



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 01:46 AM UTC

PDB ID : 3IY2 / pdb_00003iy2
EMDB ID : EMD-5107
Title : Variable domains of the computer generated model (WAM) of Fab 6 fitted into the cryoEM reconstruction of the virus-Fab 6 complex
Authors : Hafenstein, S.; Bowman, V.D.; Sun, T.; Nelson, C.D.; Palermo, L.M.; Chipman, P.R.; Battisti, A.J.; Parrish, C.R.; Rossmann, M.G.
Deposited on : 2009-04-09
Resolution : 18.00 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
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

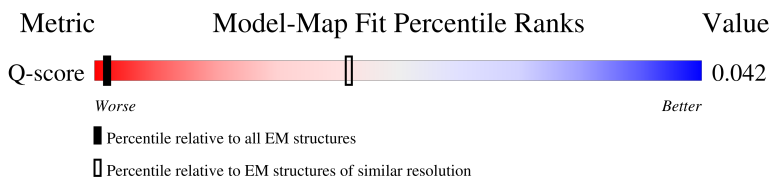
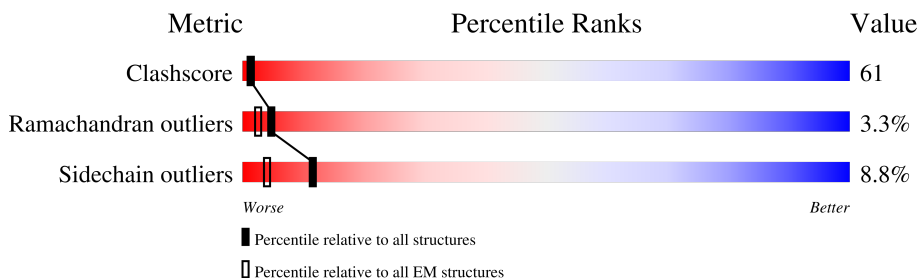
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 18.00 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	24 (17.50 - 18.30)

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

Mol	Chain	Length	Quality of chain
1	A	107	
2	B	111	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1690 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 Antibody 6, light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	107	Total	C	N	O	S	0	0
			827	524	141	159	3		

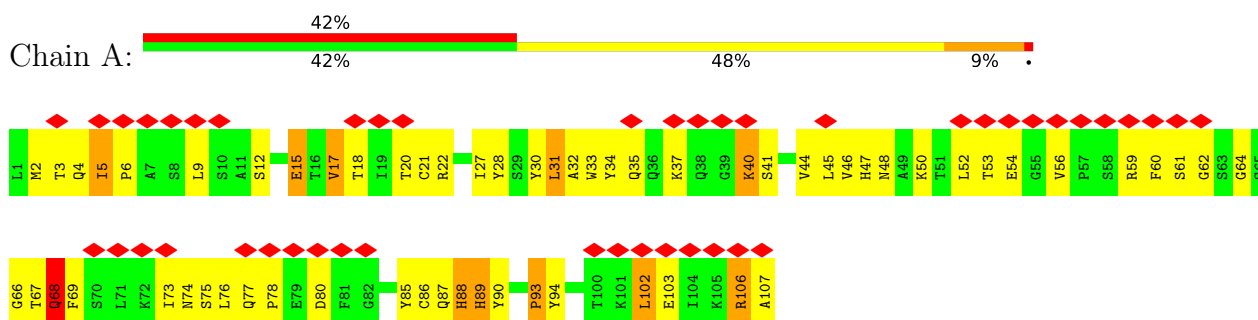
- Molecule 2 is a protein called Antibody 6, heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	111	Total	C	N	O	S	0	0
			863	550	141	167	5		

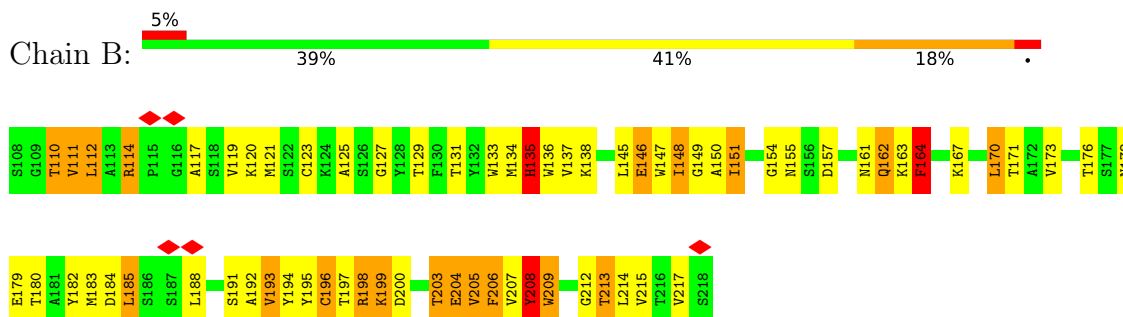
3 Residue-property plots

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: Antibody 6, light chain



- Molecule 2: Antibody 6, heavy chain



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	2520	Depositor
Resolution determination method	FSC 0.5 CUT-OFF	Depositor
CTF correction method	robem	Depositor
Microscope	FEI/PHILIPS CM300FEG/T	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	37	Depositor
Minimum defocus (nm)	1.2	Depositor
Maximum defocus (nm)	4.2	Depositor
Magnification	47190	Depositor
Image detector	KODAK SO-163 FILM	Depositor
Maximum map value	6.618	Depositor
Minimum map value	-2.780	Depositor
Average map value	0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	1	Depositor
Map size ($\text{\AA}$)	531.76, 531.76, 531.76	wwPDB
Map dimensions	184, 184, 184	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	2.89, 2.89, 2.89	Depositor

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.33	5/847 (0.6%)	1.97	17/1146 (1.5%)
2	B	1.14	0/886	2.12	33/1201 (2.7%)
All	All	1.24	5/1733 (0.3%)	2.05	50/2347 (2.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	B	1	4
All	All	1	5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	67	THR	C-O	-17.25	1.02	1.24
1	A	67	THR	C-N	12.86	1.50	1.33
1	A	68	GLN	N-CA	6.46	1.53	1.46
1	A	68	GLN	CD-OE1	5.99	1.34	1.23
1	A	68	GLN	CD-NE2	-5.67	1.21	1.33

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	68	GLN	OE1-CD-NE2	-19.35	103.25	122.60
2	B	164	PHE	CD1-CE1-CZ	-18.66	86.40	120.00
2	B	135	HIS	CB-CG-CD2	-14.83	111.92	131.20
2	B	164	PHE	CZ-CE2-CD2	-14.50	93.89	120.00
1	A	88	HIS	ND1-CE1-NE2	12.41	120.81	108.40
2	B	170	LEU	N-CA-CB	-12.31	90.34	110.41
1	A	68	GLN	CG-CD-NE2	12.25	134.78	116.40

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	135	HIS	ND1-CE1-NE2	10.96	119.36	108.40
1	A	89	HIS	ND1-CE1-NE2	10.91	119.31	108.40
1	A	47	HIS	ND1-CE1-NE2	10.43	118.83	108.40
1	A	93	PRO	CB-CA-C	-10.35	86.14	112.00
1	A	88	HIS	CE1-NE2-CD2	-9.93	99.07	109.00
2	B	135	HIS	CA-CB-CG	-9.48	104.32	113.80
1	A	89	HIS	CE1-NE2-CD2	-9.41	99.59	109.00
2	B	146	GLU	CG-CD-OE1	-9.31	96.98	118.40
2	B	205	VAL	CA-CB-CG2	-9.10	94.93	110.40
1	A	47	HIS	CE1-NE2-CD2	-9.04	99.96	109.00
2	B	135	HIS	N-CA-CB	-9.02	94.96	111.37
1	A	67	THR	CA-C-N	-8.75	110.66	122.99
1	A	67	THR	C-N-CA	-8.75	110.66	122.99
1	A	67	THR	O-C-N	8.53	132.17	122.27
2	B	135	HIS	CE1-NE2-CD2	-8.47	100.53	109.00
1	A	17	VAL	CA-CB-CG2	-7.87	97.03	110.40
2	B	146	GLU	CG-CD-OE2	7.73	136.18	118.40
2	B	170	LEU	CB-CA-C	-7.06	97.44	109.51
1	A	54	GLU	N-CA-CB	-6.81	98.70	109.19
2	B	151	ILE	CB-CA-C	-6.80	100.03	110.50
2	B	135	HIS	CG-ND1-CE1	-6.78	97.78	109.30
2	B	164	PHE	CG-CD2-CE2	-6.42	109.79	120.70
2	B	146	GLU	CB-CG-CD	-6.17	102.11	112.60
2	B	205	VAL	N-CA-CB	-6.16	101.06	111.23
1	A	89	HIS	CA-C-O	6.05	128.56	121.78
2	B	127	GLY	N-CA-C	-6.05	98.83	113.18
2	B	135	HIS	ND1-CG-CD2	6.05	112.15	106.10
2	B	151	ILE	N-CA-CB	5.89	121.10	111.44
1	A	88	HIS	CG-ND1-CE1	-5.88	99.31	109.30
2	B	164	PHE	CD1-CG-CD2	5.80	127.30	118.60
2	B	151	ILE	CA-CB-CG2	5.72	120.23	110.50
2	B	110	THR	O-C-N	5.61	130.05	122.59
2	B	203	THR	CA-CB-CG2	5.61	120.03	110.50
2	B	148	ILE	CG1-CB-CG2	-5.37	94.60	110.70
2	B	154	GLY	N-CA-C	-5.31	107.46	115.32
2	B	208	TYR	N-CA-C	5.29	122.06	110.80
1	A	17	VAL	CB-CA-C	5.24	119.15	110.30
2	B	129	THR	N-CA-C	-5.19	106.80	113.23
2	B	162	GLN	O-C-N	5.16	127.59	122.12
2	B	164	PHE	CE1-CZ-CE2	5.15	129.26	120.00
2	B	110	THR	OG1-CB-CG2	5.12	119.54	109.30
2	B	199	LYS	N-CA-C	-5.10	99.10	108.02

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	110	THR	CA-C-O	-5.07	113.26	120.51

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	208	TYR	CA

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	68	GLN	Sidechain
2	B	135	HIS	Sidechain
2	B	164	PHE	Sidechain
2	B	206	PHE	Sidechain
2	B	208	TYR	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	A	827	0	816	68	0
2	B	863	0	829	150	0
All	All	1690	0	1645	200	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 61.

All (200) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:173:VAL:HG11	2:B:176:THR:HG22	1.29	1.12
2:B:121:MET:HE3	2:B:183:MET:SD	1.89	1.11
1:A:9:LEU:HD21	1:A:17:VAL:HG21	1.35	1.09
1:A:9:LEU:HD21	1:A:17:VAL:CG2	1.86	1.04
1:A:44:VAL:HG13	1:A:53:THR:HG21	1.40	1.02
2:B:195:TYR:HB3	2:B:208:TYR:HE2	1.22	1.01

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:GLN:HB3	1:A:78:PRO:HD2	1.43	1.00
2:B:197:THR:CG2	2:B:208:TYR:HA	1.94	0.97
2:B:173:VAL:CG1	2:B:176:THR:HG22	1.96	0.95
2:B:151:ILE:HB	2:B:170:LEU:HD23	1.50	0.93
2:B:173:VAL:CG1	2:B:178:ASN:HB2	1.97	0.93
2:B:173:VAL:HG11	2:B:176:THR:CG2	1.99	0.92
1:A:31:LEU:HD22	1:A:88:HIS:HA	1.51	0.92
2:B:176:THR:HG23	2:B:178:ASN:HD22	1.34	0.91
2:B:111:VAL:HG22	2:B:112:LEU:H	1.36	0.91
2:B:163:LYS:HG2	2:B:164:PHE:HZ	1.34	0.90
2:B:197:THR:HG22	2:B:208:TYR:HA	1.52	0.90
1:A:32:ALA:HB3	2:B:205:VAL:HG21	1.53	0.90
2:B:195:TYR:HB3	2:B:208:TYR:CE2	2.08	0.88
2:B:151:ILE:HB	2:B:170:LEU:CD2	2.02	0.88
1:A:44:VAL:CG1	1:A:53:THR:HG21	2.04	0.87
2:B:120:LYS:HE3	2:B:180:THR:HG21	1.57	0.86
1:A:9:LEU:CD2	1:A:17:VAL:HG21	2.07	0.83
2:B:121:MET:HE2	2:B:215:VAL:HG21	1.61	0.82
2:B:163:LYS:HE2	2:B:164:PHE:CZ	2.15	0.82
2:B:119:VAL:HG12	2:B:183:MET:HB2	1.64	0.80
1:A:5:ILE:HD12	1:A:6:PRO:HA	1.63	0.80
2:B:173:VAL:HG11	2:B:178:ASN:HB2	1.63	0.79
1:A:93:PRO:HB3	2:B:147:TRP:CE3	2.18	0.79
2:B:123:CYS:HB3	2:B:179:GLU:HB3	1.64	0.79
2:B:133:TRP:HB2	2:B:199:LYS:HB3	1.64	0.78
2:B:197:THR:HG21	2:B:208:TYR:HA	1.67	0.77
2:B:173:VAL:HG13	2:B:173:VAL:O	1.86	0.76
2:B:136:TRP:HD1	2:B:170:LEU:CD1	1.99	0.75
1:A:37:LYS:HB2	1:A:40:LYS:HE2	1.69	0.75
2:B:136:TRP:O	2:B:148:ILE:HG22	1.87	0.73
1:A:32:ALA:CB	2:B:205:VAL:HG21	2.17	0.73
1:A:89:HIS:ND1	2:B:204:GLU:HA	2.02	0.73
2:B:121:MET:HE2	2:B:215:VAL:CG2	2.19	0.72
2:B:151:ILE:CD1	2:B:171:THR:HA	2.19	0.72
2:B:136:TRP:HB3	2:B:148:ILE:CG2	2.20	0.71
2:B:184:ASP:HB3	2:B:185:LEU:HD12	1.71	0.71
2:B:173:VAL:HG12	2:B:178:ASN:HB2	1.73	0.70
1:A:89:HIS:HB3	2:B:204:GLU:HB3	1.72	0.69
2:B:121:MET:CE	2:B:215:VAL:HG21	2.21	0.69
1:A:9:LEU:HD21	1:A:17:VAL:HG22	1.74	0.69
1:A:93:PRO:HB3	2:B:147:TRP:CZ3	2.28	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:196:CYS:O	2:B:197:THR:HG23	1.93	0.69
2:B:163:LYS:HG2	2:B:164:PHE:CZ	2.23	0.68
2:B:197:THR:HA	2:B:209:TRP:NE1	2.09	0.68
2:B:136:TRP:C	2:B:148:ILE:HG22	2.18	0.67
2:B:121:MET:CE	2:B:183:MET:SD	2.76	0.67
2:B:138:LYS:CE	2:B:163:LYS:HZ3	2.07	0.67
1:A:2:MET:HG2	1:A:86:CYS:SG	2.35	0.67
1:A:48:ASN:H	2:B:204:GLU:HG3	1.60	0.66
1:A:87:GLN:HB3	2:B:205:VAL:HG11	1.75	0.66
2:B:131:THR:HB	2:B:198:ARG:HB3	1.77	0.65
2:B:197:THR:HG21	2:B:208:TYR:CG	2.31	0.65
1:A:31:LEU:CD2	1:A:88:HIS:HA	2.26	0.65
1:A:3:THR:HB	1:A:22:ARG:HG2	1.77	0.65
2:B:151:ILE:N	2:B:170:LEU:HD21	2.12	0.65
1:A:59:ARG:HB2	1:A:74:ASN:OD1	1.96	0.65
2:B:176:THR:HG23	2:B:178:ASN:ND2	2.09	0.64
1:A:17:VAL:HG22	1:A:18:THR:N	2.13	0.63
2:B:110:THR:O	2:B:111:VAL:HG12	1.99	0.63
2:B:123:CYS:HG	2:B:196:CYS:CB	2.11	0.63
2:B:137:VAL:HB	2:B:208:TYR:OH	1.98	0.63
2:B:110:THR:HG21	2:B:213:THR:O	1.99	0.62
2:B:110:THR:HB	2:B:214:LEU:O	1.99	0.62
2:B:200:ASP:HB2	2:B:203:THR:OG1	1.98	0.62
1:A:27:ILE:HD12	1:A:31:LEU:HD23	1.81	0.62
2:B:119:VAL:CG1	2:B:183:MET:SD	2.88	0.62
2:B:135:HIS:CE1	2:B:137:VAL:HG23	2.35	0.61
2:B:135:HIS:CE1	2:B:147:TRP:HD1	2.19	0.61
2:B:135:HIS:CE1	2:B:137:VAL:CG2	2.84	0.60
2:B:195:TYR:CB	2:B:208:TYR:CE2	2.83	0.60
1:A:106:ARG:HG2	1:A:107:ALA:N	2.17	0.59
2:B:110:THR:HG21	2:B:213:THR:C	2.28	0.59
2:B:138:LYS:CE	2:B:163:LYS:NZ	2.65	0.59
2:B:110:THR:HG22	2:B:215:VAL:CG2	2.32	0.59
2:B:184:ASP:HB3	2:B:185:LEU:CD1	2.32	0.59
1:A:85:TYR:CD1	2:B:145:LEU:HD12	2.38	0.59
2:B:136:TRP:CB	2:B:148:ILE:CG2	2.81	0.59
2:B:110:THR:HG22	2:B:215:VAL:HG23	1.85	0.58
1:A:12:SER:O	1:A:15:GLU:HB2	2.02	0.58
2:B:138:LYS:HE2	2:B:163:LYS:HZ3	1.67	0.58
2:B:151:ILE:HD12	2:B:171:THR:HA	1.84	0.58
1:A:5:ILE:HD12	1:A:6:PRO:CA	2.34	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:ARG:HA	1:A:68:GLN:OE1	2.04	0.57
1:A:27:ILE:CD1	1:A:31:LEU:HD23	2.34	0.57
2:B:135:HIS:ND1	2:B:137:VAL:HG23	2.20	0.57
2:B:161:ASN:OD1	2:B:163:LYS:HB3	2.05	0.57
2:B:136:TRP:HD1	2:B:170:LEU:HD12	1.70	0.56
2:B:131:THR:CB	2:B:198:ARG:HB3	2.36	0.56
2:B:136:TRP:CB	2:B:148:ILE:HG23	2.35	0.55
2:B:197:THR:HA	2:B:209:TRP:CD1	2.40	0.55
1:A:50:LYS:HG2	1:A:62:GLY:O	2.06	0.55
2:B:135:HIS:NE2	2:B:147:TRP:CD1	2.75	0.55
2:B:136:TRP:CD1	2:B:170:LEU:CD1	2.86	0.55
2:B:111:VAL:HG22	2:B:112:LEU:N	2.14	0.55
2:B:197:THR:HA	2:B:209:TRP:HE1	1.71	0.55
1:A:45:LEU:HA	1:A:56:VAL:HG21	1.89	0.55
2:B:193:VAL:HG11	2:B:195:TYR:CZ	2.42	0.55
2:B:114:ARG:H	2:B:114:ARG:HD3	1.73	0.54
2:B:136:TRP:CD1	2:B:170:LEU:HD12	2.43	0.54
1:A:89:HIS:CG	2:B:204:GLU:HA	2.43	0.54
1:A:87:GLN:OE1	2:B:205:VAL:HB	2.09	0.53
2:B:136:TRP:HD1	2:B:170:LEU:HD11	1.74	0.53
2:B:136:TRP:HB3	2:B:148:ILE:HG21	1.90	0.52
2:B:125:ALA:HA	2:B:209:TRP:CZ3	2.45	0.52
1:A:76:LEU:CD1	1:A:80:ASP:HB2	2.39	0.52
2:B:114:ARG:H	2:B:114:ARG:CD	2.23	0.52
1:A:9:LEU:CG	1:A:17:VAL:HG21	2.39	0.52
1:A:17:VAL:HG22	1:A:18:THR:H	1.74	0.52
1:A:27:ILE:HG22	1:A:90:TYR:CD2	2.46	0.51
2:B:197:THR:HG23	2:B:208:TYR:CE2	2.46	0.51
2:B:213:THR:O	2:B:213:THR:HG23	2.11	0.51
1:A:34:TYR:OH	2:B:205:VAL:HG23	2.10	0.51
2:B:163:LYS:CE	2:B:164:PHE:CZ	2.91	0.50
2:B:146:GLU:CD	2:B:163:LYS:HZ2	2.19	0.50
2:B:173:VAL:HG12	2:B:178:ASN:CB	2.39	0.50
2:B:150:ALA:C	2:B:170:LEU:HD21	2.37	0.50
2:B:191:SER:O	2:B:192:ALA:HB2	2.11	0.50
1:A:77:GLN:HB3	1:A:78:PRO:CD	2.27	0.50
2:B:151:ILE:HD13	2:B:171:THR:HA	1.93	0.50
2:B:199:LYS:HB2	2:B:206:PHE:HD1	1.76	0.50
2:B:136:TRP:HB2	2:B:148:ILE:HG23	1.93	0.49
2:B:120:LYS:CE	2:B:180:THR:HG21	2.38	0.49
2:B:131:THR:CG2	2:B:198:ARG:HB3	2.42	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:151:ILE:HG13	2:B:157:ASP:O	2.13	0.49
1:A:68:GLN:OE1	1:A:68:GLN:HA	2.12	0.49
1:A:32:ALA:HB3	2:B:205:VAL:CG2	2.36	0.49
2:B:135:HIS:O	2:B:196:CYS:HA	2.12	0.49
1:A:59:ARG:HD3	1:A:75:SER:O	2.13	0.48
2:B:199:LYS:HB2	2:B:206:PHE:CD1	2.49	0.48
2:B:131:THR:HG21	2:B:198:ARG:HG2	1.96	0.47
2:B:197:THR:CG2	2:B:208:TYR:CD2	2.97	0.47
1:A:33:TRP:CZ3	1:A:86:CYS:HB3	2.49	0.47
1:A:102:LEU:HD12	1:A:103:GLU:N	2.29	0.47
2:B:163:LYS:HD3	2:B:164:PHE:CE1	2.50	0.47
1:A:64:GLY:HA3	1:A:69:PHE:HA	1.96	0.47
2:B:197:THR:HG21	2:B:208:TYR:CA	2.39	0.47
1:A:106:ARG:HG2	1:A:107:ALA:H	1.80	0.47
1:A:15:GLU:HA	1:A:15:GLU:OE1	2.14	0.47
2:B:114:ARG:HE	2:B:117:ALA:HB2	1.80	0.47
1:A:68:GLN:OE1	1:A:68:GLN:CA	2.63	0.46
2:B:200:ASP:HB2	2:B:203:THR:CB	2.44	0.46
2:B:162:GLN:OE1	2:B:162:GLN:HA	2.14	0.46
2:B:135:HIS:CE1	2:B:137:VAL:HG22	2.50	0.46
2:B:184:ASP:CB	2:B:185:LEU:CD1	2.93	0.46
2:B:184:ASP:CB	2:B:185:LEU:HD12	2.41	0.46
2:B:188:LEU:HD12	2:B:217:VAL:O	2.15	0.46
2:B:197:THR:HG21	2:B:208:TYR:CD1	2.50	0.46
1:A:94:TYR:CD2	2:B:206:PHE:CE2	3.04	0.46
2:B:205:VAL:H	2:B:205:VAL:HG22	0.88	0.46
1:A:32:ALA:CB	2:B:205:VAL:CG2	2.91	0.45
1:A:94:TYR:HD2	2:B:206:PHE:CE2	2.33	0.45
2:B:212:GLY:O	2:B:213:THR:HB	2.17	0.45
2:B:173:VAL:CG1	2:B:173:VAL:O	2.61	0.45
2:B:167:LYS:HB3	2:B:167:LYS:HE2	1.86	0.45
2:B:135:HIS:HE1	2:B:137:VAL:HG22	1.82	0.45
2:B:135:HIS:NE2	2:B:147:TRP:HD1	2.13	0.45
2:B:138:LYS:NZ	2:B:164:PHE:CE2	2.81	0.45
2:B:125:ALA:HA	2:B:209:TRP:HZ3	1.81	0.44
1:A:35:GLN:HB2	1:A:45:LEU:HD11	1.98	0.44
2:B:123:CYS:HB2	2:B:136:TRP:CH2	2.52	0.44
2:B:125:ALA:HB2	2:B:209:TRP:CH2	2.53	0.44
2:B:198:ARG:O	2:B:207:VAL:HG22	2.18	0.44
2:B:110:THR:O	2:B:111:VAL:CG1	2.66	0.43
2:B:193:VAL:CG1	2:B:195:TYR:CE2	3.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:148:ILE:HG21	2:B:148:ILE:HD13	1.61	0.43
2:B:173:VAL:HG22	2:B:176:THR:H	1.83	0.43
2:B:196:CYS:O	2:B:197:THR:CG2	2.65	0.43
2:B:196:CYS:C	2:B:197:THR:HG23	2.43	0.43
1:A:76:LEU:HD11	1:A:80:ASP:HB2	1.99	0.43
2:B:135:HIS:CD2	2:B:149:GLY:O	2.71	0.43
1:A:28:TYR:HB2	1:A:30:TYR:CE2	2.53	0.43
1:A:59:ARG:O	1:A:73:ILE:HA	2.19	0.43
2:B:123:CYS:SG	2:B:196:CYS:CB	3.04	0.43
1:A:4:GLN:HG3	1:A:21:CYS:SG	2.59	0.43
2:B:193:VAL:HG23	2:B:214:LEU:HD13	1.99	0.43
1:A:94:TYR:CD2	2:B:206:PHE:CZ	3.07	0.43
1:A:22:ARG:HB2	1:A:68:GLN:OE1	2.18	0.42
2:B:148:ILE:HD12	2:B:194:TYR:HE2	1.83	0.42
2:B:185:LEU:HD12	2:B:185:LEU:N	2.35	0.42
2:B:182:TYR:CE2	2:B:184:ASP:OD1	2.73	0.41
1:A:20:THR:HG23	1:A:69:PHE:O	2.20	0.41
2:B:199:LYS:HD2	2:B:200:ASP:O	2.20	0.41
1:A:59:ARG:HA	1:A:74:ASN:OD1	2.21	0.41
2:B:131:THR:HG21	2:B:198:ARG:HB3	2.01	0.41
2:B:173:VAL:CG2	2:B:176:THR:HG22	2.51	0.41
1:A:44:VAL:HG22	1:A:53:THR:OG1	2.21	0.41
1:A:85:TYR:CE1	2:B:145:LEU:HD12	2.55	0.41
2:B:134:MET:SD	2:B:179:GLU:OE1	2.79	0.40
1:A:22:ARG:CB	1:A:68:GLN:OE1	2.69	0.40
1:A:76:LEU:HD12	1:A:80:ASP:HB2	2.02	0.40
1:A:102:LEU:HD12	1:A:102:LEU:C	2.46	0.40
1:A:15:GLU:OE1	1:A:15:GLU:CA	2.69	0.40
1:A:52:LEU:HD21	1:A:60:PHE:O	2.20	0.40
2:B:136:TRP:CD1	2:B:170:LEU:HD11	2.53	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	A	105/107 (98%)	98 (93%)	6 (6%)	1 (1%)	12	49
2	B	109/111 (98%)	93 (85%)	10 (9%)	6 (6%)	1	15
All	All	214/218 (98%)	191 (89%)	16 (8%)	7 (3%)	5	21

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	111	VAL
2	B	208	TYR
2	B	213	THR
2	B	155	ASN
2	B	209	TRP
2	B	204	GLU
1	A	66	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	90/90 (100%)	80 (89%)	10 (11%)	6	20
2	B	92/92 (100%)	86 (94%)	6 (6%)	15	37
All	All	182/182 (100%)	166 (91%)	16 (9%)	11	28

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ILE
1	A	15	GLU
1	A	31	LEU
1	A	40	LYS
1	A	41	SER
1	A	46	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	61	SER
1	A	68	GLN
1	A	102	LEU
1	A	106	ARG
2	B	112	LEU
2	B	114	ARG
2	B	185	LEU
2	B	193	VAL
2	B	196	CYS
2	B	198	ARG

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

Mol	Chain	Res	Type
1	A	43	GLN
2	B	143	GLN
2	B	178	ASN

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](#)

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 ⓘ

There are no chain breaks in this entry.

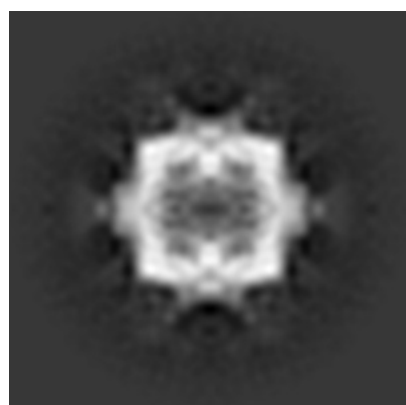
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-5107. These allow visual inspection of the internal detail of the map and identification of artifacts.

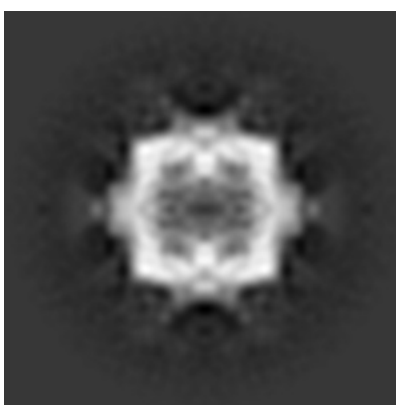
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

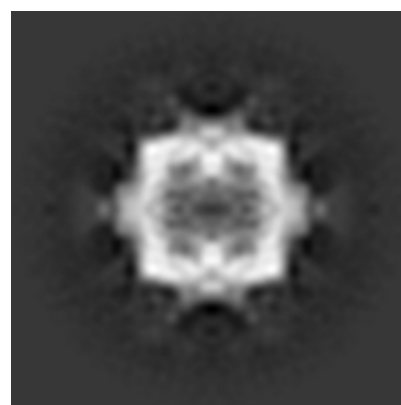
6.1.1 Primary map



X



Y

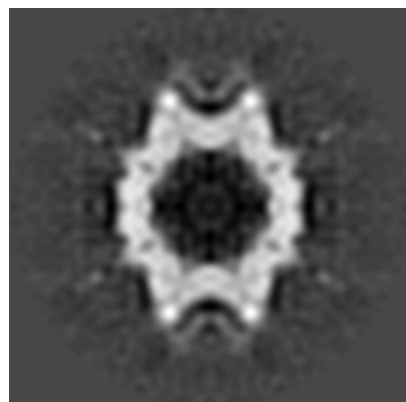


Z

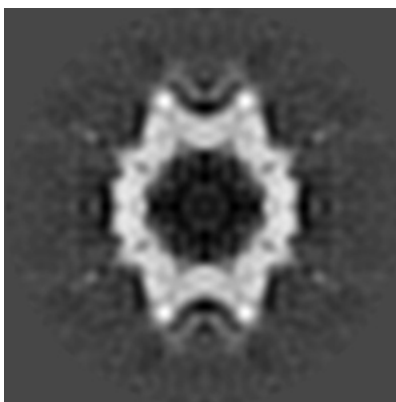
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

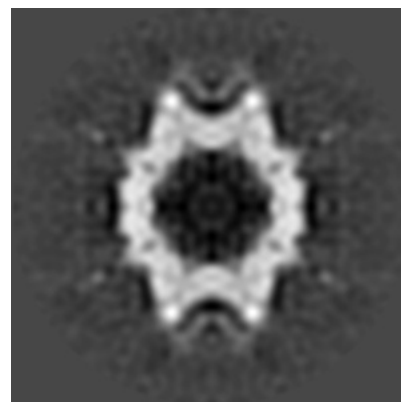
6.2.1 Primary map



X Index: 92



Y Index: 92

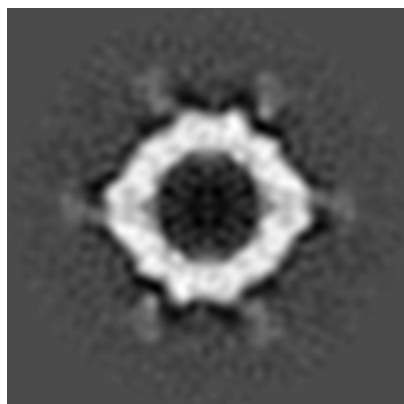


Z Index: 92

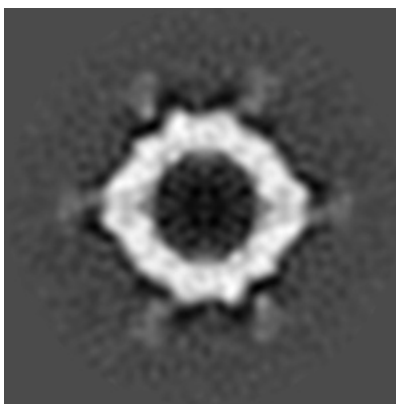
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

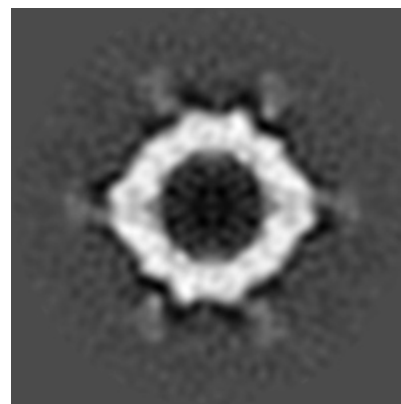
6.3.1 Primary map



X Index: 81



Y Index: 103

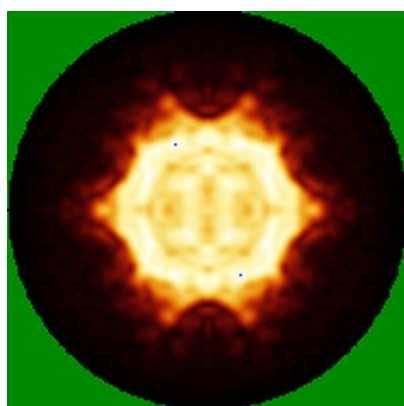


Z Index: 81

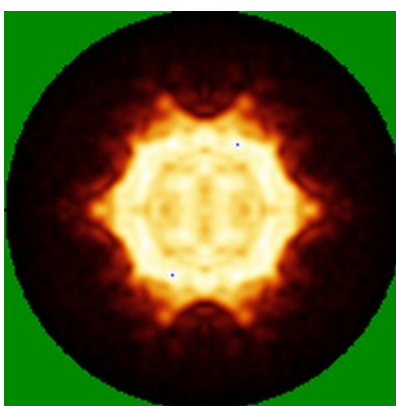
The images above show the largest variance slices of the map in three orthogonal directions.

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

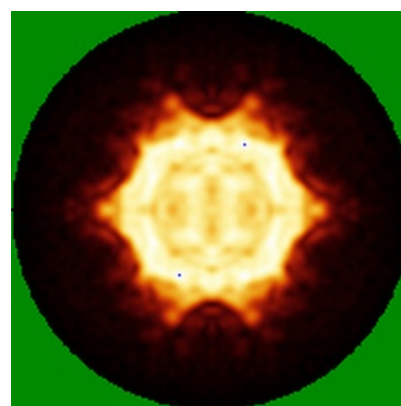
6.4.1 Primary map



X



Y

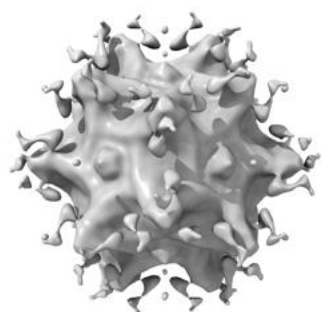


Z

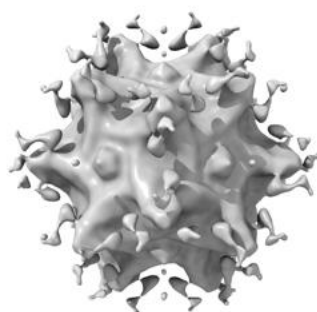
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

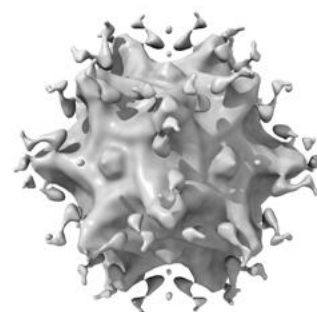
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 1.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

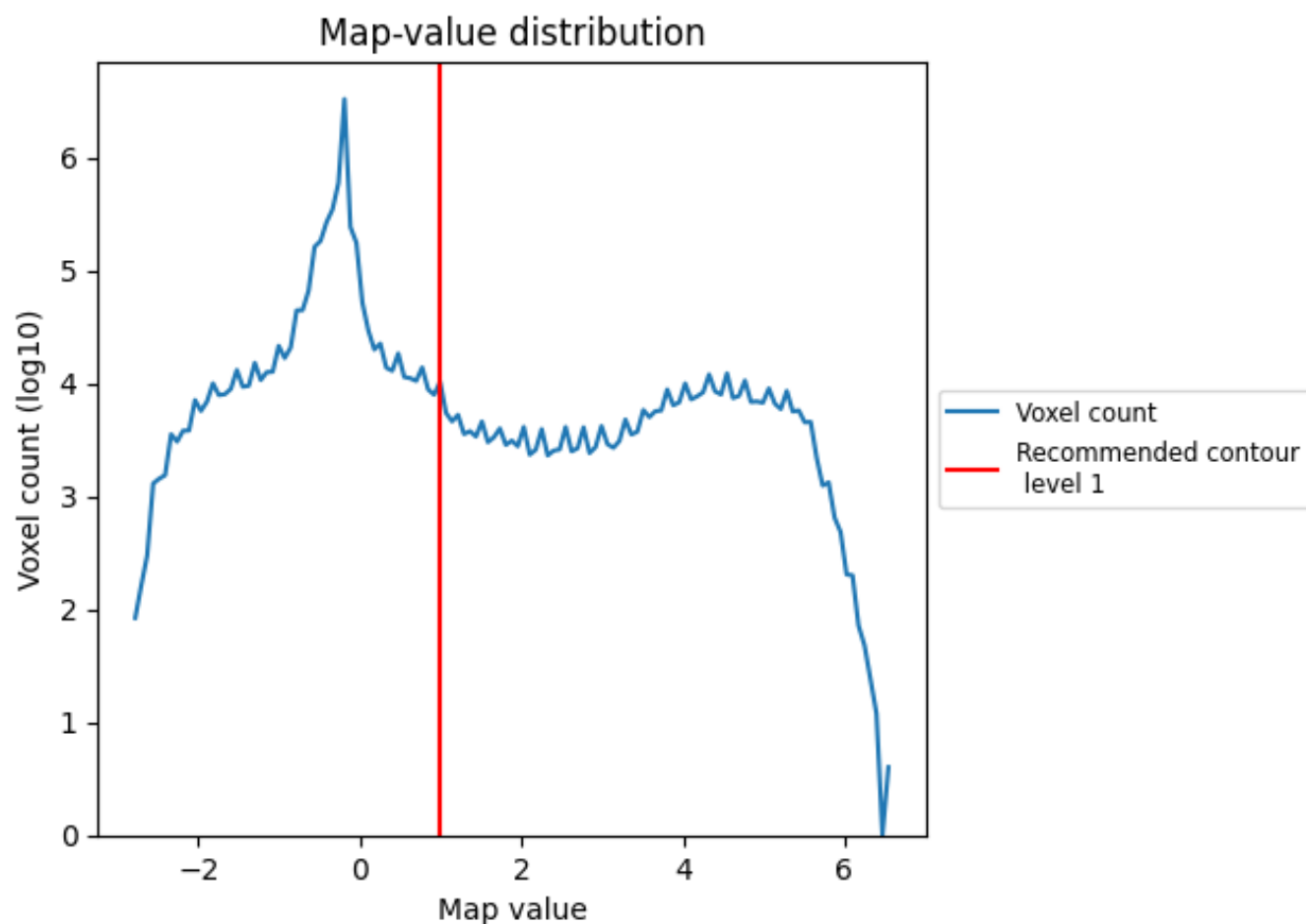
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

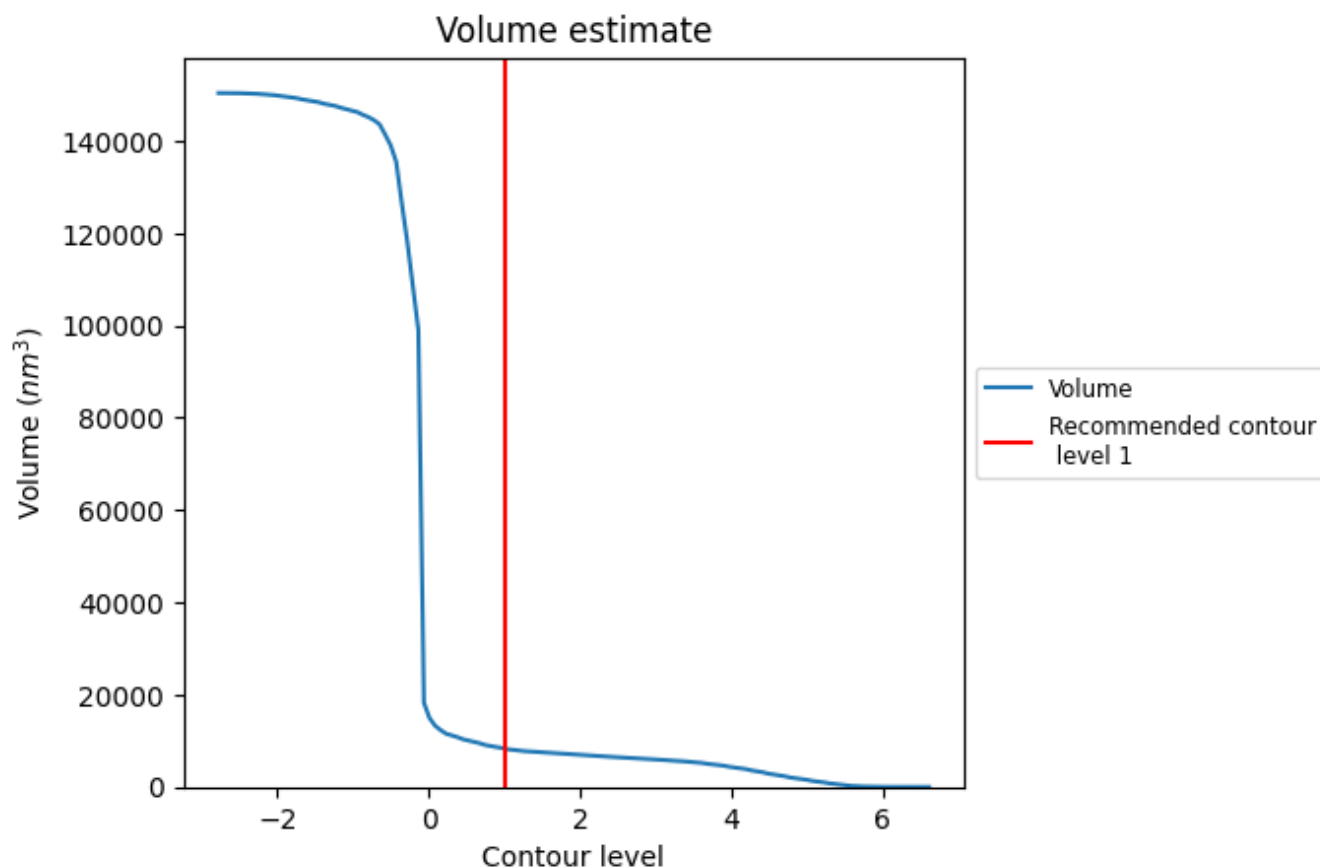
This section contains the results of statistical analysis of the map.

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.

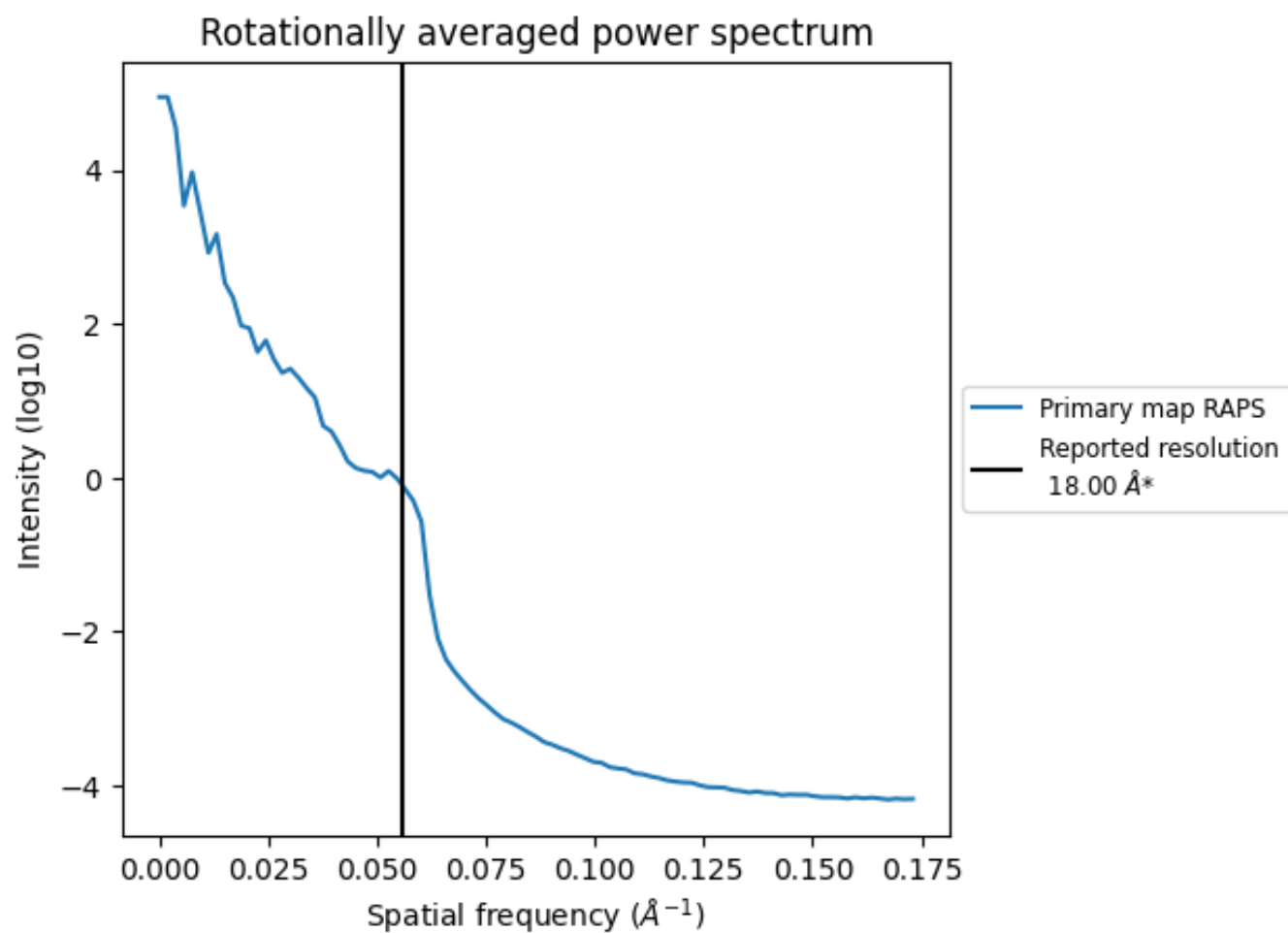
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 8324 nm³; this corresponds to an approximate mass of 7519 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 0.056 Å⁻¹

8 Fourier-Shell correlation ⓘ

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

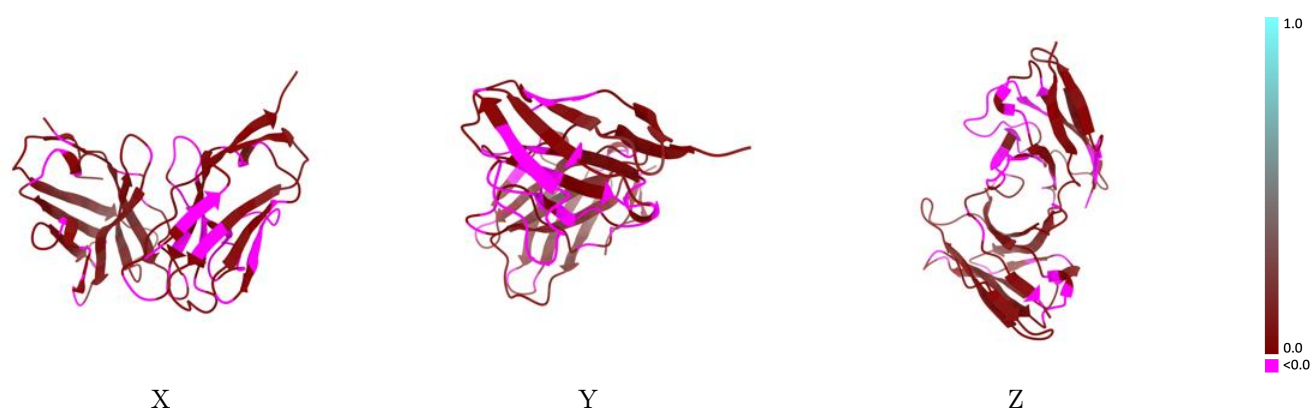
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-5107 and PDB model 3IY2. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlay [i](#)

This section was not generated.

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

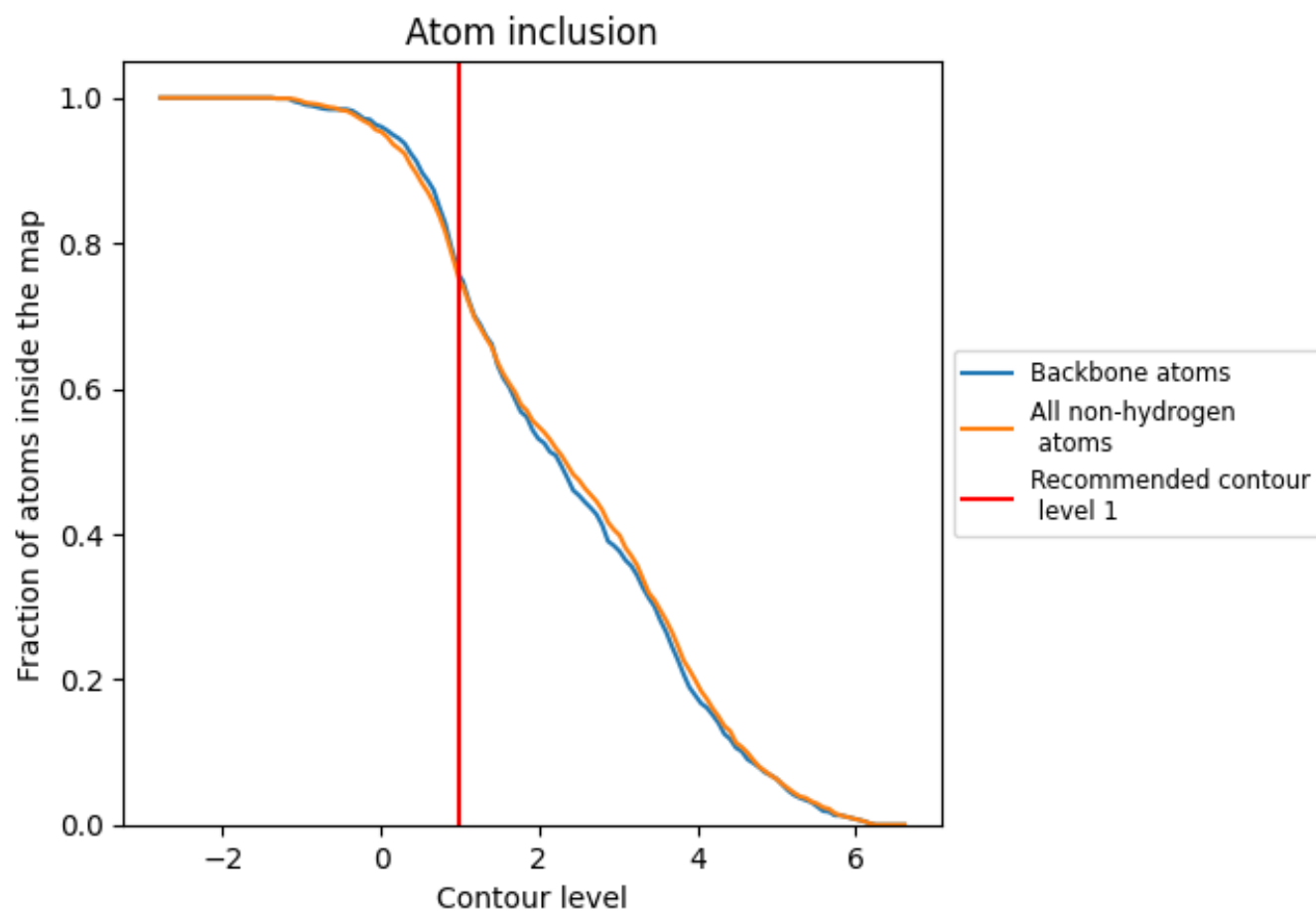


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 75% 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 (1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7470	0.0420
A	0.5500	0.0180
B	0.9360	0.0650

