



Full wwPDB EM Validation Report ⓘ

Oct 6, 2024 – 12:33 PM JST

PDB ID : 8Y8A
EMDB ID : EMD-39039
Title : Structure of HCoV-HKU1C spike in the functionally anchored-3up conformation with 3TMPRSS2
Authors : Lu, Y.C.; Zhang, X.; Wang, H.F.; Liu, X.C.; Sun, L.; Yang, H.T.
Deposited on : 2024-02-06
Resolution : 3.19 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.dev113
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

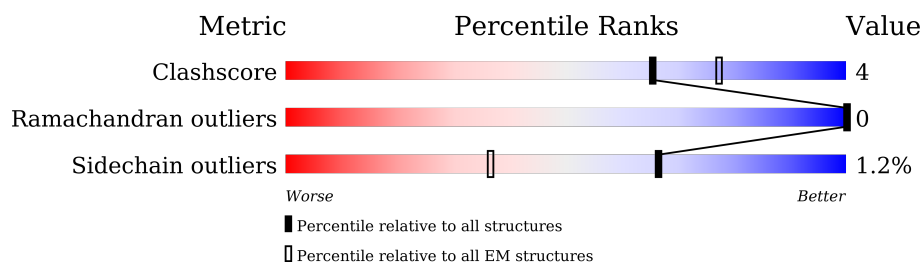
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.19 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	G	383	
1	H	383	
1	T	383	
2	A	1263	
2	B	1263	
2	C	1263	
3	D	2	
3	E	2	

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Mol	Chain	Length	Quality of chain
3	F	2	 100%
3	J	2	 100%
3	L	2	 100%
3	M	2	 100%
3	N	2	 100%
3	O	2	 100%
3	P	2	 100%
3	Q	2	 100%
3	R	2	 100%
4	I	3	 100%
4	K	3	 100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 33387 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 Transmembrane protease serine 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	T	237	Total	C	N	O	S	0	0
			1725	1106	292	314	13		
1	H	237	Total	C	N	O	S	0	0
			1727	1108	292	314	13		
1	G	237	Total	C	N	O	S	0	0
			1727	1108	292	314	13		

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
T	251	ASP	SER	engineered mutation	UNP O15393
T	252	ASP	SER	engineered mutation	UNP O15393
T	?	-	ARG	deletion	UNP O15393
T	253	ASP	GLN	engineered mutation	UNP O15393
T	254	ASP	SER	engineered mutation	UNP O15393
T	255	LYS	ARG	engineered mutation	UNP O15393
H	251	ASP	SER	engineered mutation	UNP O15393
H	252	ASP	SER	engineered mutation	UNP O15393
H	?	-	ARG	deletion	UNP O15393
H	253	ASP	GLN	engineered mutation	UNP O15393
H	254	ASP	SER	engineered mutation	UNP O15393
H	255	LYS	ARG	engineered mutation	UNP O15393
G	251	ASP	SER	engineered mutation	UNP O15393
G	252	ASP	SER	engineered mutation	UNP O15393
G	?	-	ARG	deletion	UNP O15393
G	253	ASP	GLN	engineered mutation	UNP O15393
G	254	ASP	SER	engineered mutation	UNP O15393
G	255	LYS	ARG	engineered mutation	UNP O15393

- Molecule 2 is a protein called Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	1185	Total	C	N	O	S	0	0
			9191	5871	1507	1757	56		
2	C	1194	Total	C	N	O	S	0	0
			9236	5890	1515	1774	57		
2	B	1194	Total	C	N	O	S	0	0
			9255	5907	1517	1774	57		

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	752	GLY	ARG	engineered mutation	UNP Q0ZME7
A	753	GLY	ARG	engineered mutation	UNP Q0ZME7
A	754	SER	LYS	engineered mutation	UNP Q0ZME7
A	755	GLY	ARG	engineered mutation	UNP Q0ZME7
A	756	SER	ARG	engineered mutation	UNP Q0ZME7
A	902	PRO	LEU	engineered mutation	UNP Q0ZME7
A	980	PRO	SER	engineered mutation	UNP Q0ZME7
A	1023	PRO	ASN	engineered mutation	UNP Q0ZME7
A	1067	PRO	ASN	engineered mutation	UNP Q0ZME7
A	1068	PRO	LEU	engineered mutation	UNP Q0ZME7
C	752	GLY	ARG	engineered mutation	UNP Q0ZME7
C	753	GLY	ARG	engineered mutation	UNP Q0ZME7
C	754	SER	LYS	engineered mutation	UNP Q0ZME7
C	755	GLY	ARG	engineered mutation	UNP Q0ZME7
C	756	SER	ARG	engineered mutation	UNP Q0ZME7
C	902	PRO	LEU	engineered mutation	UNP Q0ZME7
C	980	PRO	SER	engineered mutation	UNP Q0ZME7
C	1023	PRO	ASN	engineered mutation	UNP Q0ZME7
C	1067	PRO	ASN	engineered mutation	UNP Q0ZME7
C	1068	PRO	LEU	engineered mutation	UNP Q0ZME7
B	752	GLY	ARG	engineered mutation	UNP Q0ZME7
B	753	GLY	ARG	engineered mutation	UNP Q0ZME7
B	754	SER	LYS	engineered mutation	UNP Q0ZME7
B	755	GLY	ARG	engineered mutation	UNP Q0ZME7
B	756	SER	ARG	engineered mutation	UNP Q0ZME7
B	902	PRO	LEU	engineered mutation	UNP Q0ZME7
B	980	PRO	SER	engineered mutation	UNP Q0ZME7
B	1023	PRO	ASN	engineered mutation	UNP Q0ZME7
B	1067	PRO	ASN	engineered mutation	UNP Q0ZME7
B	1068	PRO	LEU	engineered mutation	UNP Q0ZME7

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



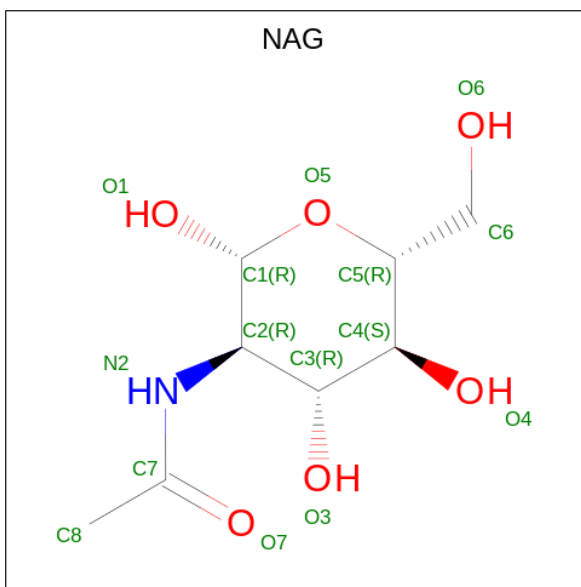
Mol	Chain	Residues	Atoms				AltConf	Trace
3	D	2	Total	C	N	O	0	0
			28	16	2	10		
3	E	2	Total	C	N	O	0	0
			28	16	2	10		
3	F	2	Total	C	N	O	0	0
			28	16	2	10		
3	J	2	Total	C	N	O	0	0
			28	16	2	10		
3	L	2	Total	C	N	O	0	0
			28	16	2	10		
3	M	2	Total	C	N	O	0	0
			28	16	2	10		
3	N	2	Total	C	N	O	0	0
			28	16	2	10		
3	O	2	Total	C	N	O	0	0
			28	16	2	10		
3	P	2	Total	C	N	O	0	0
			28	16	2	10		
3	Q	2	Total	C	N	O	0	0
			28	16	2	10		
3	R	2	Total	C	N	O	0	0
			28	16	2	10		

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

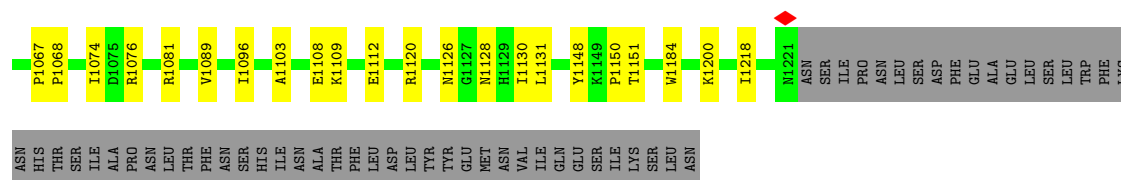


Mol	Chain	Residues	Atoms				AltConf	Trace
4	I	3	Total	C	N	O	0	0
			39	22	2	15		
4	K	3	Total	C	N	O	0	0
			39	22	2	15		

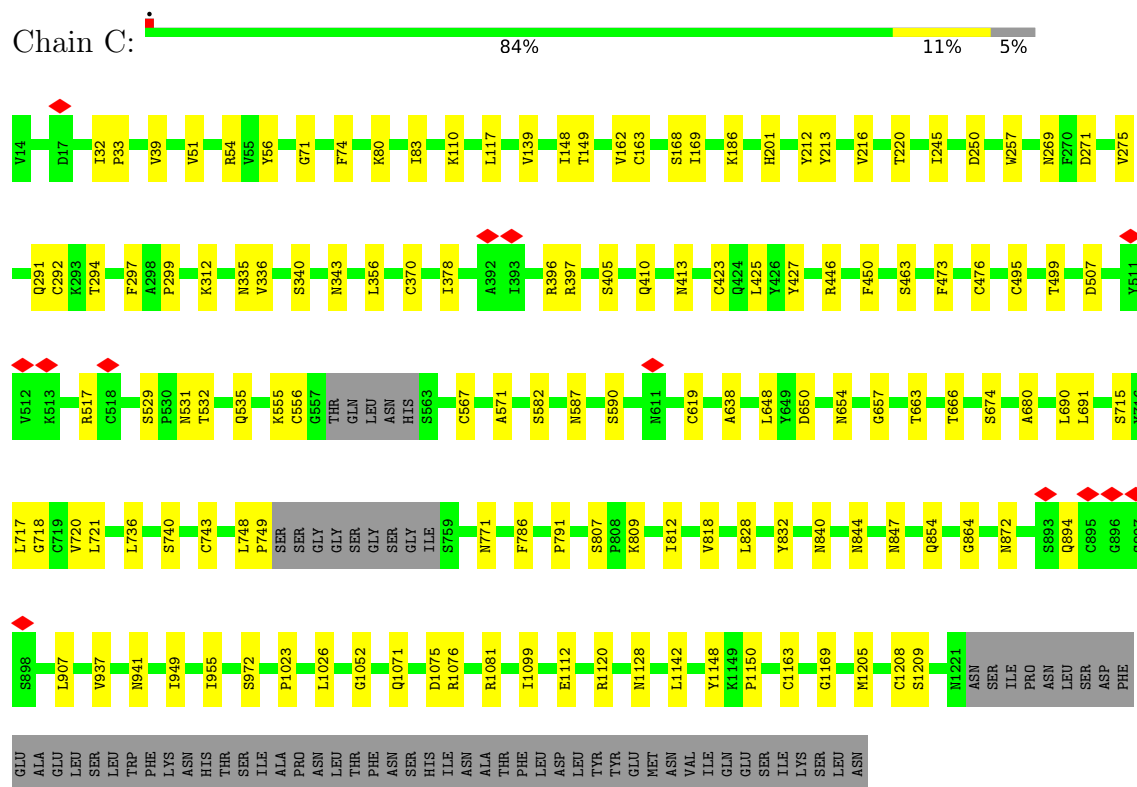
- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



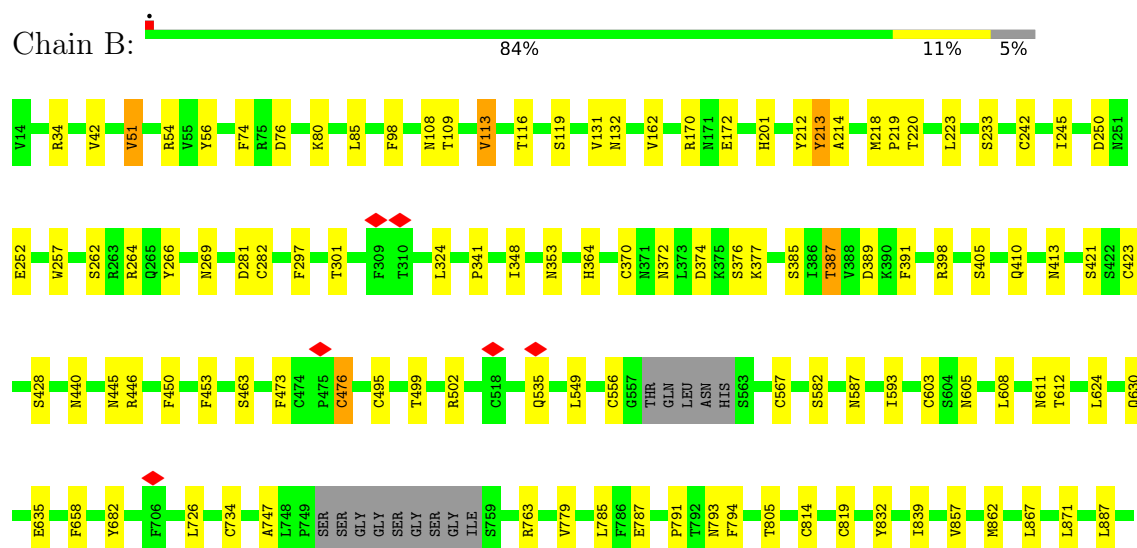
Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5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

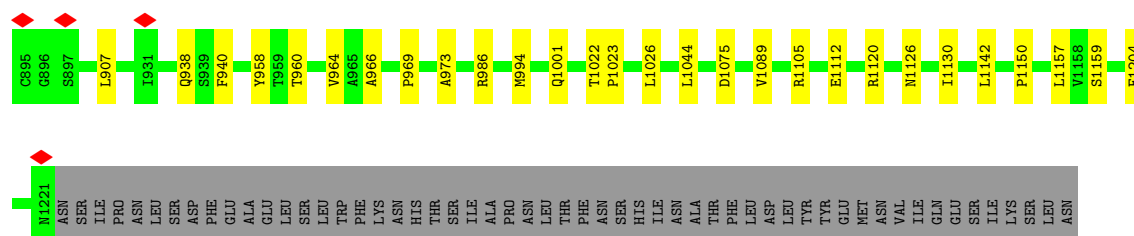


• Molecule 2: Spike glycoprotein



• Molecule 2: Spike glycoprotein





- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain R:  100%



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

Chain I:  100%



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

Chain K:

100%



MAG1
MAG2
EM13

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	142943	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.870	Depositor
Minimum map value	-0.002	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.032	Depositor
Recommended contour level	0.006	Depositor
Map size (Å)	328.0, 328.0, 328.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.82, 0.82, 0.82	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: BMA, 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	G	0.23	0/1776	0.42	0/2441
1	H	0.23	0/1776	0.42	0/2441
1	T	0.23	0/1774	0.42	0/2438
2	A	0.24	0/9416	0.41	0/12833
2	B	0.24	0/9482	0.41	0/12923
2	C	0.24	0/9462	0.41	0/12901
All	All	0.24	0/33686	0.41	0/45977

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 [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	G	1727	0	1605	14	0
1	H	1727	0	1605	10	0
1	T	1725	0	1598	18	0
2	A	9191	0	8797	77	0
2	B	9255	0	8847	75	0
2	C	9236	0	8839	71	0
3	D	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	28	0	25	0	0
3	F	28	0	25	0	0
3	J	28	0	25	0	0
3	L	28	0	25	0	0
3	M	28	0	25	0	0
3	N	28	0	25	0	0
3	O	28	0	25	0	0
3	P	28	0	25	0	0
3	Q	28	0	25	0	0
3	R	28	0	25	0	0
4	I	39	0	34	0	0
4	K	39	0	34	0	0
5	A	56	0	52	0	0
5	B	42	0	39	0	0
5	C	42	0	39	0	0
All	All	33387	0	31764	248	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 4.

All (248) 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:A:201:HIS:HB2	2:A:212:TYR:HB2	1.73	0.70
2:B:440:ASN:H	2:B:446:ARG:HH21	1.38	0.70
2:B:162:VAL:HG22	2:B:172:GLU:HB2	1.77	0.66
2:A:701:LEU:HA	2:A:705:SER:HB3	1.78	0.65
2:A:1120:ARG:NH2	2:C:1112:GLU:OE2	2.28	0.65
2:B:445:ASN:HD21	2:B:549:LEU:H	1.43	0.65
2:C:201:HIS:HB2	2:C:212:TYR:HB2	1.78	0.64
2:C:556:CYS:HA	2:C:567:CYS:HA	1.79	0.64
2:B:201:HIS:HB2	2:B:212:TYR:HB2	1.79	0.64
2:C:80:LYS:HE3	2:C:245:ILE:HB	1.79	0.63
2:A:871:LEU:HD13	2:B:779:VAL:HG21	1.81	0.63
2:A:691:LEU:HD13	2:A:720:VAL:HG22	1.80	0.62
2:A:1112:GLU:OE1	2:A:1120:ARG:NH1	2.34	0.60
2:B:556:CYS:HA	2:B:567:CYS:HA	1.83	0.60
2:A:370:CYS:HA	2:A:423:CYS:HA	1.84	0.60
2:B:1112:GLU:OE1	2:B:1120:ARG:NH2	2.35	0.59
1:H:445:LEU:HD23	1:H:457:GLY:HA3	1.84	0.59
1:G:298:VAL:HG11	1:G:331:VAL:HG11	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:64:THR:HB	2:A:265:GLN:HG2	1.85	0.58
2:C:413:ASN:HD21	2:C:463:SER:H	1.51	0.58
2:A:425:LEU:HB3	2:A:590:SER:HB3	1.86	0.58
2:C:405:SER:O	2:C:410:GLN:NE2	2.37	0.57
2:C:832:TYR:OH	2:C:1075:ASP:OD1	2.22	0.57
2:C:691:LEU:HD13	2:C:720:VAL:HG22	1.87	0.57
2:A:74:PHE:HB3	2:A:257:TRP:HB3	1.86	0.57
2:C:937:VAL:O	2:C:941:ASN:ND2	2.37	0.57
2:C:854:GLN:HE21	2:C:955:ILE:HD11	1.70	0.57
2:B:374:ASP:OD1	2:B:377:LYS:NZ	2.35	0.57
2:A:556:CYS:HA	2:A:567:CYS:HA	1.85	0.57
1:H:311:PHE:HB3	1:H:314:ILE:HD12	1.87	0.57
2:C:162:VAL:HG13	2:C:168:SER:HB3	1.86	0.56
1:T:268:PRO:HB2	1:T:362:LYS:H	1.70	0.56
2:C:907:LEU:HD11	2:C:1142:LEU:HD22	1.86	0.56
2:B:495:CYS:SG	2:B:499:THR:OG1	2.63	0.56
2:A:351:ASN:HD22	2:A:601:THR:HG1	1.51	0.56
2:C:495:CYS:SG	2:C:499:THR:OG1	2.63	0.56
2:A:214:ALA:HB2	2:A:220:THR:HA	1.88	0.56
2:C:840:ASN:O	2:C:844:ASN:ND2	2.38	0.56
1:T:445:LEU:HD23	1:T:457:GLY:HA3	1.87	0.56
1:T:463:SER:OG	1:T:470:ARG:NH1	2.38	0.56
1:H:268:PRO:HB2	1:H:362:LYS:H	1.70	0.56
2:A:801:GLU:OE1	2:A:1109:LYS:NZ	2.39	0.56
2:A:1130:ILE:HG22	2:A:1131:LEU:HG	1.88	0.56
2:C:370:CYS:HA	2:C:423:CYS:HA	1.87	0.56
2:C:51:VAL:HG21	2:C:56:TYR:HE2	1.71	0.55
2:A:80:LYS:NZ	2:A:81:GLY:O	2.39	0.55
2:A:709:GLN:HG3	2:A:721:LEU:HD23	1.88	0.55
2:B:324:LEU:O	2:B:353:ASN:ND2	2.38	0.55
1:T:431:GLN:HB3	2:C:531:ASN:HD21	1.71	0.55
2:A:818:VAL:O	2:A:1081:ARG:NH1	2.40	0.55
2:A:109:THR:HG21	2:A:254:LEU:HD23	1.89	0.55
1:T:311:PHE:HB3	1:T:314:ILE:HD12	1.88	0.55
2:C:1023:PRO:HG2	2:C:1026:LEU:HD23	1.89	0.54
2:B:726:LEU:HD22	2:B:763:ARG:HD2	1.89	0.54
2:B:862:MET:HE2	2:B:958:TYR:HA	1.89	0.54
1:G:445:LEU:HD23	1:G:457:GLY:HA3	1.89	0.53
2:A:306:LEU:HD12	2:A:677:VAL:HB	1.91	0.53
2:C:680:ALA:HB1	2:C:736:LEU:HD13	1.90	0.53
2:A:839:ILE:HG23	2:A:1089:VAL:HG11	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1184:TRP:HB2	2:A:1218:ILE:HD11	1.89	0.53
2:A:132:ASN:HB3	2:A:154:THR:HG22	1.91	0.53
2:C:818:VAL:O	2:C:1081:ARG:NH1	2.42	0.53
2:A:495:CYS:SG	2:A:499:THR:OG1	2.67	0.52
2:A:1055:SER:OG	2:A:1060:GLU:OE2	2.26	0.52
2:B:446:ARG:HA	2:B:450:PHE:HB3	1.92	0.52
2:C:425:LEU:HB3	2:C:590:SER:HB3	1.90	0.52
2:B:986:ARG:NH1	2:B:1130:ILE:O	2.39	0.52
1:G:311:PHE:HB3	1:G:314:ILE:HD12	1.92	0.52
2:A:335:ASN:ND2	2:A:337:SER:OG	2.43	0.52
2:A:505:ASP:OD1	1:G:470:ARG:NH2	2.42	0.52
2:B:80:LYS:HE2	2:B:245:ILE:HB	1.91	0.52
2:B:832:TYR:OH	2:B:1075:ASP:OD1	2.24	0.51
2:C:715:SER:OG	2:C:718:GLY:O	2.26	0.51
2:B:213:TYR:HB3	2:B:223:LEU:HD11	1.92	0.51
2:B:348:ILE:HD11	2:B:385:SER:HB3	1.91	0.51
2:B:387:THR:HG23	2:B:593:ILE:HB	1.92	0.51
2:A:102:ILE:HG13	2:A:258:VAL:HG13	1.93	0.51
2:B:113:VAL:O	2:B:116:THR:OG1	2.21	0.51
2:C:748:LEU:HD12	2:C:749:PRO:HD2	1.91	0.51
2:A:604:SER:OG	2:A:606:ASP:OD1	2.29	0.51
2:B:966:ALA:HB2	2:B:973:ALA:HB3	1.94	0.50
2:B:80:LYS:NZ	2:B:242:CYS:O	2.45	0.50
2:A:1062:LEU:HD11	2:A:1074:ILE:HD11	1.93	0.50
2:A:582:SER:OG	2:A:587:ASN:ND2	2.36	0.50
2:A:167:GLY:O	2:A:243:ASN:ND2	2.44	0.50
2:A:353:ASN:HB3	2:A:605:ASN:HD21	1.76	0.50
2:A:1037:ALA:O	2:A:1041:ASN:ND2	2.32	0.50
2:B:85:LEU:HG	2:B:170:ARG:HH12	1.76	0.50
2:C:74:PHE:HB3	2:C:257:TRP:HB3	1.93	0.50
2:C:555:LYS:HB2	2:C:571:ALA:HB2	1.93	0.50
2:C:674:SER:OG	2:B:940:PHE:O	2.29	0.49
2:B:374:ASP:OD2	2:B:376:SER:OG	2.30	0.49
2:B:473:PHE:HZ	2:B:535:GLN:HE21	1.60	0.49
2:C:828:LEU:HD11	2:C:1071:GLN:HG3	1.95	0.49
1:T:431:GLN:HG3	2:C:529:SER:HA	1.95	0.49
2:A:162:VAL:HG22	2:A:172:GLU:HB2	1.94	0.49
2:C:1128:ASN:HB3	2:C:1148:TYR:HB3	1.94	0.49
2:B:867:LEU:HD23	2:B:871:LEU:HD11	1.94	0.49
2:B:1023:PRO:HD2	2:B:1026:LEU:HB3	1.93	0.49
2:A:351:ASN:ND2	2:A:601:THR:OG1	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:412:SER:HB3	1:T:415:VAL:HG12	1.95	0.48
2:C:1112:GLU:OE1	2:C:1120:ARG:NH1	2.46	0.48
2:B:791:PRO:HB3	2:B:1150:PRO:HB3	1.95	0.48
1:T:463:SER:OG	2:C:507:ASP:OD2	2.28	0.48
2:B:364:HIS:ND1	2:B:428:SER:OG	2.46	0.48
2:A:131:VAL:HG22	2:A:132:ASN:H	1.78	0.48
2:C:529:SER:HG	2:C:532:THR:HG1	1.61	0.48
1:G:272:SER:HA	1:G:282:GLY:HA2	1.95	0.48
1:T:291:ILE:N	1:T:349:MET:O	2.47	0.48
2:A:821:ASN:HD21	2:B:630:GLN:HE22	1.62	0.48
2:B:109:THR:HG23	2:B:252:GLU:HG2	1.96	0.47
1:G:268:PRO:HB2	1:G:362:LYS:H	1.79	0.47
2:A:400:ASP:O	2:A:410:GLN:NE2	2.47	0.47
2:A:109:THR:OG1	2:A:120:GLU:O	2.32	0.47
2:A:131:VAL:HG13	2:A:133:THR:H	1.78	0.47
2:A:846:VAL:HG13	2:A:1096:ILE:HD12	1.96	0.47
2:A:1023:PRO:HD2	2:A:1026:LEU:HB3	1.97	0.47
2:B:51:VAL:HG11	2:B:54:ARG:HD2	1.97	0.47
2:C:807:SER:HB2	2:C:1099:ILE:HD13	1.96	0.47
1:H:298:VAL:HG11	1:H:331:VAL:HG11	1.96	0.47
2:B:214:ALA:HB2	2:B:220:THR:HA	1.97	0.47
2:C:294:THR:HG21	2:C:299:PRO:HB3	1.95	0.46
2:C:812:ILE:O	2:C:840:ASN:ND2	2.47	0.46
2:B:582:SER:OG	2:B:587:ASN:ND2	2.35	0.46
2:A:507:ASP:OD2	1:G:463:SER:OG	2.34	0.46
1:H:291:ILE:N	1:H:349:MET:O	2.48	0.46
2:A:473:PHE:HZ	2:A:535:GLN:HE21	1.63	0.46
2:C:271:ASP:OD1	2:C:275:VAL:N	2.39	0.46
2:A:476:CYS:O	2:A:502:ARG:N	2.48	0.46
2:C:648:LEU:HB3	2:C:657:GLY:HA3	1.97	0.46
1:H:458:ASP:O	1:H:475:GLY:N	2.36	0.46
2:C:110:LYS:HD2	2:C:117:LEU:HD11	1.98	0.46
2:A:432:VAL:HG22	2:A:433:ASN:HD22	1.81	0.46
2:A:737:ARG:HD3	2:C:949:ILE:HD13	1.98	0.46
2:B:787:GLU:HB3	2:B:1157:LEU:HD23	1.97	0.46
1:T:416:TYR:OH	1:T:471:PRO:O	2.33	0.46
2:C:1052:GLY:O	2:C:1076:ARG:NE	2.48	0.46
2:B:131:VAL:HG22	2:B:132:ASN:H	1.81	0.46
2:B:74:PHE:HB3	2:B:257:TRP:HB3	1.99	0.45
2:A:332:TRP:O	2:A:390:LYS:NZ	2.40	0.45
1:G:425:ILE:HG23	1:G:476:ASN:HB2	1.96	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:839:ILE:HG23	2:B:1089:VAL:HG11	1.98	0.45
1:H:446:VAL:HB	1:H:453:TRP:HB3	1.98	0.45
2:A:636:VAL:HG21	2:A:668:THR:HG23	1.99	0.45
2:A:805:THR:HG22	2:A:1103:ALA:HA	1.99	0.45
2:B:871:LEU:HD12	2:B:871:LEU:HA	1.83	0.45
2:A:427:TYR:HE1	2:A:590:SER:HB2	1.82	0.45
2:C:864:GLY:O	2:C:972:SER:OG	2.34	0.45
2:B:476:CYS:O	2:B:502:ARG:N	2.48	0.45
2:A:343:ASN:ND2	2:A:439:PHE:O	2.42	0.45
2:C:39:VAL:HB	2:C:71:GLY:HA2	1.99	0.45
2:B:262:SER:OG	2:B:264:ARG:NH1	2.49	0.45
2:A:340:SER:OG	2:A:343:ASN:OD1	2.34	0.44
2:C:51:VAL:HG11	2:C:54:ARG:HD2	1.99	0.44
2:C:690:LEU:HB2	2:C:721:LEU:HB2	1.99	0.44
1:G:291:ILE:N	1:G:349:MET:O	2.50	0.44
2:A:168:SER:HA	2:A:243:ASN:HD21	1.81	0.44
1:T:415:VAL:O	2:C:517:ARG:NH2	2.40	0.44
2:A:1128:ASN:HB3	2:A:1148:TYR:HB3	1.99	0.44
1:T:257:VAL:HG11	1:T:466:ALA:HB2	1.99	0.44
1:H:435:ASP:HB3	1:H:466:ALA:H	1.83	0.44
1:T:413:ARG:HA	1:T:417:ASP:HA	2.00	0.44
1:H:272:SER:HA	1:H:282:GLY:HA2	1.99	0.44
2:A:51:VAL:HG11	2:A:54:ARG:HD2	2.00	0.44
2:A:793:ASN:HB3	2:A:1151:THR:HB	2.00	0.44
2:C:250:ASP:OD1	2:C:250:ASP:N	2.51	0.44
2:C:473:PHE:HZ	2:C:535:GLN:HE21	1.65	0.44
2:C:791:PRO:HB3	2:C:1150:PRO:HB3	2.00	0.44
2:B:281:ASP:OD1	2:B:282:CYS:N	2.42	0.44
2:B:405:SER:H	2:B:410:GLN:HE22	1.66	0.44
2:A:655:ILE:HG12	2:A:669:ILE:HD12	2.00	0.44
2:B:793:ASN:OD1	2:B:794:PHE:N	2.51	0.44
2:C:356:LEU:HD11	2:C:378:ILE:HG22	2.00	0.43
2:B:372:ASN:HB2	2:B:421:SER:HA	2.00	0.43
2:B:389:ASP:OD1	2:B:389:ASP:N	2.51	0.43
2:C:340:SER:OG	2:C:343:ASN:OD1	2.35	0.43
2:B:218:MET:HB2	2:B:219:PRO:HD3	1.99	0.43
2:B:605:ASN:HB2	2:B:608:LEU:HD11	1.99	0.43
2:B:611:ASN:OD1	2:B:612:THR:N	2.51	0.43
1:G:378:LEU:HD11	1:G:401:LYS:HD2	1.99	0.43
2:B:405:SER:O	2:B:410:GLN:NE2	2.50	0.43
2:A:1108:GLU:OE2	2:B:1105:ARG:NH1	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:269:ASN:HB2	2:B:297:PHE:CE1	2.53	0.43
2:A:52:LEU:HD12	2:A:52:LEU:H	1.84	0.43
2:C:446:ARG:HA	2:C:450:PHE:HB3	2.00	0.43
2:C:650:ASP:OD1	2:C:654:ASN:N	2.51	0.43
2:A:80:LYS:NZ	2:A:242:CYS:H	2.16	0.43
2:A:529:SER:HA	1:G:431:GLN:HG3	2.00	0.43
2:C:1205:MET:HE3	2:B:994:MET:HB3	2.01	0.43
2:C:582:SER:OG	2:C:587:ASN:ND2	2.42	0.42
2:B:907:LEU:HD11	2:B:1142:LEU:HD22	2.00	0.42
2:C:872:ASN:HB2	2:C:894:GLN:HE22	1.84	0.42
2:C:1163:CYS:HA	2:C:1169:GLY:HA2	2.01	0.42
2:B:108:ASN:HB2	2:B:119:SER:HB2	2.00	0.42
2:B:785:LEU:HD22	2:B:1159:SER:HB2	2.00	0.42
2:B:1126:ASN:OD1	2:B:1126:ASN:N	2.52	0.42
2:A:413:ASN:HD21	2:A:463:SER:H	1.67	0.42
2:C:809:LYS:HE3	2:C:847:ASN:HB3	2.00	0.42
1:T:443:GLY:HA2	1:T:444:PRO:HD3	1.89	0.42
2:B:51:VAL:HG21	2:B:56:TYR:CE1	2.55	0.42
2:A:938:GLN:HG2	2:A:1044:LEU:HD22	2.02	0.42
2:A:1200:LYS:NZ	2:B:1204:PHE:HB2	2.35	0.42
2:C:32:ILE:HD12	2:C:33:PRO:HD2	2.01	0.42
1:G:406:GLU:HG3	1:G:409:ARG:H	1.85	0.42
1:T:272:SER:HA	1:T:282:GLY:HA2	2.02	0.42
2:C:139:VAL:HG22	2:C:148:ILE:HG12	2.01	0.42
2:C:291:GLN:HA	2:C:294:THR:HG22	2.02	0.42
2:B:301:THR:HG23	2:B:682:TYR:HA	2.02	0.42
2:A:387:THR:HG23	2:A:593:ILE:HB	2.02	0.41
2:C:396:ARG:HB2	2:C:397:ARG:CZ	2.50	0.41
2:C:717:LEU:HD11	2:C:740:SER:HB2	2.02	0.41
2:B:250:ASP:N	2:B:250:ASP:OD1	2.53	0.41
2:B:341:PRO:HA	2:B:391:PHE:HB2	2.01	0.41
1:G:435:ASP:HB3	1:G:466:ALA:H	1.84	0.41
2:B:98:PHE:N	2:B:233:SER:O	2.42	0.41
2:B:42:VAL:HG11	2:B:266:TYR:HE2	1.85	0.41
2:C:1209:SER:HB2	2:B:1001:GLN:HE22	1.86	0.41
2:B:370:CYS:HA	2:B:423:CYS:HA	2.02	0.41
2:B:624:LEU:HD13	2:B:658:PHE:CD1	2.55	0.41
2:A:791:PRO:HB3	2:A:1150:PRO:HB3	2.02	0.41
2:A:799:HIS:CE1	2:A:1126:ASN:HD21	2.39	0.41
2:A:1052:GLY:O	2:A:1076:ARG:NH1	2.53	0.41
2:C:149:THR:HG22	2:C:186:LYS:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:405:SER:H	2:C:410:GLN:HE22	1.68	0.41
2:C:786:PHE:HD2	2:B:969:PRO:HG3	1.85	0.41
2:B:938:GLN:HG2	2:B:1044:LEU:HD22	2.03	0.41
1:T:334:HIS:HE1	1:T:336:ASN:HB2	1.86	0.41
2:A:680:ALA:HB1	2:A:736:LEU:HD13	2.02	0.41
2:C:335:ASN:OD1	2:C:336:VAL:N	2.54	0.41
2:B:960:THR:O	2:B:964:VAL:HG23	2.20	0.41
1:T:415:VAL:HG22	1:T:416:TYR:CD2	2.55	0.41
2:A:212:TYR:HB3	2:A:220:THR:HB	2.02	0.41
2:B:34:ARG:HB3	2:B:76:ASP:HB3	2.02	0.41
2:B:413:ASN:HD21	2:B:463:SER:H	1.69	0.41
2:A:113:VAL:O	2:A:116:THR:OG1	2.23	0.41
2:C:83:ILE:HG22	2:C:169:ILE:HG21	2.02	0.41
2:C:638:ALA:HB3	2:C:666:THR:HG21	2.02	0.41
2:A:42:VAL:HG21	2:A:266:TYR:HE2	1.86	0.40
2:B:805:THR:HG21	2:B:857:VAL:HG21	2.03	0.40
1:H:275:VAL:HG22	1:H:276:GLN:HG2	2.02	0.40
2:A:700:VAL:HA	2:A:704:ILE:HG22	2.03	0.40
1:T:275:VAL:HG22	1:T:276:GLN:HG2	2.02	0.40
2:A:106:VAL:HB	2:A:198:LEU:HB2	2.04	0.40
2:C:269:ASN:HB2	2:C:297:PHE:CE1	2.56	0.40
2:C:427:TYR:HE1	2:C:590:SER:HB2	1.86	0.40
2:B:747:ALA:HB2	2:B:763:ARG:NH1	2.37	0.40
1:G:305:PRO:HB3	1:G:331:VAL:HG23	2.04	0.40
2:A:1067:PRO:N	2:A:1068:PRO:HD2	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	G	235/383 (61%)	217 (92%)	18 (8%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	235/383 (61%)	215 (92%)	20 (8%)	0	100	100
1	T	235/383 (61%)	215 (92%)	20 (8%)	0	100	100
2	A	1175/1263 (93%)	1147 (98%)	28 (2%)	0	100	100
2	B	1188/1263 (94%)	1152 (97%)	36 (3%)	0	100	100
2	C	1188/1263 (94%)	1160 (98%)	28 (2%)	0	100	100
All	All	4256/4938 (86%)	4106 (96%)	150 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	171/325 (53%)	170 (99%)	1 (1%)	84	92
1	H	171/325 (53%)	169 (99%)	2 (1%)	67	85
1	T	170/325 (52%)	168 (99%)	2 (1%)	67	85
2	A	1039/1132 (92%)	1028 (99%)	11 (1%)	70	86
2	B	1047/1132 (92%)	1033 (99%)	14 (1%)	65	83
2	C	1048/1132 (93%)	1036 (99%)	12 (1%)	70	86
All	All	3646/4371 (83%)	3604 (99%)	42 (1%)	66	85

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	T	355	LEU
1	T	387	THR
1	H	355	LEU
1	H	387	THR
2	A	20	CYS
2	A	64	THR
2	A	163	CYS
2	A	213	TYR

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Mol	Chain	Res	Type
2	A	242	CYS
2	A	474	CYS
2	A	476	CYS
2	A	682	TYR
2	A	734	CYS
2	A	836	CYS
2	A	1022	THR
2	C	163	CYS
2	C	213	TYR
2	C	216	VAL
2	C	220	THR
2	C	292	CYS
2	C	312	LYS
2	C	476	CYS
2	C	619	CYS
2	C	663	THR
2	C	743	CYS
2	C	771	ASN
2	C	1208	CYS
2	B	51	VAL
2	B	113	VAL
2	B	213	TYR
2	B	387	THR
2	B	398	ARG
2	B	453	PHE
2	B	476	CYS
2	B	603	CYS
2	B	635	GLU
2	B	734	CYS
2	B	814	CYS
2	B	819	CYS
2	B	887	LEU
2	B	1022	THR
1	G	387	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (56) such sidechains are listed below:

Mol	Chain	Res	Type
1	T	276	GLN
1	H	276	GLN
1	H	408	GLN
2	A	142	HIS

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Mol	Chain	Res	Type
2	A	201	HIS
2	A	243	ASN
2	A	273	HIS
2	A	291	GLN
2	A	335	ASN
2	A	351	ASN
2	A	433	ASN
2	A	587	ASN
2	A	605	ASN
2	A	709	GLN
2	A	799	HIS
2	A	847	ASN
2	A	863	GLN
2	A	1046	GLN
2	A	1059	GLN
2	A	1071	GLN
2	C	108	ASN
2	C	243	ASN
2	C	410	GLN
2	C	587	ASN
2	C	645	GLN
2	C	652	ASN
2	C	683	GLN
2	C	722	ASN
2	C	840	ASN
2	C	844	ASN
2	C	854	GLN
2	C	881	ASN
2	C	988	ASN
2	C	1038	GLN
2	C	1045	GLN
2	C	1059	GLN
2	C	1091	GLN
2	C	1211	ASN
2	B	204	GLN
2	B	265	GLN
2	B	269	ASN
2	B	291	GLN
2	B	371	ASN
2	B	402	GLN
2	B	410	GLN
2	B	445	ASN

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Mol	Chain	Res	Type
2	B	535	GLN
2	B	587	ASN
2	B	630	GLN
2	B	789	GLN
2	B	941	ASN
2	B	1001	GLN
2	B	1091	GLN
2	B	1128	ASN
2	B	1135	GLN
1	G	276	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

28 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	$\# Z > 2$
3	NAG	D	1	3,2	14,14,15	0.19	0	17,19,21	0.39	0
3	NAG	D	2	3	14,14,15	0.24	0	17,19,21	0.42	0
3	NAG	E	1	3,2	14,14,15	0.22	0	17,19,21	0.45	0
3	NAG	E	2	3	14,14,15	0.23	0	17,19,21	0.41	0
3	NAG	F	1	3,2	14,14,15	0.21	0	17,19,21	0.42	0
3	NAG	F	2	3	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	I	1	4,2	14,14,15	0.24	0	17,19,21	0.44	0
4	NAG	I	2	4	14,14,15	0.21	0	17,19,21	0.39	0
4	BMA	I	3	4	11,11,12	0.59	0	15,15,17	0.79	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	J	1	3,2	14,14,15	0.23	0	17,19,21	0.46	0
3	NAG	J	2	3	14,14,15	0.23	0	17,19,21	0.41	0
4	NAG	K	1	4,2	14,14,15	0.24	0	17,19,21	0.46	0
4	NAG	K	2	4	14,14,15	0.22	0	17,19,21	0.39	0
4	BMA	K	3	4	11,11,12	0.58	0	15,15,17	0.77	0
3	NAG	L	1	3,2	14,14,15	0.24	0	17,19,21	0.46	0
3	NAG	L	2	3	14,14,15	0.25	0	17,19,21	0.42	0
3	NAG	M	1	3,2	14,14,15	0.21	0	17,19,21	0.43	0
3	NAG	M	2	3	14,14,15	0.24	0	17,19,21	0.43	0
3	NAG	N	1	3,2	14,14,15	0.21	0	17,19,21	0.40	0
3	NAG	N	2	3	14,14,15	0.24	0	17,19,21	0.42	0
3	NAG	O	1	3,2	14,14,15	0.18	0	17,19,21	0.42	0
3	NAG	O	2	3	14,14,15	0.25	0	17,19,21	0.44	0
3	NAG	P	1	3,2	14,14,15	0.23	0	17,19,21	0.42	0
3	NAG	P	2	3	14,14,15	0.23	0	17,19,21	0.42	0
3	NAG	Q	1	3,2	14,14,15	0.22	0	17,19,21	0.45	0
3	NAG	Q	2	3	14,14,15	0.24	0	17,19,21	0.41	0
3	NAG	R	1	3,2	14,14,15	0.22	0	17,19,21	0.43	0
3	NAG	R	2	3	14,14,15	0.23	0	17,19,21	0.41	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	D	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
3	NAG	E	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	NAG	F	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
4	NAG	I	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	1/6/23/26	0/1/1/1
4	BMA	I	3	4	-	0/2/19/22	0/1/1/1
3	NAG	J	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1
4	NAG	K	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	K	2	4	-	0/6/23/26	0/1/1/1
4	BMA	K	3	4	-	0/2/19/22	0/1/1/1
3	NAG	L	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	L	2	3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	M	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	M	2	3	-	0/6/23/26	0/1/1/1
3	NAG	N	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	N	2	3	-	0/6/23/26	0/1/1/1
3	NAG	O	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	O	2	3	-	0/6/23/26	0/1/1/1
3	NAG	P	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	P	2	3	-	0/6/23/26	0/1/1/1
3	NAG	Q	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	Q	2	3	-	2/6/23/26	0/1/1/1
3	NAG	R	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	R	2	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

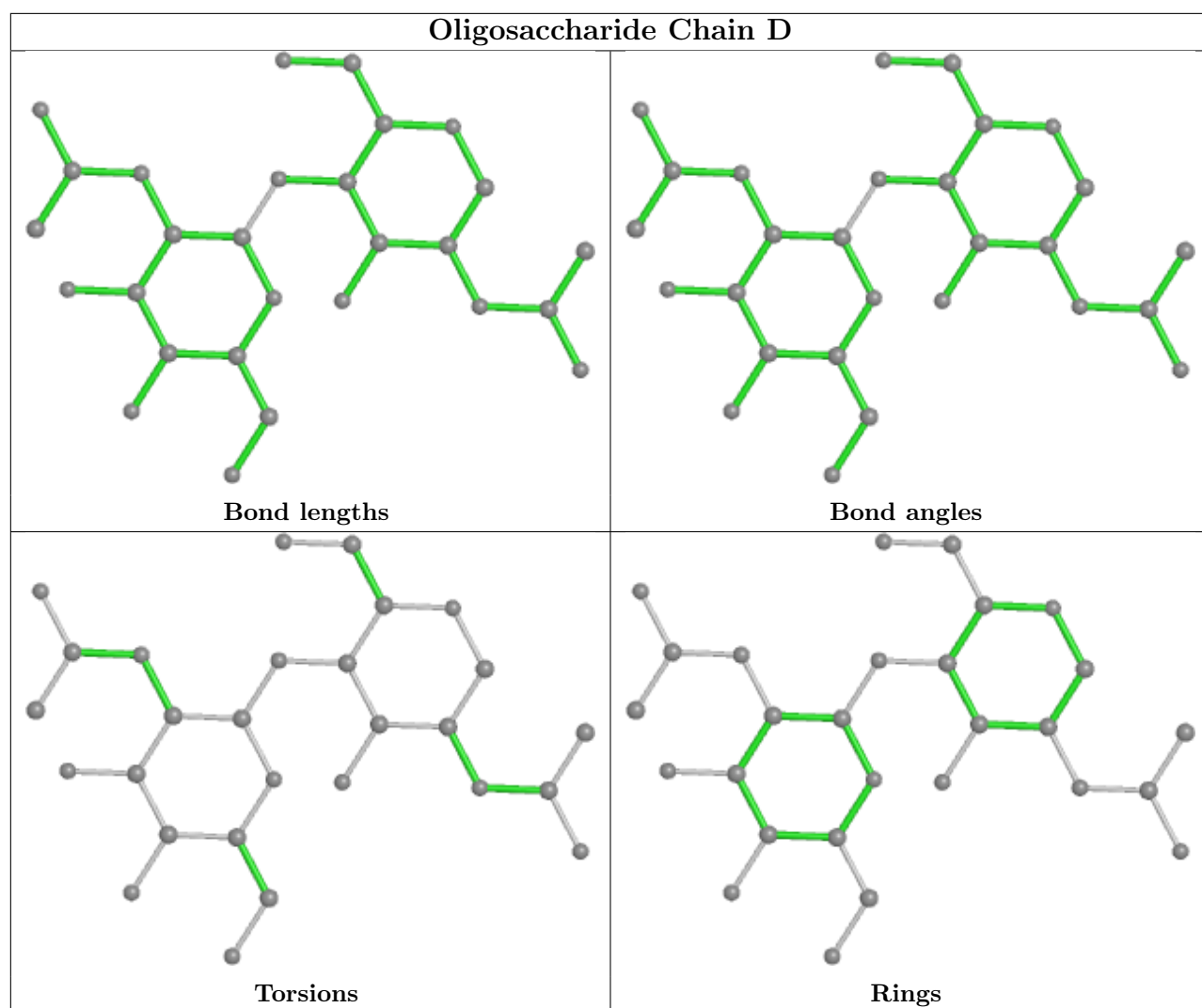
All (13) torsion outliers are listed below:

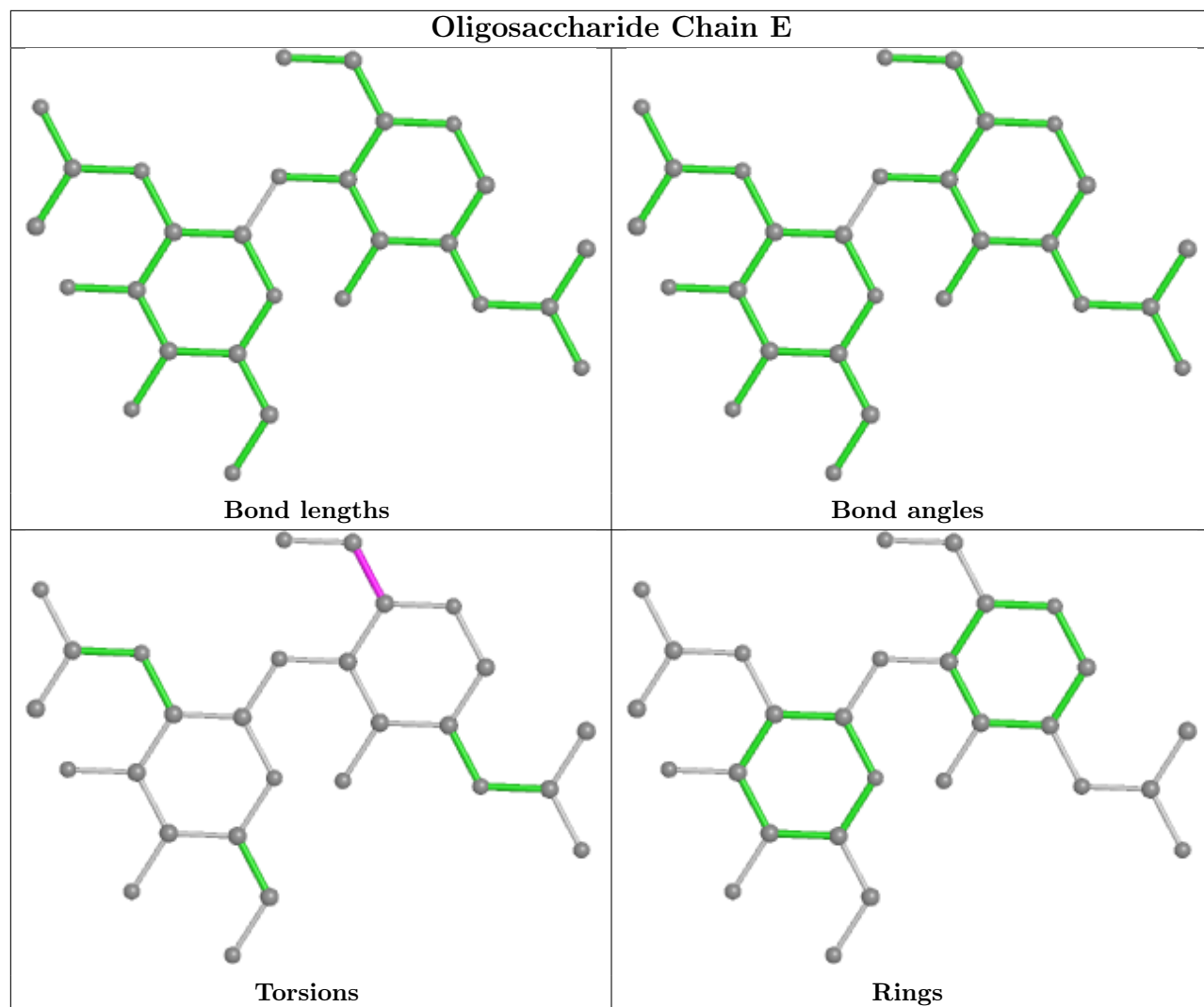
Mol	Chain	Res	Type	Atoms
3	F	2	NAG	O5-C5-C6-O6
3	O	1	NAG	C8-C7-N2-C2
3	O	1	NAG	O7-C7-N2-C2
3	Q	1	NAG	O5-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6
3	Q	1	NAG	C4-C5-C6-O6
3	Q	2	NAG	O5-C5-C6-O6
3	Q	2	NAG	C4-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
3	M	1	NAG	C4-C5-C6-O6
4	I	2	NAG	C4-C5-C6-O6
3	M	1	NAG	O5-C5-C6-O6

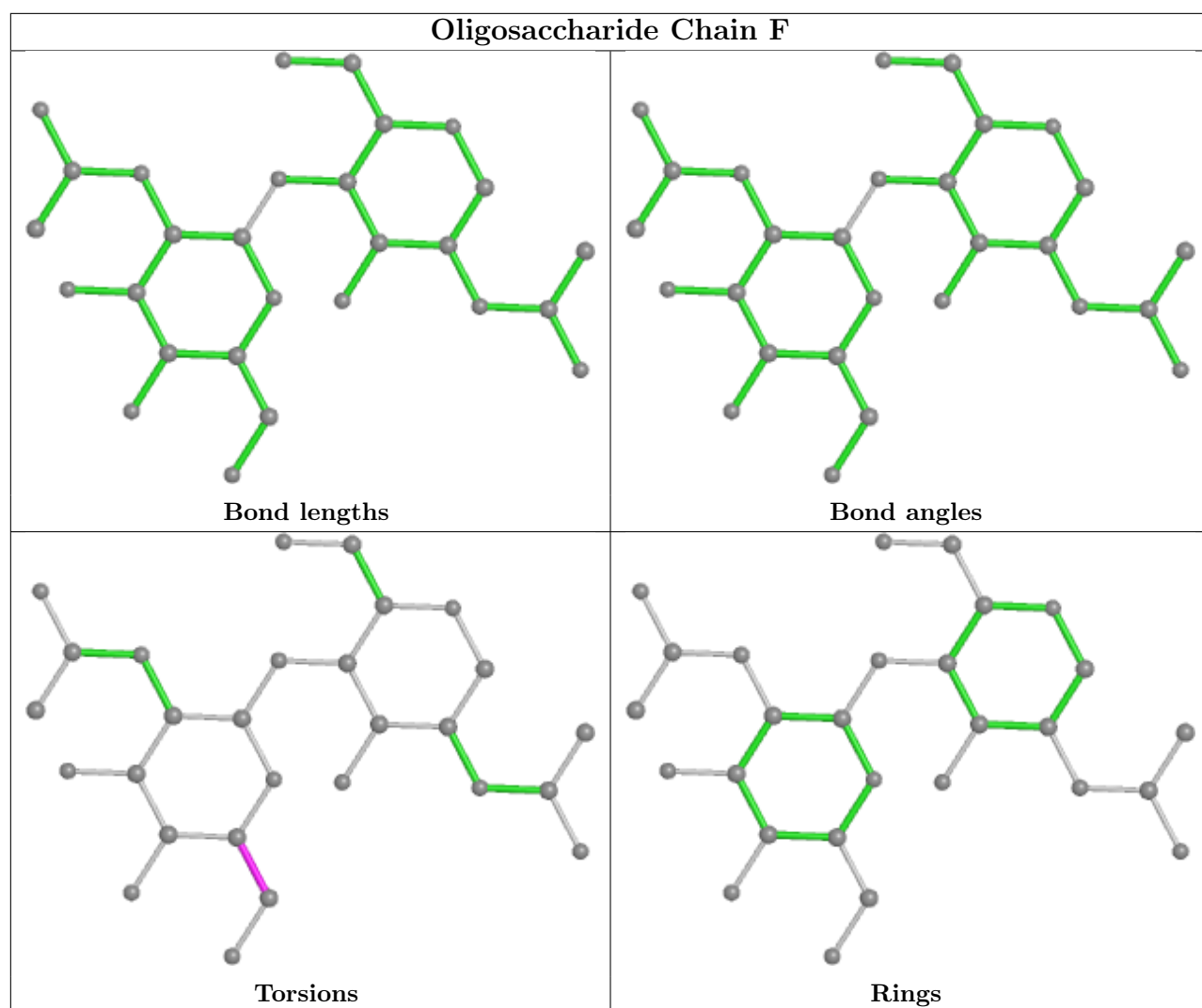
There are no ring outliers.

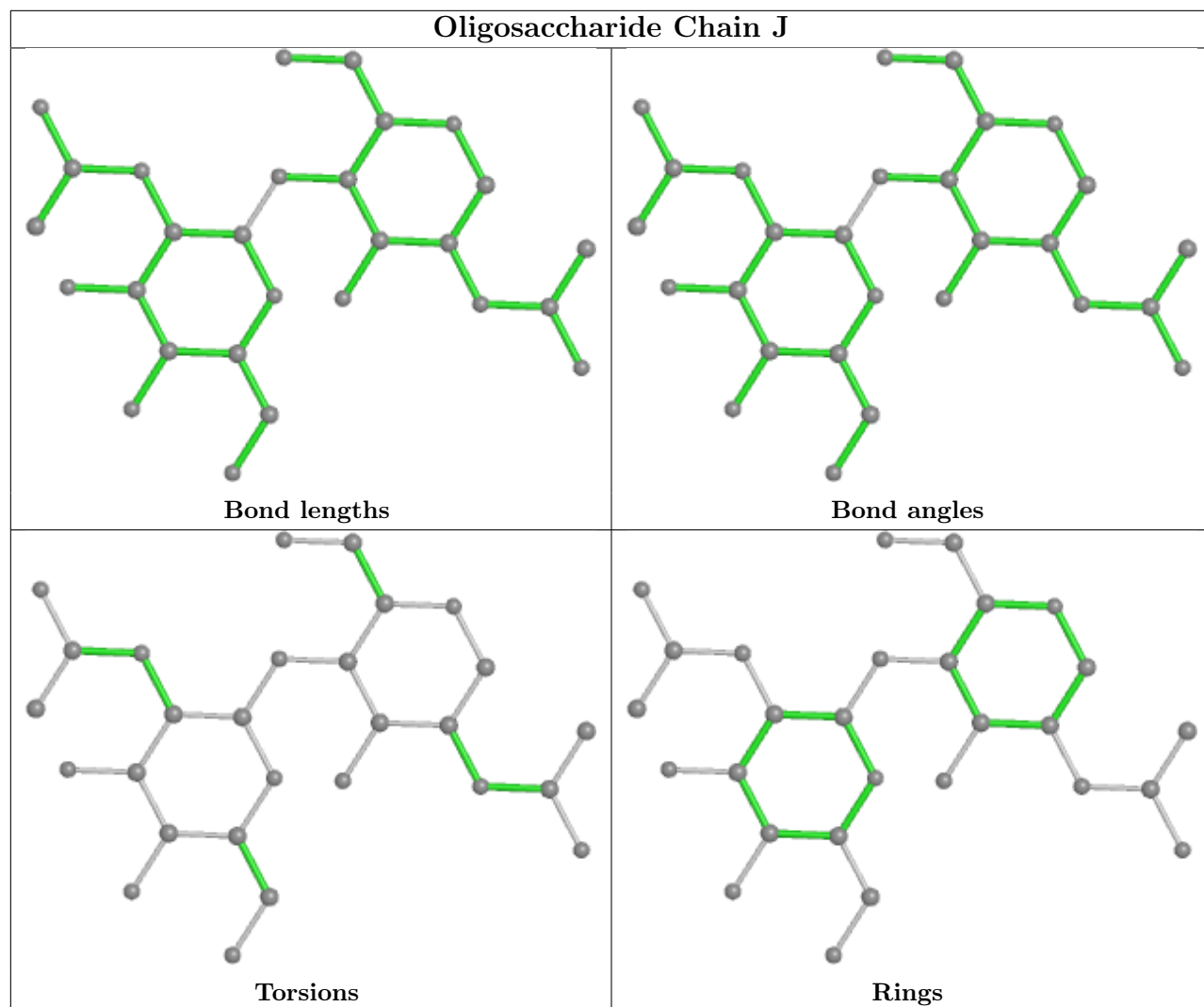
No monomer is involved in short contacts.

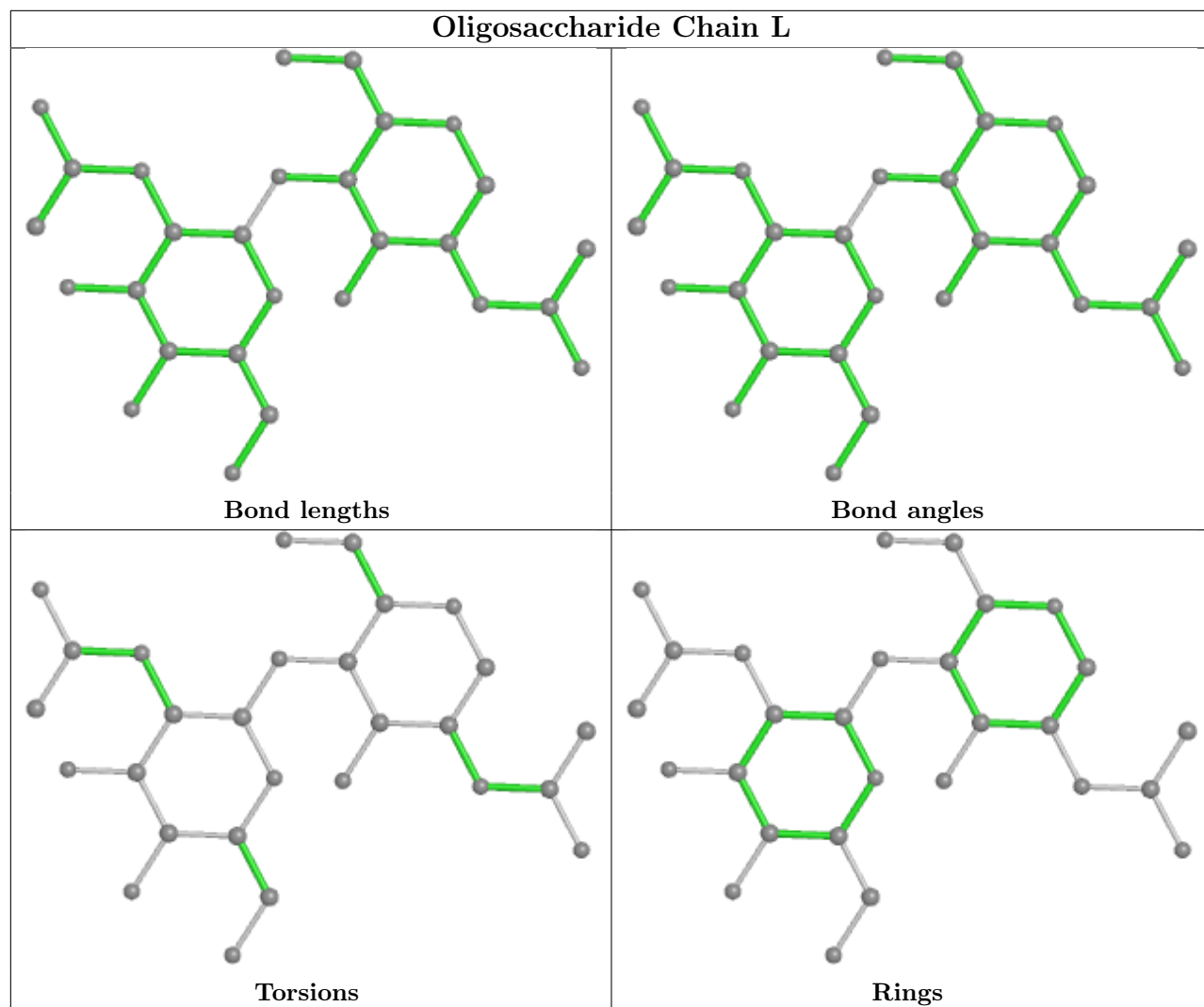
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

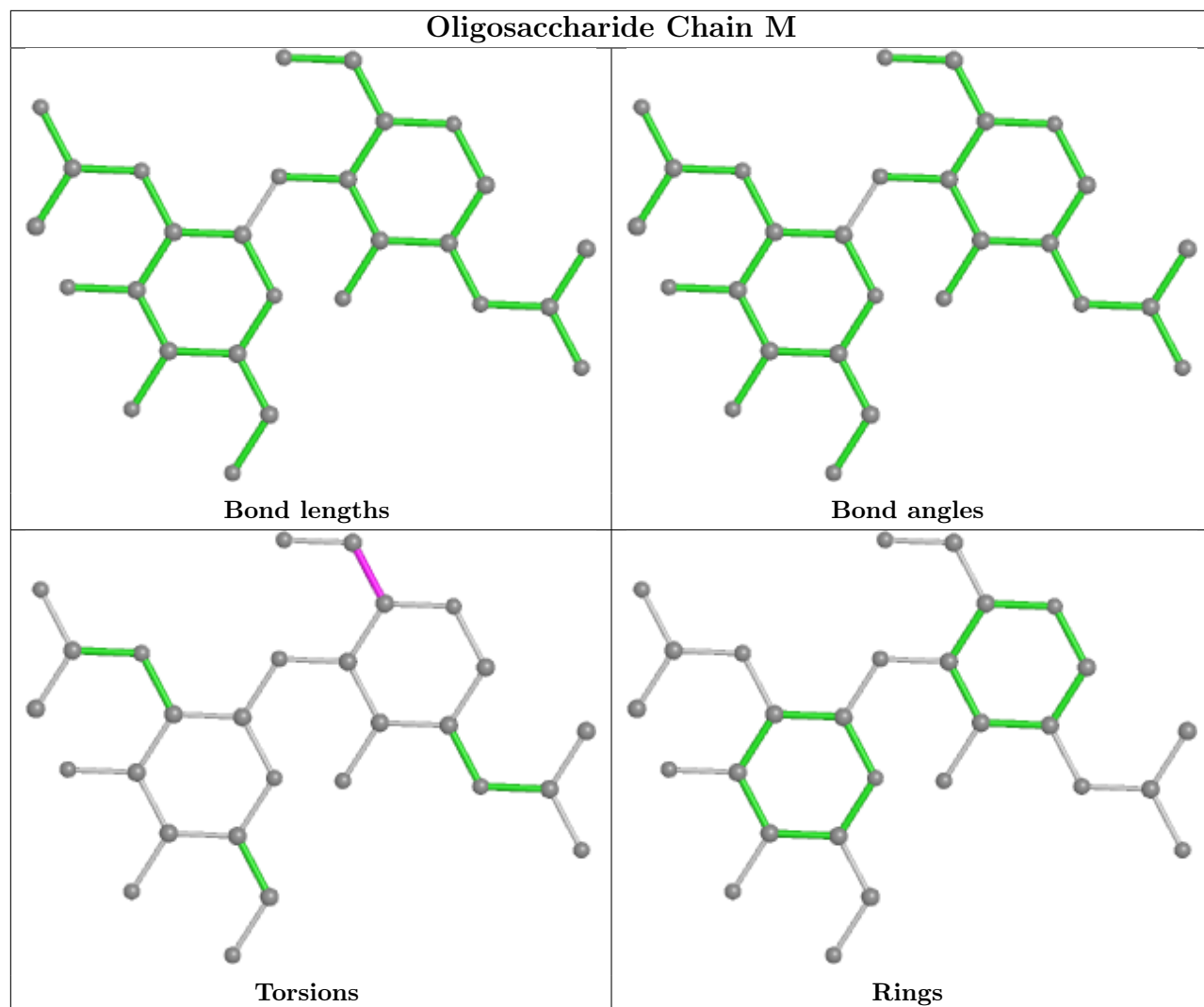


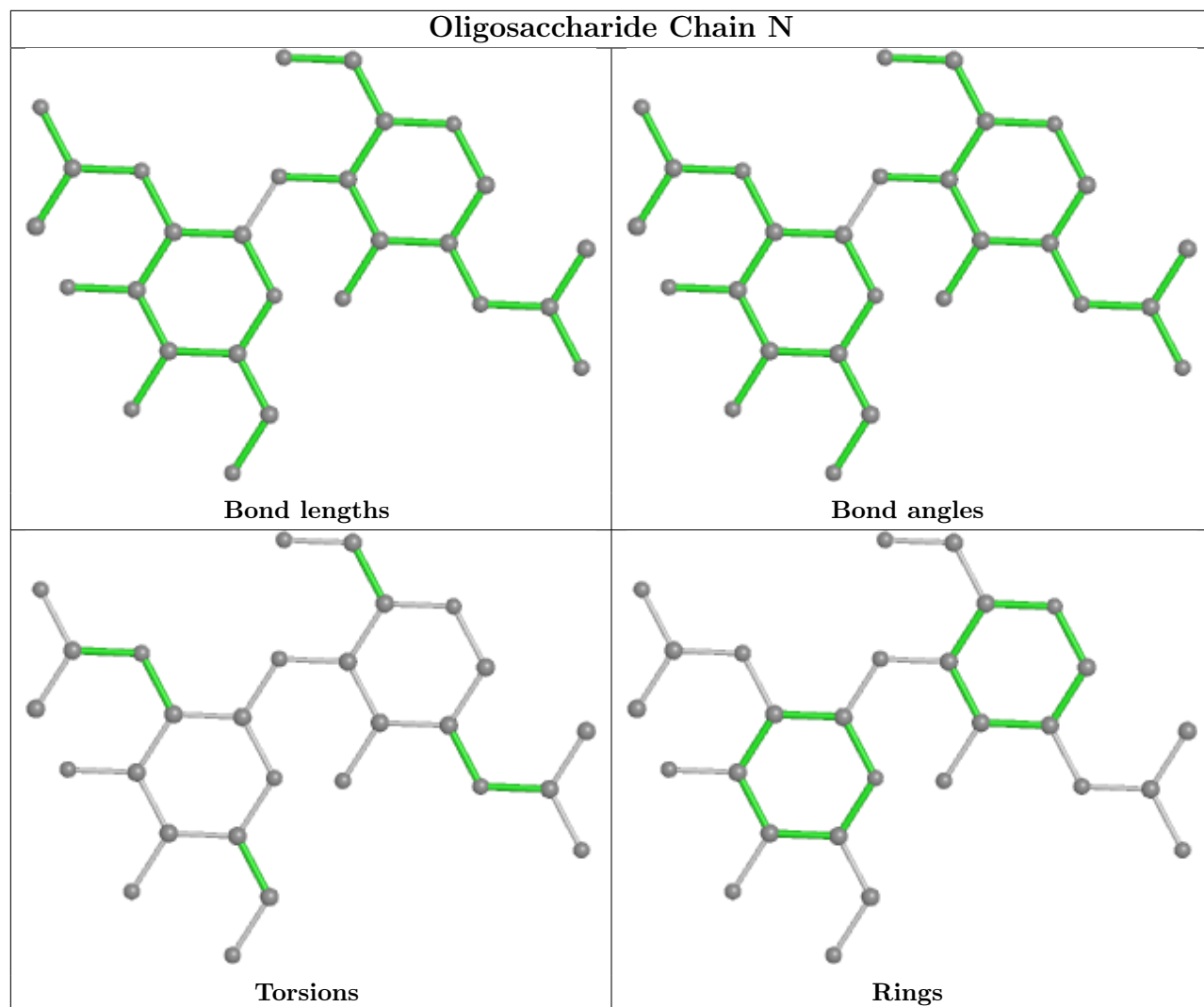


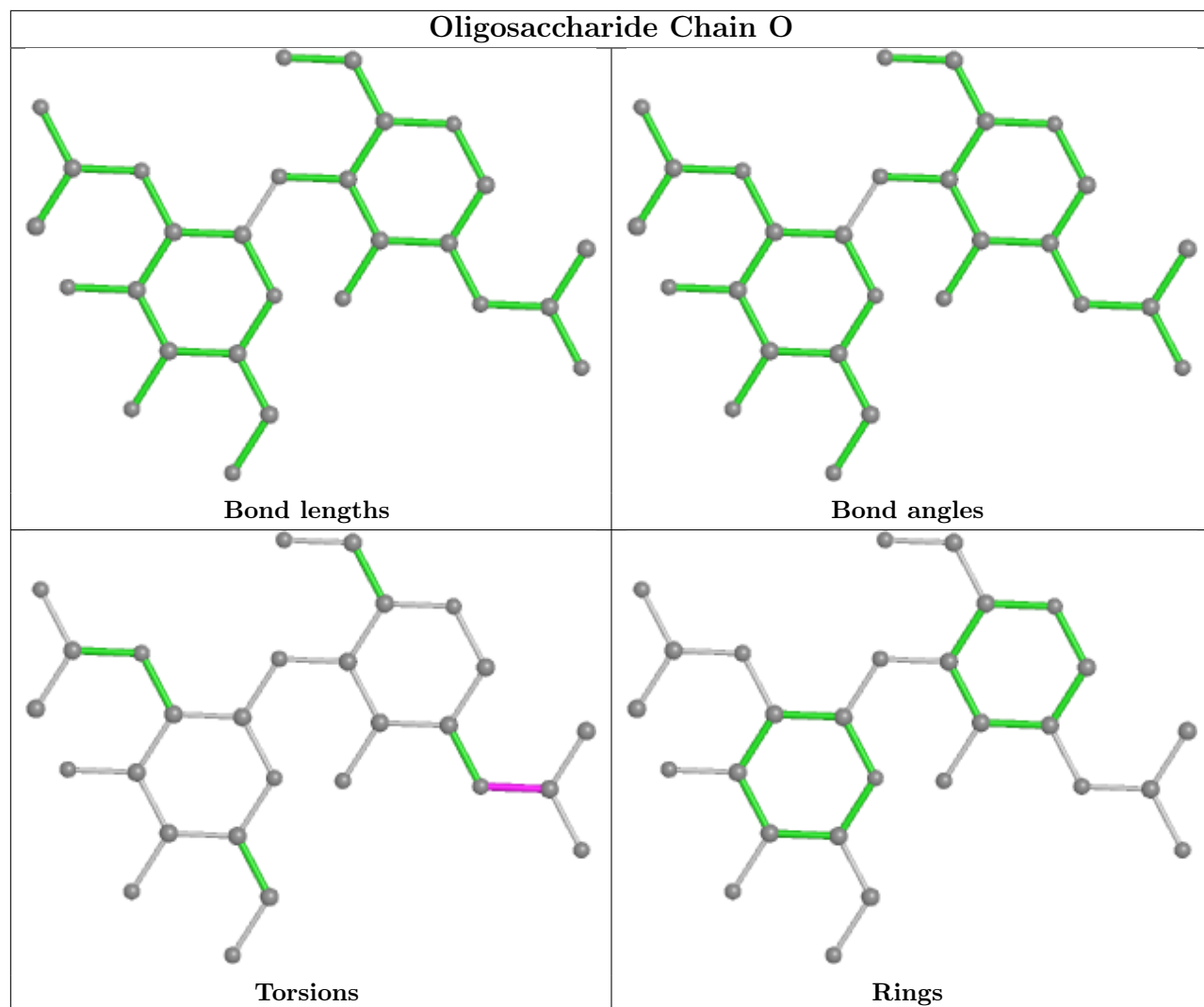


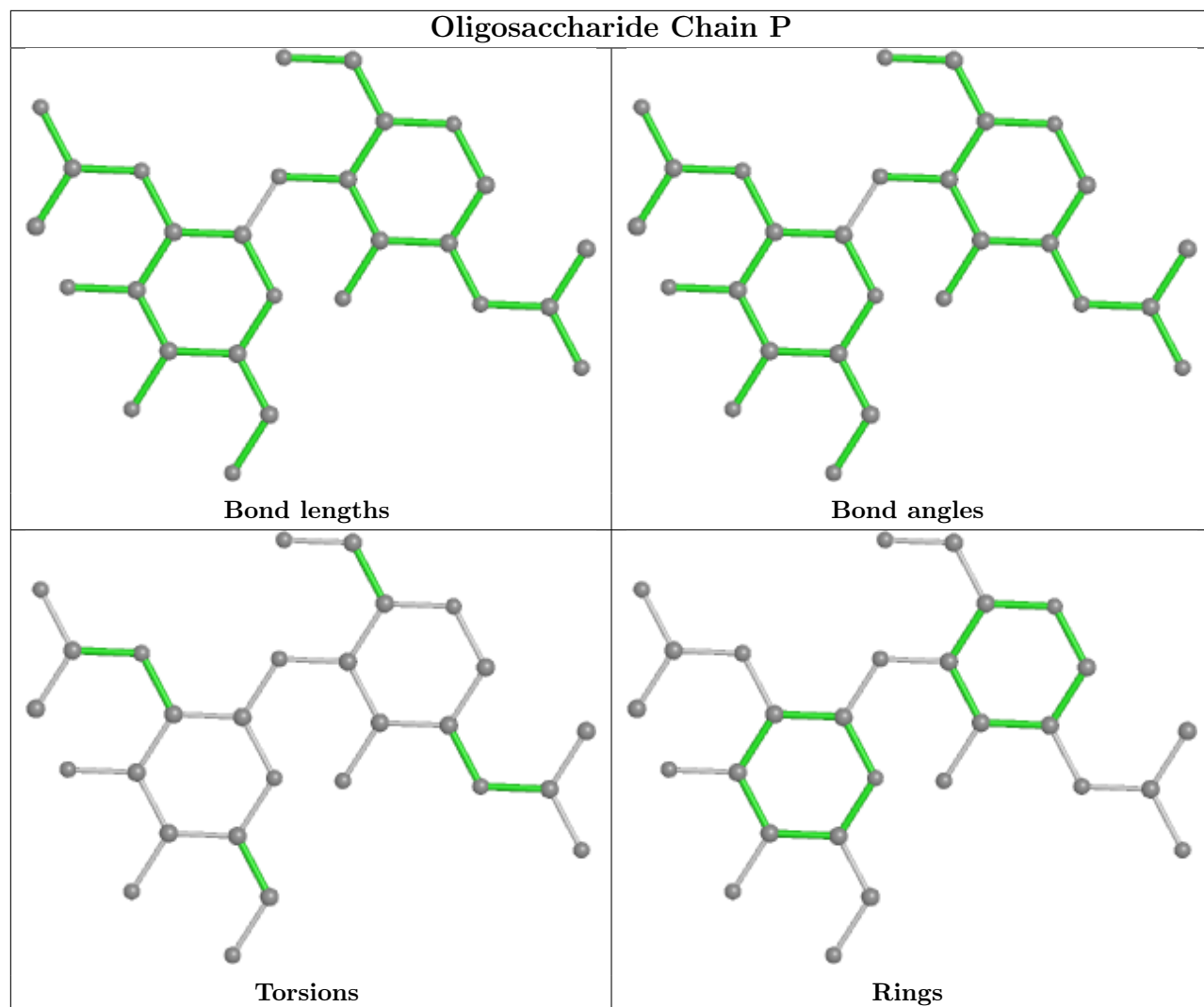


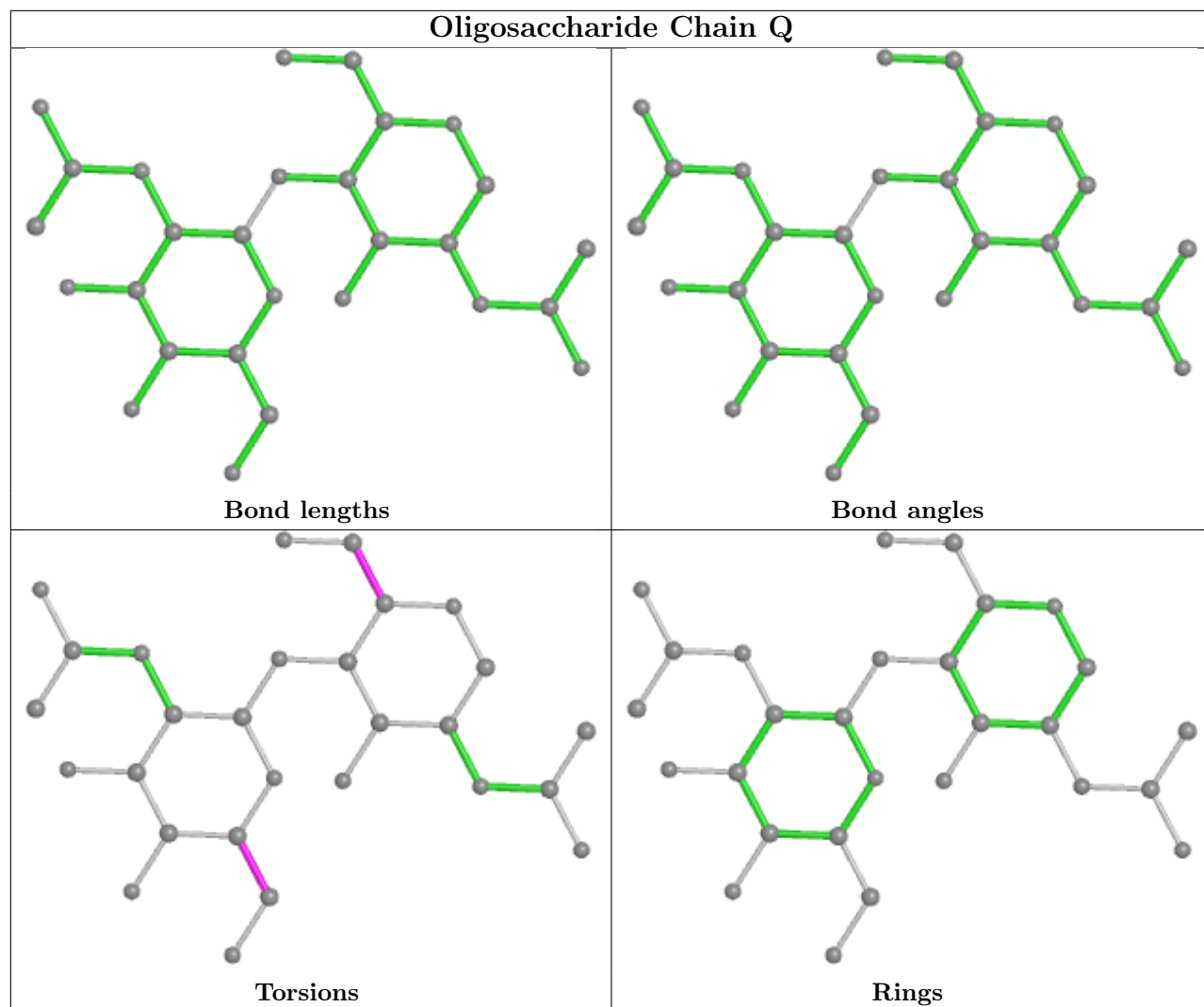


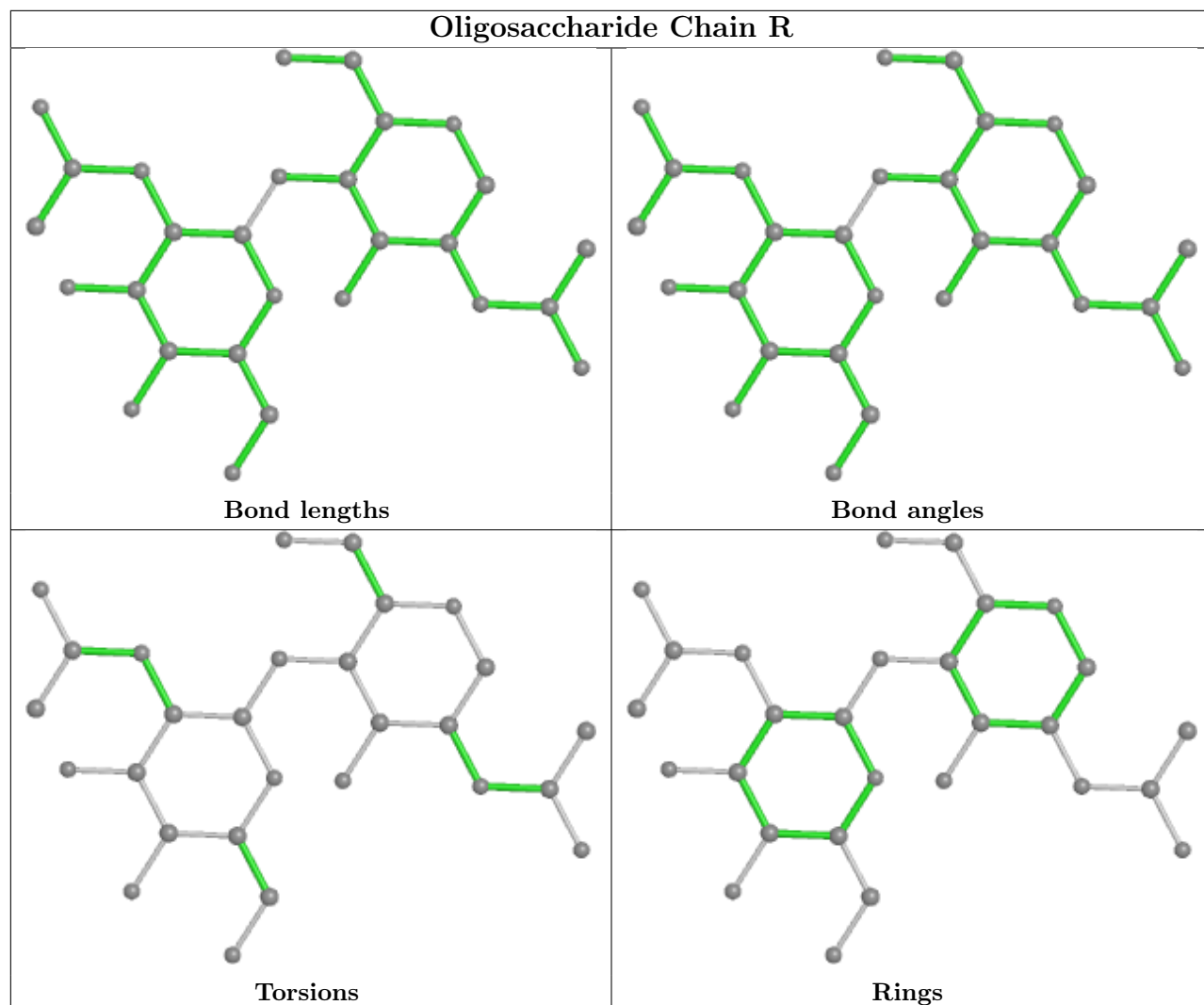


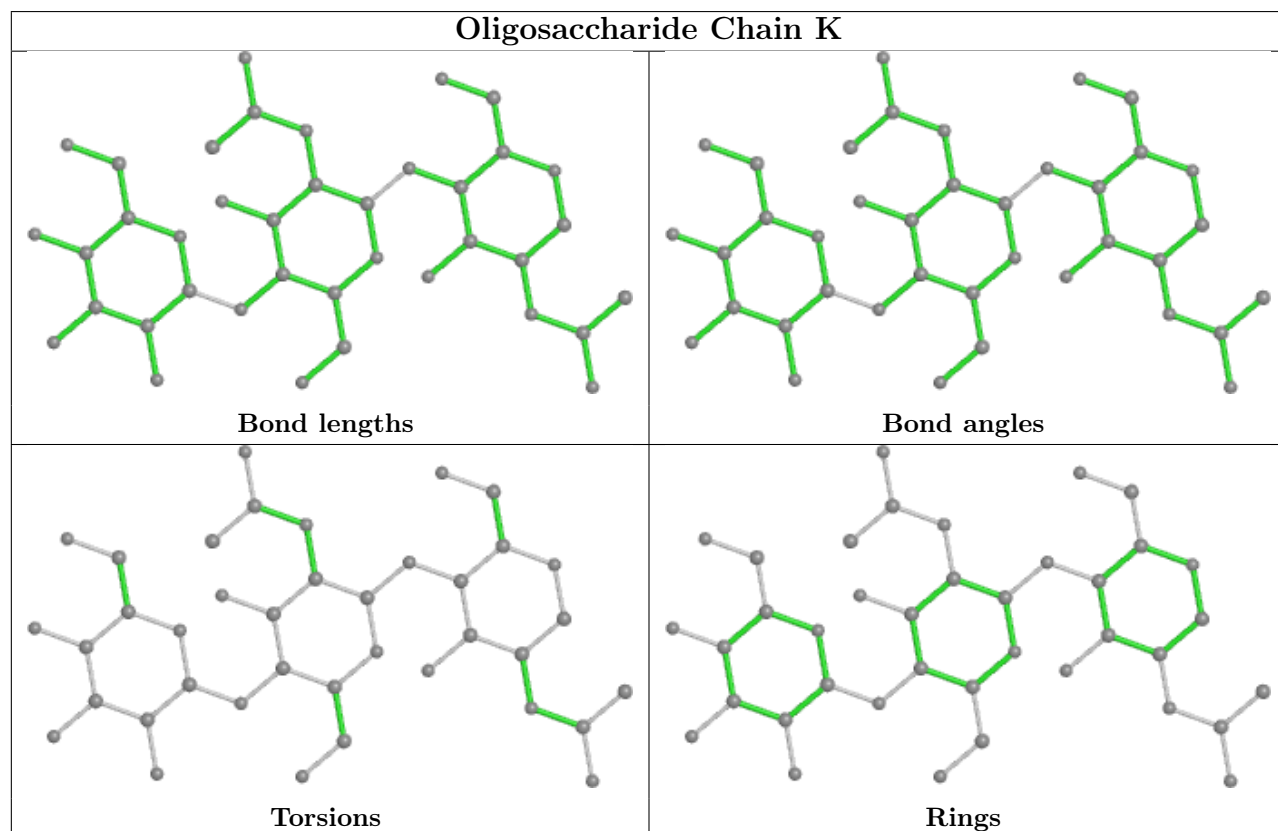
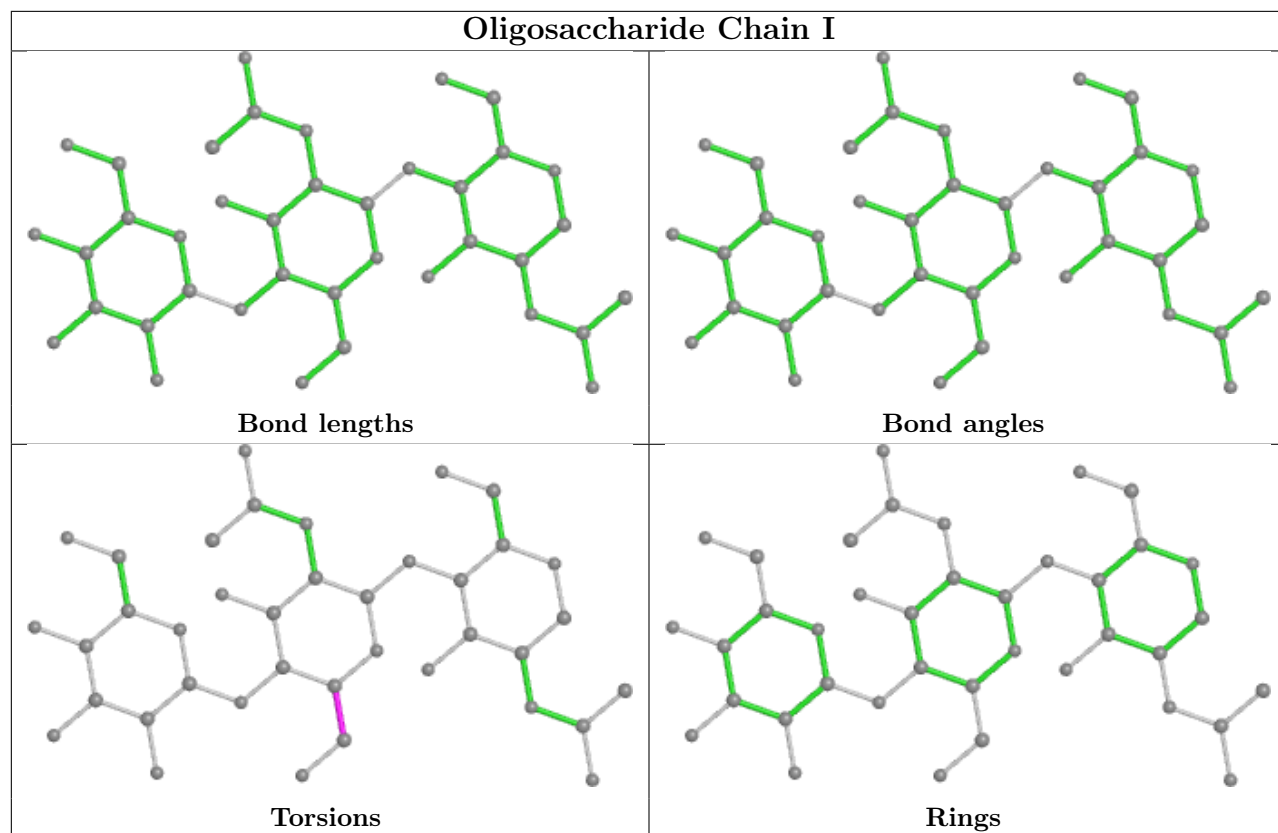












5.6 Ligand geometry

10 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	$\# Z > 2$
5	NAG	C	1302	2	14,14,15	0.23	0	17,19,21	0.42	0
5	NAG	B	1302	2	14,14,15	0.23	0	17,19,21	0.42	0
5	NAG	A	1302	2	14,14,15	0.23	0	17,19,21	0.43	0
5	NAG	B	1301	2	14,14,15	0.21	0	17,19,21	0.42	0
5	NAG	A	1303	2	14,14,15	0.20	0	17,19,21	0.42	0
5	NAG	B	1303	2	14,14,15	0.22	0	17,19,21	0.40	0
5	NAG	A	1304	2	14,14,15	0.22	0	17,19,21	0.42	0
5	NAG	A	1301	2	14,14,15	0.22	0	17,19,21	0.42	0
5	NAG	C	1303	2	14,14,15	0.22	0	17,19,21	0.44	0
5	NAG	C	1301	2	14,14,15	0.22	0	17,19,21	0.42	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
5	NAG	C	1302	2	-	0/6/23/26	0/1/1/1
5	NAG	B	1302	2	-	0/6/23/26	0/1/1/1
5	NAG	A	1302	2	-	2/6/23/26	0/1/1/1
5	NAG	B	1301	2	-	0/6/23/26	0/1/1/1
5	NAG	A	1303	2	-	0/6/23/26	0/1/1/1
5	NAG	B	1303	2	-	2/6/23/26	0/1/1/1
5	NAG	A	1304	2	-	0/6/23/26	0/1/1/1
5	NAG	A	1301	2	-	2/6/23/26	0/1/1/1
5	NAG	C	1303	2	-	2/6/23/26	0/1/1/1
5	NAG	C	1301	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1303	NAG	C4-C5-C6-O6
5	A	1301	NAG	C4-C5-C6-O6
5	B	1303	NAG	O5-C5-C6-O6
5	A	1301	NAG	O5-C5-C6-O6
5	C	1301	NAG	O5-C5-C6-O6
5	C	1301	NAG	C4-C5-C6-O6
5	A	1302	NAG	C4-C5-C6-O6
5	A	1302	NAG	O5-C5-C6-O6
5	C	1303	NAG	C4-C5-C6-O6
5	C	1303	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

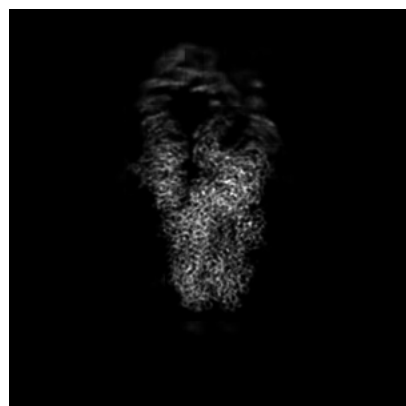
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-39039. These allow visual inspection of the internal detail of the map and identification of artifacts.

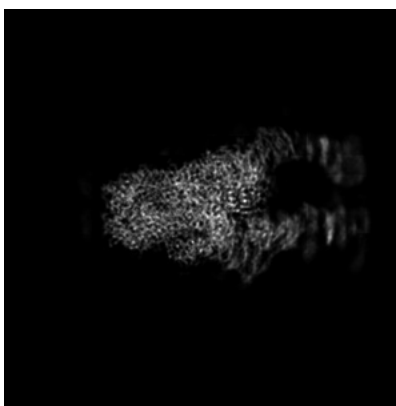
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

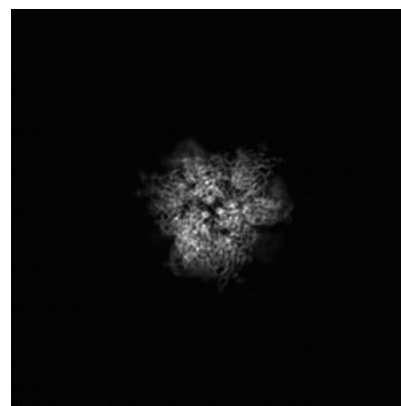
6.1.1 Primary map



X

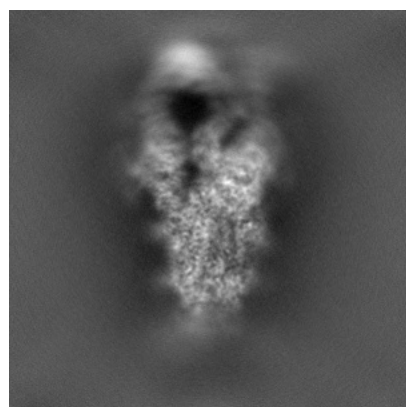


Y

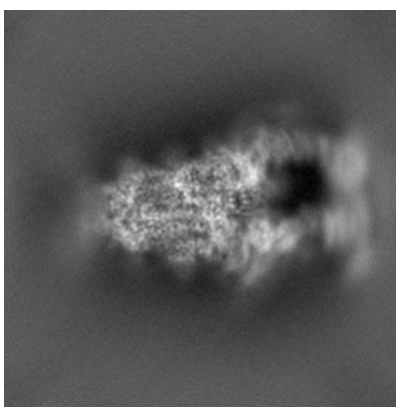


Z

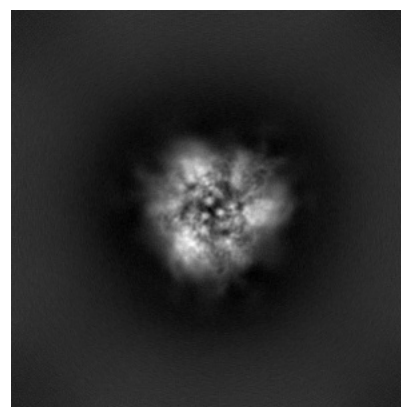
6.1.2 Raw map



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: 200

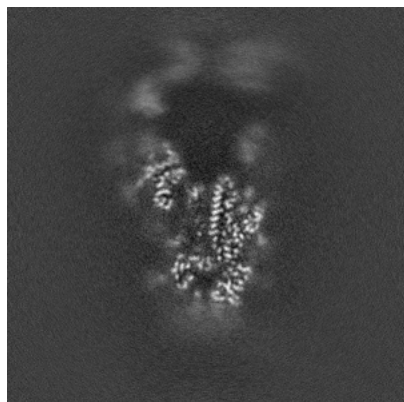


Y Index: 200

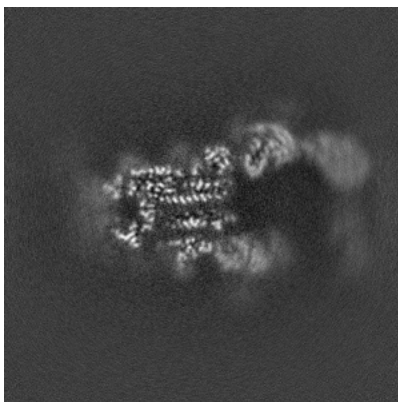


Z Index: 200

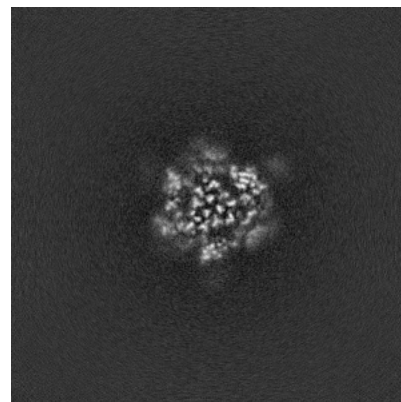
6.2.2 Raw map



X Index: 200



Y Index: 200



Z Index: 200

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: 222

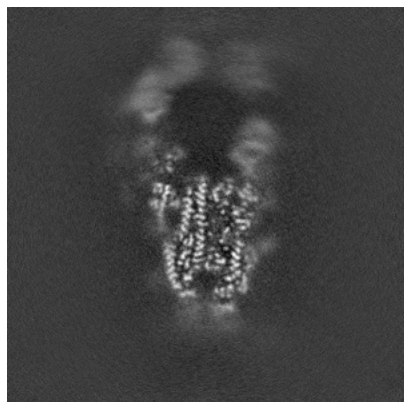


Y Index: 188

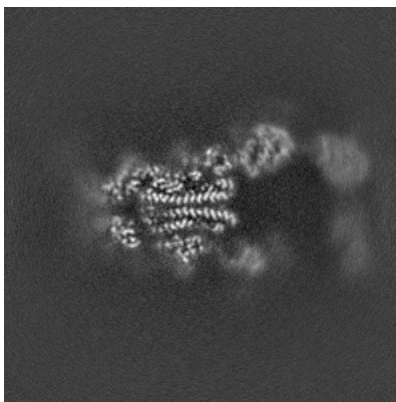


Z Index: 215

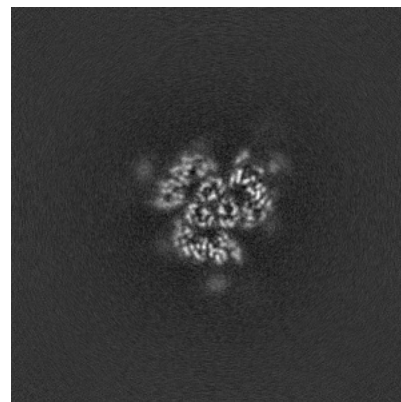
6.3.2 Raw map



X Index: 192



Y Index: 196



Z Index: 214

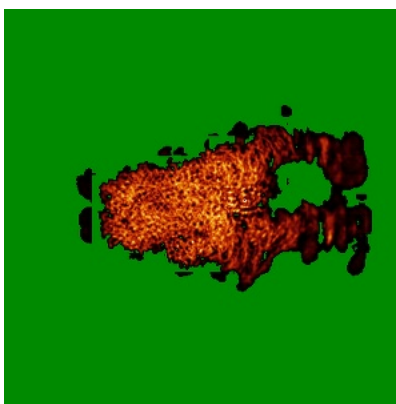
The images above show the largest variance slices of the map in three orthogonal directions.

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

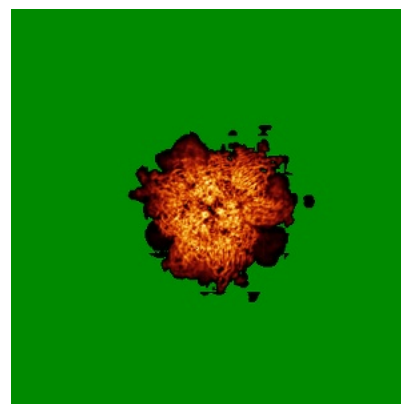
6.4.1 Primary map



X

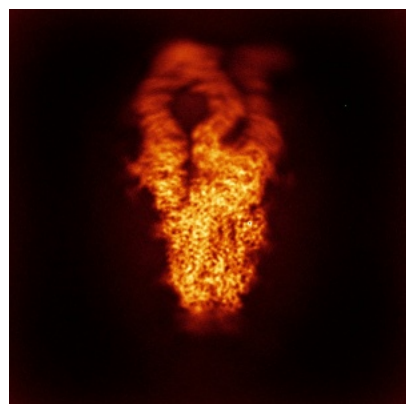


Y

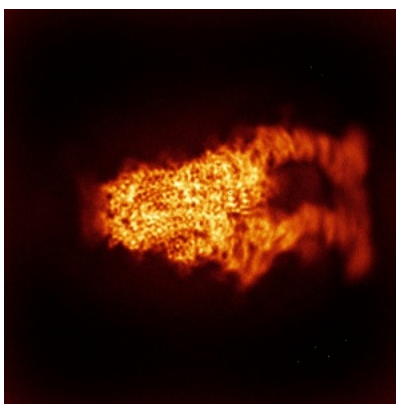


Z

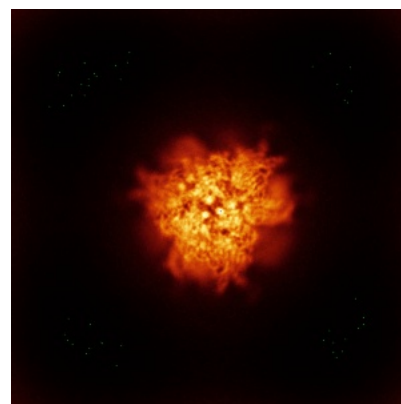
6.4.2 Raw map



X



Y

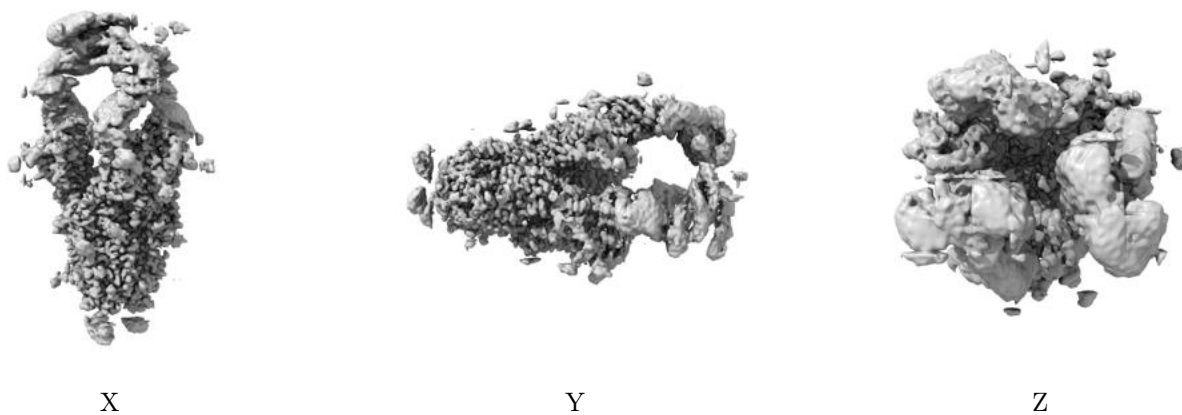


Z

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.

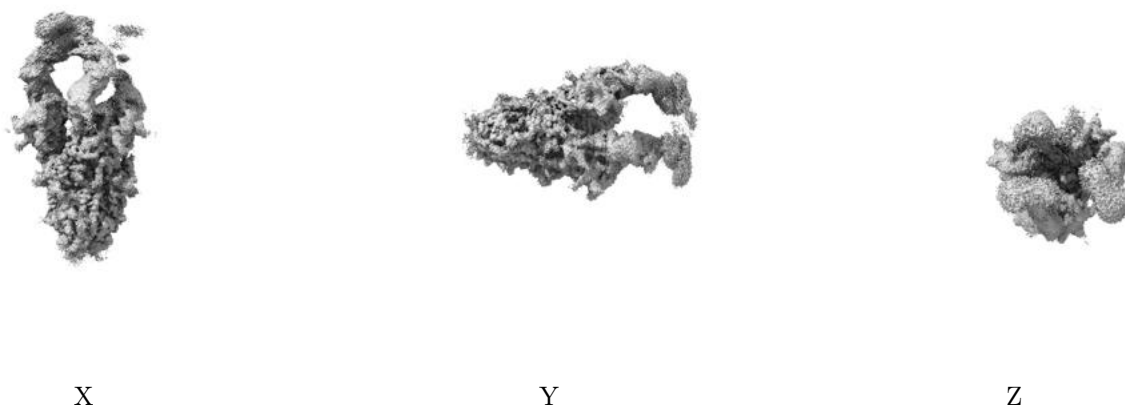
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.006. 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.

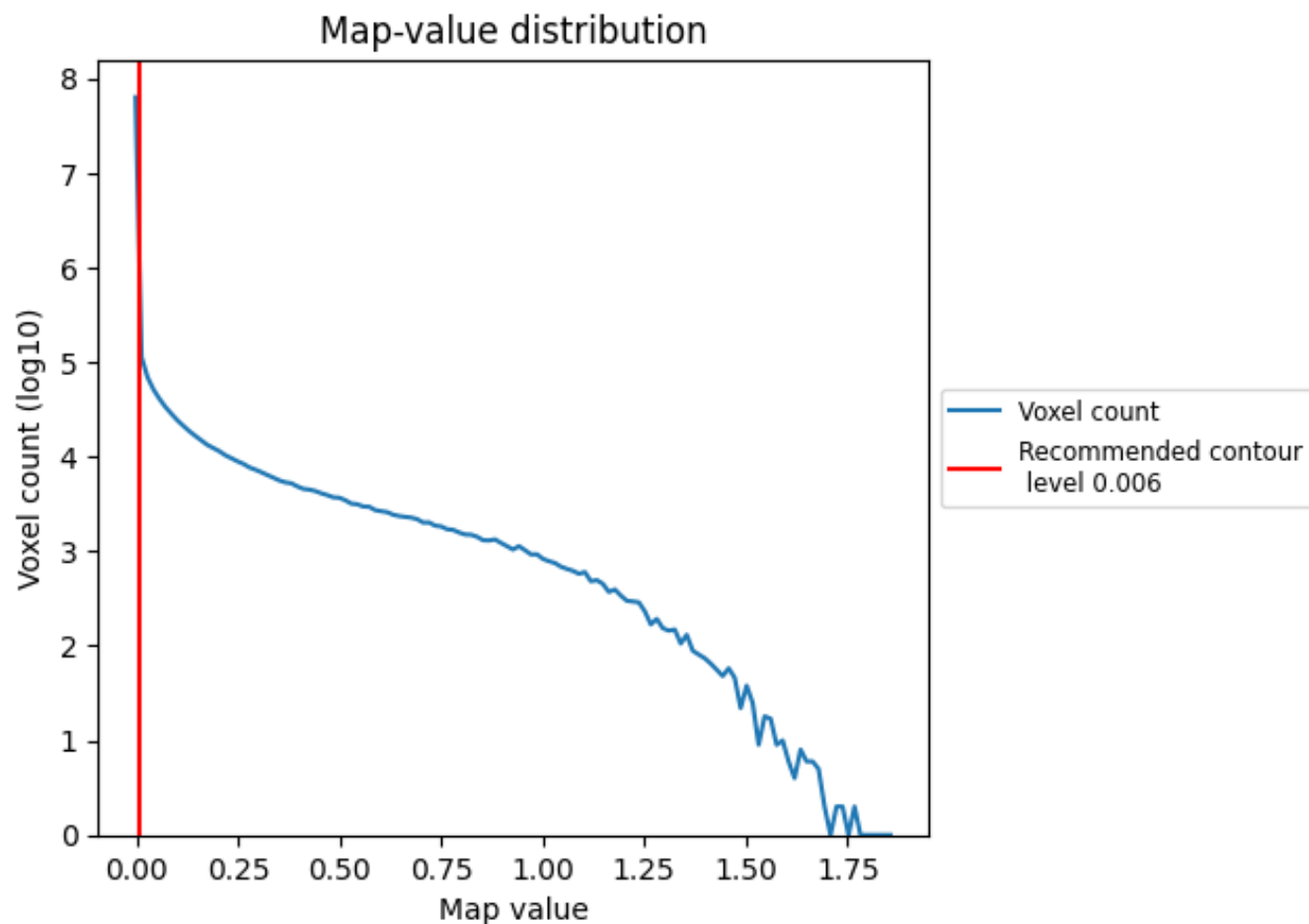
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

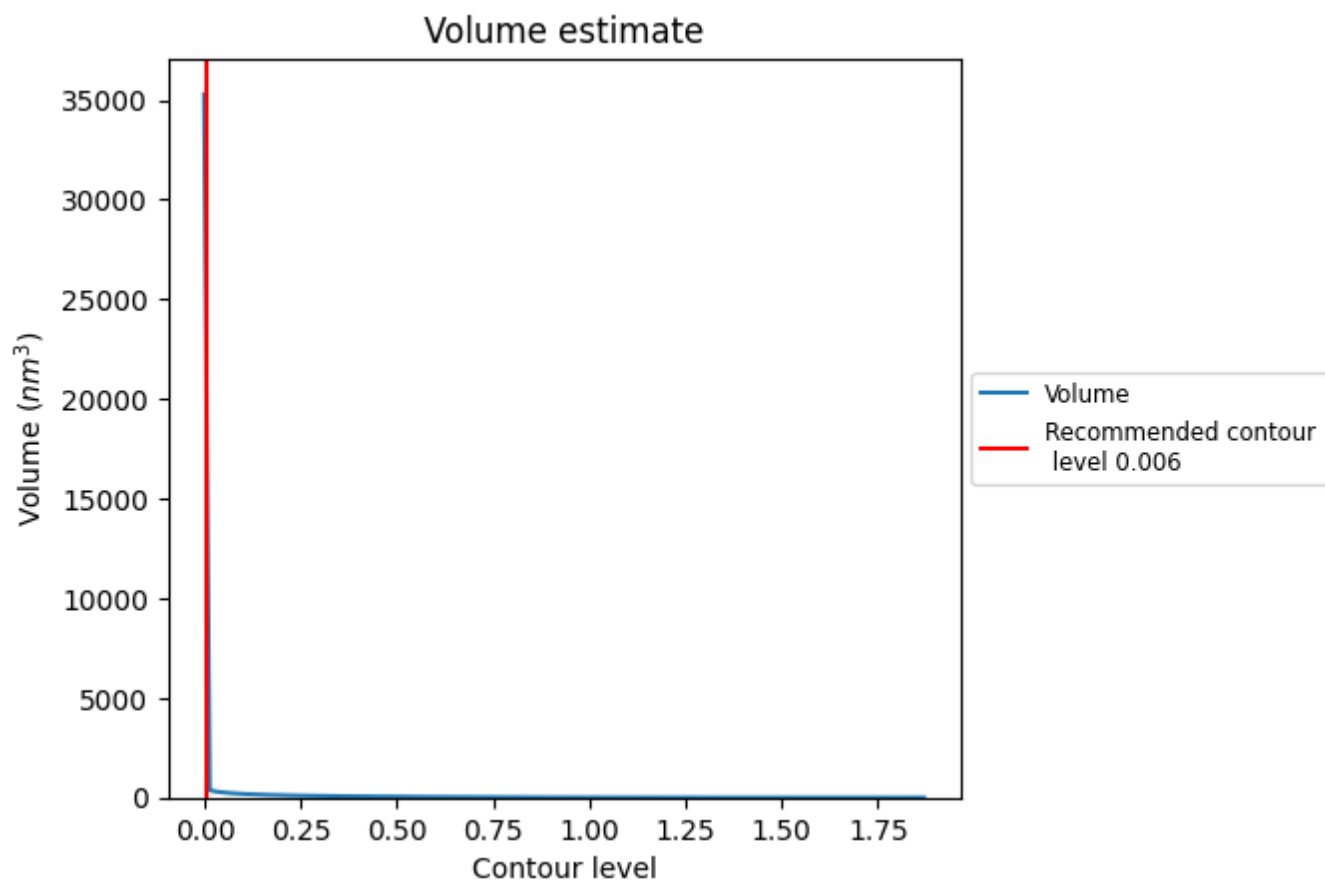
This section contains the results of statistical analysis of the map.

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.

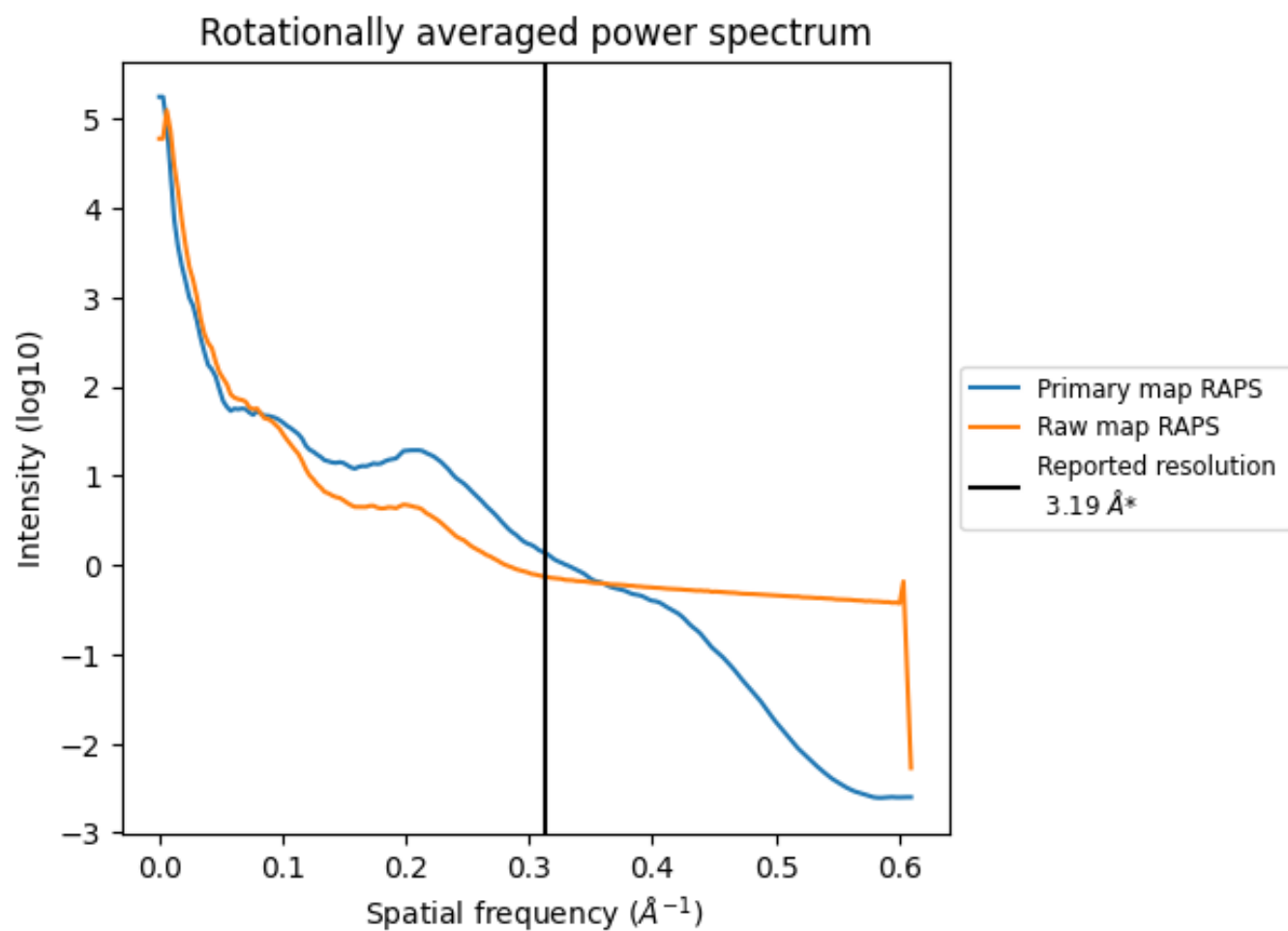
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 16960 nm^3 ; this corresponds to an approximate mass of 15320 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 ⓘ

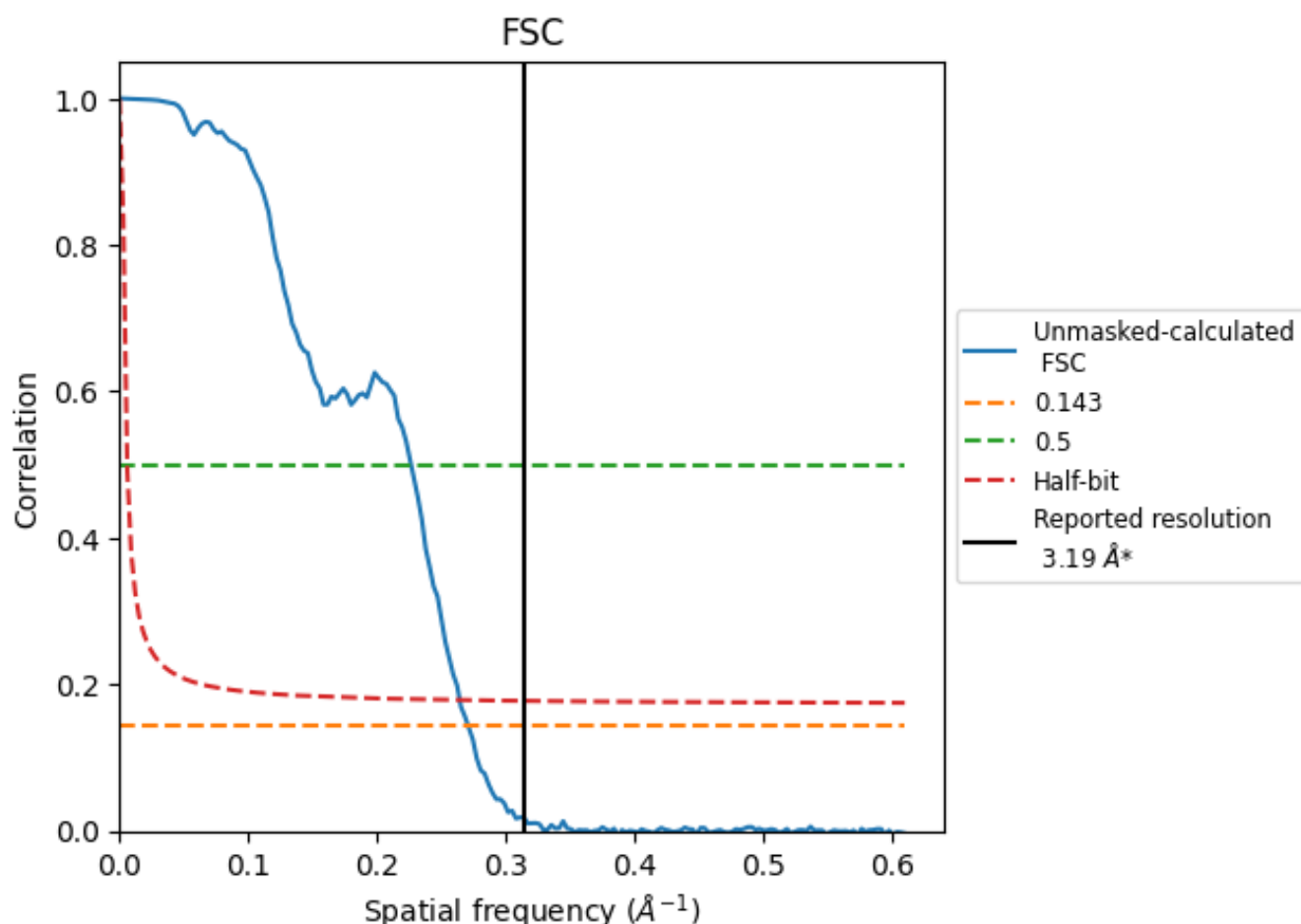


*Reported resolution corresponds to spatial frequency of 0.313 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.313 Å⁻¹

8.2 Resolution estimates [i](#)

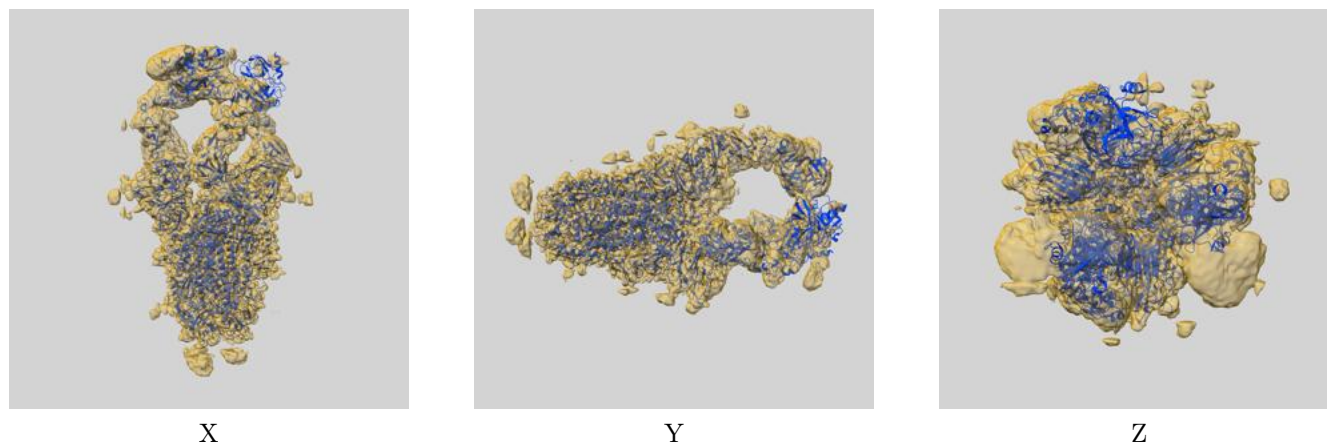
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.19	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.69	4.42	3.78

*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 3.69 differs from the reported value 3.19 by more than 10 %

9 Map-model fit [i](#)

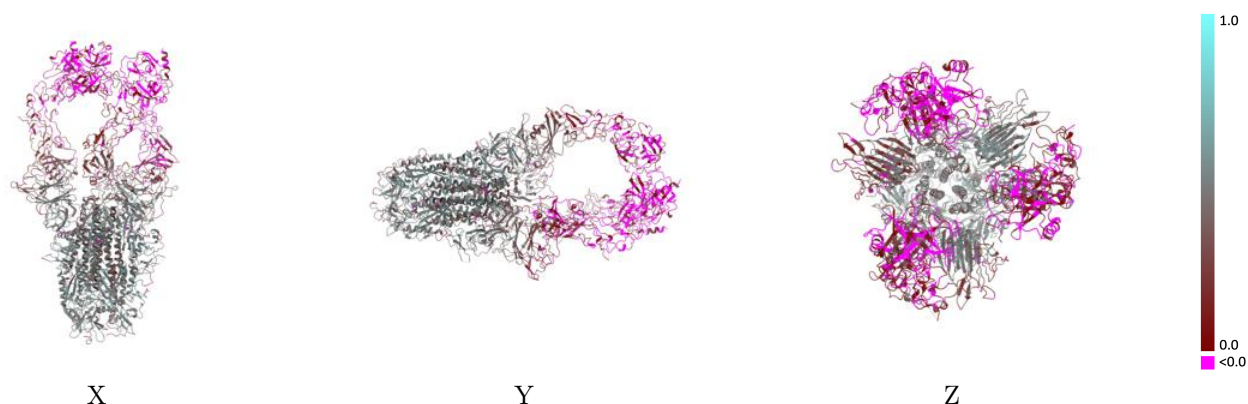
This section contains information regarding the fit between EMDB map EMD-39039 and PDB model 8Y8A. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



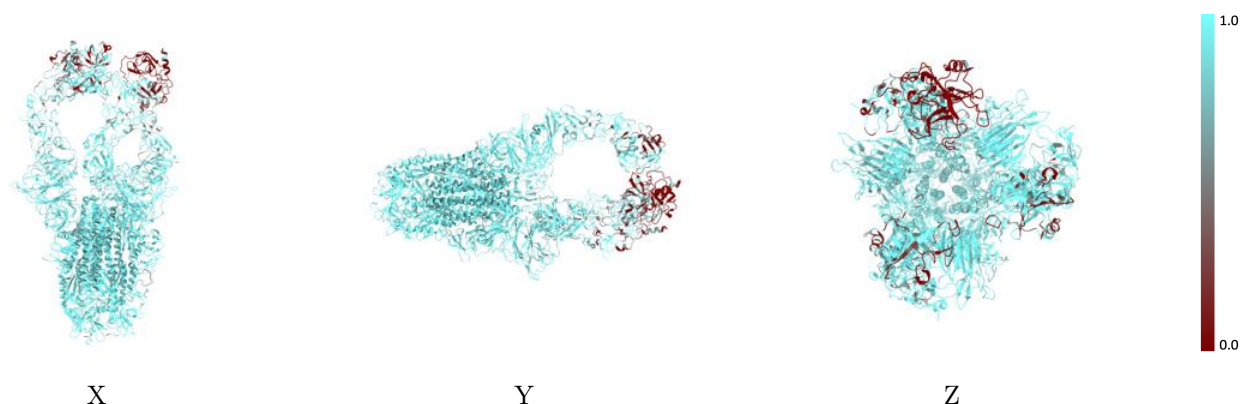
The images above show the 3D surface view of the map at the recommended contour level 0.006 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



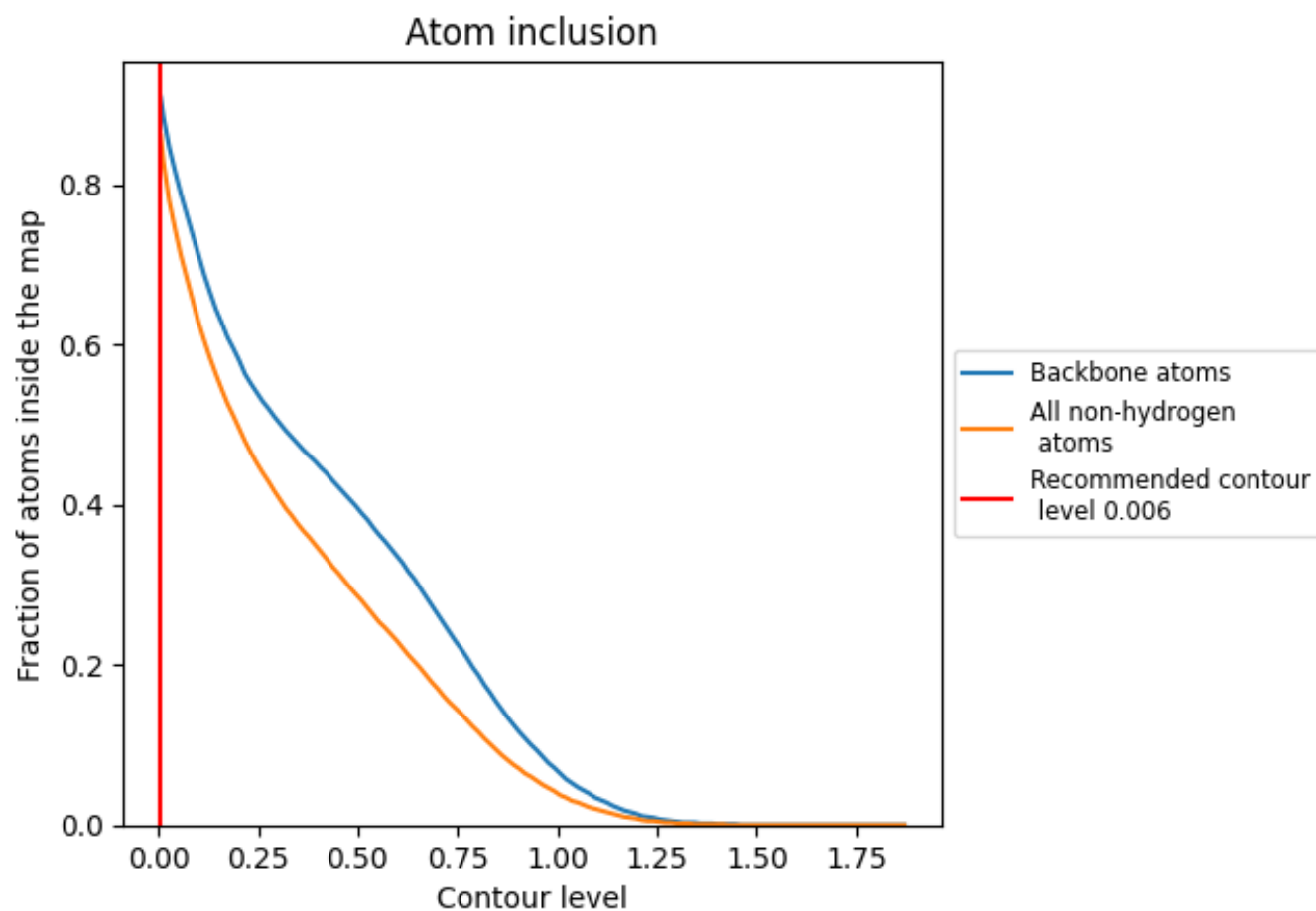
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.006).









































9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 86% 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 (0.006) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8610	 0.3080
A	 0.9150	 0.3580
B	 0.9340	 0.3890
C	 0.9250	 0.3600
D	 0.8570	 0.3210
E	 0.8570	 0.3120
F	 0.8570	 0.2520
G	 0.2290	 -0.0380
H	 0.6860	 0.0080
I	 0.9490	 0.3750
J	 0.8930	 0.2670
K	 0.8210	 0.2810
L	 0.6430	 0.2440
M	 0.8210	 0.2320
N	 0.8570	 0.2900
O	 0.7500	 0.1040
P	 1.0000	 0.3390
Q	 0.7500	 0.2730
R	 0.7500	 0.2300
T	 0.6520	 -0.0150

