



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 09:17 AM UTC

PDB ID : 8SAK / pdb_00008sak
EMDB ID : EMD-40272
Title : BtCoV-422 in complex with neutralizing antibody JC57-11
Authors : McFadden, E.; McLellan, J.S.
Deposited on : 2023-04-01
Resolution : 3.00 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:

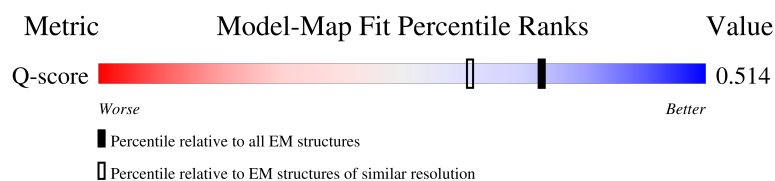
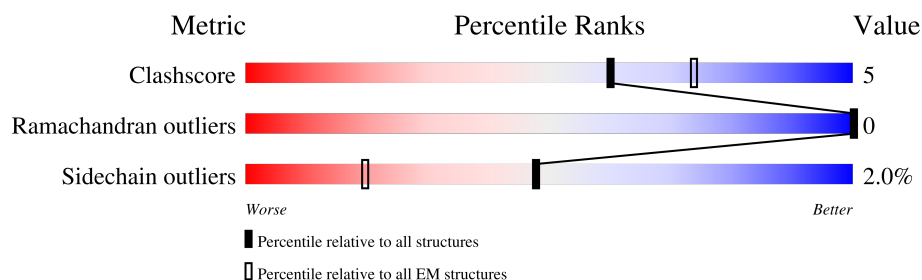
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 3.00 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.



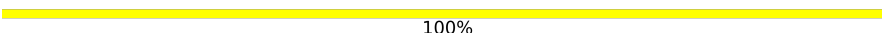
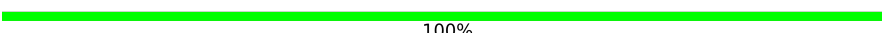
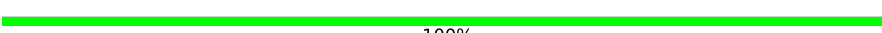
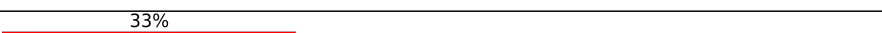
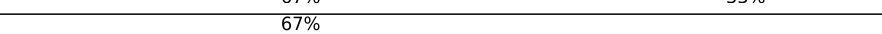
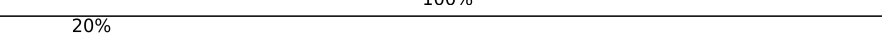
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14081 (2.50 - 3.50)

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	1289	
1	B	1289	
1	C	1289	
2	H	237	

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Mol	Chain	Length	Quality of chain
3	L	216	
4	5	2	
4	7	2	
4	E	2	
4	K	2	
4	U	2	
4	X	2	
4	c	2	
4	f	2	
4	l	2	
4	n	2	
4	p	2	
4	y	2	
5	1	3	
5	Q	3	
5	h	3	
6	t	5	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 28520 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 Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1121	Total	C	N	O	S	0	0
			8662	5511	1428	1670	53		
1	B	1124	Total	C	N	O	S	0	0
			8684	5523	1432	1676	53		
1	C	1117	Total	C	N	O	S	0	0
			8634	5494	1424	1664	52		

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	885	PRO	ALA	engineered mutation	UNP A0A2R4KP93
A	1056	PRO	ALA	engineered mutation	UNP A0A2R4KP93
A	1057	PRO	VAL	engineered mutation	UNP A0A2R4KP93
B	885	PRO	ALA	engineered mutation	UNP A0A2R4KP93
B	1056	PRO	ALA	engineered mutation	UNP A0A2R4KP93
B	1057	PRO	VAL	engineered mutation	UNP A0A2R4KP93
C	885	PRO	ALA	engineered mutation	UNP A0A2R4KP93
C	1056	PRO	ALA	engineered mutation	UNP A0A2R4KP93
C	1057	PRO	VAL	engineered mutation	UNP A0A2R4KP93

- Molecule 2 is a protein called JC57-11 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	131	Total	C	N	O	S	0	0
			996	629	167	196	4		

- Molecule 3 is a protein called JC57-11 Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	L	105	Total	C	N	O	S	0	0
			781	484	137	157	3		

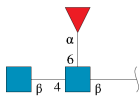
- Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-a

cetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
4	E	2	Total	C	N	O	0	0
			28	16	2	10		
4	K	2	Total	C	N	O	0	0
			28	16	2	10		
4	U	2	Total	C	N	O	0	0
			28	16	2	10		
4	X	2	Total	C	N	O	0	0
			28	16	2	10		
4	c	2	Total	C	N	O	0	0
			28	16	2	10		
4	f	2	Total	C	N	O	0	0
			28	16	2	10		
4	l	2	Total	C	N	O	0	0
			28	16	2	10		
4	n	2	Total	C	N	O	0	0
			28	16	2	10		
4	p	2	Total	C	N	O	0	0
			28	16	2	10		
4	y	2	Total	C	N	O	0	0
			28	16	2	10		
4	5	2	Total	C	N	O	0	0
			28	16	2	10		
4	7	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 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.



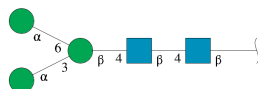
Mol	Chain	Residues	Atoms				AltConf	Trace
5	Q	3	Total	C	N	O	0	0
			38	22	2	14		
5	h	3	Total	C	N	O	0	0
			38	22	2	14		

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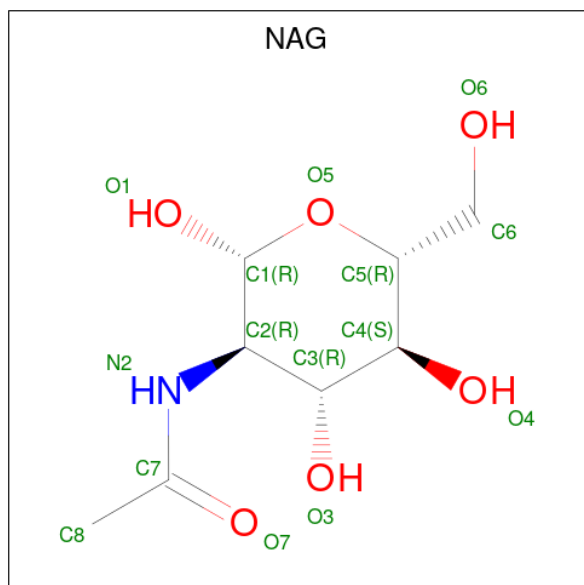
Mol	Chain	Residues	Atoms				AltConf	Trace
5	1	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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
6	t	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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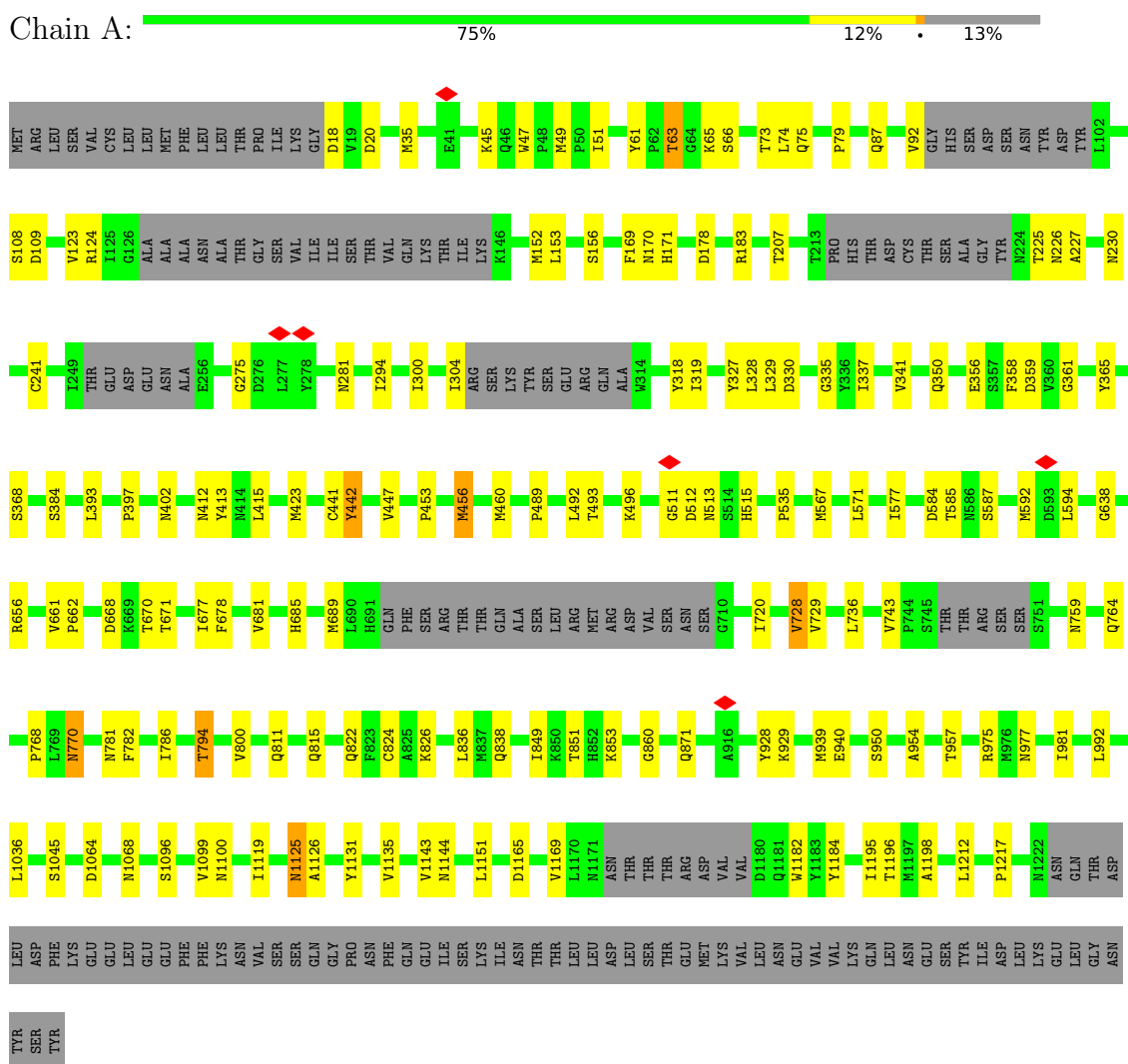
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Mol	Chain	Residues	Atoms				AltConf
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C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

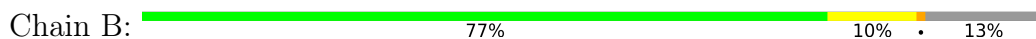
3 Residue-property plots [i](#)

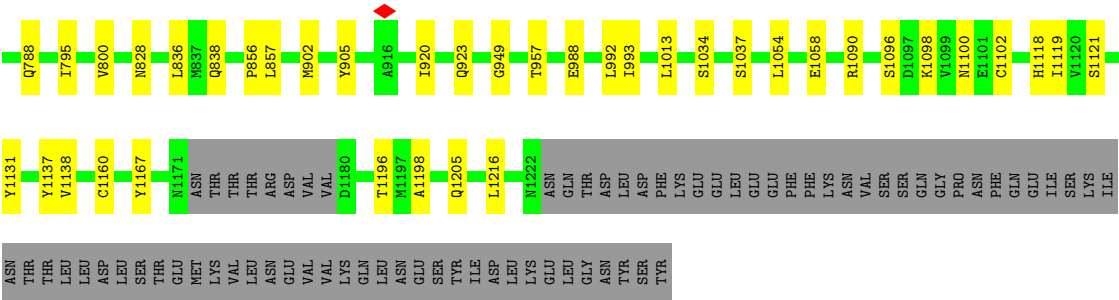
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Spike glycoprotein

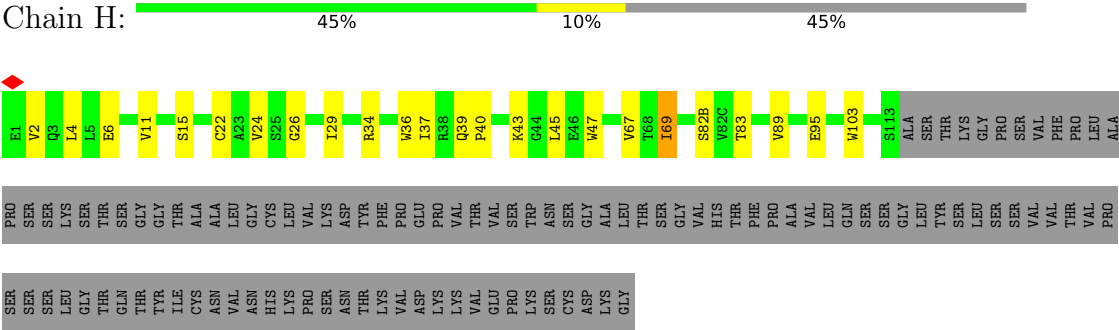


• Molecule 1: Spike glycoprotein

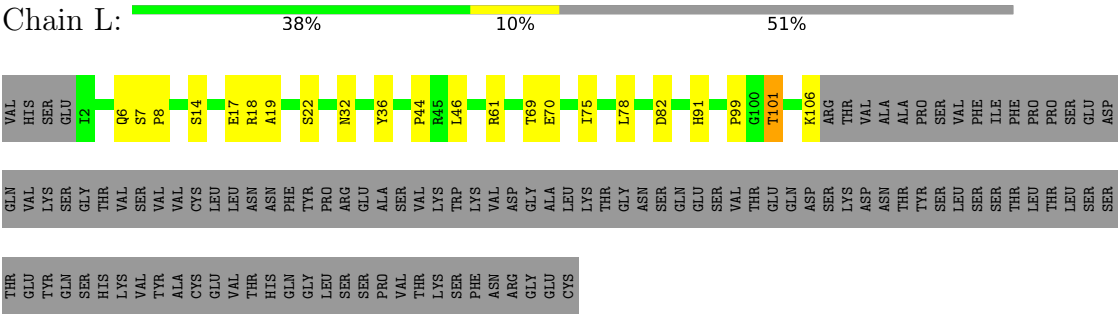




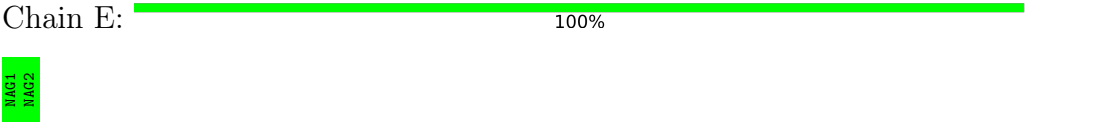
• Molecule 2: JC57-11 Fab heavy chain



• Molecule 3: JC57-11 Fab light chain



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



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



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

Chain U:  100%



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

Chain X:  100%



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

Chain c:  100%



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

Chain f:  100%



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

Chain l:  50%  50%



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

Chain n:  100%



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

Chain p:  50%  50%



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

Chain y:



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

Chain 5:



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

Chain 7:



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain h:



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 1:



- Molecule 6: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-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	57821	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 ($e^-/\text{\AA}^2$)	80.5	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	158.847	Depositor
Minimum map value	-7.304	Depositor
Average map value	0.010	Depositor
Map value standard deviation	1.190	Depositor
Recommended contour level	1	Depositor
Map size (Å)	333.28, 333.28, 333.28	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8332, 0.8332, 0.8332	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, MAN, BMA, FUC

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	0.12	0/8866	0.24	0/12055
1	B	0.13	0/8888	0.24	0/12085
1	C	0.12	0/8837	0.24	0/12015
2	H	0.26	0/1020	0.37	0/1389
3	L	0.19	0/798	0.33	0/1082
All	All	0.13	0/28409	0.25	0/38626

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	A	8662	0	8331	88	0
1	B	8684	0	8350	77	0
1	C	8634	0	8306	79	0
2	H	996	0	971	14	0
3	L	781	0	757	13	0
4	5	28	0	25	1	0
4	7	28	0	25	1	0
4	E	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	K	28	0	25	0	0
4	U	28	0	25	0	0
4	X	28	0	25	0	0
4	c	28	0	25	0	0
4	f	28	0	25	1	0
4	l	28	0	25	0	0
4	n	28	0	25	0	0
4	p	28	0	25	0	0
4	y	28	0	25	1	0
5	l	38	0	34	0	0
5	Q	38	0	34	0	0
5	h	38	0	34	0	0
6	t	61	0	52	1	0
7	A	126	0	117	1	0
7	B	70	0	65	0	0
7	C	56	0	52	0	0
All	All	28520	0	27403	253	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 5.

All (253) 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:B:866:ASN:H	1:B:998:ASN:HD21	1.35	0.73
2:H:39:GLN:HB2	2:H:45:LEU:HD23	1.71	0.72
1:C:472:ASN:HD21	1:C:503:ILE:H	1.37	0.71
3:L:6:GLN:NE2	3:L:101:THR:OG1	2.22	0.71
1:B:862:ASN:HD22	1:B:867:LEU:HD12	1.55	0.71
1:C:162:ASN:ND2	1:C:203:SER:O	2.25	0.69
3:L:6:GLN:HG3	3:L:99:PRO:HD2	1.75	0.68
1:A:1064:ASP:OD1	1:A:1068:ASN:ND2	2.27	0.68
1:A:853:LYS:HG2	1:C:764:GLN:HB2	1.76	0.66
1:C:1096:SER:O	1:C:1100:ASN:ND2	2.28	0.66
1:A:794:THR:OG1	1:A:838:GLN:NE2	2.29	0.66
1:A:350:GLN:NE2	1:A:368:SER:OG	2.30	0.64
1:B:496:LYS:HG2	1:B:571:LEU:HB2	1.78	0.64
1:C:677:ILE:HG12	1:C:719:VAL:HG12	1.79	0.64
1:A:74:LEU:HD13	1:A:329:LEU:HD12	1.79	0.64
1:A:226:ASN:O	1:A:230:ASN:ND2	2.24	0.63
1:A:153:LEU:HD23	1:A:294:ILE:HG21	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:262:GLN:HE22	1:B:294:ILE:H	1.47	0.63
1:B:661:VAL:HG13	1:B:681:VAL:HG11	1.80	0.63
1:B:263:ASN:HD22	1:B:264:THR:H	1.46	0.63
1:A:35:MET:HE2	1:A:207:THR:HG23	1.80	0.62
1:C:800:VAL:O	1:C:828:ASN:ND2	2.32	0.62
3:L:14:SER:HB3	3:L:106:LYS:HD2	1.82	0.62
2:H:36:TRP:HD1	2:H:69:ILE:HD12	1.65	0.61
1:A:384:SER:OG	1:A:412:ASN:OD1	2.18	0.61
1:B:768:PRO:HD3	1:C:856:PRO:HD2	1.81	0.61
1:A:453:PRO:HG2	1:A:456:MET:HG3	1.82	0.61
1:B:791:ILE:HD11	1:B:1098:LYS:HD2	1.83	0.60
1:A:786:ILE:HG12	1:A:1135:VAL:HG22	1.83	0.60
1:A:764:GLN:HB2	1:B:853:LYS:HG2	1.83	0.60
1:A:18:ASP:N	1:A:241:CYS:O	2.36	0.59
1:B:190:VAL:HG23	1:B:239:GLN:HG3	1.85	0.59
1:C:401:TYR:HB3	1:C:472:ASN:HD22	1.68	0.59
2:H:15:SER:HA	2:H:82(B):SER:HA	1.84	0.58
1:A:109:ASP:HA	1:A:207:THR:HG21	1.85	0.58
1:A:512:ASP:OD1	1:A:513:ASN:N	2.34	0.58
1:A:849:ILE:HG23	1:A:1099:VAL:HG11	1.86	0.58
1:B:233:LYS:HG2	1:B:238:LEU:HD13	1.86	0.58
1:A:496:LYS:HB3	1:A:567:MET:HE2	1.85	0.57
1:A:656:ARG:NH1	1:B:912:GLY:O	2.38	0.57
1:C:109:ASP:OD1	1:C:112:THR:OG1	2.22	0.57
1:A:75:GLN:HG3	1:B:818:ARG:HD2	1.87	0.57
1:B:953:GLY:HA3	1:B:957:THR:HG21	1.87	0.56
2:H:2:VAL:HG22	2:H:4:LEU:HD21	1.86	0.56
1:A:49:MET:HB3	1:A:124:ARG:HH11	1.70	0.56
1:B:715:ALA:HA	1:C:902:MET:HE2	1.87	0.56
3:L:19:ALA:N	3:L:75:ILE:O	2.38	0.56
1:A:156:SER:HB3	1:A:169:PHE:HB2	1.87	0.56
1:B:631:GLN:HE22	1:C:274:LYS:HG2	1.70	0.56
1:C:174:LEU:HB3	1:C:176:LEU:HD13	1.88	0.56
1:A:397:PRO:HG2	1:A:571:LEU:HD21	1.87	0.56
1:C:644:GLY:HA2	1:C:654:CYS:HA	1.88	0.56
1:B:979:ILE:HD13	1:B:1119:ILE:HD11	1.89	0.55
1:C:413:TYR:HE2	1:C:415:LEU:HD23	1.72	0.54
3:L:17:GLU:O	3:L:18:ARG:NH1	2.40	0.54
1:A:584:ASP:OD1	1:A:585:THR:N	2.41	0.54
1:B:976:MET:HE1	1:B:993:ILE:HG21	1.89	0.54
1:B:435:ASP:OD2	1:B:439:ARG:NH1	2.40	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:787:THR:HG23	1:B:1134:HIS:HB2	1.90	0.53
1:B:1064:ASP:OD1	1:B:1068:ASN:ND2	2.41	0.53
1:A:638:GLY:HA2	1:B:72:LEU:HD21	1.90	0.53
1:A:729:VAL:HG11	1:A:743:VAL:HG21	1.88	0.53
1:B:845:LEU:O	1:B:849:ILE:HG12	2.09	0.53
1:B:1064:ASP:O	1:B:1068:ASN:ND2	2.40	0.52
1:C:256:GLU:OE1	1:C:273:ARG:NH1	2.40	0.52
1:C:364:VAL:HG22	1:C:666:VAL:HG13	1.91	0.52
7:A:1304:NAG:O6	1:C:532:ASN:O	2.21	0.52
1:A:1096:SER:O	1:A:1100:ASN:ND2	2.37	0.52
1:B:786:ILE:HG12	1:B:1135:VAL:HG22	1.92	0.52
3:L:61:ARG:NH1	3:L:82:ASP:OD2	2.42	0.52
1:B:384:SER:HB2	1:B:412:ASN:HD22	1.75	0.52
1:C:501:SER:HB3	1:C:538:PHE:HD2	1.75	0.52
1:B:782:PHE:O	1:B:996:LYS:NZ	2.40	0.51
1:B:584:ASP:OD1	1:B:585:THR:N	2.43	0.51
1:C:397:PRO:HG3	1:C:571:LEU:HD21	1.93	0.51
2:H:4:LEU:HD22	2:H:24:VAL:HG12	1.90	0.51
1:C:109:ASP:OD2	1:C:113:LYS:NZ	2.39	0.51
1:B:972:ILE:HG22	1:B:976:MET:HE2	1.92	0.51
1:B:392:LEU:HD11	1:B:575:PHE:HE2	1.74	0.51
1:A:393:LEU:O	1:A:493:THR:OG1	2.22	0.51
1:C:148:TYR:O	1:C:302:ARG:NH2	2.38	0.51
1:A:73:THR:HG23	1:A:356:GLU:HG3	1.92	0.51
1:B:123:VAL:HG22	1:B:319:ILE:HG12	1.92	0.51
1:C:782:PHE:HZ	1:C:993:ILE:HG12	1.75	0.51
1:B:259:GLY:HA3	1:B:282:MET:HE1	1.93	0.50
1:B:425:ASN:OD1	1:B:457:LYS:NZ	2.42	0.50
1:B:866:ASN:H	1:B:998:ASN:ND2	2.06	0.50
1:B:357:SER:OG	1:B:359:ASP:OD1	2.30	0.50
3:L:7:SER:OG	3:L:22:SER:OG	2.17	0.50
1:A:781:ASN:OD1	1:A:782:PHE:N	2.44	0.50
1:C:123:VAL:HG22	1:C:319:ILE:HG12	1.92	0.50
1:C:74:LEU:HD12	1:C:329:LEU:HD12	1.93	0.50
1:B:147:ILE:HD11	1:B:302:ARG:HG2	1.93	0.50
1:B:393:LEU:O	1:B:493:THR:OG1	2.30	0.49
1:B:845:LEU:HD12	1:B:1096:SER:HB3	1.94	0.49
1:C:330:ASP:HB2	1:C:358:PHE:CE1	2.48	0.49
1:C:20:ASP:HB2	4:y:2:NAG:H3	1.94	0.49
1:A:728:VAL:HG12	1:A:759:ASN:HB2	1.93	0.49
1:B:800:VAL:HG12	1:B:928:TYR:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:678:PHE:HE1	1:A:720:ILE:HG12	1.78	0.49
1:C:402:ASN:ND2	1:C:535:PRO:O	2.45	0.49
1:A:447:VAL:HG22	1:A:577:ILE:HG12	1.95	0.49
1:B:892:ASN:ND2	1:B:892:ASN:O	2.45	0.49
1:A:662:PRO:HB3	1:B:929:LYS:HE2	1.94	0.49
1:C:108:SER:OG	1:C:109:ASP:N	2.46	0.49
1:C:351:LEU:HD11	1:C:365:TYR:CD1	2.48	0.48
1:B:679:GLY:O	1:B:717:GLY:HA3	2.13	0.48
1:A:92:VAL:HG13	1:A:304:ILE:HG13	1.93	0.48
1:A:413:TYR:OH	1:A:442:TYR:OH	2.28	0.48
1:A:800:VAL:HG12	1:A:928:TYR:HB3	1.95	0.48
1:A:123:VAL:HG22	1:A:319:ILE:HG12	1.96	0.48
1:B:476:SER:HB3	1:B:524:PRO:HB2	1.96	0.48
1:B:384:SER:HB2	1:B:412:ASN:ND2	2.29	0.48
1:A:977:ASN:ND2	1:A:981:ILE:O	2.47	0.47
1:A:275:GLY:HA2	1:C:631:GLN:HE22	1.78	0.47
1:A:413:TYR:HE2	1:A:415:LEU:HD13	1.79	0.47
1:C:795:ILE:O	1:C:838:GLN:NE2	2.32	0.47
2:H:2:VAL:HG22	2:H:4:LEU:CD2	2.43	0.47
1:A:489:PRO:HB2	1:A:492:LEU:HD23	1.97	0.47
1:C:584:ASP:OD1	1:C:585:THR:N	2.48	0.47
1:A:328:LEU:HB3	1:A:341:VAL:HG23	1.95	0.47
1:A:1126:ALA:HB2	1:A:1131:TYR:HB2	1.96	0.47
1:B:904:GLY:HA2	1:B:917:ARG:HH22	1.79	0.47
1:A:1184:TYR:HB2	1:A:1195:ILE:HD13	1.96	0.47
1:C:1034:SER:O	1:C:1037:SER:OG	2.33	0.47
4:7:1:NAG:O6	4:7:2:NAG:O3	2.31	0.47
1:C:686:ILE:HG23	1:C:690:LEU:HD22	1.97	0.47
1:C:1119:ILE:HD11	1:C:1137:TYR:HB2	1.97	0.46
1:C:788:GLN:OE1	1:C:1131:TYR:OH	2.21	0.46
1:C:1098:LYS:HG3	1:C:1102:CYS:SG	2.55	0.46
1:C:1118:HIS:NE2	1:C:1121:SER:OG	2.43	0.46
3:L:32:ASN:HB3	3:L:91:HIS:NE2	2.30	0.46
1:A:512:ASP:H	1:A:515:HIS:CD2	2.33	0.46
1:B:1051:LEU:HD11	1:B:1063:ILE:HD11	1.97	0.46
1:B:1164:ILE:O	1:B:1184:TYR:OH	2.26	0.46
1:C:372:VAL:HG13	1:C:660:SER:HB3	1.97	0.46
1:A:423:MET:HE3	1:C:456:MET:HE2	1.97	0.46
1:A:178:ASP:OD2	1:A:227:ALA:N	2.49	0.46
1:C:111:ALA:HB3	1:C:207:THR:HG22	1.97	0.46
1:B:413:TYR:HH	1:B:442:TYR:HH	1.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:685:HIS:HA	1:A:689:MET:HE2	1.97	0.46
1:A:957:THR:HG21	1:C:1167:TYR:CZ	2.51	0.45
1:B:799:THR:HA	1:B:831:LEU:HD13	1.98	0.45
1:B:1087:THR:HG23	1:C:1090:ARG:HH12	1.81	0.45
1:A:860:GLY:H	1:A:871:GLN:HE22	1.65	0.45
1:B:170:ASN:C	1:B:190:VAL:HG12	2.42	0.45
1:C:75:GLN:NE2	1:C:326:THR:OG1	2.49	0.45
2:H:36:TRP:CD1	2:H:69:ILE:HD12	2.47	0.45
1:C:664:SER:HB2	1:C:677:ILE:HB	1.98	0.45
3:L:69:THR:OG1	3:L:70:GLU:OE1	2.33	0.45
1:A:677:ILE:HD11	1:A:736:LEU:HD22	1.99	0.45
1:A:768:PRO:HD3	1:B:856:PRO:HD2	1.98	0.45
1:C:426:GLU:OE1	1:C:457:LYS:NZ	2.45	0.45
6:t:1:NAG:H62	6:t:2:NAG:C7	2.47	0.45
1:A:361:GLY:O	1:A:365:TYR:OH	2.30	0.45
1:B:74:LEU:HD12	1:B:329:LEU:HD12	1.98	0.45
1:B:772:SER:OG	1:B:1210:GLU:OE2	2.27	0.44
1:A:1064:ASP:O	1:A:1068:ASN:ND2	2.49	0.44
1:B:178:ASP:OD1	1:B:179:GLY:N	2.40	0.44
1:C:512:ASP:O	1:C:513:ASN:ND2	2.51	0.44
4:5:1:NAG:H4	4:5:2:NAG:H2	1.78	0.44
1:B:1015:PHE:O	1:B:1018:VAL:HG12	2.18	0.44
1:C:573:MET:HE3	1:C:573:MET:HB3	1.69	0.44
1:A:47:TRP:O	1:A:318:TYR:OH	2.34	0.44
1:A:152:MET:HE3	1:A:300:ILE:HD11	1.99	0.44
1:B:549:LEU:HB3	1:B:553:GLU:HB3	1.99	0.44
1:B:681:VAL:HG12	1:C:905:TYR:CG	2.53	0.44
1:A:61:TYR:OH	1:A:335:GLY:O	2.26	0.44
1:A:63:THR:OG1	1:C:583:THR:O	2.28	0.44
1:A:511:GLY:HA3	1:A:515:HIS:CD2	2.52	0.44
1:B:123:VAL:HG11	1:B:149:PRO:HG2	1.99	0.44
1:C:387:CYS:HB2	1:C:413:TYR:HB3	1.99	0.44
1:B:152:MET:HE3	1:B:168:TYR:CE1	2.53	0.43
1:C:67:TYR:HB3	1:C:70:ILE:HD11	1.99	0.43
2:H:6:GLU:HA	2:H:22:CYS:HA	1.99	0.43
2:H:40:PRO:HB2	2:H:43:LYS:HB2	2.00	0.43
3:L:36:TYR:CZ	3:L:46:LEU:HD13	2.53	0.43
1:B:397:PRO:HG3	1:B:571:LEU:HD21	2.00	0.43
1:B:44:ASN:OD1	1:B:46:GLN:NE2	2.40	0.43
1:B:590:PRO:HG2	1:B:592:MET:HG2	2.00	0.43
1:C:342:ASP:O	1:C:349:SER:OG	2.29	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:ASN:HD21	1:A:337:ILE:HB	1.83	0.43
1:A:1036:LEU:O	1:A:1045:SER:OG	2.37	0.43
1:B:123:VAL:N	1:B:258:PHE:O	2.35	0.43
1:B:714:THR:OG1	1:B:717:GLY:O	2.34	0.43
1:B:377:GLN:O	1:B:603:LYS:NZ	2.51	0.43
1:A:1165:ASP:OD1	1:A:1184:TYR:OH	2.36	0.43
1:B:159:ASN:OD1	1:B:159:ASN:N	2.52	0.43
1:B:800:VAL:HA	1:B:928:TYR:HA	2.01	0.43
2:H:2:VAL:CG2	2:H:4:LEU:HD21	2.49	0.43
1:A:822:GLN:O	1:A:826:LYS:HG2	2.19	0.43
1:C:920:ILE:HA	1:C:923:GLN:HE21	1.84	0.43
3:L:7:SER:HG	3:L:8:PRO:HD3	1.84	0.43
1:A:402:ASN:ND2	1:A:535:PRO:O	2.52	0.43
1:B:1001:LEU:O	1:B:1004:MET:HG2	2.19	0.43
1:C:380:GLU:OE2	1:C:647:SER:OG	2.34	0.43
1:C:413:TYR:HH	1:C:442:TYR:HH	1.62	0.43
1:C:413:TYR:OH	1:C:442:TYR:OH	2.29	0.43
1:A:1195:ILE:HB	1:A:1217:PRO:HB3	2.00	0.42
3:L:106:LYS:HB3	3:L:106:LYS:HE2	1.83	0.42
1:A:45:LYS:O	1:A:87:GLN:NE2	2.27	0.42
1:A:183:ARG:HD2	1:A:225:THR:HB	2.01	0.42
1:B:887:GLU:OE2	1:B:1125:ASN:HB2	2.19	0.42
1:C:19:VAL:HG21	1:C:243:PHE:HA	2.01	0.42
1:C:123:VAL:HG11	1:C:149:PRO:HG2	2.00	0.42
1:C:600:ILE:HD11	1:C:608:VAL:HG21	2.00	0.42
1:A:975:ARG:HD3	1:A:1119:ILE:O	2.19	0.42
1:B:29:SER:OG	1:B:197:SER:OG	2.30	0.42
1:C:347:ASP:OD1	1:C:367:VAL:HG21	2.19	0.42
1:A:170:ASN:HB3	1:A:171:HIS:H	1.63	0.42
1:C:376:GLY:HA3	1:C:603:LYS:HD2	2.02	0.42
2:H:37:ILE:HD13	2:H:47:TRP:HA	2.02	0.42
1:A:1182:TRP:HE1	1:A:1212:LEU:HD21	1.83	0.42
1:B:568:THR:OG1	1:B:570:THR:O	2.37	0.42
4:f:1:NAG:H61	4:f:2:NAG:C7	2.50	0.42
1:A:66:SER:HB2	1:C:632:GLN:HB2	2.00	0.42
1:A:592:MET:HE3	1:A:594:LEU:HD11	2.01	0.42
1:A:1151:LEU:HD22	1:A:1212:LEU:HD11	2.02	0.42
1:C:49:MET:O	1:C:124:ARG:NH1	2.53	0.42
1:C:607:CYS:HA	1:C:620:VAL:HG12	2.01	0.42
1:B:473:TYR:CE1	1:B:576:ILE:HG13	2.55	0.42
1:C:730:ASP:OD1	1:C:759:ASN:ND2	2.43	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:940:GLU:HB3	1:A:1125:ASN:HD21	1.84	0.41
1:C:1054:LEU:HD13	1:C:1058:GLU:HG2	2.02	0.41
1:A:811:GLN:NE2	1:A:815:GLN:OE1	2.46	0.41
1:B:611:ASN:ND2	1:B:616:SER:HB3	2.35	0.41
2:H:2:VAL:HA	2:H:26:GLY:HA3	2.01	0.41
2:H:34:ARG:NE	2:H:95:GLU:OE1	2.36	0.41
1:A:79:PRO:O	1:A:327:TYR:OH	2.30	0.41
1:A:330:ASP:HB2	1:A:358:PHE:CE1	2.56	0.41
1:A:851:THR:HG21	1:A:950:SER:HB2	2.02	0.41
1:A:929:LYS:HD3	1:C:662:PRO:HB3	2.02	0.41
1:A:954:ALA:HB2	1:C:1167:TYR:HE1	1.85	0.41
1:A:668:ASP:HB3	1:A:671:THR:HG22	2.03	0.41
1:B:193:PRO:CB	1:B:200:PRO:HB2	2.51	0.41
1:A:65:LYS:HG2	1:C:636:TYR:CE1	2.56	0.41
1:A:441:CYS:HB3	1:A:587:SER:HB3	2.01	0.41
1:A:939:MET:SD	1:C:760:TYR:HB2	2.61	0.41
1:A:1196:THR:HG23	1:A:1198:ALA:H	1.86	0.41
1:C:1160:CYS:SG	1:C:1205:GLN:HA	2.61	0.41
2:H:103:TRP:CE3	3:L:44:PRO:HD2	2.56	0.41
1:B:274:LYS:HD3	1:B:274:LYS:HA	1.84	0.41
1:B:728:VAL:HG12	1:B:759:ASN:HB2	2.01	0.41
1:A:770:ASN:O	1:A:770:ASN:ND2	2.54	0.40
1:A:108:SER:OG	1:A:109:ASP:N	2.54	0.40
1:C:568:THR:OG1	1:C:570:THR:O	2.38	0.40
1:C:857:LEU:HD11	1:C:949:GLY:HA3	2.02	0.40
1:C:1196:THR:HG23	1:C:1198:ALA:H	1.86	0.40
1:A:20:ASP:OD1	1:A:20:ASP:N	2.51	0.40
1:C:148:TYR:H	1:C:302:ARG:NH2	2.20	0.40
1:C:591:MET:SD	1:C:591:MET:N	2.93	0.40
1:C:1216:LEU:H	1:C:1216:LEU:HD12	1.87	0.40
1:C:193:PRO:HB3	1:C:200:PRO: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	1103/1289 (86%)	1081 (98%)	22 (2%)	0	100	100
1	B	1106/1289 (86%)	1078 (98%)	28 (2%)	0	100	100
1	C	1097/1289 (85%)	1078 (98%)	19 (2%)	0	100	100
2	H	129/237 (54%)	121 (94%)	8 (6%)	0	100	100
3	L	103/216 (48%)	95 (92%)	8 (8%)	0	100	100
All	All	3538/4320 (82%)	3453 (98%)	85 (2%)	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	A	967/1121 (86%)	948 (98%)	19 (2%)	48	76
1	B	969/1121 (86%)	953 (98%)	16 (2%)	53	78
1	C	964/1121 (86%)	944 (98%)	20 (2%)	47	75
2	H	112/203 (55%)	106 (95%)	6 (5%)	20	53
3	L	87/188 (46%)	85 (98%)	2 (2%)	44	74
All	All	3099/3754 (83%)	3036 (98%)	63 (2%)	48	76

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	ILE
1	A	63	THR
1	A	359	ASP
1	A	442	TYR
1	A	456	MET
1	A	460	MET
1	A	661	VAL

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Mol	Chain	Res	Type
1	A	670	THR
1	A	681	VAL
1	A	728	VAL
1	A	770	ASN
1	A	794	THR
1	A	824	CYS
1	A	836	LEU
1	A	992	LEU
1	A	1125	ASN
1	A	1143	VAL
1	A	1144	ASN
1	A	1169	VAL
1	B	73	THR
1	B	115	VAL
1	B	159	ASN
1	B	189	CYS
1	B	263	ASN
1	B	302	ARG
1	B	348	LEU
1	B	396	THR
1	B	442	TYR
1	B	681	VAL
1	B	732	CYS
1	B	787	THR
1	B	892	ASN
1	B	895	THR
1	B	957	THR
1	B	1052	LYS
1	C	72	LEU
1	C	90	TYR
1	C	207	THR
1	C	341	VAL
1	C	372	VAL
1	C	381	GLN
1	C	396	THR
1	C	456	MET
1	C	515	HIS
1	C	573	MET
1	C	604	LEU
1	C	620	VAL
1	C	624	CYS
1	C	761	THR

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Mol	Chain	Res	Type
1	C	836	LEU
1	C	957	THR
1	C	988	GLU
1	C	992	LEU
1	C	1013	LEU
1	C	1138	VAL
2	H	11	VAL
2	H	29	ILE
2	H	67	VAL
2	H	69	ILE
2	H	83	THR
2	H	89	VAL
3	L	78	LEU
3	L	101	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (60) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	HIS
1	A	350	GLN
1	A	515	HIS
1	A	540	GLN
1	A	548	GLN
1	A	651	ASN
1	A	733	GLN
1	A	792	GLN
1	A	804	GLN
1	A	838	GLN
1	A	871	GLN
1	A	970	GLN
1	A	1023	ASN
1	A	1059	GLN
1	A	1125	ASN
1	A	1142	HIS
1	A	1181	GLN
1	B	118	ASN
1	B	119	ASN
1	B	262	GLN
1	B	263	ASN
1	B	377	GLN
1	B	412	ASN
1	B	513	ASN

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Mol	Chain	Res	Type
1	B	611	ASN
1	B	631	GLN
1	B	641	ASN
1	B	752	GLN
1	B	777	GLN
1	B	788	GLN
1	B	804	GLN
1	B	828	ASN
1	B	911	GLN
1	B	998	ASN
1	B	1141	ASN
1	B	1191	ASN
1	C	75	GLN
1	C	81	HIS
1	C	87	GLN
1	C	162	ASN
1	C	323	HIS
1	C	425	ASN
1	C	472	ASN
1	C	513	ASN
1	C	631	GLN
1	C	777	GLN
1	C	804	GLN
1	C	815	GLN
1	C	828	ASN
1	C	903	GLN
1	C	983	GLN
1	C	984	GLN
1	C	1027	GLN
1	C	1093	GLN
1	C	1191	ASN
1	C	1205	GLN
2	H	3	GLN
2	H	100(K)	ASN
3	L	6	GLN
3	L	37	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 ⓘ

38 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	$\# Z > 2$
5	NAG	1	1	5,1	14,14,15	0.26	0	17,19,21	0.46	0
5	NAG	1	2	5	14,14,15	0.26	0	17,19,21	0.41	0
5	FUC	1	3	5	10,10,11	0.58	0	14,14,16	0.74	0
4	NAG	5	1	4,1	14,14,15	0.66	0	17,19,21	1.24	1 (5%)
4	NAG	5	2	4	14,14,15	0.44	0	17,19,21	0.53	0
4	NAG	7	1	4,1	14,14,15	0.51	0	17,19,21	0.55	0
4	NAG	7	2	4	14,14,15	0.38	0	17,19,21	0.37	0
4	NAG	E	1	4,1	14,14,15	0.25	0	17,19,21	0.53	0
4	NAG	E	2	4	14,14,15	0.21	0	17,19,21	0.40	0
4	NAG	K	1	4,1	14,14,15	0.47	0	17,19,21	0.75	1 (5%)
4	NAG	K	2	4	14,14,15	0.25	0	17,19,21	0.50	0
5	NAG	Q	1	5,1	14,14,15	0.76	1 (7%)	17,19,21	1.38	1 (5%)
5	NAG	Q	2	5	14,14,15	0.23	0	17,19,21	0.43	0
5	FUC	Q	3	5	10,10,11	0.89	0	14,14,16	0.89	0
4	NAG	U	1	4,1	14,14,15	0.32	0	17,19,21	0.41	0
4	NAG	U	2	4	14,14,15	0.20	0	17,19,21	0.41	0
4	NAG	X	1	4,1	14,14,15	0.28	0	17,19,21	0.62	0
4	NAG	X	2	4	14,14,15	0.24	0	17,19,21	0.40	0
4	NAG	c	1	4,1	14,14,15	0.16	0	17,19,21	0.44	0
4	NAG	c	2	4	14,14,15	0.24	0	17,19,21	0.45	0
4	NAG	f	1	4,1	14,14,15	0.31	0	17,19,21	0.47	0
4	NAG	f	2	4	14,14,15	0.24	0	17,19,21	0.44	0
5	NAG	h	1	5,1	14,14,15	0.28	0	17,19,21	0.42	0
5	NAG	h	2	5	14,14,15	0.25	0	17,19,21	0.42	0
5	FUC	h	3	5	10,10,11	0.59	0	14,14,16	0.73	0
4	NAG	l	1	4,1	14,14,15	0.25	0	17,19,21	0.46	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	l	2	4	14,14,15	0.43	0	17,19,21	0.97	2 (11%)
4	NAG	n	1	4,1	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	n	2	4	14,14,15	0.19	0	17,19,21	0.45	0
4	NAG	p	1	4,1	14,14,15	0.44	0	17,19,21	0.78	1 (5%)
4	NAG	p	2	4	14,14,15	0.22	0	17,19,21	0.47	0
6	NAG	t	1	6,1	14,14,15	0.22	0	17,19,21	0.56	0
6	NAG	t	2	6	14,14,15	0.23	0	17,19,21	0.59	0
6	BMA	t	3	6	11,11,12	0.72	0	15,15,17	1.05	1 (6%)
6	MAN	t	4	6	11,11,12	0.88	0	15,15,17	1.08	2 (13%)
6	MAN	t	5	6	11,11,12	1.47	3 (27%)	15,15,17	2.16	2 (13%)
4	NAG	y	1	4,1	14,14,15	0.24	0	17,19,21	0.40	0
4	NAG	y	2	4	14,14,15	0.26	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	1	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	1	2	5	-	0/6/23/26	0/1/1/1
5	FUC	1	3	5	-	-	0/1/1/1
4	NAG	5	1	4,1	-	1/6/23/26	0/1/1/1
4	NAG	5	2	4	-	2/6/23/26	0/1/1/1
4	NAG	7	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	7	2	4	-	2/6/23/26	0/1/1/1
4	NAG	E	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	4/6/23/26	0/1/1/1
4	NAG	K	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	K	2	4	-	0/6/23/26	0/1/1/1
5	NAG	Q	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	Q	2	5	-	2/6/23/26	0/1/1/1
5	FUC	Q	3	5	-	-	0/1/1/1
4	NAG	U	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	U	2	4	-	3/6/23/26	0/1/1/1
4	NAG	X	1	4,1	-	4/6/23/26	0/1/1/1
4	NAG	X	2	4	-	4/6/23/26	0/1/1/1
4	NAG	c	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	c	2	4	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	f	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	f	2	4	-	4/6/23/26	0/1/1/1
5	NAG	h	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	h	2	5	-	0/6/23/26	0/1/1/1
5	FUC	h	3	5	-	-	0/1/1/1
4	NAG	l	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	l	2	4	-	4/6/23/26	0/1/1/1
4	NAG	n	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	n	2	4	-	0/6/23/26	0/1/1/1
4	NAG	p	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	p	2	4	-	1/6/23/26	0/1/1/1
6	NAG	t	1	6,1	-	4/6/23/26	0/1/1/1
6	NAG	t	2	6	-	2/6/23/26	0/1/1/1
6	BMA	t	3	6	-	1/2/19/22	0/1/1/1
6	MAN	t	4	6	-	1/2/19/22	0/1/1/1
6	MAN	t	5	6	-	2/2/19/22	0/1/1/1
4	NAG	y	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	y	2	4	-	2/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	t	5	MAN	O5-C1	2.92	1.48	1.43
6	t	5	MAN	C1-C2	2.87	1.59	1.52
5	Q	1	NAG	O5-C1	2.74	1.48	1.43
6	t	5	MAN	O5-C5	2.40	1.48	1.43

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	t	5	MAN	C1-O5-C5	7.25	121.90	112.19
5	Q	1	NAG	C1-O5-C5	5.46	119.50	112.19
4	5	1	NAG	C1-O5-C5	4.85	118.69	112.19
6	t	3	BMA	C1-O5-C5	2.68	115.78	112.19
4	l	2	NAG	C1-O5-C5	2.68	115.77	112.19
6	t	4	MAN	C1-O5-C5	2.53	115.58	112.19
6	t	5	MAN	O2-C2-C3	-2.30	105.38	110.15
4	K	1	NAG	C1-O5-C5	2.23	115.17	112.19
4	l	2	NAG	C3-C4-C5	2.23	114.27	110.23
4	p	1	NAG	C1-O5-C5	2.21	115.14	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	t	4	MAN	O2-C2-C3	-2.15	105.71	110.15

There are no chirality outliers.

All (61) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	7	2	NAG	O5-C5-C6-O6
4	f	2	NAG	C4-C5-C6-O6
4	p	1	NAG	O5-C5-C6-O6
4	7	2	NAG	C4-C5-C6-O6
5	Q	1	NAG	O5-C5-C6-O6
4	l	2	NAG	O5-C5-C6-O6
4	f	2	NAG	O5-C5-C6-O6
4	y	1	NAG	O5-C5-C6-O6
4	l	2	NAG	C4-C5-C6-O6
4	y	1	NAG	C4-C5-C6-O6
5	Q	1	NAG	C4-C5-C6-O6
4	n	1	NAG	O5-C5-C6-O6
6	t	5	MAN	O5-C5-C6-O6
4	E	2	NAG	C8-C7-N2-C2
4	E	2	NAG	O7-C7-N2-C2
4	U	2	NAG	C8-C7-N2-C2
4	U	2	NAG	O7-C7-N2-C2
4	X	1	NAG	C8-C7-N2-C2
4	X	1	NAG	O7-C7-N2-C2
4	X	2	NAG	C8-C7-N2-C2
4	X	2	NAG	O7-C7-N2-C2
4	f	2	NAG	C8-C7-N2-C2
4	f	2	NAG	O7-C7-N2-C2
4	l	2	NAG	C8-C7-N2-C2
4	l	2	NAG	O7-C7-N2-C2
4	7	1	NAG	C8-C7-N2-C2
4	7	1	NAG	O7-C7-N2-C2
5	Q	1	NAG	C8-C7-N2-C2
5	Q	1	NAG	O7-C7-N2-C2
5	1	1	NAG	C8-C7-N2-C2
5	1	1	NAG	O7-C7-N2-C2
6	t	1	NAG	C8-C7-N2-C2
6	t	1	NAG	O7-C7-N2-C2
4	E	1	NAG	O5-C5-C6-O6
4	X	1	NAG	O5-C5-C6-O6
4	X	2	NAG	C4-C5-C6-O6

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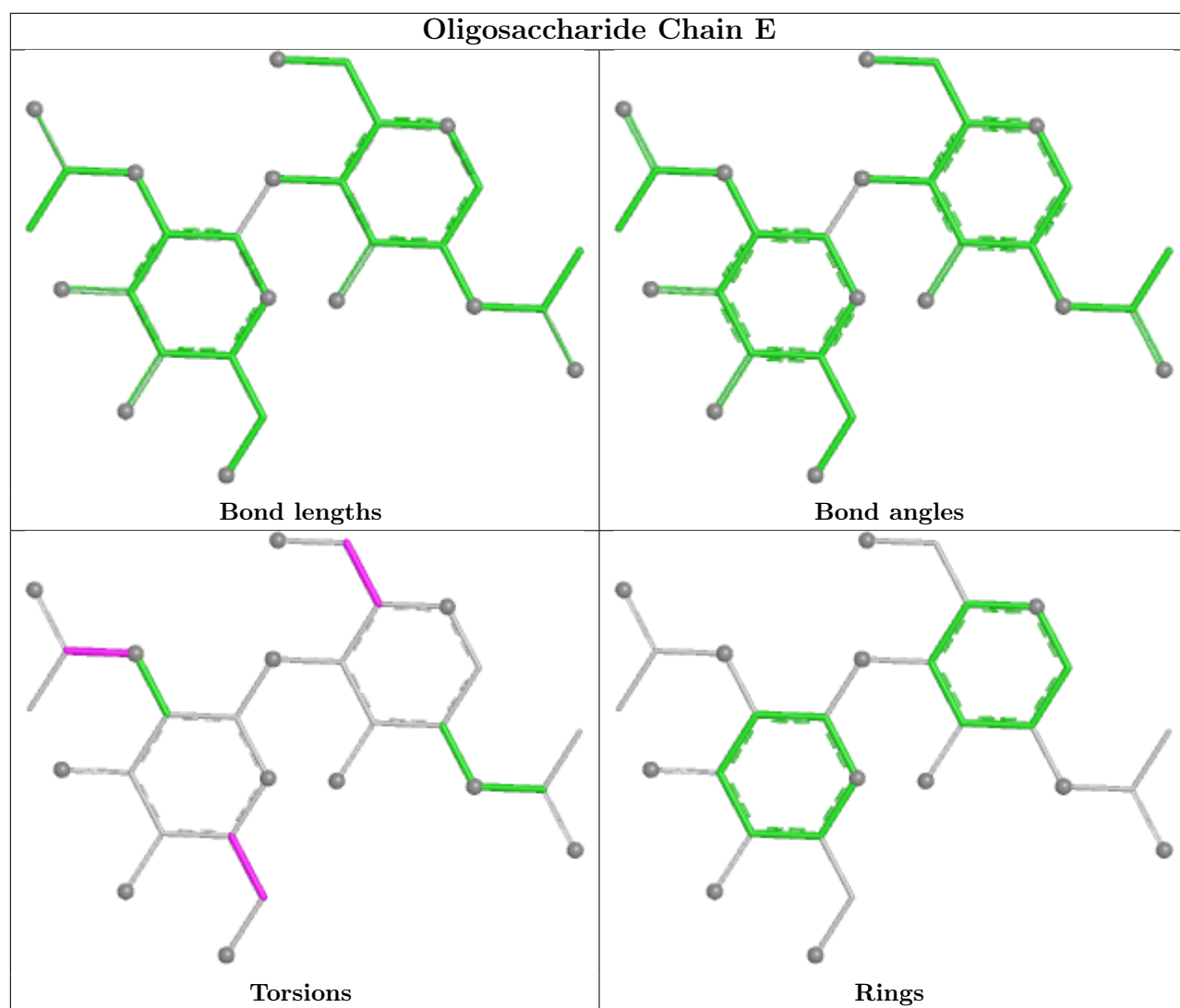
Mol	Chain	Res	Type	Atoms
4	E	2	NAG	O5-C5-C6-O6
4	n	1	NAG	C4-C5-C6-O6
4	E	1	NAG	C4-C5-C6-O6
4	X	1	NAG	C4-C5-C6-O6
4	p	1	NAG	C4-C5-C6-O6
4	X	2	NAG	O5-C5-C6-O6
6	t	1	NAG	C4-C5-C6-O6
4	c	2	NAG	C4-C5-C6-O6
4	c	2	NAG	O5-C5-C6-O6
6	t	3	BMA	O5-C5-C6-O6
4	5	1	NAG	O5-C5-C6-O6
6	t	4	MAN	O5-C5-C6-O6
6	t	1	NAG	O5-C5-C6-O6
4	5	2	NAG	C1-C2-N2-C7
4	E	2	NAG	C4-C5-C6-O6
6	t	5	MAN	C4-C5-C6-O6
5	Q	2	NAG	C4-C5-C6-O6
4	5	2	NAG	C3-C2-N2-C7
6	t	2	NAG	C3-C2-N2-C7
5	Q	2	NAG	O5-C5-C6-O6
6	t	2	NAG	C1-C2-N2-C7
4	y	2	NAG	C4-C5-C6-O6
4	y	2	NAG	O5-C5-C6-O6
4	U	2	NAG	C4-C5-C6-O6
4	p	2	NAG	C4-C5-C6-O6

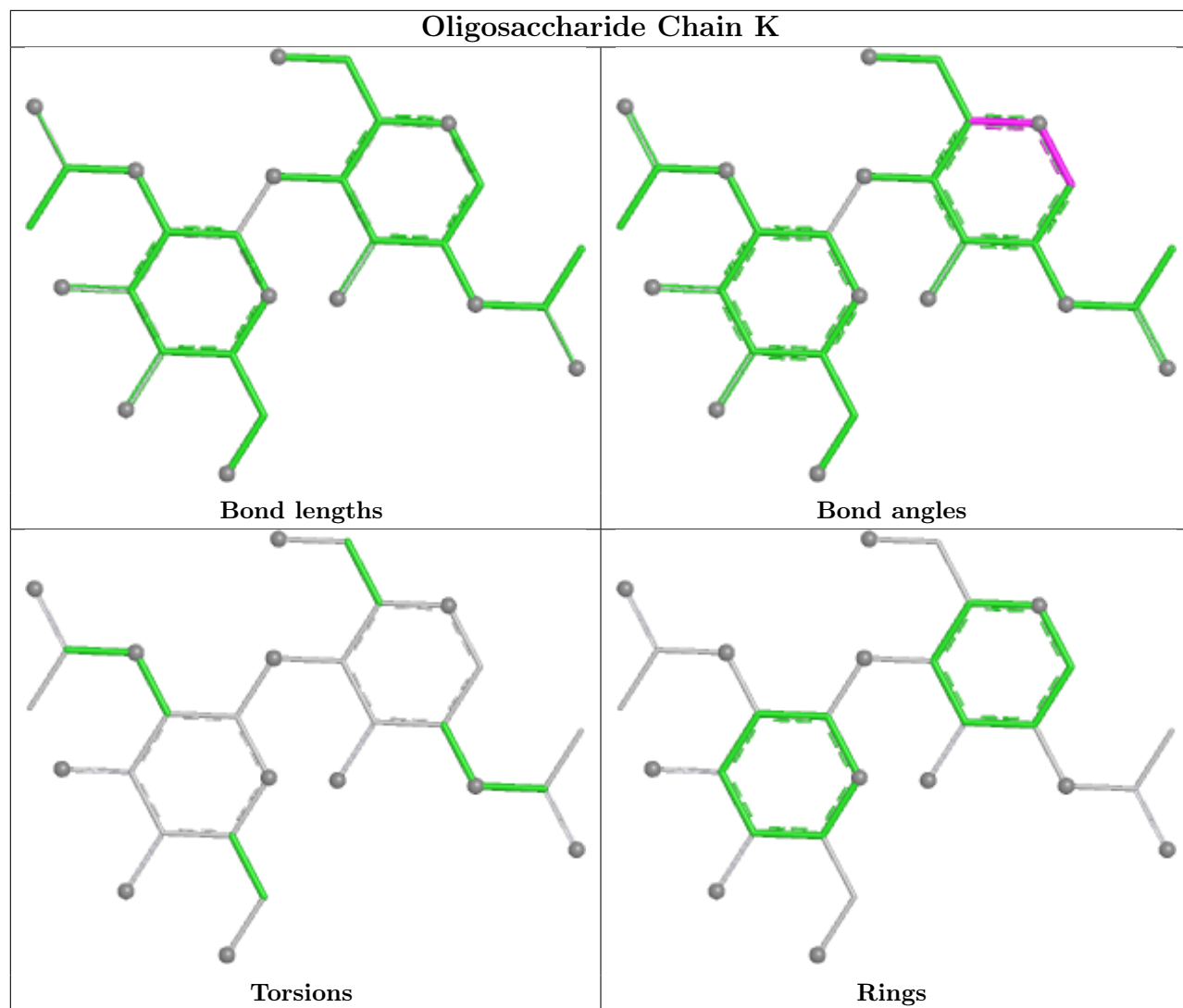
There are no ring outliers.

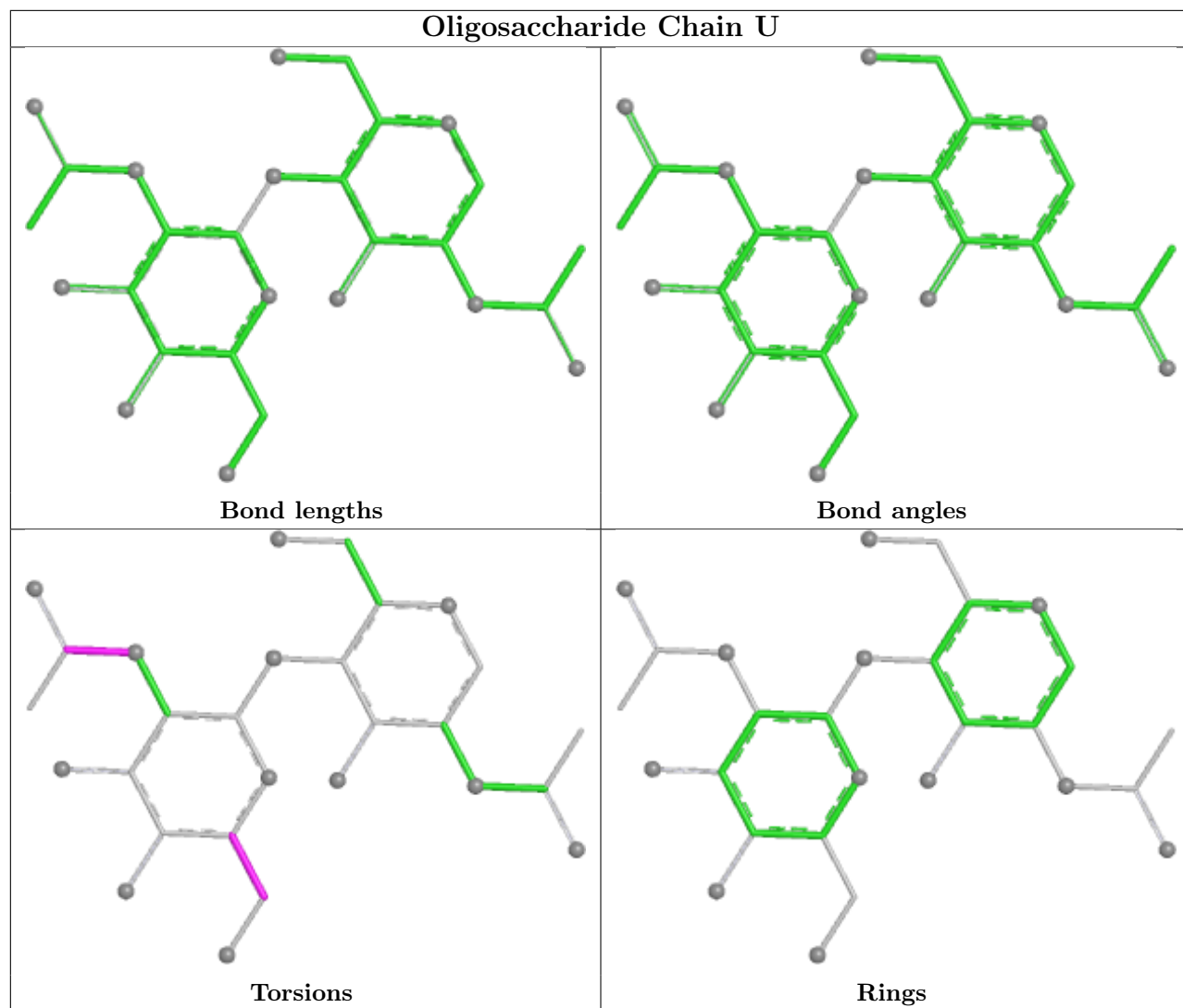
9 monomers are involved in 5 short contacts:

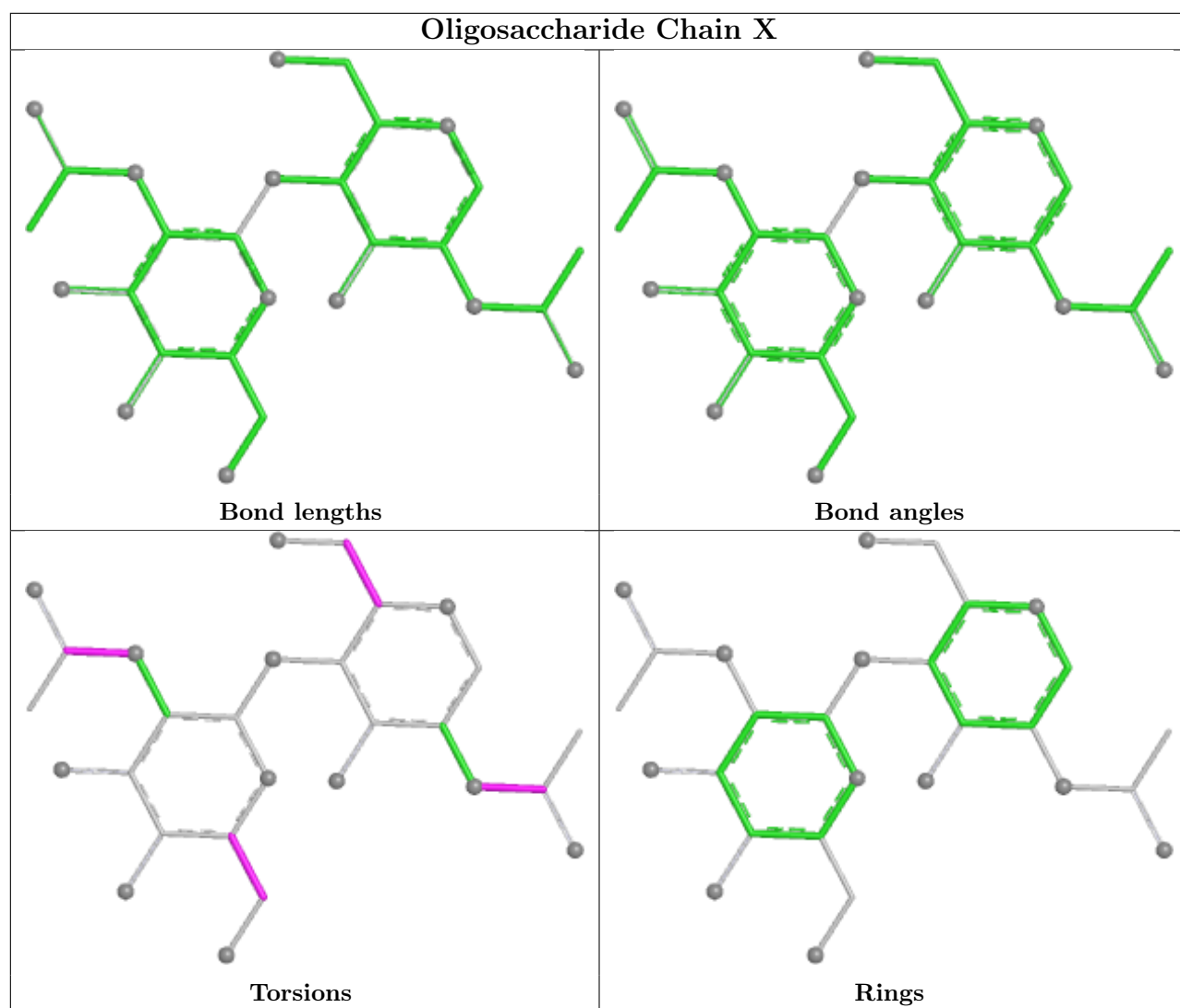
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	t	2	NAG	1	0
4	f	2	NAG	1	0
6	t	1	NAG	1	0
4	f	1	NAG	1	0
4	7	2	NAG	1	0
4	5	1	NAG	1	0
4	y	2	NAG	1	0
4	7	1	NAG	1	0
4	5	2	NAG	1	0

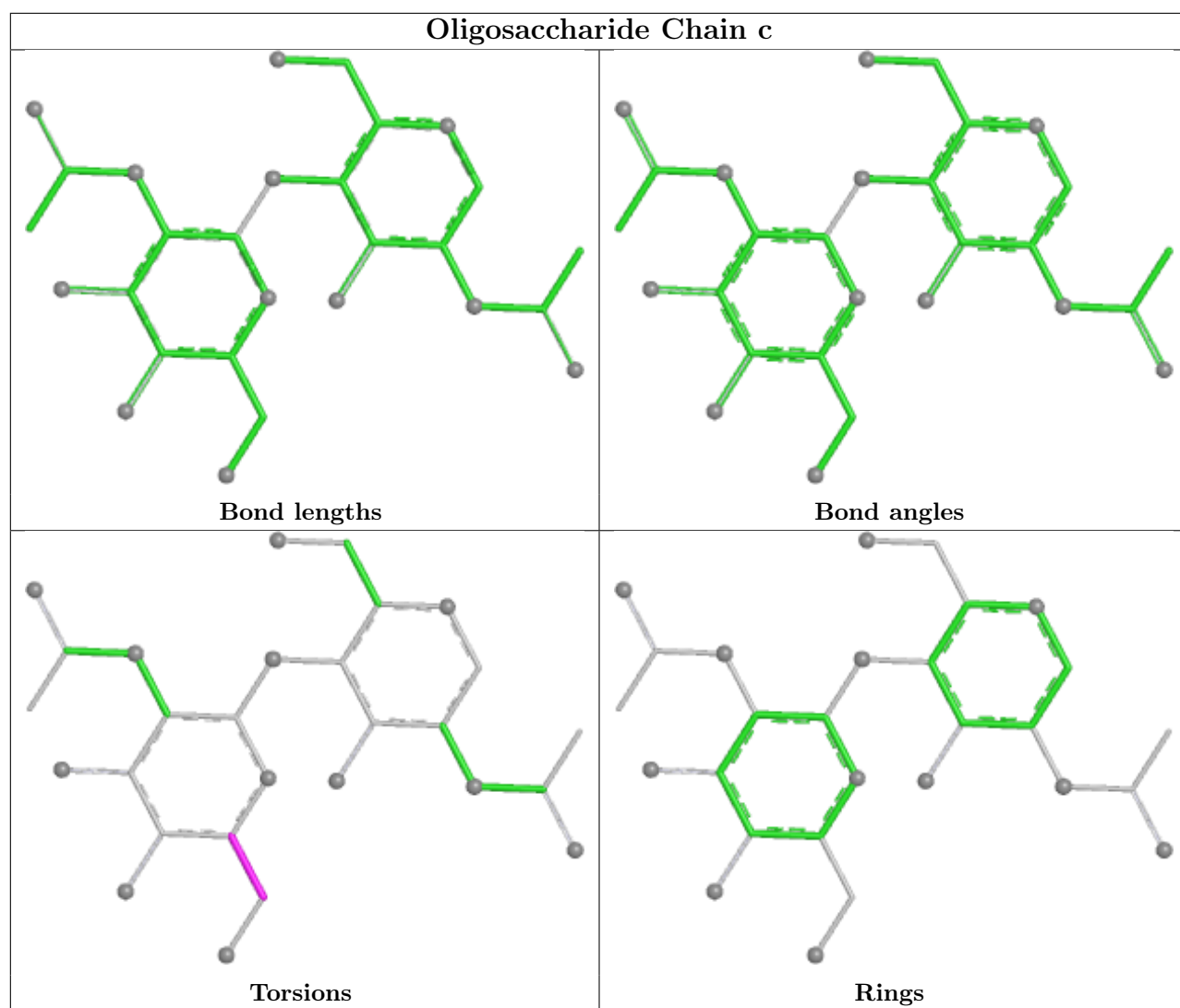
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

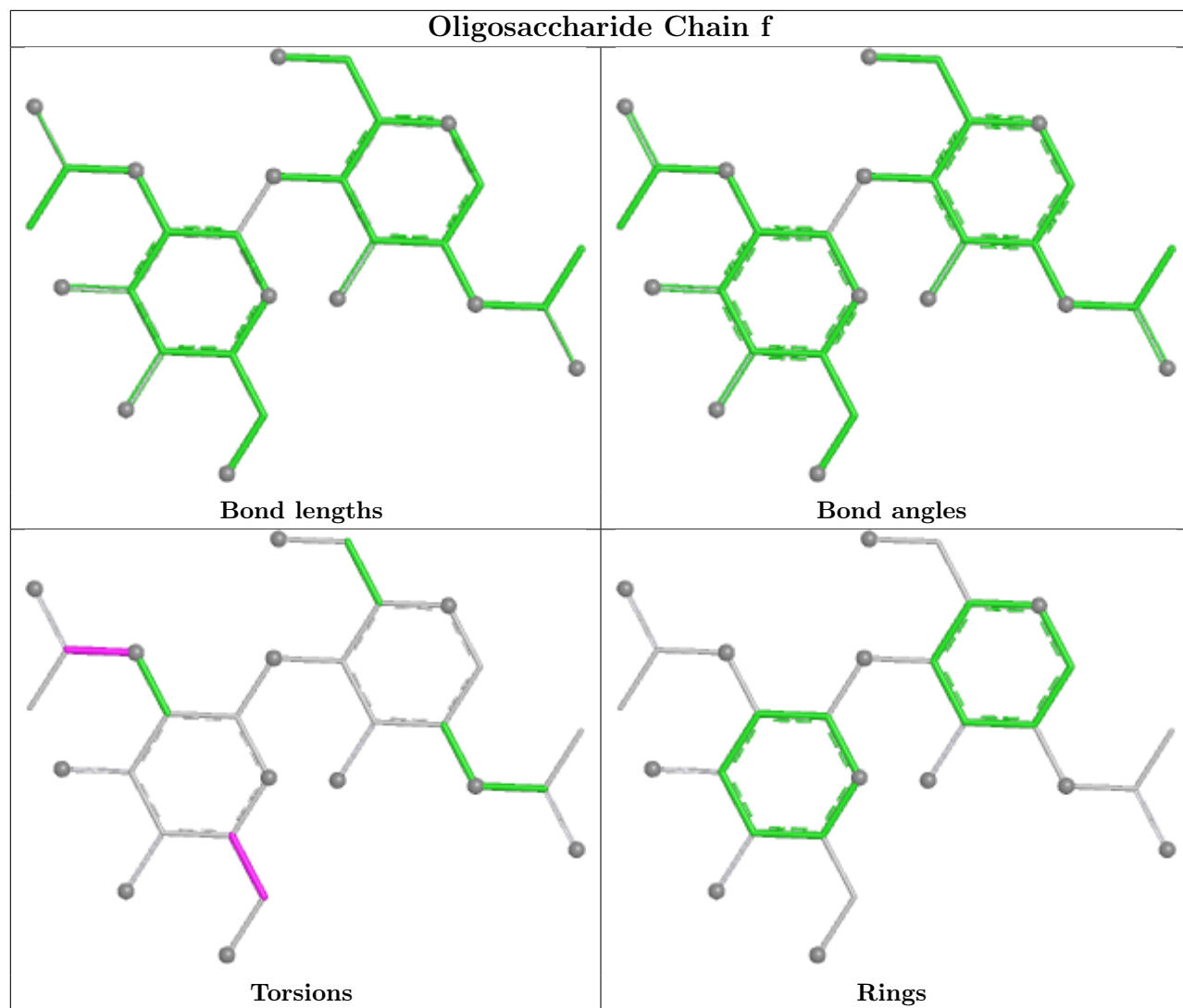


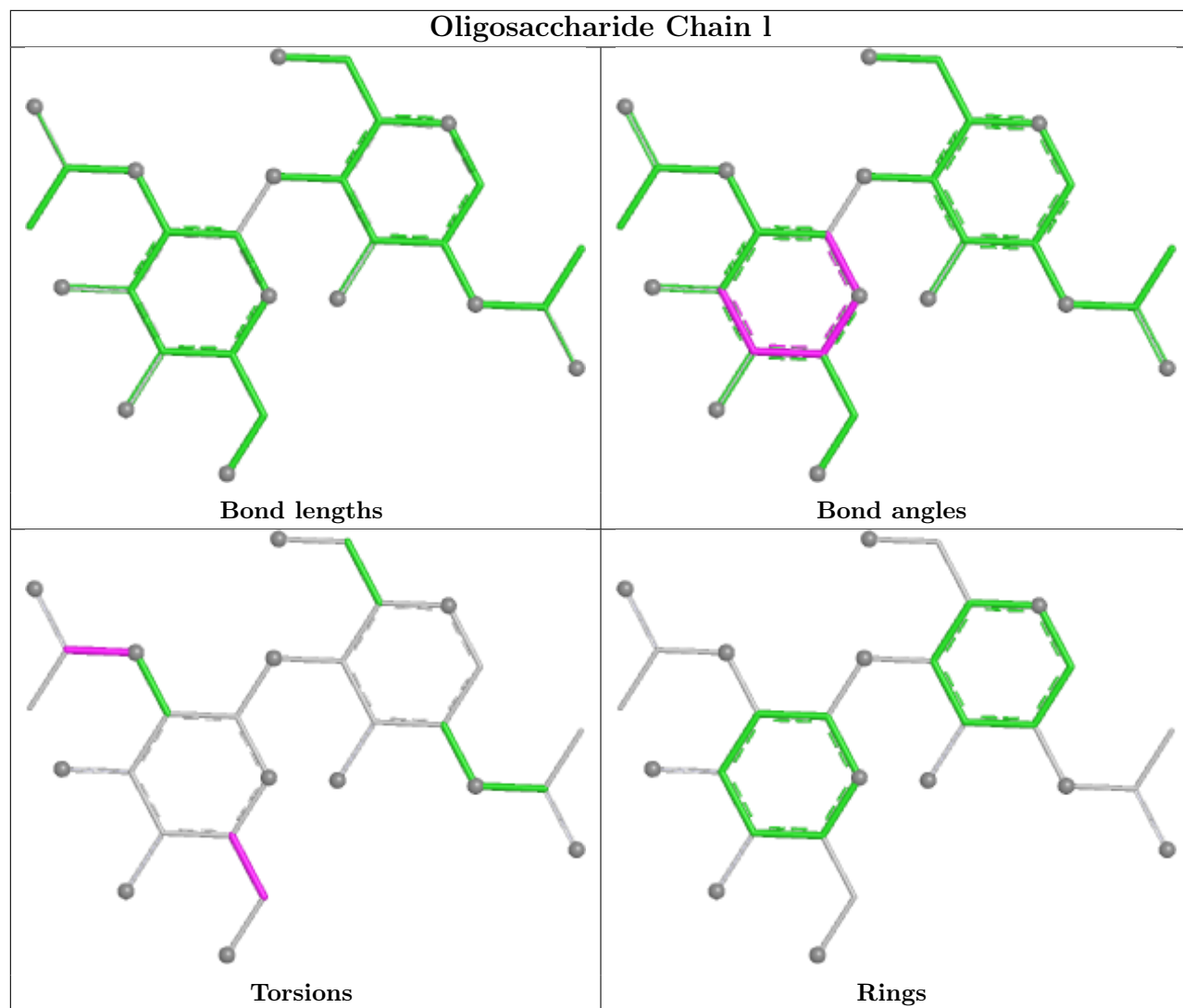


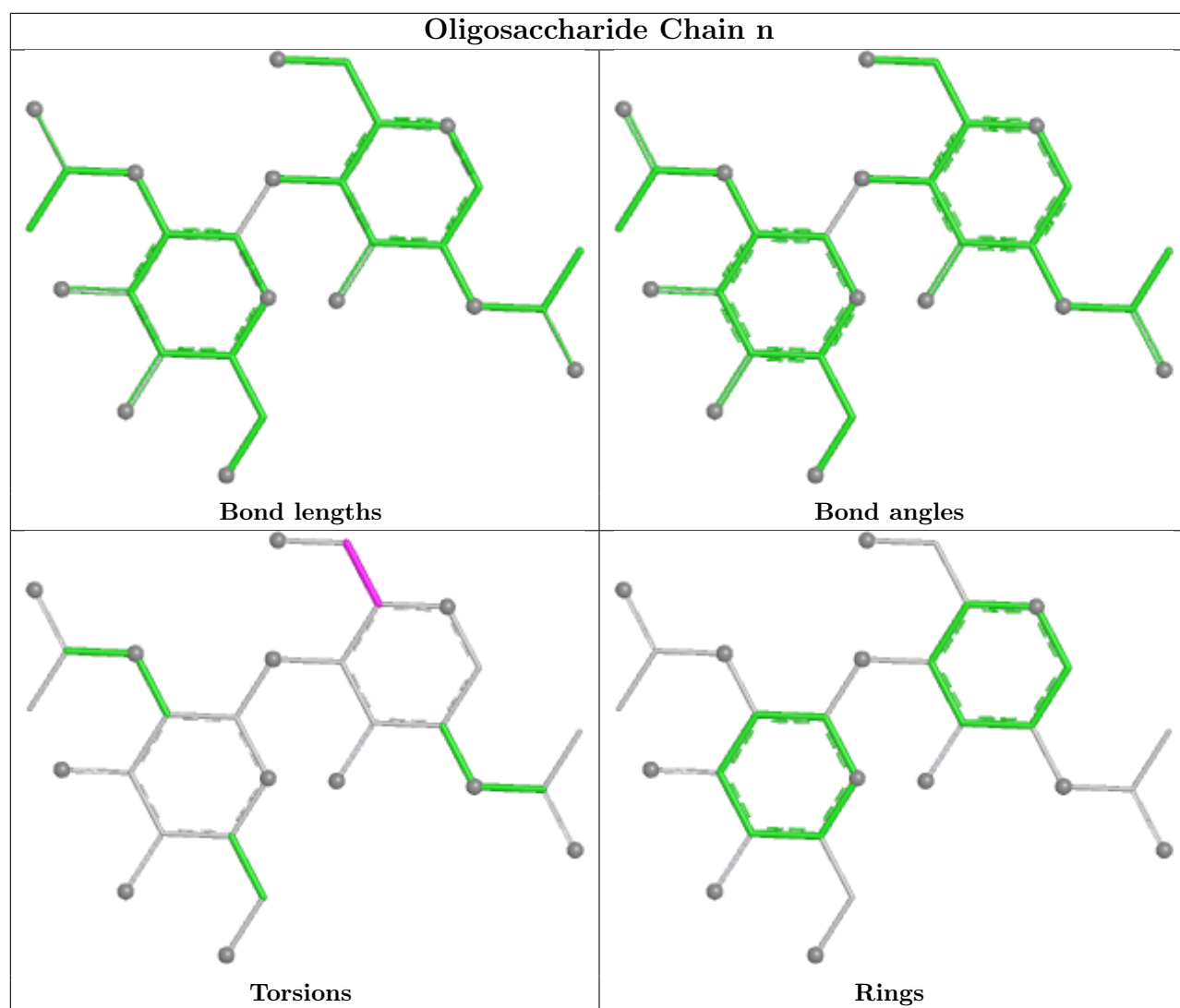


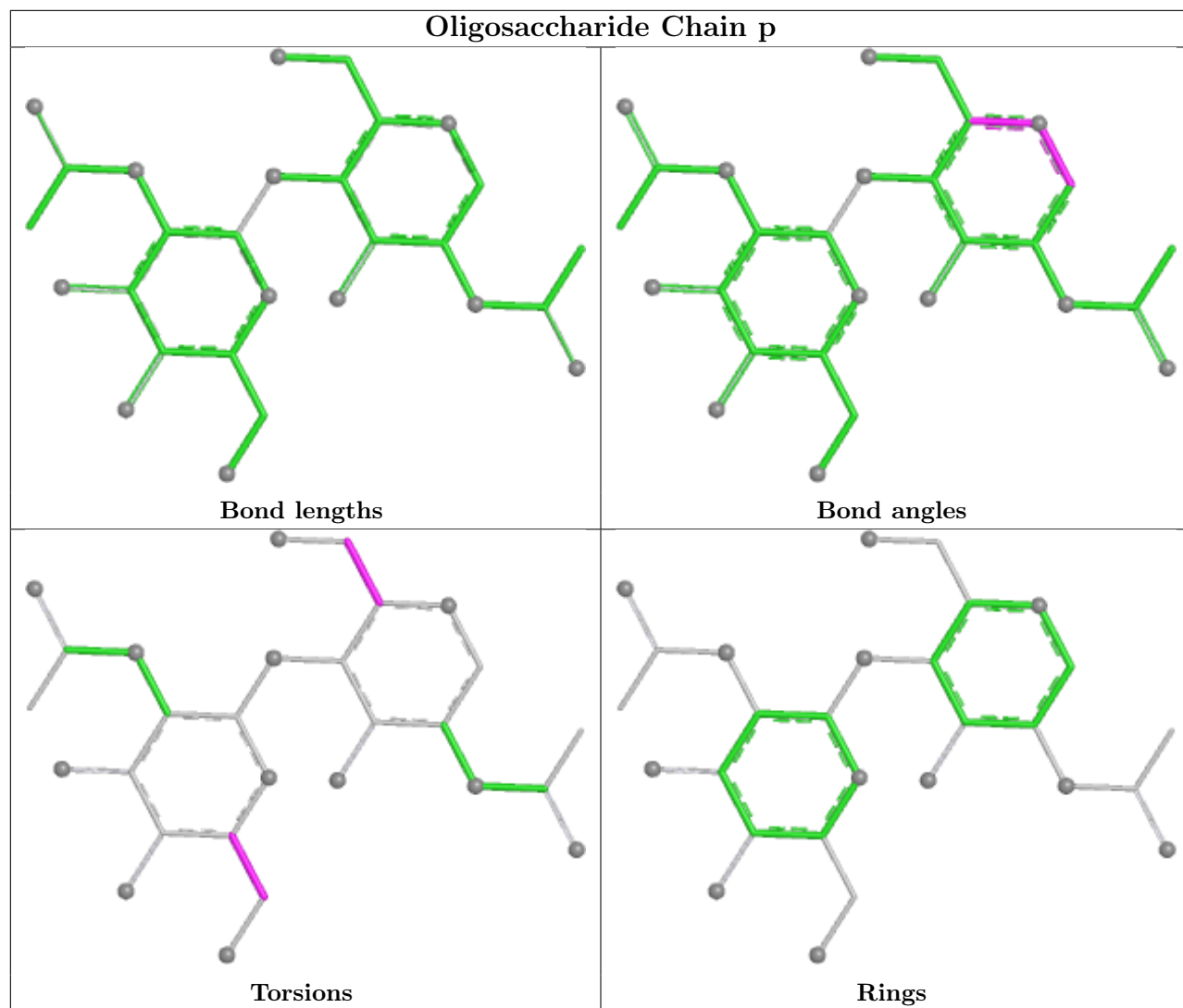


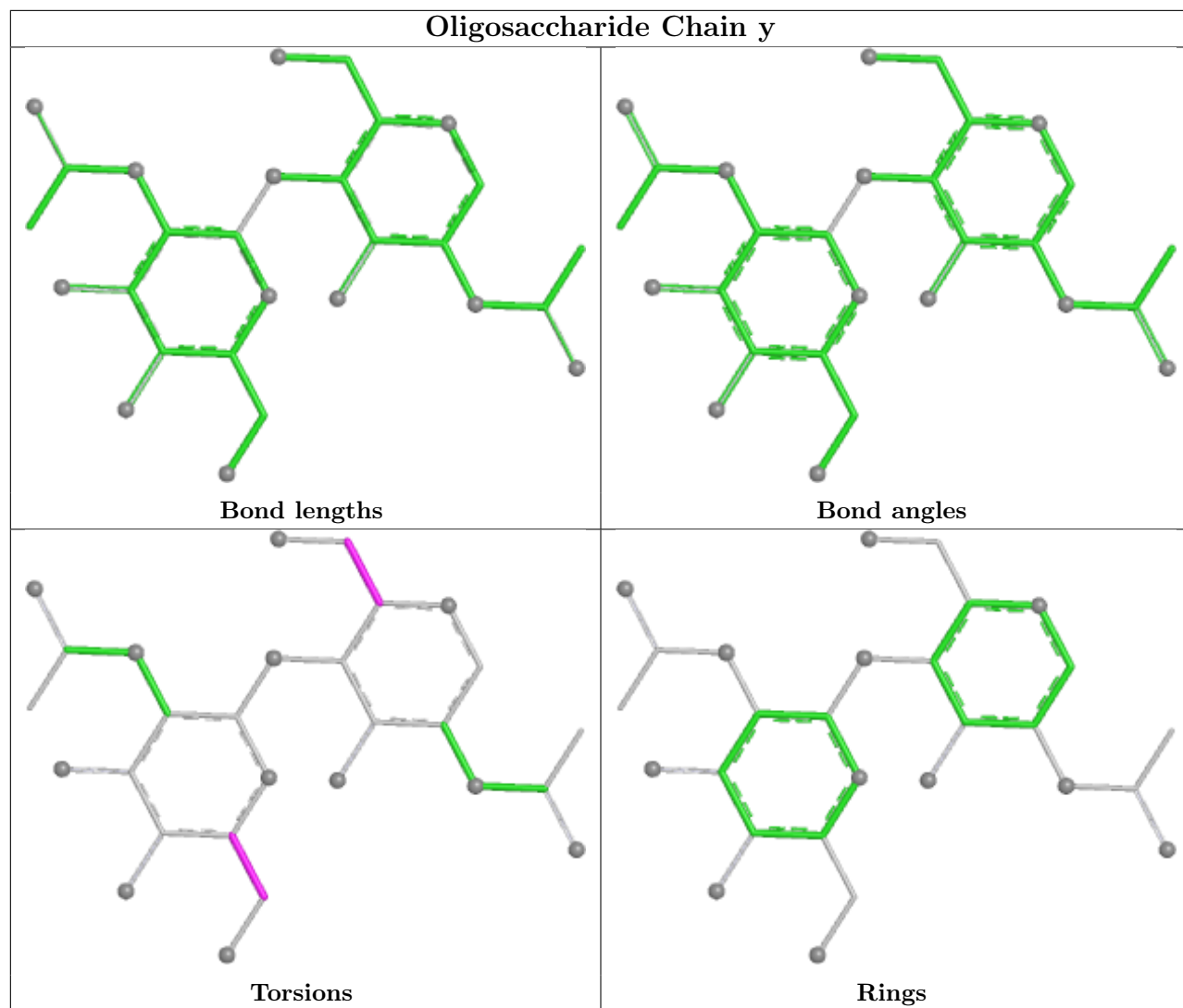


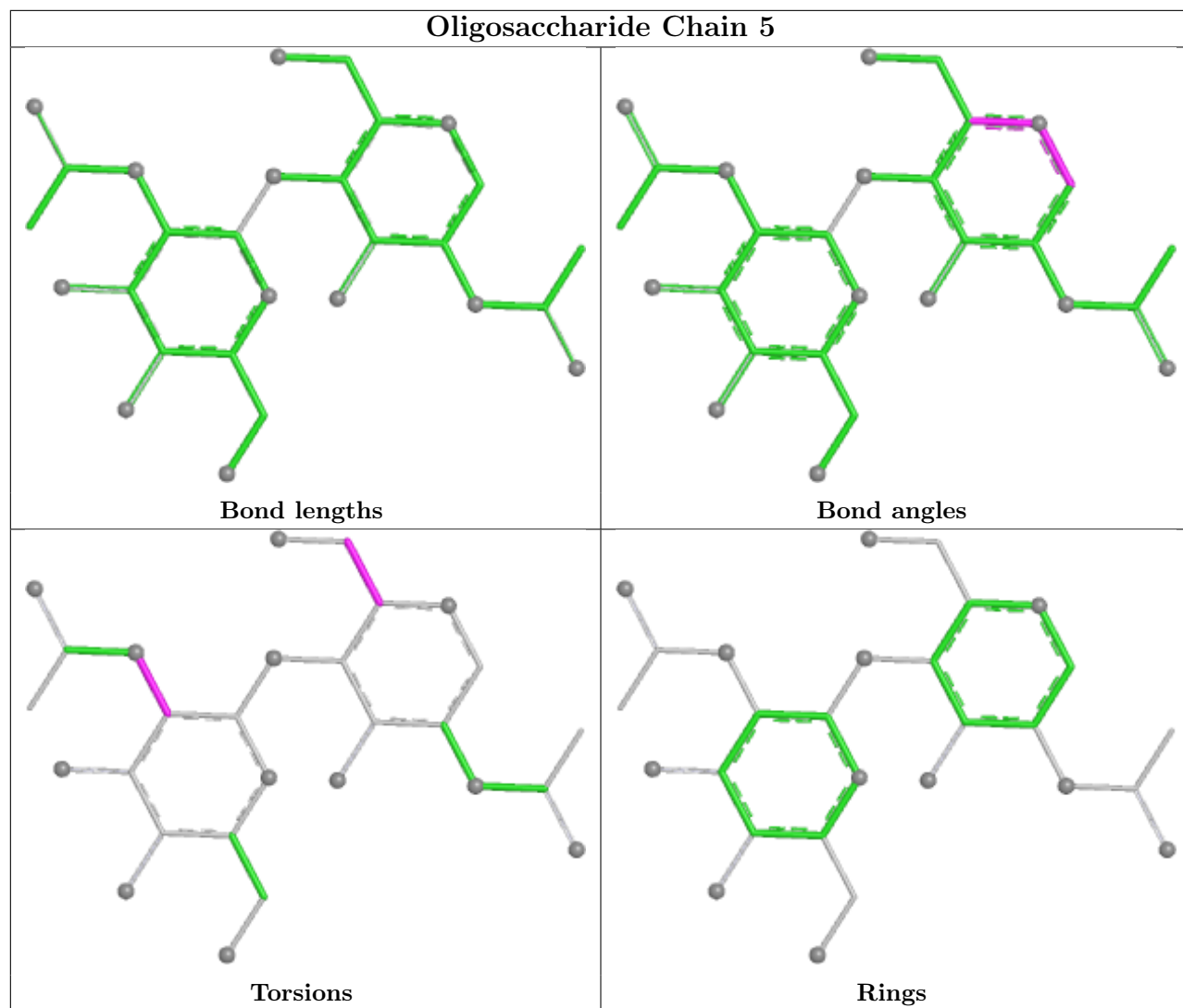


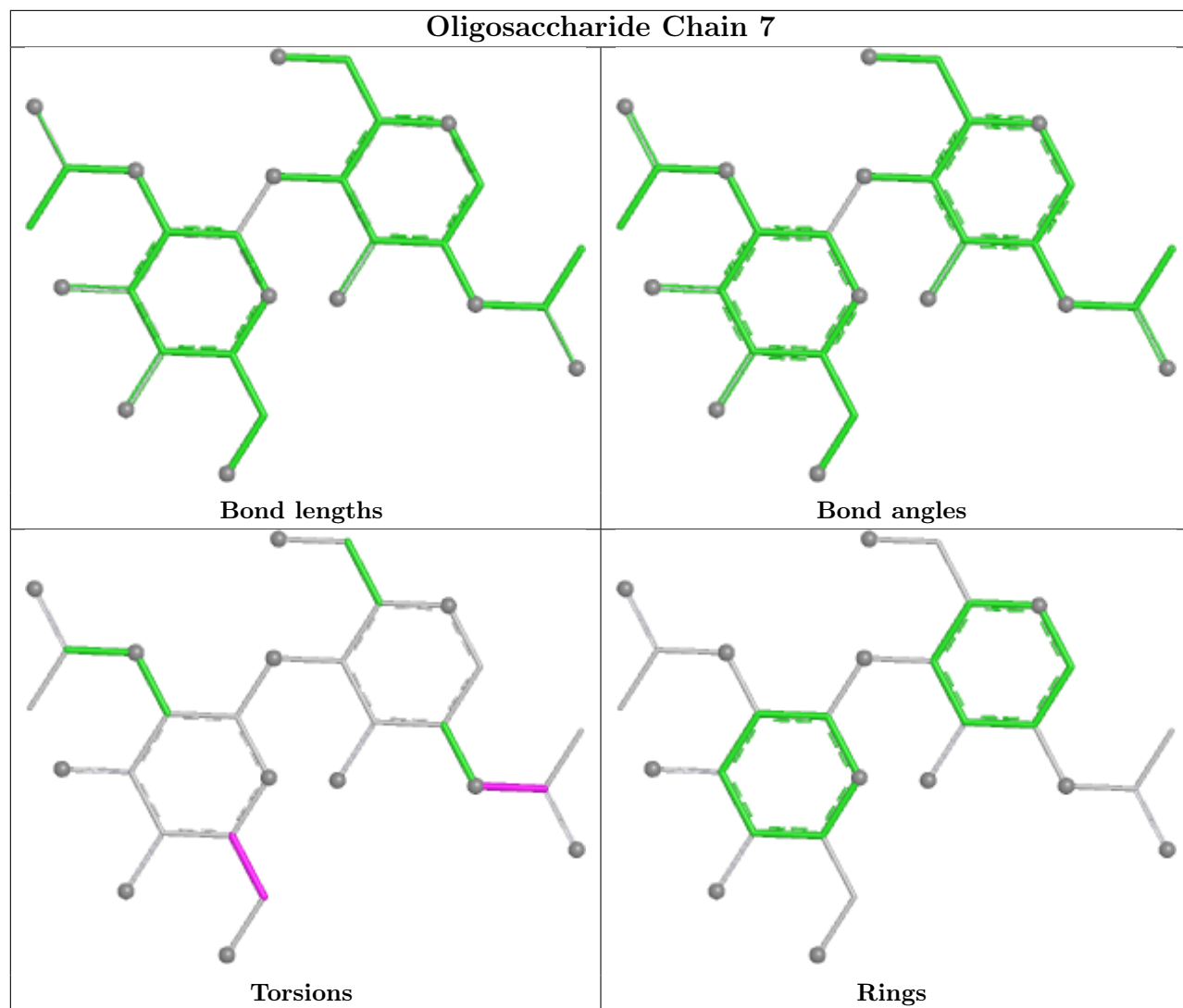


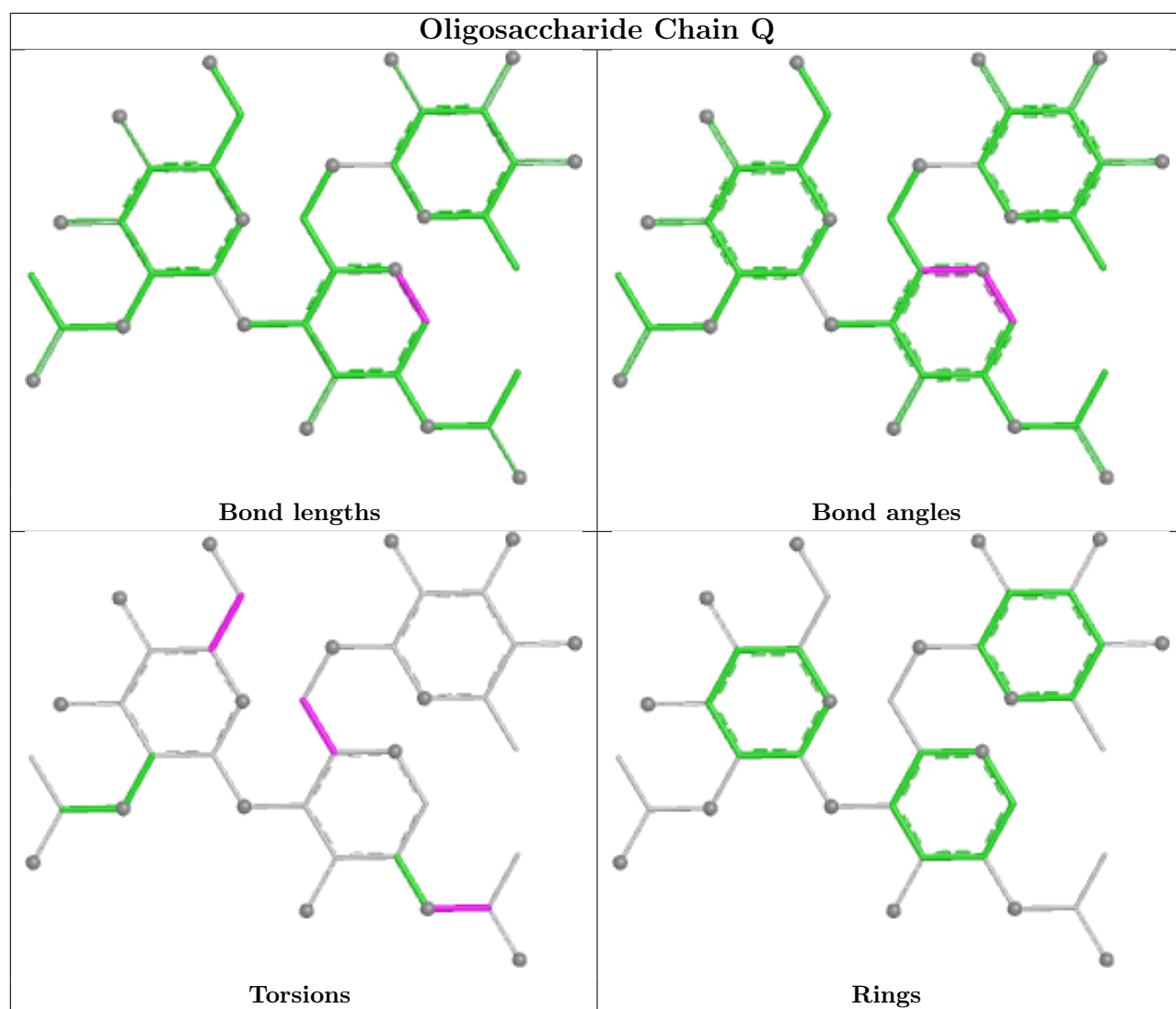


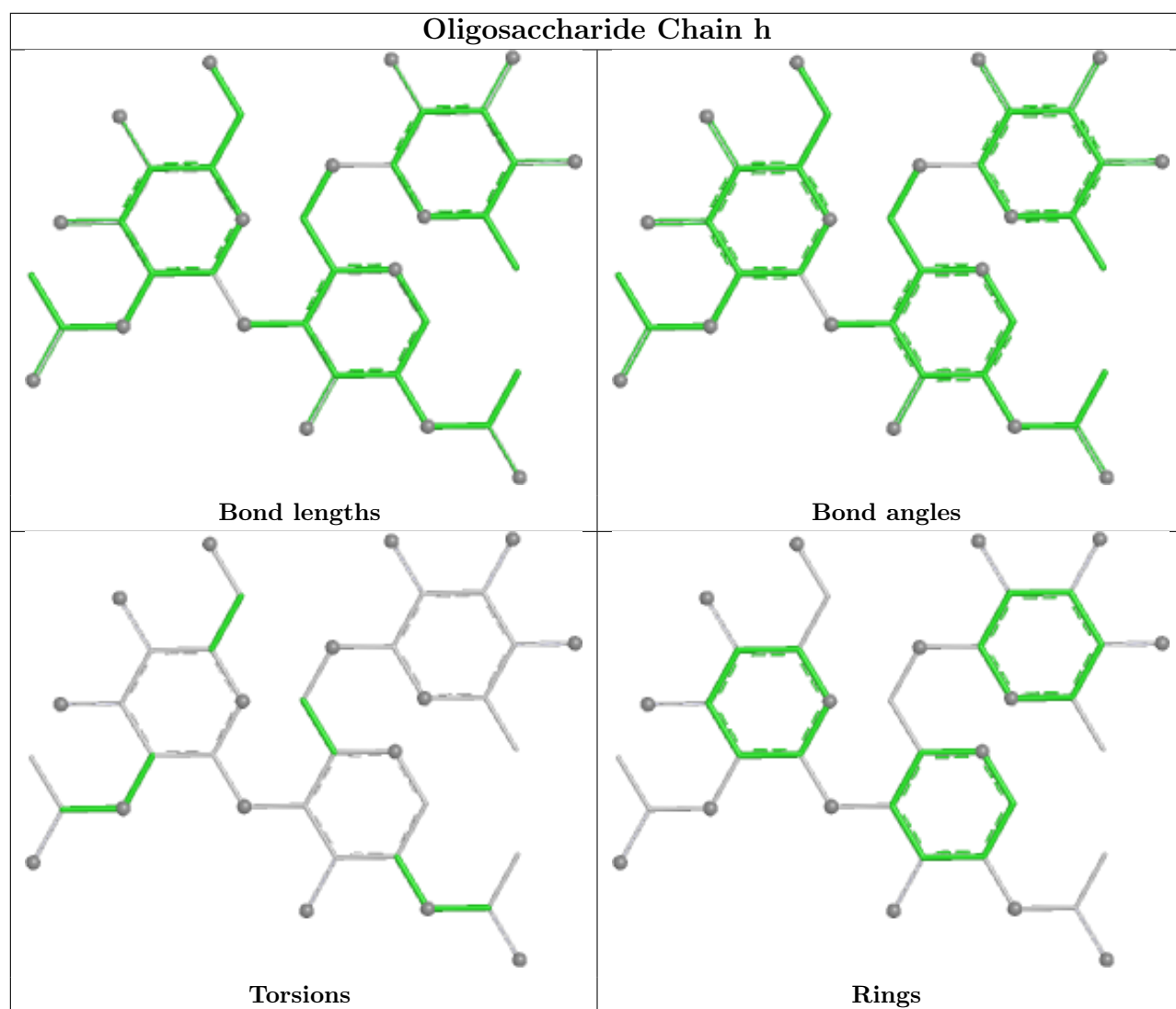


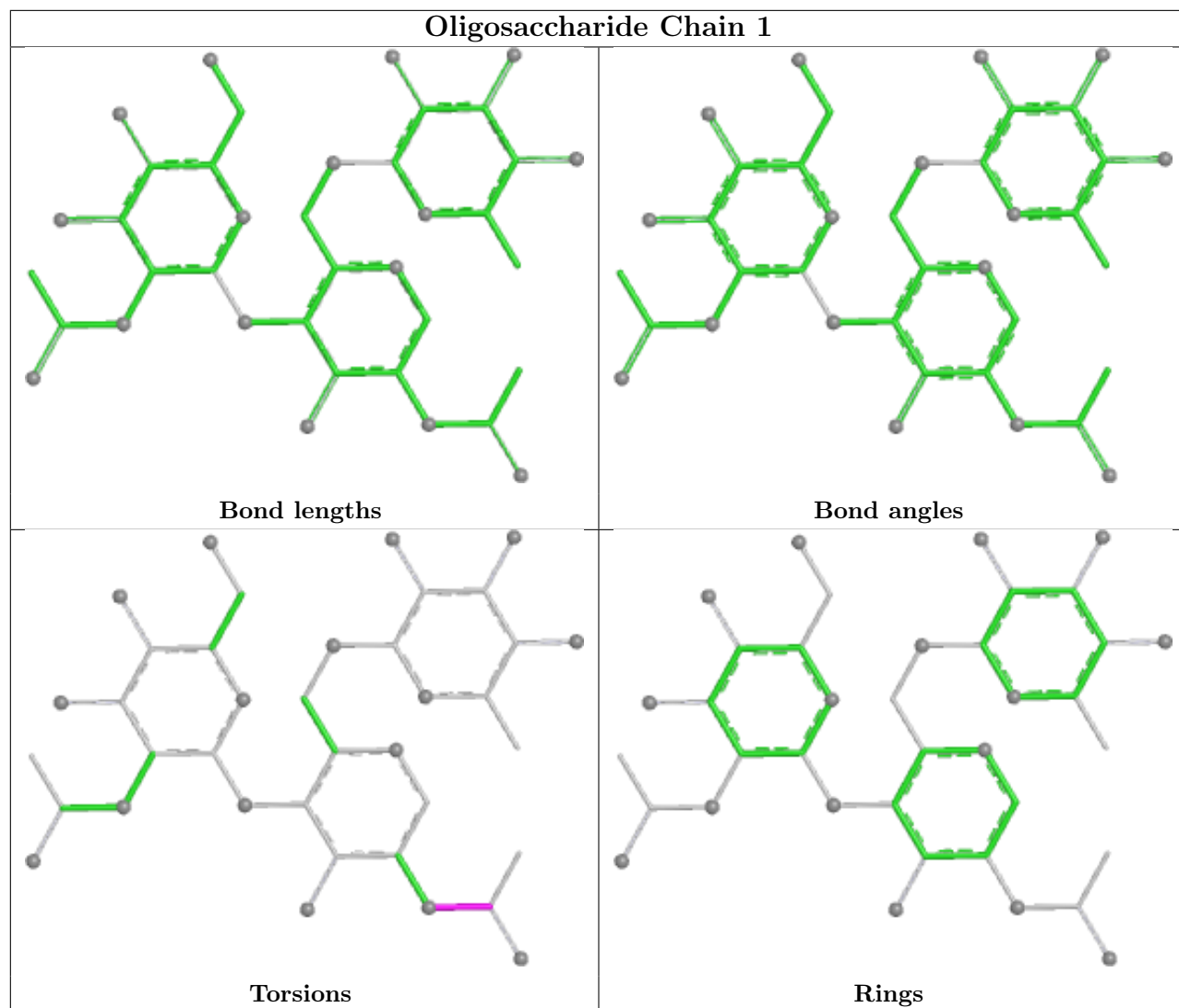


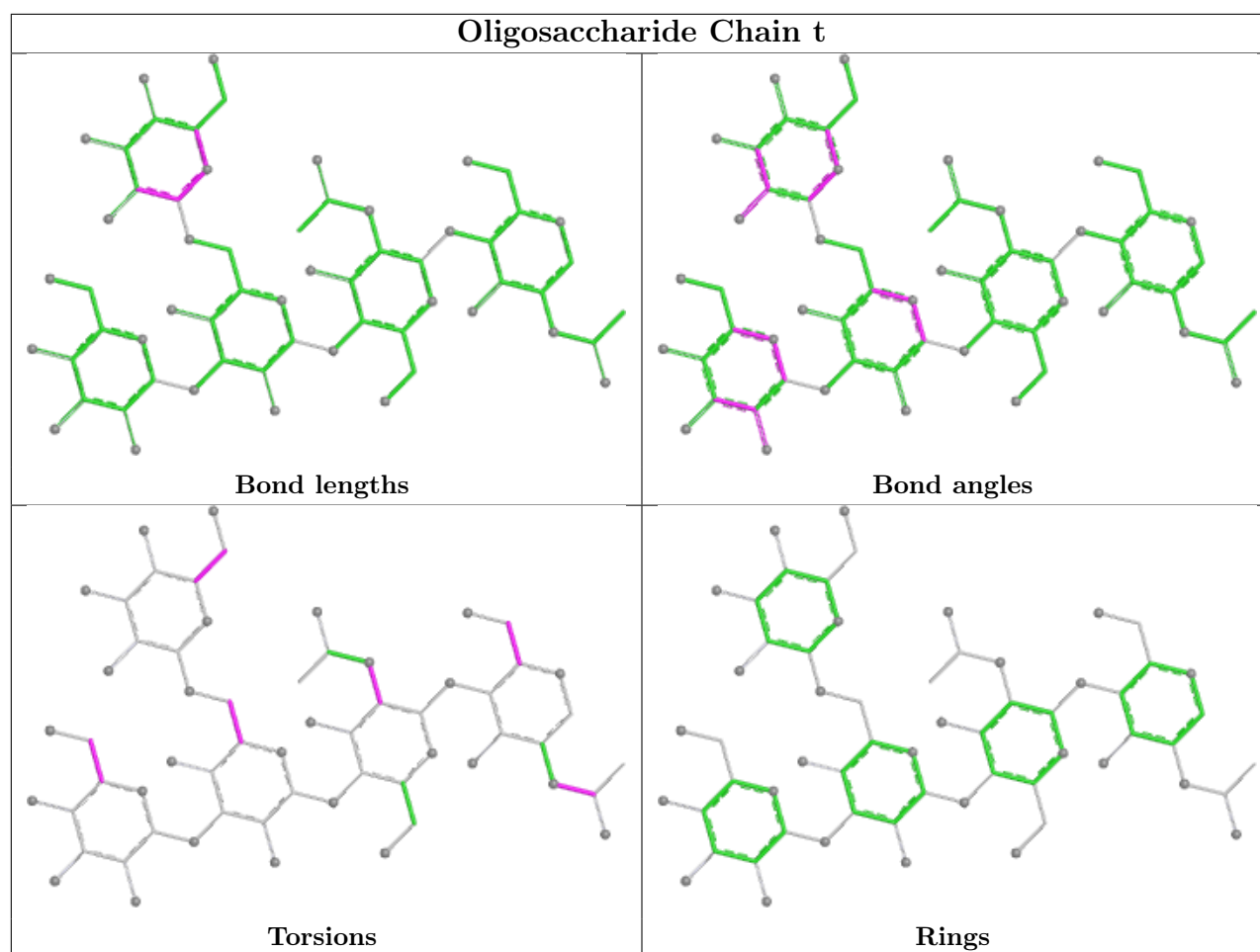












5.6 Ligand geometry [i](#)

18 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	$\# Z > 2$
7	NAG	A	1309	1	14,14,15	0.25	0	17,19,21	0.48	0
7	NAG	C	1303	1	14,14,15	0.19	0	17,19,21	0.43	0
7	NAG	A	1303	1	14,14,15	0.42	0	17,19,21	0.48	0
7	NAG	A	1301	1	14,14,15	0.28	0	17,19,21	0.55	0
7	NAG	C	1302	1	14,14,15	0.20	0	17,19,21	0.37	0
7	NAG	C	1304	1	14,14,15	0.25	0	17,19,21	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NAG	A	1305	1	14,14,15	0.22	0	17,19,21	0.43	0
7	NAG	B	1305	1	14,14,15	0.19	0	17,19,21	0.44	0
7	NAG	A	1307	1	14,14,15	0.19	0	17,19,21	0.45	0
7	NAG	B	1302	1	14,14,15	0.20	0	17,19,21	0.46	0
7	NAG	B	1303	1	14,14,15	0.23	0	17,19,21	0.49	0
7	NAG	B	1304	1	14,14,15	0.24	0	17,19,21	0.55	0
7	NAG	A	1302	1	14,14,15	0.39	0	17,19,21	0.55	0
7	NAG	A	1308	1	14,14,15	0.25	0	17,19,21	0.46	0
7	NAG	C	1301	1	14,14,15	0.20	0	17,19,21	0.46	0
7	NAG	A	1304	1	14,14,15	0.27	0	17,19,21	0.35	0
7	NAG	A	1306	1	14,14,15	0.21	0	17,19,21	0.42	0
7	NAG	B	1301	1	14,14,15	0.21	0	17,19,21	0.47	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
7	NAG	A	1309	1	-	0/6/23/26	0/1/1/1
7	NAG	C	1303	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1303	1	-	4/6/23/26	0/1/1/1
7	NAG	A	1301	1	-	2/6/23/26	0/1/1/1
7	NAG	C	1302	1	-	1/6/23/26	0/1/1/1
7	NAG	C	1304	1	-	1/6/23/26	0/1/1/1
7	NAG	A	1305	1	-	1/6/23/26	0/1/1/1
7	NAG	B	1305	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1307	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1302	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1303	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1304	1	-	4/6/23/26	0/1/1/1
7	NAG	A	1302	1	-	3/6/23/26	0/1/1/1
7	NAG	A	1308	1	-	2/6/23/26	0/1/1/1
7	NAG	C	1301	1	-	3/6/23/26	0/1/1/1
7	NAG	A	1304	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1306	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1301	1	-	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 (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1302	NAG	C1-C2-N2-C7
7	A	1302	NAG	C8-C7-N2-C2
7	A	1302	NAG	O7-C7-N2-C2
7	B	1305	NAG	C4-C5-C6-O6
7	B	1303	NAG	O5-C5-C6-O6
7	B	1305	NAG	O5-C5-C6-O6
7	B	1302	NAG	O5-C5-C6-O6
7	A	1308	NAG	O5-C5-C6-O6
7	A	1303	NAG	O5-C5-C6-O6
7	A	1304	NAG	O5-C5-C6-O6
7	B	1303	NAG	C4-C5-C6-O6
7	A	1304	NAG	C4-C5-C6-O6
7	A	1308	NAG	C4-C5-C6-O6
7	C	1303	NAG	C4-C5-C6-O6
7	A	1307	NAG	O5-C5-C6-O6
7	B	1302	NAG	C4-C5-C6-O6
7	A	1303	NAG	C4-C5-C6-O6
7	C	1301	NAG	C8-C7-N2-C2
7	C	1301	NAG	O7-C7-N2-C2
7	A	1306	NAG	O5-C5-C6-O6
7	C	1302	NAG	O5-C5-C6-O6
7	A	1307	NAG	C4-C5-C6-O6
7	C	1303	NAG	O5-C5-C6-O6
7	A	1305	NAG	O5-C5-C6-O6
7	C	1304	NAG	O5-C5-C6-O6
7	B	1304	NAG	C1-C2-N2-C7
7	A	1301	NAG	C3-C2-N2-C7
7	A	1303	NAG	C3-C2-N2-C7
7	B	1304	NAG	C3-C2-N2-C7
7	C	1301	NAG	O5-C5-C6-O6
7	A	1301	NAG	C1-C2-N2-C7
7	A	1303	NAG	C1-C2-N2-C7
7	B	1304	NAG	C4-C5-C6-O6
7	A	1306	NAG	C4-C5-C6-O6
7	B	1304	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1304	NAG	1	0

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-40272. These allow visual inspection of the internal detail of the map and identification of artifacts.

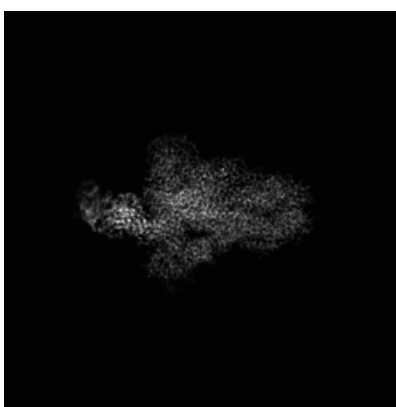
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

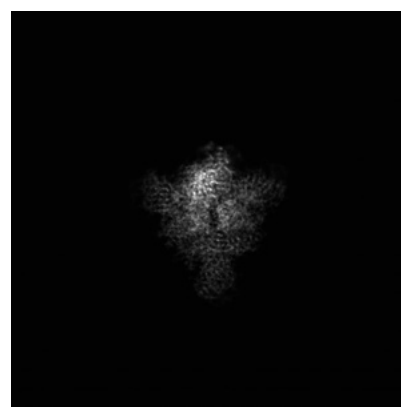
6.1.1 Primary map



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: 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](#)

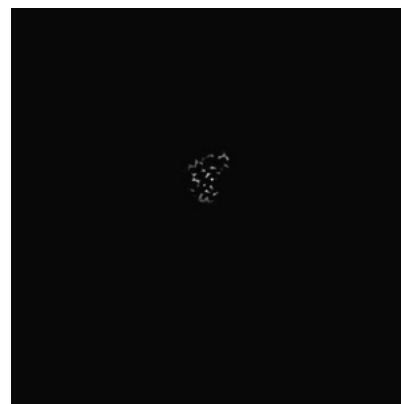
6.3.1 Primary map



X Index: 196



Y Index: 227

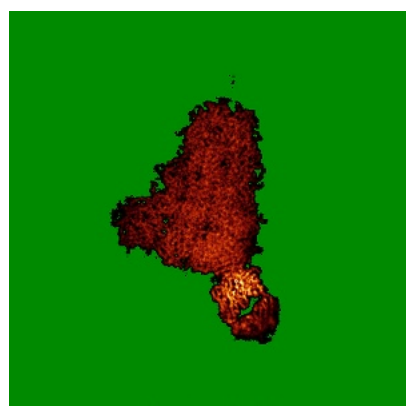


Z Index: 123

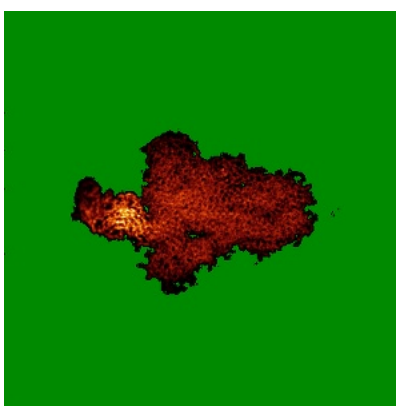
The images above show the largest variance slices of the map in three orthogonal directions.

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

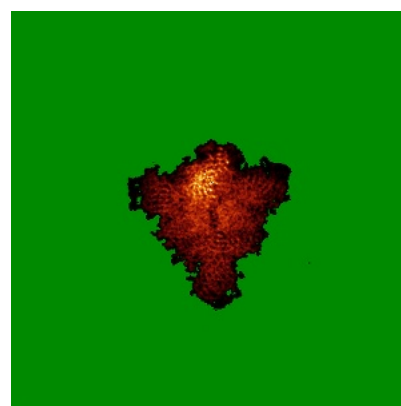
6.4.1 Primary map



X



Y

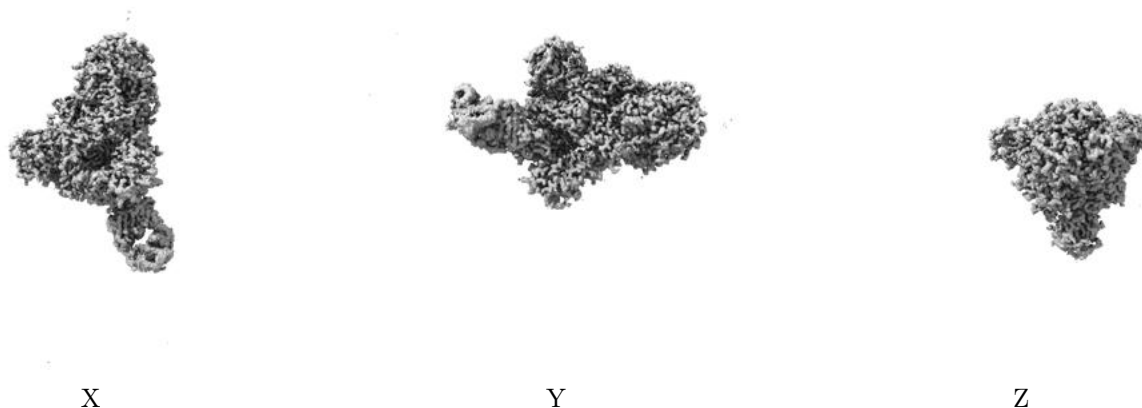


Z

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 [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 1.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

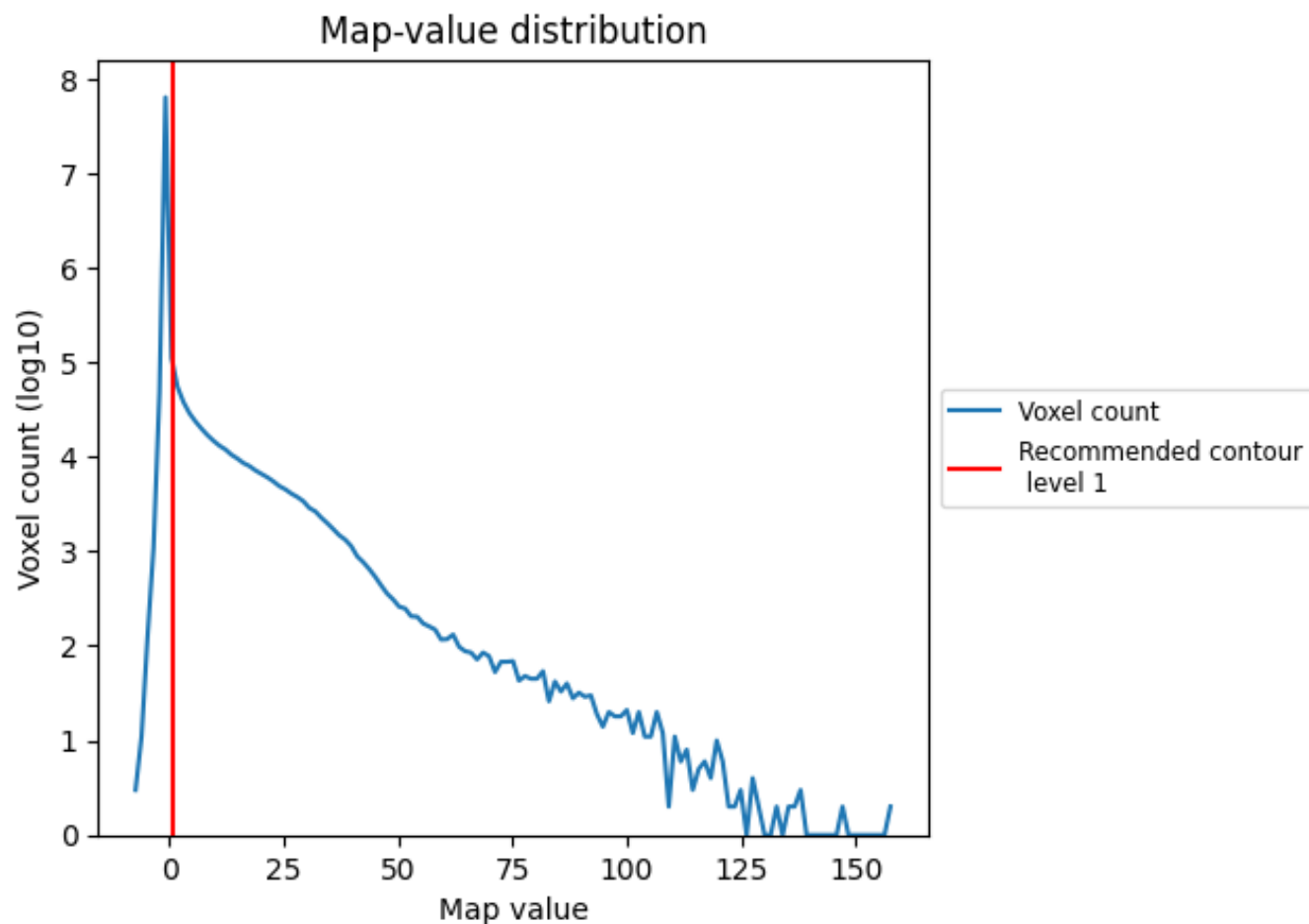
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

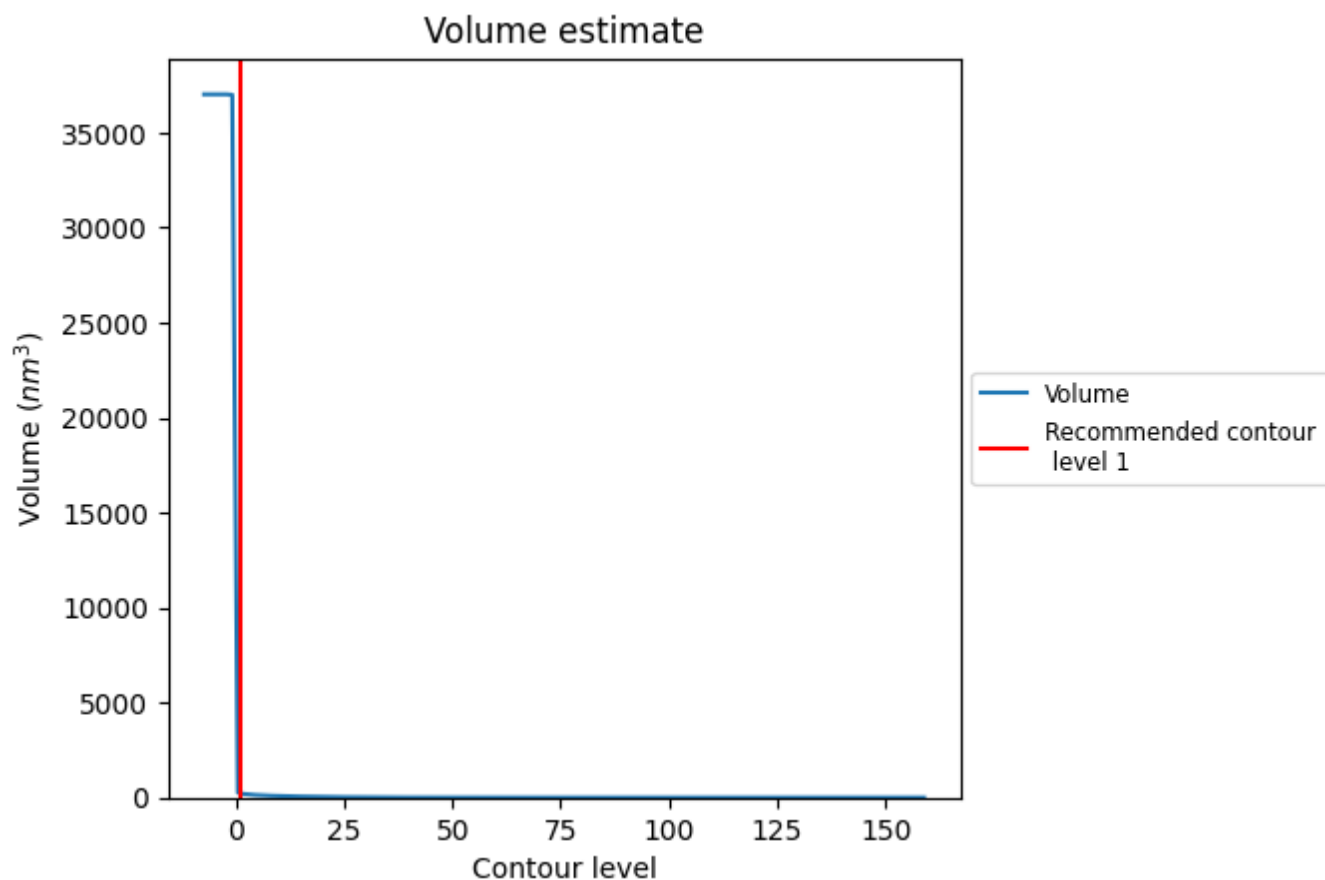
This section contains the results of statistical analysis of the map.

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.

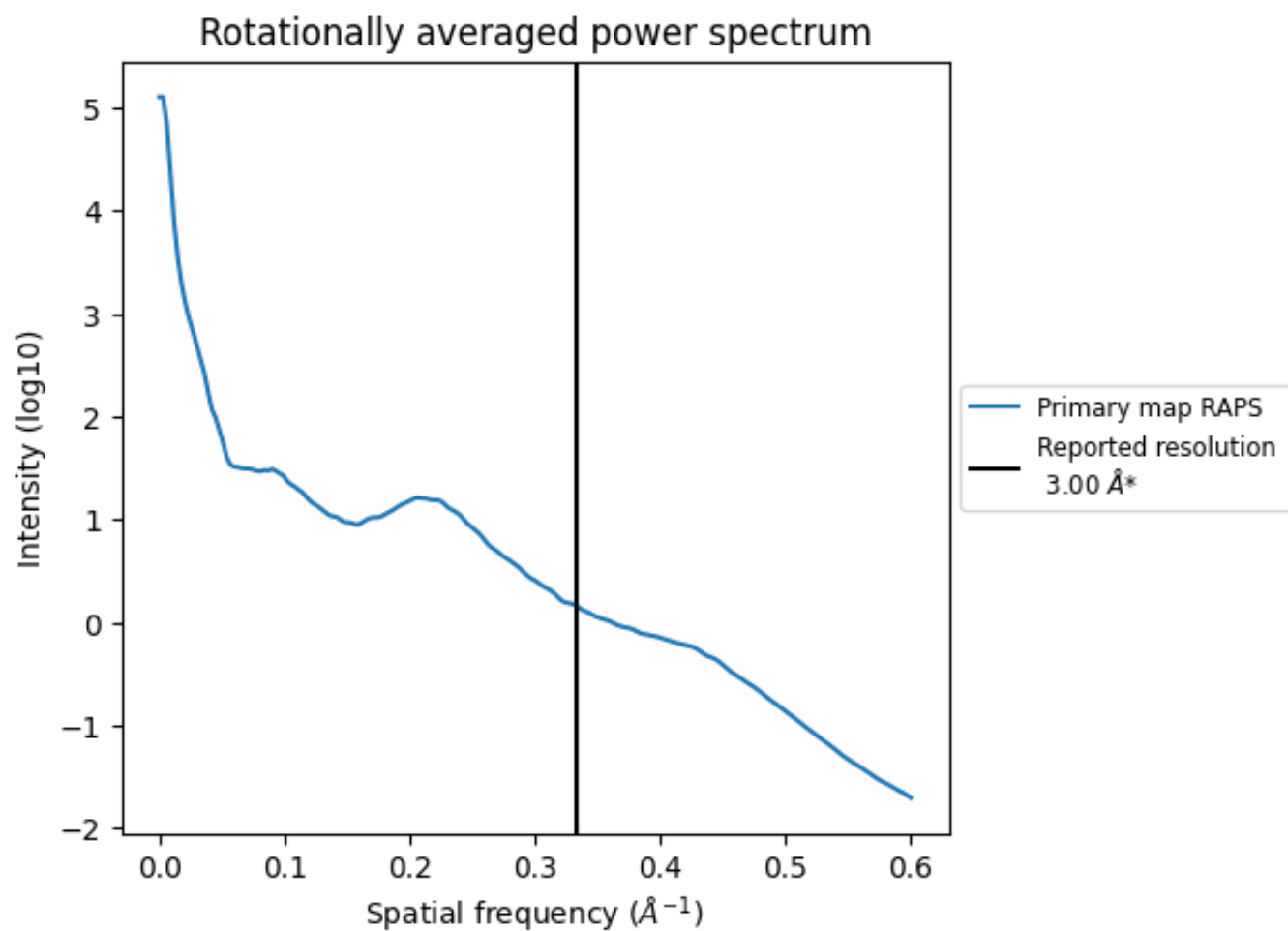
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 233 nm³; this corresponds to an approximate mass of 210 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 ⓘ

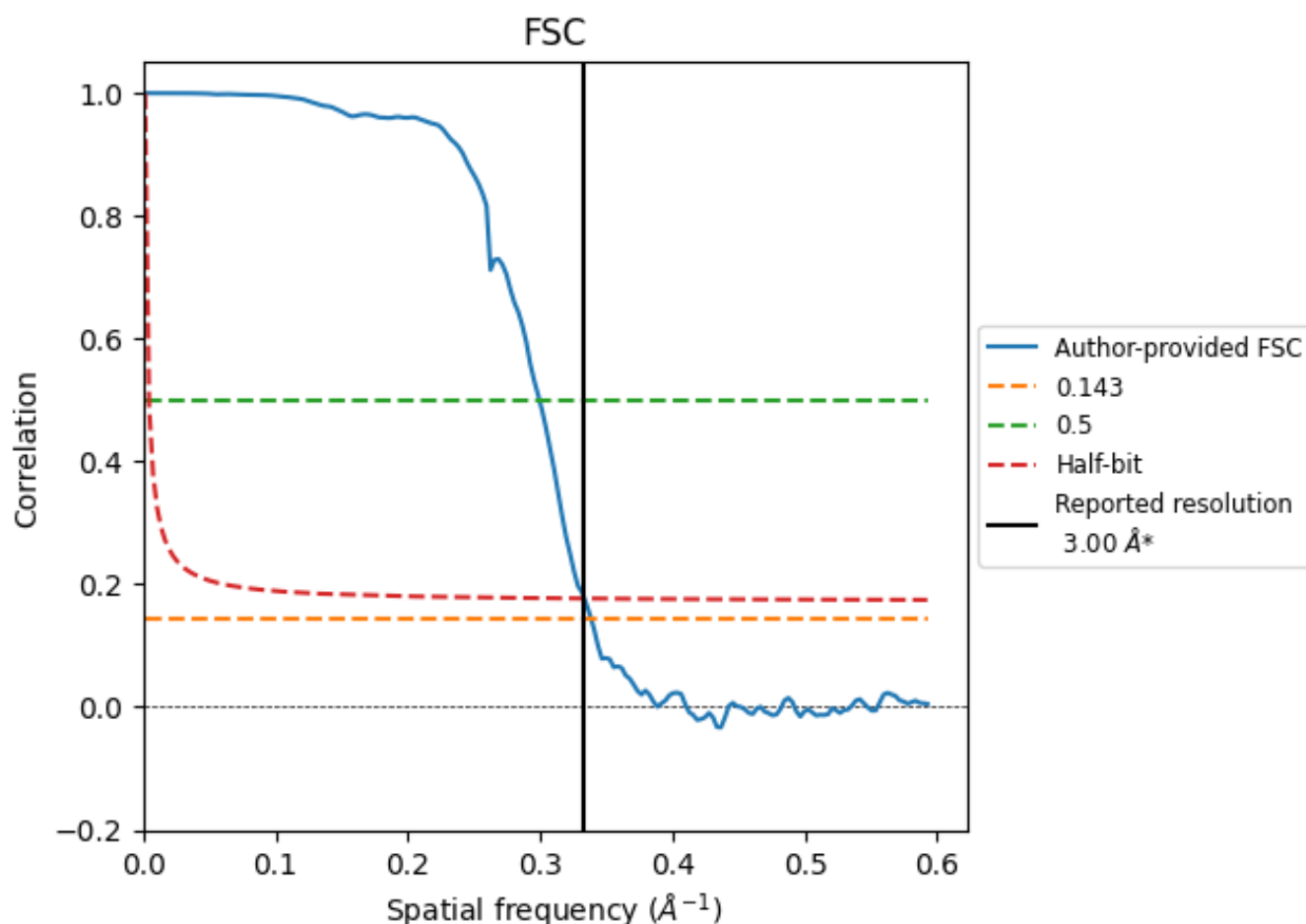


*Reported resolution corresponds to spatial frequency of 0.333 Å⁻¹

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 0.333 Å⁻¹

8.2 Resolution estimates [i](#)

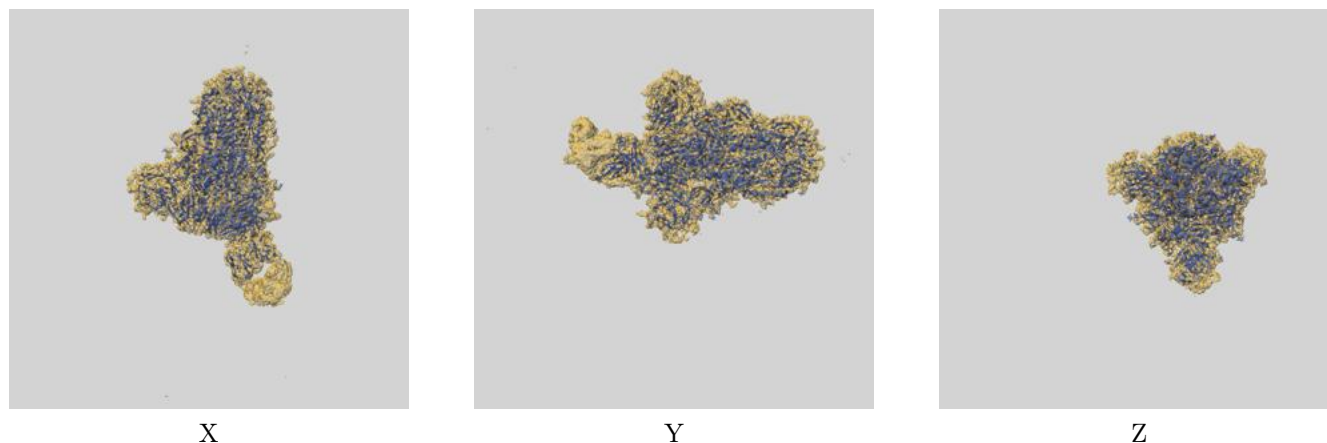
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	2.95	3.34	3.00
Unmasked-calculated*	-	-	-

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

9 Map-model fit [i](#)

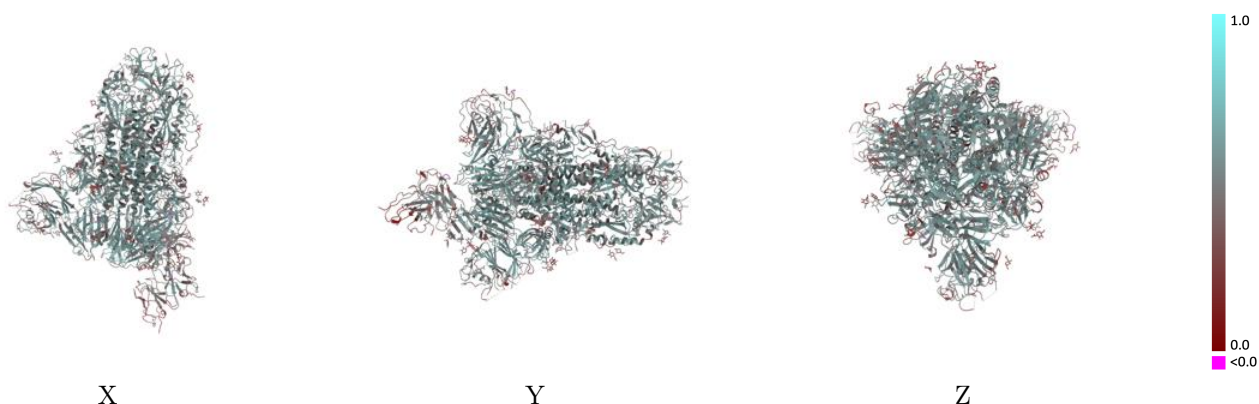
This section contains information regarding the fit between EMDB map EMD-40272 and PDB model 8SAK. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



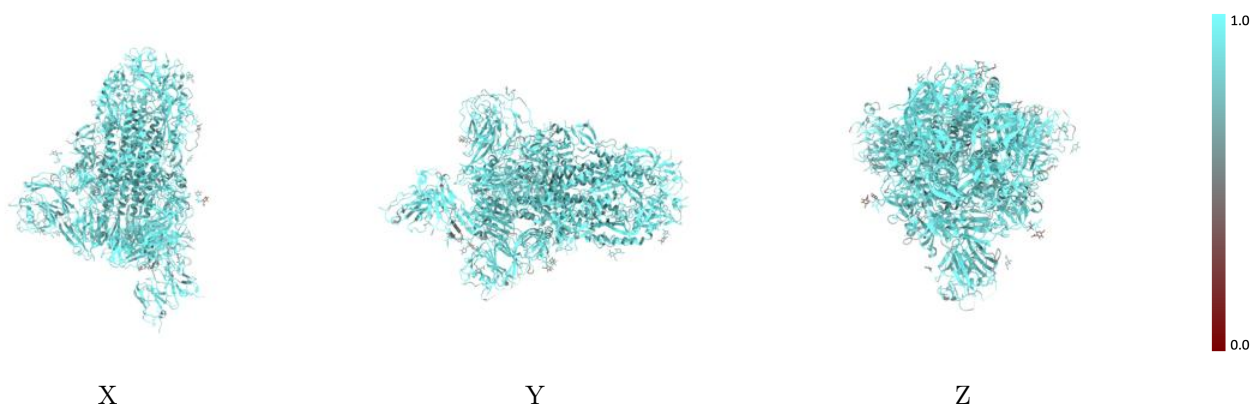
The images above show the 3D surface view of the map at the recommended contour level 1.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



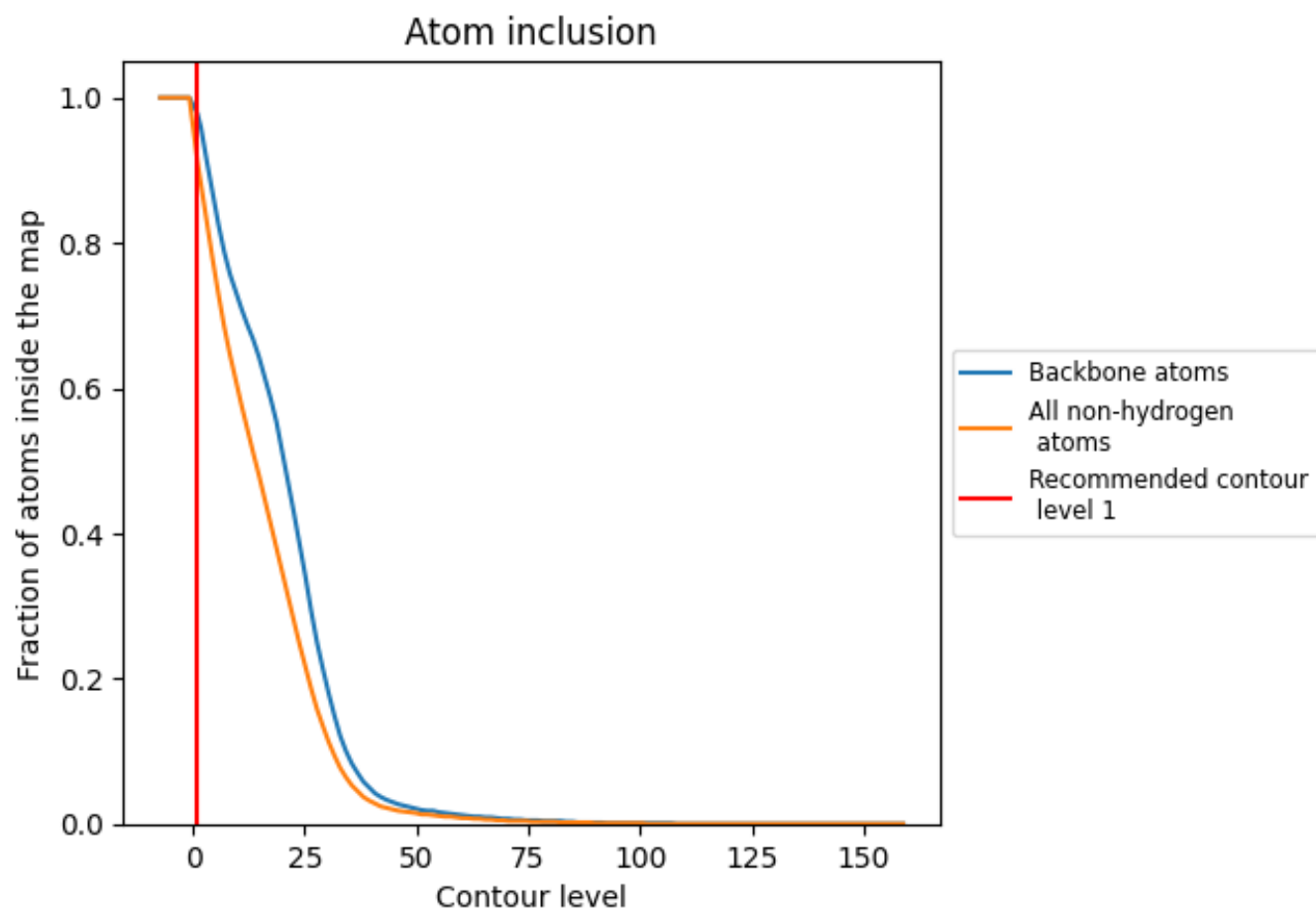
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (1).

9.4 Atom inclusion ⓘ



At the recommended contour level, 98% of all backbone atoms, 91% 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 (1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9120	 0.5140
1	 0.6320	 0.3070
5	 0.7500	 0.3240
7	 0.7500	 0.2870
A	 0.9120	 0.5220
B	 0.9210	 0.5280
C	 0.9180	 0.5280
E	 0.7500	 0.4150
H	 0.9240	 0.4540
K	 0.6790	 0.1870
L	 0.8770	 0.3280
Q	 0.6050	 0.2800
U	 0.6790	 0.3310
X	 0.7500	 0.3840
c	 0.6430	 0.2820
f	 0.5710	 0.2280
h	 0.5530	 0.2890
l	 0.8210	 0.3290
n	 0.6430	 0.3180
p	 0.7500	 0.3190
t	 0.8030	 0.3310
y	 0.6790	 0.3330

