



Full wwPDB EM Validation Report ⓘ

Apr 12, 2026 – 10:54 AM UTC

PDB ID : 9FYQ / pdb_00009fyq
EMDB ID : EMD-50889
Title : Cryo-EM structure of native SV2A in complex with TeNT-Hc, gangliosides and Pro-Macrobody 5
Authors : Schenck, S.; Brunner, J.D.
Deposited on : 2024-07-03
Resolution : 3.25 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

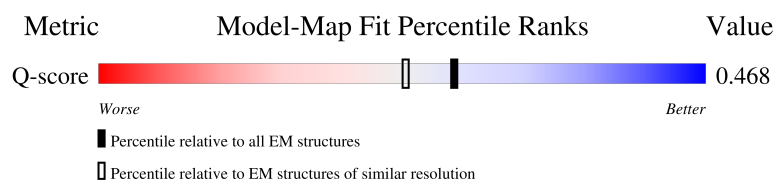
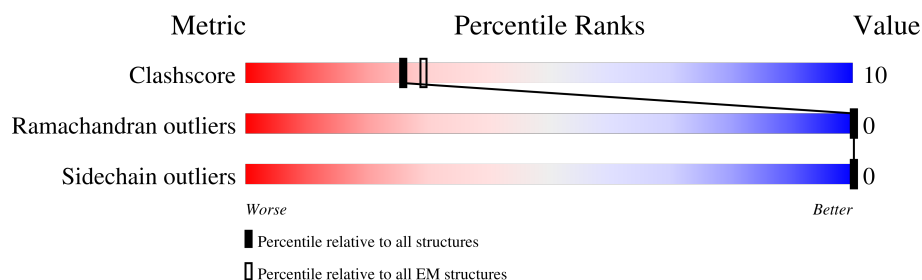
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.25 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	14599 (2.75 - 3.75)

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	742	 8% 65% 11% 23%
2	C	469	 8% 58% 33% 9%
3	D	490	 23% 75%
4	F	3	 33% 33% 67%

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Mol	Chain	Length	Quality of chain
4	I	3	33%33%67%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 14368 atoms, of which 5187 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 Synaptic vesicle glycoprotein 2A.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	570	Total	C	H	N	O	S	0	0
			8590	2929	4111	734	782	34		

- Molecule 2 is a protein called Tetanus toxin heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	429	Total	C	N	O	S	0	0
			3468	2225	581	653	9		

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	847	MET	-	initiating methionine	UNP P04958
C	848	GLY	-	expression tag	UNP P04958
C	849	LYS	-	expression tag	UNP P04958
C	850	HIS	-	expression tag	UNP P04958
C	851	HIS	-	expression tag	UNP P04958
C	852	HIS	-	expression tag	UNP P04958
C	853	HIS	-	expression tag	UNP P04958
C	854	HIS	-	expression tag	UNP P04958
C	855	HIS	-	expression tag	UNP P04958
C	856	HIS	-	expression tag	UNP P04958
C	857	HIS	-	expression tag	UNP P04958
C	858	HIS	-	expression tag	UNP P04958
C	859	HIS	-	expression tag	UNP P04958
C	860	GLY	-	expression tag	UNP P04958
C	861	SER	-	expression tag	UNP P04958
C	862	ALA	-	expression tag	UNP P04958
C	863	SER	-	expression tag	UNP P04958
C	864	LEU	-	expression tag	UNP P04958
C	865	GLU	-	expression tag	UNP P04958
C	866	VAL	-	expression tag	UNP P04958
C	867	LEU	-	expression tag	UNP P04958

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Chain	Residue	Modelled	Actual	Comment	Reference
C	868	PHE	-	expression tag	UNP P04958
C	869	GLN	-	expression tag	UNP P04958
C	870	GLY	-	expression tag	UNP P04958
C	871	PRO	-	expression tag	UNP P04958
C	872	SER	-	expression tag	UNP P04958
C	873	HIS	-	expression tag	UNP P04958
C	874	MET	-	expression tag	UNP P04958

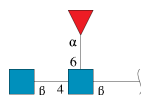
- Molecule 3 is a protein called Pro-Macrobody 5,Maltose/maltodextrin-binding periplasmic protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	D	122	Total	C	H	N	O	S	0	0
			1836	595	889	162	185	5		

There are 3 discrepancies between the modelled and reference sequences:

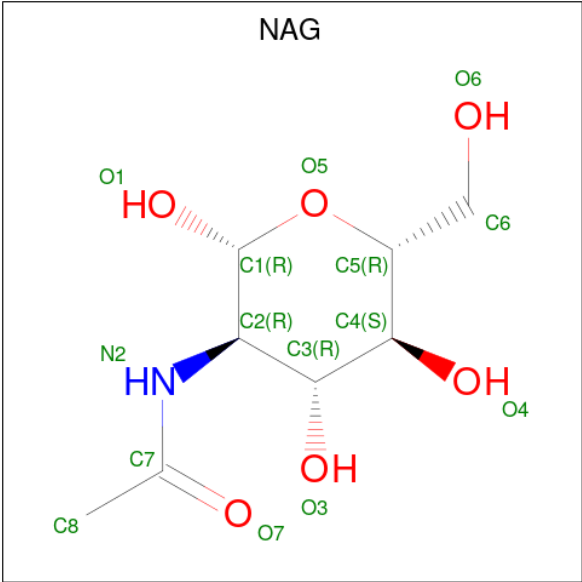
Chain	Residue	Modelled	Actual	Comment	Reference
D	487	PRO	-	expression tag	UNP P0AEX9
D	488	GLY	-	expression tag	UNP P0AEX9
D	489	ALA	-	expression tag	UNP P0AEX9

- Molecule 4 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.



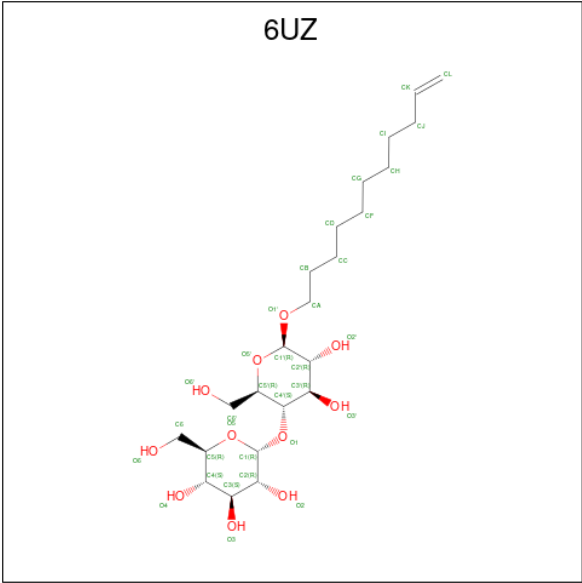
Mol	Chain	Residues	Atoms				AltConf	Trace
4	F	3	Total	C	N	O	0	0
			38	22	2	14		
4	I	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	

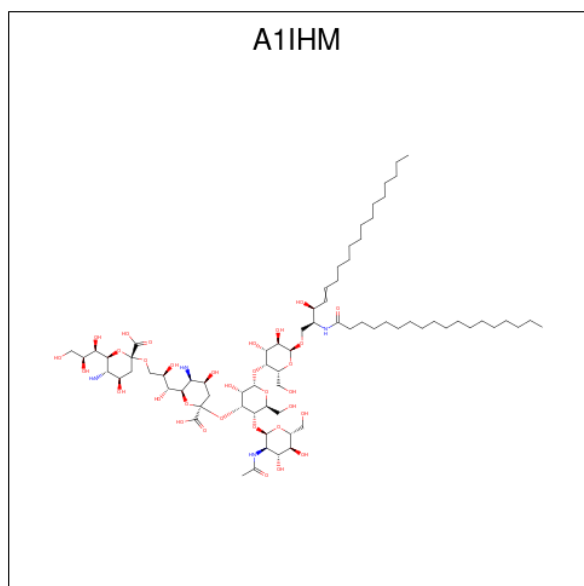
- Molecule 6 is omega-undecylenyl-beta-D-maltopyranoside (CCD ID: 6UZ) (formula: $C_{23}H_{42}O_{11}$).



Mol	Chain	Residues	Atoms				AltConf
6	A	1	Total	C	H	O	0
			76	23	42	11	

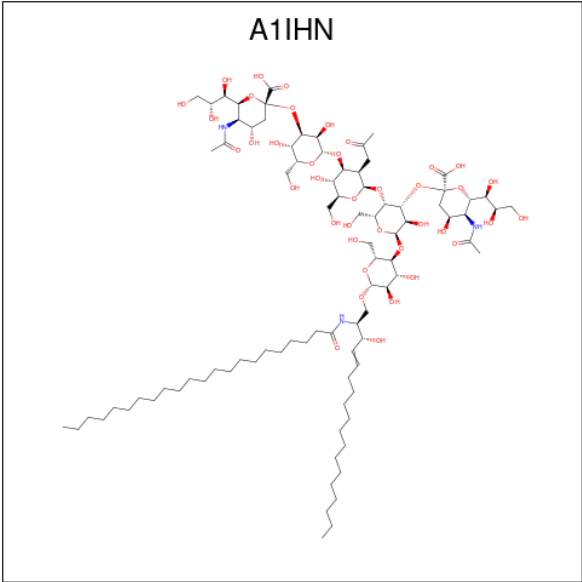
- Molecule 7 is (2 {R},4 {R},5 {S},6 {S})-2-[(2 {R},3 {R})-3-[(2 {R},3 {S},4 {S},6 {S})-6-[(2 {S},3 {S},4 {R},5 {S},6 {S})-3-[(2 {R},3 {R},4 {R},5 {S},6 {R})-3-acetamido-6-(hydroxy

methyl)-4,5-bis(oxidanyl)oxan-2-yl]oxy-2-(hydroxymethyl)-6-[(2 {R},3 {R},4 {R},5 {R},6 {S})-2-(hydroxymethyl)-6-[2-(octadecanoylamino)-3-oxidanyl-octadec-4-enoy]-4,5-bis(oxidanyl)oxan-3-yl]oxy-5-oxidanyl-oxan-4-yl]oxy-3-azanyl-6-carboxy-4-oxidanyl-oxan-2-yl]-2,3-bis(oxidanyl)propoxy]-5-azanyl-4-oxidanyl-6-[(1 {R},2 {S})-1,2,3-tris(oxidanyl)propyl]oxane-2-carboxylic acid (CCD ID: A1IHM) (formula: $C_{74}H_{134}N_4O_{32}$) (labeled as "Ligand of Interest" by depositor).

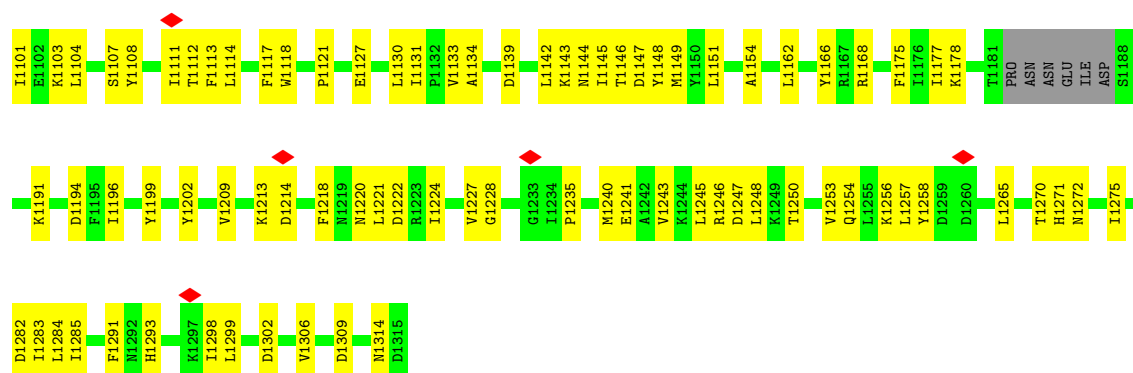


Mol	Chain	Residues	Atoms					AltConf
7	C	1	Total	C	H	N	O	0
			138	40	65	3	30	

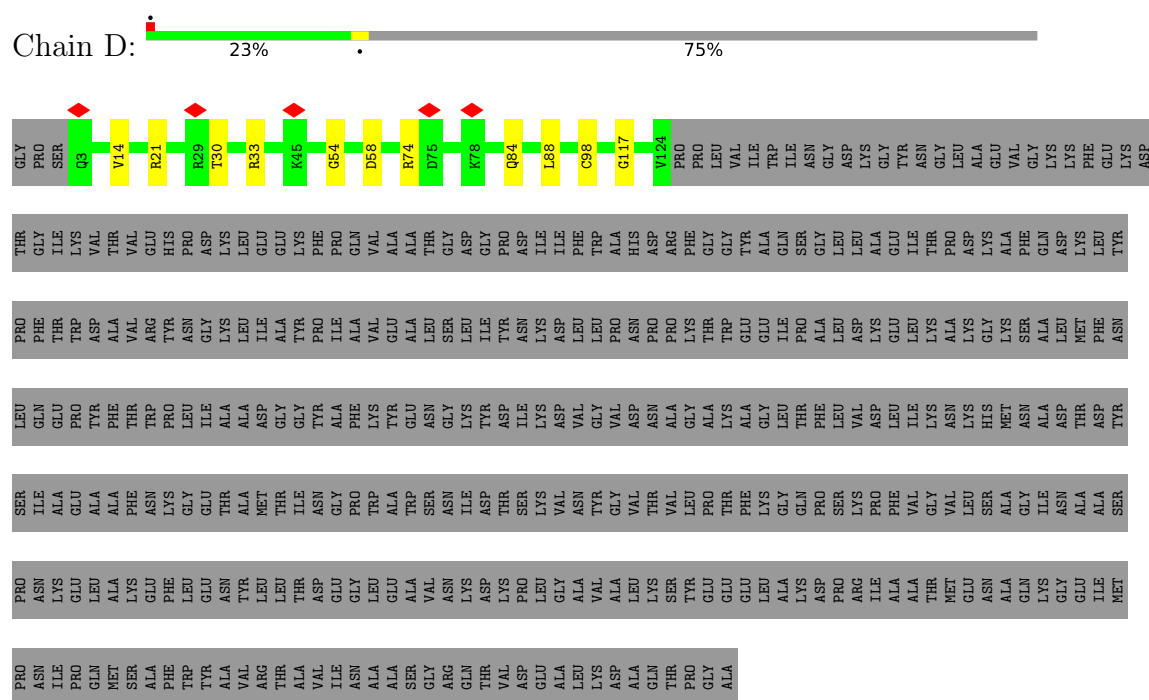
- Molecule 8 is (2 {S},4 {S},5 {R},6 {S})-5-acetamido-2-[(2 {R},3 {R},4 {R},5 {S},6 {R})-2-[(2 {R},3 {R},4 {S},5 {R},6 {S})-2-[(2 {R},3 {S},4 {R},5 {R},6 {R})-4-[(2 {S},4 {S},5 {S},6 {S})-5-acetamido-2-carboxy-4-oxidanyl-6-[(1 {S},2 {R})-1,2,3-tris(oxidanyl)propyl]oxan-2-yl]oxy-6-[(2 {R},3 {S},4 {R},5 {R},6 {R})-6-[2-(docosanoylamino)-3-oxidanyl-octadec-4-enoy]-2-(hydroxymethyl)-4,5-bis(oxidanyl)oxan-3-yl]oxy-2-(hydroxymethyl)-5-oxidanyl-oxan-3-yl]oxy-6-(hydroxymethyl)-5-oxidanyl-3-(2-oxidanylidene-propyl)oxan-4-yl]oxy-6-(hydroxymethyl)-3,5-bis(oxidanyl)oxan-4-yl]oxy-4-oxidanyl-6-[(1 {R},2 {R})-1,2,3-tris(oxidanyl)propyl]oxane-2-carboxylic acid (CCD ID: A1IHN) (formula: $C_{89}H_{157}N_3O_{39}$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	H	N	O	
8	C	1	170	51	80	2	37	0



- Molecule 3: Pro-Macrobody 5, Maltose/maltodextrin-binding periplasmic protein



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



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



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	89713	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	JEOL CRYO ARM 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	60000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.330	Depositor
Minimum map value	-0.161	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.0657	Depositor
Map size (Å)	318.08, 318.08, 318.08	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.71, 0.71, 0.71	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: A1IHM, A1IHN, FUC, NAG, 6UZ

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	0.20	0/4598	0.37	0/6222
2	C	0.18	0/3545	0.36	0/4806
3	D	0.21	0/969	0.40	0/1309
All	All	0.19	0/9112	0.37	0/12337

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4479	4111	4426	53	0
2	C	3468	0	3429	120	0
3	D	947	889	888	8	0
4	F	38	0	34	0	0
4	I	38	0	34	1	0
5	A	14	0	13	0	0
6	A	34	42	0	1	0
7	C	73	65	0	2	0
8	C	90	80	0	0	0
All	All	9181	5187	8824	184	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 10.

All (184) 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:C:881:LEU:HD21	2:C:1100:GLU:HB3	1.56	0.85
2:C:1256:LYS:HE2	2:C:1291:PHE:HZ	1.44	0.82
1:A:539:VAL:HG12	1:A:559:THR:HG22	1.60	0.82
2:C:1247:ASP:HB3	2:C:1250:THR:HG22	1.67	0.77
2:C:1271:HIS:HB2	2:C:1285:ILE:HD13	1.65	0.77
2:C:1035:THR:HG21	2:C:1101:ILE:HD12	1.70	0.71
2:C:981:LYS:HA	2:C:1067:ASP:HB2	1.70	0.71
2:C:913:GLN:HB2	2:C:924:HIS:HB3	1.71	0.71
2:C:1029:ASN:ND2	2:C:1309:ASP:OD2	2.26	0.69
2:C:1213:LYS:HB2	2:C:1235:PRO:HG2	1.74	0.69
2:C:1254:GLN:NE2	2:C:1302:ASP:OD1	2.25	0.66
1:A:560:ASP:O	1:A:565:LYS:NZ	2.27	0.66
2:C:886:ILE:HD11	2:C:1093:CYS:HB3	1.78	0.65
1:A:168:LEU:HD13	1:A:354:GLN:OE1	1.97	0.65
2:C:1079:ASN:HB3	2:C:1082:GLN:HG2	1.79	0.65
2:C:1079:ASN:HB3	2:C:1082:GLN:CG	2.27	0.64
2:C:878:ASP:HB2	2:C:1103:LYS:HE3	1.77	0.64
2:C:1246:ARG:NH1	2:C:1302:ASP:OD2	2.31	0.64
1:A:562:PHE:H	1:A:565:LYS:HG3	1.62	0.63
2:C:1030:LYS:NZ	2:C:1117:PHE:O	2.31	0.63
1:A:242:ASN:HD22	1:A:269:ILE:HD13	1.65	0.61
2:C:1145:ILE:O	2:C:1146:THR:OG1	2.14	0.61
1:A:344:SER:O	1:A:348:ILE:HD12	2.00	0.61
2:C:1112:THR:HG23	2:C:1113:PHE:HD1	1.66	0.61
1:A:242:ASN:ND2	1:A:269:ILE:HD13	2.16	0.60
2:C:928:ASN:HD21	2:C:930:SER:HB2	1.66	0.60
2:C:1112:THR:HG23	2:C:1113:PHE:CD1	2.36	0.60
1:A:369:GLU:N	1:A:369:GLU:OE1	2.34	0.60
2:C:1047:TYR:CE1	2:C:1052:LEU:HD13	2.37	0.60
1:A:547:ARG:HA	1:A:567:VAL:O	2.00	0.60
2:C:1178:LYS:HD3	2:C:1199:TYR:CE2	2.36	0.60
2:C:1096:LEU:HB2	2:C:1101:ILE:HD11	1.83	0.59
3:D:54:GLY:O	3:D:74:ARG:NH1	2.34	0.59
2:C:889:LEU:HD13	2:C:1089:PHE:HB3	1.85	0.58
2:C:911:ASP:OD2	2:C:928:ASN:ND2	2.36	0.58
1:A:226:ALA:HB3	1:A:683:THR:HG21	1.84	0.58
2:C:1130:LEU:HD13	2:C:1175:PHE:CD1	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:252:VAL:CG1	1:A:257:THR:HG23	2.34	0.58
2:C:1241:GLU:HG2	2:C:1243:VAL:HG13	1.86	0.58
2:C:1097:ASN:HB2	2:C:1098:PRO:HD2	1.85	0.57
1:A:470:ASP:HB3	1:A:708:VAL:HG21	1.86	0.57
2:C:1047:TYR:HE1	2:C:1052:LEU:HD13	1.69	0.57
1:A:524:VAL:O	1:A:544:THR:HA	2.05	0.57
2:C:945:ASP:HB3	2:C:1066:GLU:OE1	2.04	0.57
3:D:14:VAL:HG11	3:D:88:LEU:CD1	2.35	0.56
1:A:645:GLY:O	1:A:712:LYS:NZ	2.38	0.56
2:C:1220:ASN:C	2:C:1221:LEU:HD12	2.31	0.56
1:A:528:ASP:OD1	1:A:548:ASN:ND2	2.27	0.56
4:I:1:NAG:H61	4:I:2:NAG:C7	2.36	0.56
1:A:479:ASP:OD1	1:A:483:ARG:NH2	2.38	0.55
2:C:1046:LEU:O	2:C:1052:LEU:HD12	2.07	0.55
2:C:1146:THR:HG22	2:C:1228:GLY:O	2.06	0.55
2:C:991:TRP:HB3	2:C:1004:LEU:HD23	1.89	0.55
1:A:621:LYS:O	1:A:737:ARG:NH1	2.40	0.55
1:A:553:ASN:OD1	1:A:573:ASN:ND2	2.40	0.55
2:C:977:ILE:HD13	2:C:1089:PHE:CE1	2.42	0.54
2:C:993:VAL:HG22	2:C:1002:TRP:HE3	1.72	0.54
1:A:263:LEU:O	1:A:267:VAL:HG23	2.08	0.54
2:C:975:SER:HA	2:C:994:SER:CB	2.37	0.54
2:C:1002:TRP:HZ2	2:C:1038:ASN:HD21	1.55	0.54
2:C:1224:ILE:HD13	2:C:1275:ILE:HD11	1.90	0.54
2:C:1145:ILE:HG22	2:C:1146:THR:HG23	1.89	0.53
1:A:470:ASP:CB	1:A:708:VAL:HG21	2.38	0.53
1:A:176:LEU:CD2	1:A:273:ILE:HD13	2.38	0.53
1:A:248:PHE:O	1:A:252:VAL:HG23	2.08	0.53
2:C:927:ASN:ND2	2:C:1079:ASN:O	2.41	0.53
2:C:996:LYS:O	2:C:999:ASN:HB2	2.09	0.53
2:C:1022:LYS:HB2	2:C:1166:TYR:HB3	1.90	0.53
2:C:1133:VAL:HB	2:C:1302:ASP:HB2	1.90	0.53
2:C:1148:TYR:CD2	2:C:1283:ILE:HD12	2.43	0.53
1:A:674:VAL:O	1:A:682:ARG:NH1	2.38	0.53
2:C:1131:ILE:HD12	2:C:1306:VAL:HG21	1.89	0.52
2:C:1218:PHE:HB2	2:C:1222:ASP:HB2	1.90	0.52
1:A:558:ASN:HA	1:A:578:HIS:O	2.09	0.52
3:D:30:THR:HG23	3:D:33:ARG:HD3	1.92	0.52
1:A:163:ARG:HB2	1:A:384:ALA:HB2	1.91	0.52
2:C:1177:ILE:HG23	2:C:1196:ILE:HD12	1.90	0.52
2:C:1067:ASP:OD1	2:C:1068:ASN:N	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1097:ASN:O	2:C:1101:ILE:HG12	2.08	0.52
2:C:886:ILE:CD1	2:C:1093:CYS:HB3	2.40	0.52
2:C:915:VAL:HG13	2:C:924:HIS:HB2	1.92	0.52
2:C:1117:PHE:HB2	2:C:1306:VAL:HG22	1.91	0.52
2:C:1071:THR:HG22	2:C:1073:LYS:HG2	1.92	0.51
2:C:1240:MET:HG2	2:C:1257:LEU:HD13	1.92	0.51
7:C:1401:A1IHM:O28	7:C:1401:A1IHM:O24	2.28	0.51
2:C:1178:LYS:HD3	2:C:1199:TYR:CD2	2.46	0.51
1:A:493:GLU:O	1:A:495:VAL:HG23	2.11	0.51
2:C:1209:VAL:HG22	2:C:1227:VAL:HG22	1.94	0.50
2:C:1029:ASN:HB3	2:C:1118:TRP:CZ3	2.46	0.50
3:D:14:VAL:HG11	3:D:88:LEU:HD12	1.93	0.50
2:C:1148:TYR:CE2	2:C:1283:ILE:HD12	2.47	0.50
2:C:1151:LEU:HD11	2:C:1270:THR:HG21	1.93	0.50
1:A:466:VAL:HG11	1:A:701:ILE:HG12	1.93	0.50
1:A:605:THR:O	1:A:608:VAL:HG22	2.12	0.50
2:C:1114:LEU:HD22	2:C:1253:VAL:HG22	1.93	0.50
2:C:1191:LYS:O	2:C:1194:ASP:HB2	2.11	0.50
2:C:1058:ILE:O	2:C:1058:ILE:HG13	2.11	0.49
2:C:1293:HIS:O	2:C:1299:LEU:HD22	2.11	0.49
2:C:1077:CYS:SG	2:C:1082:GLN:HG3	2.52	0.49
2:C:1104:LEU:HG	2:C:1108:TYR:HE2	1.77	0.49
2:C:1079:ASN:HB3	2:C:1082:GLN:HG3	1.94	0.49
2:C:1088:LYS:HD2	2:C:1108:TYR:HE1	1.78	0.49
1:A:462:TYR:HA	1:A:465:THR:HG22	1.95	0.48
2:C:1097:ASN:OD1	2:C:1100:GLU:HG3	2.13	0.48
2:C:1247:ASP:HB3	2:C:1250:THR:CG2	2.40	0.48
2:C:972:ASN:O	2:C:996:LYS:HG3	2.13	0.48
2:C:977:ILE:HD13	2:C:1089:PHE:HE1	1.76	0.48
3:D:21:ARG:NH2	3:D:84:GLN:OE1	2.46	0.48
1:A:172:LEU:HB3	1:A:276:VAL:HG13	1.94	0.48
2:C:1257:LEU:O	2:C:1265:LEU:N	2.46	0.48
1:A:168:LEU:HA	1:A:171:VAL:HG22	1.94	0.48
2:C:1097:ASN:CG	2:C:1100:GLU:HG3	2.38	0.48
2:C:1272:ASN:HB2	2:C:1282:ASP:OD1	2.12	0.48
1:A:585:LEU:HG	1:A:588:THR:HG21	1.96	0.48
1:A:252:VAL:HG13	1:A:257:THR:HG23	1.96	0.48
2:C:1134:ALA:HA	2:C:1298:ILE:HD11	1.95	0.48
3:D:98:CYS:O	3:D:117:GLY:N	2.46	0.48
1:A:575:THR:HG22	1:A:577:LEU:HG	1.95	0.47
2:C:928:ASN:ND2	2:C:930:SER:HB2	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1096:LEU:HD22	2:C:1100:GLU:OE1	2.15	0.47
2:C:915:VAL:HG11	2:C:924:HIS:CD2	2.49	0.47
2:C:1213:LYS:HD2	2:C:1235:PRO:CG	2.45	0.47
2:C:957:ARG:HB3	2:C:1085:SER:HB2	1.97	0.47
2:C:1142:LEU:HD21	2:C:1144:ASN:C	2.40	0.47
2:C:943:TYR:CD1	2:C:946:MET:HE3	2.50	0.46
2:C:972:ASN:OD1	2:C:973:GLU:N	2.48	0.46
2:C:1111:ILE:HG13	2:C:1248:LEU:HD13	1.98	0.46
1:A:467:TRP:CZ3	1:A:642:LEU:HD22	2.50	0.46
2:C:944:ASN:O	2:C:948:ASN:HB2	2.15	0.46
2:C:1145:ILE:H	2:C:1145:ILE:HD12	1.80	0.46
1:A:227:ASP:OD2	1:A:684:THR:N	2.49	0.46
2:C:949:ASN:CG	2:C:1039:ASP:HA	2.40	0.46
2:C:878:ASP:HB2	2:C:1103:LYS:CE	2.43	0.46
2:C:1224:ILE:HD13	2:C:1275:ILE:CD1	2.46	0.46
1:A:171:VAL:HG21	1:A:354:GLN:HG3	1.98	0.46
2:C:1047:TYR:HD1	2:C:1052:LEU:HA	1.81	0.46
1:A:460:SER:OG	1:A:722:ALA:HB3	2.17	0.45
2:C:1149:MET:HE3	2:C:1284:LEU:HD22	1.98	0.45
2:C:974:TYR:CE2	2:C:995:LEU:HD12	2.51	0.45
2:C:922:ALA:CB	2:C:1087:ASP:HB2	2.45	0.45
2:C:1256:LYS:HD2	2:C:1258:TYR:CZ	2.51	0.45
2:C:1022:LYS:HA	2:C:1022:LYS:HD3	1.62	0.45
6:A:802:6UZ:O3	6:A:802:6UZ:O6	2.35	0.45
1:A:176:LEU:HD22	1:A:273:ILE:HD13	1.98	0.45
2:C:943:TYR:HB3	2:C:946:MET:HE3	1.99	0.45
2:C:1114:LEU:O	2:C:1121:PRO:HA	2.16	0.45
2:C:1168:ARG:HD2	2:C:1314:ASN:HB3	1.98	0.45
2:C:937:LYS:HB2	2:C:942:GLU:HG2	1.99	0.45
2:C:1127:GLU:HA	2:C:1175:PHE:O	2.16	0.45
2:C:1097:ASN:ND2	2:C:1100:GLU:HG3	2.32	0.44
2:C:934:ILE:HD13	2:C:1071:THR:HA	1.99	0.44
1:A:519:LEU:O	1:A:539:VAL:HA	2.17	0.44
1:A:620:ASP:OD1	1:A:624:ARG:NH1	2.50	0.43
2:C:876:ASP:O	2:C:880:ILE:HG12	2.18	0.43
2:C:967:GLU:OE1	2:C:1162:LEU:HD21	2.18	0.43
2:C:1213:LYS:HB3	2:C:1235:PRO:HD2	2.00	0.43
2:C:1104:LEU:O	2:C:1107:SER:HB3	2.19	0.43
1:A:196:ASP:OD2	1:A:331:HIS:ND1	2.52	0.43
2:C:938:ALA:O	2:C:942:GLU:HG3	2.18	0.43
2:C:935:VAL:O	2:C:1069:ASN:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:951:THR:HB	2:C:1092:PHE:HB2	2.00	0.43
1:A:650:ALA:O	1:A:654:LEU:HD13	2.18	0.42
1:A:599:PHE:CZ	1:A:603:LEU:HD11	2.54	0.42
1:A:538:ASP:OD1	1:A:558:ASN:HB3	2.19	0.42
1:A:576:PHE:HB3	1:A:579:ASN:OD1	2.20	0.42
1:A:670:ASP:O	1:A:674:VAL:HG23	2.20	0.42
2:C:1071:THR:CG2	2:C:1073:LYS:HG2	2.50	0.42
1:A:161:HIS:CE1	1:A:165:GLN:HB3	2.55	0.42
1:A:183:VAL:O	1:A:183:VAL:HG22	2.20	0.42
2:C:918:ILE:HG13	2:C:955:TRP:CD2	2.55	0.42
1:A:541:SER:HB2	1:A:544:THR:HB	2.01	0.41
2:C:922:ALA:HB2	2:C:1087:ASP:HB2	2.01	0.41
7:C:1401:A1IHM:O3	7:C:1401:A1IHM:O4	2.39	0.41
2:C:932:GLU:HG2	2:C:1073:LYS:CD	2.51	0.41
2:C:1139:ASP:CG	2:C:1154:ALA:HB2	2.46	0.41
1:A:735:GLU:O	1:A:739:GLN:NE2	2.54	0.41
1:A:568:ASN:H	3:D:58:ASP:HB3	1.86	0.41
2:C:1145:ILE:HD11	2:C:1202:TYR:HE2	1.86	0.41
1:A:539:VAL:CG1	1:A:559:THR:HG22	2.42	0.41
2:C:1213:LYS:HG2	2:C:1214:ASP:OD1	2.21	0.41
2:C:1143:LYS:HD3	2:C:1147:ASP:OD2	2.21	0.40
2:C:933:VAL:HB	2:C:1072:LEU:HB2	2.03	0.40
2:C:1243:VAL:CG2	2:C:1245:LEU:HD13	2.51	0.40
3:D:14:VAL:HG11	3:D:88:LEU:HD13	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	564/742 (76%)	547 (97%)	17 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	423/469 (90%)	403 (95%)	20 (5%)	0	100	100
3	D	120/490 (24%)	115 (96%)	5 (4%)	0	100	100
All	All	1107/1701 (65%)	1065 (96%)	42 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	476/604 (79%)	476 (100%)	0	100	100
2	C	385/421 (91%)	385 (100%)	0	100	100
3	D	97/390 (25%)	97 (100%)	0	100	100
All	All	958/1415 (68%)	958 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	292	HIS
1	A	612	ASN
2	C	928	ASN
2	C	948	ASN
2	C	998	ASN
2	C	1293	HIS

5.3.3 RNA [i](#)

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 ⓘ

6 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	F	1	1,4	14,14,15	0.70	0	17,19,21	1.17	0
4	NAG	F	2	4	14,14,15	0.72	0	17,19,21	0.94	1 (5%)
4	FUC	F	3	4	10,10,11	0.78	0	14,14,16	2.00	4 (28%)
4	NAG	I	1	1,4	14,14,15	0.79	0	17,19,21	0.78	0
4	NAG	I	2	4	14,14,15	0.72	0	17,19,21	0.80	0
4	FUC	I	3	4	10,10,11	0.74	0	14,14,16	0.97	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	F	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	F	2	4	-	1/6/23/26	0/1/1/1
4	FUC	F	3	4	-	-	0/1/1/1
4	NAG	I	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	I	2	4	-	0/6/23/26	0/1/1/1
4	FUC	I	3	4	-	-	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	3	FUC	C1-C2-C3	4.40	116.05	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	3	FUC	C2-C3-C4	3.22	116.52	110.86
4	F	3	FUC	O5-C5-C4	3.16	115.23	109.55
4	F	2	NAG	C1-C2-N2	-2.37	106.70	110.43
4	F	3	FUC	C3-C4-C5	2.13	113.05	109.81

There are no chirality outliers.

All (5) torsion outliers are listed below:

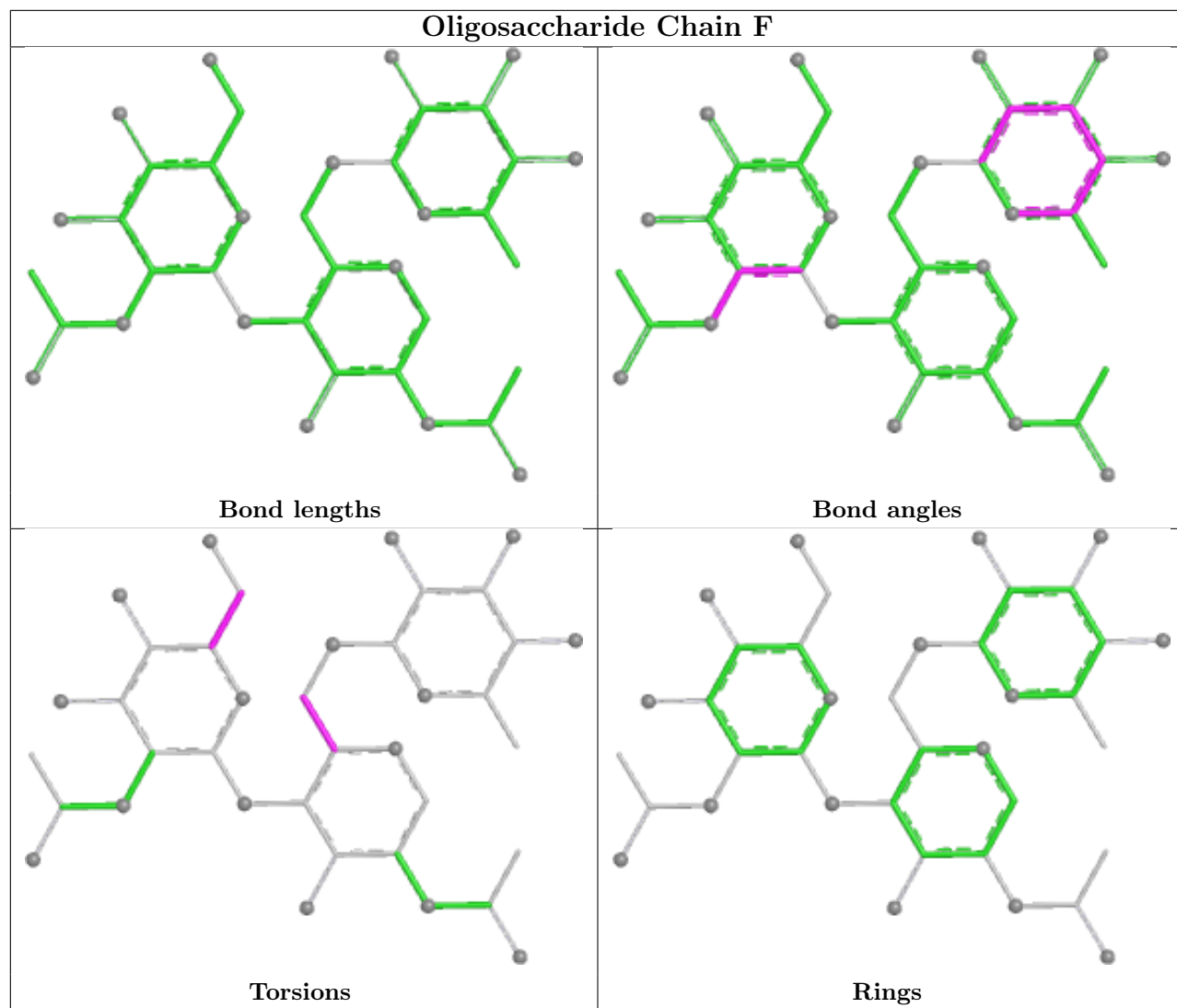
Mol	Chain	Res	Type	Atoms
4	I	1	NAG	C4-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6
4	F	1	NAG	C4-C5-C6-O6
4	F	2	NAG	O5-C5-C6-O6
4	F	1	NAG	O5-C5-C6-O6

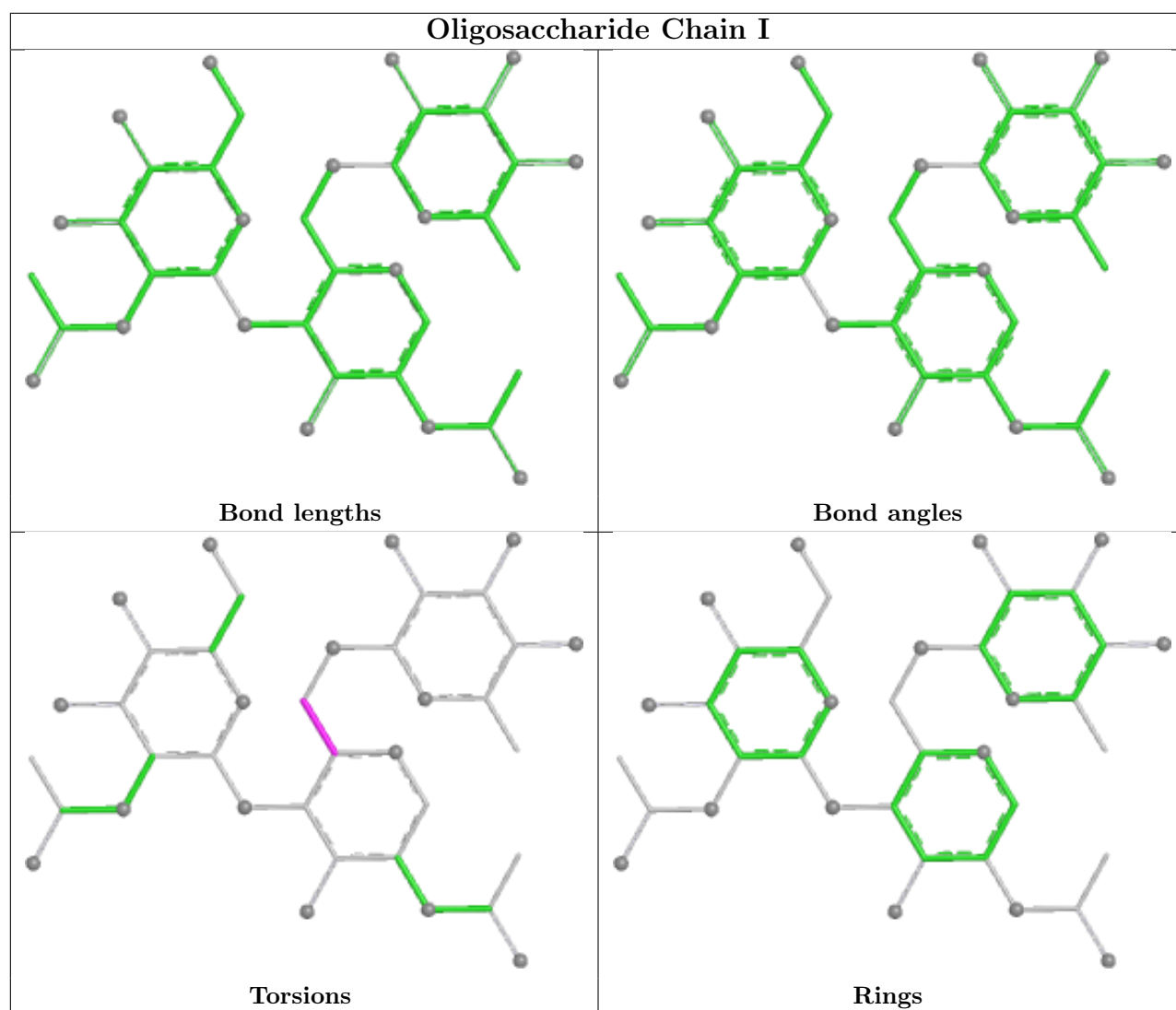
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	I	2	NAG	1	0
4	I	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.





5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
8	A1IHN	C	1402	-	93,95,136	1.39	13 (13%)	128,141,185	1.56	17 (13%)
7	A1IHM	C	1401	-	76,77,114	1.42	12 (15%)	108,115,155	1.36	14 (12%)
5	NAG	A	801	1	14,14,15	0.72	0	17,19,21	0.78	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	6UZ	A	802	-	35,35,35	1.75	11 (31%)	46,46,46	1.00	2 (4%)

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	A1IHN	C	1402	-	-	27/75/193/239	0/6/6/6
7	A1IHM	C	1401	-	-	27/61/158/200	1/5/5/5
5	NAG	A	801	1	-	0/6/23/26	0/1/1/1
6	6UZ	A	802	-	-	7/20/60/60	0/2/2/2

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	C	1402	A1IHN	O27-C41	-4.19	1.36	1.42
6	A	802	6UZ	O5-C5	4.13	1.54	1.44
6	A	802	6UZ	O5'-C5'	4.07	1.54	1.44
7	C	1401	A1IHM	C21-N1	3.87	1.46	1.34
7	C	1401	A1IHM	C30-C29	3.71	1.56	1.51
7	C	1401	A1IHM	C28-C27	3.54	1.57	1.52
8	C	1402	A1IHN	C35-N1	3.51	1.45	1.34
8	C	1402	A1IHN	C49-N2	3.44	1.45	1.34
7	C	1401	A1IHM	O25-C25	-3.21	1.36	1.43
6	A	802	6UZ	C6'-C5'	-2.92	1.42	1.51
7	C	1401	A1IHM	O14-C27	-2.88	1.39	1.44
8	C	1402	A1IHN	O27-C45	-2.82	1.39	1.44
6	A	802	6UZ	C6-C5	-2.70	1.42	1.51
7	C	1401	A1IHM	O14-C23	-2.67	1.38	1.42
6	A	802	6UZ	C3-C2	-2.50	1.45	1.52
7	C	1401	A1IHM	O23-C29	-2.45	1.38	1.43
6	A	802	6UZ	O5-C1	2.42	1.48	1.41
7	C	1401	A1IHM	O15-C31	2.39	1.44	1.40
8	C	1402	A1IHN	C43-C44	2.38	1.55	1.53
8	C	1402	A1IHN	O8-C19	-2.37	1.38	1.44
8	C	1402	A1IHN	O12-C25	-2.36	1.38	1.44
7	C	1401	A1IHM	O5-C13	-2.36	1.38	1.44
8	C	1402	A1IHN	O35-C10	-2.35	1.37	1.43
7	C	1401	A1IHM	C24-C25	2.33	1.56	1.53
6	A	802	6UZ	C3'-C2'	-2.30	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	802	6UZ	O3-C3	2.30	1.48	1.43
7	C	1401	A1IHM	O28-C10	-2.20	1.37	1.43
6	A	802	6UZ	O5'-C1'	2.16	1.47	1.41
8	C	1402	A1IHN	O21-C29	-2.12	1.38	1.43
8	C	1402	A1IHN	C31-C30	2.08	1.56	1.53
8	C	1402	A1IHN	O12-C21	-2.06	1.36	1.41
6	A	802	6UZ	O3'-C3'	2.05	1.48	1.43
6	A	802	6UZ	O2'-C2'	2.04	1.48	1.43
8	C	1402	A1IHN	O31-C49	-2.02	1.18	1.23
8	C	1402	A1IHN	C32-C31	2.02	1.55	1.52
7	C	1401	A1IHM	O5-C9	-2.01	1.36	1.41

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	1402	A1IHN	C27-O16-C31	8.32	133.91	114.36
7	C	1401	A1IHM	C23-O14-C27	5.45	127.16	114.36
7	C	1401	A1IHM	C3-O2-C7	4.98	123.44	113.72
8	C	1402	A1IHN	C15-O8-C19	-4.65	104.65	113.72
8	C	1402	A1IHN	C21-O12-C25	4.56	122.62	113.72
8	C	1402	A1IHN	C3-O2-C7	4.54	122.59	113.72
7	C	1401	A1IHM	C15-O8-C19	3.62	120.80	113.72
8	C	1402	A1IHN	C41-O27-C45	3.33	122.18	114.36
8	C	1402	A1IHN	O11-C21-C22	3.23	116.03	108.09
8	C	1402	A1IHN	C9-C10-C11	3.14	115.58	110.11
6	A	802	6UZ	O1-C4'-C3'	2.96	114.76	107.23
7	C	1401	A1IHM	O8-C15-C16	-2.73	105.44	110.59
7	C	1401	A1IHM	C23-C24-C25	2.70	116.19	110.73
8	C	1402	A1IHN	O16-C31-C30	2.57	112.20	109.84
7	C	1401	A1IHM	O4-C6-C5	2.51	113.62	107.23
7	C	1401	A1IHM	O16-C31-C32	-2.49	108.00	111.35
8	C	1402	A1IHN	O26-C11-C10	-2.48	104.82	109.82
8	C	1402	A1IHN	C9-O5-C13	2.47	118.55	113.72
7	C	1401	A1IHM	C2-O1-C3	2.45	119.27	113.91
8	C	1402	A1IHN	C50-C49-N2	2.37	120.05	116.12
8	C	1402	A1IHN	C31-C30-N1	2.34	114.64	110.91
6	A	802	6UZ	C3-C4-C5	2.33	114.45	110.23
8	C	1402	A1IHN	C36-C35-N1	2.32	119.97	116.12
8	C	1402	A1IHN	O26-C41-C42	2.30	112.97	107.23
7	C	1401	A1IHM	C32-C31-C39	-2.26	106.29	111.39
8	C	1402	A1IHN	O27-C45-C46	2.21	110.11	106.65
7	C	1401	A1IHM	O7-C12-C11	2.17	112.93	107.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	1401	A1IHM	O14-C23-C24	-2.14	108.47	111.35
8	C	1402	A1IHN	O8-C15-C16	-2.12	109.57	112.23
7	C	1401	A1IHM	C15-C16-N1	-2.08	107.42	110.92
8	C	1402	A1IHN	C11-C12-C13	-2.08	106.32	110.58
7	C	1401	A1IHM	O7-C12-C13	2.07	114.90	109.48
7	C	1401	A1IHM	C22-C21-N1	2.03	119.48	116.12

There are no chirality outliers.

All (61) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	C	1401	A1IHM	C26-C27-C28-O24
7	C	1401	A1IHM	O14-C27-C28-O24
7	C	1401	A1IHM	C24-C23-O13-C11
7	C	1401	A1IHM	O14-C23-O13-C11
7	C	1401	A1IHM	O16-C35-C36-C37
7	C	1401	A1IHM	C35-C36-C37-O18
7	C	1401	A1IHM	O19-C36-C37-C38
7	C	1401	A1IHM	O19-C36-C37-O18
8	C	1402	A1IHN	O16-C27-C37-O23
8	C	1402	A1IHN	O26-C41-C51-O33
8	C	1402	A1IHN	C44-C45-C46-C47
8	C	1402	A1IHN	C44-C45-C46-O30
8	C	1402	A1IHN	O27-C45-C46-C47
8	C	1402	A1IHN	O27-C45-C46-O30
8	C	1402	A1IHN	C22-C21-O11-C17
6	A	802	6UZ	C3'-C4'-O1-C1
7	C	1401	A1IHM	C13-C12-O7-C15
7	C	1401	A1IHM	O5-C9-O4-C6
6	A	802	6UZ	C4-C5-C6-O6
8	C	1402	A1IHN	O8-C19-C20-O9
7	C	1401	A1IHM	C35-C36-C37-C38
7	C	1401	A1IHM	O8-C19-C20-O9
8	C	1402	A1IHN	C12-C13-C14-O6
8	C	1402	A1IHN	O5-C13-C14-O6
6	A	802	6UZ	O5-C5-C6-O6
7	C	1401	A1IHM	C12-C13-C14-O6
8	C	1402	A1IHN	O18-C33-C34-O17
8	C	1402	A1IHN	C18-C19-C20-O9
6	A	802	6UZ	O1'-CA-CB-CC
7	C	1401	A1IHM	C4-C3-O1-C2
7	C	1401	A1IHM	O5-C13-C14-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
7	C	1401	A1IHM	O2-C3-O1-C2
7	C	1401	A1IHM	C29-C30-O15-C31
8	C	1402	A1IHN	O12-C25-C26-O13
8	C	1402	A1IHN	O19-C32-C33-C34
8	C	1402	A1IHN	O19-C32-C33-O18
8	C	1402	A1IHN	C31-C32-C33-C34
8	C	1402	A1IHN	C31-C32-C33-O18
7	C	1401	A1IHM	O2-C7-C8-O3
6	A	802	6UZ	CH-CI-CJ-CK
8	C	1402	A1IHN	C32-C33-C34-O17
7	C	1401	A1IHM	C7-C6-O4-C9
7	C	1401	A1IHM	C5-C6-O4-C9
8	C	1402	A1IHN	C28-C27-C37-O22
8	C	1402	A1IHN	C24-C25-C26-O13
8	C	1402	A1IHN	C1-C2-O1-C3
7	C	1401	A1IHM	C1-C2-O1-C3
8	C	1402	A1IHN	C16-C38-C39-C40
7	C	1401	A1IHM	C18-C19-C20-O9
7	C	1401	A1IHM	O14-C27-C28-C29
7	C	1401	A1IHM	C26-C27-C28-C29
7	C	1401	A1IHM	C34-C35-C36-C37
8	C	1402	A1IHN	C30-C31-C32-C33
8	C	1402	A1IHN	C16-C38-C39-O25
8	C	1402	A1IHN	O12-C21-O11-C17
7	C	1401	A1IHM	C24-C23-C40-O26
6	A	802	6UZ	CC-CD-CF-CG
7	C	1401	A1IHM	O18-C37-C38-O17
8	C	1402	A1IHN	C30-C31-C32-O19
6	A	802	6UZ	CI-CJ-CK-CL
8	C	1402	A1IHN	C29-C30-N1-C35

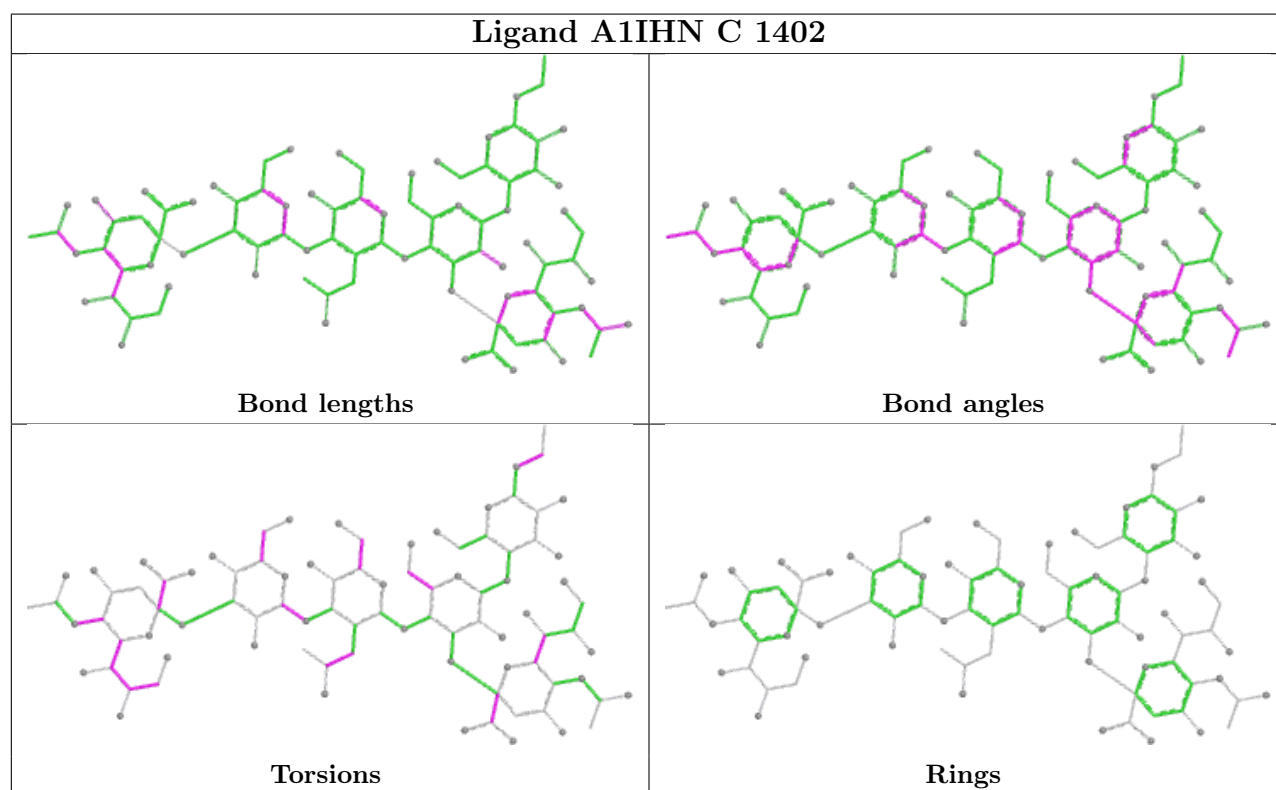
All (1) ring outliers are listed below:

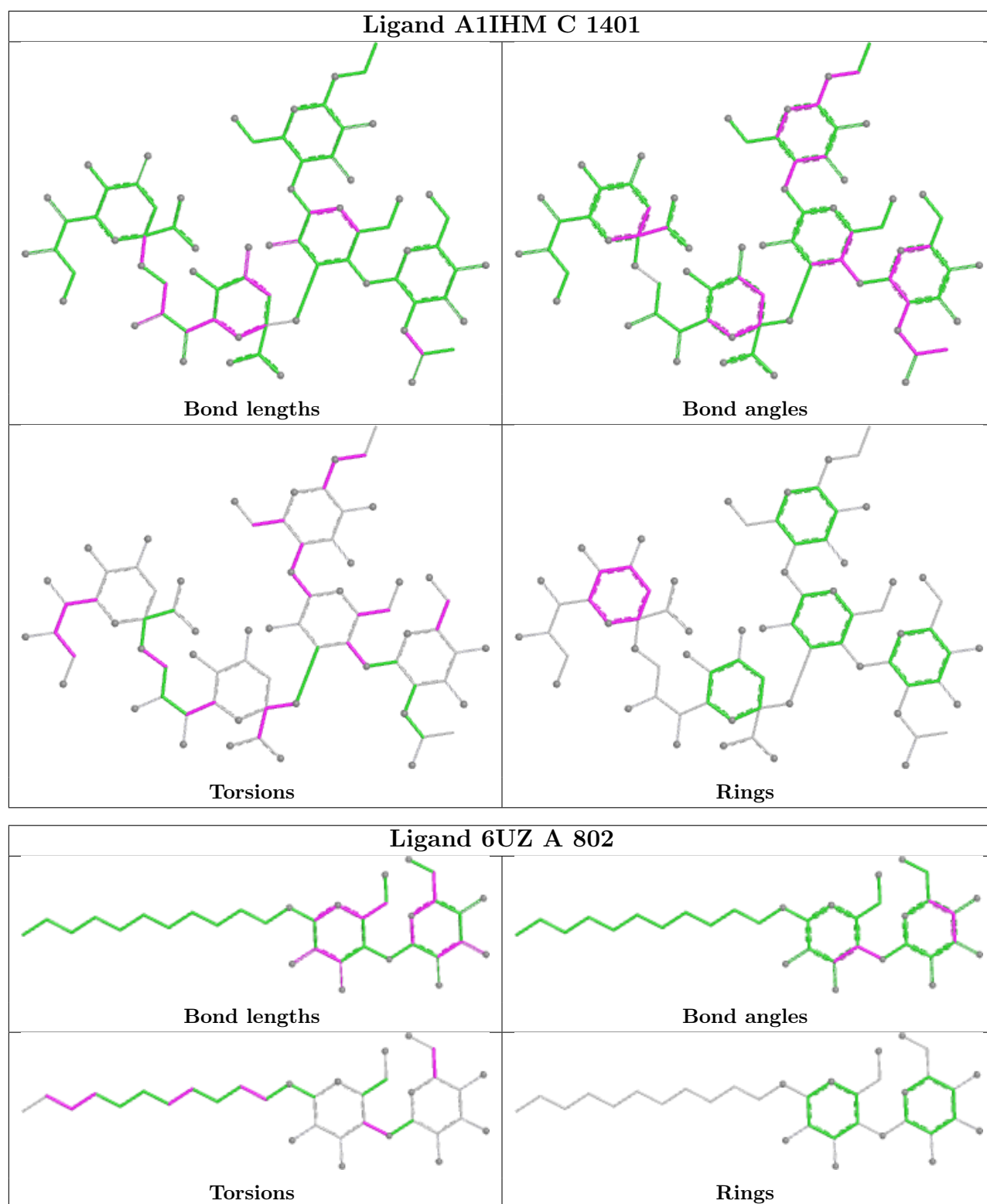
Mol	Chain	Res	Type	Atoms
7	C	1401	A1IHM	C31-C32-C33-C34-C35-O16

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	1401	A1IHM	2	0
6	A	802	6UZ	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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.

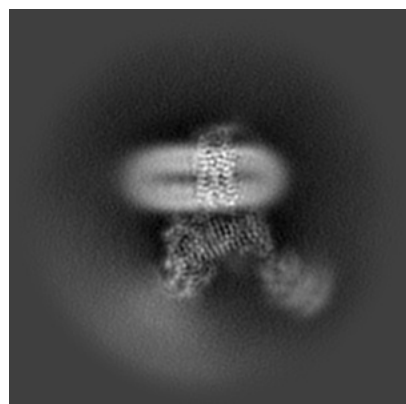
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-50889. These allow visual inspection of the internal detail of the map and identification of artifacts.

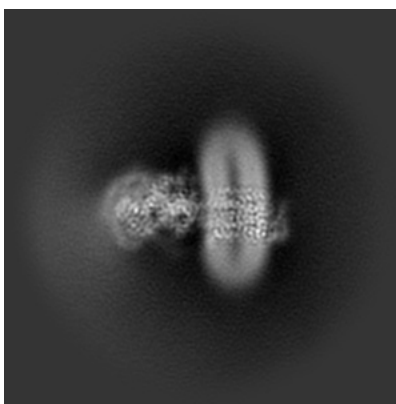
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

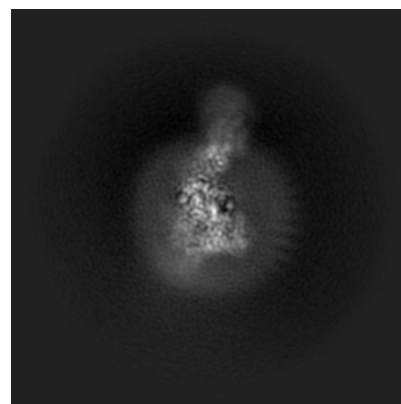
6.1.1 Primary map



X

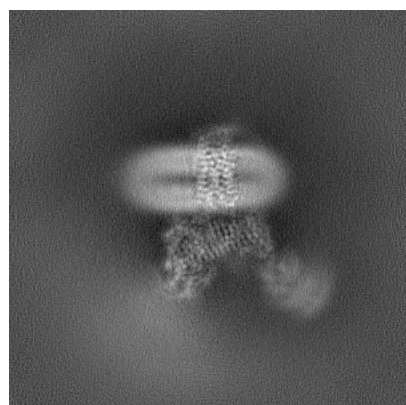


Y

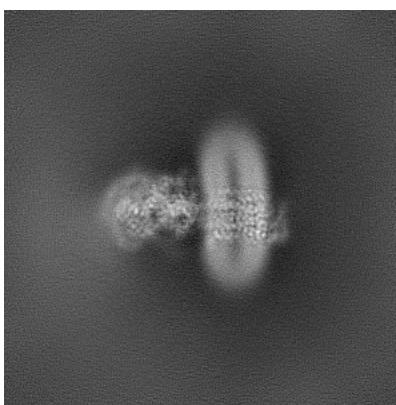


Z

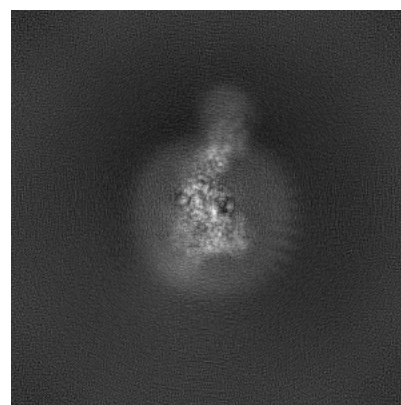
6.1.2 Raw map



X



Y

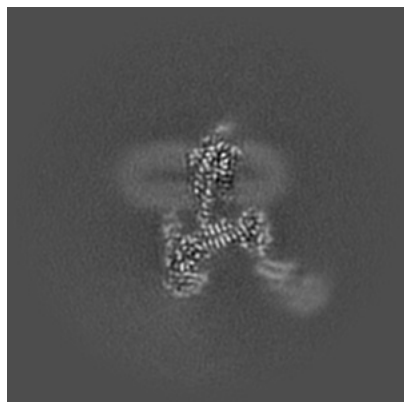


Z

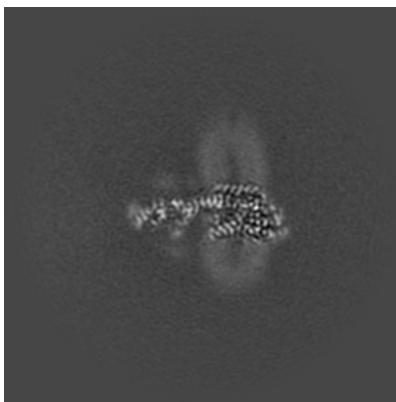
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

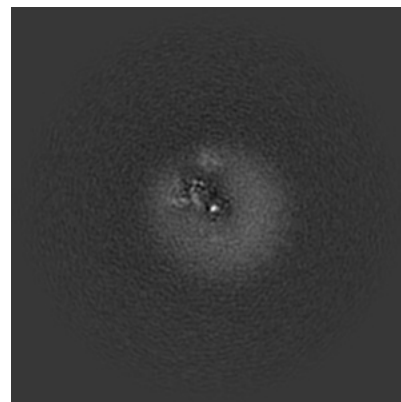
6.2.1 Primary map



X Index: 224

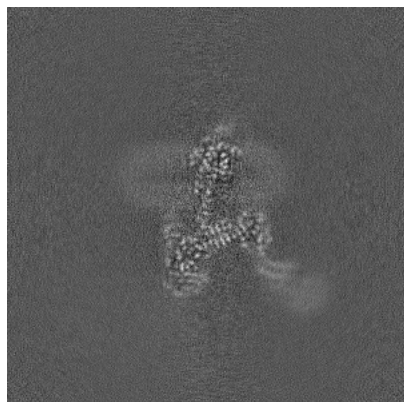


Y Index: 224

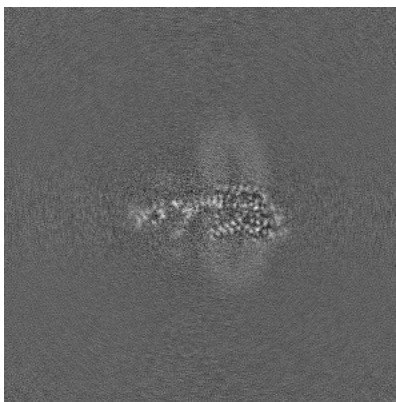


Z Index: 224

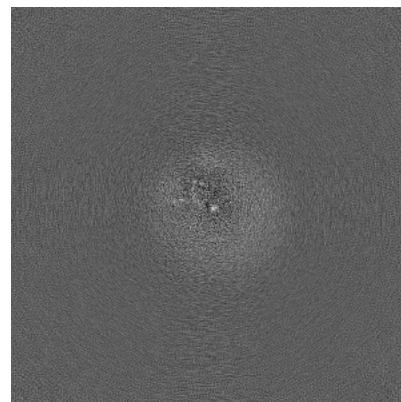
6.2.2 Raw map



X Index: 224



Y Index: 224

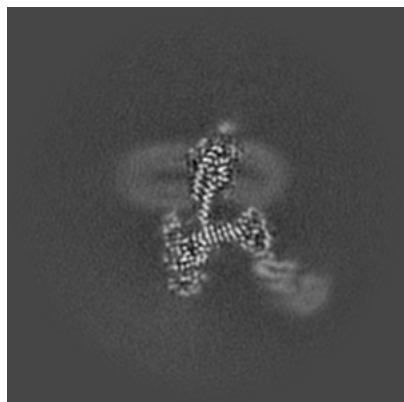


Z Index: 224

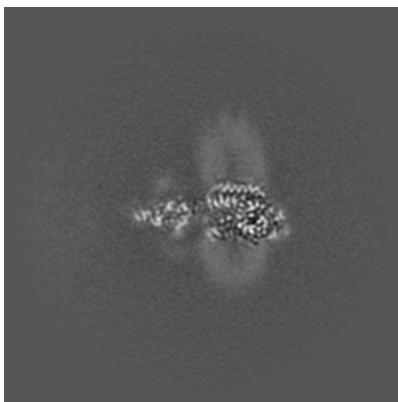
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

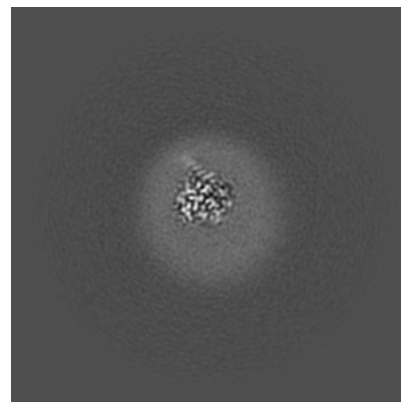
6.3.1 Primary map



X Index: 227



Y Index: 227

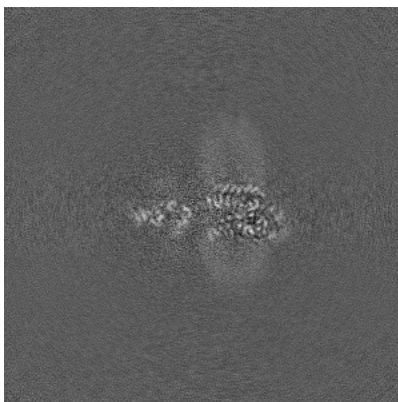


Z Index: 284

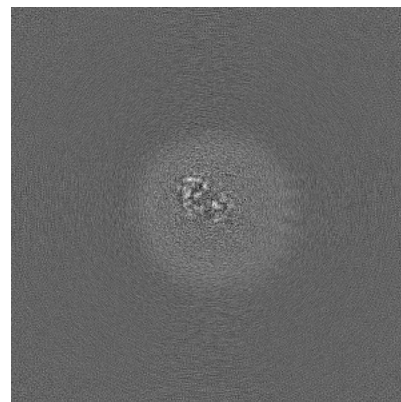
6.3.2 Raw map



X Index: 227



Y Index: 226

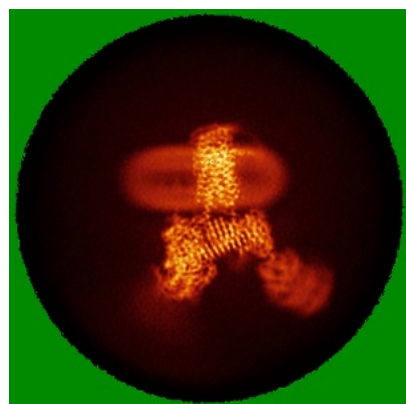


Z Index: 237

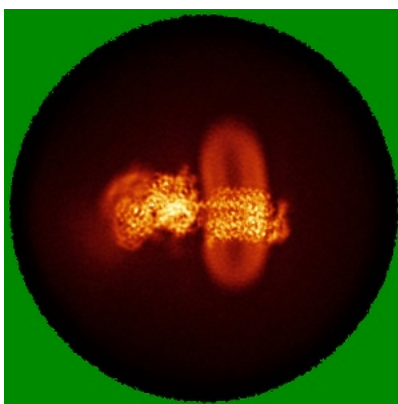
The images above show the largest variance slices of the map in three orthogonal directions.

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

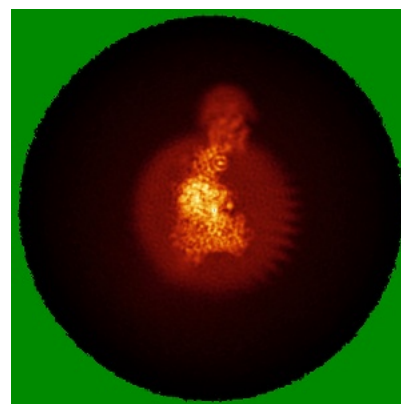
6.4.1 Primary map



X

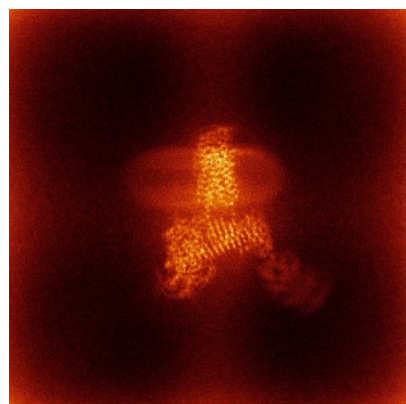


Y

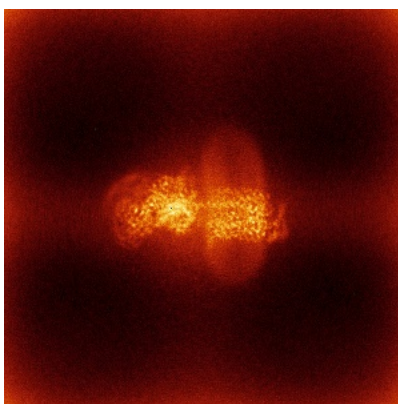


Z

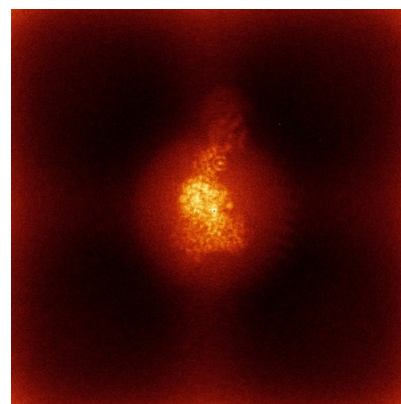
6.4.2 Raw map



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.

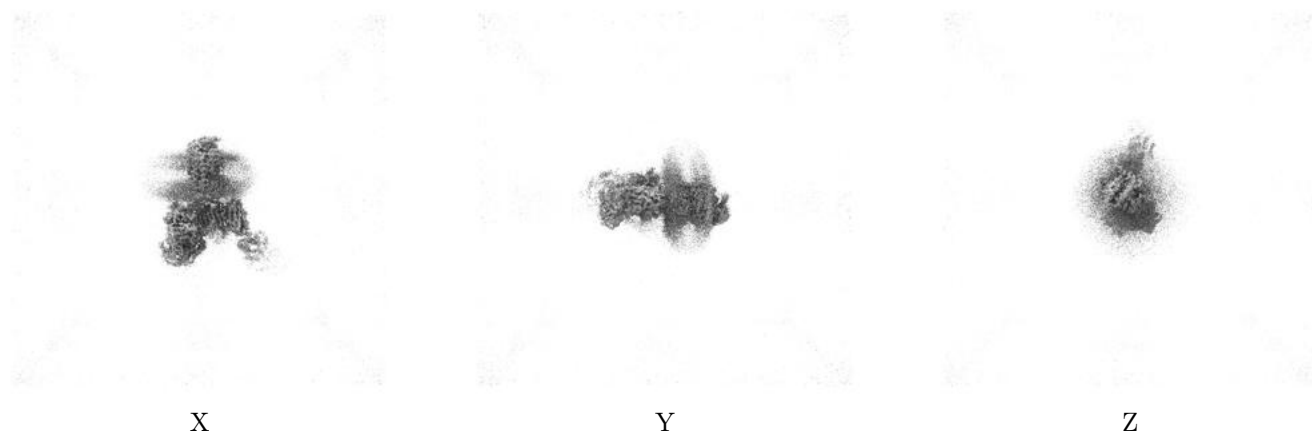
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0657. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

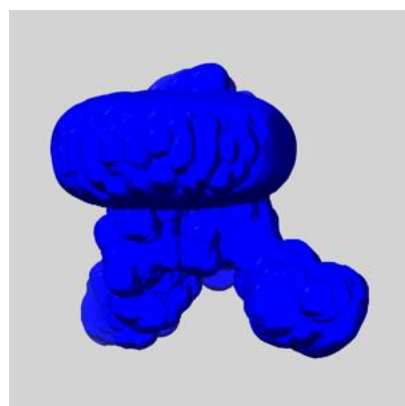
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

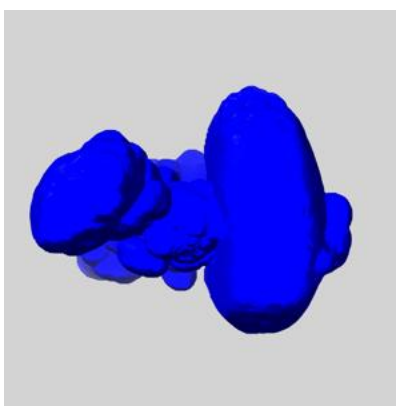
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

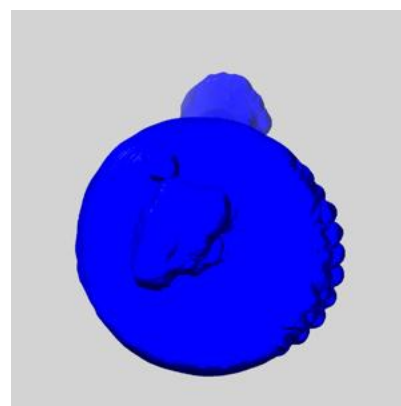
6.6.1 emd_50889_msk_1.map [i](#)



X



Y

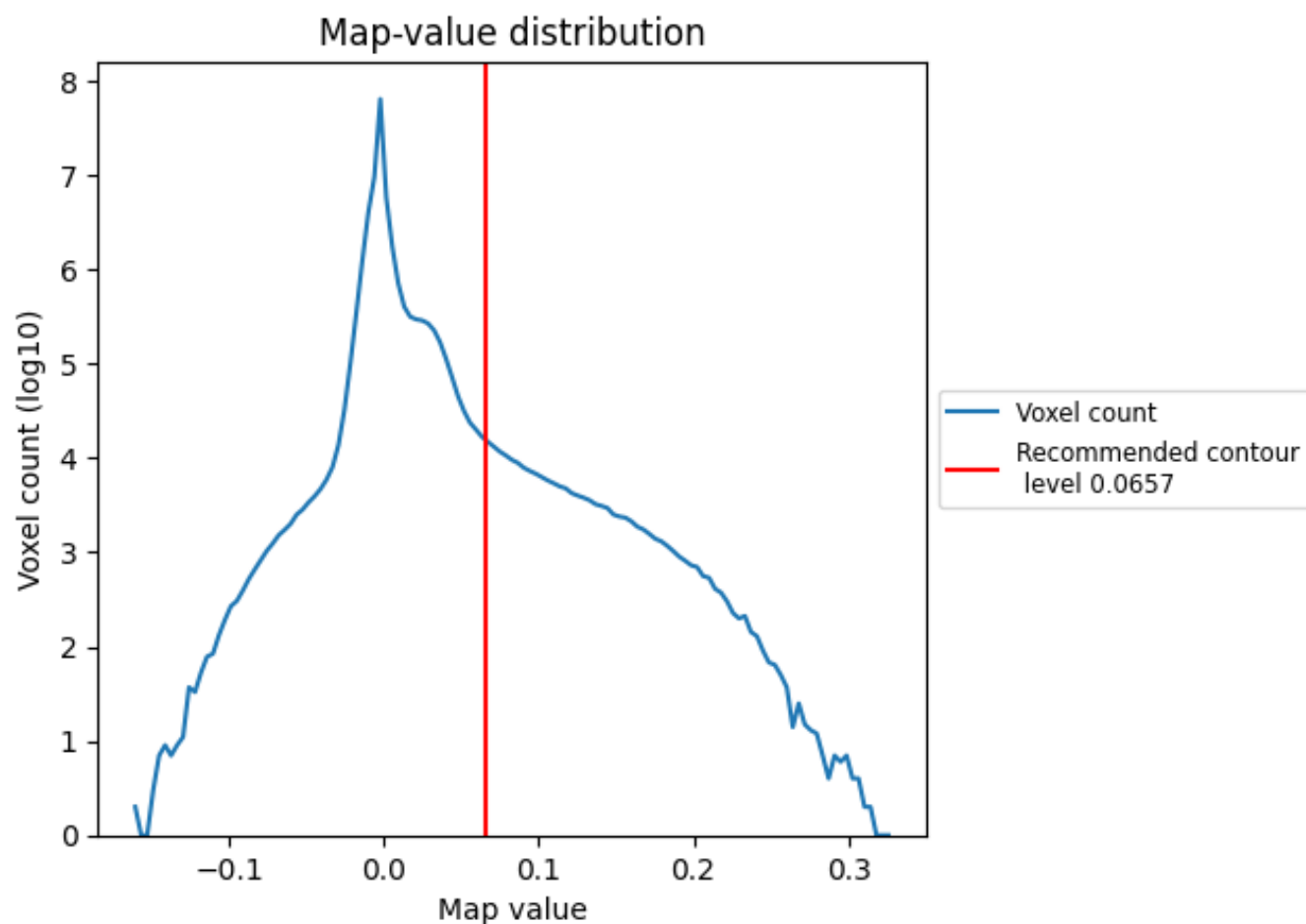


Z

7 Map analysis [i](#)

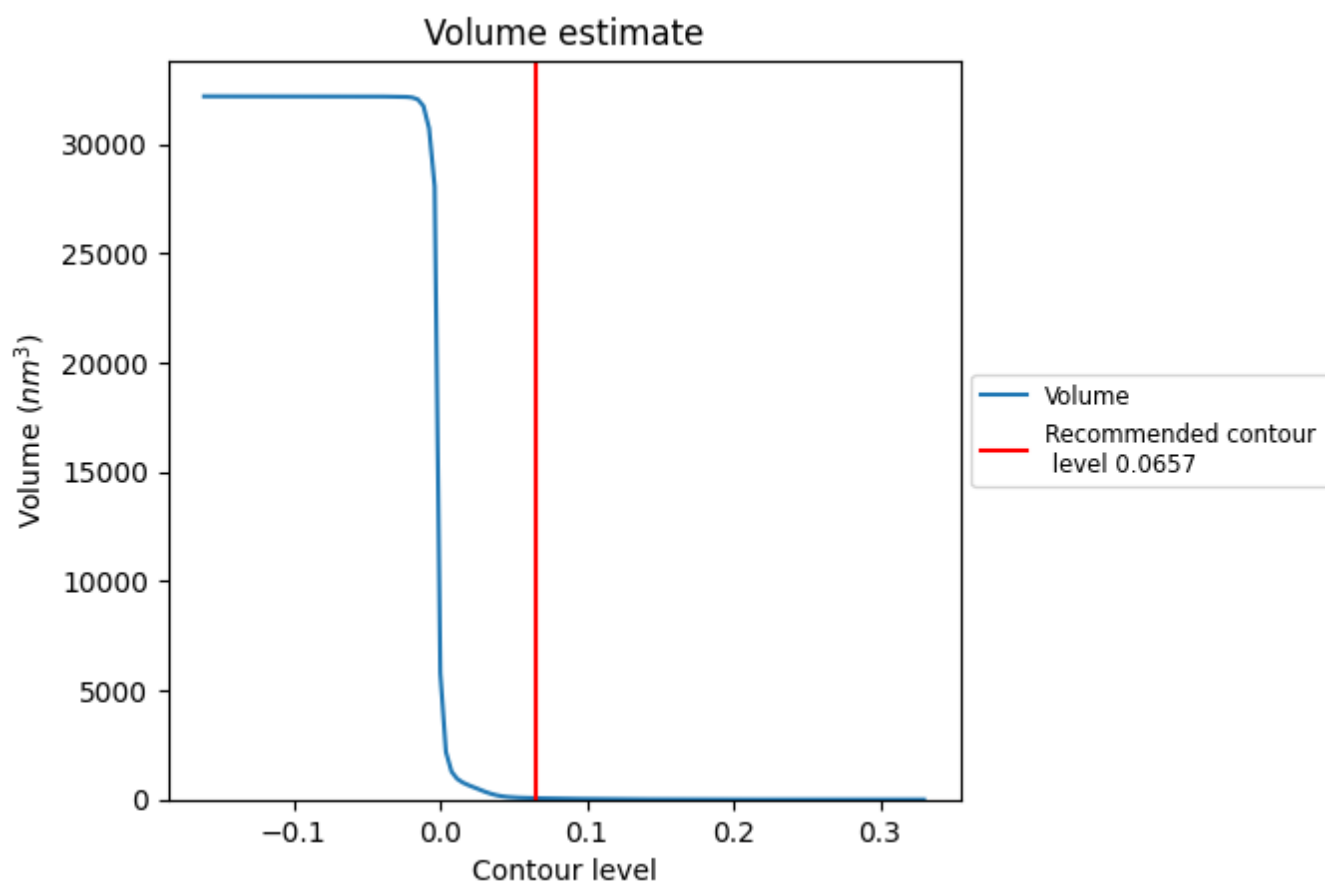
This section contains the results of statistical analysis of the map.

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.

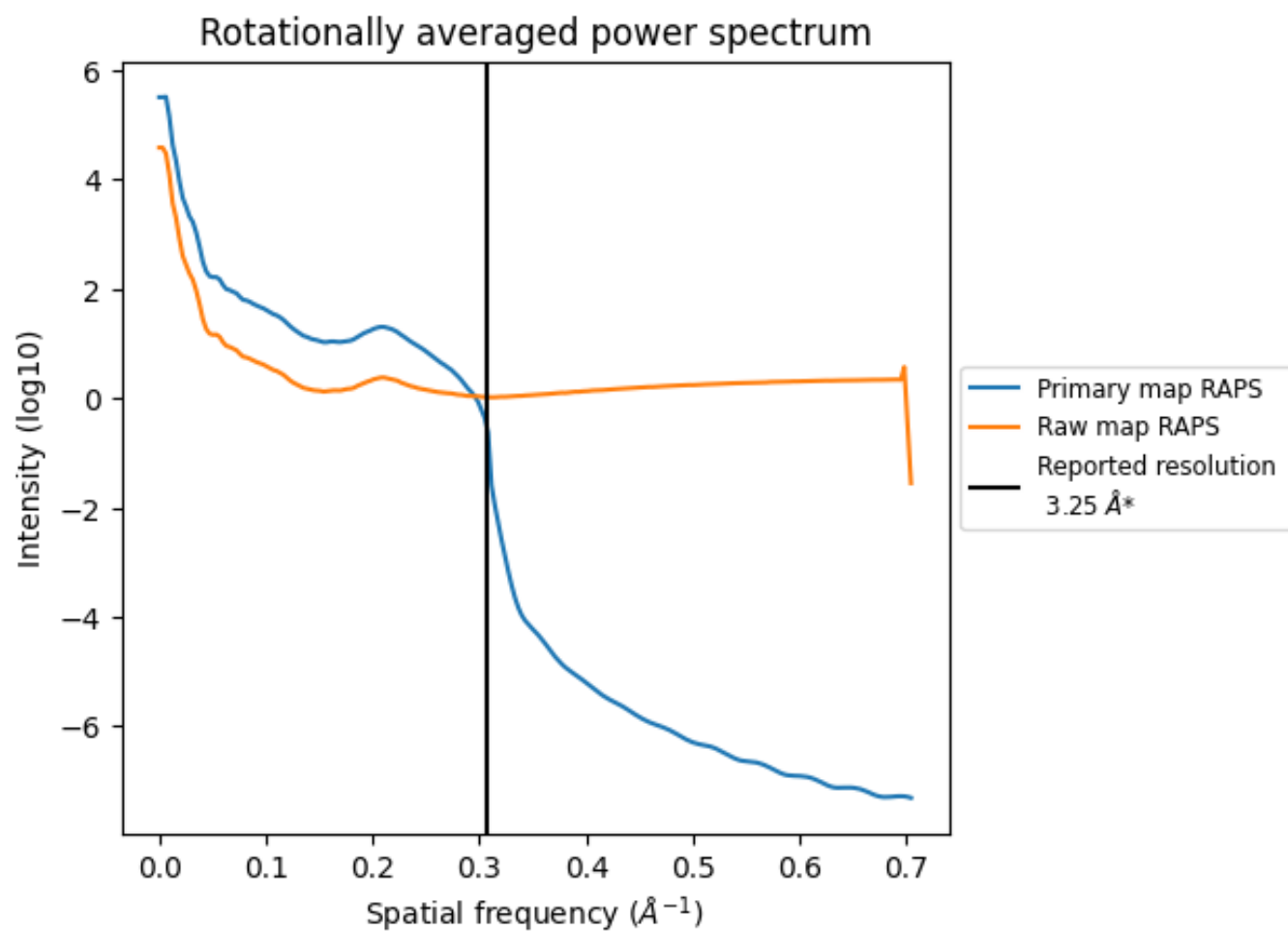
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 63 nm³; this corresponds to an approximate mass of 57 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 ⓘ

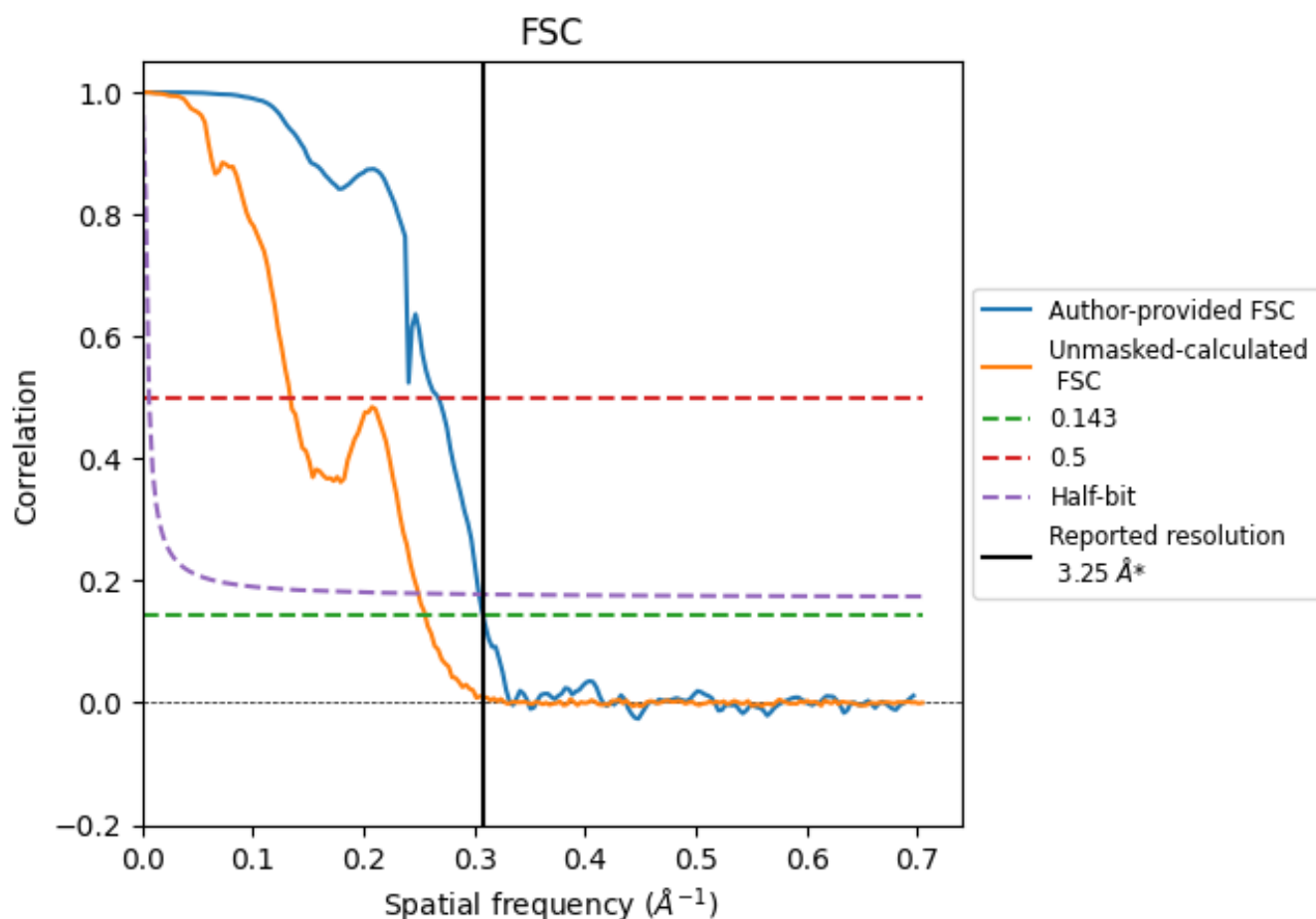


*Reported resolution corresponds to spatial frequency of 0.308 $\text{\AA}^{-1}$

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 0.308 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

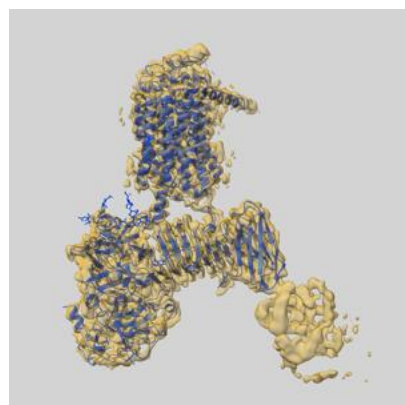
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.25	-	-
Author-provided FSC curve	3.25	3.75	3.29
Unmasked-calculated*	3.91	7.48	4.01

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.91 differs from the reported value 3.25 by more than 10 %

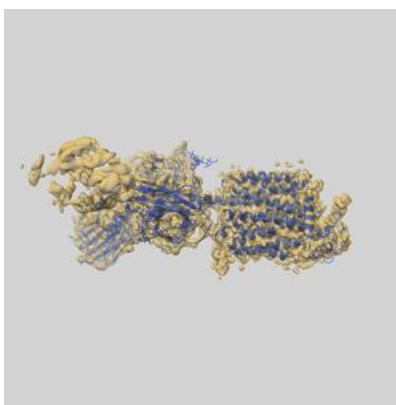
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-50889 and PDB model 9FYQ. Per-residue inclusion information can be found in section 3 on page 9.

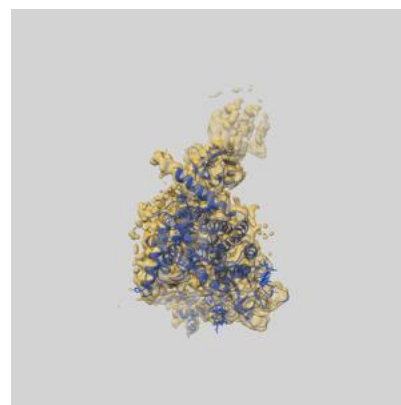
9.1 Map-model overlay [i](#)



X



Y



Z

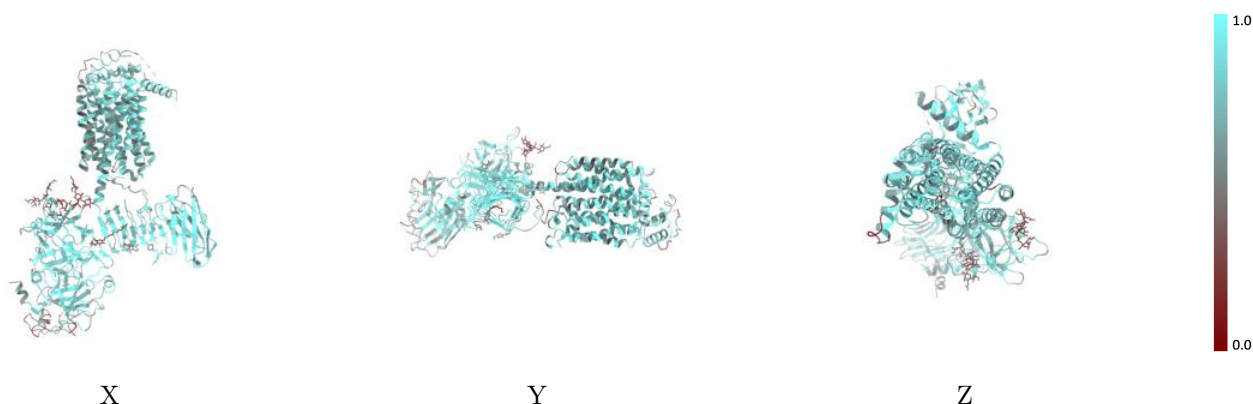
The images above show the 3D surface view of the map at the recommended contour level 0.0657 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



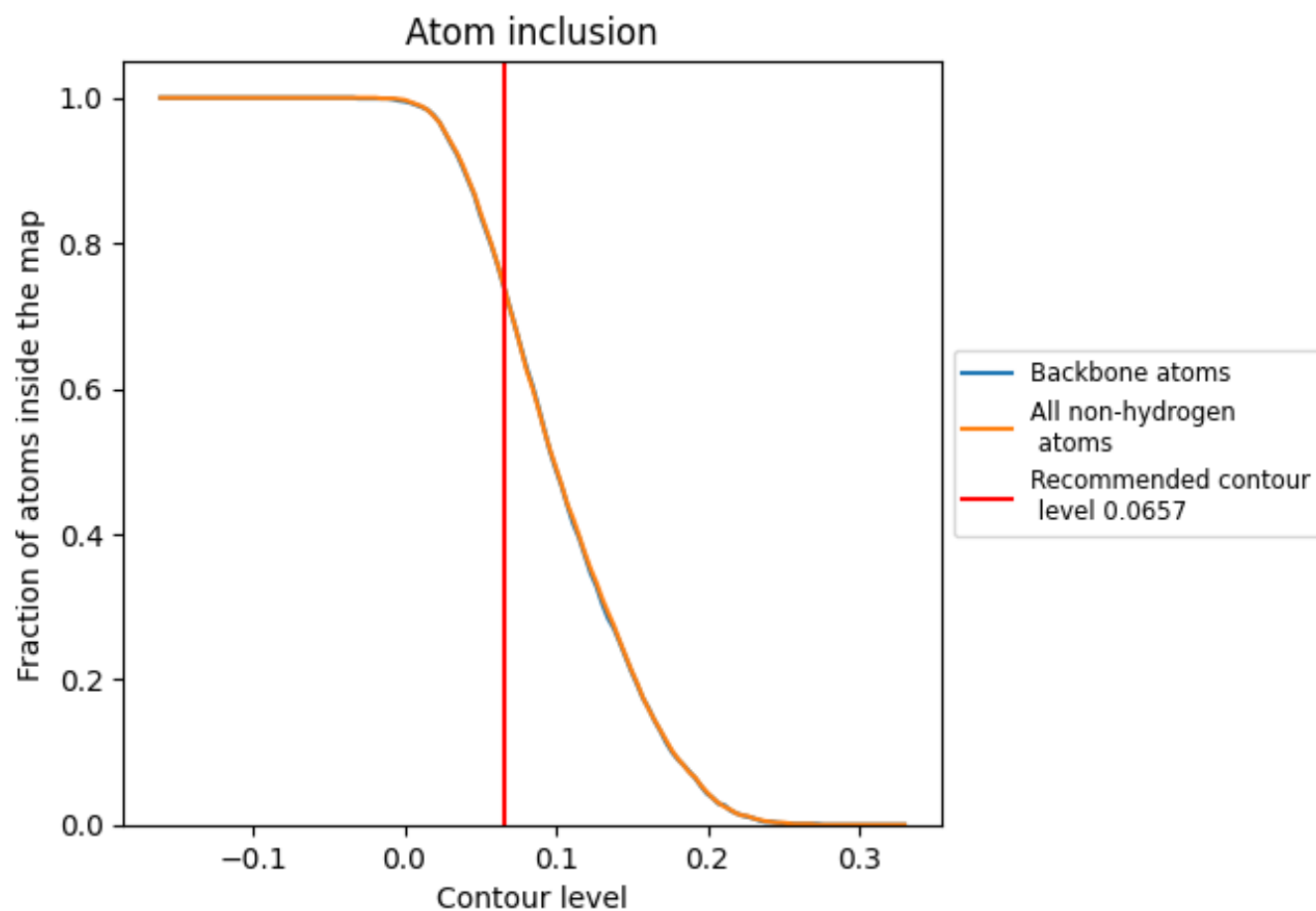
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0657).

9.4 Atom inclusion [i](#)



At the recommended contour level, 74% of all backbone atoms, 74% 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 (0.0657) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7410	0.4680
A	0.7570	0.4770
C	0.7190	0.4540
D	0.8130	0.4760
F	0.7630	0.4660
I	0.5260	0.4490

