



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 8, 2026 – 05:23 PM JST

PDB ID : 9X9Z / pdb_00009x9z
Title : Structure of the Omicron BA.2.86 variant Spike N-terminal domain(NTD)in complex with the antibody Fab fragment, C1717
Authors : Zhou, J.J.; Gao, G.F.
Deposited on : 2025-10-22
Resolution : 2.55 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

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:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

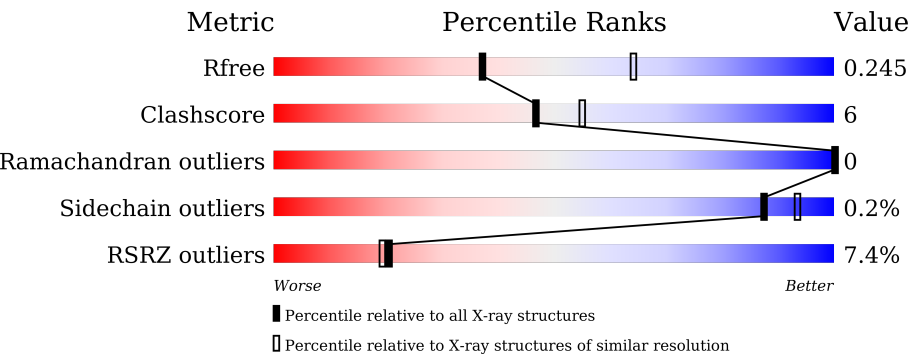
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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.



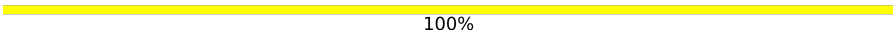
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	295	8% 80% 18% .
2	H	233	4% 85% 11% .
3	L	216	10% 88% 11% .
4	B	3	33% 67%
4	E	3	67% 33%
5	C	4	25% 75%

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Mol	Chain	Length	Quality of chain
6	D	2	 50%50%
6	F	2	 100%
6	G	2	 50%50%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5903 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2316	1496	375	436	9			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	MET	-	insertion	UNP P0DTC2
A	18	PRO	-	insertion	UNP P0DTC2
A	19	LEU	-	insertion	UNP P0DTC2
A	20	PHE	-	insertion	UNP P0DTC2
A	23	ILE	THR	variant	UNP P0DTC2
A	25	THR	ARG	conflict	UNP P0DTC2
A	?	-	LEU	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	28	SER	ALA	variant	UNP P0DTC2
A	51	LEU	SER	conflict	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	126	PHE	VAL	conflict	UNP P0DTC2
A	141	ASP	GLY	variant	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	155	SER	PHE	conflict	UNP P0DTC2
A	156	GLY	ARG	conflict	UNP P0DTC2
A	?	-	ASN	deletion	UNP P0DTC2
A	209	ILE	LEU	variant	UNP P0DTC2
A	210	GLY	VAL	variant	UNP P0DTC2
A	213	PHE	LEU	conflict	UNP P0DTC2
A	242	ASN	HIS	conflict	UNP P0DTC2
A	261	ASP	ALA	conflict	UNP P0DTC2
A	303	HIS	-	expression tag	UNP P0DTC2
A	304	HIS	-	expression tag	UNP P0DTC2
A	305	HIS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	306	HIS	-	expression tag	UNP P0DTC2
A	307	HIS	-	expression tag	UNP P0DTC2
A	308	HIS	-	expression tag	UNP P0DTC2

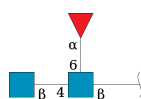
- Molecule 2 is a protein called C1717 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	224	Total	C	N	O	S	0	0	0
			1673	1057	282	327	7			

- Molecule 3 is a protein called C1717 light chain.

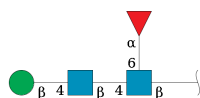
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total	C	N	O	S	0	0	0
			1570	973	269	323	5			

- Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	B	3	Total	C	N	O	0	0	0
			38	22	2	14			
4	E	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



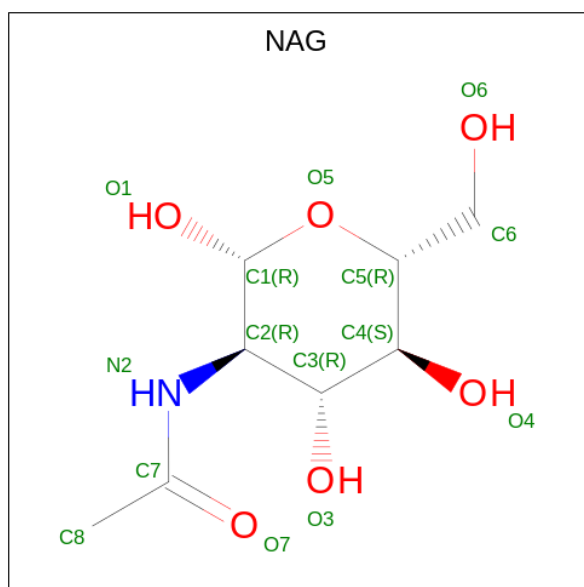
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	C	4	Total	C	N	O	0	0	0
			49	28	2	19			

- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
6	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
6	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	66	Total	O	0	0
			66	66		
8	H	37	Total	O	0	0
			37	37		

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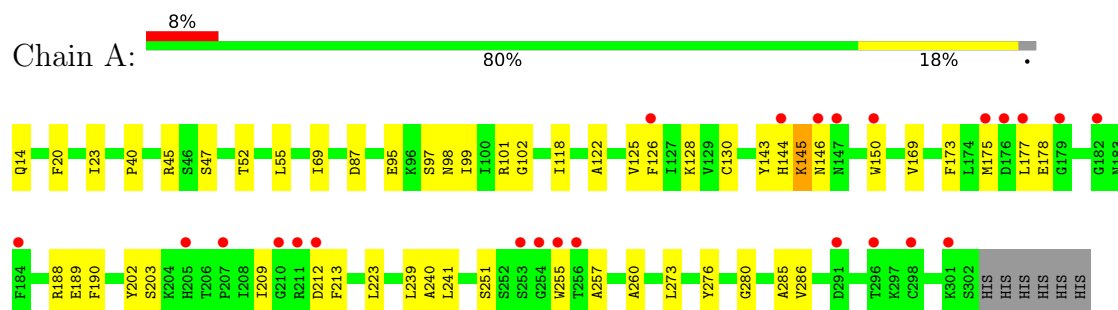
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	18	Total	O	0	0
			18	18		

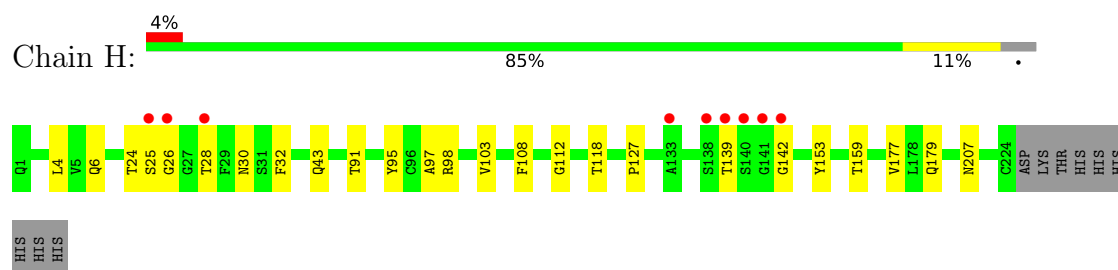
3 Residue-property plots

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 electron 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 dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

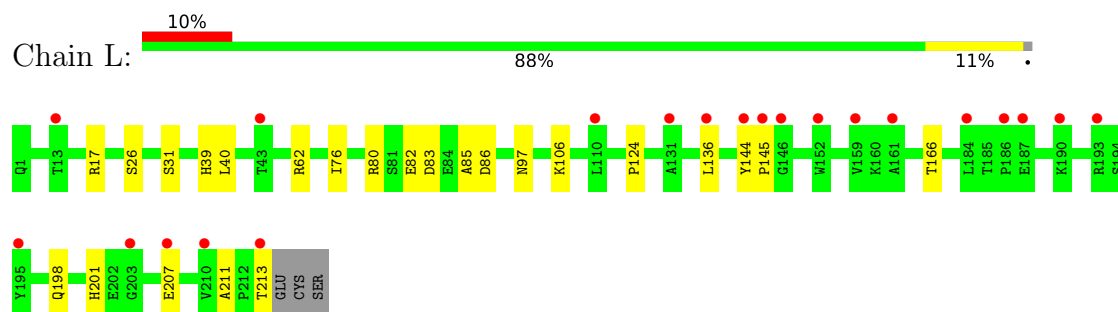
- Molecule 1: Spike protein S1



- Molecule 2: C1717 heavy chain



- Molecule 3: C1717 light chain



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



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

Chain E:  67% 33%

NAG1
NAG2
FUC3

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

Chain C:  25% 75%

NAG1
NAG2
BMA3
FUC4

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  50% 50%

NAG1
NAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%

NAG1
NAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  50% 50%

NAG1
NAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.79Å 58.17Å 112.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.11 – 2.55 42.11 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.6 (42.11-2.55) 99.7 (42.11-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.197 , 0.247 0.197 , 0.245	Depositor DCC
R_{free} test set	1546 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	37.7	Xtriage
Anisotropy	0.857	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5903	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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, 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	0.14	0/2380	0.37	0/3237
2	H	0.10	0/1714	0.32	0/2338
3	L	0.21	0/1607	0.37	0/2194
All	All	0.15	0/5701	0.35	0/7769

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	2316	0	2227	38	0
2	H	1673	0	1654	16	0
3	L	1570	0	1524	14	0
4	B	38	0	34	0	0
4	E	38	0	34	0	0
5	C	49	0	43	0	0
6	D	28	0	25	0	0
6	F	28	0	25	0	0
6	G	28	0	25	3	0
7	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	66	0	0	0	0
8	H	37	0	0	0	0
8	L	18	0	0	0	0
All	All	5903	0	5604	65	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (65) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:TYR:HD2	1:A:145:LYS:HB3	1.57	0.70
2:H:28:THR:HG22	2:H:98:ARG:HH11	1.61	0.65
2:H:127:PRO:HB3	2:H:153:TYR:HB3	1.80	0.61
3:L:40:LEU:HD23	3:L:85:ALA:HB2	1.82	0.61
2:H:6:GLN:HE21	2:H:112:GLY:HA3	1.66	0.59
1:A:97:SER:HB2	1:A:260:ALA:HA	1.85	0.58
1:A:212:ASP:OD1	1:A:213:PHE:N	2.36	0.57
3:L:198:GLN:HG2	3:L:207:GLU:HG3	1.88	0.56
1:A:143:TYR:CD2	1:A:145:LYS:HB3	2.37	0.56
1:A:178:GLU:HG3	6:G:1:NAG:H82	1.88	0.56
3:L:62:ARG:NH2	3:L:83:ASP:OD2	2.40	0.55
1:A:98:ASN:HD22	1:A:178:GLU:HA	1.71	0.55
2:H:139:THR:HG23	2:H:142:GLY:H	1.72	0.55
1:A:144:HIS:CD2	6:G:1:NAG:H3	2.42	0.55
1:A:101:ARG:HG3	1:A:240:ALA:HB2	1.89	0.54
2:H:97:ALA:HB1	2:H:108:PHE:HB3	1.91	0.53
1:A:99:ILE:HG23	1:A:239:LEU:HB3	1.90	0.52
1:A:98:ASN:OD1	1:A:98:ASN:N	2.42	0.51
2:H:30:ASN:OD1	2:H:30:ASN:N	2.41	0.51
1:A:144:HIS:NE2	6:G:1:NAG:H3	2.27	0.50
1:A:40:PRO:HG2	1:A:52:THR:HG21	1.94	0.50
1:A:128:LYS:HG2	1:A:130:CYS:SG	2.51	0.50
1:A:145:LYS:HG3	1:A:146:ASN:N	2.27	0.50
2:H:25:SER:OG	2:H:26:GLY:N	2.45	0.49
1:A:23:ILE:HG23	1:A:69:ILE:HG23	1.95	0.48
1:A:190:PHE:HA	1:A:202:TYR:O	2.14	0.48
1:A:286:VAL:HG22	2:H:103:VAL:HG22	1.93	0.48
3:L:26:SER:O	3:L:31:SER:OG	2.27	0.47
2:H:91:THR:HG23	2:H:118:THR:HA	1.97	0.46
3:L:144:TYR:HB3	3:L:145:PRO:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:TRP:H	1:A:150:TRP:HE3	1.64	0.46
1:A:20:PHE:HB2	1:A:241:LEU:HD13	1.99	0.45
2:H:159:THR:OG1	2:H:207:ASN:HB3	2.16	0.45
1:A:20:PHE:CE1	1:A:257:ALA:HB2	2.52	0.45
1:A:189:GLU:O	1:A:203:SER:HA	2.18	0.44
3:L:80:ARG:NH1	3:L:82:GLU:OE1	2.50	0.44
1:A:125:VAL:HG21	1:A:173:PHE:CZ	2.52	0.44
1:A:145:LYS:HZ3	1:A:146:ASN:HB3	1.82	0.44
2:H:43:GLN:H	2:H:43:GLN:HG2	1.60	0.44
2:H:177:VAL:HB	3:L:166:THR:HG22	1.99	0.44
3:L:86:ASP:OD1	3:L:106:LYS:HG2	2.18	0.44
3:L:124:PRO:HD3	3:L:136:LEU:HD22	1.99	0.44
1:A:251:SER:HB3	1:A:255:TRP:CZ2	2.53	0.43
3:L:17:ARG:HG3	3:L:76:ILE:O	2.18	0.43
3:L:145:PRO:HD2	3:L:201:HIS:NE2	2.33	0.43
1:A:209:ILE:HD12	3:L:97:ASN:HB2	2.00	0.43
3:L:39:HIS:O	3:L:85:ALA:HB1	2.18	0.43
1:A:45:ARG:O	1:A:280:GLY:HA2	2.19	0.43
1:A:145:LYS:HG3	1:A:145:LYS:HZ3	1.68	0.43
3:L:211:ALA:O	3:L:213:THR:N	2.52	0.43
1:A:273:LEU:O	1:A:285:ALA:HA	2.19	0.42
1:A:173:PHE:CD2	1:A:223:LEU:HD21	2.53	0.42
1:A:122:ALA:N	1:A:177:LEU:HD11	2.34	0.42
1:A:47:SER:HA	1:A:276:TYR:O	2.19	0.42
1:A:102:GLY:HA3	1:A:118:ILE:O	2.19	0.42
1:A:101:ARG:NE	1:A:177:LEU:HD22	2.34	0.42
2:H:28:THR:HA	2:H:32:PHE:CD1	2.55	0.42
1:A:95:GLU:HG3	1:A:188:ARG:NH2	2.35	0.41
2:H:6:GLN:HE22	2:H:95:TYR:HA	1.85	0.41
1:A:14:GLN:N	1:A:14:GLN:OE1	2.54	0.41
1:A:126:PHE:CD1	1:A:169:VAL:HG22	2.55	0.41
2:H:4:LEU:HD23	2:H:24:THR:HA	2.02	0.41
1:A:175:MET:HE3	1:A:175:MET:HB2	1.98	0.41
1:A:55:LEU:HD13	1:A:87:ASP:O	2.21	0.41
2:H:179:GLN:H	2:H:179:GLN:HG2	1.72	0.41

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 X-ray entries followed by that with respect to entries of similar resolution.

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	287/295 (97%)	269 (94%)	18 (6%)	0	100	100
2	H	222/233 (95%)	213 (96%)	9 (4%)	0	100	100
3	L	211/216 (98%)	195 (92%)	16 (8%)	0	100	100
All	All	720/744 (97%)	677 (94%)	43 (6%)	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 X-ray entries followed by that with respect to entries of similar resolution.

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	260/266 (98%)	259 (100%)	1 (0%)	84	90
2	H	189/198 (96%)	189 (100%)	0	100	100
3	L	178/181 (98%)	178 (100%)	0	100	100
All	All	627/645 (97%)	626 (100%)	1 (0%)	87	93

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	53	GLN
1	A	136	ASN
1	A	186	ASN
1	A	205	HIS
1	A	236	GLN
2	H	6	GLN
2	H	207	ASN
3	L	198	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 ⓘ

16 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	B	1	1,4	14,14,15	0.70	0	17,19,21	1.32	2 (11%)
4	NAG	B	2	4	14,14,15	0.71	0	17,19,21	1.21	1 (5%)
4	FUC	B	3	4	10,10,11	0.83	0	14,14,16	0.93	0
5	NAG	C	1	1,5	14,14,15	0.84	0	17,19,21	1.14	1 (5%)
5	NAG	C	2	5	14,14,15	0.71	0	17,19,21	1.21	1 (5%)
5	BMA	C	3	5	11,11,12	0.87	0	15,15,17	2.23	4 (26%)
5	FUC	C	4	5	10,10,11	0.77	0	14,14,16	1.05	0
6	NAG	D	1	1,6	14,14,15	0.71	0	17,19,21	1.13	1 (5%)
6	NAG	D	2	6	14,14,15	0.69	0	17,19,21	0.92	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	E	1	1,4	14,14,15	0.75	0	17,19,21	1.20	1 (5%)
4	NAG	E	2	4	14,14,15	0.72	0	17,19,21	0.83	0
4	FUC	E	3	4	10,10,11	0.82	0	14,14,16	0.94	0
6	NAG	F	1	1,6	14,14,15	0.71	0	17,19,21	1.61	2 (11%)
6	NAG	F	2	6	14,14,15	0.79	1 (7%)	17,19,21	2.04	4 (23%)
6	NAG	G	1	1,6	14,14,15	0.72	0	17,19,21	1.18	2 (11%)
6	NAG	G	2	6	14,14,15	0.71	0	17,19,21	0.93	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	1	1,4	-	4/6/23/26	0/1/1/1
4	NAG	B	2	4	-	3/6/23/26	0/1/1/1
4	FUC	B	3	4	-	-	0/1/1/1
5	NAG	C	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	C	2	5	-	1/6/23/26	0/1/1/1
5	BMA	C	3	5	-	2/2/19/22	0/1/1/1
5	FUC	C	4	5	-	-	0/1/1/1
6	NAG	D	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	D	2	6	-	2/6/23/26	0/1/1/1
4	NAG	E	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	FUC	E	3	4	-	-	0/1/1/1
6	NAG	F	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	F	2	6	-	2/6/23/26	0/1/1/1
6	NAG	G	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	G	2	6	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	2	NAG	C1-C2	2.17	1.55	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	3	BMA	C1-O5-C5	6.61	121.15	112.19
6	F	2	NAG	C1-O5-C5	6.61	121.14	112.19
6	F	1	NAG	C1-O5-C5	4.51	118.31	112.19
4	E	1	NAG	C1-O5-C5	3.58	117.04	112.19
4	B	1	NAG	C2-N2-C7	3.27	127.57	122.90
5	C	2	NAG	C2-N2-C7	3.26	127.54	122.90
4	B	2	NAG	C2-N2-C7	3.22	127.48	122.90
6	F	1	NAG	O4-C4-C5	3.16	117.14	109.30
6	F	2	NAG	C3-C4-C5	-2.93	105.01	110.24
6	G	1	NAG	C1-O5-C5	2.92	116.16	112.19
6	D	1	NAG	O5-C1-C2	-2.83	106.82	111.29
5	C	3	BMA	C3-C4-C5	2.61	114.90	110.24
5	C	3	BMA	C2-C3-C4	2.41	115.07	110.89
6	F	2	NAG	O5-C1-C2	2.37	115.03	111.29
6	F	2	NAG	C2-N2-C7	2.34	126.24	122.90
4	B	1	NAG	O5-C1-C2	-2.30	107.66	111.29
6	G	1	NAG	C2-N2-C7	2.27	126.13	122.90
5	C	1	NAG	C2-N2-C7	2.26	126.12	122.90
5	C	3	BMA	O4-C4-C3	-2.02	105.69	110.35

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	3	BMA	O5-C5-C6-O6
6	D	2	NAG	C8-C7-N2-C2
6	D	2	NAG	O7-C7-N2-C2
6	F	1	NAG	C8-C7-N2-C2
6	F	1	NAG	O7-C7-N2-C2
6	G	1	NAG	C8-C7-N2-C2
6	G	1	NAG	O7-C7-N2-C2
6	G	2	NAG	C8-C7-N2-C2
6	G	2	NAG	O7-C7-N2-C2
5	C	3	BMA	C4-C5-C6-O6
4	B	1	NAG	O5-C5-C6-O6
4	B	2	NAG	O5-C5-C6-O6
4	B	1	NAG	C4-C5-C6-O6
4	B	1	NAG	C1-C2-N2-C7
4	B	1	NAG	C3-C2-N2-C7
5	C	2	NAG	C3-C2-N2-C7
4	B	2	NAG	C1-C2-N2-C7
6	F	2	NAG	C1-C2-N2-C7
4	B	2	NAG	C3-C2-N2-C7

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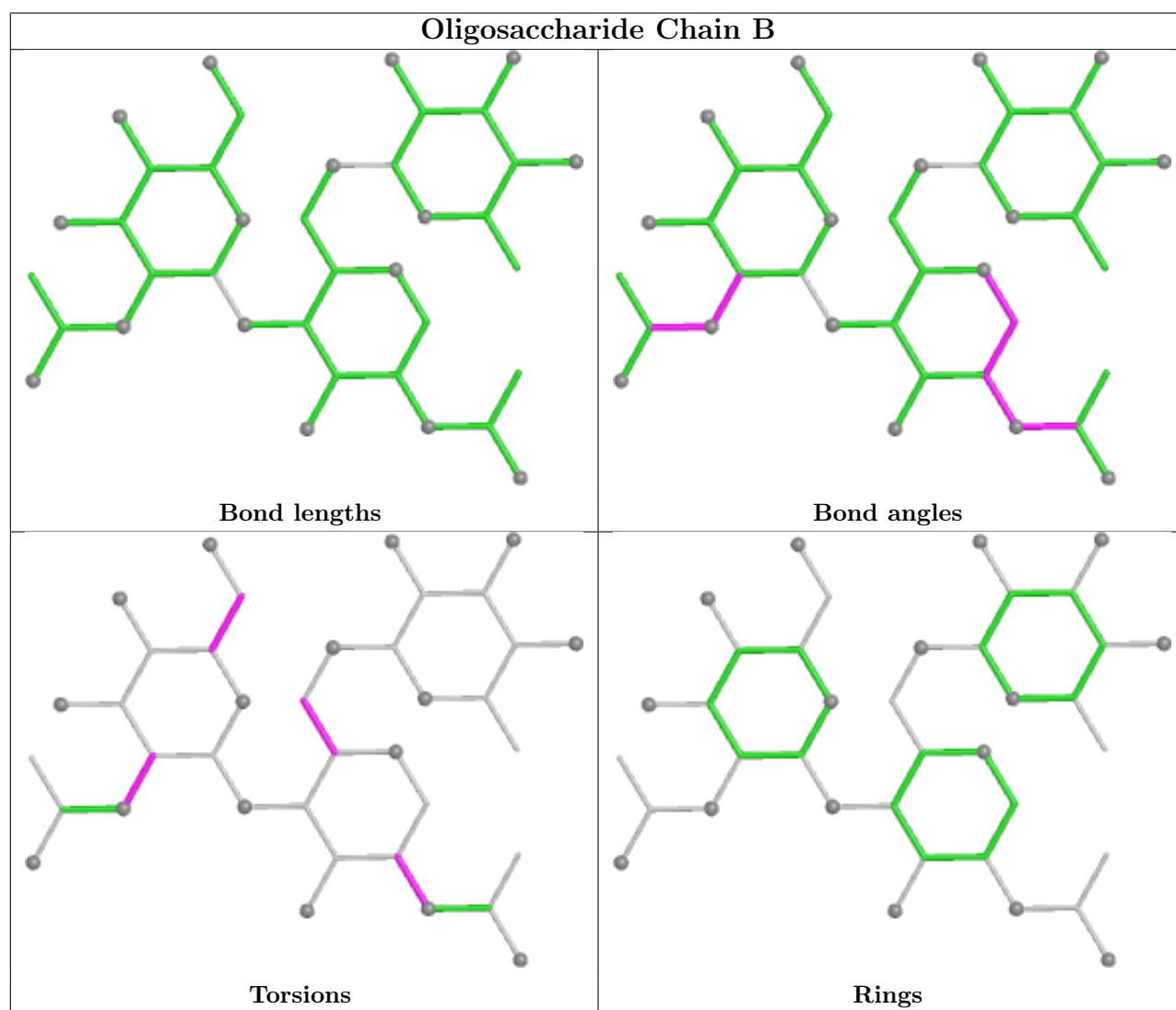
Mol	Chain	Res	Type	Atoms
6	F	2	NAG	C3-C2-N2-C7

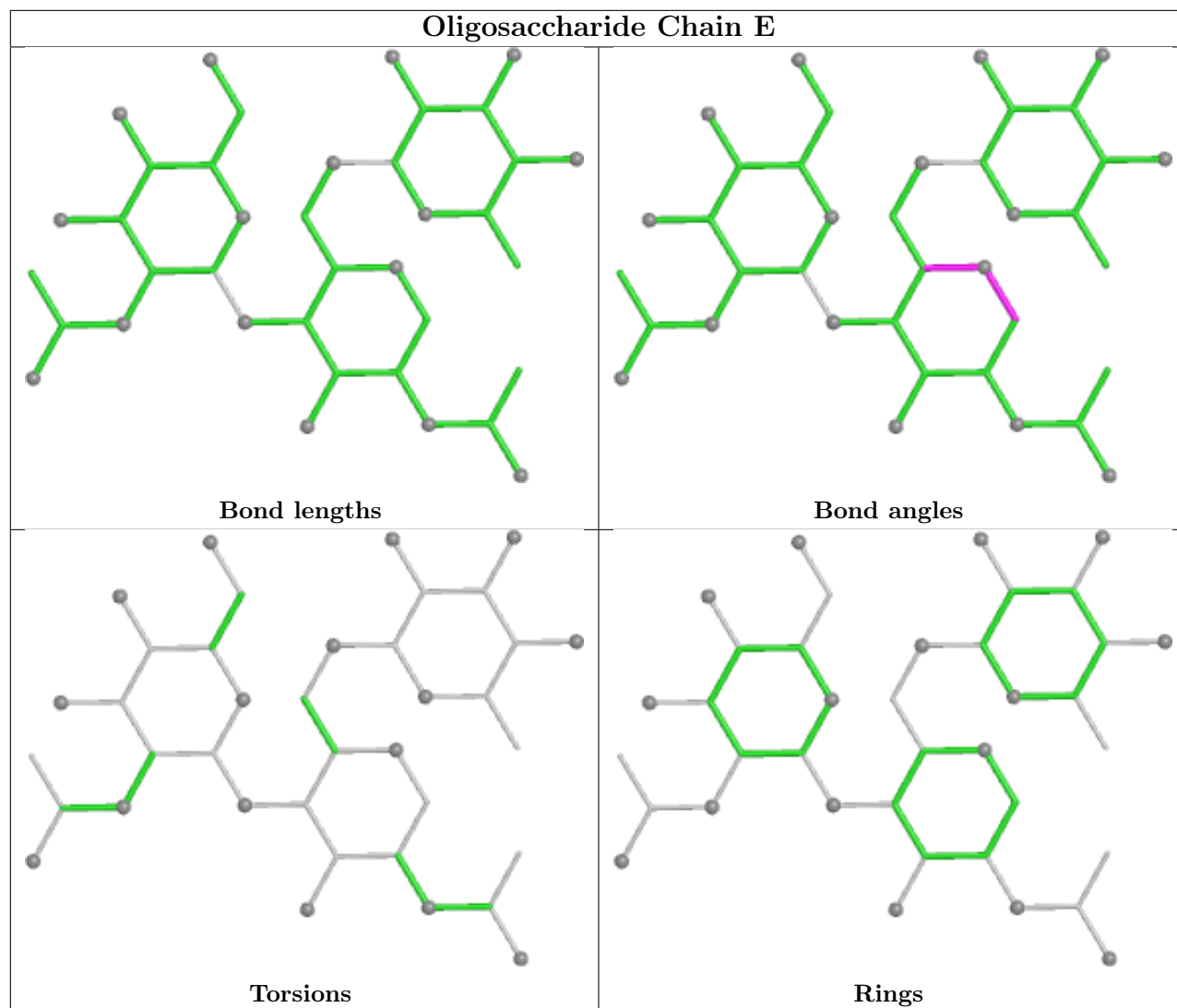
There are no ring outliers.

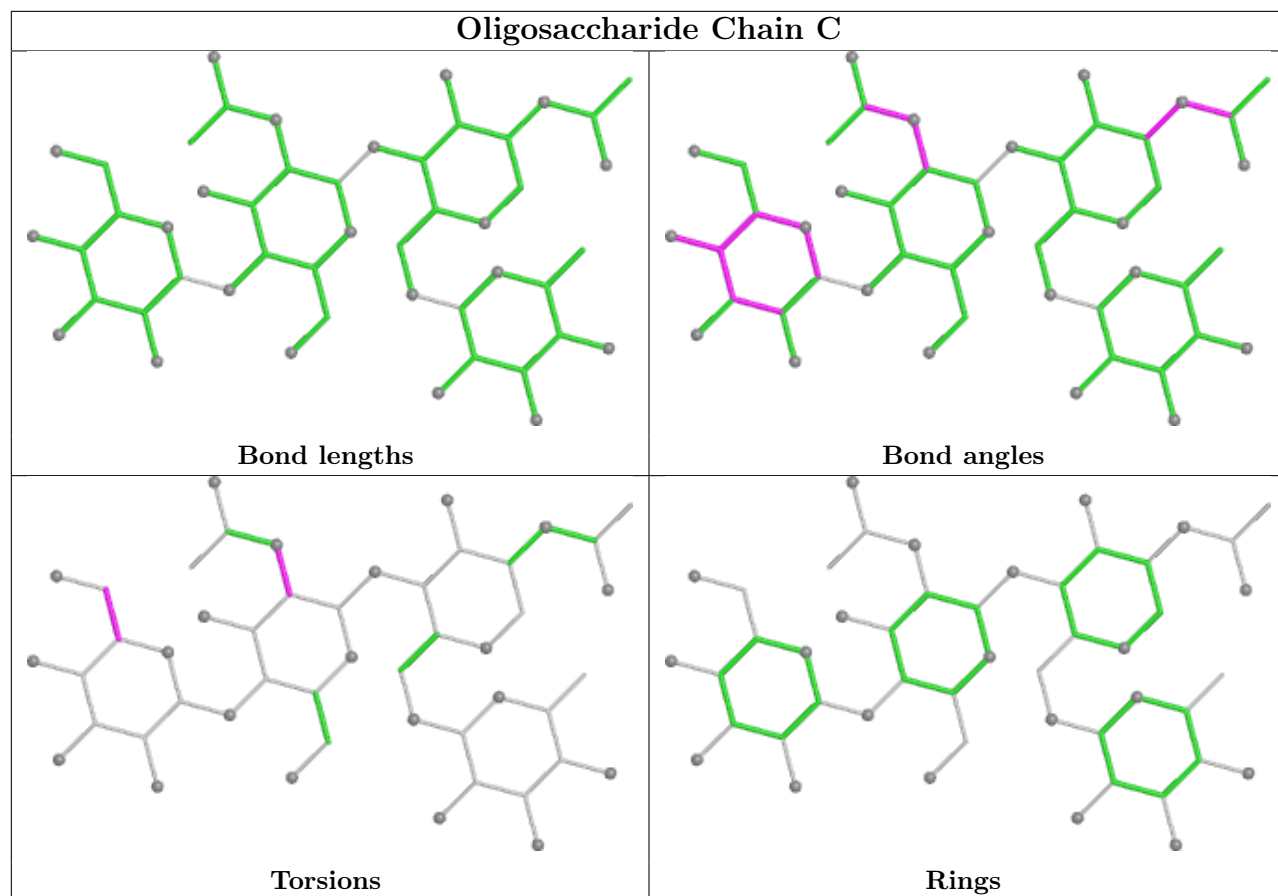
1 monomer is involved in 3 short contacts:

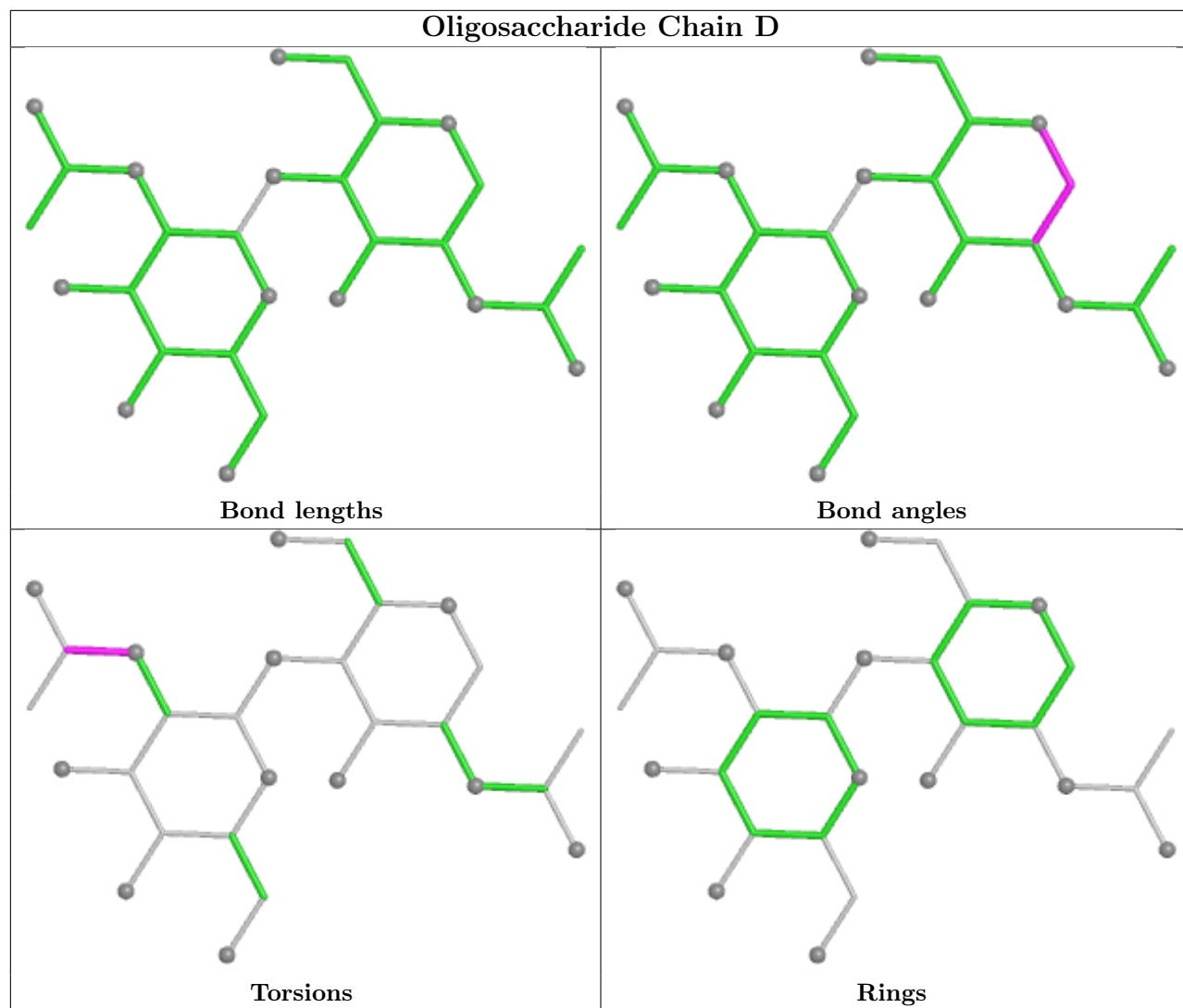
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	G	1	NAG	3	0

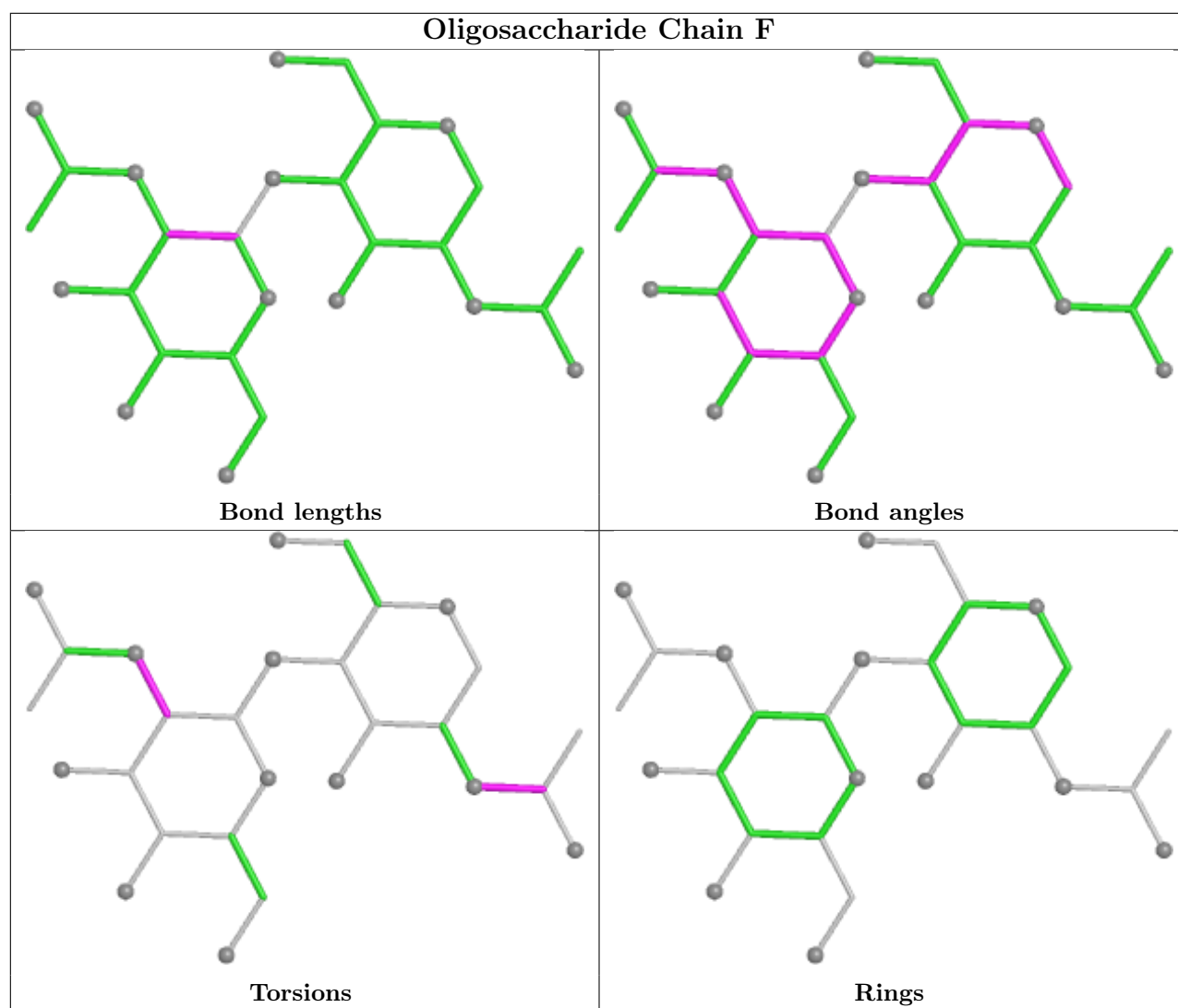
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

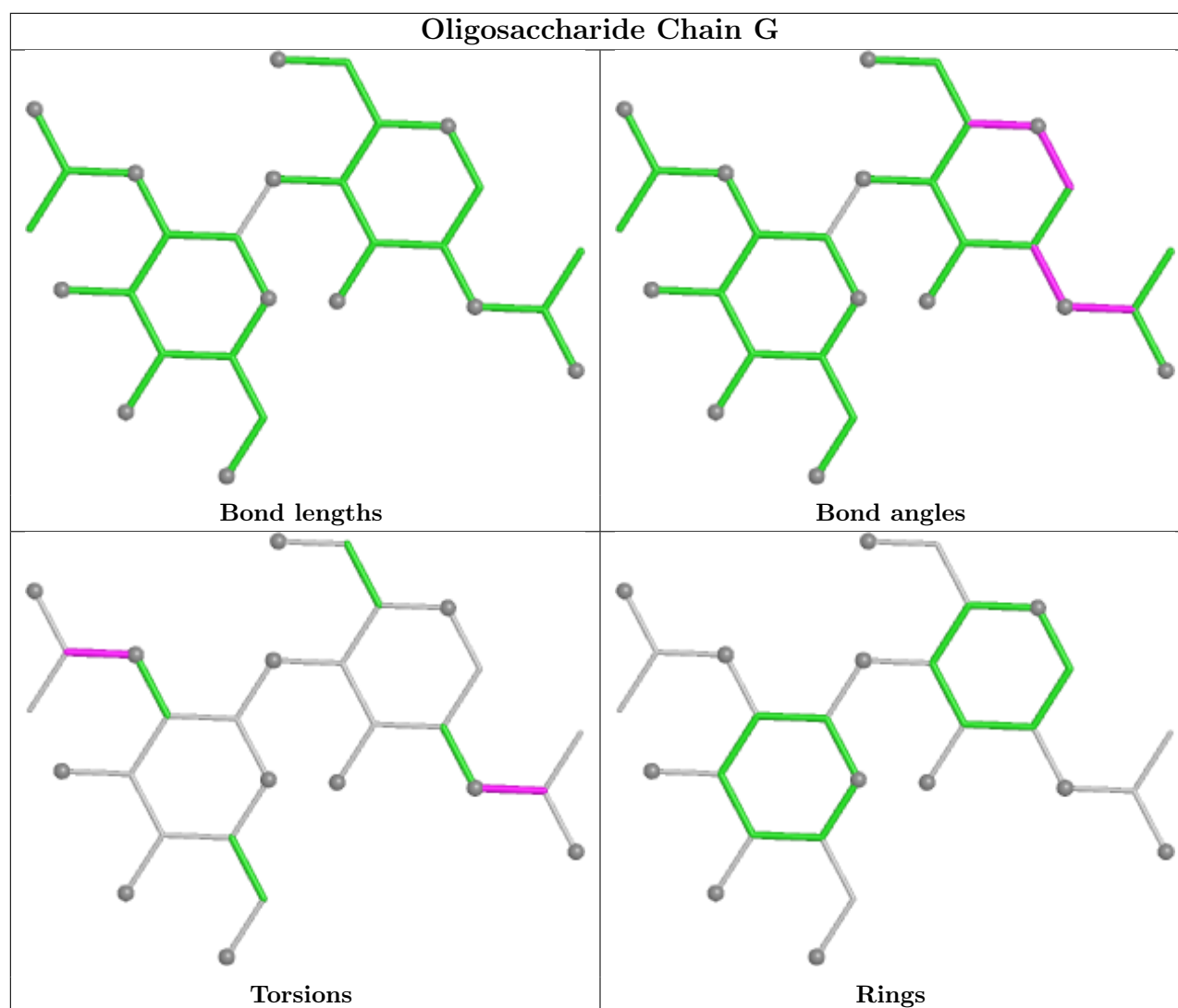












5.6 Ligand geometry [i](#)

1 ligand is 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	401	1	14,14,15	0.73	0	17,19,21	0.92	1 (5%)

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	401	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	401	NAG	C2-N2-C7	2.02	125.78	122.90

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	401	NAG	C8-C7-N2-C2
7	A	401	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	289/295 (97%)	0.38	24 (8%)	17 17	23, 39, 119, 150	0
2	H	224/233 (96%)	0.23	9 (4%)	42 43	25, 46, 81, 150	0
3	L	213/216 (98%)	0.83	21 (9%)	13 12	32, 66, 106, 127	0
All	All	726/744 (97%)	0.46	54 (7%)	20 19	23, 48, 108, 150	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	150	TRP	4.8
1	A	255	TRP	4.1
3	L	152	TRP	4.0
1	A	184	PHE	4.0
3	L	144	TYR	3.9
1	A	254	GLY	3.8
2	H	139	THR	3.7
3	L	210	VAL	3.6
3	L	186	PRO	3.6
3	L	213	THR	3.6
2	H	140	SER	3.6
3	L	146	GLY	3.2
1	A	177	LEU	3.1
1	A	212	ASP	3.0
1	A	147	ASN	2.9
1	A	210	GLY	2.9
2	H	141	GLY	2.9
3	L	184	LEU	2.9
1	A	146	ASN	2.8
3	L	159	VAL	2.8
2	H	25	SER	2.7
1	A	296	THR	2.7
2	H	28	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	256	THR	2.6
1	A	176	ASP	2.6
3	L	193	ARG	2.6
3	L	161	ALA	2.5
1	A	291	ASP	2.5
1	A	205	HIS	2.5
1	A	253	SER	2.5
1	A	182	GLY	2.4
3	L	203	GLY	2.4
3	L	190	LYS	2.4
1	A	126	PHE	2.4
3	L	145	PRO	2.4
2	H	138	SER	2.3
1	A	207	PRO	2.3
2	H	26	GLY	2.3
3	L	43	THR	2.2
1	A	179	GLY	2.2
3	L	207	GLU	2.2
1	A	144	HIS	2.2
3	L	195	TYR	2.2
3	L	13	THR	2.2
1	A	298	CYS	2.2
2	H	142	GLY	2.2
3	L	187	GLU	2.2
1	A	175	MET	2.1
3	L	131	ALA	2.1
1	A	211	ARG	2.1
1	A	301	LYS	2.1
2	H	133	ALA	2.1
3	L	110	LEU	2.1
3	L	136	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

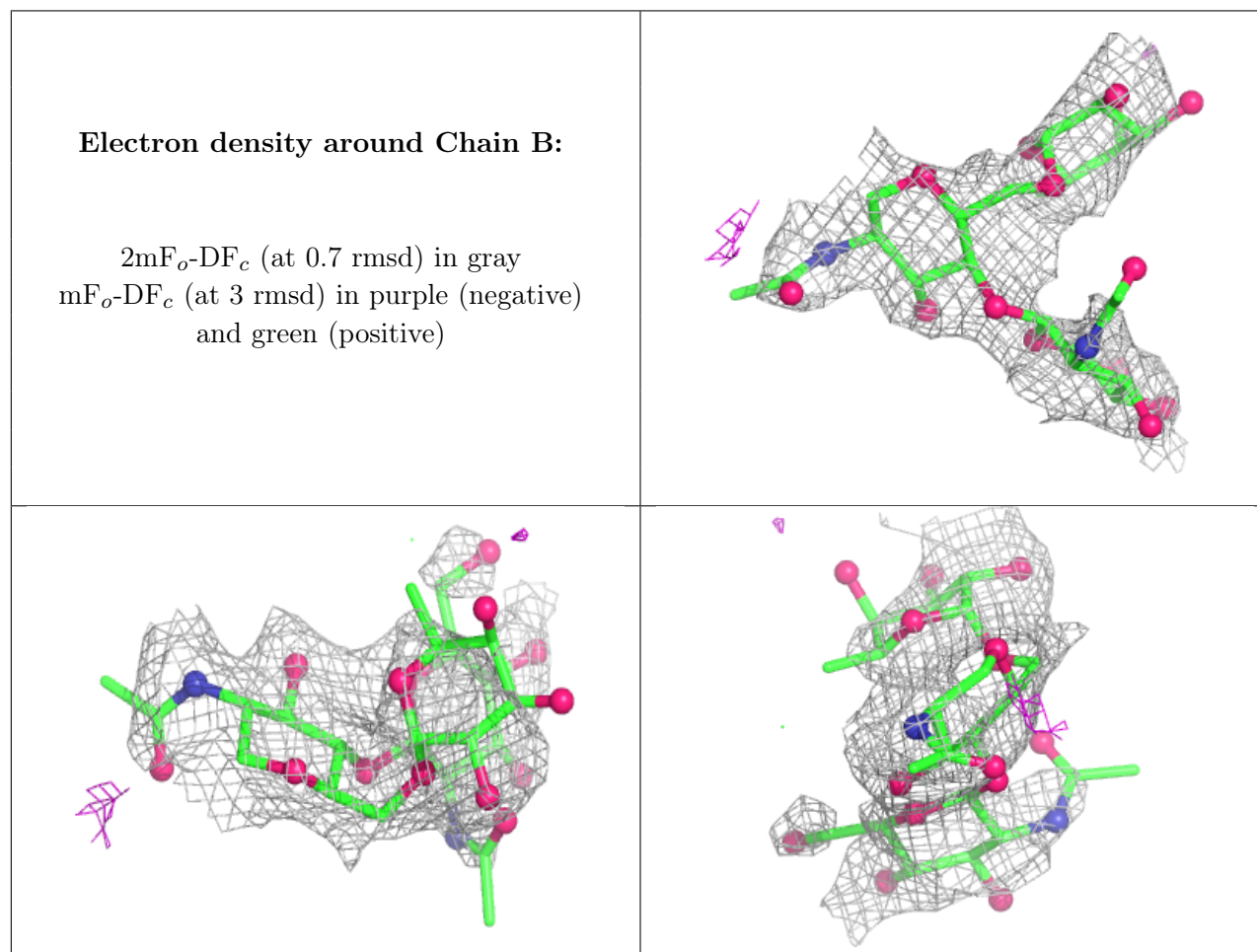
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

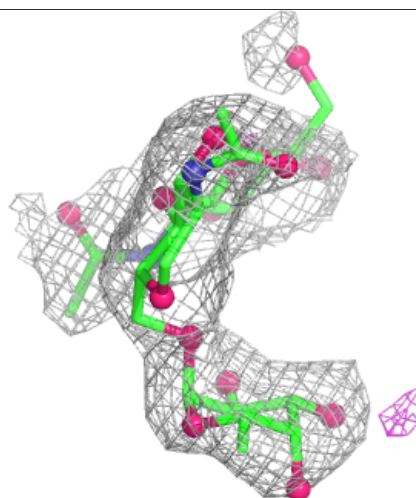
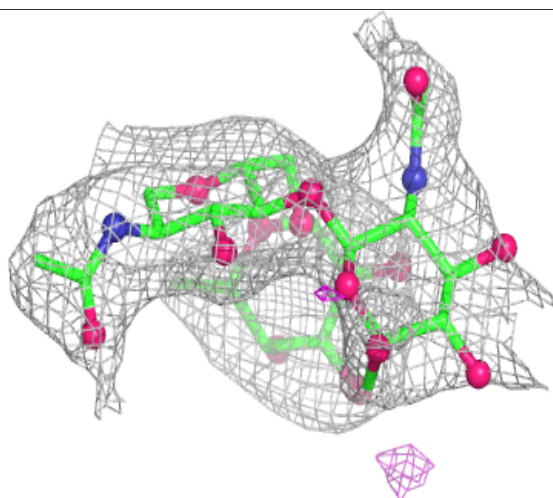
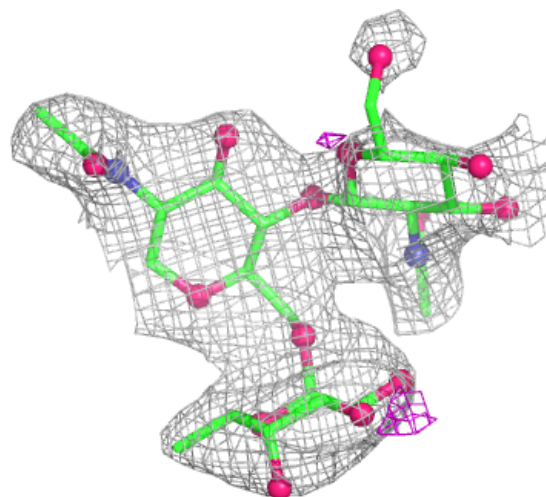
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	B	1	14/15	-	-	78,93,110,120	0
4	NAG	B	2	14/15	-	-	96,111,115,116	0
4	FUC	B	3	10/11	-	-	103,117,133,135	0
6	NAG	G	2	14/15	0.56	0.17	126,145,156,158	0
6	NAG	F	2	14/15	0.57	0.19	122,149,158,158	0
4	FUC	E	3	10/11	-	-	63,90,99,100	0
5	NAG	C	1	14/15	-	-	38,61,87,87	0
5	NAG	C	2	14/15	-	-	93,99,109,123	0
5	BMA	C	3	11/12	-	-	107,115,122,123	0
5	FUC	C	4	10/11	-	-	71,103,116,121	0
6	NAG	D	2	14/15	0.63	0.17	78,90,96,110	0
4	NAG	E	2	14/15	0.67	0.15	67,99,107,118	0
6	NAG	D	1	14/15	0.78	0.14	53,66,78,80	0
6	NAG	F	1	14/15	0.84	0.12	91,117,133,141	0
4	NAG	E	1	14/15	0.89	0.11	43,66,89,92	0
6	NAG	G	1	14/15	0.93	0.09	112,130,150,150	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



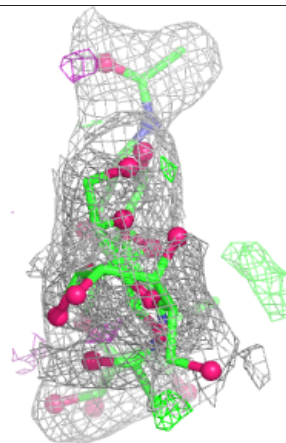
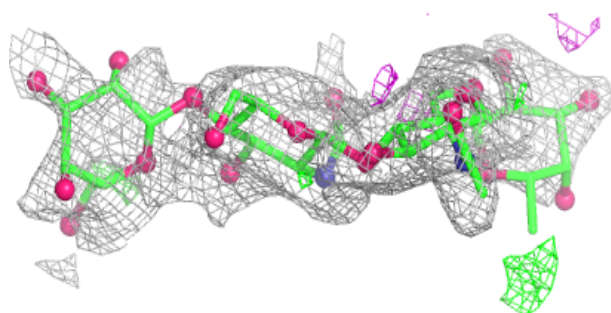
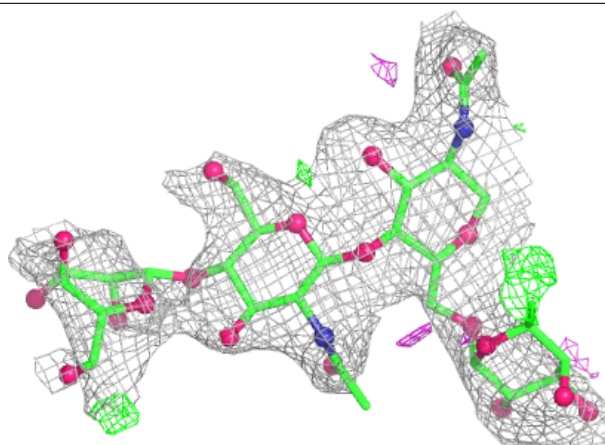
Electron density around Chain E:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



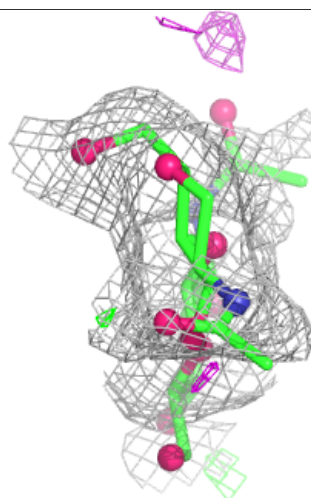
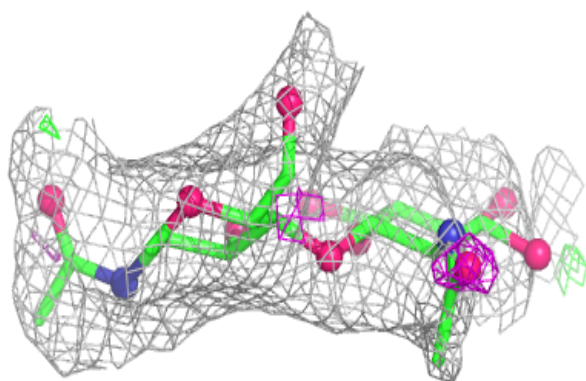
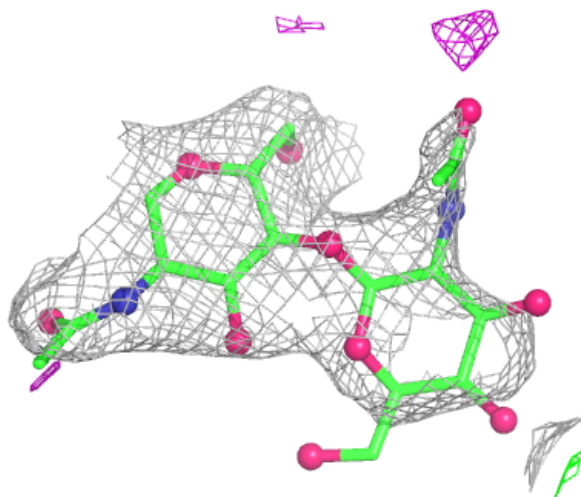
Electron density around Chain C:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



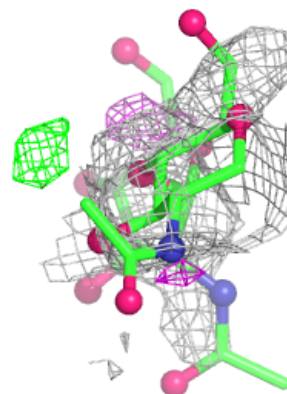
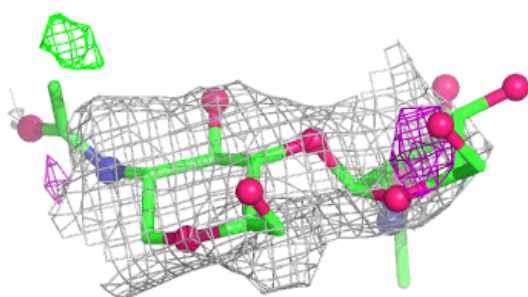
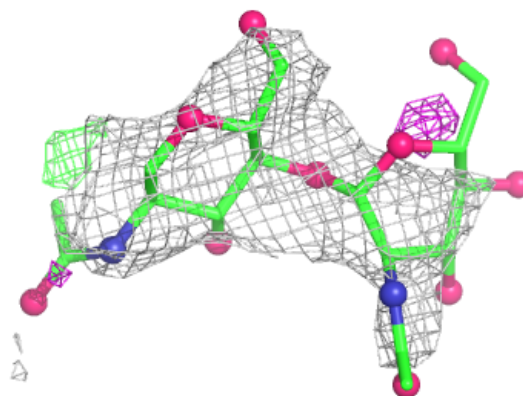
Electron density around Chain D:

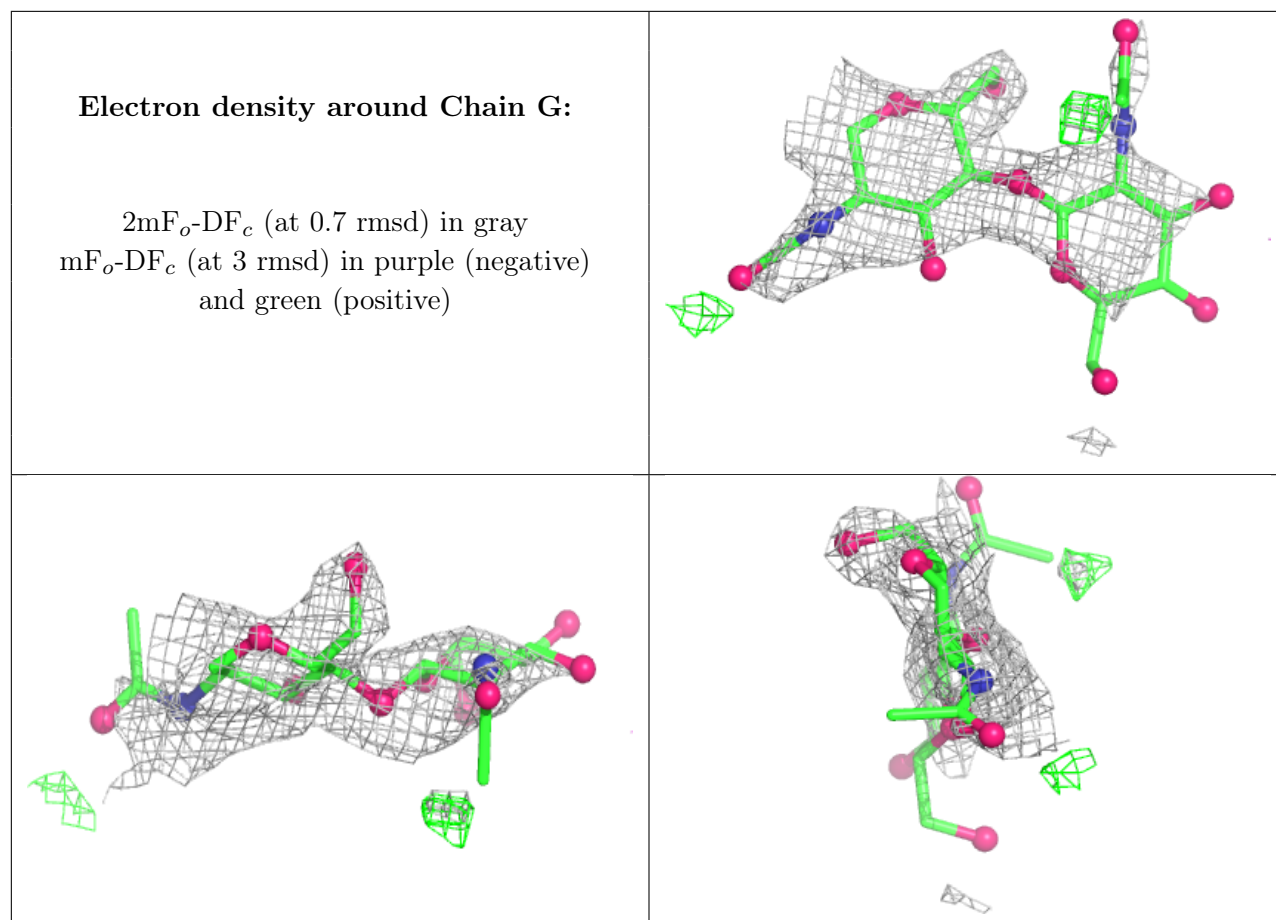
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain F:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)





6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	NAG	A	401	14/15	0.51	0.19	105,122,139,146	0

6.5 Other polymers [i](#)

There are no such residues in this entry.