



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 11:13 AM UTC

PDB ID : 7LKH / pdb_00007lkh
EMDB ID : EMD-23408
Title : Chicken Scap D435V L1-L7 domain / Fab complex focused map
Authors : Kober, D.L.; Radhakrishnan, A.; Goldstein, J.L.; Brown, M.S.; Clark, L.D.;
Bai, X.-C.; Rosenbaum, D.M.
Deposited on : 2021-02-02
Resolution : 3.50 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
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

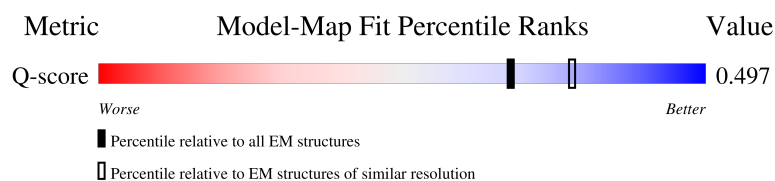
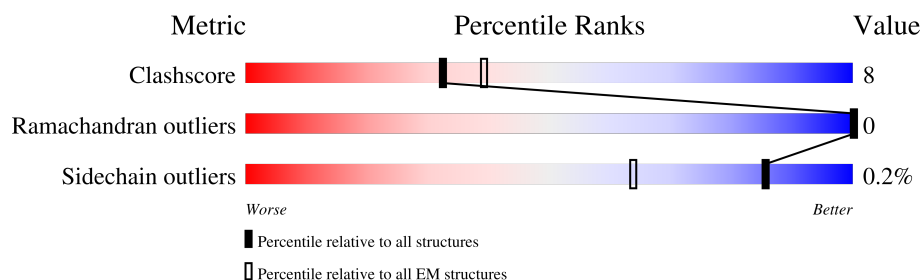
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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	13950 (3.00 - 4.00)

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	1346	
2	H	231	
3	L	214	
4	B	2	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4179 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 Sterol regulatory element-binding protein cleavage-activating protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	286	2338	1509	408	418	3	0	0

There are 31 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	PRO	-	expression tag	UNP A0A3Q3ANV4
A	0	ALA	-	expression tag	UNP A0A3Q3ANV4
A	435	VAL	ASP	engineered mutation	UNP A0A3Q3ANV4
A	1324	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1325	TYR	-	expression tag	UNP A0A3Q3ANV4
A	1326	LYS	-	expression tag	UNP A0A3Q3ANV4
A	1327	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1328	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1329	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1330	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1331	LYS	-	expression tag	UNP A0A3Q3ANV4
A	1332	GLY	-	expression tag	UNP A0A3Q3ANV4
A	1333	SER	-	expression tag	UNP A0A3Q3ANV4
A	1334	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1335	TYR	-	expression tag	UNP A0A3Q3ANV4
A	1336	LYS	-	expression tag	UNP A0A3Q3ANV4
A	1337	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1338	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1339	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1340	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1341	LYS	-	expression tag	UNP A0A3Q3ANV4
A	1342	GLY	-	expression tag	UNP A0A3Q3ANV4
A	1343	SER	-	expression tag	UNP A0A3Q3ANV4
A	1344	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1345	TYR	-	expression tag	UNP A0A3Q3ANV4
A	1346	LYS	-	expression tag	UNP A0A3Q3ANV4
A	1347	ASP	-	expression tag	UNP A0A3Q3ANV4

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	1348	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1349	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1350	ASP	-	expression tag	UNP A0A3Q3ANV4
A	1351	LYS	-	expression tag	UNP A0A3Q3ANV4

- Molecule 2 is a protein called 4G10 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	121	Total	C	N	O	S	0	0
			959	605	160	192	2		

- Molecule 3 is a protein called 4G10 Fab kappa chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	L	109	Total	C	N	O	S	0	0
			854	531	147	173	3		

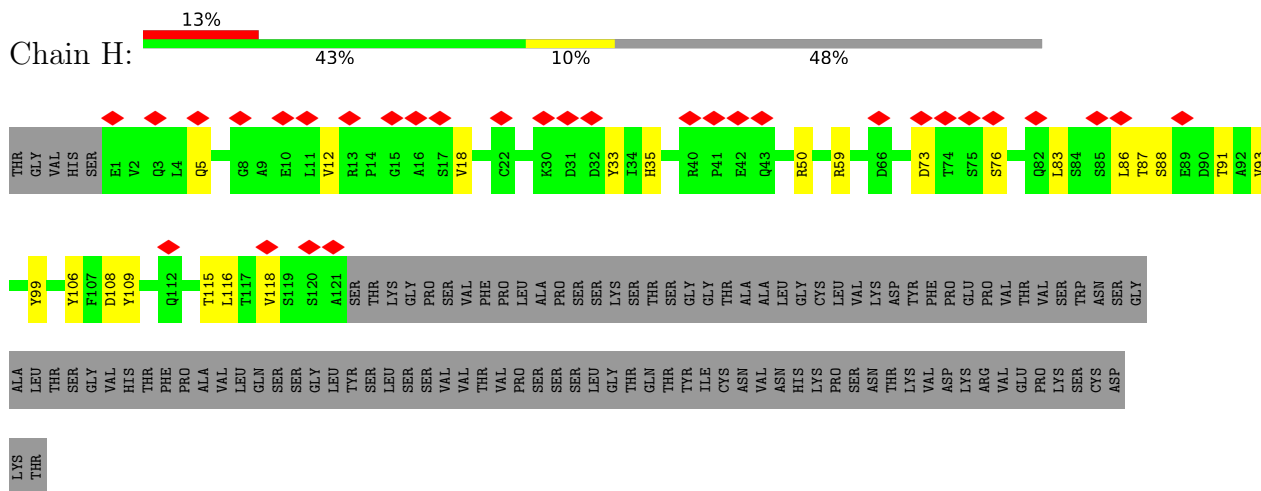
- Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



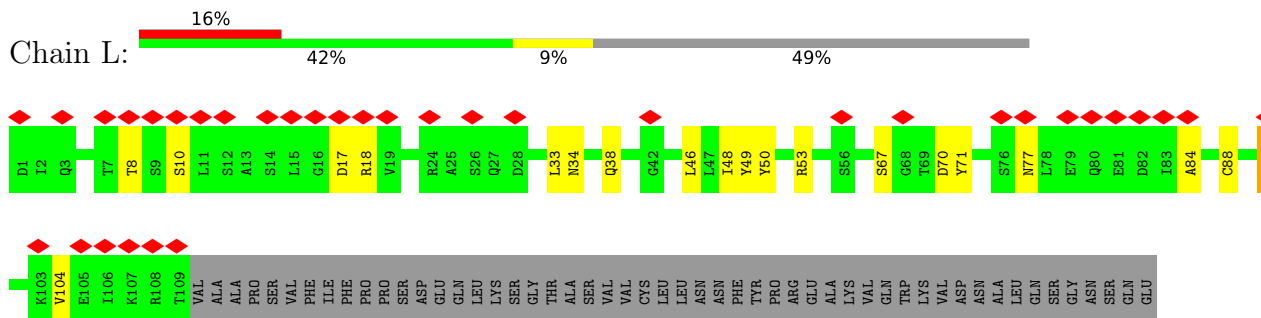
Mol	Chain	Residues	Atoms				AltConf	Trace
4	B	2	Total	C	N	O	0	0
			28	16	2	10		

[illegible]

- Molecule 2: 4G10 Fab heavy chain



- Molecule 3: 4G10 Fab kappa chain



SER	VAL	THR	GLU	GLN	ASP	SER	LYS	ASP	SER	THR	TYR	SER	LEU	SER	THR	LEU	THR	LEU	SER	LYS	ALA	ASP	TYR	GLU	LYS	HIS	LYS	VAL	TYR	ALA	CYS	GLU	VAL	THR	HIS	GLN	GLY	LEU	SER	SER	PRO	VAL	THR	LYS	SER	PHE	ASN	ARG	GLY	GLU	CYS
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

● Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	155187	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1.6	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.270	Depositor
Minimum map value	-0.310	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.005	Depositor
Recommended contour level	0.0561	Depositor
Map size (Å)	291.6, 291.6, 291.6	wwPDB
Map dimensions	270, 270, 270	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	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: NAG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.32	0/2403	0.62	0/3271
2	H	0.24	0/983	0.52	0/1338
3	L	0.24	0/870	0.59	2/1180 (0.2%)
All	All	0.29	0/4256	0.59	2/5789 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	94	LEU	CA-C-N	-5.17	113.38	119.84
3	L	94	LEU	C-N-CA	-5.17	113.38	119.84

There are no chirality outliers.

There are no planarity outliers.

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	2338	0	2334	32	0
2	H	959	0	905	19	0
3	L	854	0	830	19	0
4	B	28	0	25	3	0
All	All	4179	0	4094	70	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 8.

All (70) 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)
4:B:1:NAG:O4	4:B:2:NAG:O7	1.90	0.89
2:H:18:VAL:HG23	2:H:86:LEU:HD11	1.62	0.82
1:A:163:LEU:HD23	1:A:209:LEU:HD21	1.70	0.73
3:L:33:LEU:HD13	3:L:71:TYR:HD2	1.56	0.71
2:H:73:ASP:OD2	2:H:76:SER:OG	2.06	0.70
1:A:87:GLY:O	1:A:674:TYR:OH	2.10	0.69
2:H:12:VAL:HG11	2:H:18:VAL:HG22	1.79	0.64
3:L:46:LEU:HD21	3:L:49:TYR:HB3	1.81	0.62
1:A:696:ARG:NH2	1:A:701:ALA:HB2	2.15	0.61
2:H:88:SER:O	2:H:91:THR:HG23	2.02	0.59
2:H:118:VAL:O	2:H:118:VAL:HG23	2.03	0.59
1:A:252:ARG:HG2	1:A:256:LYS:HE2	1.86	0.57
2:H:91:THR:HG22	2:H:118:VAL:HG22	1.86	0.57
1:A:222:GLY:O	1:A:227:ASN:ND2	2.34	0.57
3:L:33:LEU:HD13	3:L:71:TYR:CD2	2.40	0.56
2:H:5:GLN:N	2:H:5:GLN:OE1	2.34	0.55
1:A:103:SER:OG	1:A:104:PRO:HD3	2.08	0.54
3:L:77:ASN:O	3:L:77:ASN:OD1	2.25	0.54
1:A:276:PHE:O	1:A:671:ALA:O	2.28	0.52
3:L:50:TYR:HB3	3:L:53:ARG:HG3	1.92	0.52
1:A:117:SER:OG	1:A:118:PRO:HD3	2.11	0.51
2:H:87:THR:O	2:H:118:VAL:HG21	2.10	0.51
1:A:703:ARG:NH1	1:A:704:TYR:OH	2.43	0.51
2:H:106:TYR:HB3	3:L:49:TYR:HB2	1.93	0.51
3:L:48:ILE:O	3:L:49:TYR:C	2.55	0.50
1:A:163:LEU:HD23	1:A:209:LEU:CD2	2.39	0.50
3:L:34:ASN:OD1	3:L:49:TYR:O	2.29	0.50
1:A:697:HIS:O	1:A:700:GLU:HG2	2.12	0.49
4:B:2:NAG:O7	4:B:2:NAG:C1	2.62	0.48
3:L:33:LEU:HD21	3:L:88:CYS:HB2	1.96	0.48
1:A:641:VAL:HG11	1:A:684:THR:CG2	2.44	0.48
1:A:65:PHE:CE2	1:A:272:VAL:HG11	2.49	0.47
1:A:252:ARG:HG2	1:A:256:LYS:CE	2.44	0.47
1:A:192:ILE:CD1	1:A:225:LEU:HD11	2.43	0.47
1:A:81:ARG:N	1:A:82:PRO:CD	2.77	0.47
3:L:33:LEU:CD1	3:L:71:TYR:HD2	2.27	0.47
3:L:48:ILE:O	3:L:48:ILE:HG22	2.14	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:70:ASP:O	3:L:71:TYR:HD1	1.97	0.47
1:A:65:PHE:CB	1:A:66:PRO:HD3	2.46	0.46
1:A:65:PHE:HB2	1:A:66:PRO:HD3	1.97	0.46
2:H:59:ARG:HD3	3:L:94:LEU:HD21	1.98	0.46
2:H:35:HIS:CE1	2:H:99:TYR:CE2	3.03	0.46
2:H:33:TYR:HD2	2:H:99:TYR:HH	1.64	0.46
1:A:278:GLU:OE1	1:A:278:GLU:C	2.59	0.46
3:L:38:GLN:O	3:L:84:ALA:HB1	2.15	0.46
1:A:157:LEU:O	1:A:157:LEU:HG	2.16	0.45
2:H:93:VAL:HG22	2:H:115:THR:HG22	2.00	0.44
1:A:65:PHE:CZ	1:A:272:VAL:HG11	2.53	0.44
1:A:97:PHE:HE1	1:A:216:LEU:HD21	1.83	0.44
1:A:107:LYS:O	1:A:108:ASN:HB3	2.19	0.43
2:H:83:LEU:HB3	2:H:86:LEU:HD21	2.01	0.43
2:H:108:ASP:OD1	2:H:109:TYR:N	2.52	0.43
3:L:67:SER:CA	3:L:71:TYR:HE1	2.32	0.43
1:A:150:VAL:HG21	1:A:666:TYR:CD1	2.54	0.43
1:A:144:GLU:OE2	1:A:169:CYS:SG	2.77	0.43
3:L:17:ASP:OD1	3:L:18:ARG:N	2.51	0.42
4:B:1:NAG:O3	4:B:2:NAG:H2	2.19	0.42
1:A:156:GLY:O	1:A:157:LEU:HB3	2.20	0.42
2:H:118:VAL:O	2:H:118:VAL:CG2	2.67	0.42
1:A:264:CYS:SG	1:A:682:PRO:HG3	2.60	0.42
2:H:12:VAL:O	2:H:118:VAL:HA	2.20	0.41
3:L:8:THR:OG1	3:L:10:SER:O	2.37	0.41
2:H:18:VAL:HG21	2:H:116:LEU:HD11	2.03	0.41
1:A:67:PRO:O	1:A:84:TRP:NE1	2.51	0.41
3:L:8:THR:O	3:L:8:THR:HG23	2.21	0.41
1:A:136:ASP:OD1	2:H:50:ARG:NH2	2.53	0.41
3:L:84:ALA:O	3:L:104:VAL:HG22	2.21	0.41
1:A:162:ASN:HB3	1:A:179:GLN:HE21	1.87	0.40
1:A:696:ARG:HH22	1:A:701:ALA:HB2	1.82	0.40
1:A:699:GLN:O	1:A:703:ARG:HB2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	280/1346 (21%)	262 (94%)	18 (6%)	0	100	100
2	H	119/231 (52%)	117 (98%)	2 (2%)	0	100	100
3	L	107/214 (50%)	99 (92%)	8 (8%)	0	100	100
All	All	506/1791 (28%)	478 (94%)	28 (6%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	261/1165 (22%)	260 (100%)	1 (0%)	84	81
2	H	102/198 (52%)	102 (100%)	0	100	100
3	L	97/191 (51%)	97 (100%)	0	100	100
All	All	460/1554 (30%)	459 (100%)	1 (0%)	85	85

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	PHE

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

Mol	Chain	Res	Type
1	A	132	HIS
1	A	196	HIS
1	A	259	HIS
1	A	658	HIS
3	L	34	ASN
3	L	55	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

2 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	B	1	1,4	14,14,15	0.74	1 (7%)	17,19,21	0.45	0
4	NAG	B	2	4	14,14,15	0.25	0	17,19,21	0.66	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
4	NAG	B	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	B	2	4	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1	NAG	O5-C1	-2.42	1.39	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

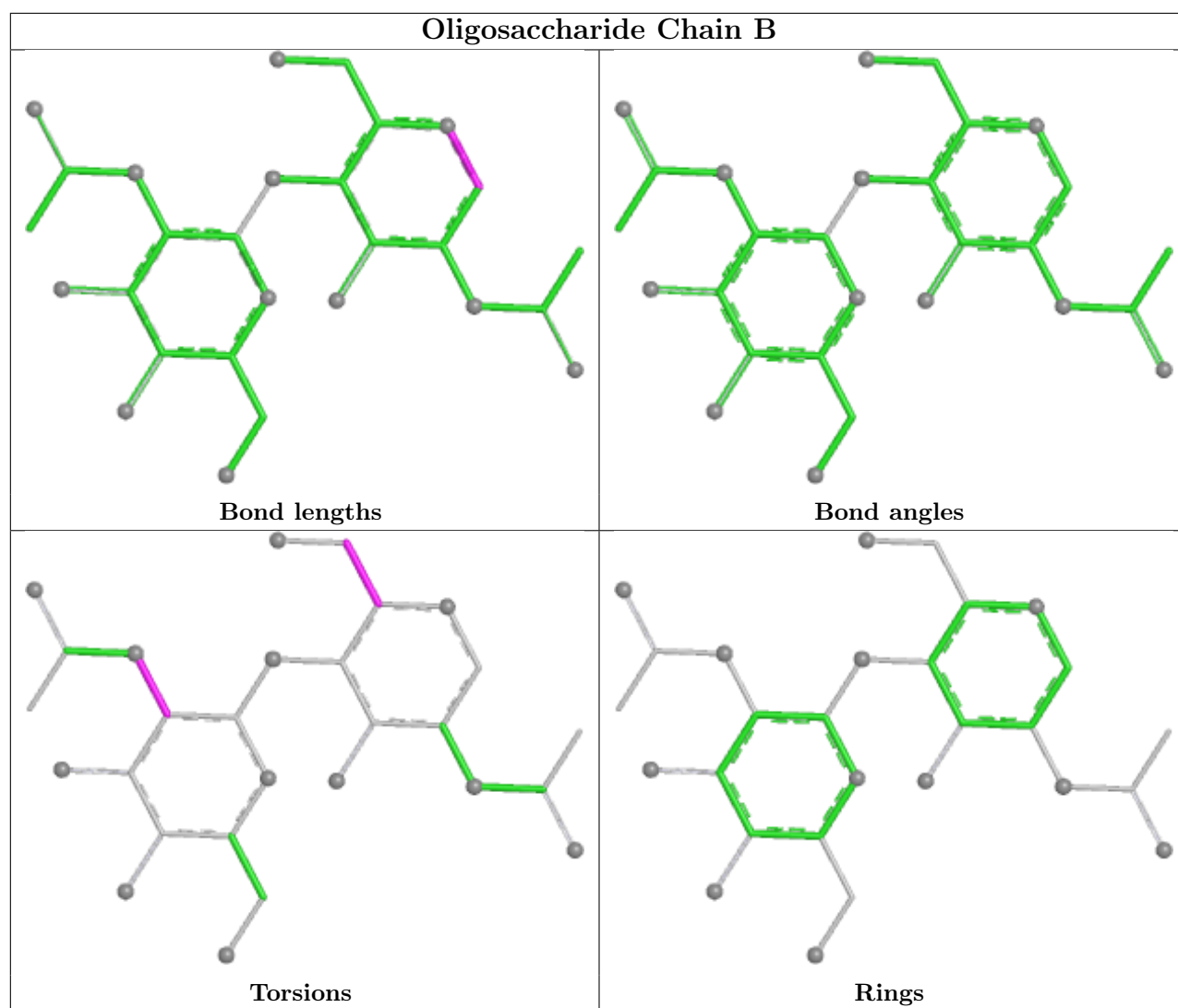
Mol	Chain	Res	Type	Atoms
4	B	2	NAG	C1-C2-N2-C7
4	B	1	NAG	O5-C5-C6-O6
4	B	1	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2	NAG	3	0
4	B	1	NAG	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



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.

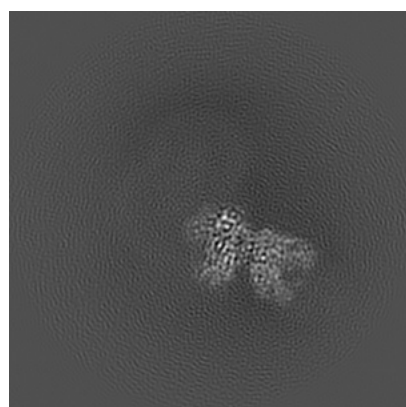
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23408. These allow visual inspection of the internal detail of the map and identification of artifacts.

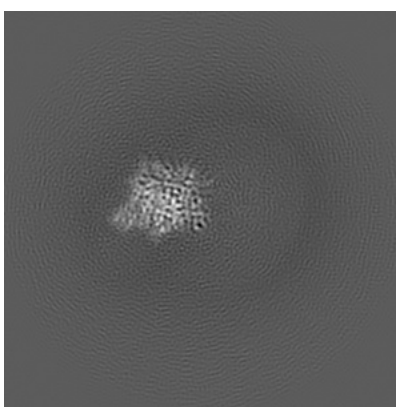
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

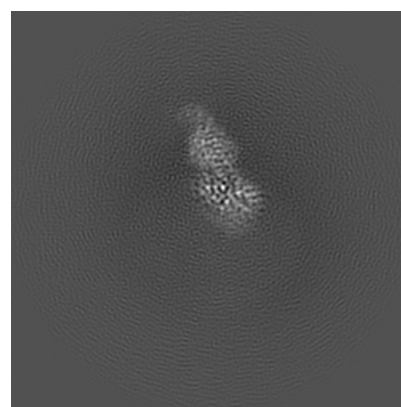
6.1.1 Primary map



X



Y

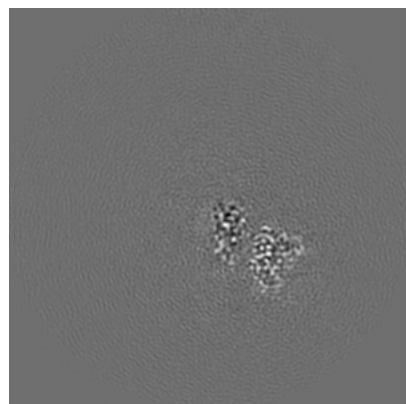


Z

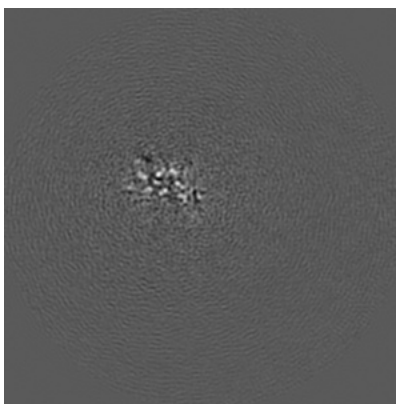
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

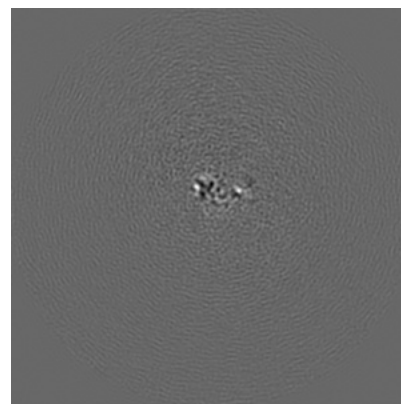
6.2.1 Primary map



X Index: 135



Y Index: 135

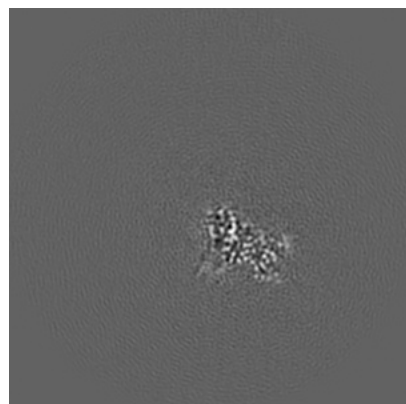


Z Index: 135

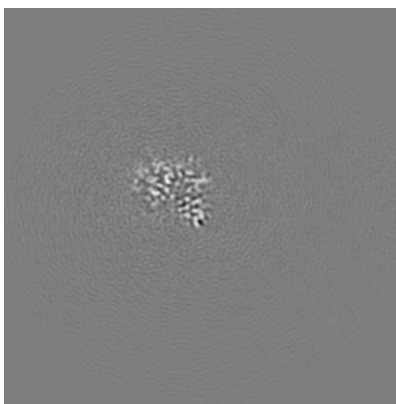
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

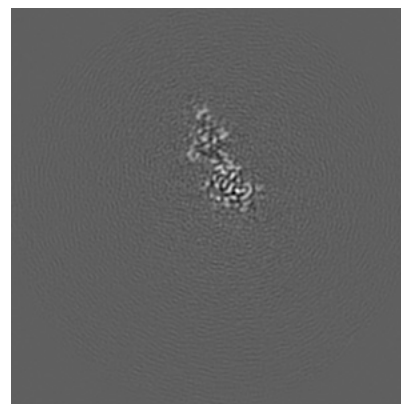
6.3.1 Primary map



X Index: 142



Y Index: 146

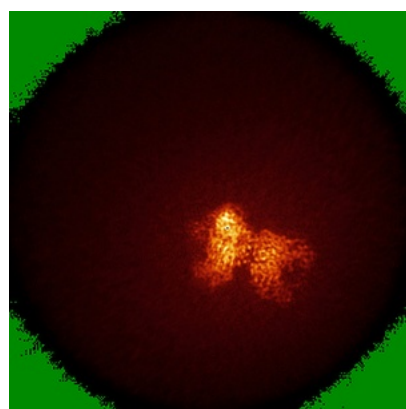


Z Index: 110

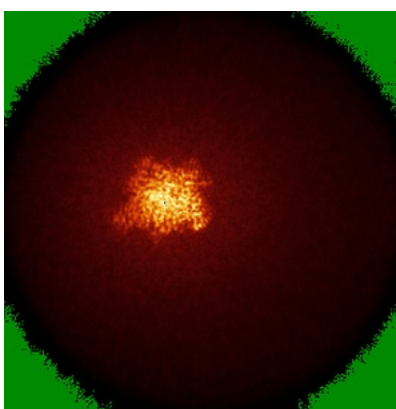
The images above show the largest variance slices of the map in three orthogonal directions.

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

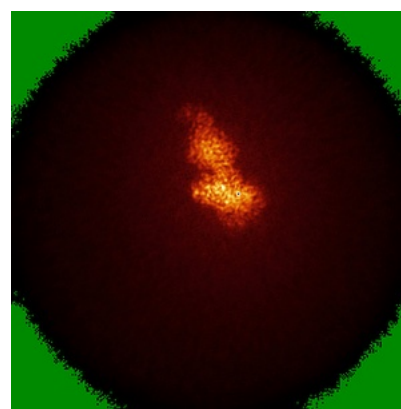
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0561. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

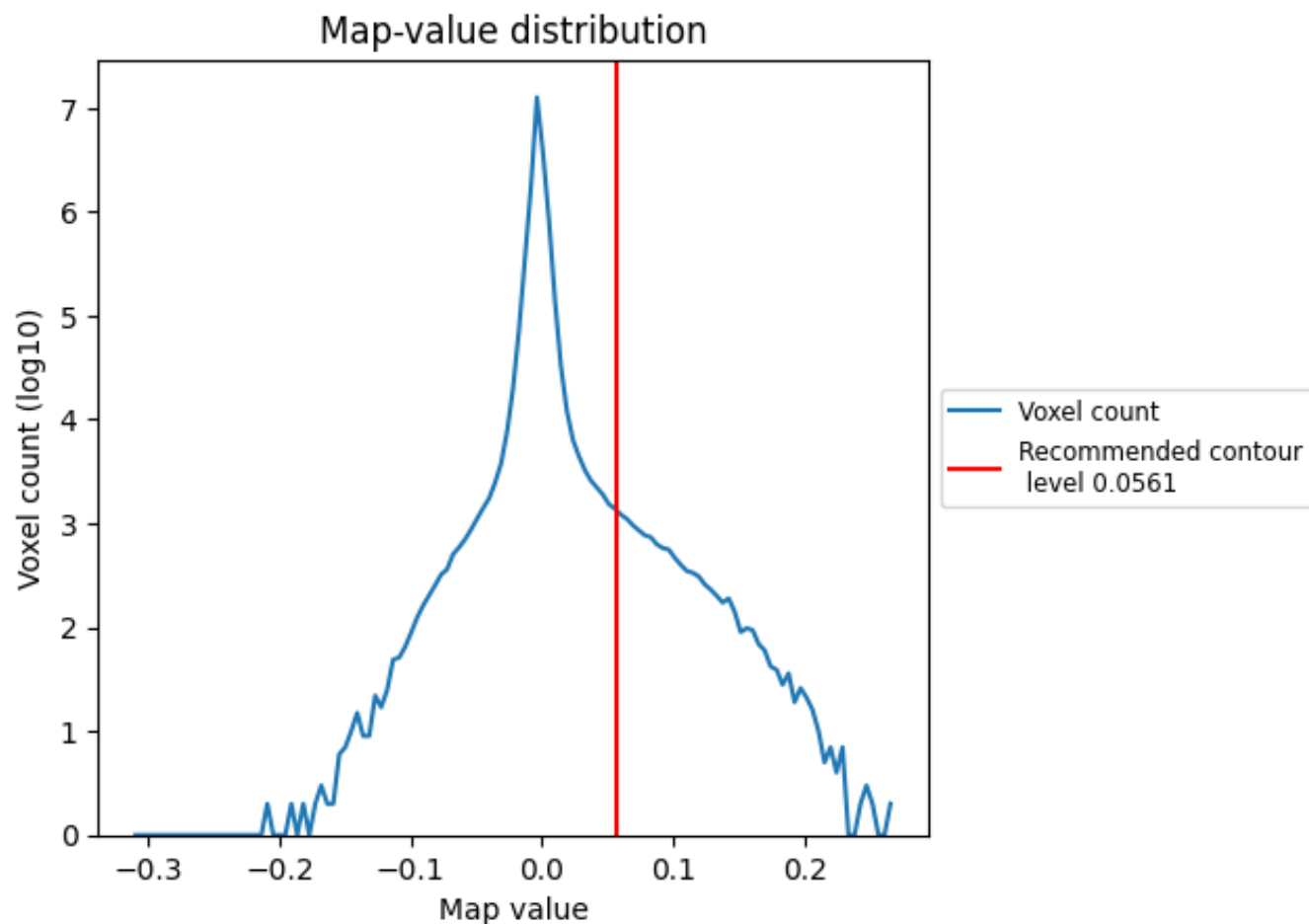
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

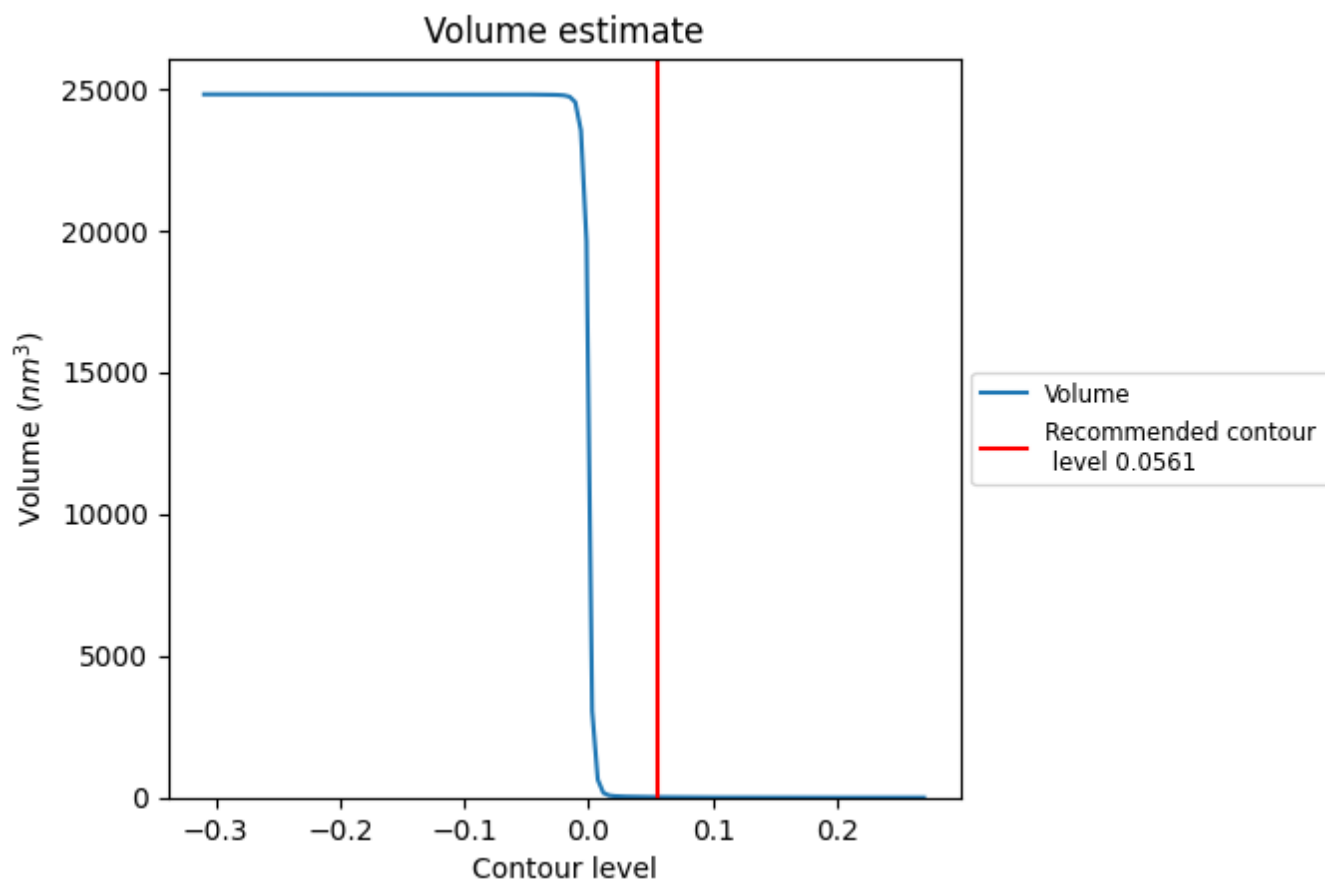
This section contains the results of statistical analysis of the map.

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.

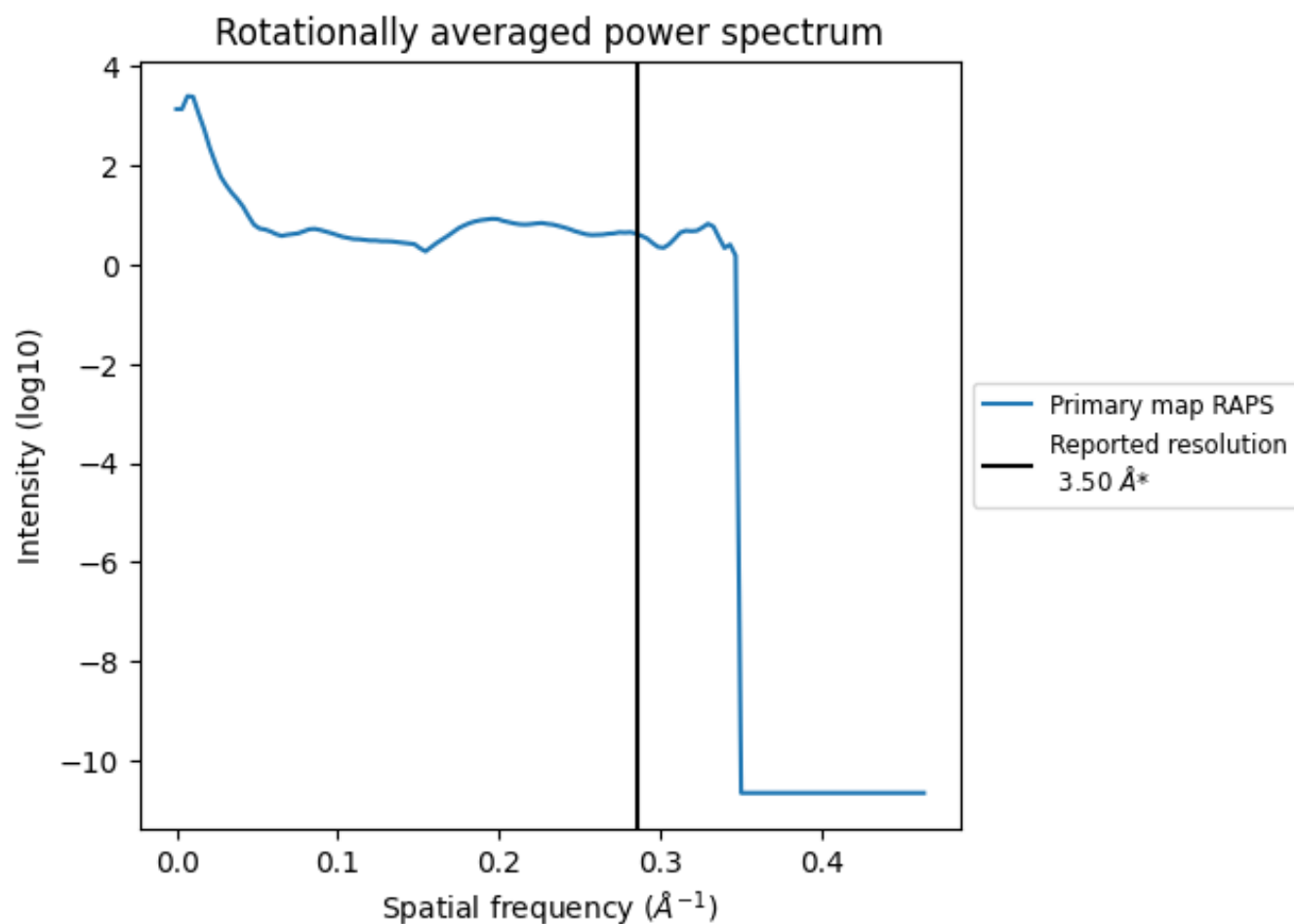
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 16 nm³; this corresponds to an approximate mass of 14 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 ⓘ

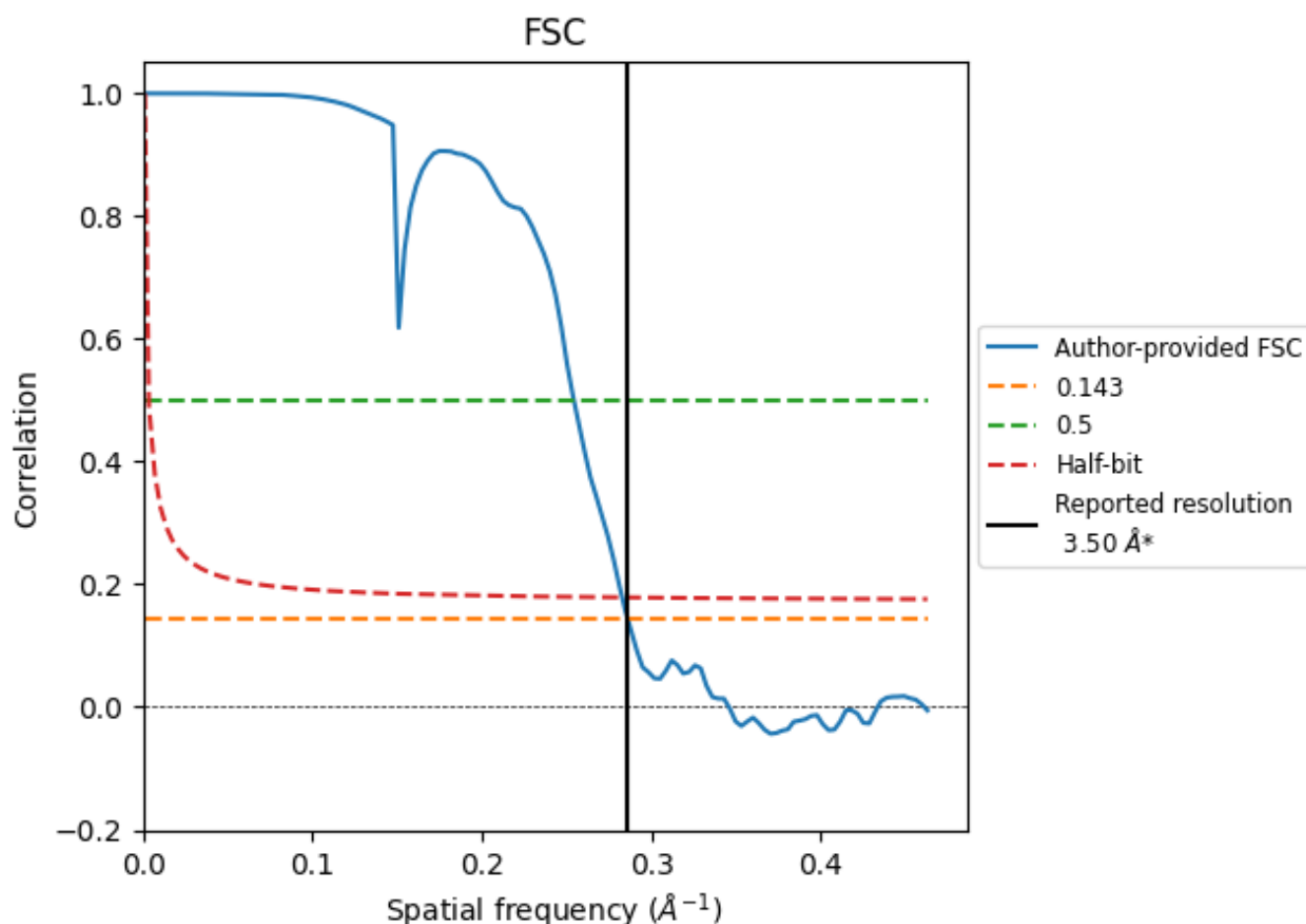


*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

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 0.286 Å⁻¹

8.2 Resolution estimates [i](#)

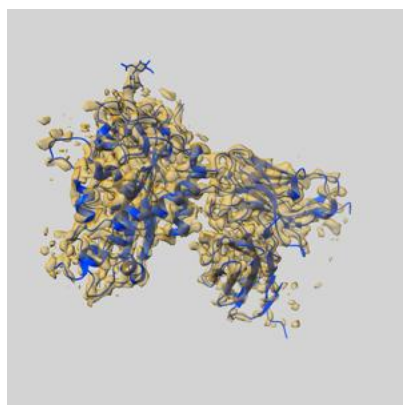
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	3.49	3.93	3.54
Unmasked-calculated*	-	-	-

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

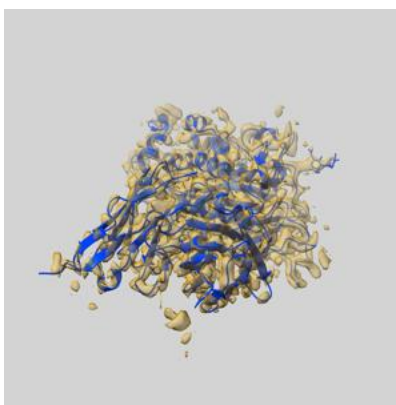
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23408 and PDB model 7LKH. Per-residue inclusion information can be found in section 3 on page 5.

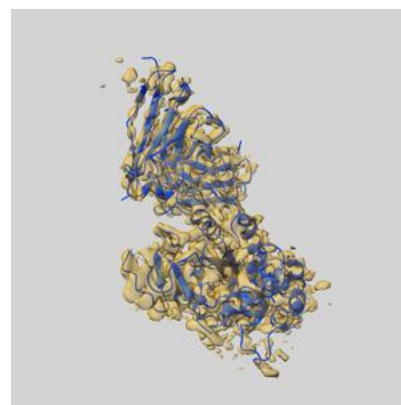
9.1 Map-model overlay [i](#)



X



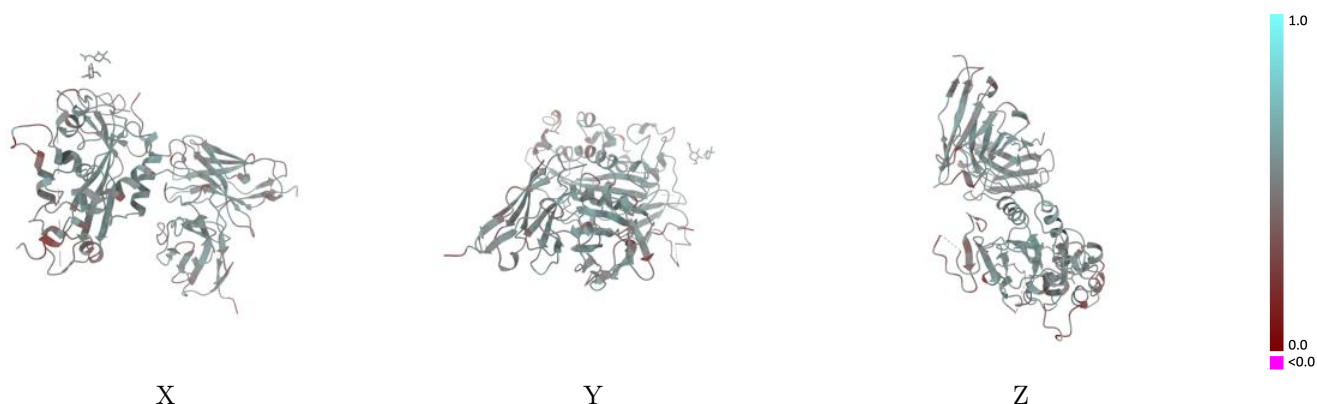
Y



Z

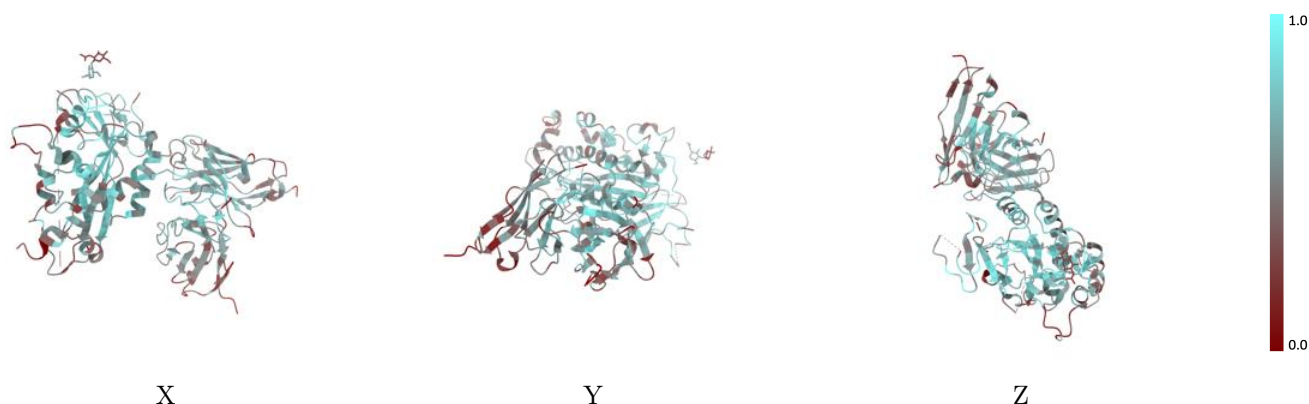
The images above show the 3D surface view of the map at the recommended contour level 0.0561 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



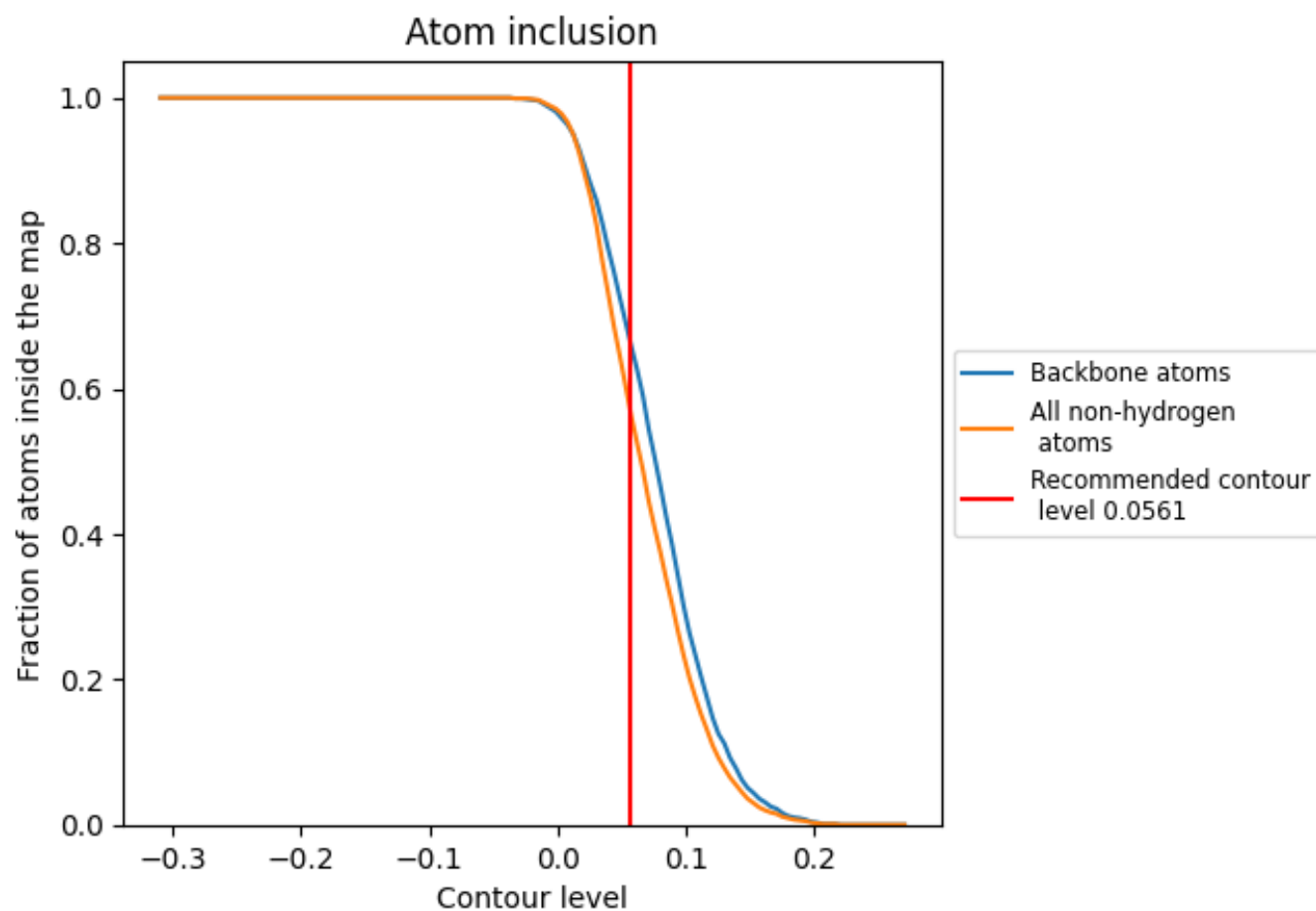
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0561).

9.4 Atom inclusion [i](#)



At the recommended contour level, 67% of all backbone atoms, 57% 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 (0.0561) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5730	0.4970
A	0.6120	0.4940
B	0.3930	0.5010
H	0.5540	0.5060
L	0.4920	0.4960

