



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 04:48 PM UTC

PDB ID : 6OEU / pdb_00006oeu
EMDB ID : EMD-7795
Title : Structure of human Patched1
Authors : Qi, X.; Li, X.; Wang, J.
Deposited on : 2019-03-27
Resolution : 3.50 Å(reported)

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

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with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

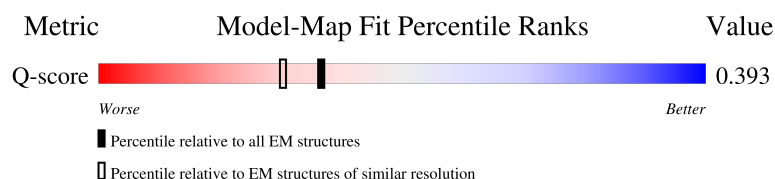
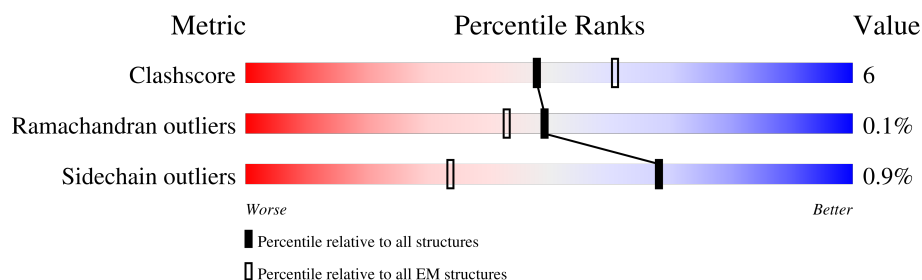
EMDB validation analysis	:	0.0.1.dev132
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13950 (3.00 - 4.00)

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

Mol	Chain	Length	Quality of chain
1	A	1455	30% 57% 9% 34%
2	B	2	100% 50% 50%
2	C	2	100% 50% 50%
2	D	2	100% 50% 50%

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Mol	Chain	Length	Quality of chain
2	E	2	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	D	1	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7310 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	964	7198	4670	1207	1287	34	0	0

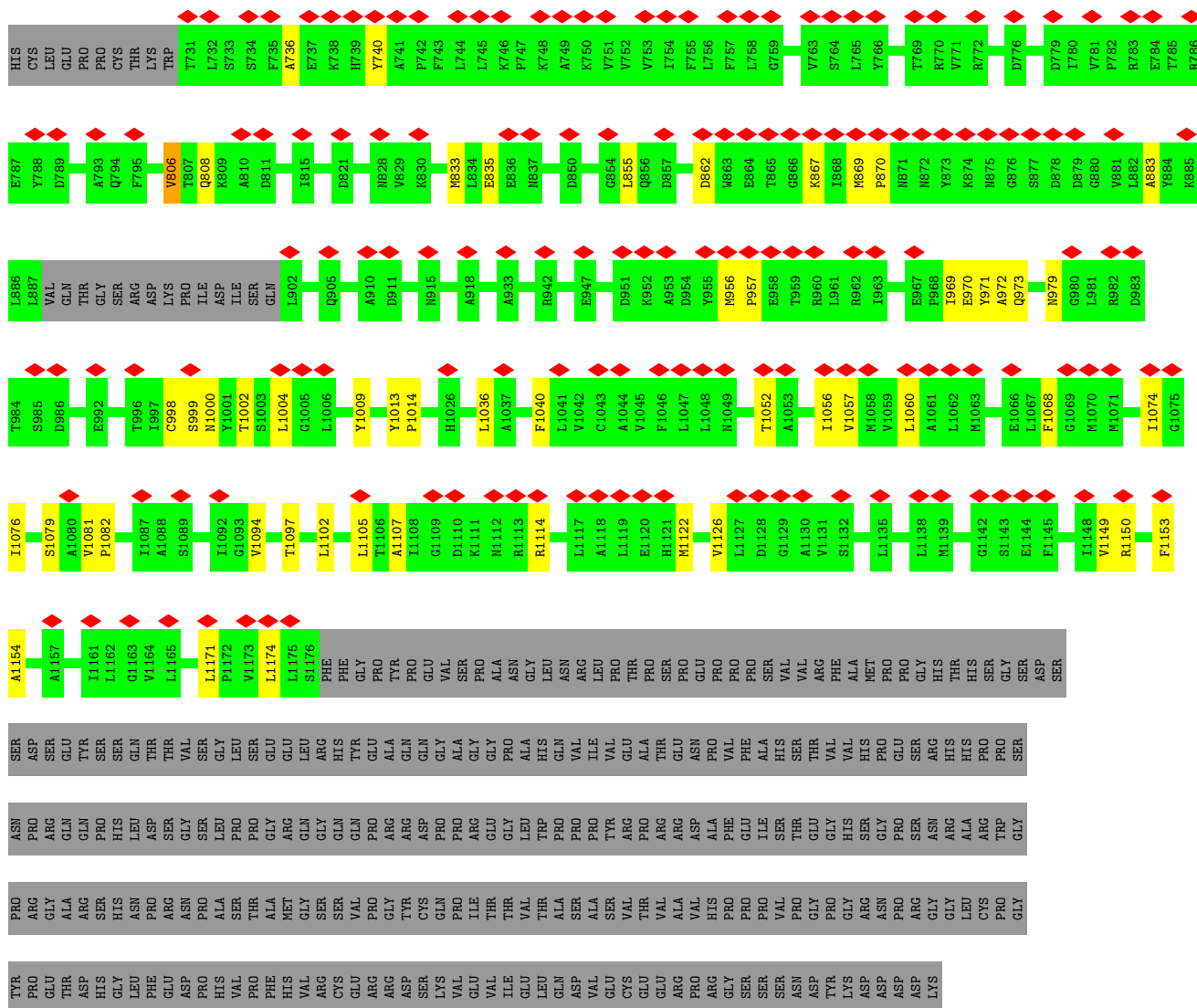
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1448	ASP	-	expression tag	UNP Q13635
A	1449	TYR	-	expression tag	UNP Q13635
A	1450	LYS	-	expression tag	UNP Q13635
A	1451	ASP	-	expression tag	UNP Q13635
A	1452	ASP	-	expression tag	UNP Q13635
A	1453	ASP	-	expression tag	UNP Q13635
A	1454	ASP	-	expression tag	UNP Q13635
A	1455	LYS	-	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
2	C	2	28	16	2	10	0	0
2	D	2	28	16	2	10	0	0
2	E	2	28	16	2	10	0	0



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



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



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



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



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	789118	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1.6	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	3.830	Depositor
Minimum map value	-2.511	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.103	Depositor
Recommended contour level	0.8	Depositor
Map size ($\text{\AA}$)	256.0, 256.0, 256.0	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.0, 1.0, 1.0	Depositor

5 Model quality

5.1 Standard geometry

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	1.10	1/7354 (0.0%)	1.56	18/10020 (0.2%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	342	ILE	N-CA	5.01	1.50	1.46

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	415	SER	N-CA-C	-13.17	91.14	110.23
1	A	121	LEU	N-CA-C	-8.86	93.00	108.69
1	A	416	THR	N-CA-C	-8.12	101.16	112.45
1	A	415	SER	CB-CA-C	7.68	122.46	109.72
1	A	252	PRO	N-CA-CB	7.45	110.31	103.08
1	A	261	PRO	N-CA-CB	7.27	110.88	103.25
1	A	231	PRO	N-CA-CB	6.88	109.88	103.41
1	A	253	PRO	N-CA-CB	6.68	110.26	103.25
1	A	870	PRO	N-CA-CB	6.53	110.45	103.52
1	A	957	PRO	N-CA-CB	6.05	110.08	103.30
1	A	956	MET	CA-C-N	6.02	126.45	119.47
1	A	956	MET	C-N-CA	6.02	126.45	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	PRO	N-CA-CB	5.77	109.67	103.33
1	A	869	MET	CA-C-N	5.70	125.23	119.19
1	A	869	MET	C-N-CA	5.70	125.23	119.19
1	A	247	TYR	N-CA-C	-5.20	101.21	108.54
1	A	1068	PHE	CA-C-N	5.11	125.51	119.94
1	A	1068	PHE	C-N-CA	5.11	125.51	119.94

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	313	SER	Peptide

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	7198	0	6945	85	0
2	B	28	0	25	2	0
2	C	28	0	25	2	0
2	D	28	0	25	12	0
2	E	28	0	25	5	0
All	All	7310	0	7045	85	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 (85) 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:1000:ASN:HD21	2:E:1:NAG:C1	1.29	1.43
1:A:1000:ASN:ND2	2:E:1:NAG:C1	1.89	1.35
1:A:141:ASN:ND2	2:B:1:NAG:C1	2.16	1.09
1:A:414:ASN:ND2	2:D:1:NAG:H2	1.73	1.02
1:A:1000:ASN:OD1	2:E:1:NAG:H83	1.60	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:ASN:HD22	2:B:1:NAG:C1	1.75	0.97
1:A:414:ASN:ND2	2:D:1:NAG:C2	2.40	0.85
1:A:412:ALA:HB2	2:D:1:NAG:H83	1.57	0.84
1:A:312:ASN:HD22	2:C:1:NAG:C1	1.92	0.83
1:A:414:ASN:HD21	2:D:1:NAG:H4	1.47	0.80
1:A:414:ASN:HD21	2:D:1:NAG:C4	1.96	0.77
1:A:414:ASN:HD22	2:D:1:NAG:H2	1.50	0.74
1:A:476:LEU:HD11	1:A:595:ILE:HG13	1.75	0.69
1:A:999:SER:HA	1:A:1002:THR:HG22	1.74	0.68
1:A:1150:ARG:O	1:A:1154:ALA:HB3	1.94	0.67
1:A:999:SER:HA	1:A:1002:THR:CG2	2.25	0.65
1:A:163:LYS:HD2	1:A:416:THR:HG23	1.77	0.65
1:A:1000:ASN:CG	2:E:1:NAG:C1	2.67	0.64
1:A:95:GLN:HE21	1:A:597:SER:HB2	1.63	0.64
1:A:156:GLN:HE21	1:A:400:GLN:HE22	1.46	0.63
1:A:414:ASN:ND2	2:D:1:NAG:O5	2.32	0.62
1:A:1107:ALA:HB2	1:A:1114:ARG:HB2	1.83	0.60
1:A:1000:ASN:ND2	2:E:1:NAG:O5	2.14	0.60
1:A:547:SER:HB3	1:A:1102:LEU:HD22	1.84	0.59
1:A:449:MET:HG3	1:A:508:LEU:HD21	1.84	0.59
1:A:999:SER:CA	1:A:1002:THR:HG22	2.33	0.59
1:A:414:ASN:ND2	2:D:1:NAG:H4	2.17	0.59
1:A:411:VAL:HG13	1:A:419:VAL:HG22	1.86	0.58
1:A:338:GLN:HE21	1:A:835:GLU:HB3	1.68	0.57
1:A:414:ASN:HD21	2:D:1:NAG:C5	2.21	0.54
1:A:330:LEU:HG	1:A:337:TRP:HH2	1.73	0.54
1:A:164:GLU:HB2	1:A:167:ALA:HB2	1.90	0.53
1:A:411:VAL:CG1	1:A:419:VAL:HG22	2.38	0.53
1:A:414:ASN:ND2	2:D:1:NAG:C1	2.72	0.53
1:A:312:ASN:ND2	2:C:1:NAG:C1	2.68	0.52
1:A:808:GLN:HA	1:A:970:GLU:O	2.10	0.52
1:A:1081:VAL:HB	1:A:1082:PRO:HD3	1.91	0.52
1:A:998:CYS:O	1:A:1002:THR:HG22	2.11	0.51
1:A:736:ALA:O	1:A:740:TYR:HB2	2.10	0.51
1:A:91:GLY:HA3	1:A:538:GLY:HA2	1.92	0.50
1:A:441:ARG:NH2	1:A:501:GLN:OE1	2.45	0.50
1:A:1122:MET:O	1:A:1126:VAL:HG23	2.11	0.50
1:A:969:ILE:HD11	1:A:972:ALA:HB2	1.93	0.49
1:A:862:ASP:O	1:A:867:LYS:N	2.44	0.49
1:A:1149:VAL:O	1:A:1153:PHE:HB3	2.13	0.48
1:A:163:LYS:HE3	1:A:416:THR:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:282:LEU:HG	1:A:287:VAL:HG13	1.96	0.48
1:A:1013:TYR:N	1:A:1014:PRO:HD2	2.29	0.48
1:A:1079:SER:O	1:A:1082:PRO:HD2	2.14	0.47
1:A:76:PRO:HG2	1:A:77:LEU:HD12	1.97	0.47
1:A:296:CYS:HB3	1:A:309:PRO:HG2	1.97	0.47
1:A:308:ALA:HB3	1:A:309:PRO:HD3	1.97	0.47
1:A:87:LEU:HD21	1:A:545:GLY:HA3	1.97	0.47
1:A:736:ALA:O	1:A:740:TYR:CB	2.63	0.46
1:A:140:LEU:HA	1:A:143:THR:HG22	1.97	0.46
1:A:189:HIS:HB3	1:A:199:LEU:HB3	1.96	0.46
1:A:806:VAL:HG21	1:A:971:TYR:CE1	2.51	0.45
1:A:485:LEU:HD21	1:A:503:LEU:HD23	1.99	0.45
1:A:414:ASN:ND2	2:D:1:NAG:C4	2.73	0.44
1:A:1056:ILE:HG12	1:A:1097:THR:HG22	2.00	0.44
1:A:349:ASN:HB3	1:A:354:LEU:HD22	2.00	0.44
1:A:806:VAL:HA	1:A:973:GLN:HA	2.00	0.43
1:A:414:ASN:O	1:A:415:SER:C	2.58	0.43
1:A:162:PRO:HA	1:A:417:GLN:HA	2.01	0.43
1:A:296:CYS:O	1:A:297:LEU:HG	2.19	0.43
1:A:531:ILE:HB	1:A:532:PRO:HD3	2.01	0.43
1:A:434:PHE:O	1:A:501:GLN:NE2	2.46	0.43
1:A:168:ASN:HD21	1:A:355:VAL:HG22	1.85	0.42
1:A:139:GLU:OE2	1:A:1009:TYR:HB2	2.18	0.42
1:A:833:MET:HG3	1:A:979:ASN:HD22	1.84	0.42
1:A:1060:LEU:HD11	1:A:1094:VAL:HG22	2.00	0.42
1:A:414:ASN:HD21	2:D:1:NAG:C2	2.29	0.42
1:A:157:LEU:HG	1:A:363:MET:HG2	2.02	0.42
1:A:496:ASN:HD22	1:A:498:ALA:HB3	1.84	0.42
1:A:199:LEU:HD23	1:A:199:LEU:H	1.84	0.42
1:A:587:MET:O	1:A:591:ILE:HB	2.20	0.42
1:A:1040:PHE:HB2	1:A:1057:VAL:HG11	2.02	0.41
1:A:1074:ILE:HG23	1:A:1076:ILE:HG12	2.02	0.41
1:A:1171:LEU:HA	1:A:1174:LEU:HB2	2.03	0.41
1:A:496:ASN:HD21	1:A:569:ALA:HB1	1.85	0.41
1:A:203:CYS:HA	1:A:366:LEU:HA	2.03	0.41
1:A:87:LEU:HD11	1:A:541:LEU:HD12	2.03	0.41
1:A:518:LEU:HD12	1:A:588:VAL:HG22	2.03	0.41
1:A:1036:LEU:HD23	1:A:1036:LEU:HA	1.96	0.40
1:A:855:LEU:HG	1:A:883:ALA:HB2	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	958/1455 (66%)	858 (90%)	99 (10%)	1 (0%)	48 79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	261	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	704/1238 (57%)	698 (99%)	6 (1%)	70 76

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	508	LEU
1	A	583	PHE
1	A	806	VAL
1	A	1004	LEU
1	A	1052	THR
1	A	1105	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	GLN
1	A	124	ASN
1	A	141	ASN
1	A	154	ASN
1	A	156	GLN
1	A	283	ASN
1	A	312	ASN
1	A	328	HIS
1	A	338	GLN
1	A	375	HIS
1	A	414	ASN
1	A	496	ASN
1	A	739	HIS
1	A	808	GLN
1	A	1000	ASN
1	A	1099	HIS
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](#)

8 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	14,14,15	0.89	1 (7%)	17,19,21	1.80	5 (29%)
2	NAG	B	2	2	14,14,15	1.01	1 (7%)	17,19,21	1.79	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	1	2	14,14,15	0.79	1 (7%)	17,19,21	2.13	8 (47%)
2	NAG	C	2	2	14,14,15	0.82	0	17,19,21	1.31	2 (11%)
2	NAG	D	1	2	14,14,15	0.99	1 (7%)	17,19,21	1.59	2 (11%)
2	NAG	D	2	2	14,14,15	0.81	1 (7%)	17,19,21	1.58	2 (11%)
2	NAG	E	1	2	14,14,15	0.61	0	17,19,21	1.50	3 (17%)
2	NAG	E	2	2	14,14,15	0.58	0	17,19,21	1.43	2 (11%)

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	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	NAG	C	1	2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	NAG	D	1	2	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1
2	NAG	E	1	2	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	NAG	C1-C2	3.42	1.57	1.52
2	D	2	NAG	C1-C2	2.42	1.55	1.52
2	D	1	NAG	C1-C2	2.34	1.55	1.52
2	B	1	NAG	O5-C1	-2.19	1.40	1.43
2	C	1	NAG	C1-C2	2.16	1.55	1.52

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	C1-O5-C5	5.38	119.40	112.19
2	D	2	NAG	C2-N2-C7	4.55	128.99	122.90
2	D	1	NAG	C2-N2-C7	4.01	128.28	122.90
2	C	1	NAG	C2-N2-C7	3.99	128.24	122.90
2	E	2	NAG	C1-O5-C5	3.74	117.20	112.19
2	B	1	NAG	O4-C4-C5	3.73	118.51	109.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C4-C3-C2	-3.54	105.83	111.02
2	E	1	NAG	C1-C2-N2	-3.36	105.14	110.43
2	C	1	NAG	C1-O5-C5	3.36	116.69	112.19
2	C	2	NAG	C2-N2-C7	3.19	127.17	122.90
2	B	1	NAG	O5-C1-C2	3.12	116.11	111.29
2	B	1	NAG	C2-N2-C7	3.04	126.97	122.90
2	E	1	NAG	C2-N2-C7	2.98	126.90	122.90
2	C	1	NAG	O5-C5-C4	-2.84	103.92	110.83
2	C	1	NAG	O4-C4-C5	2.78	116.17	109.32
2	C	1	NAG	O7-C7-N2	2.61	126.60	121.98
2	B	1	NAG	O5-C5-C4	-2.48	104.79	110.83
2	E	2	NAG	C2-N2-C7	2.39	126.10	122.90
2	B	1	NAG	O5-C5-C6	-2.36	103.06	107.66
2	B	2	NAG	C2-N2-C7	2.36	126.06	122.90
2	E	1	NAG	C4-C3-C2	-2.32	107.62	111.02
2	D	1	NAG	O5-C5-C6	-2.28	103.22	107.66
2	C	2	NAG	O5-C5-C4	-2.26	105.33	110.83
2	B	2	NAG	O5-C1-C2	-2.26	107.80	111.29
2	C	1	NAG	C8-C7-N2	-2.19	112.49	116.12
2	C	1	NAG	C3-C4-C5	-2.13	106.36	110.23
2	D	2	NAG	C1-O5-C5	2.07	114.96	112.19

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1	NAG	C1-C2-N2-C7
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
2	C	1	NAG	C3-C2-N2-C7
2	D	1	NAG	C3-C2-N2-C7
2	C	1	NAG	C1-C2-N2-C7
2	D	2	NAG	C1-C2-N2-C7
2	D	2	NAG	C3-C2-N2-C7

There are no ring outliers.

4 monomers are involved in 21 short contacts:

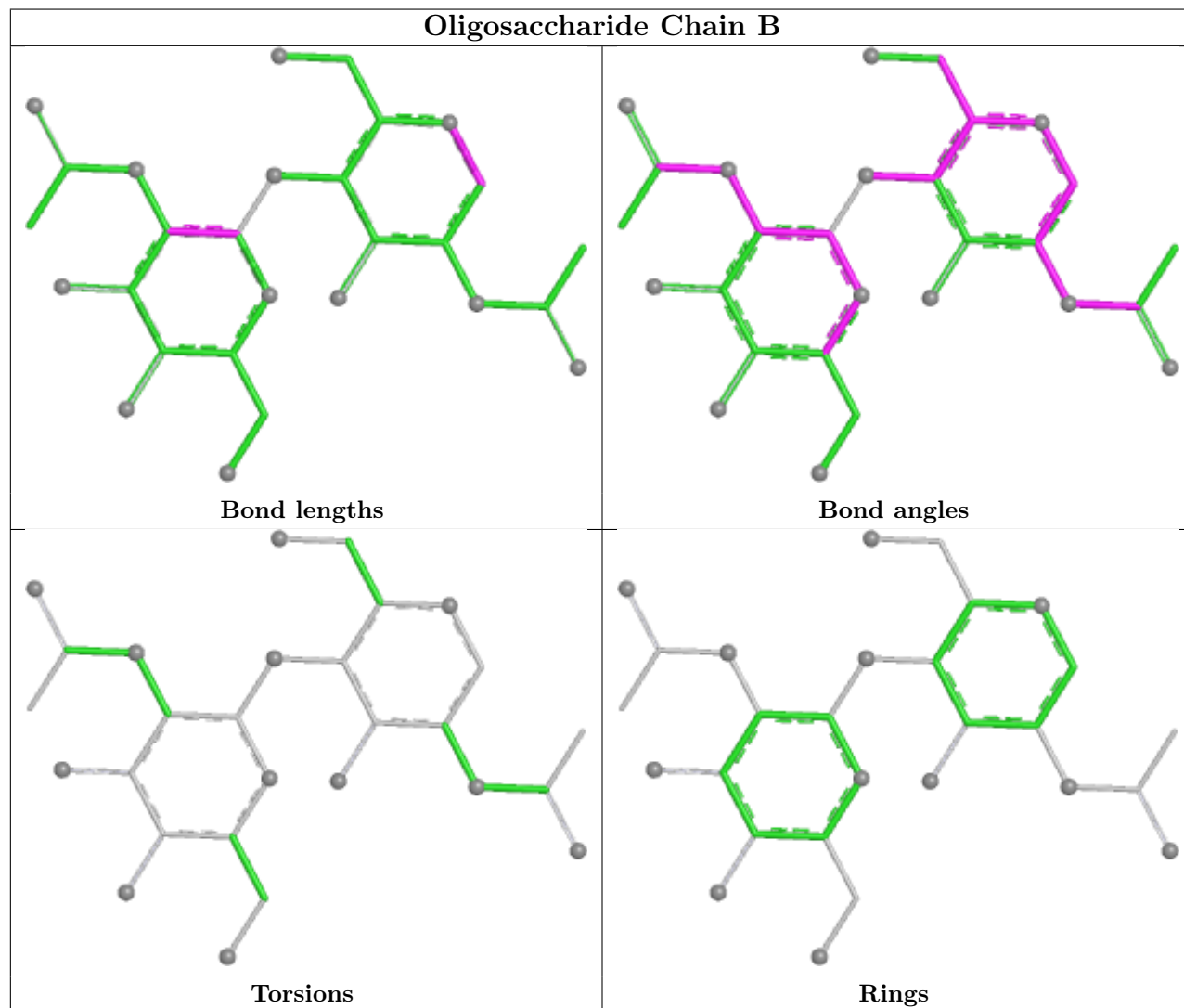
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	NAG	2	0
2	C	1	NAG	2	0

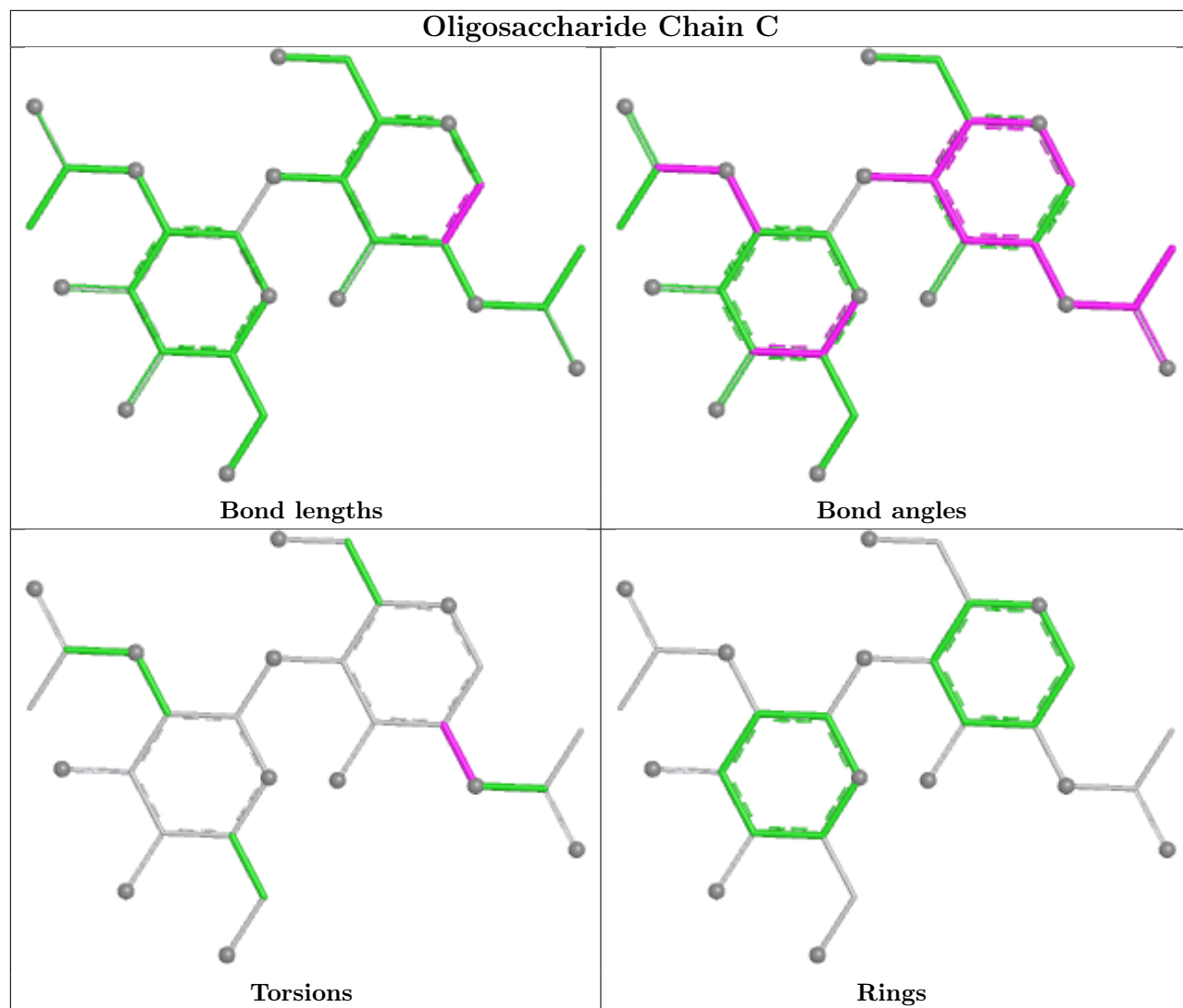
Continued on next page...

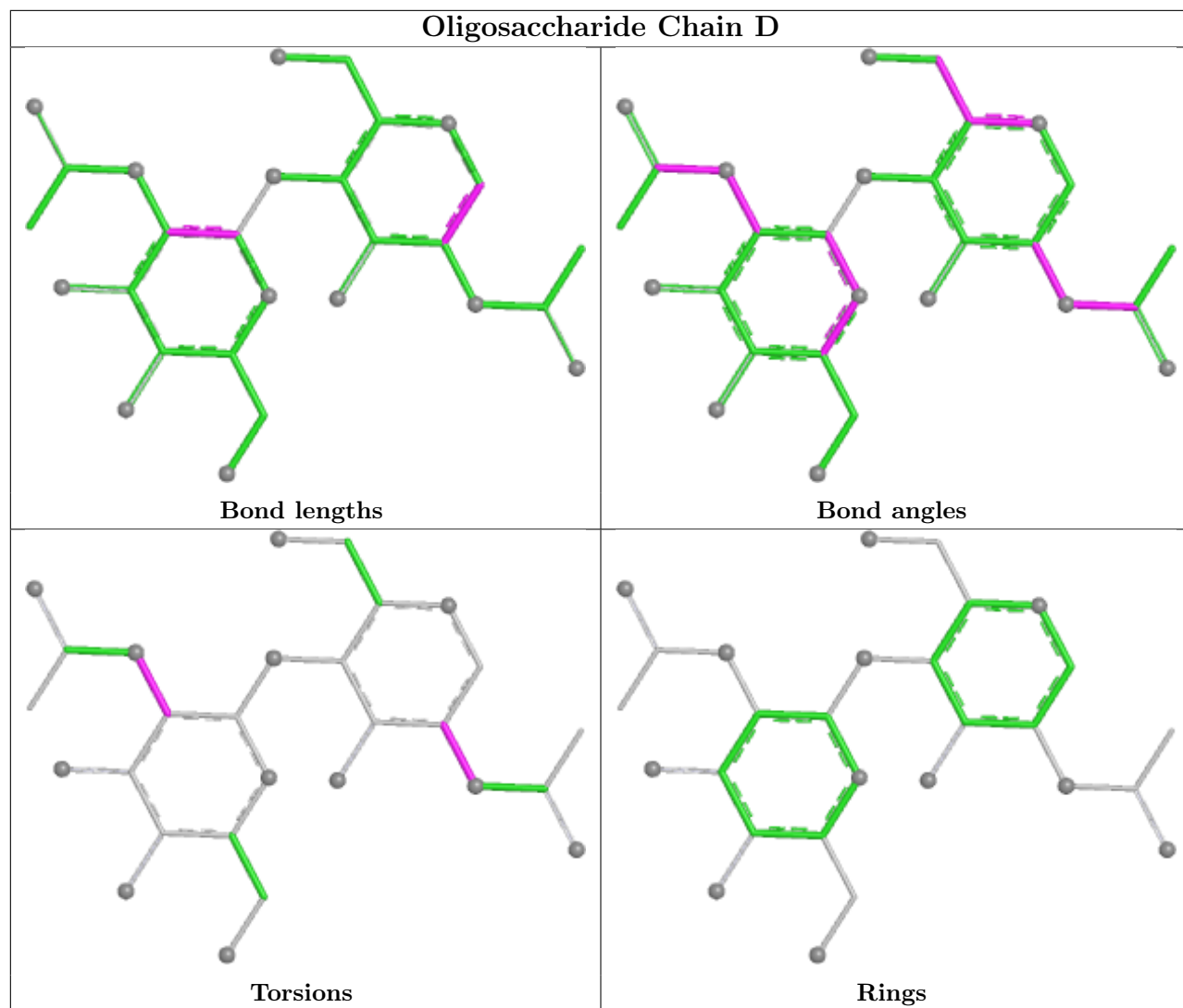
Continued from previous page...

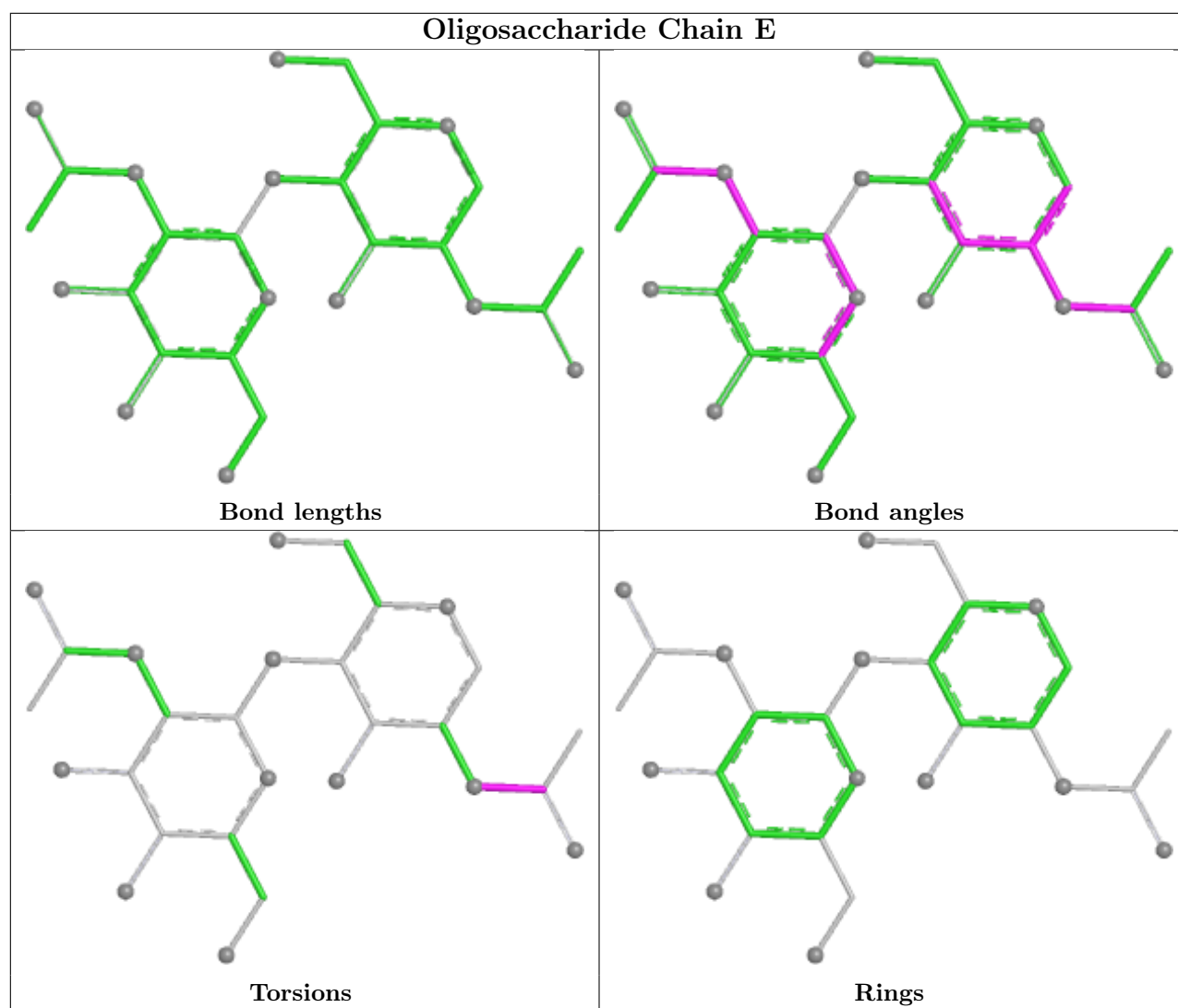
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	1	NAG	5	0
2	D	1	NAG	12	0

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









5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

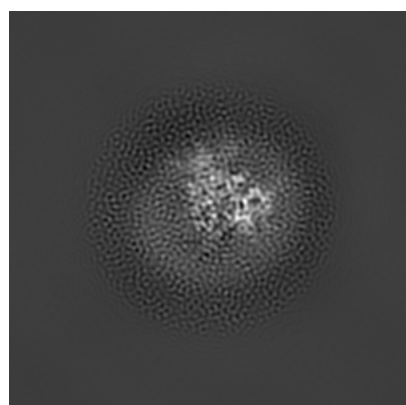
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-7795. These allow visual inspection of the internal detail of the map and identification of artifacts.

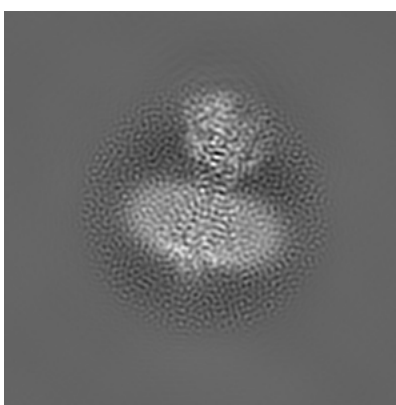
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

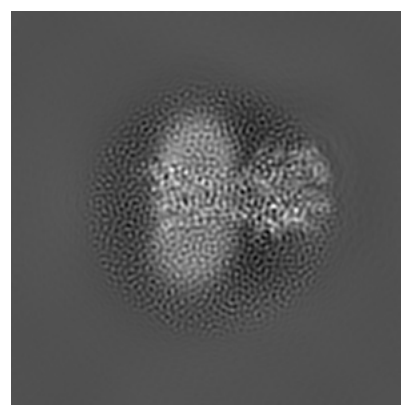
6.1.1 Primary map



X



Y

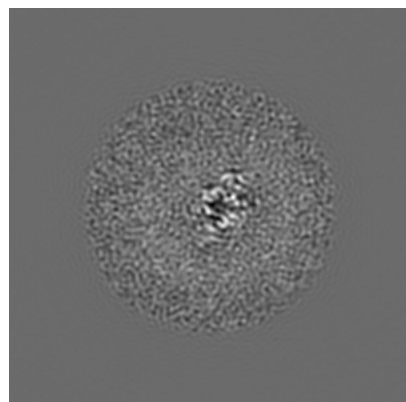


Z

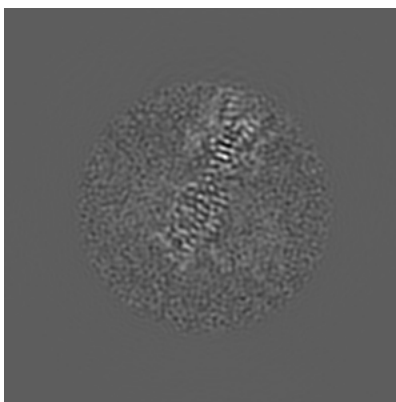
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

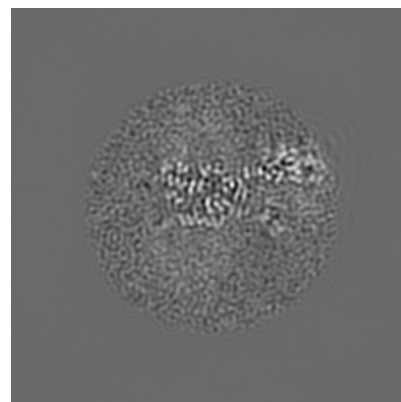
6.2.1 Primary map



X Index: 128



Y Index: 128

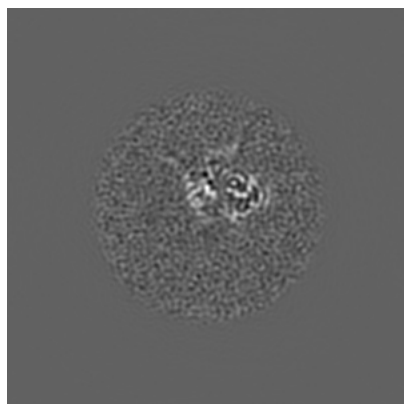


Z Index: 128

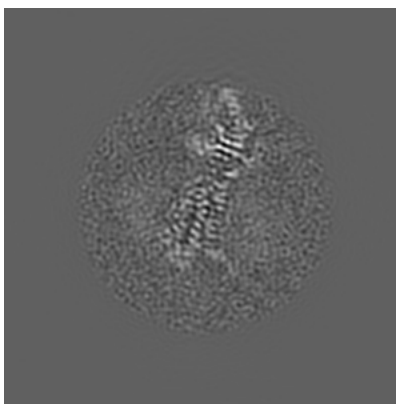
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

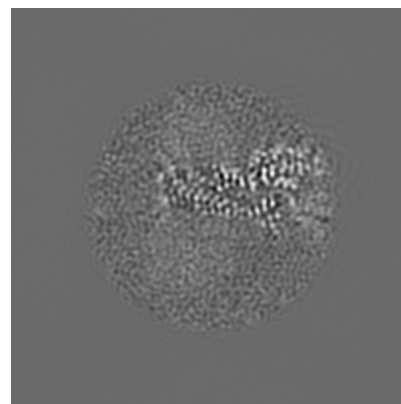
6.3.1 Primary map



X Index: 161



Y Index: 130

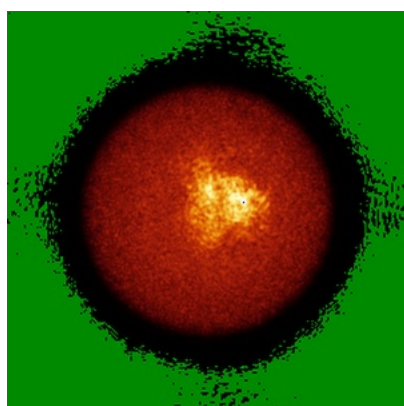


Z Index: 136

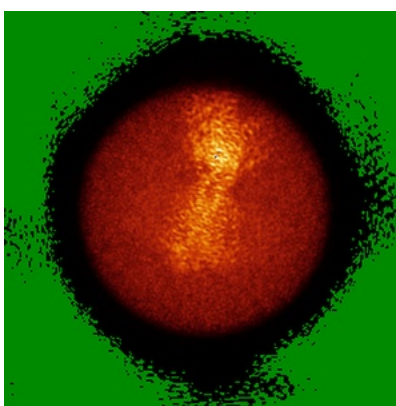
The images above show the largest variance slices of the map in three orthogonal directions.

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

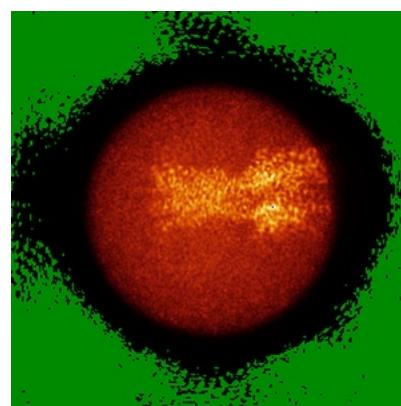
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views

This section was not generated.

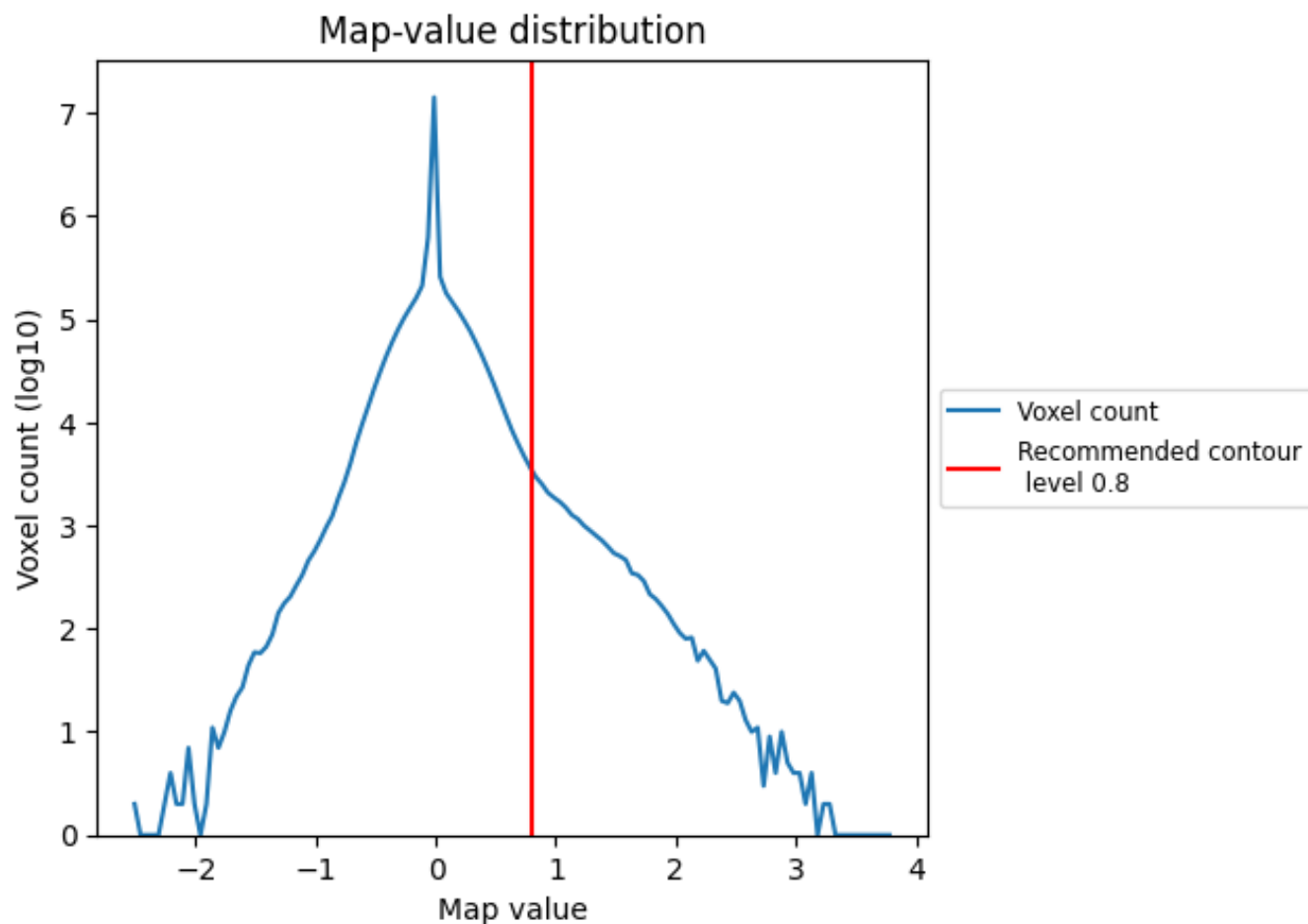
6.6 Mask visualisation

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

7 Map analysis [i](#)

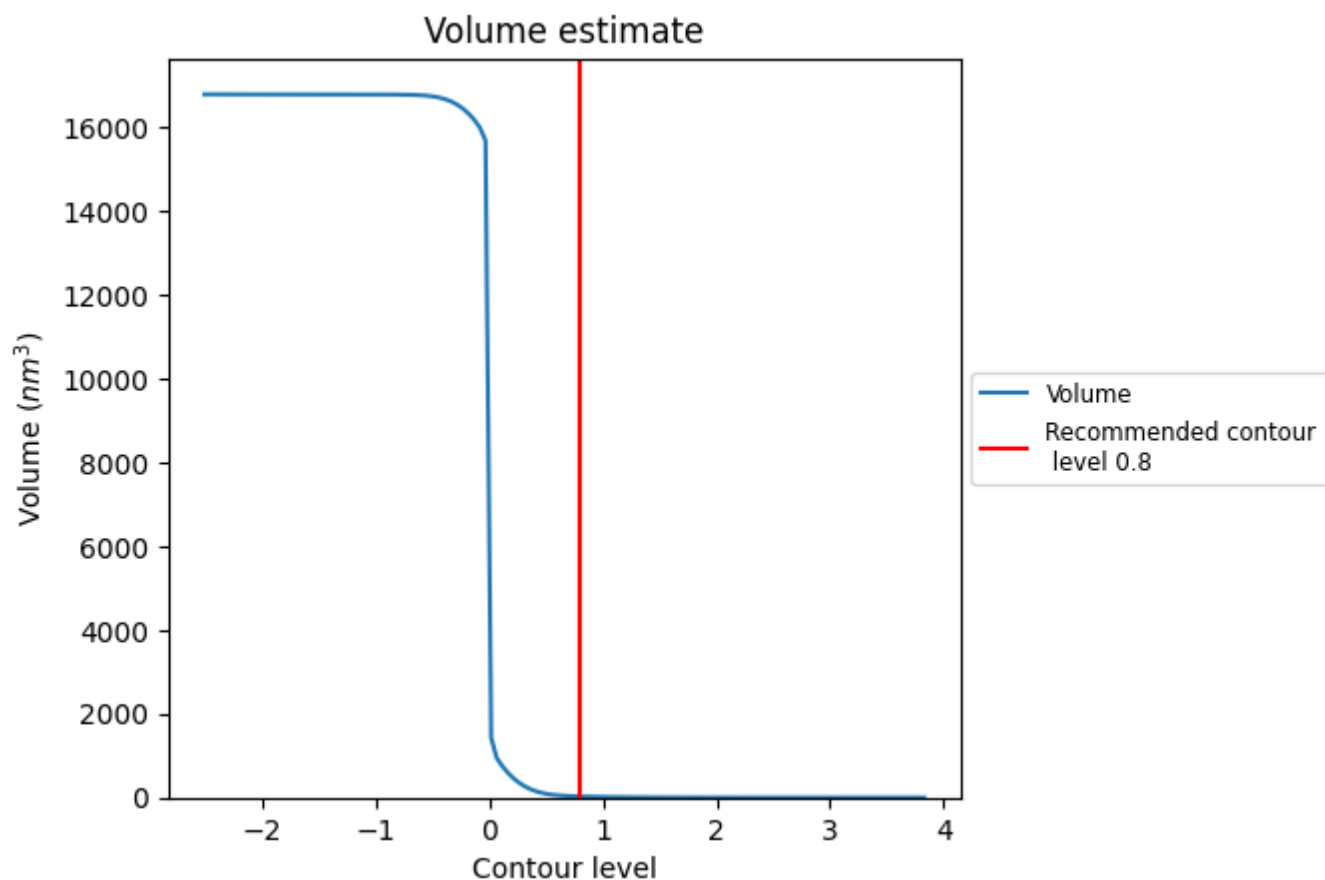
This section contains the results of statistical analysis of the map.

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.

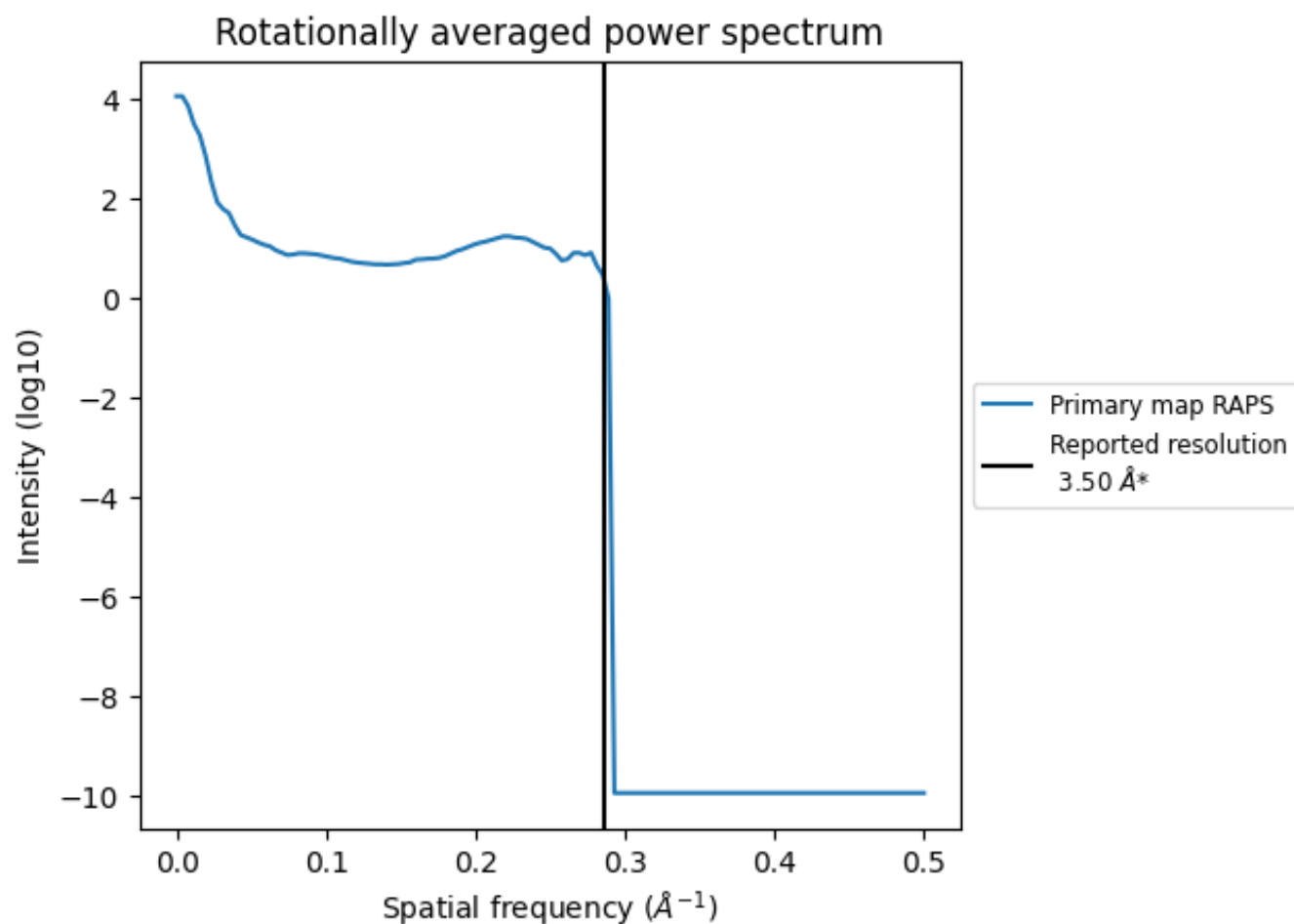
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 25 nm³; this corresponds to an approximate mass of 23 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



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

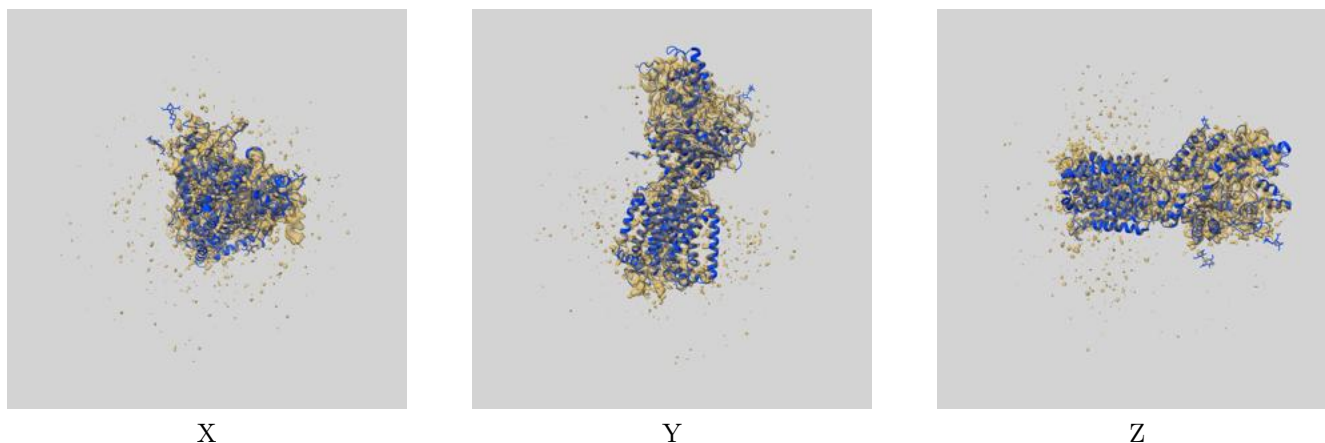
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

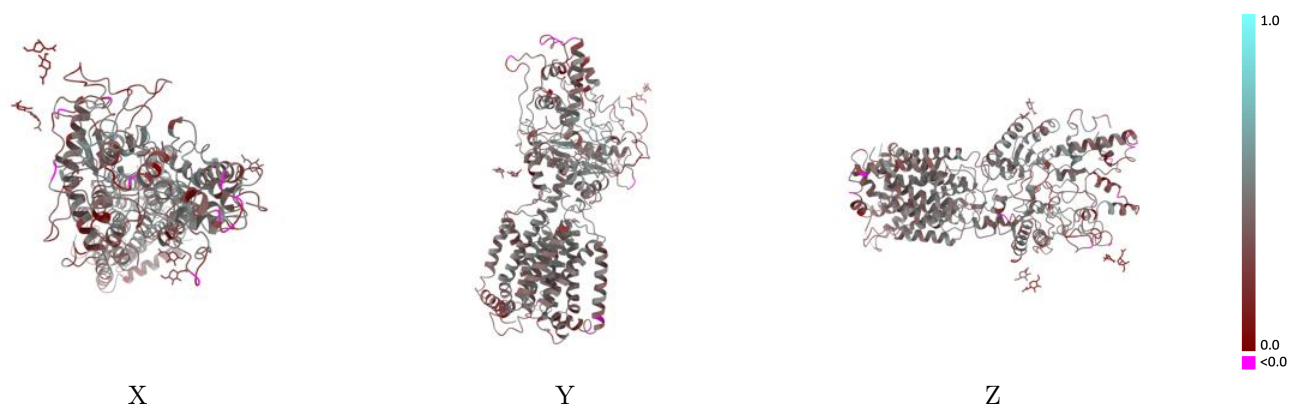
This section contains information regarding the fit between EMDB map EMD-7795 and PDB model 6OEU. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



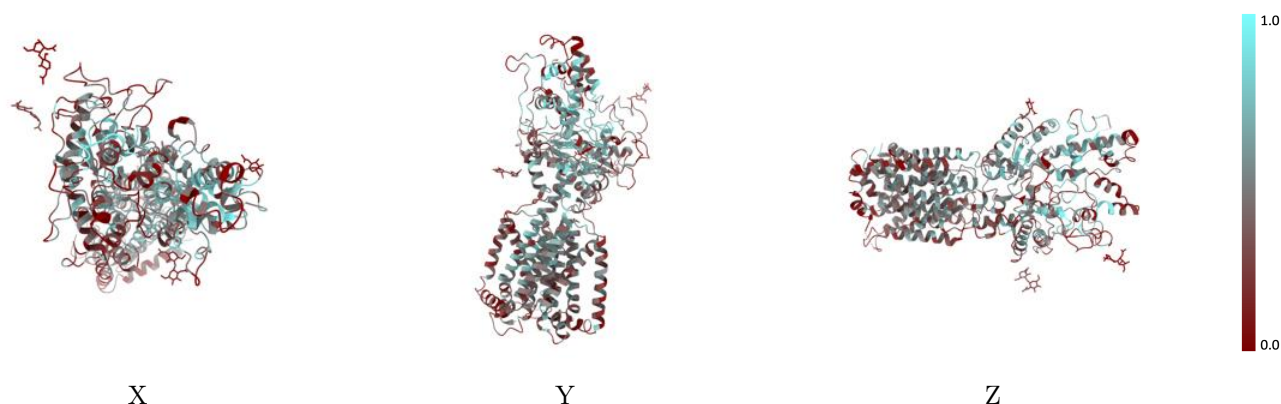
The images above show the 3D surface view of the map at the recommended contour level 0.8 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



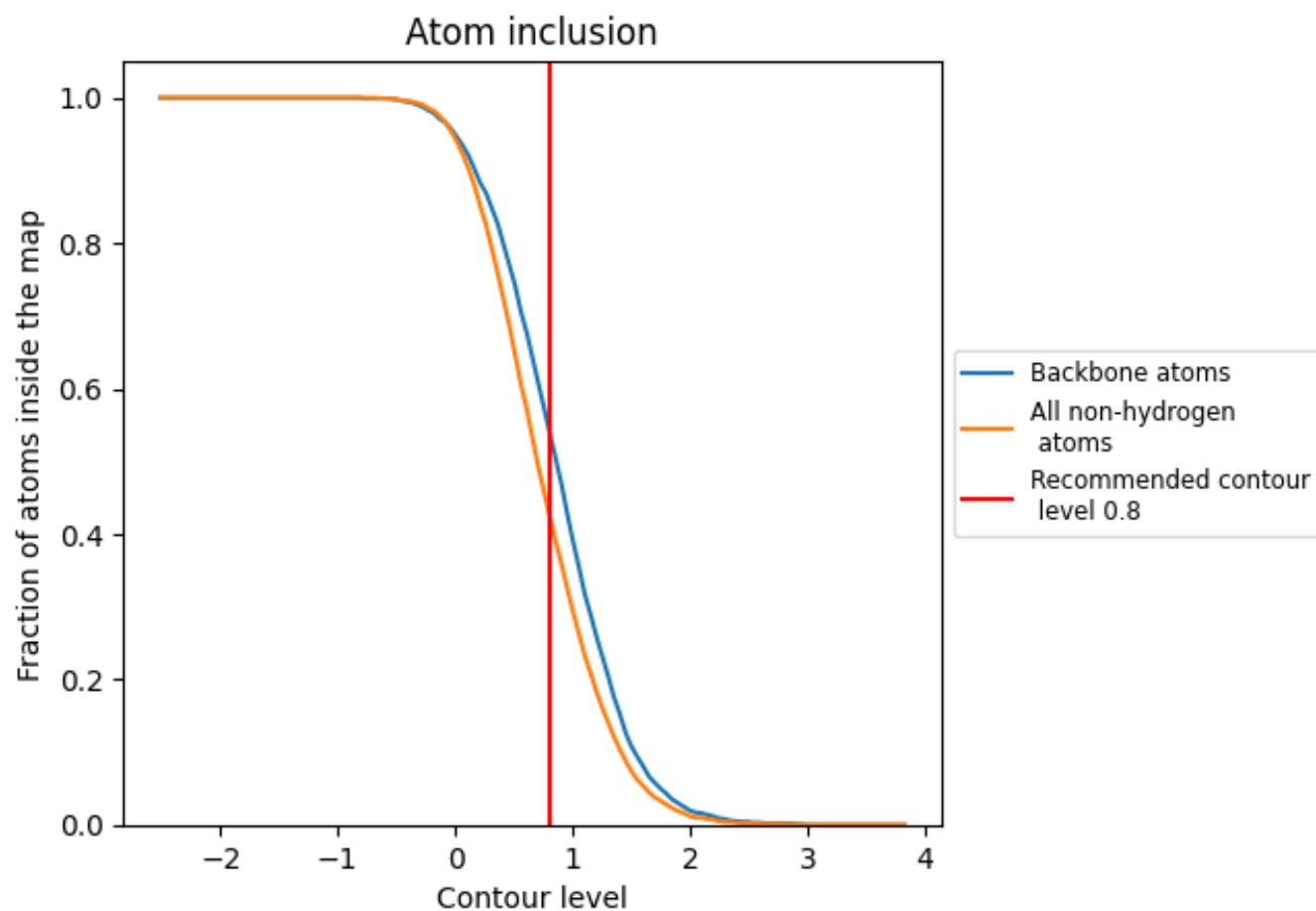
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.8).

9.4 Atom inclusion [i](#)



At the recommended contour level, 54% of all backbone atoms, 43% 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 (0.8) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.4280	0.3930
A	0.4320	0.3950
B	0.1790	0.2630
C	0.0000	0.1400
D	0.2500	0.2270
E	0.1430	0.3460

