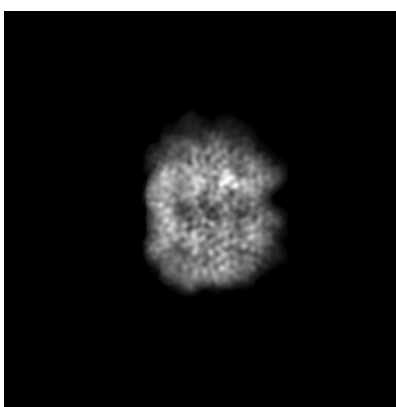
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

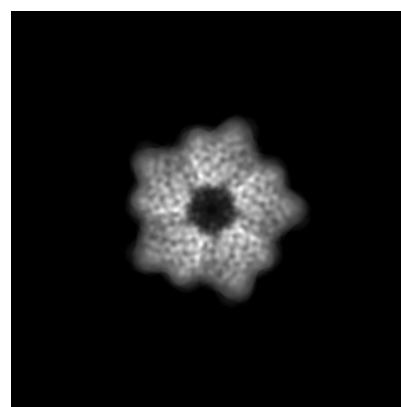
#### 6.1.1 Primary map



X



Y

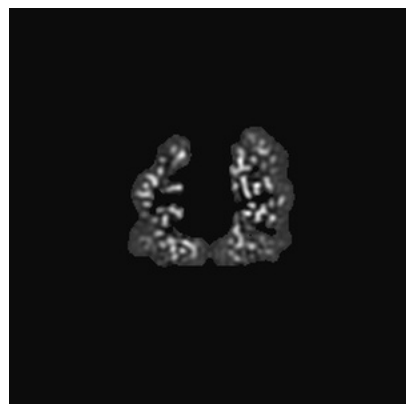


Z

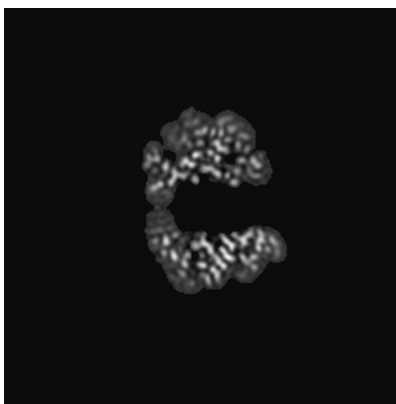
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

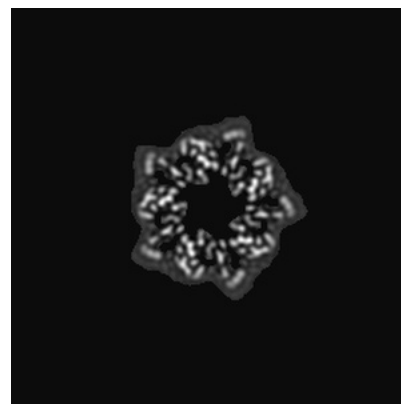
#### 6.2.1 Primary map



X Index: 128



Y Index: 128

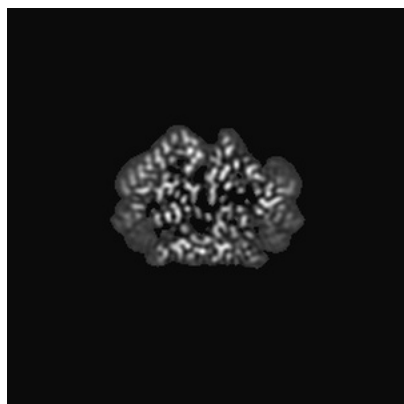


Z Index: 128

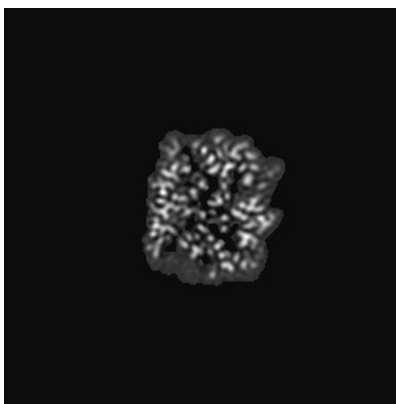
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

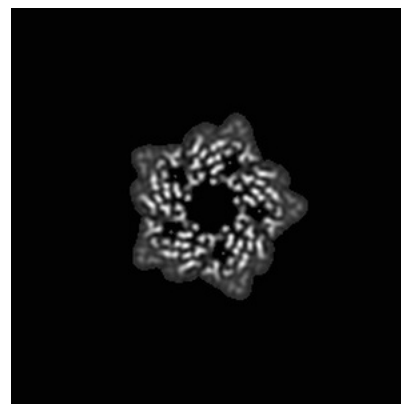
### 6.3.1 Primary map



X Index: 145



Y Index: 147

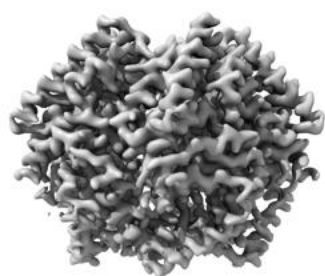


Z Index: 141

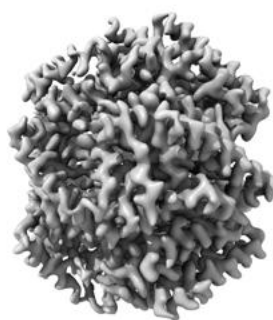
The images above show the largest variance slices of the map in three orthogonal directions.

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

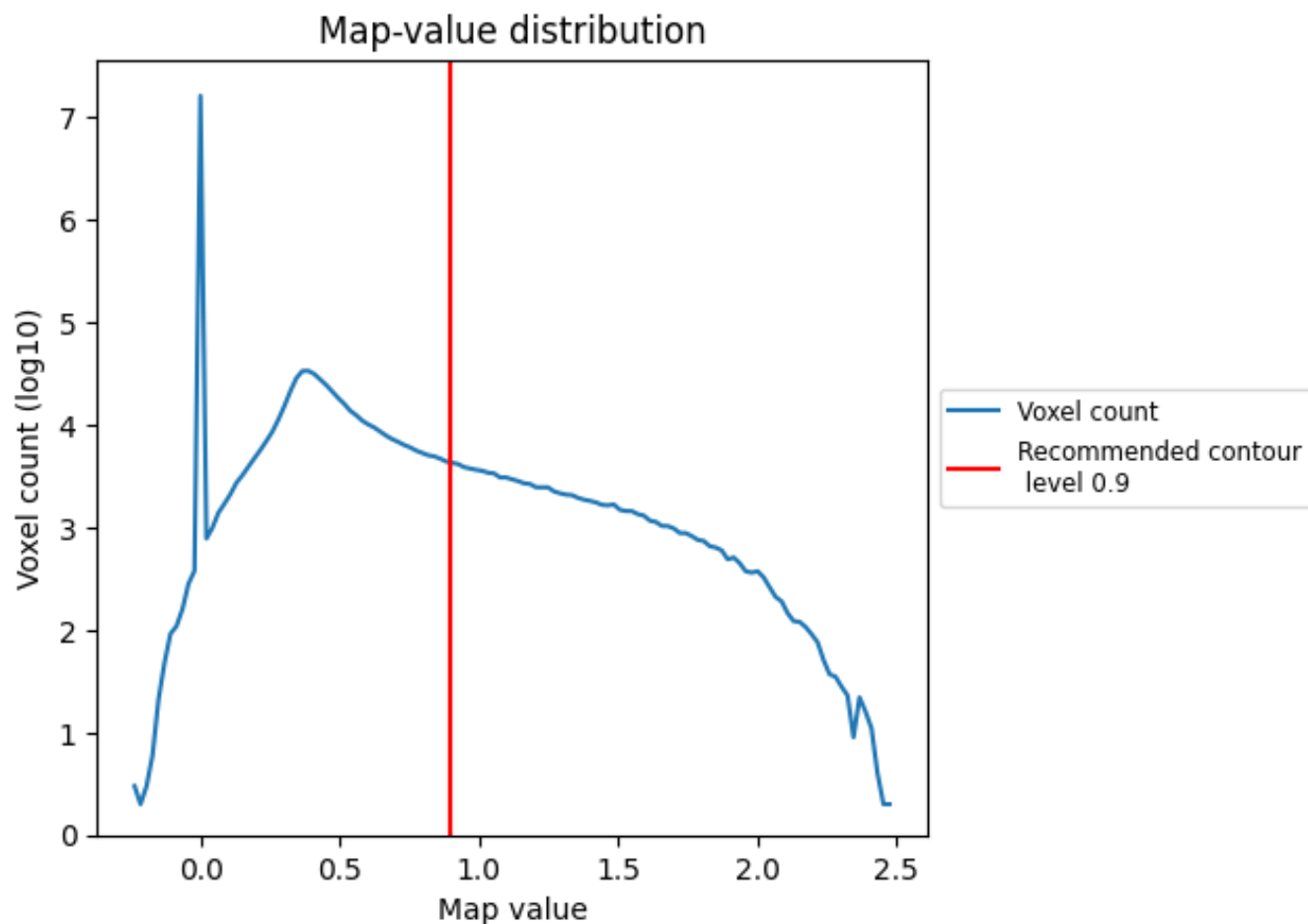
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

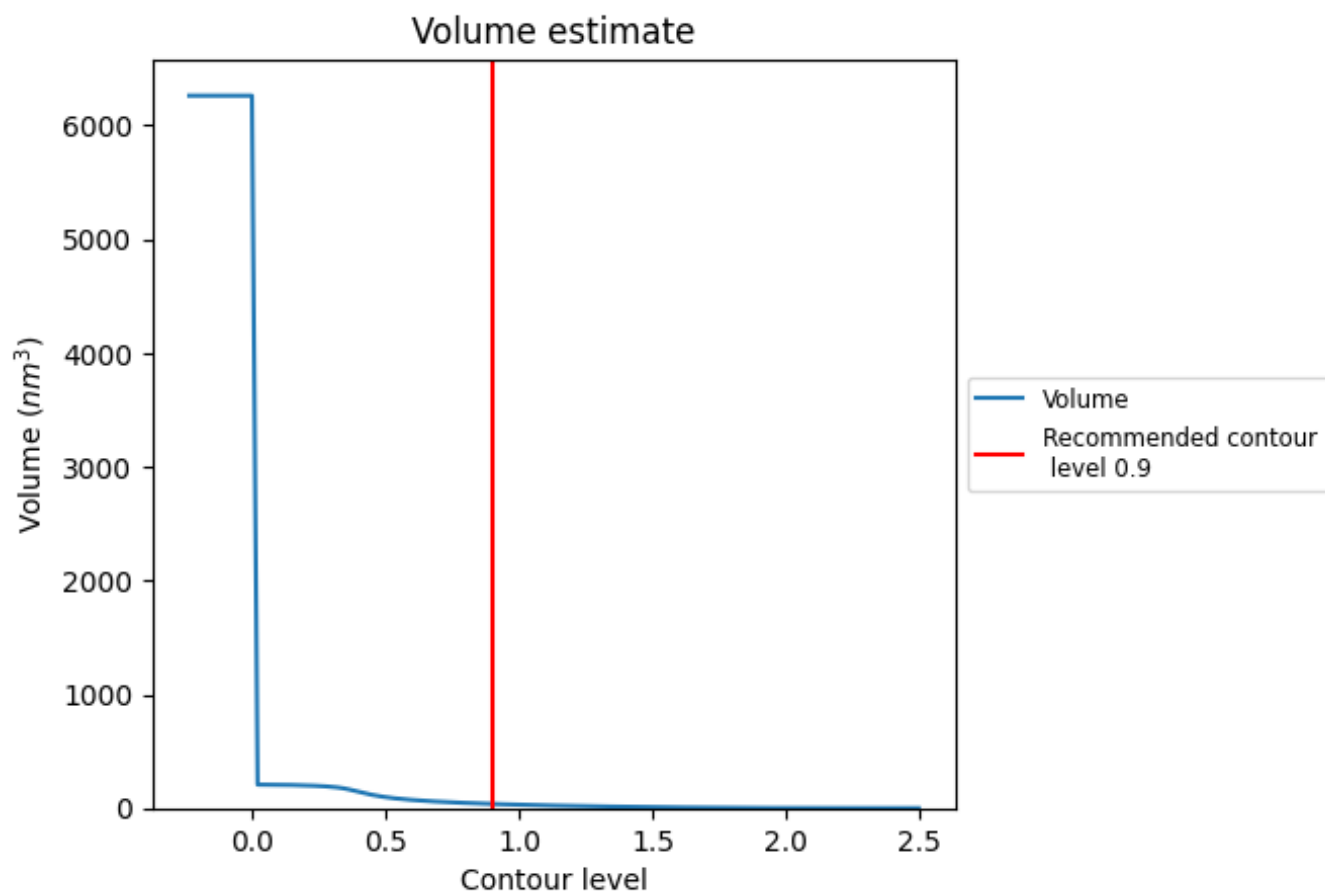
This section contains the results of statistical analysis of the map.

### 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.

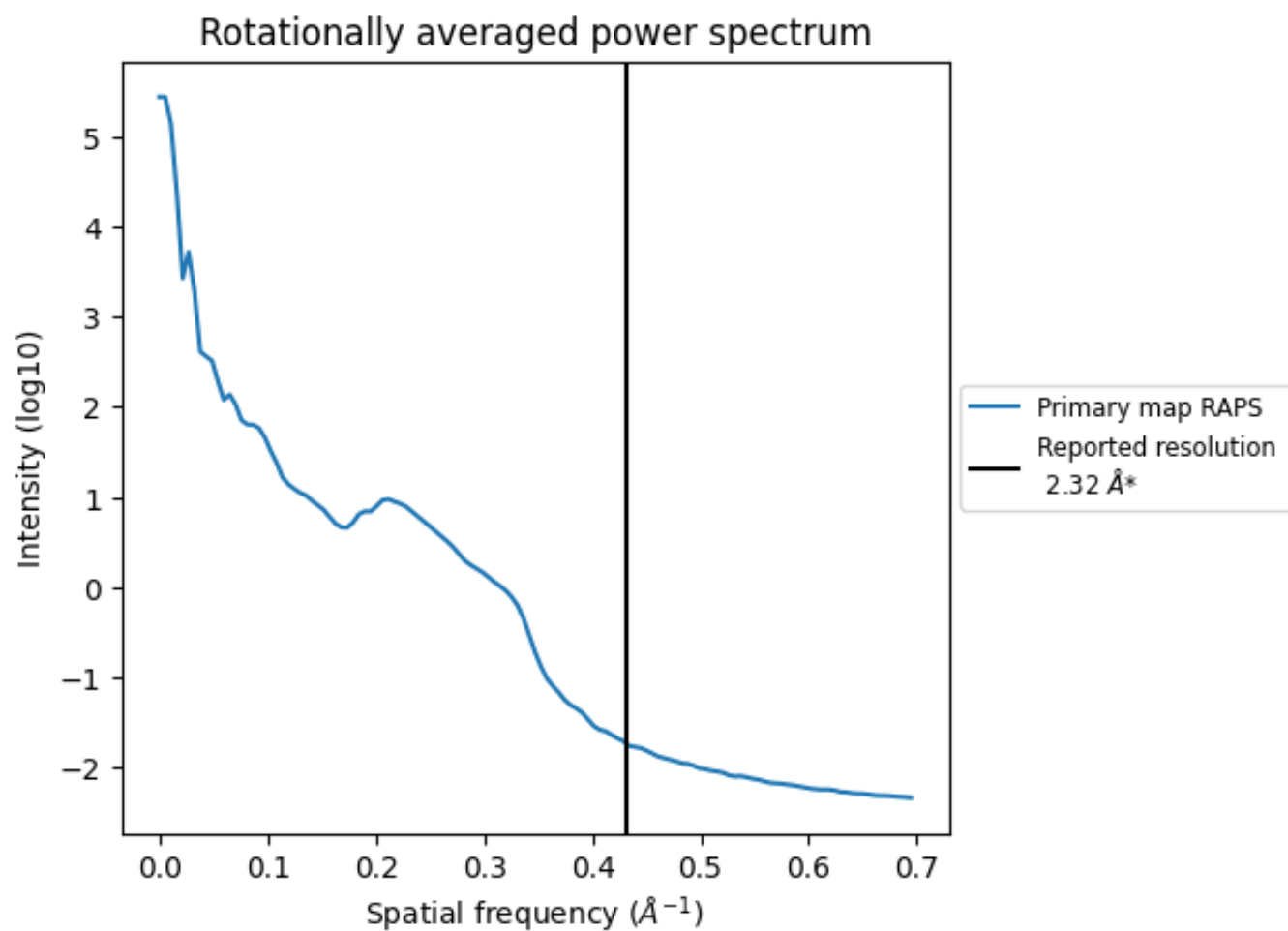
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 37  $\text{nm}^3$ ; this corresponds to an approximate mass of 33 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 0.431 Å<sup>-1</sup>

## 8 Fourier-Shell correlation ⓘ

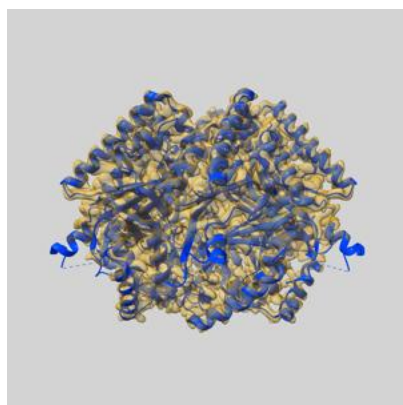
This section was not generated. No FSC curve or half-maps provided.



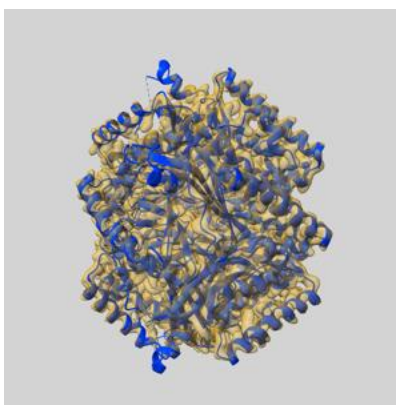
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-21373 and PDB model 6VSA. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

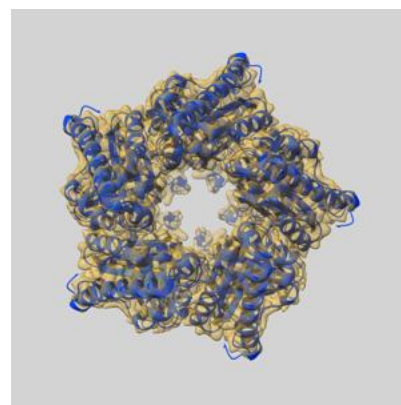
### 9.1 Map-model overlay [i](#)



X



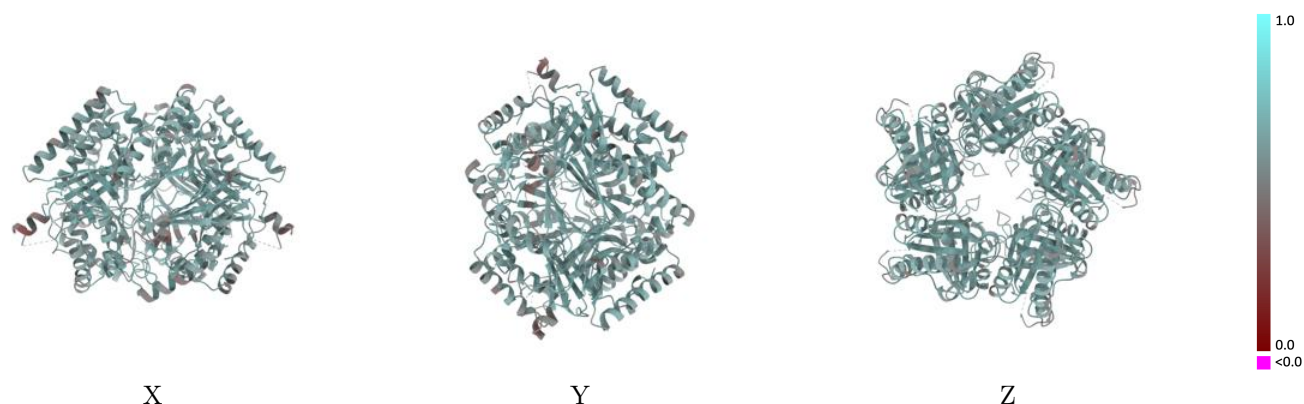
Y



Z

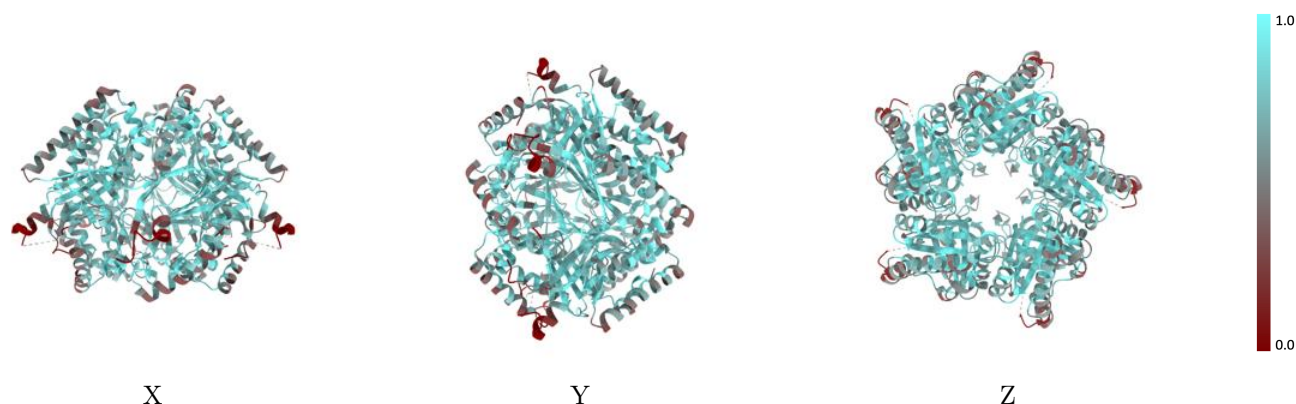
The images above show the 3D surface view of the map at the recommended contour level 0.9 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



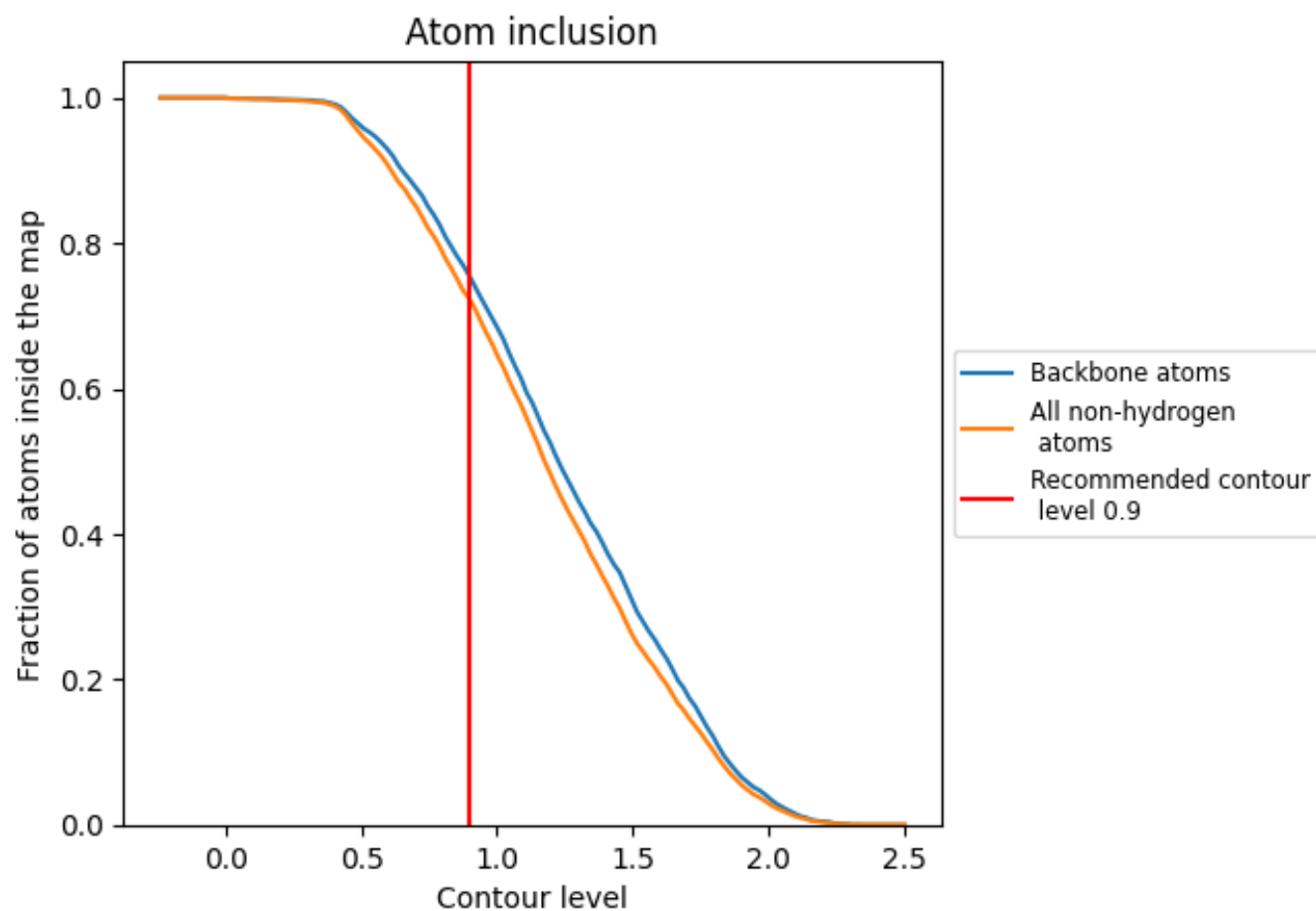
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.9).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 72% 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 (0.9) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.7222	<div></div> 0.6050
V	<div></div> 0.7291	<div></div> 0.6050
W	<div></div> 0.7296	<div></div> 0.6060
X	<div></div> 0.7322	<div></div> 0.6040
Y	<div></div> 0.7285	<div></div> 0.6040
Z	<div></div> 0.7328	<div></div> 0.6040

