Full wwPDB EM Validation Report ⓘ

Nov 29, 2022 – 08:36 PM JST

PDB ID	:	7X1M
EMDB ID	:	EMD-32944
Title	:	The complex structure of Omicron BA.1 RBD with BD604, S309,and S304
Authors	:	Huang, M.; Xie, Y.F.; Qi, J.X.
Deposited on	:	2022-02-24
Resolution	:	2.74 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:

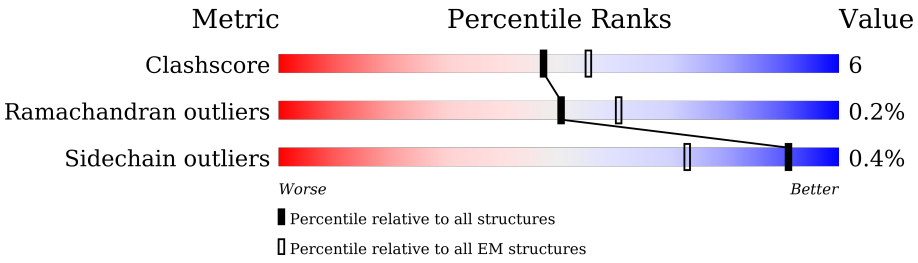
EMDB validation analysis	:	0.0.1.dev43
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

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 2.74 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	A	230	30% 83% 14%
2	B	214	35% 84% 15%
3	C	222	30% 89% 10%
4	D	214	44% 76% 22%
5	H	223	9% 42% 12% 47%
6	L	215	41% 9% 50%
7	M	209	9% 81% 12% 7%
8	E	3	67% 33%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 9908 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 S309 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	225	1701	1074	287	333	7	0	0

- Molecule 2 is a protein called S309 Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	212	1615	1006	276	329	4	0	0

- Molecule 3 is a protein called BD-604 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	219	1615	1016	270	322	7	0	0

- Molecule 4 is a protein called BD-604 Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	213	1637	1024	273	336	4	1	0

- Molecule 5 is a protein called S304 Fab haavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	H	119	921	582	152	183	4	0	0

- Molecule 6 is a protein called S304 Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	L	108	816	516	133	164	3	0	0

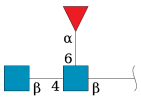
- Molecule 7 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	M	195	Total	C	N	O	S	0	0
			1565	1009	264	284	8		

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	339	ASP	GLY	variant	UNP P0DTC2
M	371	LEU	SER	variant	UNP P0DTC2
M	373	PRO	SER	variant	UNP P0DTC2
M	375	PHE	SER	variant	UNP P0DTC2
M	417	ASN	LYS	variant	UNP P0DTC2
M	440	LYS	ASN	variant	UNP P0DTC2
M	446	SER	GLY	variant	UNP P0DTC2
M	477	ASN	SER	variant	UNP P0DTC2
M	478	LYS	THR	variant	UNP P0DTC2
M	484	ALA	GLU	variant	UNP P0DTC2
M	493	ARG	GLN	variant	UNP P0DTC2
M	496	SER	GLY	variant	UNP P0DTC2
M	498	ARG	GLN	variant	UNP P0DTC2
M	501	TYR	ASN	variant	UNP P0DTC2
M	505	HIS	TYR	variant	UNP P0DTC2

- Molecule 8 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



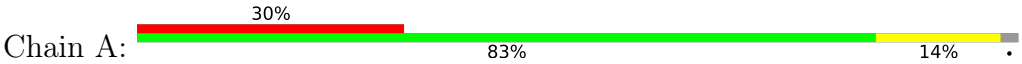
Mol	Chain	Residues	Atoms				AltConf	Trace
8	E	3	Total	C	N	O	0	0
			38	22	2	14		

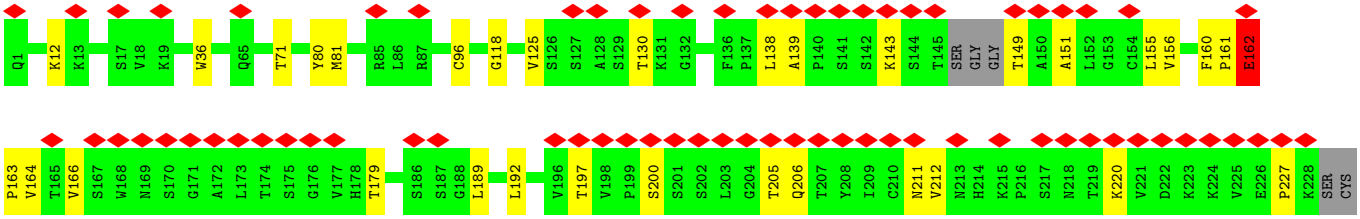
3 Residue-property plots (i)

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: S309 Fab heavy chain

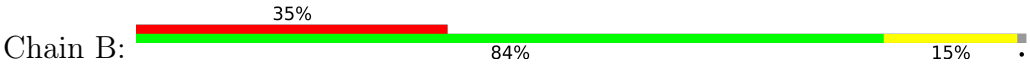
Chain A:





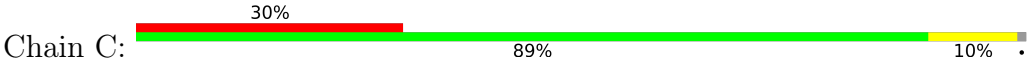
• Molecule 2: S309 Fab light chain

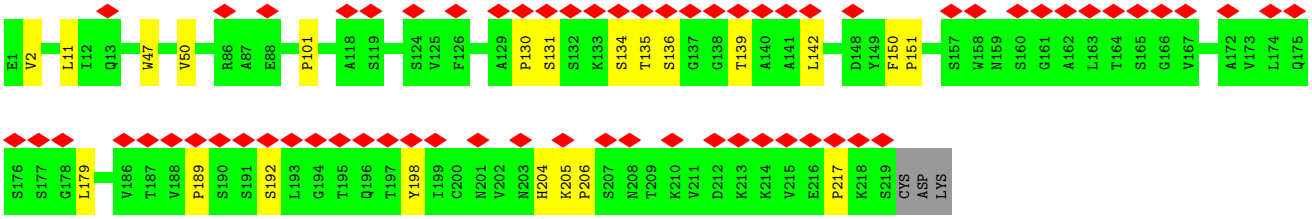
Chain B:



• Molecule 3: BD-604 Fab heavy chain

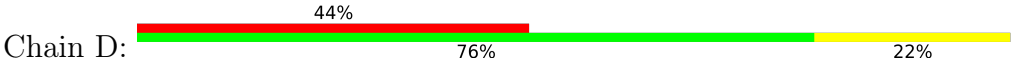
Chain C:

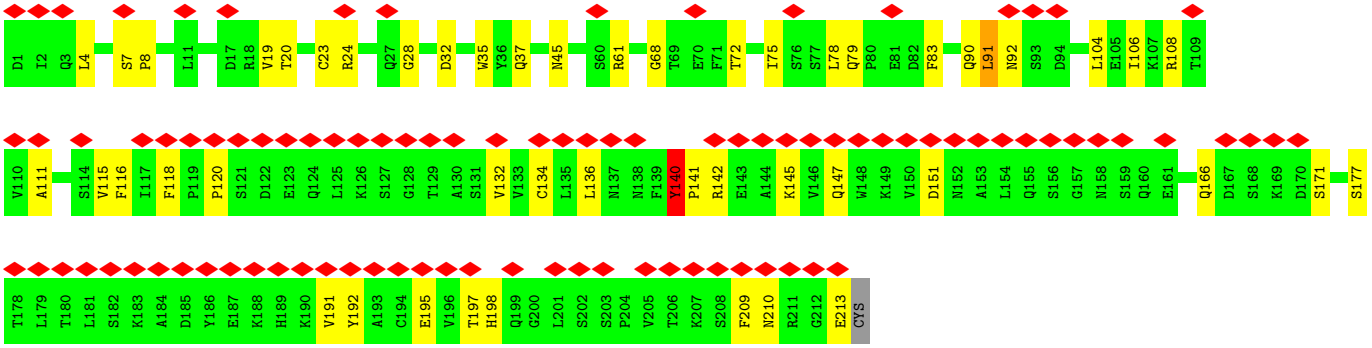




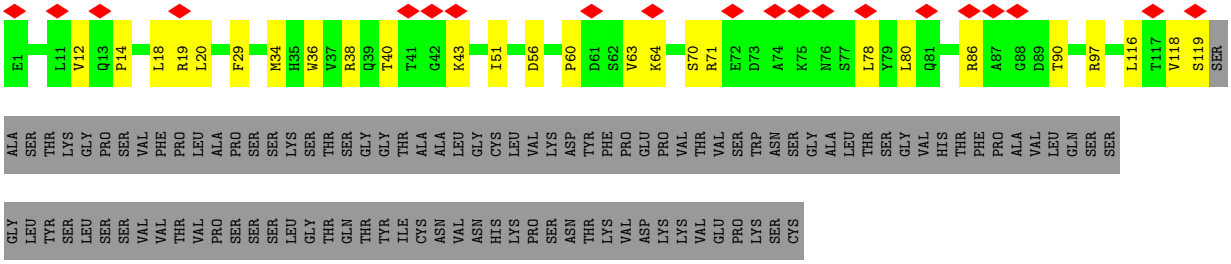
• Molecule 4: BD-604 Fab light chain

Chain D:

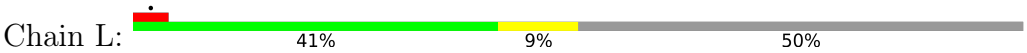


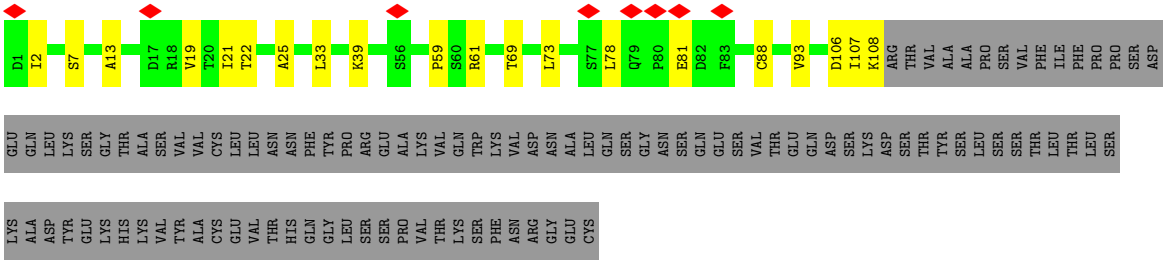


• Molecule 5: S304 Fab haavy chain

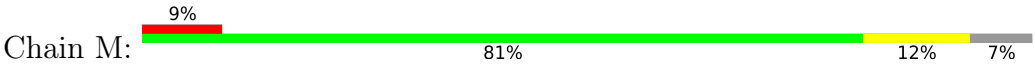


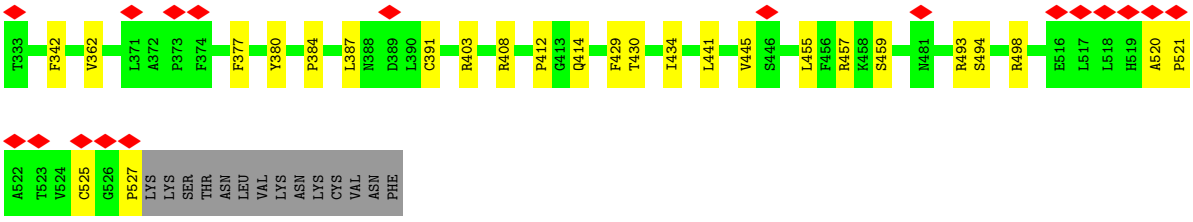
• Molecule 6: S304 Fab light chain





• Molecule 7: Spike protein S1





• Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	553923	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 ($e^-/Å^2$)	60.0	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.386	Depositor
Minimum map value	-2.215	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.079	Depositor
Recommended contour level	0.7	Depositor
Map size (Å)	306.0, 306.0, 306.0	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.85, 0.85, 0.85	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FUC, 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	0.31	0/1743	0.55	0/2374
2	B	0.30	0/1648	0.53	0/2238
3	C	0.32	0/1651	0.50	0/2249
4	D	0.28	0/1673	0.56	1/2272 (0.0%)
5	H	0.37	0/944	0.56	0/1280
6	L	0.33	0/835	0.51	0/1135
7	M	0.36	0/1612	0.53	0/2195
All	All	0.32	0/10106	0.54	1/13743 (0.0%)

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
4	D	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	D	91	LEU	CA-CB-CG	5.56	128.10	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	162	GLU	Peptide
4	D	140	TYR	Peptide

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	1701	0	1661	19	0
2	B	1615	0	1576	19	0
3	C	1615	0	1596	16	0
4	D	1637	0	1592	31	0
5	H	921	0	862	17	0
6	L	816	0	799	12	0
7	M	1565	0	1493	18	0
8	E	38	0	34	1	0
All	All	9908	0	9613	125	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 6.

All (125) 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:A:149:THR:N	1:A:200:SER:HG	1.73	0.85
2:B:108:ARG:NH1	2:B:170:ASP:O	2.23	0.70
4:D:4:LEU:HD11	4:D:90:GLN:HB2	1.73	0.70
6:L:21:ILE:HD11	6:L:73:LEU:HD23	1.75	0.69
4:D:78:LEU:HD12	4:D:83:PHE:HE1	1.58	0.67
6:L:78:LEU:HD21	6:L:107:ILE:HD12	1.75	0.67
4:D:7:SER:OG	4:D:24:ARG:NH1	2.30	0.65
4:D:7:SER:HB2	4:D:8:PRO:HD3	1.79	0.65
7:M:380:TYR:O	7:M:430:THR:HA	1.96	0.64
3:C:11:LEU:HD11	3:C:151:PRO:HG3	1.80	0.63
1:A:138:LEU:HD11	1:A:155:LEU:HB2	1.81	0.63
7:M:342:PHE:HB2	8:E:1:NAG:H82	1.80	0.62
4:D:32:ASP:HB3	4:D:91:LEU:HD21	1.81	0.62
7:M:362:VAL:HG13	7:M:527:PRO:HD2	1.81	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:130:PRO:HG2	3:C:217:PRO:HG3	1.83	0.60
3:C:131:SER:H	3:C:134:SER:HB2	1.65	0.60
1:A:130:THR:HG23	1:A:161:PRO:HD3	1.85	0.59
3:C:204:HIS:HD2	3:C:206:PRO:HD2	1.69	0.57
4:D:92:ASN:OD1	7:M:403:ARG:NH2	2.37	0.57
4:D:120:PRO:HD3	4:D:132:VAL:HG22	1.87	0.57
2:B:13:LEU:O	2:B:107:LYS:N	2.31	0.57
4:D:61:ARG:NH1	4:D:79:GLN:OE1	2.37	0.57
5:H:12:VAL:HG11	5:H:18:LEU:HD11	1.86	0.57
4:D:147:GLN:HB2	4:D:195:GLU:HB3	1.89	0.55
7:M:377:PHE:HD1	7:M:434:ILE:HG13	1.71	0.55
3:C:136:SER:HB3	4:D:116:PHE:HE1	1.71	0.55
4:D:145:LYS:HB2	4:D:197:THR:HB	1.88	0.54
5:H:12:VAL:HG11	5:H:18:LEU:HD21	1.89	0.54
4:D:78:LEU:HD12	4:D:83:PHE:CE1	2.42	0.54
5:H:36:TRP:HE1	5:H:78:LEU:CD2	2.21	0.53
2:B:32:THR:HG22	7:M:441:LEU:HD22	1.91	0.53
1:A:156:VAL:HG11	1:A:164:VAL:HG21	1.89	0.53
5:H:90:THR:OG1	5:H:116:LEU:O	2.19	0.53
7:M:457:ARG:NH1	7:M:459:SER:O	2.37	0.53
1:A:205:THR:OG1	1:A:206:GLN:OE1	2.25	0.53
1:A:162:GLU:HG3	1:A:163:PRO:HD3	1.92	0.52
2:B:125:LEU:O	2:B:183:LYS:NZ	2.34	0.51
3:C:192:SER:O	3:C:198:TYR:OH	2.26	0.51
5:H:70:SER:O	5:H:78:LEU:HG	2.11	0.51
4:D:19:VAL:HG12	4:D:75:ILE:HB	1.93	0.50
4:D:141:PRO:HD2	4:D:198:HIS:CE1	2.47	0.50
6:L:13:ALA:O	6:L:108:LYS:N	2.45	0.50
2:B:113:PRO:HD3	2:B:198:HIS:ND1	2.27	0.50
4:D:108:ARG:HH12	4:D:111:ALA:HB2	1.76	0.50
2:B:114:SER:HB3	2:B:137:ASN:HB2	1.95	0.49
5:H:51:ILE:HG12	5:H:71:ARG:HG3	1.95	0.48
5:H:19:ARG:HA	5:H:80:LEU:O	2.13	0.48
6:L:19:VAL:HG21	6:L:78:LEU:HD12	1.95	0.48
6:L:93:VAL:HG22	7:M:380:TYR:CD2	2.49	0.48
4:D:151:ASP:OD1	4:D:191:VAL:HB	2.14	0.47
7:M:493:ARG:HG3	7:M:494:SER:N	2.29	0.47
3:C:134:SER:HG	4:D:118:PHE:HE1	1.61	0.47
2:B:195:GLU:OE2	2:B:204:PRO:HB3	2.15	0.47
2:B:105:GLU:OE1	2:B:173:TYR:OH	2.33	0.47
4:D:134:CYS:HB3	4:D:177:SER:OG	2.14	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:M:391:CYS:HA	7:M:525:CYS:HA	1.97	0.47
4:D:78:LEU:HD11	4:D:104:LEU:HD21	1.96	0.47
6:L:25:ALA:O	6:L:69:THR:OG1	2.27	0.47
7:M:408:ARG:NH1	7:M:414:GLN:OE1	2.40	0.47
4:D:90:GLN:CD	4:D:91:LEU:H	2.19	0.46
5:H:60:PRO:HD2	5:H:63:VAL:HB	1.97	0.46
4:D:83:PHE:HE2	4:D:106:ILE:HG12	1.80	0.46
1:A:151:ALA:HB2	1:A:197:THR:HG22	1.98	0.46
7:M:384:PRO:HA	7:M:387:LEU:HD23	1.97	0.46
3:C:101:PRO:HD2	7:M:455:LEU:HD21	1.98	0.46
6:L:59:PRO:HB3	6:L:61:ARG:CZ	2.46	0.45
6:L:33:LEU:HD12	6:L:88:CYS:HB2	1.98	0.45
5:H:56:ASP:N	5:H:56:ASP:OD1	2.49	0.45
3:C:131:SER:O	3:C:135:THR:HG23	2.16	0.45
6:L:39:LYS:HE2	6:L:81:GLU:O	2.17	0.45
1:A:36:TRP:CE2	1:A:81:MET:HB2	2.52	0.45
1:A:96:CYS:O	1:A:118:GLY:N	2.50	0.45
1:A:212:VAL:O	1:A:220:LYS:HA	2.16	0.45
5:H:34:MET:HE2	5:H:97:ARG:HA	1.99	0.45
3:C:139:THR:HG22	3:C:189:PRO:HA	1.98	0.45
5:H:29:PHE:HZ	5:H:78:LEU:HB2	1.81	0.45
3:C:205:LYS:HB3	3:C:206:PRO:HD3	1.99	0.44
1:A:160:PHE:HB2	1:A:189:LEU:HD12	1.99	0.44
1:A:166:VAL:HA	1:A:211:ASN:O	2.17	0.44
4:D:210:ASN:O	4:D:213:GLU:HG2	2.18	0.44
2:B:80:GLU:N	2:B:80:GLU:OE1	2.51	0.44
3:C:150:PHE:HA	3:C:151:PRO:HA	1.83	0.44
2:B:120:PRO:HB2	2:B:125:LEU:HD11	1.99	0.43
7:M:445:VAL:O	7:M:498:ARG:NH1	2.52	0.43
4:D:20:THR:HG23	4:D:72:THR:HG23	2.01	0.43
4:D:108:ARG:HB3	4:D:140:TYR:CE2	2.54	0.43
4:D:166:GLN:HG2	4:D:171:SER:HA	2.01	0.43
5:H:86:ARG:O	5:H:118:VAL:HG21	2.19	0.43
1:A:71:THR:OG1	1:A:80:TYR:HB2	2.18	0.43
5:H:51:ILE:HG21	5:H:78:LEU:HD11	2.01	0.43
2:B:39:GLN:O	2:B:85:ALA:HB1	2.19	0.42
6:L:2:ILE:HD11	6:L:93:VAL:HG21	2.00	0.42
2:B:190:LYS:HE3	2:B:210:ASN:HB2	2.00	0.42
1:A:160:PHE:CD1	1:A:161:PRO:HA	2.54	0.42
1:A:139:ALA:HB1	1:A:227:PRO:HA	2.01	0.42
2:B:134:CYS:HB2	2:B:148:TRP:CZ2	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:192:SER:HG	3:C:198:TYR:HH	1.60	0.42
2:B:32:THR:CG2	7:M:441:LEU:HD22	2.50	0.42
1:A:143:LYS:H	1:A:143:LYS:HG2	1.70	0.41
1:A:179:THR:HG23	1:A:192:LEU:HD21	2.02	0.41
3:C:150:PHE:HB2	3:C:179:LEU:HD22	2.01	0.41
4:D:115:VAL:HG22	4:D:136:LEU:HD23	2.02	0.41
2:B:147:GLN:HB3	2:B:154:LEU:HD22	2.02	0.41
3:C:47:TRP:HZ2	3:C:50:VAL:HG12	1.85	0.41
5:H:38:ARG:HH21	5:H:43:LYS:NZ	2.18	0.41
2:B:150:VAL:HG13	2:B:192:TYR:HE1	1.84	0.41
1:A:162:GLU:HG3	1:A:163:PRO:CD	2.50	0.41
3:C:130:PRO:HD3	3:C:142:LEU:HB3	2.03	0.41
5:H:20:LEU:H	5:H:20:LEU:HD23	1.85	0.41
4:D:23:CYS:HB2	4:D:35:TRP:CH2	2.55	0.41
5:H:14:PRO:HD3	5:H:119:SER:C	2.41	0.41
4:D:28:GLY:HA2	4:D:68:GLY:O	2.19	0.41
4:D:83:PHE:CE2	4:D:106:ILE:HG12	2.55	0.41
5:H:40:THR:OG1	5:H:43:LYS:HG2	2.21	0.41
2:B:165:GLU:HG2	2:B:166:GLN:H	1.85	0.41
7:M:412:PRO:HG3	7:M:429:PHE:HB2	2.02	0.41
7:M:493:ARG:HG3	7:M:494:SER:H	1.86	0.41
1:A:12:LYS:O	1:A:125:VAL:HA	2.21	0.40
2:B:90:GLN:NE2	2:B:96:LEU:HD21	2.36	0.40
2:B:147:GLN:HB2	2:B:195:GLU:HB3	2.03	0.40
6:L:106:ASP:OD1	6:L:107:ILE:N	2.54	0.40
4:D:192:TYR:HB2	4:D:209:PHE:CE1	2.56	0.40
6:L:7:SER:OG	6:L:22:THR:OG1	2.39	0.40
7:M:520:ALA:HB1	7:M:521:PRO:HD2	2.04	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	221/230 (96%)	200 (90%)	20 (9%)	1 (0%)	29 48
2	B	210/214 (98%)	192 (91%)	18 (9%)	0	100 100
3	C	217/222 (98%)	201 (93%)	15 (7%)	1 (0%)	29 48
4	D	212/214 (99%)	188 (89%)	23 (11%)	1 (0%)	29 48
5	H	117/223 (52%)	113 (97%)	4 (3%)	0	100 100
6	L	106/215 (49%)	98 (92%)	8 (8%)	0	100 100
7	M	193/209 (92%)	182 (94%)	11 (6%)	0	100 100
All	All	1276/1527 (84%)	1174 (92%)	99 (8%)	3 (0%)	50 69

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	2	VAL
4	D	140	TYR
1	A	162	GLU

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	A	189/192 (98%)	189 (100%)	0	100 100
2	B	183/185 (99%)	182 (100%)	1 (0%)	88 92
3	C	183/186 (98%)	183 (100%)	0	100 100
4	D	187/187 (100%)	184 (98%)	3 (2%)	62 78
5	H	96/186 (52%)	95 (99%)	1 (1%)	76 85
6	L	92/188 (49%)	92 (100%)	0	100 100
7	M	170/184 (92%)	170 (100%)	0	100 100
All	All	1100/1308 (84%)	1095 (100%)	5 (0%)	91 92

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	78	ARG
4	D	45[A]	ASN
4	D	45[B]	ASN
4	D	142	ARG
5	H	64	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	211	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)

3 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	$\#\|Z\| > 2$
8	NAG	E	1	8,7	14,14,15	0.43	0	17,19,21	0.56	0
8	NAG	E	2	8	14,14,15	0.20	0	17,19,21	0.45	0
8	FUC	E	3	8	10,10,11	0.72	0	14,14,16	0.74	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
8	NAG	E	1	8,7	-	0/6/23/26	0/1/1/1
8	NAG	E	2	8	-	2/6/23/26	0/1/1/1
8	FUC	E	3	8	-	-	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	E	2	NAG	O5-C5-C6-O6
8	E	2	NAG	C4-C5-C6-O6

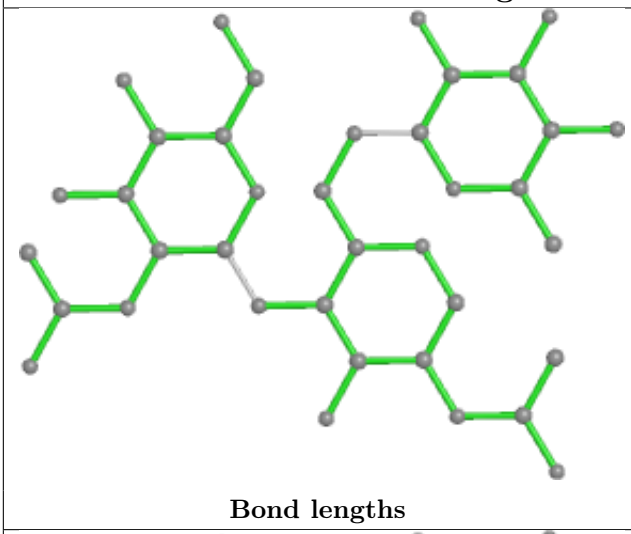
There are no ring outliers.

1 monomer is involved in 1 short contact:

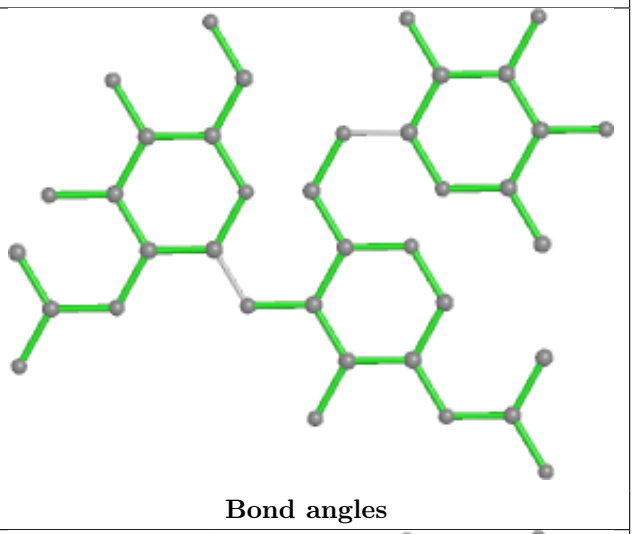
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	E	1	NAG	1	0

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

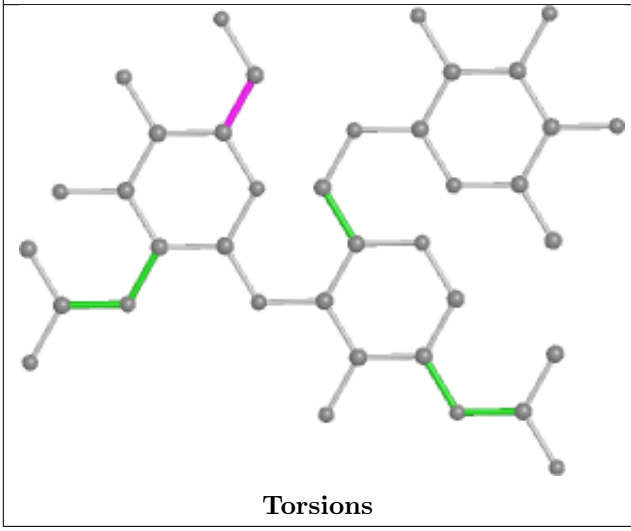
Oligosaccharide Chain E



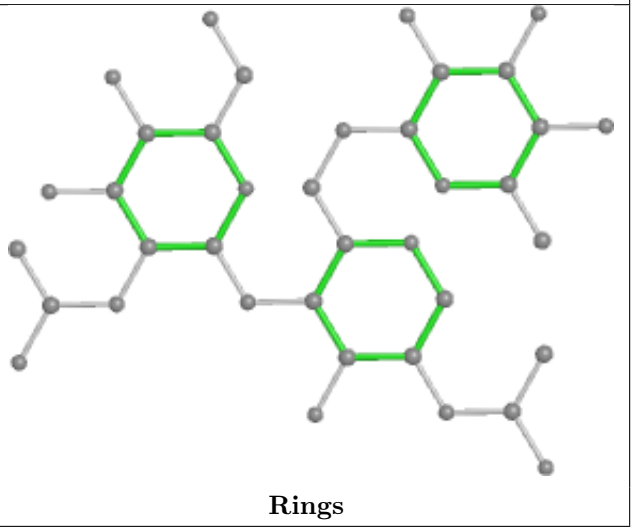
Bond lengths



Bond angles



Torsions



Rings

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.

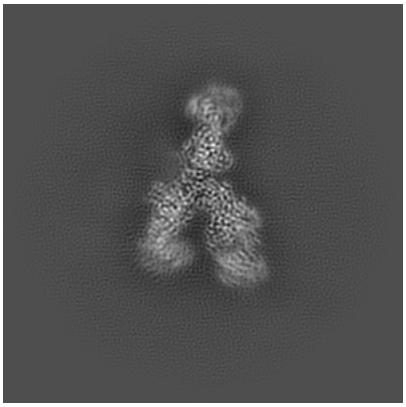
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-32944. These allow visual inspection of the internal detail of the map and identification of artifacts.

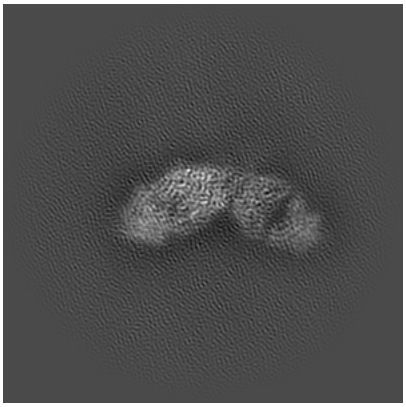
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

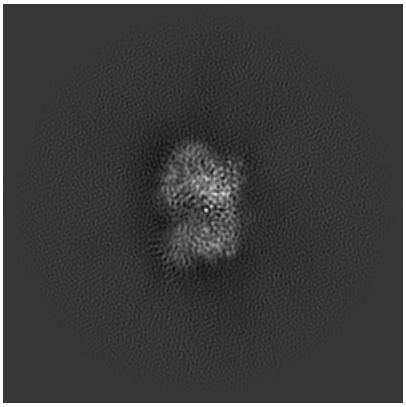
6.1.1 Primary map



X



Y

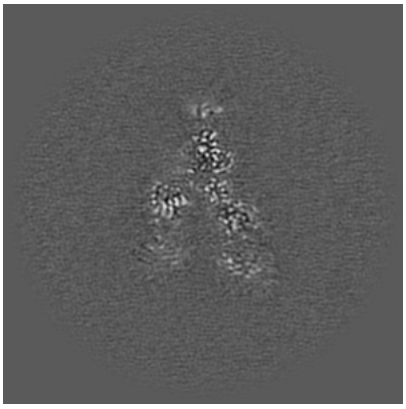


Z

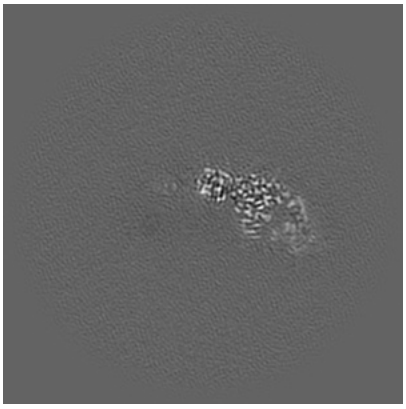
The images above show the map projected in three orthogonal directions.

6.2 Central slices

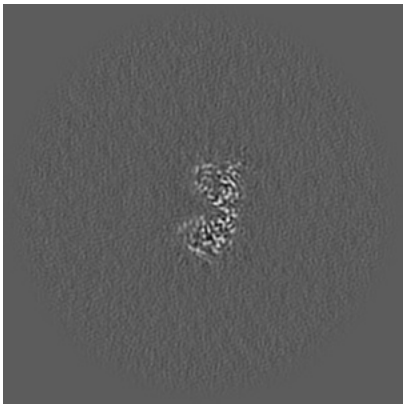
6.2.1 Primary map



X Index: 180



Y Index: 180

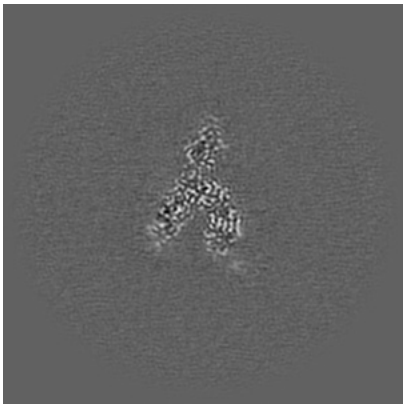


Z Index: 180

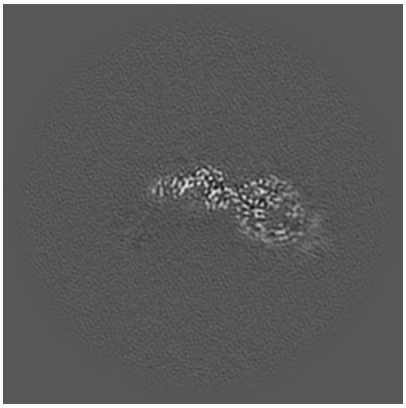
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

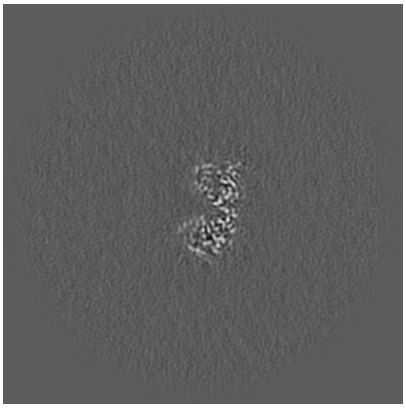
6.3.1 Primary map



X Index: 196



Y Index: 188



Z Index: 180

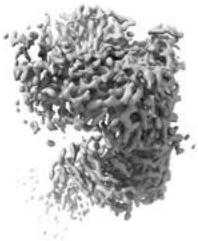
The images above show the largest variance slices of the map in three orthogonal directions.

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

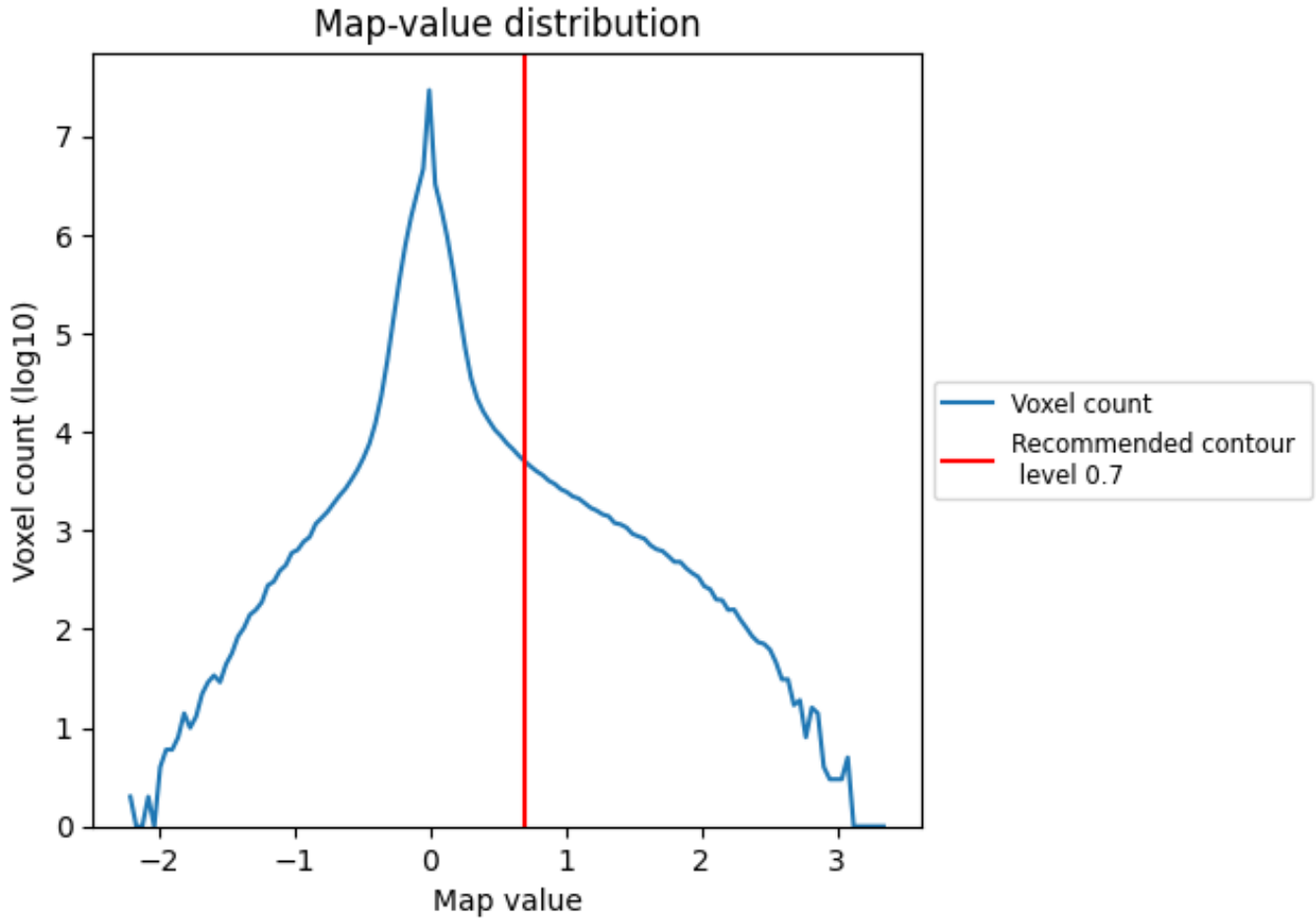
6.5 Mask visualisation (i)

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

7 Map analysis

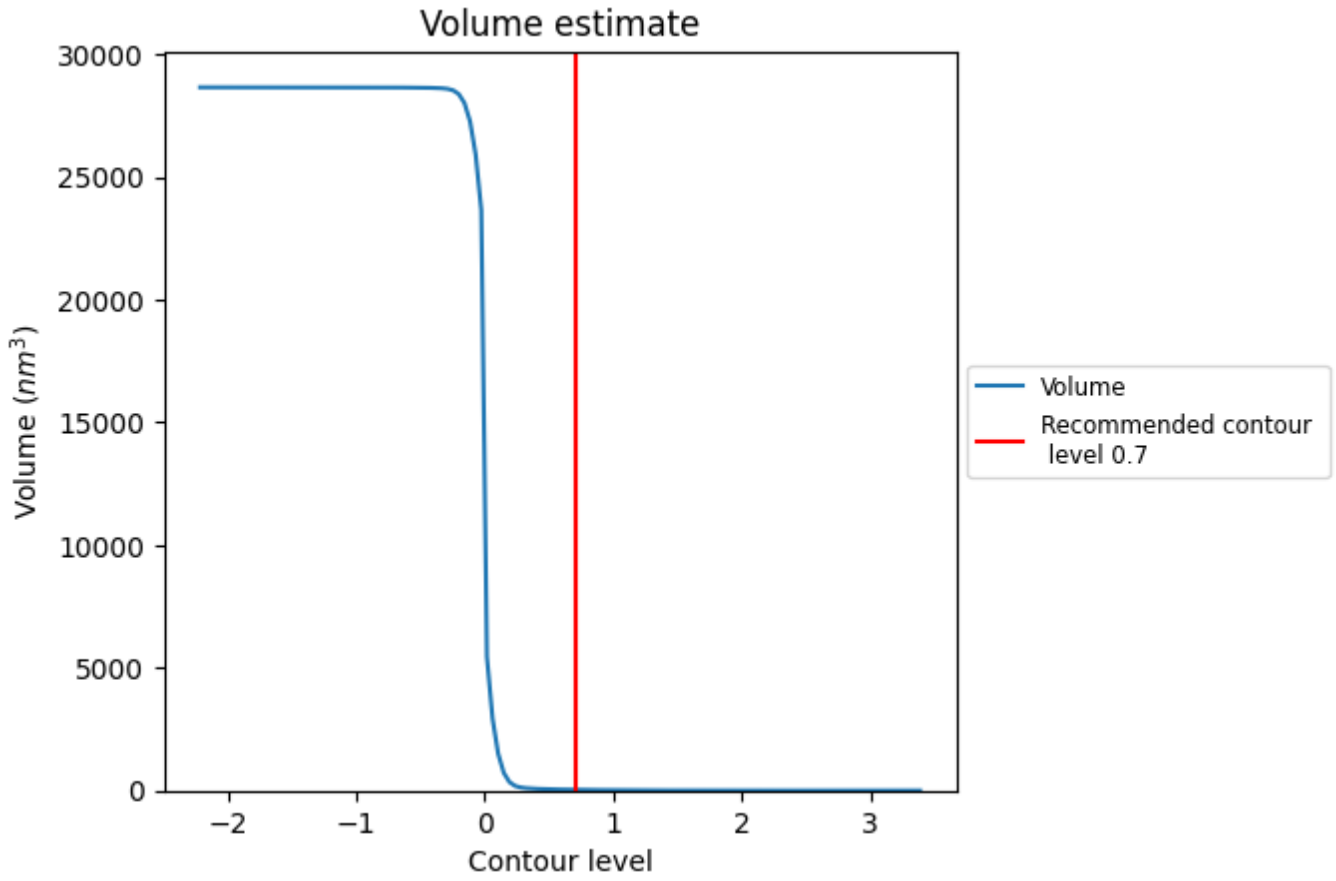
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

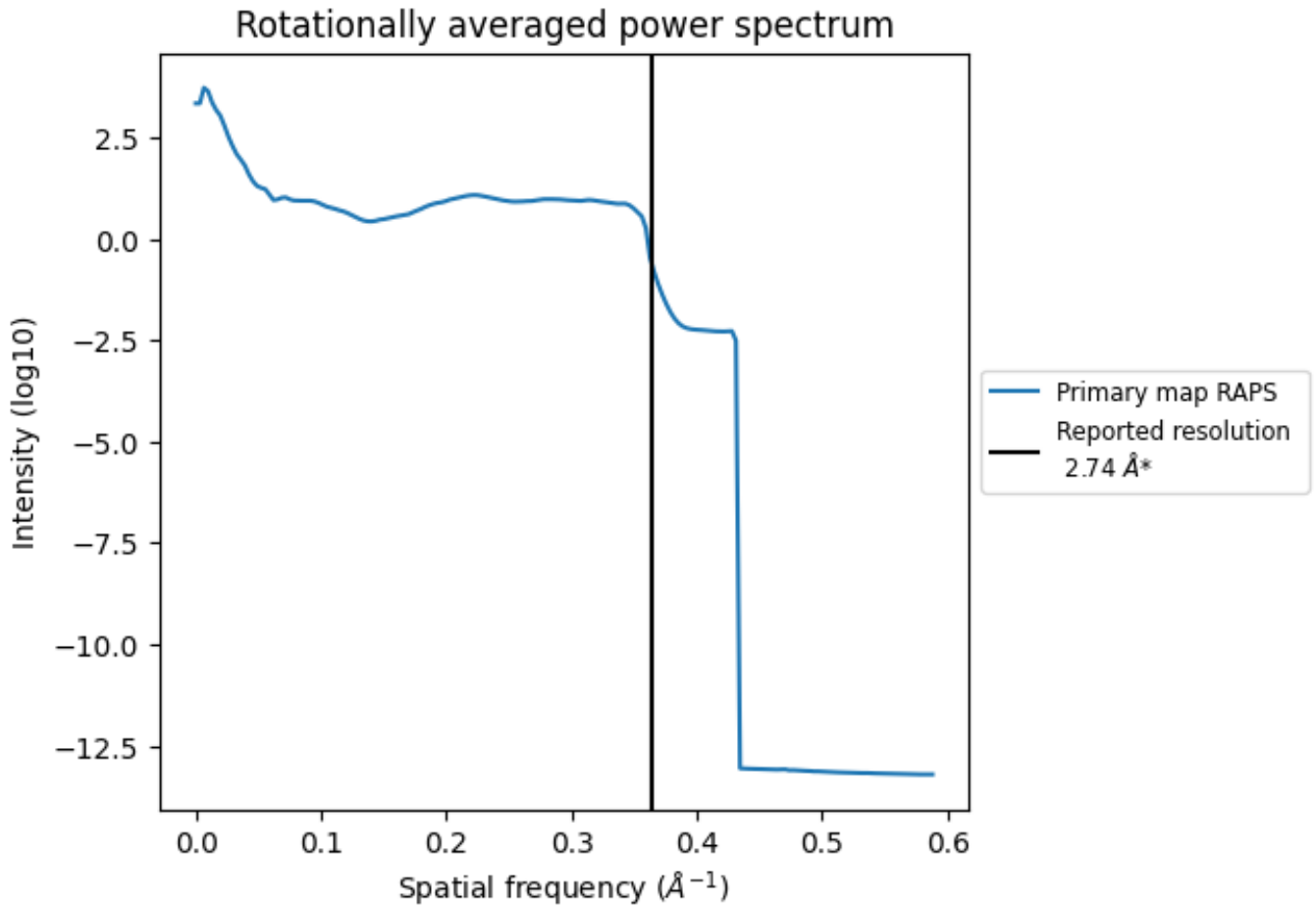
7.2 Volume estimate



The volume at the recommended contour level is 33 nm^3; this corresponds to an approximate mass of 29 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)

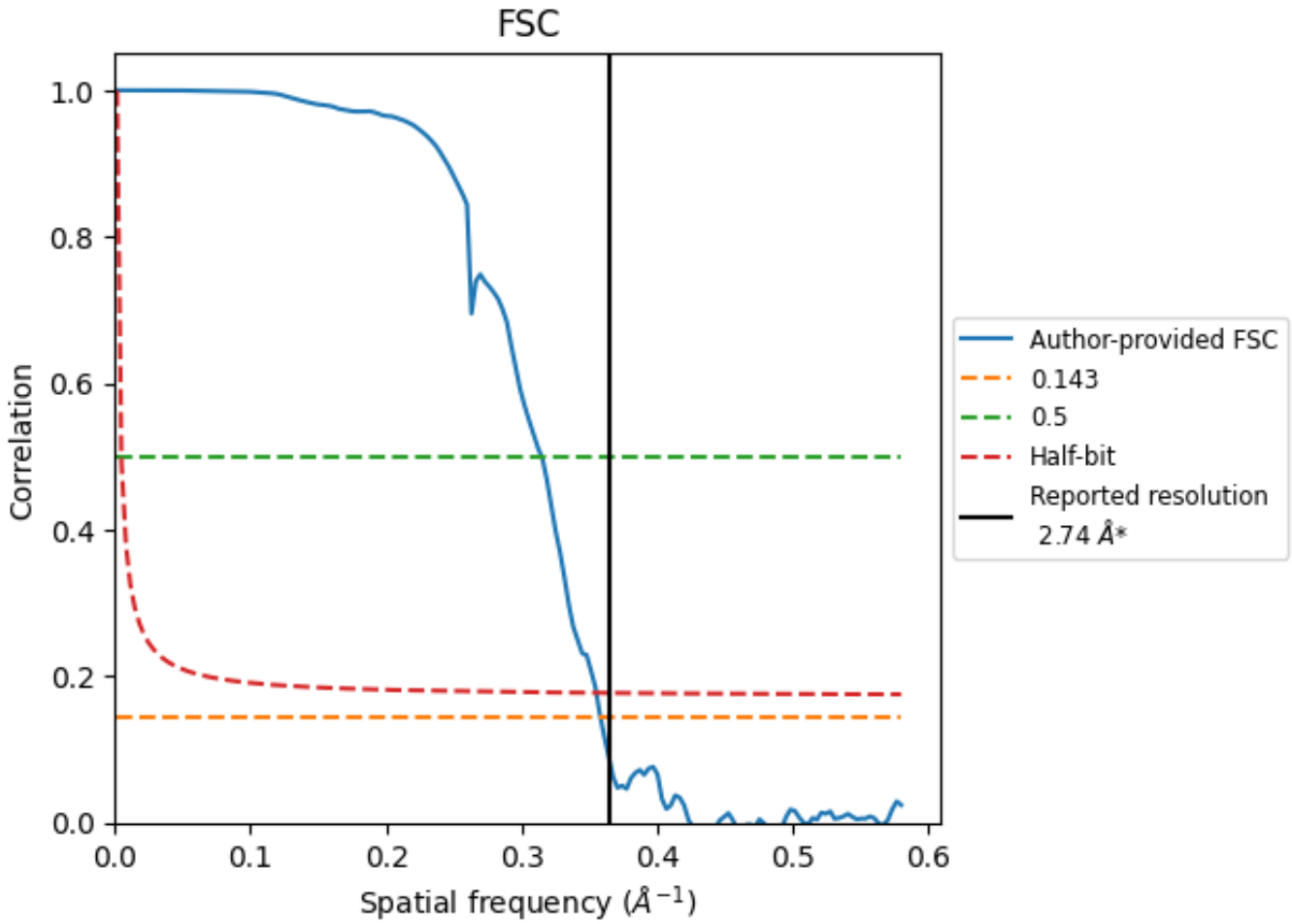


*Reported resolution corresponds to spatial frequency of 0.365 $Å^{-1}$

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.365 Å^{-1}

8.2 Resolution estimates (i)

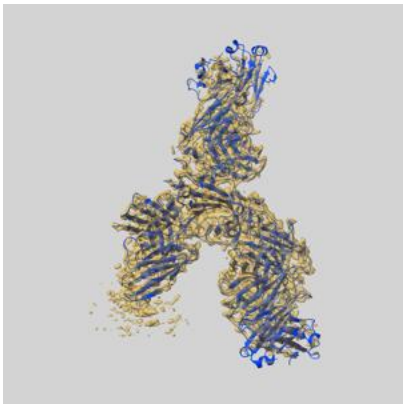
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.74	-	-
Author-provided FSC curve	2.79	3.18	2.82
Unmasked-calculated*	-	-	-

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

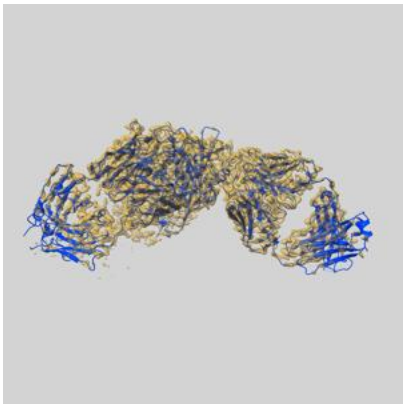
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-32944 and PDB model 7X1M. Per-residue inclusion information can be found in section 3 on page 5.

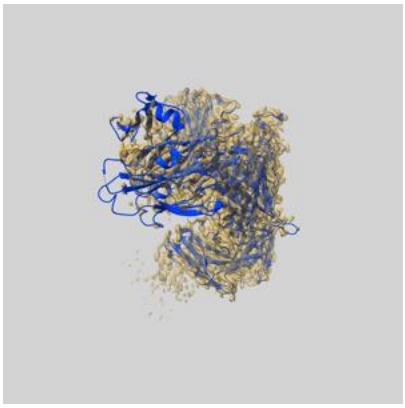
9.1 Map-model overlay



X



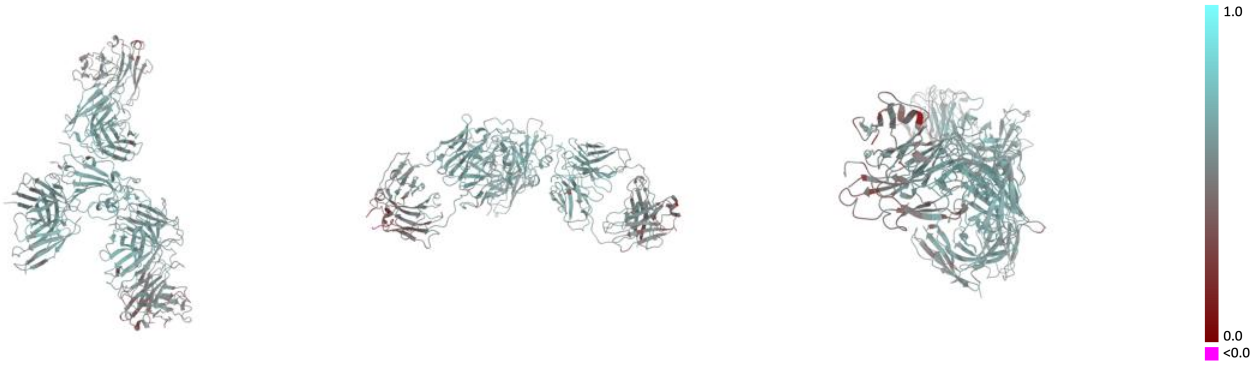
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.7 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.

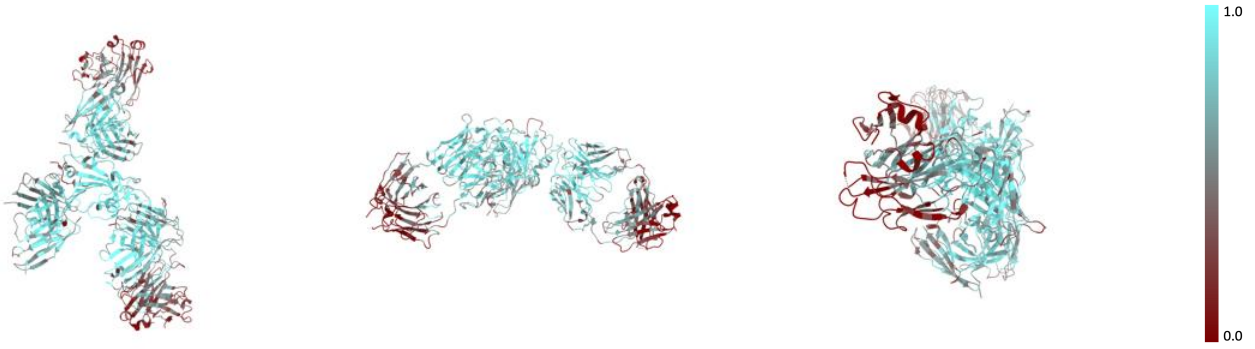
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

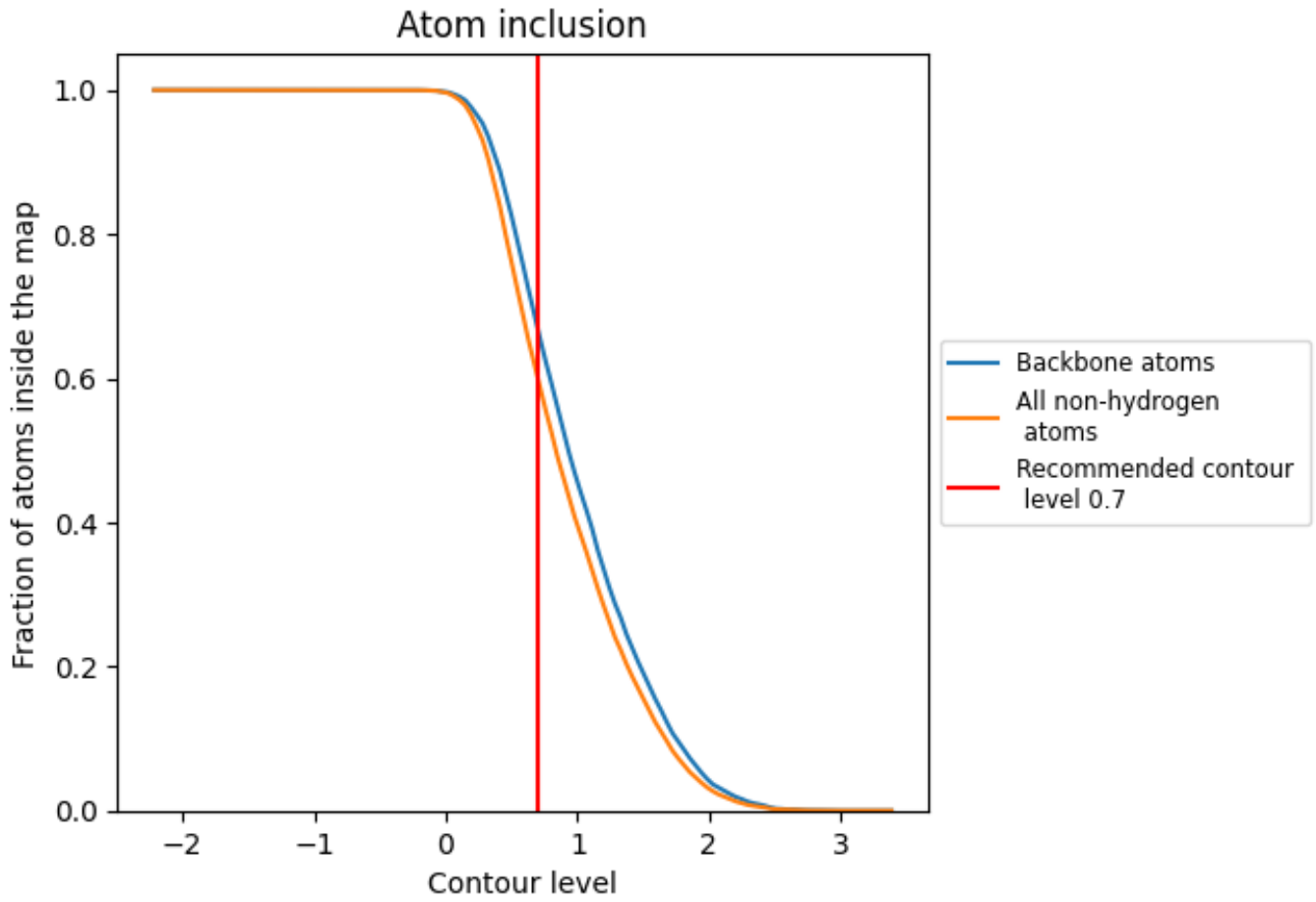
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.7).

9.4 Atom inclusion (i)



At the recommended contour level, 67% of all backbone atoms, 60% 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 (0.7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5990	0.5520
A	0.5503	0.5470
B	0.5110	0.5390
C	0.5960	0.5450
D	0.4574	0.5060
E	0.7895	0.6030
H	0.6503	0.5670
L	0.7469	0.5790
M	0.7837	0.5990