Full wwPDB EM Validation Report (i)

Nov 7, 2022 – 11:28 AM EST

PDB ID	:	6DMO
EMDB ID	:	EMD-7964
Title	:	Cryo-EM structure of human Ptch1 with three mutations L282Q/T500F/P504L
Authors	:	Yan, N.; Gong, X.; Qian, H.W.
Deposited on	:	2018-06-05
Resolution	:	4.10 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 (i) 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 (i)) were used in the production of this report:

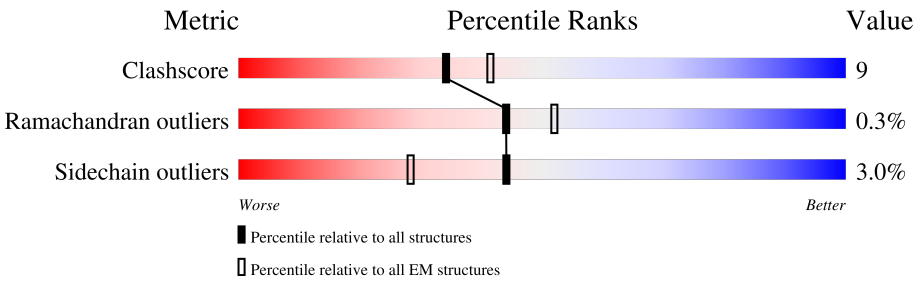
EMDB validation analysis	:	0.0.1.dev43
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 4.10 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	1349	23% 57% 14% • 27%
2	B	2	50% 50% 50%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7853 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 Protein patched homolog 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	982	7755	5061	1277	1375	42	0	0

There are 47 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP Q13635
A	-19	ALA	-	expression tag	UNP Q13635
A	-18	ASP	-	expression tag	UNP Q13635
A	-17	TYR	-	expression tag	UNP Q13635
A	-16	LYS	-	expression tag	UNP Q13635
A	-15	ASP	-	expression tag	UNP Q13635
A	-14	ASP	-	expression tag	UNP Q13635
A	-13	ASP	-	expression tag	UNP Q13635
A	-12	ASP	-	expression tag	UNP Q13635
A	-11	LYS	-	expression tag	UNP Q13635
A	-10	SER	-	expression tag	UNP Q13635
A	-9	GLY	-	expression tag	UNP Q13635
A	-8	PRO	-	expression tag	UNP Q13635
A	-7	ASP	-	expression tag	UNP Q13635
A	-6	GLU	-	expression tag	UNP Q13635
A	-5	VAL	-	expression tag	UNP Q13635
A	-4	ASP	-	expression tag	UNP Q13635
A	-3	ALA	-	expression tag	UNP Q13635
A	-2	SER	-	expression tag	UNP Q13635
A	-1	GLY	-	expression tag	UNP Q13635
A	0	ARG	-	expression tag	UNP Q13635
A	282	GLN	LEU	engineered mutation	UNP Q13635
A	500	PHE	THR	engineered mutation	UNP Q13635
A	504	LEU	PRO	engineered mutation	UNP Q13635
A	1306	LEU	-	expression tag	UNP Q13635
A	1307	GLU	-	expression tag	UNP Q13635
A	1308	GLY	-	expression tag	UNP Q13635
A	1309	SER	-	expression tag	UNP Q13635

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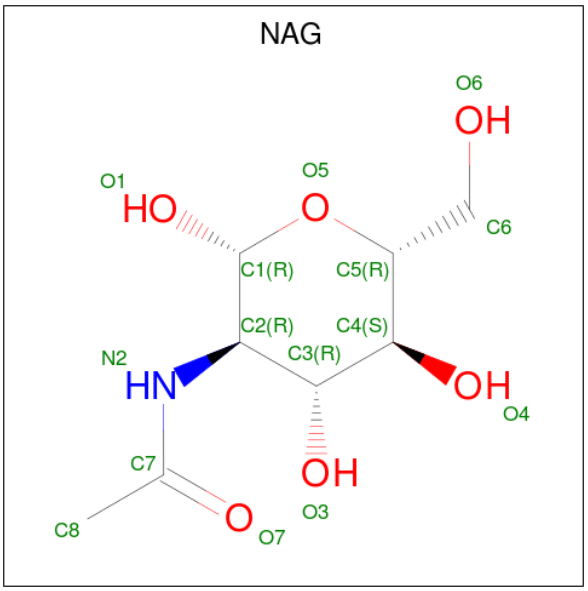
Chain	Residue	Modelled	Actual	Comment	Reference
A	1310	ASP	-	expression tag	UNP Q13635
A	1311	GLU	-	expression tag	UNP Q13635
A	1312	VAL	-	expression tag	UNP Q13635
A	1313	ASP	-	expression tag	UNP Q13635
A	1314	ALA	-	expression tag	UNP Q13635
A	1315	VAL	-	expression tag	UNP Q13635
A	1316	GLU	-	expression tag	UNP Q13635
A	1317	GLY	-	expression tag	UNP Q13635
A	1318	SER	-	expression tag	UNP Q13635
A	1319	HIS	-	expression tag	UNP Q13635
A	1320	HIS	-	expression tag	UNP Q13635
A	1321	HIS	-	expression tag	UNP Q13635
A	1322	HIS	-	expression tag	UNP Q13635
A	1323	HIS	-	expression tag	UNP Q13635
A	1324	HIS	-	expression tag	UNP Q13635
A	1325	HIS	-	expression tag	UNP Q13635
A	1326	HIS	-	expression tag	UNP Q13635
A	1327	HIS	-	expression tag	UNP Q13635
A	1328	HIS	-	expression tag	UNP Q13635

- Molecule 2 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
			Total	C	N	O		
2	B	2	28	16	2	10	0	0

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



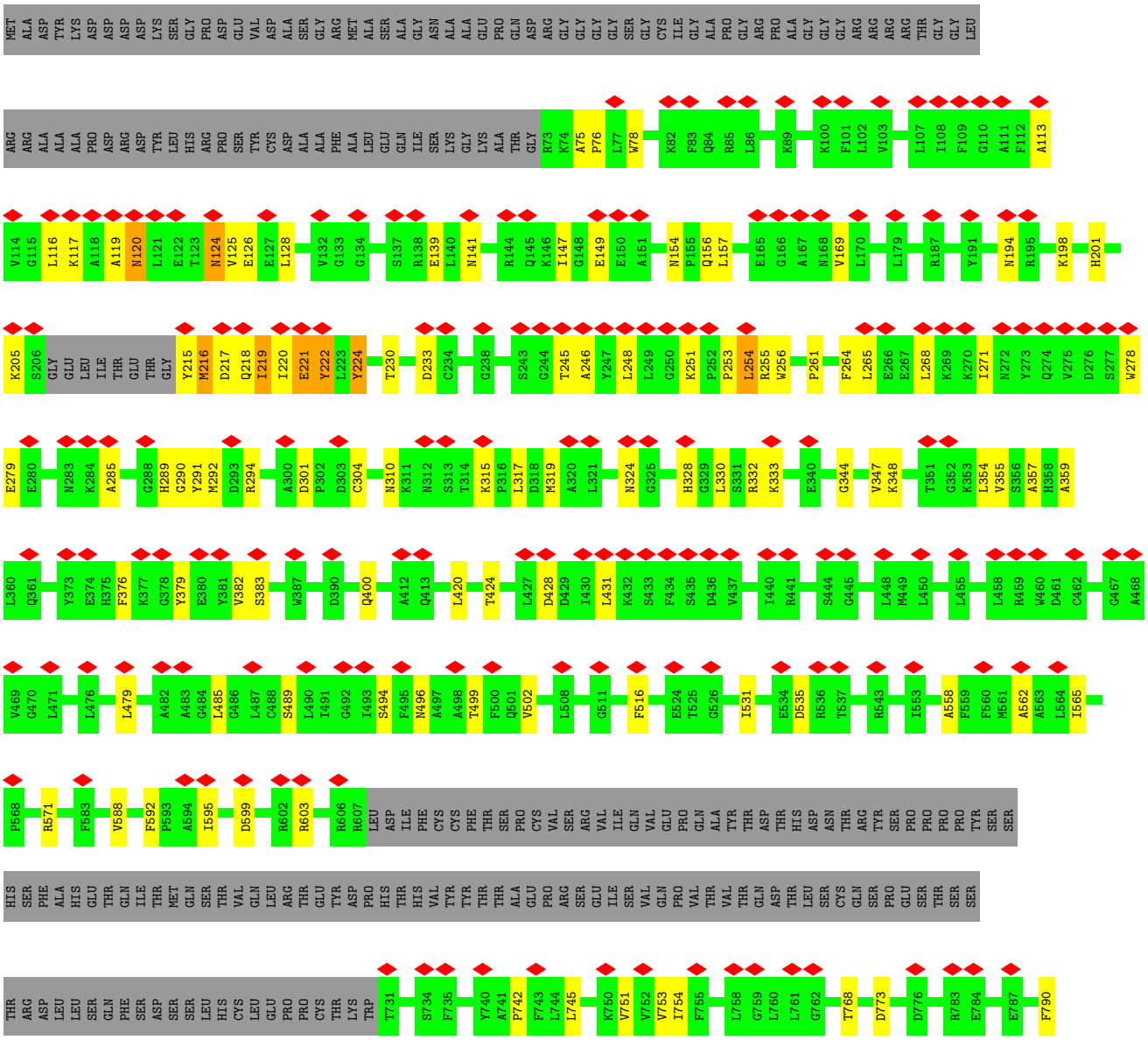
Mol	Chain	Residues	Atoms				AltConf
3	A	1	Total	C	N	O	0
			70	40	5	25	
3	A	1	Total	C	N	O	0
			70	40	5	25	
3	A	1	Total	C	N	O	0
			70	40	5	25	
3	A	1	Total	C	N	O	0
			70	40	5	25	
3	A	1	Total	C	N	O	0
			70	40	5	25	

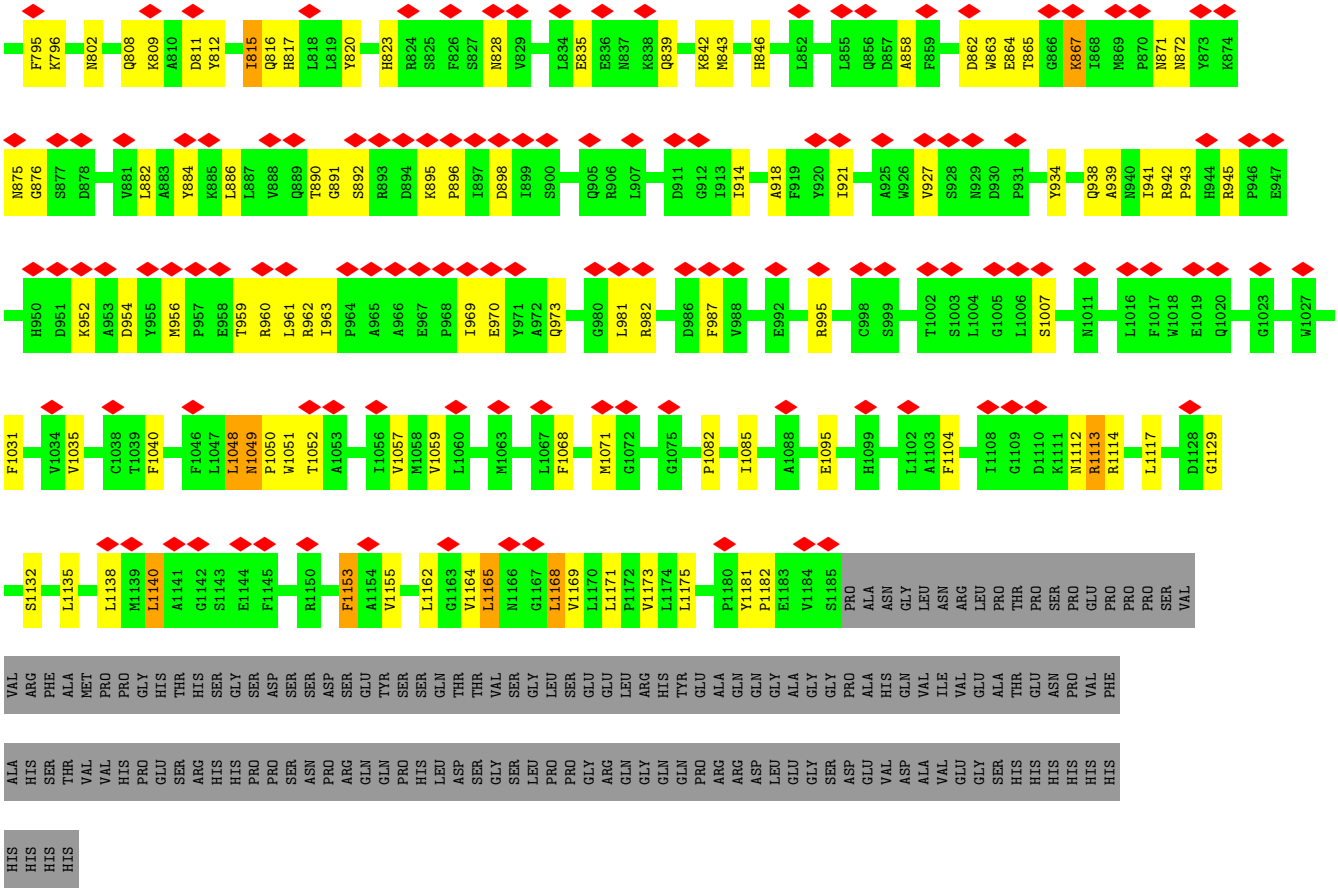
3 Residue-property plots (i)

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: Protein patched homolog 1

Chain A:





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

Chain B:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	154721	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 ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.147	Depositor
Minimum map value	-0.078	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.035	Depositor
Map size (Å)	244.38399, 244.38399, 244.38399	wwPDB
Map dimensions	224, 224, 224	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.091, 1.091, 1.091	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	0.40	0/7946	0.68	7/10804 (0.1%)

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	5

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	815	ILE	CG1-CB-CG2	-6.19	97.79	111.40
1	A	1140	LEU	CA-CB-CG	5.95	128.99	115.30
1	A	961	LEU	CA-CB-CG	5.32	127.54	115.30
1	A	224	TYR	N-CA-C	-5.27	96.78	111.00
1	A	898	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	1048	LEU	CA-CB-CG	5.08	126.99	115.30
1	A	1168	LEU	CA-CB-CG	5.06	126.94	115.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1049	ASN	Peptide
1	A	1153	PHE	Peptide
1	A	812	TYR	Peptide

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Mol	Chain	Res	Type	Group
1	A	867	LYS	Peptide
1	A	942	ARG	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	7755	0	7784	140	0
2	B	28	0	25	0	0
3	A	70	0	65	0	0
All	All	7853	0	7874	140	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 9.

All (140) 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:864:GLU:O	1:A:865:THR:HG22	1.20	1.26
1:A:864:GLU:O	1:A:865:THR:CG2	2.04	1.05
1:A:120:ASN:O	1:A:494:SER:OG	1.76	1.02
1:A:113:ALA:O	1:A:116:LEU:HD23	1.68	0.91
1:A:218:GLN:O	1:A:222:TYR:HB2	1.71	0.89
1:A:939:ALA:O	1:A:941:ILE:HD12	1.75	0.87
1:A:261:PRO:O	1:A:265:LEU:HB2	1.77	0.85
1:A:216:MET:HB3	1:A:278:TRP:CZ2	2.14	0.83
1:A:939:ALA:O	1:A:941:ILE:CD1	2.28	0.82
1:A:216:MET:HB3	1:A:278:TRP:CH2	2.16	0.80
1:A:219:ILE:HD11	1:A:246:ALA:CB	2.13	0.78
1:A:882:LEU:O	1:A:886:LEU:HB2	1.86	0.76
1:A:216:MET:SD	1:A:219:ILE:CG2	2.74	0.76
1:A:216:MET:SD	1:A:219:ILE:HG22	2.26	0.75
1:A:219:ILE:CD1	1:A:246:ALA:CB	2.66	0.74
1:A:219:ILE:HD11	1:A:246:ALA:HB3	1.70	0.73
1:A:113:ALA:O	1:A:116:LEU:CD2	2.38	0.72
1:A:943:PRO:HG2	1:A:945:ARG:HH21	1.54	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ILE:HG12	1:A:248:LEU:HD11	1.71	0.71
1:A:864:GLU:C	1:A:865:THR:HG22	2.09	0.70
1:A:941:ILE:HD12	1:A:941:ILE:N	2.06	0.70
1:A:219:ILE:HD13	1:A:246:ALA:HB1	1.73	0.69
1:A:219:ILE:HD12	1:A:219:ILE:O	1.95	0.66
1:A:939:ALA:HB1	1:A:941:ILE:HD11	1.79	0.65
1:A:154:ASN:O	1:A:156:GLN:NE2	2.29	0.65
1:A:823:HIS:ND1	1:A:843:MET:SD	2.70	0.65
1:A:943:PRO:HG2	1:A:945:ARG:NH2	2.12	0.64
1:A:941:ILE:HG23	1:A:969:ILE:HD12	1.80	0.64
1:A:219:ILE:CD1	1:A:246:ALA:HB1	2.27	0.64
1:A:938:GLN:O	1:A:973:GLN:NE2	2.32	0.62
1:A:290:GLY:O	1:A:294:ARG:NH1	2.32	0.62
1:A:216:MET:SD	1:A:219:ILE:HG21	2.40	0.61
1:A:285:ALA:HA	1:A:333:LYS:HD3	1.83	0.60
1:A:858:ALA:O	1:A:862:ASP:HB2	2.01	0.60
1:A:867:LYS:HG2	1:A:871:ASN:HD22	1.66	0.60
1:A:294:ARG:NH2	1:A:328:HIS:O	2.35	0.60
1:A:348:LYS:HA	1:A:354:LEU:HA	1.83	0.60
1:A:796:LYS:HA	1:A:982:ARG:HH22	1.65	0.60
1:A:795:PHE:O	1:A:982:ARG:NH2	2.35	0.59
1:A:876:GLY:O	1:A:952:LYS:NZ	2.35	0.59
1:A:1171:LEU:O	1:A:1175:LEU:HB2	2.02	0.58
1:A:943:PRO:HD2	1:A:969:ILE:HG22	1.86	0.57
1:A:428:ASP:N	1:A:428:ASP:OD1	2.37	0.57
1:A:1114:ARG:HH12	1:A:1181:TYR:HA	1.70	0.56
1:A:809:LYS:HA	1:A:970:GLU:HG2	1.87	0.56
1:A:169:VAL:N	1:A:357:ALA:O	2.37	0.56
1:A:376:PHE:HB3	1:A:382:VAL:HG21	1.85	0.56
1:A:599:ASP:OD2	1:A:603:ARG:NH2	2.39	0.56
1:A:939:ALA:C	1:A:941:ILE:HD12	2.26	0.56
1:A:1165:LEU:O	1:A:1169:VAL:HB	2.05	0.55
1:A:981:LEU:HB3	1:A:987:PHE:HE1	1.72	0.54
1:A:768:THR:HG22	1:A:1068:PHE:HD2	1.73	0.54
1:A:1040:PHE:HB2	1:A:1057:VAL:HG21	1.89	0.54
1:A:1068:PHE:HA	1:A:1071:MET:HE2	1.90	0.54
1:A:943:PRO:CD	1:A:969:ILE:HA	2.38	0.54
1:A:496:ASN:HB2	1:A:499:THR:HG22	1.90	0.53
1:A:124:ASN:ND2	1:A:126:GLU:OE2	2.41	0.53
1:A:400:GLN:HE22	1:A:424:THR:HG22	1.73	0.53
1:A:753:VAL:HG21	1:A:1173:VAL:HG21	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:MET:HB3	1:A:278:TRP:CE2	2.45	0.52
1:A:1049:ASN:O	1:A:1051:TRP:N	2.42	0.52
1:A:835:GLU:HG3	1:A:839:GLN:HB2	1.92	0.51
1:A:927:VAL:HG11	1:A:945:ARG:HG3	1.91	0.51
1:A:310:ASN:CG	1:A:317:LEU:HD11	2.30	0.51
1:A:891:GLY:HA3	1:A:918:ALA:HB1	1.93	0.51
1:A:125:VAL:HG23	1:A:128:LEU:HD22	1.93	0.51
1:A:808:GLN:O	1:A:1007:SER:OG	2.29	0.50
1:A:956:MET:O	1:A:959:THR:OG1	2.30	0.50
1:A:1164:VAL:O	1:A:1168:LEU:HB3	2.12	0.50
1:A:954:ASP:OD1	1:A:962:ARG:NH1	2.45	0.49
1:A:344:GLY:N	1:A:359:ALA:O	2.46	0.48
1:A:301:ASP:HB3	1:A:304:CYS:HB2	1.96	0.48
1:A:289:HIS:HB3	1:A:292:MET:HB2	1.95	0.48
1:A:941:ILE:CD1	1:A:941:ILE:N	2.73	0.47
1:A:565:ILE:O	1:A:571:ARG:NE	2.47	0.47
1:A:256:TRP:HZ2	1:A:330:LEU:HD21	1.79	0.47
1:A:817:HIS:HA	1:A:820:TYR:HD2	1.79	0.47
1:A:1113:ARG:HH22	1:A:1117:LEU:HB2	1.78	0.47
1:A:420:LEU:HB3	1:A:790:PHE:HD1	1.79	0.47
1:A:219:ILE:HD13	1:A:246:ALA:CB	2.39	0.47
1:A:941:ILE:HD12	1:A:941:ILE:H	1.79	0.47
1:A:245:THR:HG22	1:A:255:ARG:HE	1.80	0.46
1:A:943:PRO:HD2	1:A:969:ILE:HA	1.98	0.46
1:A:198:LYS:H	1:A:201:HIS:CE1	2.33	0.46
1:A:230:THR:HB	1:A:233:ASP:HB2	1.97	0.46
1:A:742:PRO:HA	1:A:745:LEU:HB2	1.97	0.46
1:A:216:MET:CE	1:A:264:PHE:HE2	2.29	0.45
1:A:558:ALA:HB1	1:A:1035:VAL:HG13	1.98	0.45
1:A:1049:ASN:ND2	1:A:1052:THR:OG1	2.49	0.45
1:A:265:LEU:HD21	1:A:279:GLU:HA	1.98	0.45
1:A:379:TYR:O	1:A:383:SER:N	2.49	0.45
1:A:139:GLU:OE2	1:A:995:ARG:NH1	2.30	0.45
1:A:1140:LEU:HB3	1:A:1153:PHE:HZ	1.81	0.45
1:A:531:ILE:HG23	1:A:535:ASP:HB3	1.99	0.45
1:A:215:TYR:CD1	1:A:215:TYR:O	2.70	0.45
1:A:1082:PRO:HA	1:A:1085:ILE:HG12	1.98	0.44
1:A:872:ASN:HD22	1:A:884:TYR:HB2	1.83	0.44
1:A:219:ILE:HD13	1:A:219:ILE:HA	1.75	0.44
1:A:268:LEU:HA	1:A:271:ILE:HG22	1.99	0.43
1:A:811:ASP:HB3	1:A:815:ILE:HD13	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:PRO:HG3	1:A:1048:LEU:HG	2.01	0.43
1:A:842:LYS:HB3	1:A:846:HIS:HB2	2.00	0.43
1:A:216:MET:HB3	1:A:278:TRP:CZ3	2.53	0.43
1:A:890:THR:HG23	1:A:892:SER:H	1.82	0.43
1:A:221:GLU:OE2	1:A:221:GLU:HA	2.19	0.43
1:A:863:TRP:HE1	1:A:876:GLY:HA3	1.83	0.43
1:A:1095:GLU:H	1:A:1095:GLU:HG3	1.67	0.43
1:A:347:VAL:HG23	1:A:355:VAL:HG23	2.00	0.43
1:A:1135:LEU:HA	1:A:1138:LEU:HD22	1.99	0.43
1:A:291:TYR:HA	1:A:294:ARG:HH11	1.84	0.43
1:A:816:GLN:HB2	1:A:914:ILE:HG22	2.00	0.43
1:A:588:VAL:HG13	1:A:592:PHE:HD2	1.82	0.43
1:A:251:LYS:HB3	1:A:254:LEU:HD23	2.01	0.42
1:A:310:ASN:ND2	1:A:317:LEU:HD11	2.33	0.42
1:A:1104:PHE:CG	1:A:1175:LEU:HD23	2.54	0.42
1:A:219:ILE:CG2	1:A:220:ILE:HG12	2.49	0.42
1:A:934:TYR:OH	1:A:941:ILE:HD13	2.18	0.42
1:A:1059:VAL:HG21	1:A:1171:LEU:HD21	2.00	0.42
1:A:499:THR:HA	1:A:502:VAL:HB	2.01	0.42
1:A:1085:ILE:HG21	1:A:1155:VAL:HG12	2.02	0.42
1:A:773:ASP:OD1	1:A:773:ASP:N	2.53	0.42
1:A:428:ASP:HA	1:A:431:LEU:HB2	2.01	0.41
1:A:751:VAL:HG13	1:A:754:ILE:HD12	2.01	0.41
1:A:75:ALA:HA	1:A:78:TRP:NE1	2.35	0.41
1:A:147:ILE:HG23	1:A:149:GLU:H	1.85	0.41
1:A:921:ILE:HG13	1:A:963:ILE:HG21	2.01	0.41
1:A:943:PRO:HD2	1:A:969:ILE:CB	2.50	0.41
1:A:485:LEU:HD23	1:A:485:LEU:HA	1.89	0.41
1:A:592:PHE:HA	1:A:595:ILE:HG22	2.03	0.41
1:A:941:ILE:CD1	1:A:941:ILE:H	2.33	0.41
1:A:116:LEU:HA	1:A:116:LEU:HD13	1.80	0.41
1:A:119:ALA:HB2	1:A:489:SER:HB2	2.03	0.41
1:A:981:LEU:HB3	1:A:987:PHE:CE1	2.54	0.41
1:A:1129:GLY:O	1:A:1132:SER:OG	2.30	0.41
1:A:205:LYS:HD2	1:A:224:TYR:HB3	2.01	0.41
1:A:261:PRO:O	1:A:265:LEU:CB	2.58	0.41
1:A:858:ALA:O	1:A:862:ASP:CB	2.68	0.40
1:A:895:LYS:N	1:A:896:PRO:HD3	2.36	0.40
1:A:562:ALA:HB1	1:A:1031:PHE:CZ	2.56	0.40
1:A:516:PHE:HD1	1:A:516:PHE:HA	1.72	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	976/1349 (72%)	865 (89%)	108 (11%)	3 (0%)	41 75

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1182	PRO
1	A	1050	PRO
1	A	253	PRO

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	826/1147 (72%)	801 (97%)	25 (3%)	41 64

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	117	LYS
1	A	120	ASN
1	A	124	ASN
1	A	141	ASN
1	A	157	LEU
1	A	194	ASN
1	A	216	MET
1	A	217	ASP

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Mol	Chain	Res	Type
1	A	219	ILE
1	A	221	GLU
1	A	222	TYR
1	A	254	LEU
1	A	315	LYS
1	A	319	MET
1	A	324	ASN
1	A	332	ARG
1	A	479	LEU
1	A	802	ASN
1	A	828	ASN
1	A	875	ASN
1	A	960	ARG
1	A	1112	ASN
1	A	1113	ARG
1	A	1162	LEU
1	A	1165	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	GLN
1	A	124	ASN
1	A	160	GLN
1	A	178	HIS
1	A	194	ASN
1	A	196	GLN
1	A	282	GLN
1	A	283	ASN
1	A	324	ASN
1	A	375	HIS
1	A	417	GLN
1	A	828	ASN
1	A	853	GLN
1	A	901	GLN
1	A	929	ASN
1	A	1121	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)

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	$\#\|Z\| > 2$
2	NAG	B	1	2,1	14,14,15	0.33	0	17,19,21	0.68	1 (5%)
2	NAG	B	2	2	14,14,15	0.32	0	17,19,21	0.53	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
2	NAG	B	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1	NAG	C1-O5-C5	2.35	115.38	112.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	NAG	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6

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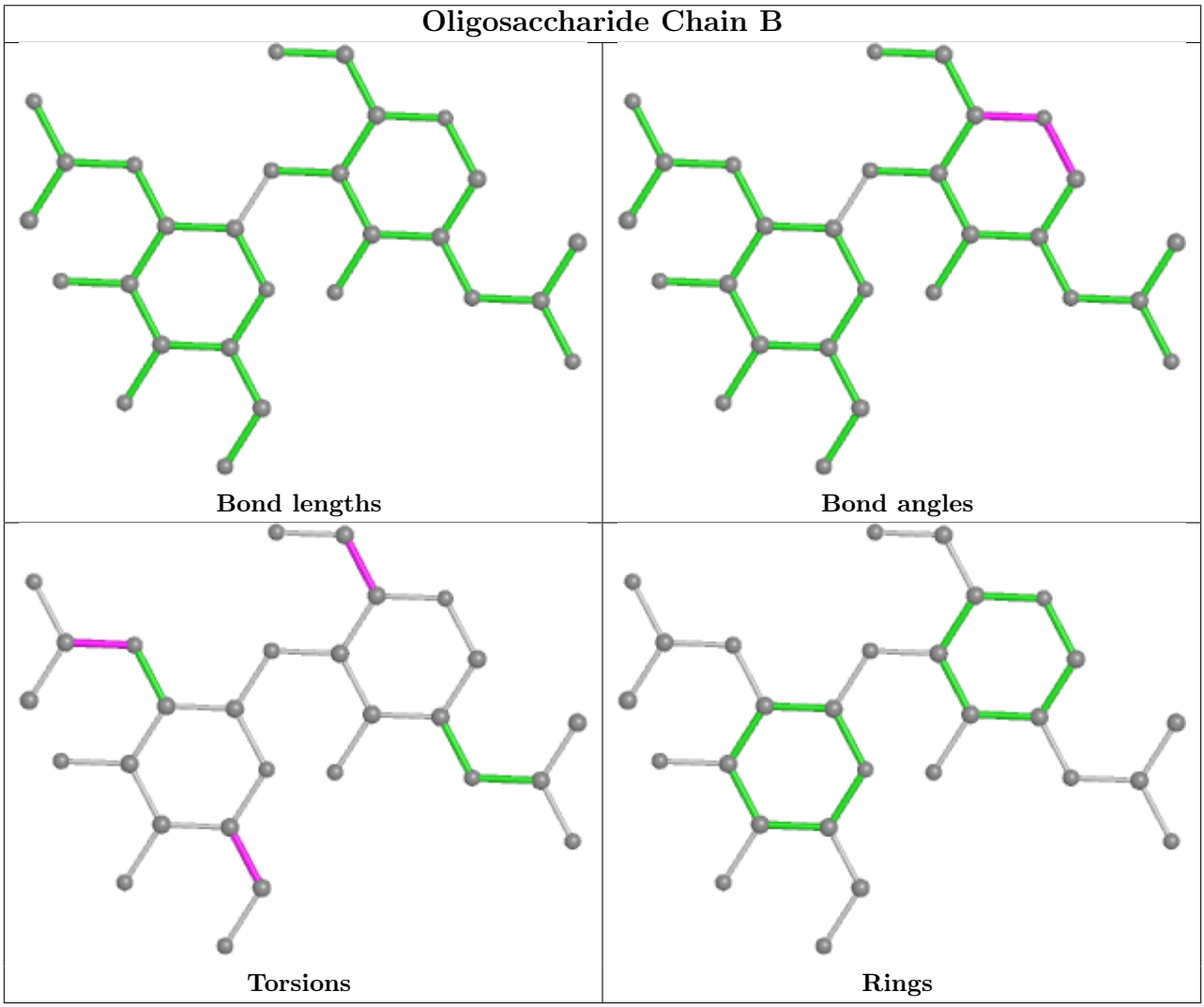
Mol	Chain	Res	Type	Atoms
2	B	2	NAG	O5-C5-C6-O6
2	B	2	NAG	C8-C7-N2-C2
2	B	2	NAG	O7-C7-N2-C2
2	B	2	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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)

5 ligands 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	$\#\|Z\| > 2$
3	NAG	A	1802	1	14,14,15	0.18	0	17,19,21	0.48	0
3	NAG	A	1801	1	14,14,15	0.62	0	17,19,21	0.78	1 (5%)
3	NAG	A	1804	1	14,14,15	0.55	0	17,19,21	0.40	0
3	NAG	A	1803	1	14,14,15	0.33	0	17,19,21	0.68	1 (5%)
3	NAG	A	1807	1	14,14,15	0.32	0	17,19,21	0.58	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	1802	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1801	1	-	4/6/23/26	0/1/1/1
3	NAG	A	1804	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1803	1	-	4/6/23/26	0/1/1/1
3	NAG	A	1807	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1801	NAG	C1-O5-C5	2.41	115.46	112.19
3	A	1803	NAG	C1-O5-C5	2.17	115.13	112.19

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1803	NAG	C4-C5-C6-O6
3	A	1801	NAG	O5-C5-C6-O6
3	A	1802	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	A	1801	NAG	C4-C5-C6-O6
3	A	1803	NAG	O5-C5-C6-O6
3	A	1802	NAG	O5-C5-C6-O6
3	A	1807	NAG	C4-C5-C6-O6
3	A	1801	NAG	C8-C7-N2-C2
3	A	1801	NAG	O7-C7-N2-C2
3	A	1803	NAG	C8-C7-N2-C2
3	A	1803	NAG	O7-C7-N2-C2
3	A	1807	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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.

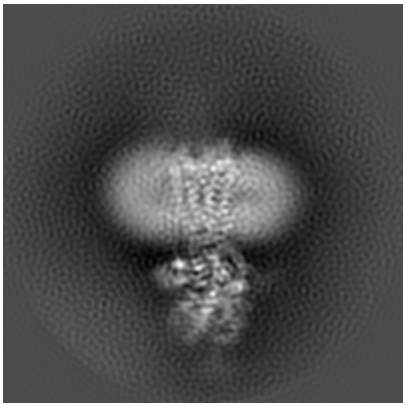
6 Map visualisation (i)

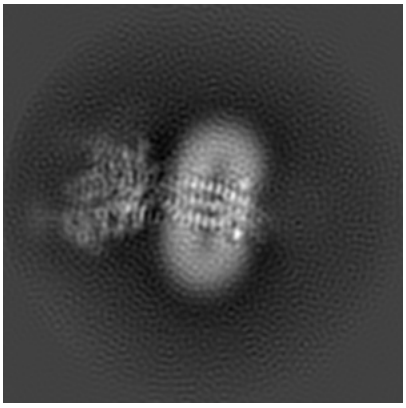
This section contains visualisations of the EMDB entry EMD-7964. These allow visual inspection of the internal detail of the map and identification of artifacts.

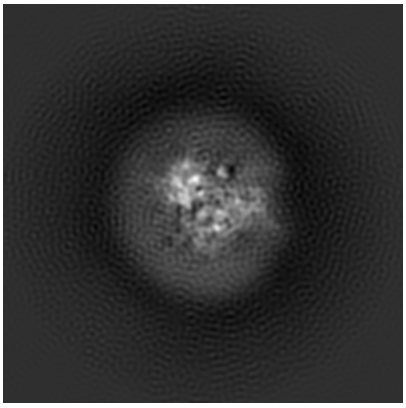
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

6.1.1 Primary map


X

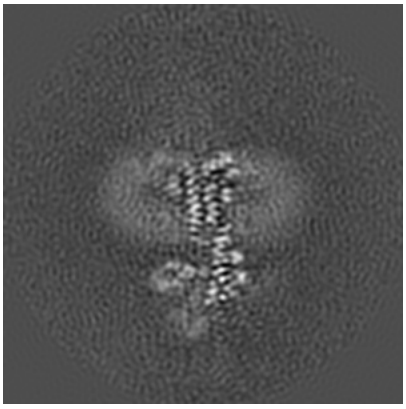

Y

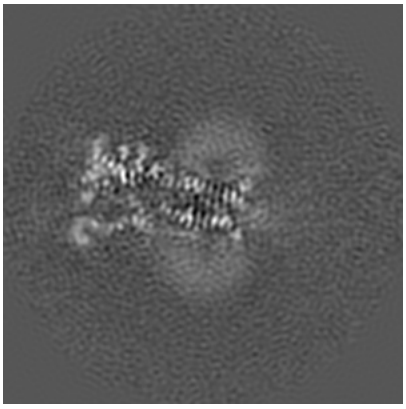

Z

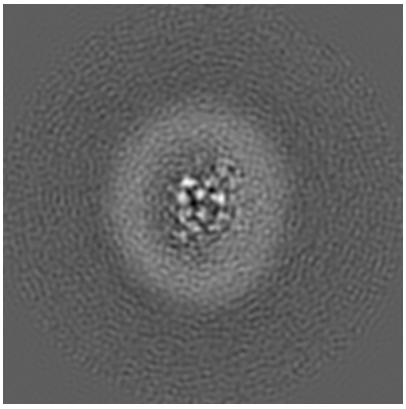
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

6.2.1 Primary map


X Index: 112


Y Index: 112

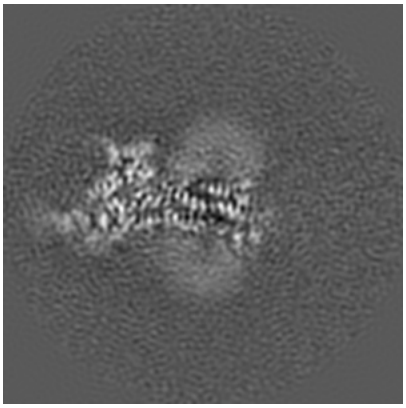

Z Index: 112

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

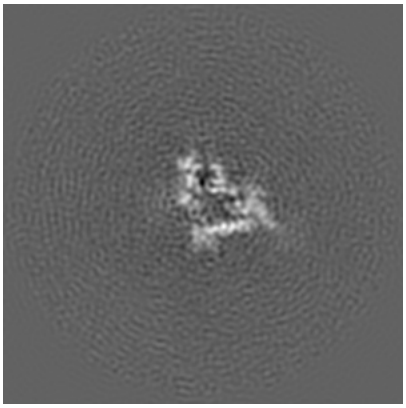
6.3 Largest variance slices (i)

6.3.1 Primary map

X Index: 108



Y Index: 118



Z Index: 77

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

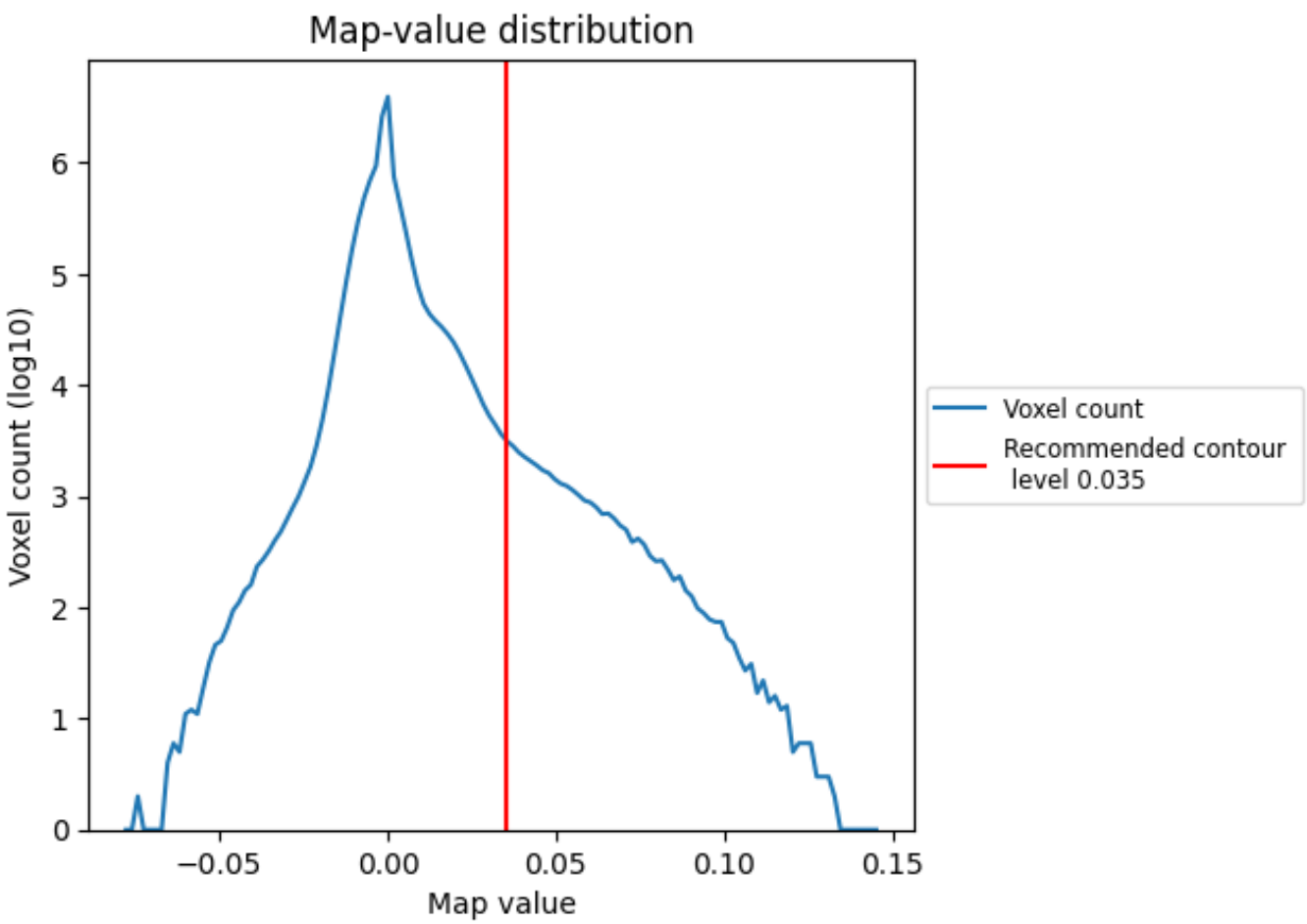
6.5 Mask visualisation (i)

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

7 Map analysis (i)

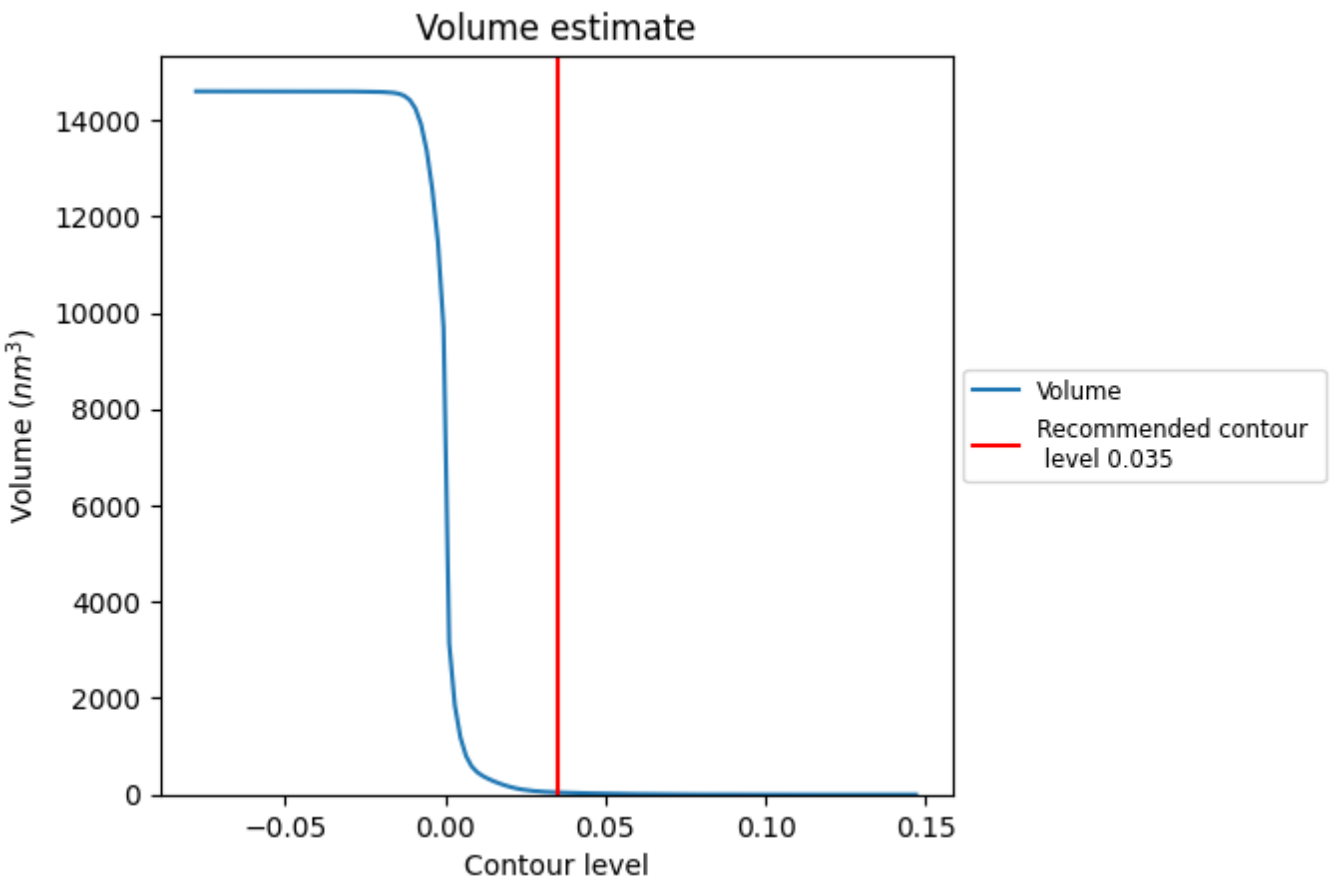
This section contains the results of statistical analysis of the map.

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.

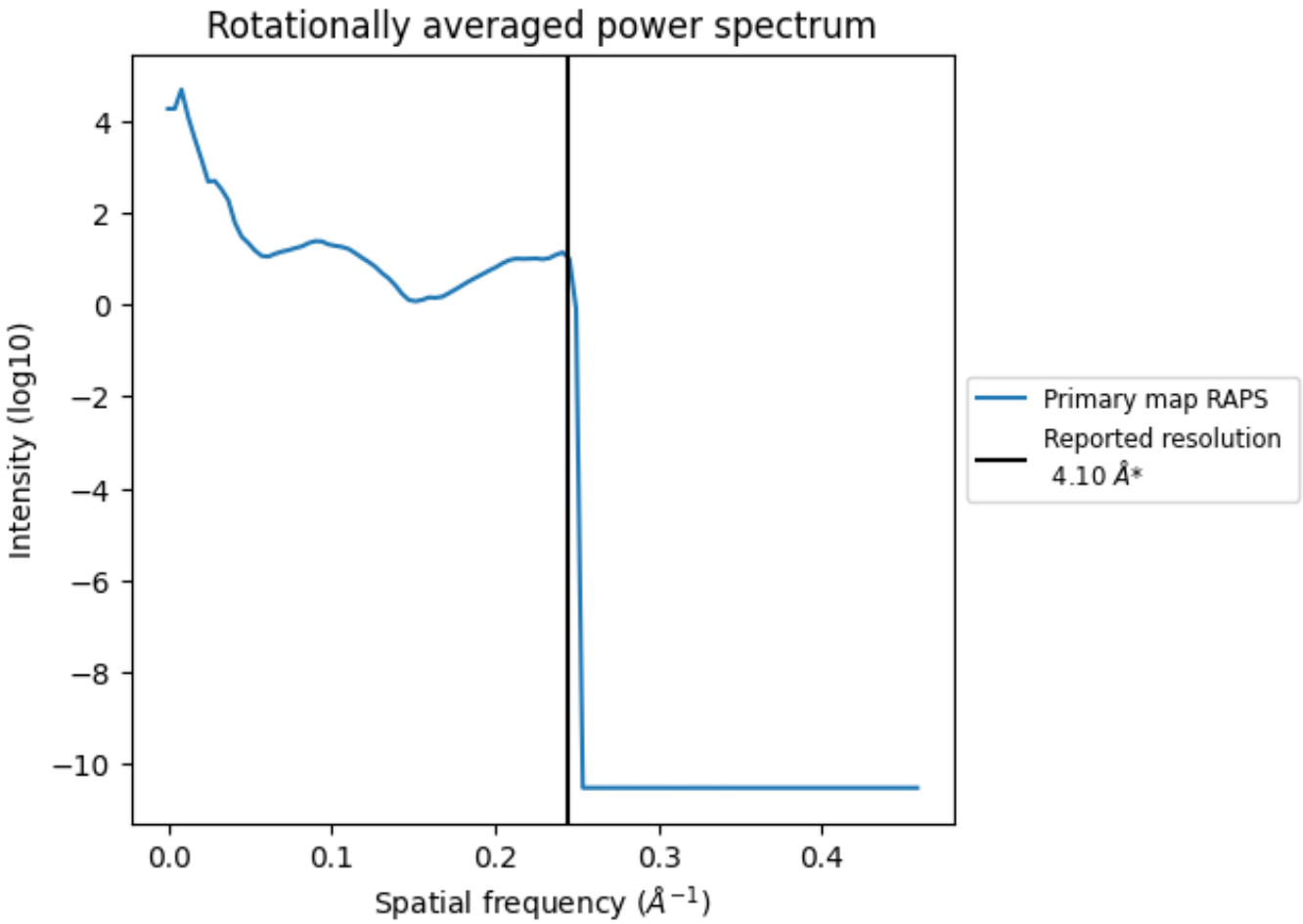
7.2 Volume estimate (i)



The volume at the recommended contour level is 45 nm^3; this corresponds to an approximate mass of 40 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 0.244 Å^{-1}

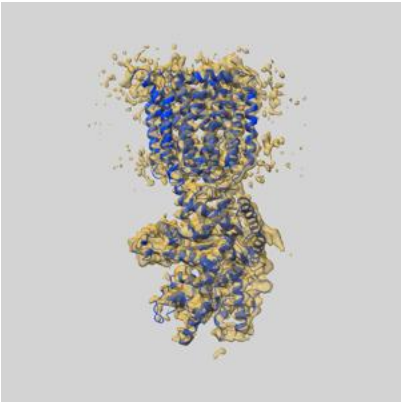
8 Fourier-Shell correlation (i)

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

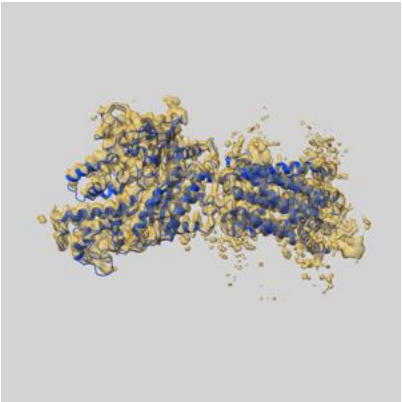
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-7964 and PDB model 6DMO. Per-residue inclusion information can be found in section 3 on page 6.

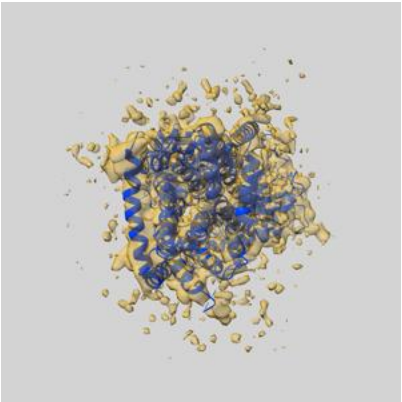
9.1 Map-model overlay (i)



X



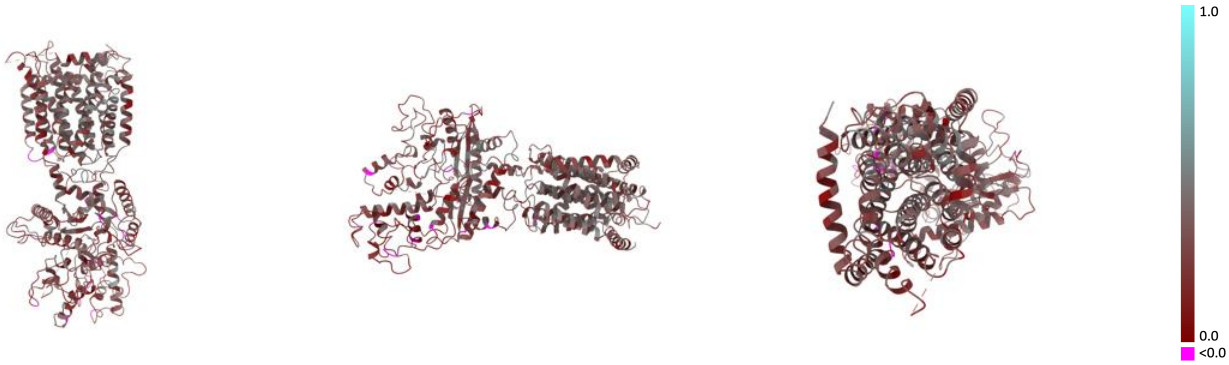
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.035 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.

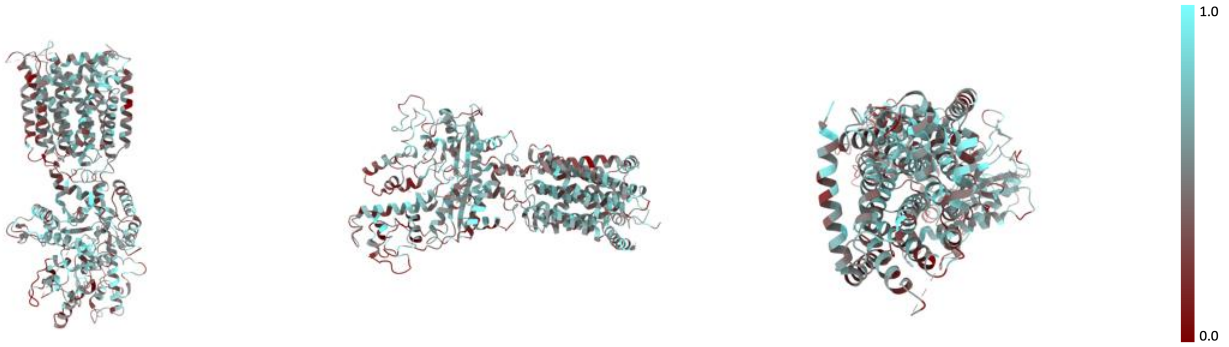
9.2 Q-score mapped to coordinate model (i)



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.

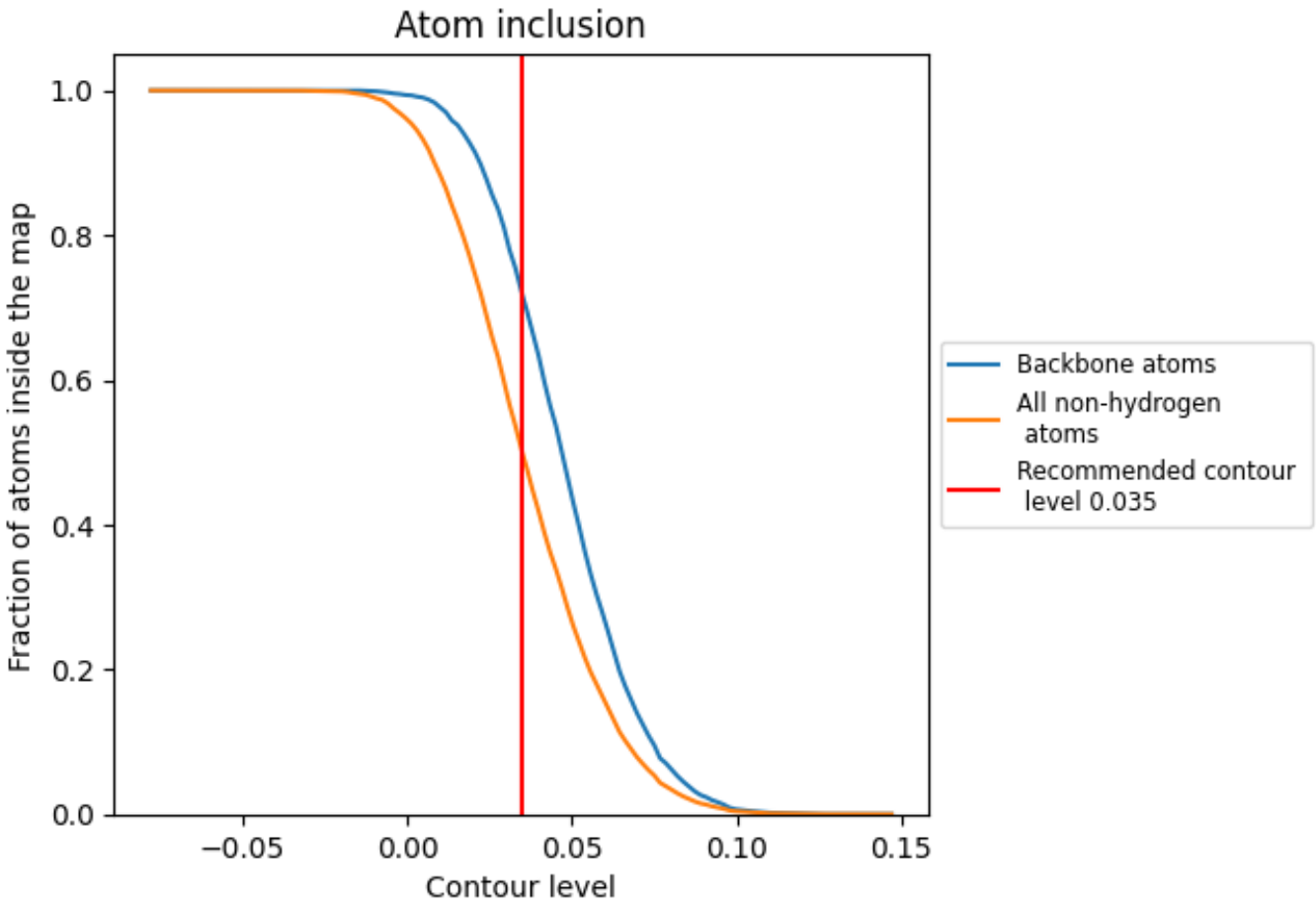
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.035).

9.4 Atom inclusion (i)



At the recommended contour level, 72% of all backbone atoms, 50% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.035) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.4987	0.2790
A	0.4991	0.2790
B	0.3929	0.2680