



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2024 – 05:11 pm BST

PDB ID : 5HNS
Title : Structure of glycosylated NPC1 luminal domain C
Authors : Zhao, Y.; Ren, J.; Harlos, K.; Stuart, D.I.
Deposited on : 2016-01-18
Resolution : 2.45 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

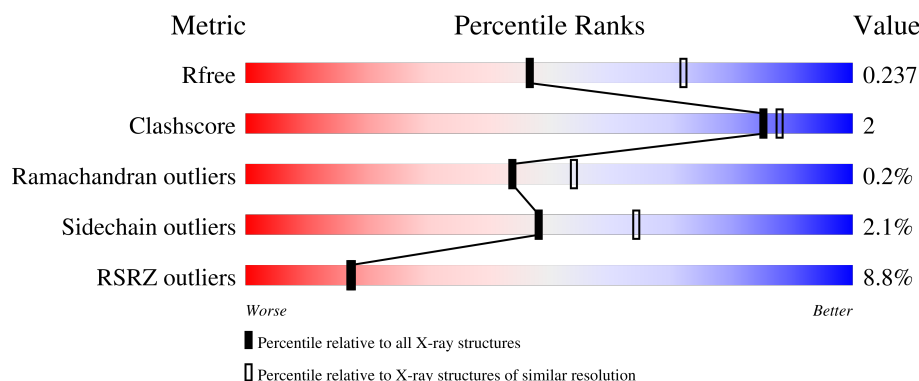
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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}	164625	1096 (2.46-2.46)
Clashscore	180529	1178 (2.46-2.46)
Ramachandran outliers	177936	1170 (2.46-2.46)
Sidechain outliers	177891	1170 (2.46-2.46)
RSRZ outliers	164620	1096 (2.46-2.46)

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	244	<div> <div>7%</div> <div>82%</div> <div>6%</div> <div>12%</div> </div>
1	B	244	<div> <div>9%</div> <div>85%</div> <div>•</div> <div>11%</div> </div>
2	C	2	<div> <div>50%</div> <div>50%</div> </div>
2	E	2	<div> <div>100%</div> </div>
2	F	2	<div> <div>100%</div> </div>

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Mol	Chain	Length	Quality of chain
2	G	2	<div><div></div><div>50%</div><div>50%</div></div>
2	H	2	<div><div></div><div>50%</div><div>50%</div></div>
3	D	4	<div><div></div><div>100%</div></div>
4	I	5	<div><div></div><div>40%</div><div>60%</div></div>
5	J	3	<div><div></div><div>33%</div><div>67%</div></div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 3943 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 Niemann-Pick C1 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	5	0
			1787	1157	288	338	4			
1	B	216	Total	C	N	O	S	0	5	0
			1796	1162	289	341	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	384	GLU	-	expression tag	UNP O15118
A	385	THR	-	expression tag	UNP O15118
A	386	GLY	-	expression tag	UNP O15118
A	619	THR	-	expression tag	UNP O15118
A	620	GLY	-	expression tag	UNP O15118
A	621	THR	-	expression tag	UNP O15118
A	622	LEU	-	expression tag	UNP O15118
A	623	GLU	-	expression tag	UNP O15118
A	624	VAL	-	expression tag	UNP O15118
A	625	LEU	-	expression tag	UNP O15118
A	626	PHE	-	expression tag	UNP O15118
A	627	GLN	-	expression tag	UNP O15118
B	384	GLU	-	expression tag	UNP O15118
B	385	THR	-	expression tag	UNP O15118
B	386	GLY	-	expression tag	UNP O15118
B	619	THR	-	expression tag	UNP O15118
B	620	GLY	-	expression tag	UNP O15118
B	621	THR	-	expression tag	UNP O15118
B	622	LEU	-	expression tag	UNP O15118
B	623	GLU	-	expression tag	UNP O15118
B	624	VAL	-	expression tag	UNP O15118
B	625	LEU	-	expression tag	UNP O15118
B	626	PHE	-	expression tag	UNP O15118
B	627	GLN	-	expression tag	UNP O15118

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



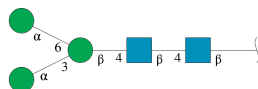
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is an oligosaccharide called 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				ZeroOcc	AltConf	Trace
3	D	4	Total	C	N	O	0	0	0
			50	28	2	20			

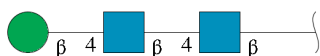
- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	I	5	Total	C	N	O	0	0	0
			61	34	2	25			

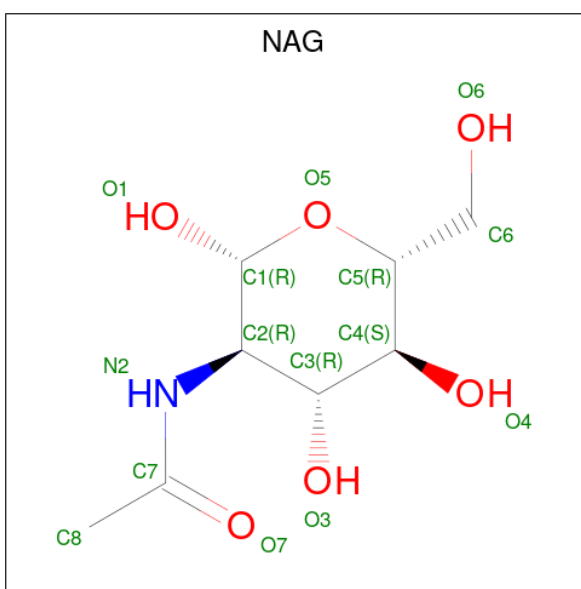
- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-b

eta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



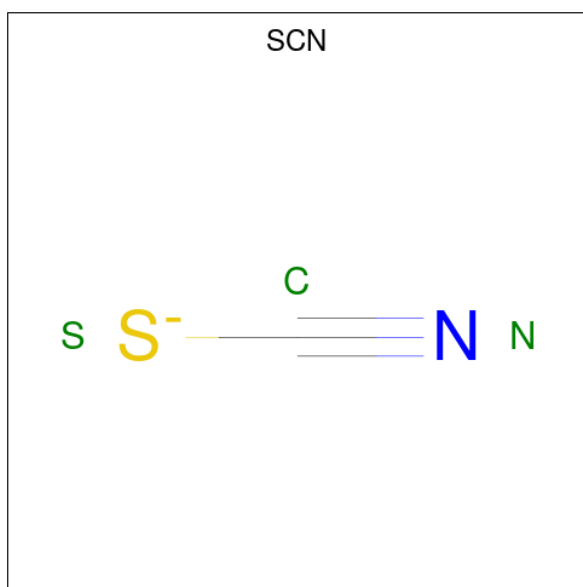
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	J	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 7 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	S	0	0
			3	1	1	1		
7	A	1	Total	C	N	S	0	0
			3	1	1	1		
7	B	1	Total	C	N	S	0	0
			3	1	1	1		
7	B	1	Total	C	N	S	0	0
			3	1	1	1		

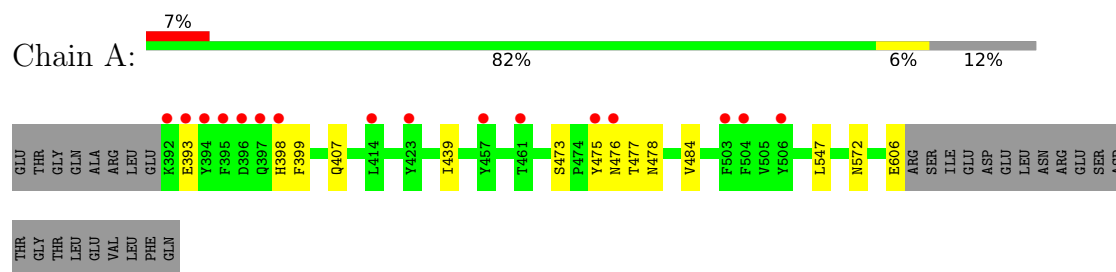
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	13	Total	O	0	0
			13	13		
8	B	31	Total	O	0	0
			31	31		

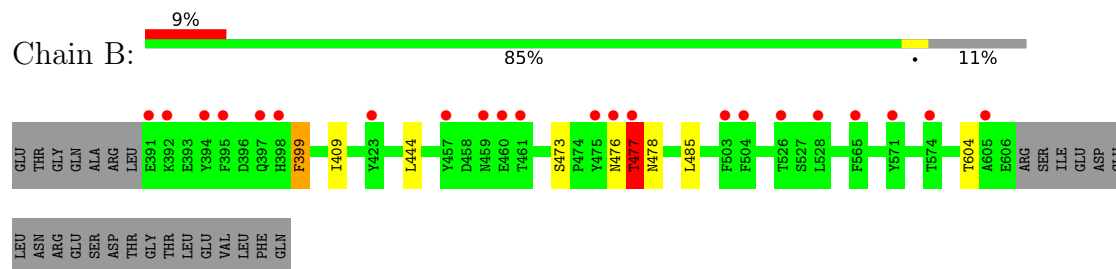
3 Residue-property plots [i](#)

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 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: Niemann-Pick C1 protein



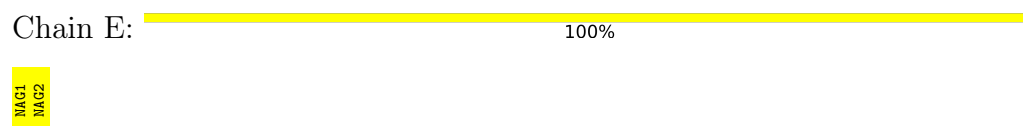
- Molecule 1: Niemann-Pick C1 protein



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



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



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

Chain F:  100%

NAG1
NAG2

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

Chain G:  50% 50%

NAG1
NAG2

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

Chain H:  50% 50%

NAG1
NAG2

- Molecule 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

Chain D:  100%

NAG1
NAG2
BMA3
MAN4

- Molecule 4: 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

Chain I:  40% 60%

NAG1
NAG2
BMA3
MAN4
MAN5

- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  33% 67%

NAG1
NAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	87.89Å 115.91Å 147.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.03 – 2.45 70.03 – 2.45	Depositor EDS
% Data completeness (in resolution range)	100.0 (70.03-2.45) 100.0 (70.03-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.224 , 0.238 0.226 , 0.237	Depositor DCC
R_{free} test set	1385 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	63.4	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3943	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.82% 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

5.1 Standard geometry

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

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	0.42	0/1844	0.63	0/2526
1	B	0.42	0/1853	0.62	1/2538 (0.0%)
All	All	0.42	0/3697	0.63	1/5064 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	399	PHE	CB-CG-CD2	-5.48	116.97	120.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	399	PHE	Peptide

5.2 Too-close contacts

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

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1787	0	1677	10	0
1	B	1796	0	1685	6	0
2	C	28	0	25	0	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	1	0
2	H	28	0	25	0	0
3	D	50	0	43	0	0
4	I	61	0	52	0	0
5	J	39	0	34	0	0
6	A	14	0	13	0	0
7	A	6	0	0	0	0
7	B	6	0	0	0	0
8	A	13	0	0	0	0
8	B	31	0	0	0	0
All	All	3943	0	3629	16	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 2.

All (16) 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:477[A]:THR:HG22	1:A:478[A]:ASN:H	1.10	1.10
1:A:475[A]:TYR:O	1:A:476[A]:ASN:O	1.86	0.92
1:A:477[A]:THR:HG22	1:A:478[A]:ASN:N	1.87	0.88
1:B:476[A]:ASN:OD1	1:B:478[A]:ASN:N	2.13	0.75
1:A:477[A]:THR:CG2	1:A:478[A]:ASN:H	1.86	0.68
1:A:475[A]:TYR:O	1:A:476[A]:ASN:C	2.34	0.65
1:A:476[B]:ASN:OD1	1:A:478[B]:ASN:N	2.15	0.59
1:A:393:GLU:HG2	1:A:407:GLN:HE22	1.70	0.56
1:B:409:ILE:CD1	1:B:604:THR:HG22	2.39	0.52
1:A:476[B]:ASN:OD1	1:A:477[B]:THR:N	2.47	0.47
1:A:477[A]:THR:CG2	1:A:478[A]:ASN:N	2.57	0.47
1:B:476[B]:ASN:O	1:B:477[B]:THR:HB	2.14	0.47
1:A:484:VAL:HG11	1:A:547:LEU:HD21	1.98	0.46
1:B:476[A]:ASN:HD21	2:G:1:NAG:H82	1.81	0.44
1:B:476[A]:ASN:OD1	1:B:477[A]:THR:N	2.51	0.43
1:B:444:LEU:HA	1:B:485:LEU:HD13	2.02	0.41

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 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	218/244 (89%)	207 (95%)	11 (5%)	0	100	100
1	B	219/244 (90%)	210 (96%)	7 (3%)	2 (1%)	14	18
All	All	437/488 (90%)	417 (95%)	18 (4%)	2 (0%)	44	32

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	477[A]	THR
1	B	477[B]	THR

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	198/219 (90%)	192 (97%)	6 (3%)	36	51
1	B	199/219 (91%)	196 (98%)	3 (2%)	60	74
All	All	397/438 (91%)	388 (98%)	9 (2%)	48	61

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	398	HIS

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Mol	Chain	Res	Type
1	A	399	PHE
1	A	439	ILE
1	A	473	SER
1	A	572	ASN
1	A	606	GLU
1	B	473	SER
1	B	477[A]	THR
1	B	477[B]	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	553	GLN
1	B	589	ASN

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 ⓘ

22 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	1,2	14,14,15	0.48	0	17,19,21	1.45	3 (17%)
2	NAG	C	2	2	14,14,15	0.40	0	17,19,21	0.97	0
3	NAG	D	1	1,3	14,14,15	0.85	1 (7%)	17,19,21	2.62	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	D	2	3	14,14,15	0.38	0	17,19,21	1.08	2 (11%)
3	BMA	D	3	3	11,11,12	0.44	0	15,15,17	1.10	1 (6%)
3	MAN	D	4	3	11,11,12	0.55	0	15,15,17	2.47	3 (20%)
2	NAG	E	1	1,2	14,14,15	0.34	0	17,19,21	1.54	3 (17%)
2	NAG	E	2	2	14,14,15	0.63	0	17,19,21	1.09	1 (5%)
2	NAG	F	1	1,2	14,14,15	0.34	0	17,19,21	0.89	1 (5%)
2	NAG	F	2	2	14,14,15	0.30	0	17,19,21	0.90	1 (5%)
2	NAG	G	1	1,2	14,14,15	0.37	0	17,19,21	0.82	0
2	NAG	G	2	2	14,14,15	0.33	0	17,19,21	0.74	0
2	NAG	H	1	1,2	14,14,15	0.44	0	17,19,21	1.24	1 (5%)
2	NAG	H	2	2	14,14,15	0.48	0	17,19,21	0.80	0
4	NAG	I	1	4,1	14,14,15	0.24	0	17,19,21	0.85	0
4	NAG	I	2	4	14,14,15	0.23	0	17,19,21	0.70	0
4	BMA	I	3	4	11,11,12	0.48	0	15,15,17	1.44	3 (20%)
4	MAN	I	4	4	11,11,12	0.34	0	15,15,17	0.90	1 (6%)
4	MAN	I	5	4	11,11,12	0.39	0	15,15,17	0.97	1 (6%)
5	NAG	J	1	5,1	14,14,15	0.31	0	17,19,21	0.97	1 (5%)
5	NAG	J	2	5	14,14,15	0.40	0	17,19,21	0.67	0
5	BMA	J	3	5	11,11,12	0.40	0	15,15,17	1.61	2 (13%)

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

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1
4	NAG	I	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	0/6/23/26	0/1/1/1
4	BMA	I	3	4	-	0/2/19/22	0/1/1/1
4	MAN	I	4	4	-	0/2/19/22	0/1/1/1
4	MAN	I	5	4	-	0/2/19/22	0/1/1/1
5	NAG	J	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	J	2	5	-	2/6/23/26	0/1/1/1
5	BMA	J	3	5	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1	NAG	C1-C2	2.70	1.56	1.52

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	NAG	C1-O5-C5	10.01	125.76	112.19
3	D	4	MAN	C1-O5-C5	7.48	122.33	112.19
5	J	3	BMA	C1-O5-C5	5.42	119.54	112.19
2	C	1	NAG	O5-C1-C2	-4.25	104.58	111.29
3	D	4	MAN	C1-C2-C3	4.14	114.75	109.67
3	D	4	MAN	O5-C1-C2	3.68	116.44	110.77
4	I	3	BMA	C3-C4-C5	3.51	116.50	110.24
2	E	1	NAG	C1-O5-C5	3.35	116.73	112.19
2	E	1	NAG	O4-C4-C3	-3.29	102.75	110.35
2	H	1	NAG	O4-C4-C3	-2.87	103.72	110.35
2	E	2	NAG	C4-C3-C2	2.82	115.15	111.02
2	F	2	NAG	C1-O5-C5	2.82	116.01	112.19
3	D	3	BMA	O5-C1-C2	-2.77	106.50	110.77
4	I	4	MAN	C1-O5-C5	2.75	115.92	112.19
2	C	1	NAG	C4-C3-C2	2.69	114.96	111.02
3	D	2	NAG	C1-O5-C5	2.69	115.83	112.19
4	I	5	MAN	C1-O5-C5	2.67	115.81	112.19
2	E	1	NAG	O5-C1-C2	-2.43	107.45	111.29
3	D	1	NAG	O5-C5-C4	2.41	116.70	110.83
2	F	1	NAG	O5-C1-C2	-2.39	107.51	111.29
2	C	1	NAG	C1-O5-C5	2.27	115.27	112.19
3	D	2	NAG	O5-C1-C2	-2.20	107.81	111.29
5	J	3	BMA	C1-C2-C3	2.20	112.37	109.67
4	I	3	BMA	O3-C3-C2	-2.19	105.81	109.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	3	BMA	C2-C3-C4	2.13	114.58	110.89
5	J	1	NAG	C1-O5-C5	2.11	115.06	112.19

There are no chirality outliers.

All (13) torsion outliers are listed below:

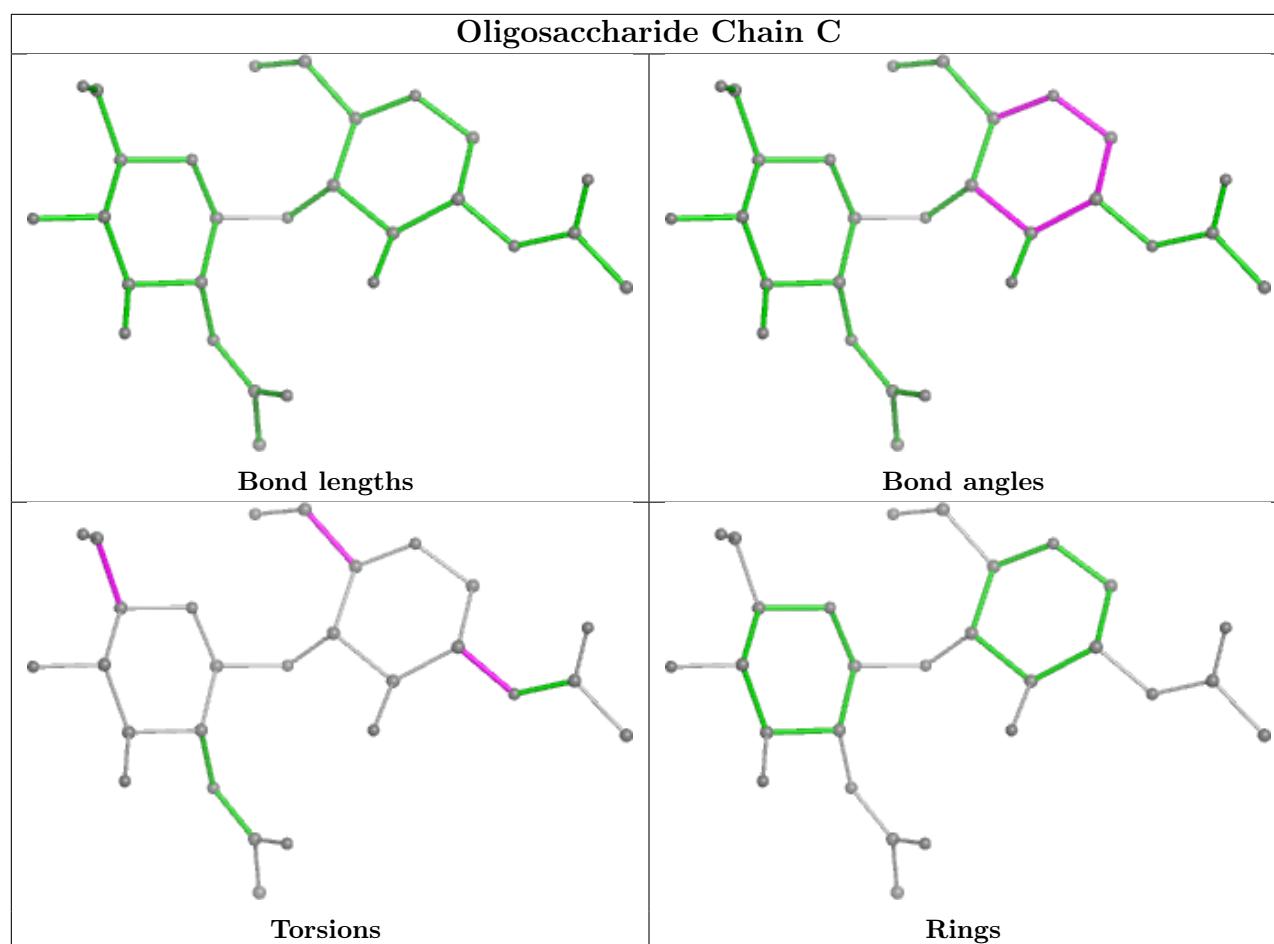
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	C4-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
5	J	2	NAG	O5-C5-C6-O6
2	E	2	NAG	C1-C2-N2-C7
2	F	2	NAG	O5-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
5	J	2	NAG	C4-C5-C6-O6
2	E	2	NAG	C3-C2-N2-C7
2	C	1	NAG	C3-C2-N2-C7
3	D	3	BMA	C4-C5-C6-O6

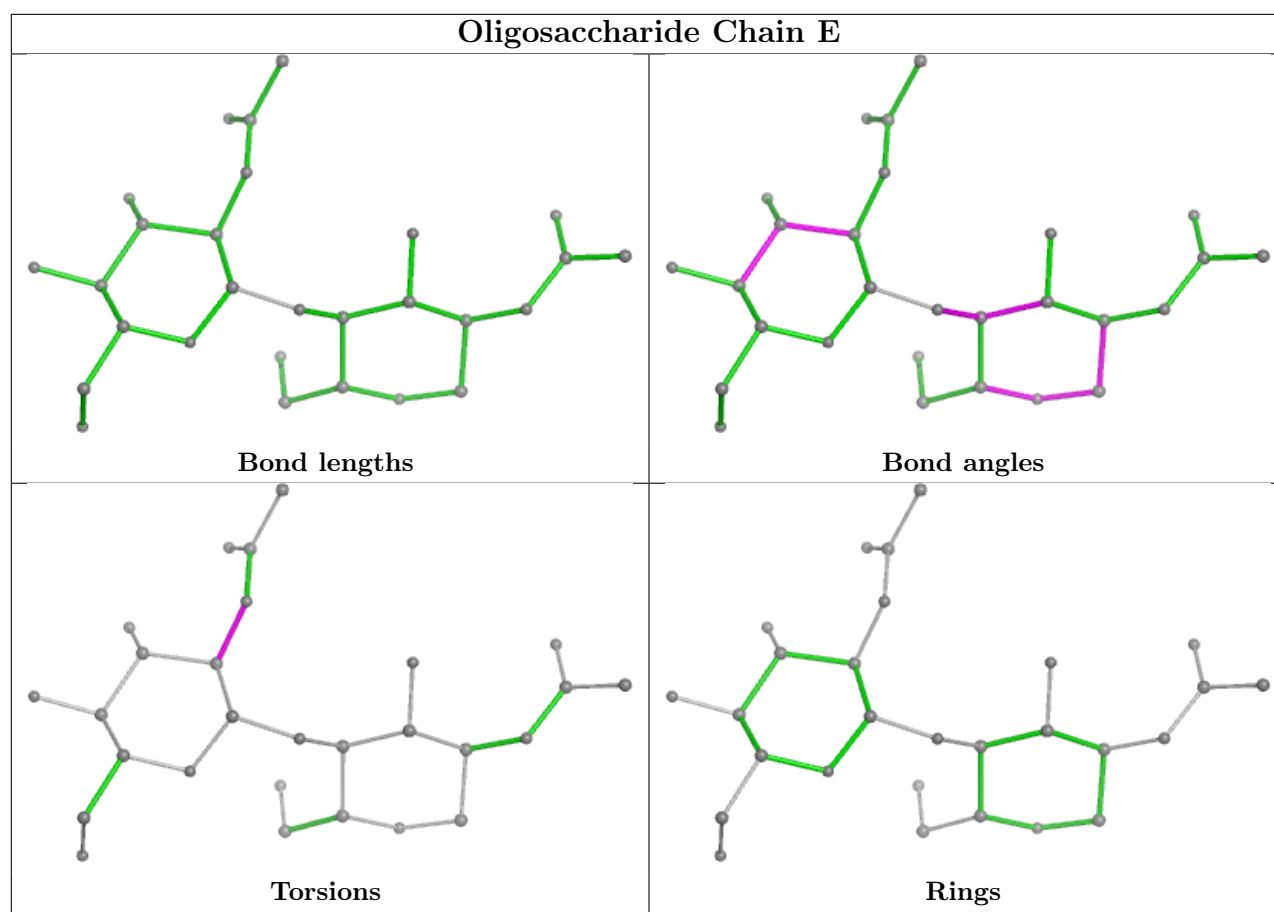
There are no ring outliers.

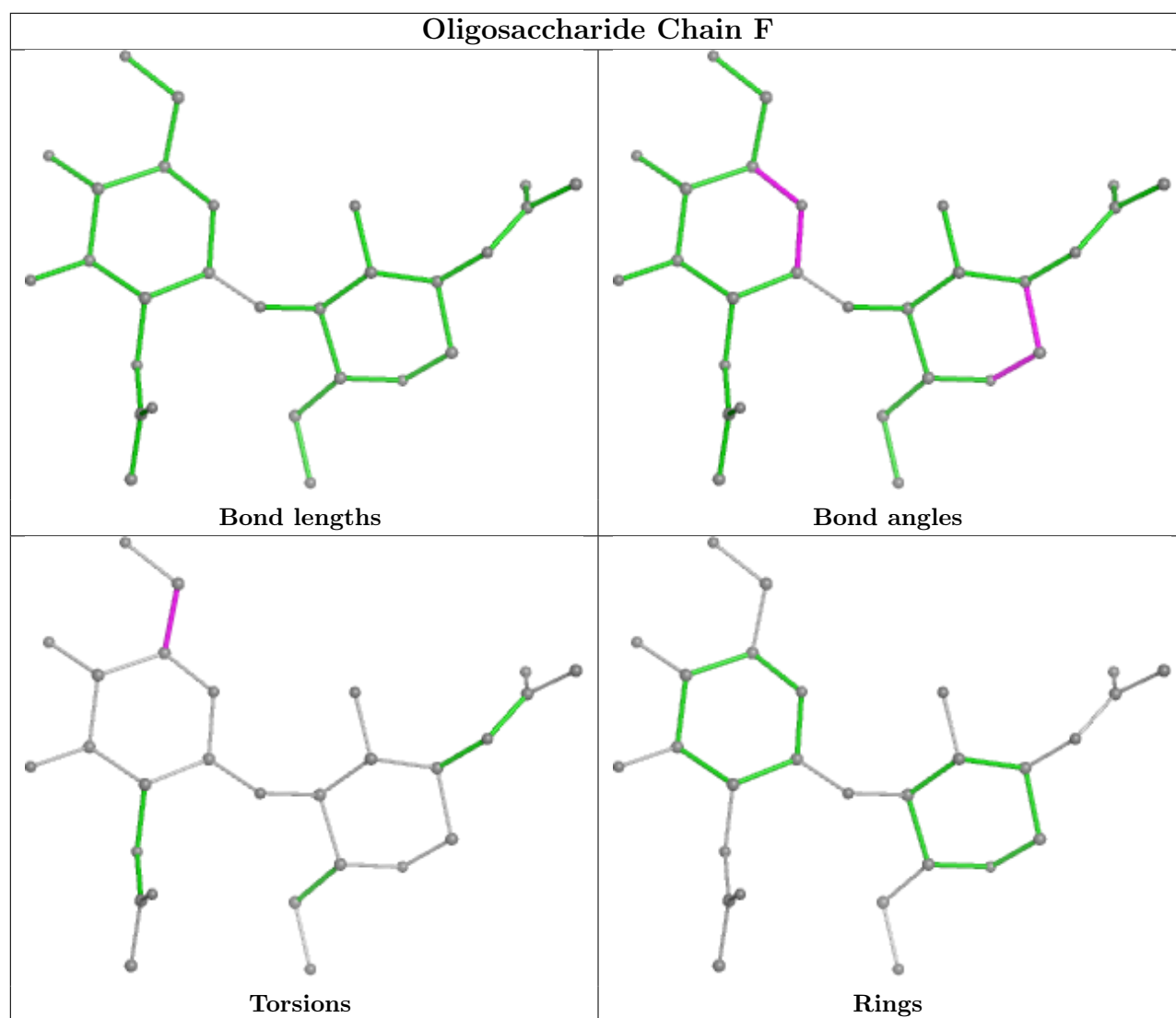
1 monomer is involved in 1 short contact:

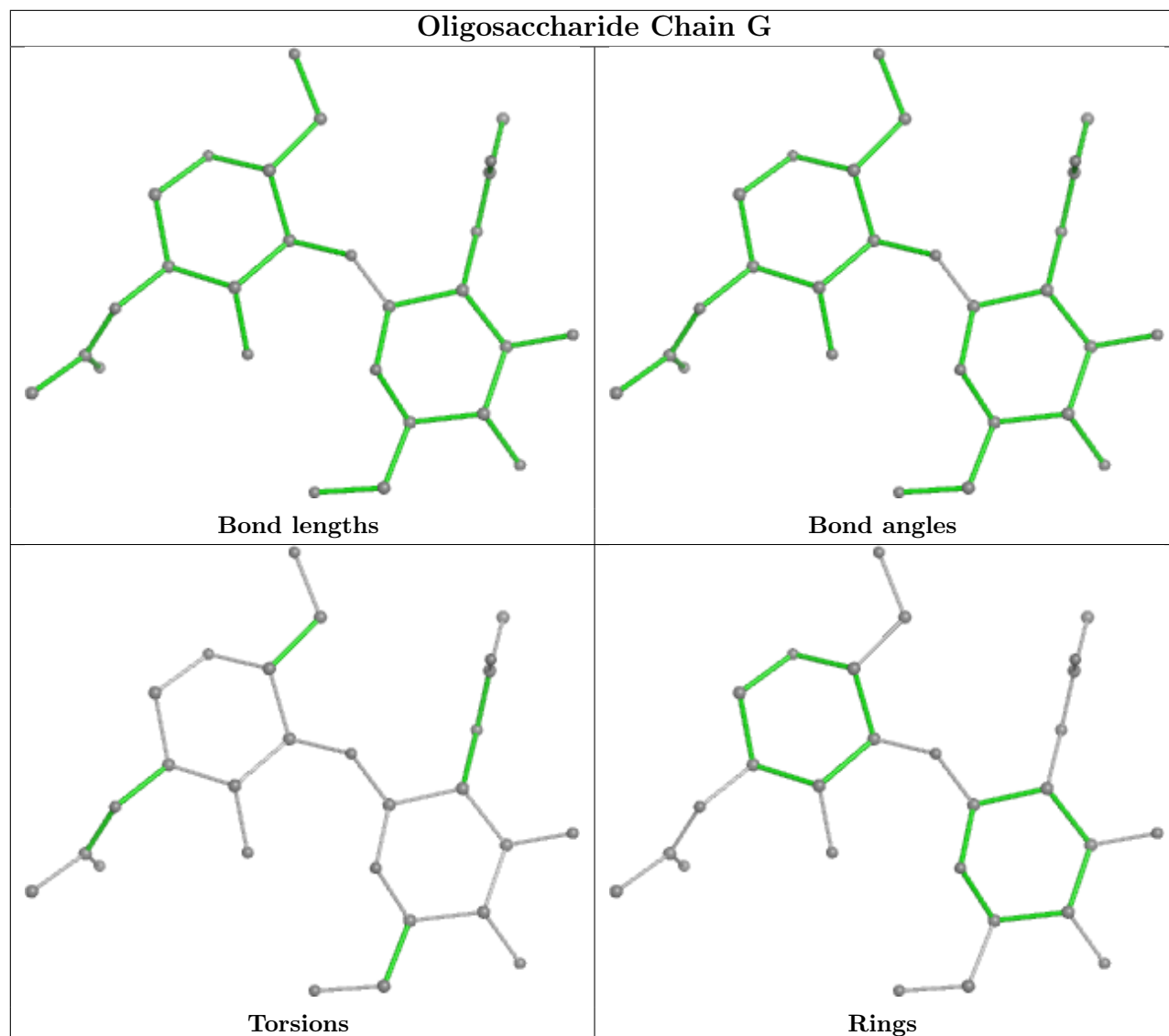
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	1	NAG	1	0

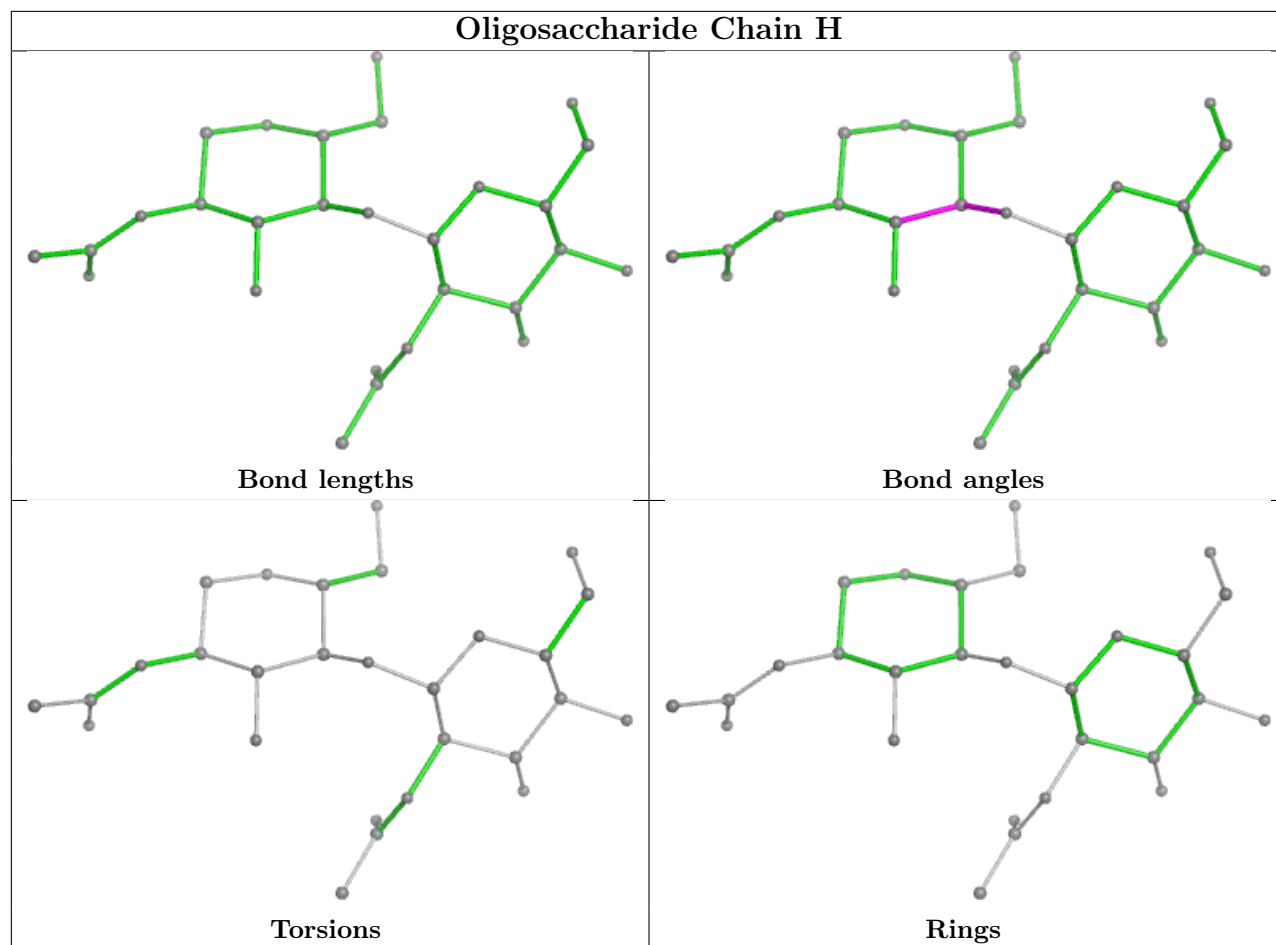
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

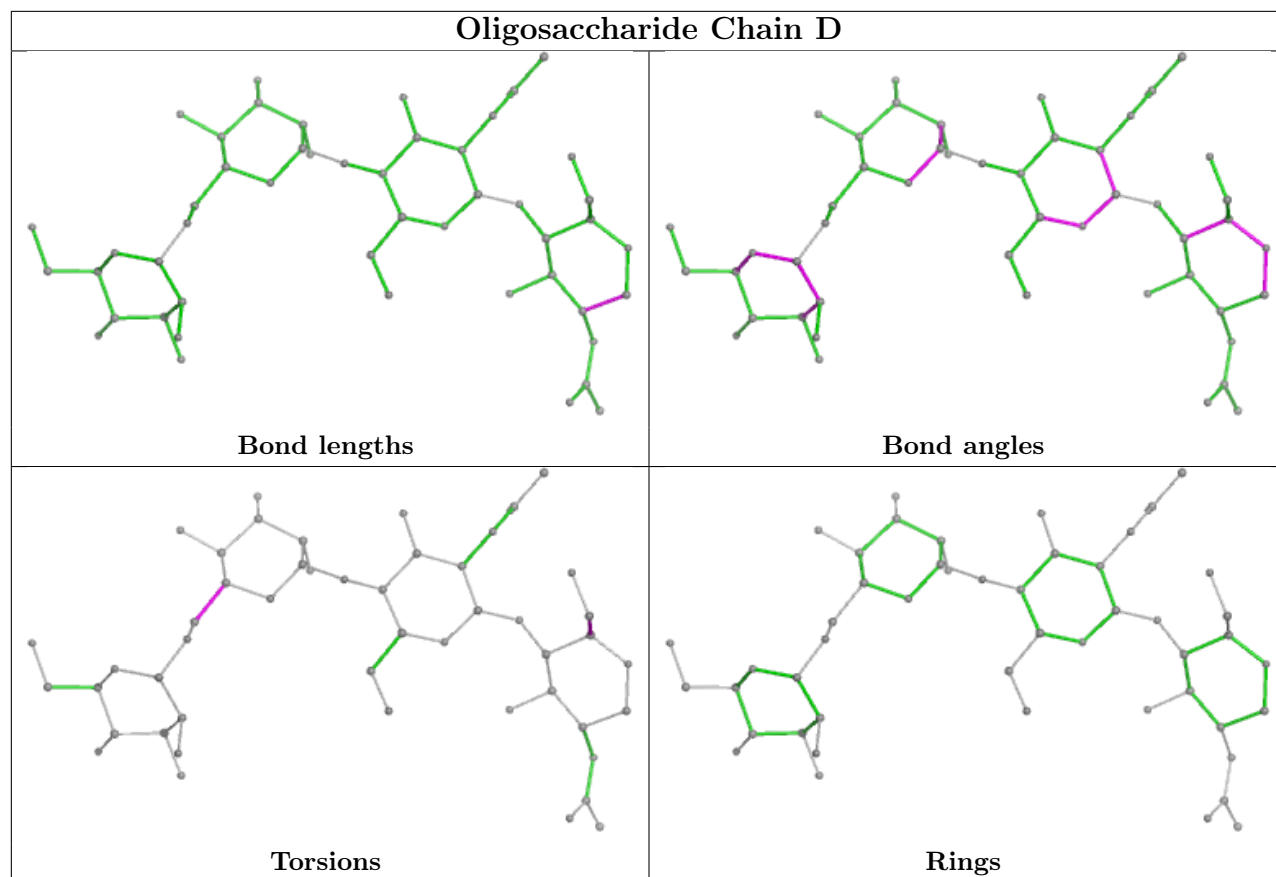


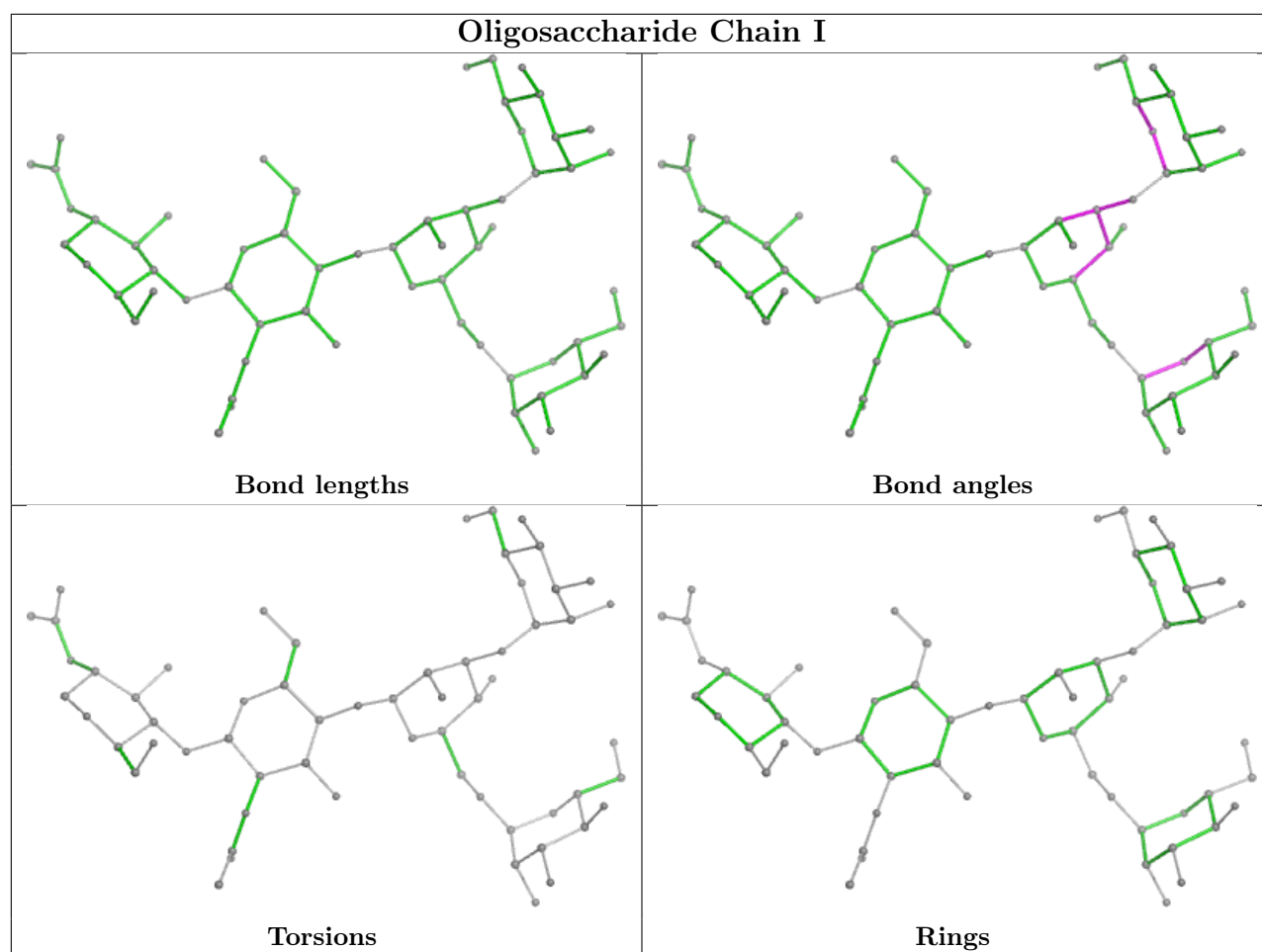


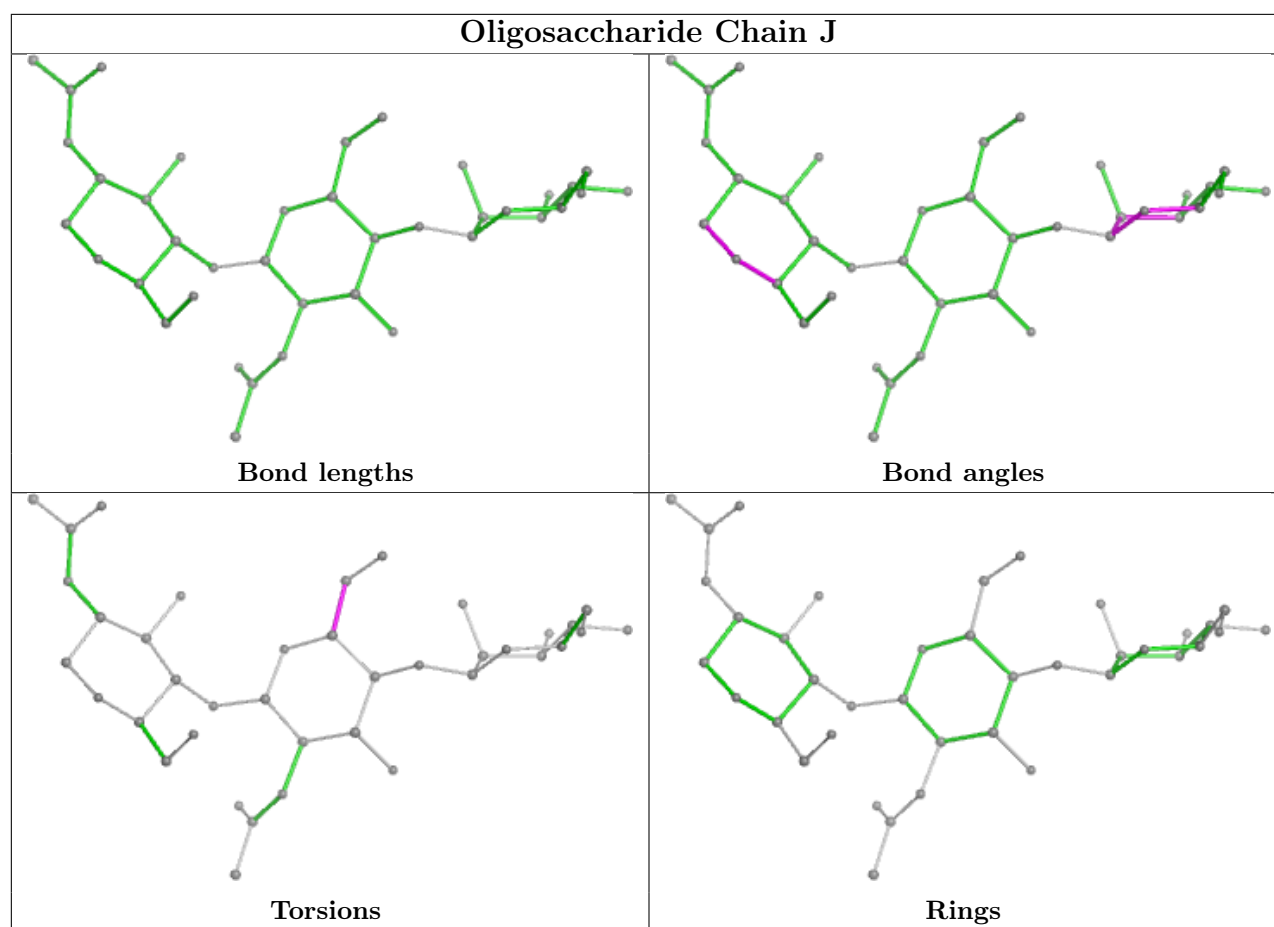












5.6 Ligand geometry [i](#)

5 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	SCN	B	714	-	1,2,2	0.29	0	0,1,1	-	-
6	NAG	A	707	1	14,14,15	0.49	0	17,19,21	1.18	0
7	SCN	B	713	-	1,2,2	0.35	0	0,1,1	-	-
7	SCN	A	712	-	1,2,2	0.40	0	0,1,1	-	-
7	SCN	A	713	-	1,2,2	0.37	0	0,1,1	-	-

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
6	NAG	A	707	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.

There are no torsion outliers.

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	215/244 (88%)	0.49	16 (7%)	22	23	40, 65, 119, 189	5 (2%)
1	B	216/244 (88%)	0.43	22 (10%)	13	14	35, 61, 113, 146	5 (2%)
All	All	431/488 (88%)	0.46	38 (8%)	17	17	35, 63, 117, 189	10 (2%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	477[A]	THR	7.6
1	A	475[A]	TYR	6.4
1	B	475[A]	TYR	5.5
1	A	476[A]	ASN	5.5
1	B	605	ALA	5.1
1	A	392	LYS	5.1
1	A	395	PHE	4.9
1	A	393	GLU	4.2
1	B	394	TYR	3.8
1	B	391	GLU	3.8
1	B	574	THR	3.6
1	B	528	LEU	3.6
1	B	503	PHE	3.6
1	A	394	TYR	3.6
1	B	392	LYS	3.4
1	A	423	TYR	3.4
1	A	503	PHE	3.4
1	B	398	HIS	3.3
1	A	398	HIS	3.3
1	A	397	GLN	3.3
1	B	397	GLN	3.2
1	A	457	TYR	3.1
1	A	506	TYR	3.0
1	B	526	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	504	PHE	2.9
1	B	565	PHE	2.9
1	B	395	PHE	2.8
1	B	457	TYR	2.8
1	B	476[A]	ASN	2.5
1	B	461	THR	2.5
1	A	414	LEU	2.4
1	B	459	ASN	2.4
1	B	571	TYR	2.3
1	B	504	PHE	2.3
1	B	460	GLU	2.2
1	A	461	THR	2.1
1	A	396	ASP	2.1
1	B	423	TYR	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.

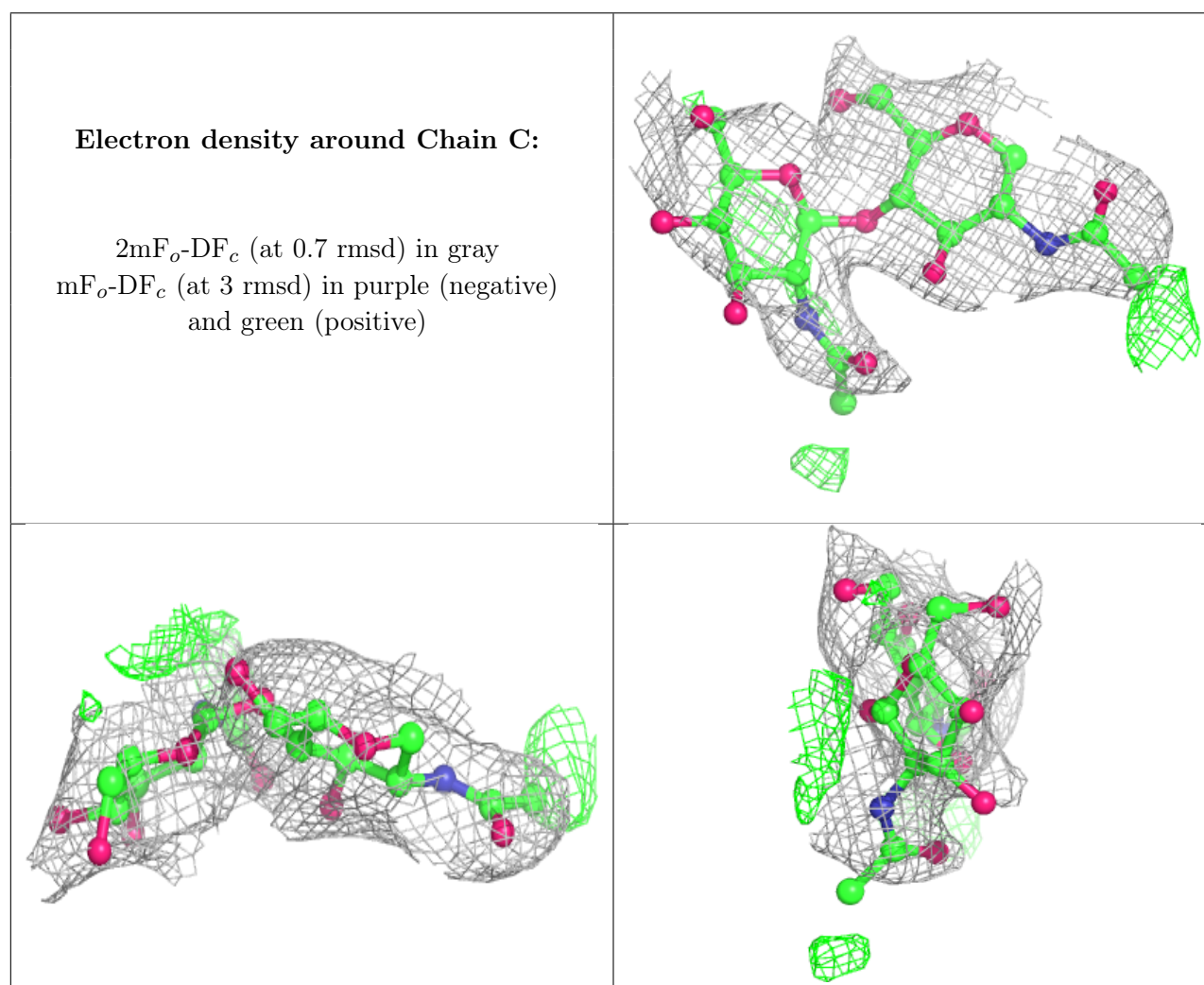
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MAN	I	4	11/12	0.28	0.18	132,143,149,150	0
3	MAN	D	4	11/12	0.36	0.17	133,136,139,142	0
4	MAN	I	5	11/12	0.37	0.17	133,136,137,139	0
5	BMA	J	3	11/12	0.39	0.22	130,135,137,137	0
2	NAG	C	2	14/15	0.46	0.18	126,129,131,131	0
2	NAG	F	2	14/15	0.51	0.17	128,133,143,146	0
3	BMA	D	3	11/12	0.54	0.14	111,117,123,128	0
2	NAG	H	2	14/15	0.58	0.16	131,134,137,137	0
2	NAG	E	2	14/15	0.61	0.18	112,117,123,124	0
2	NAG	C	1	14/15	0.64	0.14	97,102,110,122	0
4	BMA	I	3	11/12	0.67	0.14	119,126,134,139	0
5	NAG	J	2	14/15	0.76	0.17	87,95,106,120	0
2	NAG	F	1	14/15	0.76	0.17	90,105,113,121	0
2	NAG	G	1	14/15	0.77	0.16	82,98,101,108	0

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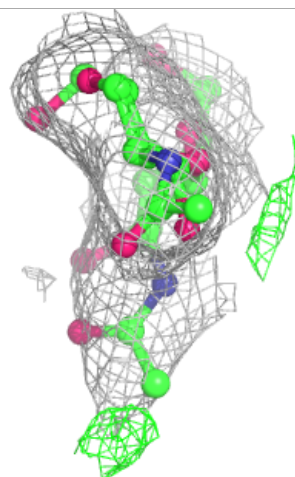
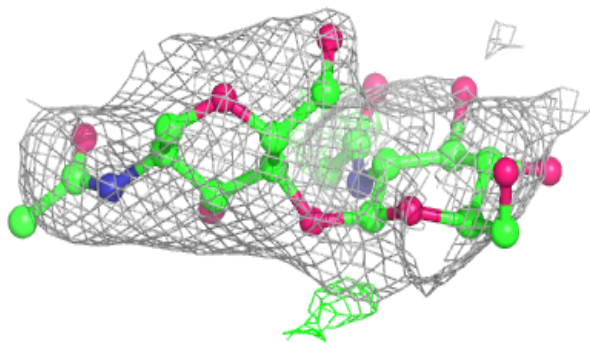
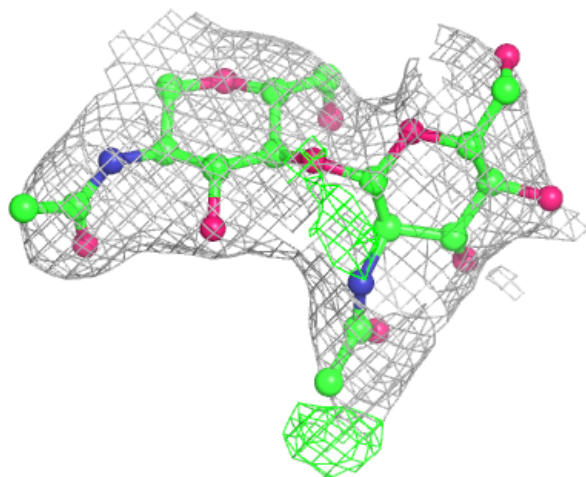
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	G	2	14/15	0.78	0.15	117,120,125,125	0
2	NAG	H	1	14/15	0.82	0.12	101,108,116,124	0
3	NAG	D	1	14/15	0.88	0.14	65,69,74,77	0
4	NAG	I	2	14/15	0.92	0.10	78,86,95,108	0
2	NAG	E	1	14/15	0.93	0.11	78,83,88,100	0
3	NAG	D	2	14/15	0.93	0.10	81,87,95,102	0
4	NAG	I	1	14/15	0.94	0.11	65,68,71,76	0
5	NAG	J	1	14/15	0.95	0.09	64,70,75,84	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.



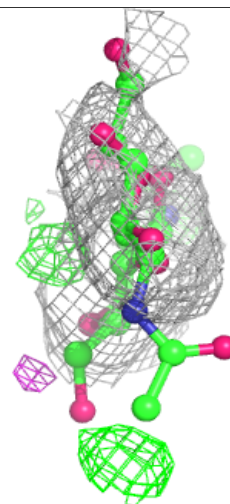
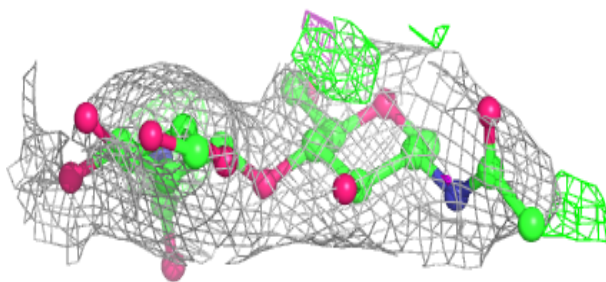
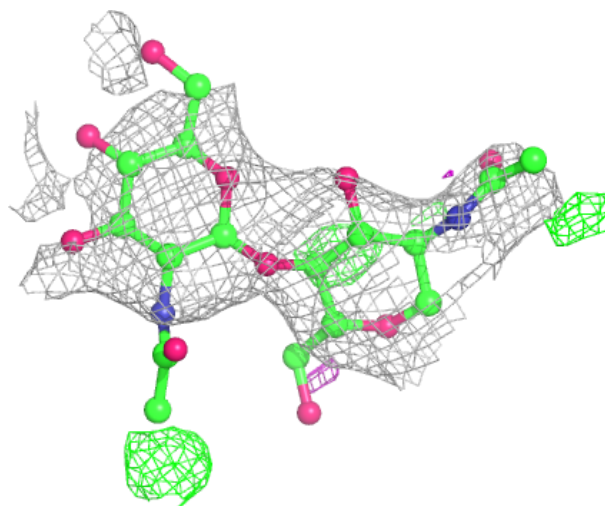
Electron density around Chain E:

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



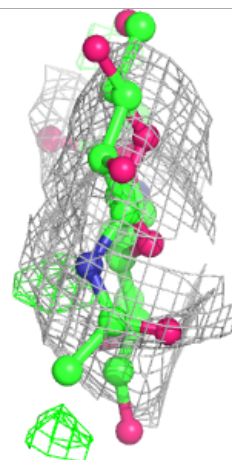
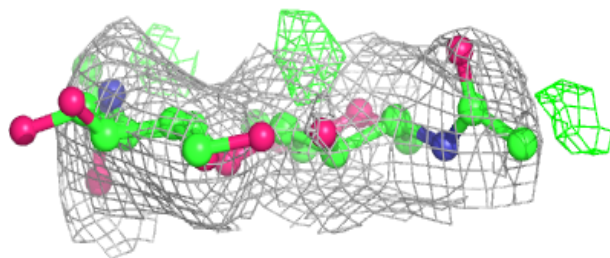
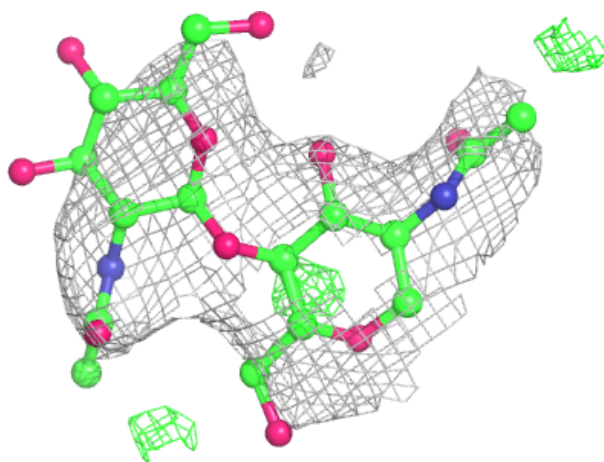
Electron density around Chain F:

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



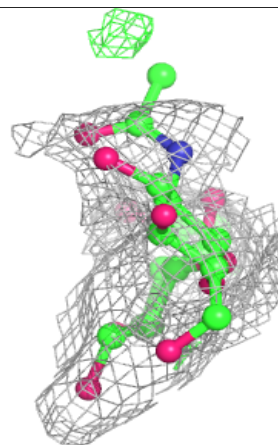
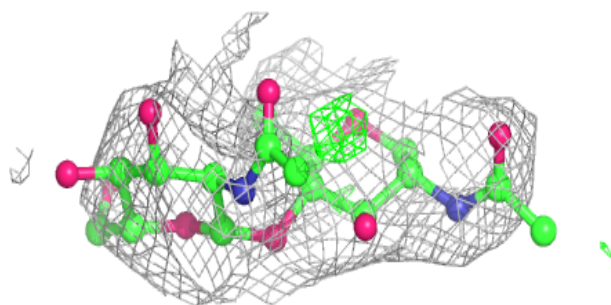
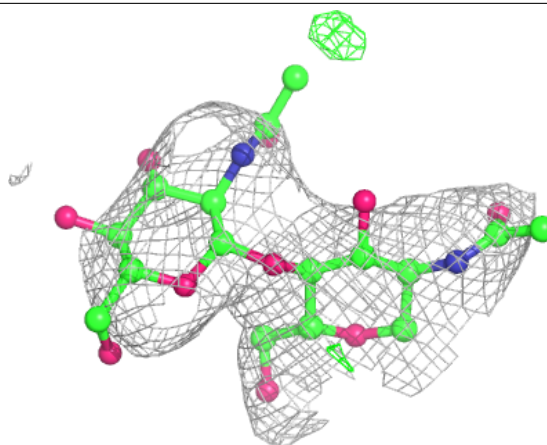
Electron density around Chain G:

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

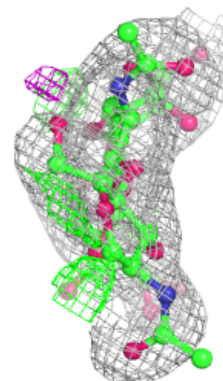
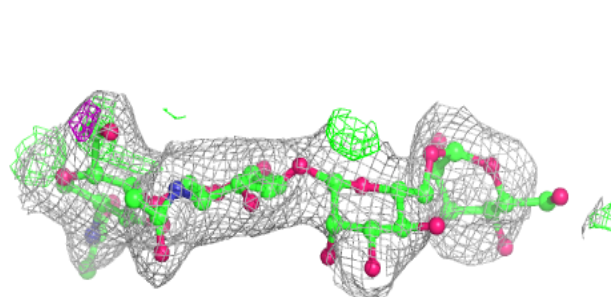
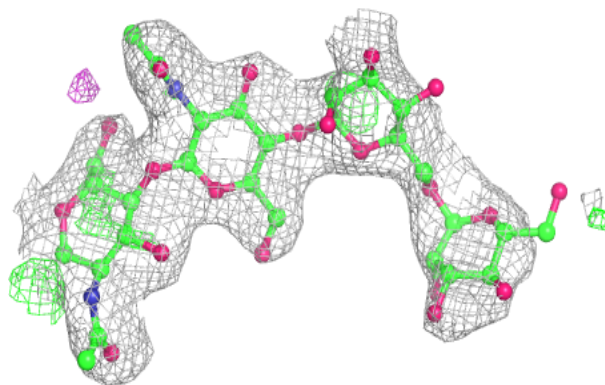


Electron density around Chain H:

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

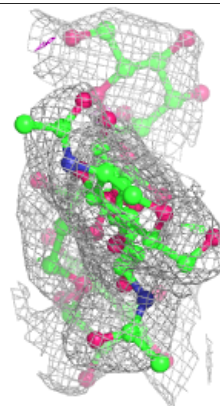
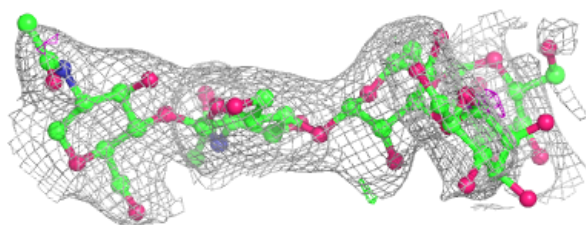
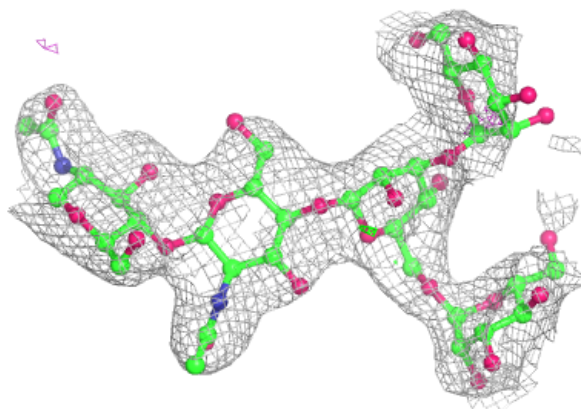
**Electron density around Chain D:**

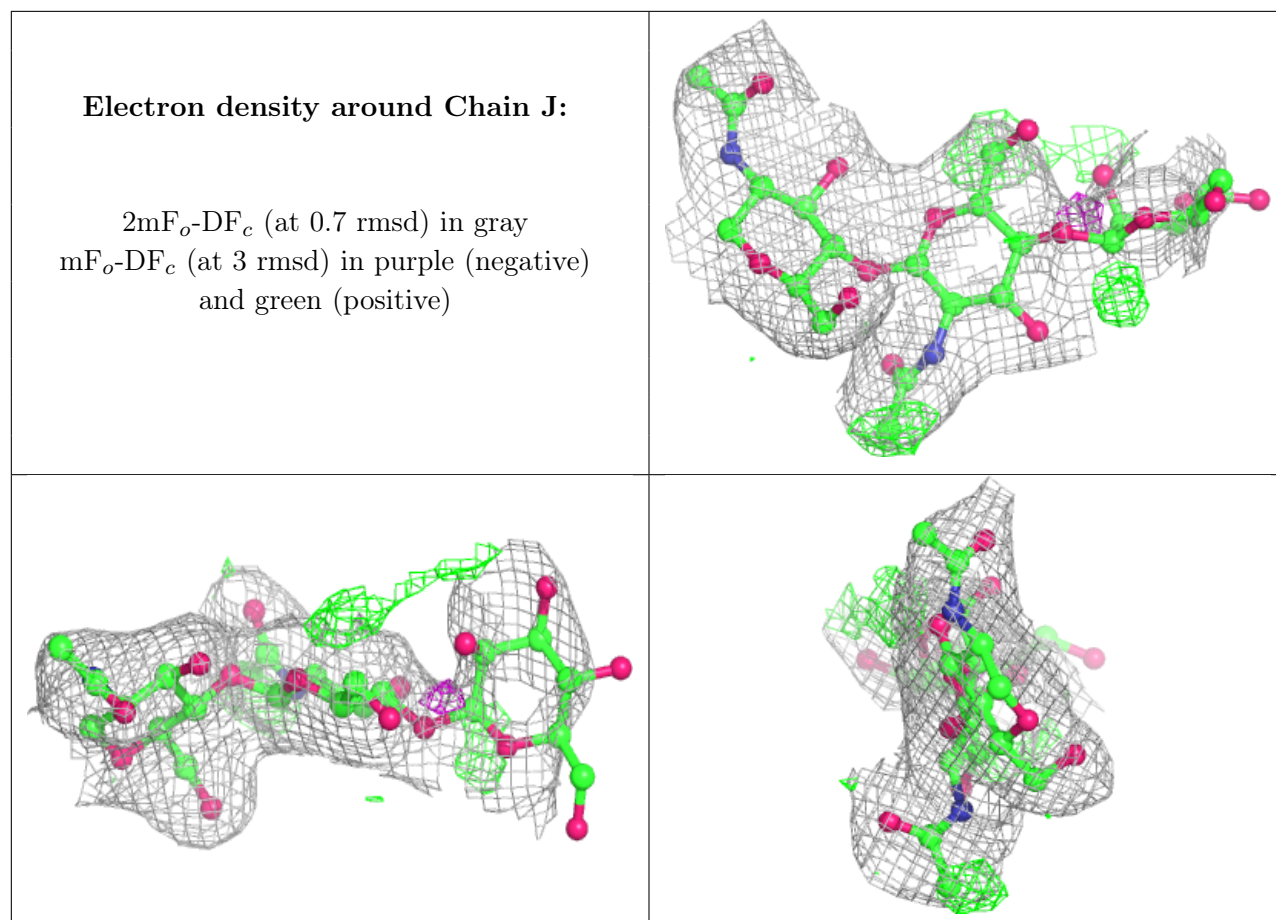
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain I:

$2mF_o-DF_c$ (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
6	NAG	A	707	14/15	0.62	0.17	112,117,125,128	0
7	SCN	A	712	3/3	0.79	0.22	94,94,99,101	0
7	SCN	A	713	3/3	0.82	0.20	88,88,94,96	0
7	SCN	B	714	3/3	0.82	0.26	102,102,105,112	0
7	SCN	B	713	3/3	0.87	0.14	98,98,99,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.