



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 28, 2026 – 03:52 AM UTC

PDB ID : 8TJ4 / pdb_00008tj4
Title : CRYSTAL STRUCTURE OF THE A/Bangkok/1/1979(H3N2) INFLUENZA VIRUS HEMAGGLUTININ WITH HUMAN RECEPTOR ANALOG 6'-SLNLN
Authors : Wu, N.C.; Zhu, X.; Wilson, I.A.
Deposited on : 2023-07-20
Resolution : 1.95 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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

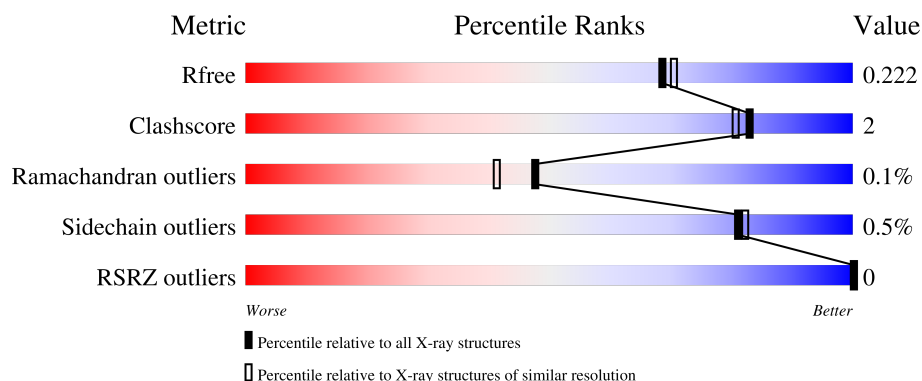
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

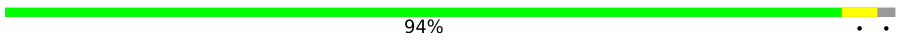
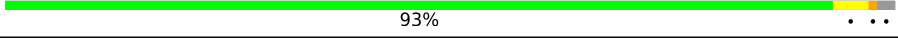
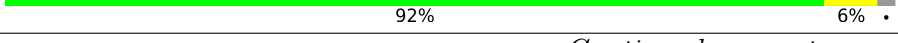
The reported resolution of this entry is 1.95 Å.

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.



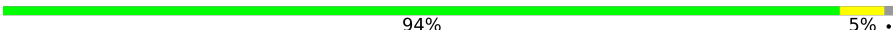
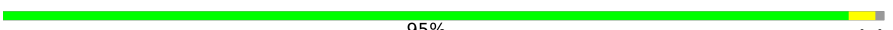
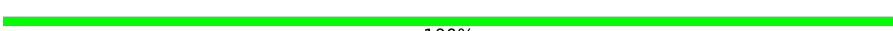
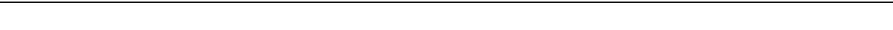
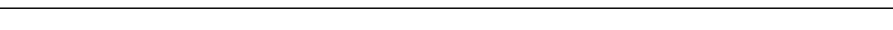
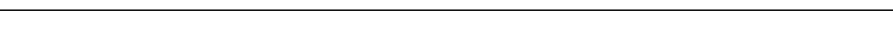
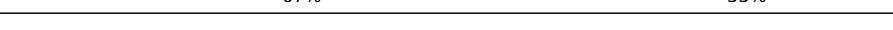
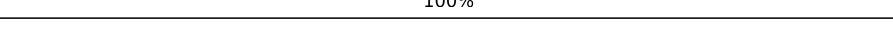
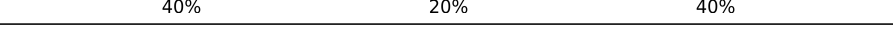
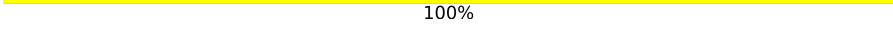
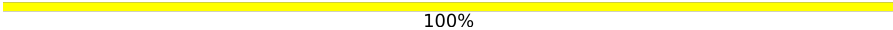
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	323	 94%
1	C	323	 89% 9%
1	E	323	 93%
1	G	323	 92% 6%

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Mol	Chain	Length	Quality of chain
2	B	174	 94% 5% .
2	D	174	 92% 7% .
2	F	174	 95% . .
2	H	174	 93% 6% ..
3	I	2	 100%
3	L	2	 50% 50%
3	O	2	 100%
3	S	2	 100%
3	T	2	 50% 50%
3	U	2	 100%
3	V	2	 50% 50%
3	W	2	 50% 50%
3	X	2	 100%
4	J	3	 67% 33%
4	M	3	 67% 33%
4	N	3	 100%
5	K	5	 60% 40%
5	Q	5	 40% 20% 40%
5	R	5	 60% 40%
6	P	4	 75% 25%
6	Y	4	 100%
7	Z	6	 33% 50% 17%
8	a	3	 33% 67%
8	c	3	 100%
9	b	2	 50% 50%

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Mol	Chain	Length	Quality of chain
9	d	2	50% 50%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 18202 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 Hemagglutinin HA1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	0	0
			2460	1543	434	471	12			
1	C	317	Total	C	N	O	S	0	0	0
			2460	1543	434	471	12			
1	E	317	Total	C	N	O	S	0	0	0
			2460	1543	434	471	12			
1	G	317	Total	C	N	O	S	0	0	0
			2460	1543	434	471	12			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	ALA	-	expression tag	UNP P03441
A	8	ASP	-	expression tag	UNP P03441
A	9	PRO	-	expression tag	UNP P03441
A	10	GLY	-	expression tag	UNP P03441
A	294	PHE	UNK	conflict	UNP P03441
C	7	ALA	-	expression tag	UNP P03441
C	8	ASP	-	expression tag	UNP P03441
C	9	PRO	-	expression tag	UNP P03441
C	10	GLY	-	expression tag	UNP P03441
C	294	PHE	UNK	conflict	UNP P03441
E	7	ALA	-	expression tag	UNP P03441
E	8	ASP	-	expression tag	UNP P03441
E	9	PRO	-	expression tag	UNP P03441
E	10	GLY	-	expression tag	UNP P03441
E	294	PHE	UNK	conflict	UNP P03441
G	7	ALA	-	expression tag	UNP P03441
G	8	ASP	-	expression tag	UNP P03441
G	9	PRO	-	expression tag	UNP P03441
G	10	GLY	-	expression tag	UNP P03441
G	294	PHE	UNK	conflict	UNP P03441

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	172	Total	C	N	O	S	0	0	0
			1386	860	243	277	6			
2	D	172	Total	C	N	O	S	0	0	0
			1386	860	243	277	6			
2	F	172	Total	C	N	O	S	0	0	0
			1392	863	246	277	6			
2	H	172	Total	C	N	O	S	0	0	0
			1392	863	246	277	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	18	VAL	UNK	conflict	UNP P03441
B	19	ASP	UNK	conflict	UNP P03441
D	18	VAL	UNK	conflict	UNP P03441
D	19	ASP	UNK	conflict	UNP P03441
F	18	VAL	UNK	conflict	UNP P03441
F	19	ASP	UNK	conflict	UNP P03441
H	18	VAL	UNK	conflict	UNP P03441
H	19	ASP	UNK	conflict	UNP P03441

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



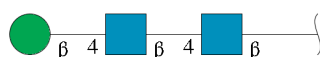
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	2	Total	C	N	O		0	0	0
			28	16	2	10				
3	L	2	Total	C	N	O		0	0	0
			28	16	2	10				
3	O	2	Total	C	N	O		0	0	0
			28	16	2	10				
3	S	2	Total	C	N	O		0	0	0
			28	16	2	10				
3	T	2	Total	C	N	O		0	0	0
			28	16	2	10				
3	U	2	Total	C	N	O		0	0	0
			28	16	2	10				

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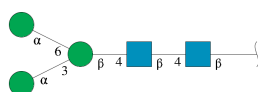
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	V	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	W	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	X	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



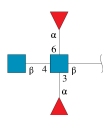
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	J	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	M	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	N	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 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.



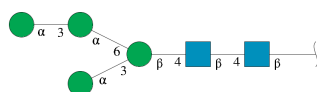
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	K	5	Total	C	N	O	0	0	0
			61	34	2	25			
5	Q	5	Total	C	N	O	0	0	0
			61	34	2	25			
5	R	5	Total	C	N	O	0	0	0
			61	34	2	25			

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



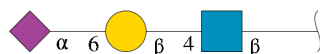
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	P	4	Total	C	N	O	0	0	0
			48	28	2	18			
6	Y	4	Total	C	N	O	0	0	0
			48	28	2	18			

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



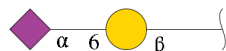
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	Z	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 8 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



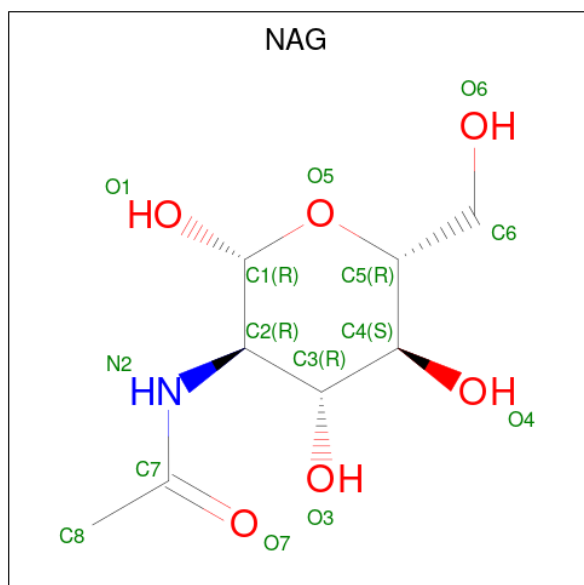
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
8	a	3	Total	C	N	O	0	0	0
			46	25	2	19			
8	c	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 9 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose.



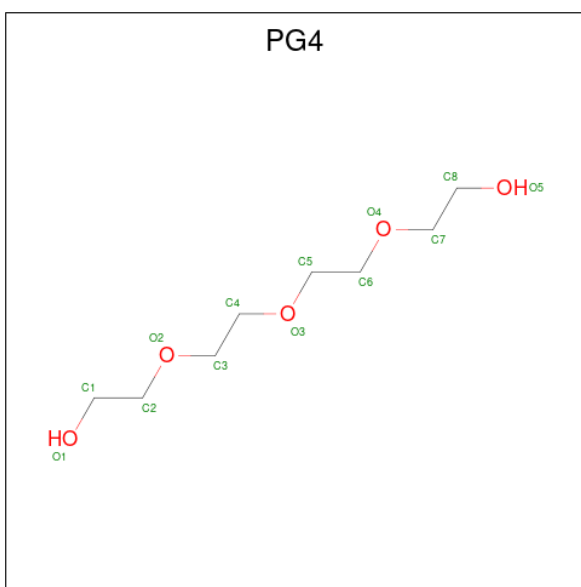
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	b	2	Total	C	N	O	0	0	0
			32	17	1	14			
9	d	2	Total	C	N	O	0	0	0
			32	17	1	14			

- Molecule 10 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).



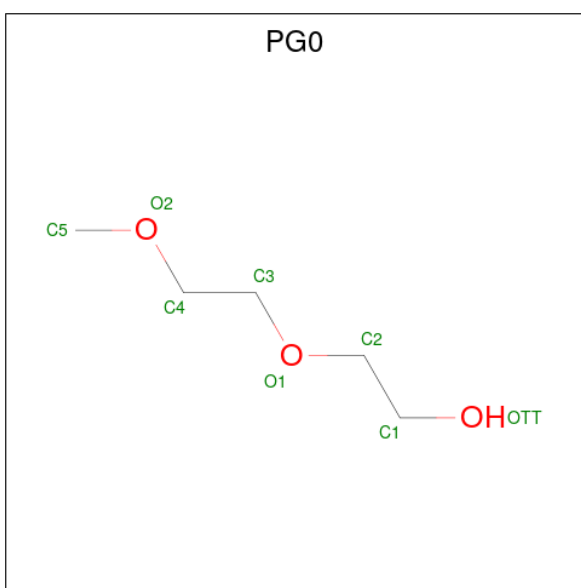
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	A	1	Total	C	N	O	0	0
			14	8	1	5		
10	B	1	Total	C	N	O	0	0
			14	8	1	5		
10	D	1	Total	C	N	O	0	0
			14	8	1	5		
10	E	1	Total	C	N	O	0	0
			14	8	1	5		
10	F	1	Total	C	N	O	0	0
			14	8	1	5		
10	G	1	Total	C	N	O	0	0
			14	8	1	5		
10	H	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 11 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



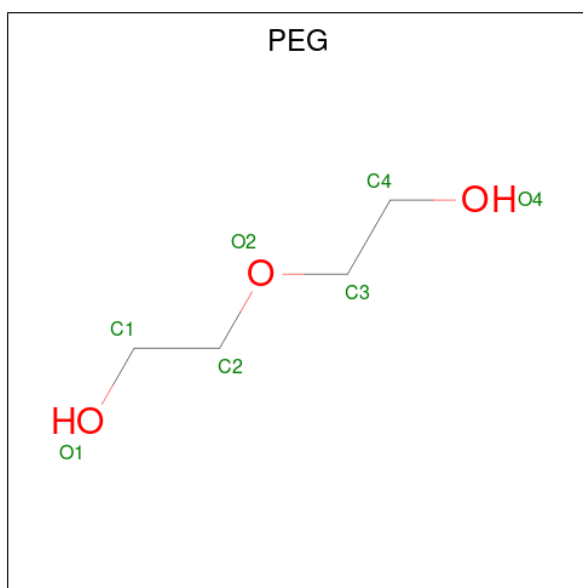
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	B	1	Total	C	O	0	0
			13	8	5		
11	F	1	Total	C	O	0	0
			13	8	5		

- Molecule 12 is 2-(2-METHOXYETHOXY)ETHANOL (CCD ID: PG0) (formula: C₅H₁₂O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	B	1	Total	C	O	0	0
			8	5	3		
12	H	1	Total	C	O	0	0
			8	5	3		

- Molecule 13 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	E	1	Total	C	O	0	0
			7	4	3		
13	F	1	Total	C	O	0	0
			7	4	3		
13	F	1	Total	C	O	0	0
			7	4	3		

- Molecule 14 is water.

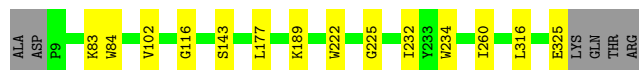
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	A	233	Total	O	0	0
			233	233		
14	B	207	Total	O	0	0
			207	207		
14	C	234	Total	O	0	0
			234	234		
14	D	216	Total	O	0	0
			216	216		
14	E	227	Total	O	0	0
			227	227		
14	F	214	Total	O	0	0
			214	214		
14	G	237	Total	O	0	0
			237	237		
14	H	201	Total	O	0	0
			201	201		

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: Hemagglutinin HA1 chain

Chain A:  94%

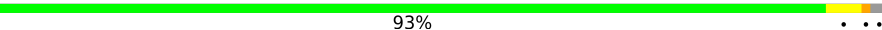


- Molecule 1: Hemagglutinin HA1 chain

Chain C:  89% 9%



- Molecule 1: Hemagglutinin HA1 chain

Chain E:  93%



- Molecule 1: Hemagglutinin HA1 chain

Chain G:  92% 6%



- Molecule 2: Hemagglutinin HA2 chain

Chain B:  94% 5%

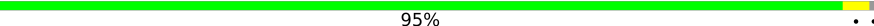


- Molecule 2: Hemagglutinin HA2 chain

Chain D:  92% 7%



- Molecule 2: Hemagglutinin HA2 chain

Chain F:  95% ..

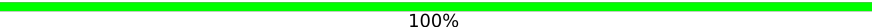


- Molecule 2: Hemagglutinin HA2 chain

Chain H:  93% 6% ..



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

Chain I:  100%



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

Chain L:  50% 50%

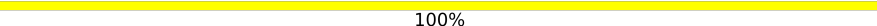


- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  100%



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

Chain S:  100%



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

Chain T:  50% 50%

MAG1
MAG2

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

Chain U:  100%

MAG1
MAG2

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

Chain V:  50% 50%

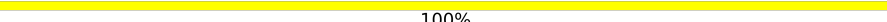
MAG1
MAG2

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

Chain W:  50% 50%

MAG1
MAG2

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

Chain X:  100%

MAG1
MAG2

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

Chain J:  67% 33%

MAG1
MAG2
BMA3

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

Chain M:  67% 33%



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

Chain N: 100%



- Molecule 5: 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 K: 60% 40%



- Molecule 5: 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 Q: 40% 20% 40%



- Molecule 5: 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 R: 60% 40%



- Molecule 6: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P: 75% 25%

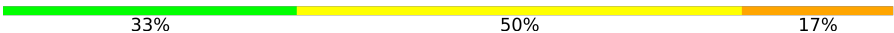


- Molecule 6: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Y: 100%

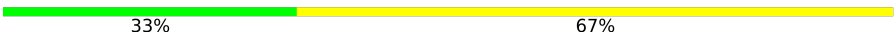
MAG1
FUC2
MAG3
FUC4

- Molecule 7: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z:  33% 50% 17%

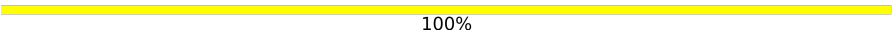
MAG1
MAG2
BMA3
MAN4
MAN5
MAN6

- Molecule 8: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain a:  33% 67%

MAG1
GAL2
STI3

- Molecule 8: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain c:  100%

MAG1
GAL2
STI3

- Molecule 9: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose

Chain b:  50% 50%

GAL3
STI2

- Molecule 9: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose

Chain d:  50% 50%

GAL3
STI2

4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	100.66Å 100.66Å 687.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.23 – 1.95 45.23 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.2 (45.23-1.95) 99.5 (45.23-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.189 , 0.223 0.189 , 0.222	Depositor DCC
R_{free} test set	9453 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.477 for -h-k,k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	18202	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.02 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.5092e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: SIA, PG4, MAN, FUC, BMA, PEG, NAG, PG0, GAL

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	0.30	0/2518	0.51	0/3423
1	C	0.32	0/2518	0.53	0/3423
1	E	0.30	0/2518	0.50	0/3423
1	G	0.33	0/2518	0.54	0/3423
2	B	0.36	0/1410	0.58	0/1896
2	D	0.43	0/1410	0.70	0/1896
2	F	0.38	0/1416	0.61	0/1903
2	H	0.47	0/1416	0.74	2/1903 (0.1%)
All	All	0.35	0/15724	0.58	2/21290 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	171	PHE	CA-C-N	5.17	131.01	121.70
2	H	171	PHE	C-N-CA	5.17	131.01	121.70

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	2460	0	2405	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2460	0	2405	15	0
1	E	2460	0	2405	9	0
1	G	2460	0	2404	10	0
2	B	1386	0	1300	7	0
2	D	1386	0	1300	7	0
2	F	1392	0	1311	4	0
2	H	1392	0	1311	8	0
3	I	28	0	25	0	0
3	L	28	0	25	0	0
3	O	28	0	25	0	0
3	S	28	0	25	1	0
3	T	28	0	25	0	0
3	U	28	0	25	0	0
3	V	28	0	25	0	0
3	W	28	0	25	1	0
3	X	28	0	25	0	0
4	J	39	0	34	0	0
4	M	39	0	34	0	0
4	N	39	0	34	0	0
5	K	61	0	52	0	0
5	Q	61	0	52	1	0
5	R	61	0	52	0	0
6	P	48	0	43	0	0
6	Y	48	0	43	2	0
7	Z	72	0	61	1	0
8	a	46	0	40	0	0
8	c	46	0	40	0	0
9	b	32	0	28	0	0
9	d	32	0	28	0	0
10	A	14	0	13	0	0
10	B	14	0	13	0	0
10	D	14	0	13	0	0
10	E	14	0	13	0	0
10	F	14	0	13	0	0
10	G	14	0	13	0	0
10	H	14	0	13	0	0
11	B	13	0	18	2	0
11	F	13	0	18	1	0
12	B	8	0	12	1	0
12	H	8	0	12	0	0
13	E	7	0	10	2	0
13	F	14	0	20	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	A	233	0	0	1	0
14	B	207	0	0	2	0
14	C	234	0	0	3	0
14	D	216	0	0	1	0
14	E	227	0	0	2	0
14	F	214	0	0	1	0
14	G	237	0	0	0	0
14	H	201	0	0	1	0
All	All	18202	0	15788	78	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 (78) 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:G:177:LEU:HB2	1:G:260:ILE:HD11	1.74	0.69
2:F:71:SER:HB3	13:F:204:PEG:H31	1.73	0.69
1:G:123:GLU:HB3	1:G:168:MET:HE3	1.77	0.67
1:A:177:LEU:HB2	1:A:260:ILE:HD11	1.76	0.66
1:C:177:LEU:HB2	1:C:260:ILE:HD11	1.79	0.65
1:G:222:TRP:CZ2	1:G:225:GLY:HA2	2.33	0.63
1:A:222:TRP:CZ2	1:A:225:GLY:HA2	2.36	0.61
1:C:222:TRP:CZ2	1:C:225:GLY:HA2	2.38	0.59
1:C:123:GLU:HB3	1:C:168:MET:HE3	1.85	0.59
1:A:102:VAL:HG22	1:A:232:ILE:HB	1.86	0.58
1:G:126:ASN:OD1	1:G:128:THR:HG23	2.02	0.58
14:E:615:HOH:O	13:F:203:PEG:H41	2.06	0.55
13:E:402:PEG:H42	13:F:203:PEG:H11	1.88	0.55
2:B:39:LYS:HE3	14:B:482:HOH:O	2.08	0.54
1:E:102:VAL:HG22	1:E:232:ILE:HB	1.89	0.54
1:G:77:ASP:OD2	1:G:141:ARG:NH1	2.42	0.53
1:E:174:PHE:C	1:E:174:PHE:CD1	2.87	0.53
2:D:51:LYS:O	2:D:55:VAL:HG23	2.09	0.52
2:B:58:LYS:HE3	2:B:59:THR:C	2.35	0.51
1:C:77:ASP:OD2	1:C:141:ARG:NH1	2.43	0.51
2:H:51:LYS:O	2:H:55:VAL:HG23	2.11	0.50
13:E:402:PEG:H11	3:S:1:NAG:H83	1.94	0.50
1:E:195:TYR:O	1:E:196:VAL:HG22	2.11	0.49
2:D:1:GLY:HA2	14:D:310:HOH:O	2.12	0.49
2:H:54:ARG:HA	2:H:57:GLU:HG3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:LYS:HE2	1:A:83:LYS:HB2	1.70	0.48
1:E:320:MET:HG3	1:E:321:ARG:O	2.14	0.48
2:F:19:ASP:OD2	14:F:301:HOH:O	2.20	0.47
1:C:126:ASN:OD1	1:C:128:THR:HG23	2.15	0.47
2:B:58:LYS:NZ	2:B:60:ASN:HA	2.30	0.47
1:G:167:THR:OG1	1:G:242:ILE:HD11	2.15	0.47
1:C:132:GLN:NE2	1:C:152:ASN:HD21	2.14	0.46
11:B:202:PG4:H41	11:B:202:PG4:H22	1.76	0.46
7:Z:3:BMA:H62	7:Z:4:MAN:H2	1.70	0.46
1:C:92:LYS:NZ	14:C:411:HOH:O	2.48	0.46
1:A:234:TRP:NE1	14:A:502:HOH:O	2.23	0.46
2:B:115:MET:HE3	2:B:115:MET:HA	1.98	0.46
1:E:32:ASP:OD1	1:E:33:GLN:HG2	2.17	0.45
5:Q:3:BMA:H62	5:Q:5:MAN:H2	1.75	0.45
11:F:202:PG4:H21	11:F:202:PG4:H41	1.57	0.45
2:H:17:MET:HE1	2:H:23:GLY:HA3	1.99	0.45
1:A:143:SER:HB2	2:D:161:VAL:CG1	2.47	0.45
2:B:39:LYS:HE2	14:B:458:HOH:O	2.17	0.45
1:C:27:LYS:HG2	1:C:32:ASP:O	2.17	0.45
12:B:203:PG0:H31	12:B:203:PG0:H11	1.42	0.45
2:H:171:PHE:O	2:H:172:GLN:HB2	2.18	0.44
1:E:22:ASN:OD1	1:E:22:ASN:N	2.50	0.44
2:D:17:MET:HE1	2:D:23:GLY:HA3	2.00	0.43
1:E:234:TRP:NE1	14:E:504:HOH:O	2.37	0.43
1:C:237:VAL:HG21	1:C:243:LEU:HB2	1.99	0.43
1:G:322:ASN:O	1:G:324:PRO:HD3	2.19	0.43
1:C:324:PRO:O	1:C:325:GLU:HB2	2.19	0.42
2:D:171:PHE:O	2:D:172:GLN:HB2	2.19	0.42
2:H:82:LYS:NZ	14:H:313:HOH:O	2.53	0.42
2:B:171:PHE:O	2:B:172:GLN:HB2	2.19	0.42
1:C:207:LYS:HD2	1:C:241:ASP:HA	2.02	0.42
2:H:54:ARG:HG2	2:H:57:GLU:OE2	2.19	0.42
1:E:207:LYS:HG3	1:E:241:ASP:OD1	2.19	0.42
2:F:171:PHE:O	2:F:172:GLN:HB2	2.19	0.42
1:A:189:LYS:HB2	1:A:189:LYS:HE3	1.77	0.42
1:G:177:LEU:HB2	1:G:260:ILE:CD1	2.46	0.42
3:W:1:NAG:H3	6:Y:2:FUC:O4	2.20	0.42
1:C:166:VAL:HG12	1:C:245:ILE:HB	2.02	0.41
2:H:4:GLY:O	2:H:8:GLY:HA3	2.21	0.41
1:C:238:LYS:NZ	14:C:403:HOH:O	2.43	0.41
2:D:60:ASN:HB2	2:D:62:LYS:HE2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:29:ILE:HD13	1:E:29:ILE:HA	1.93	0.41
1:C:197:ARG:NH1	1:C:248:ASN:O	2.54	0.41
1:G:186:SER:HA	1:G:218:GLY:O	2.20	0.41
1:C:321:ARG:HD2	14:C:553:HOH:O	2.20	0.41
2:D:115:MET:HE3	2:D:115:MET:HA	2.02	0.41
6:Y:2:FUC:H5	6:Y:3:NAG:O5	2.21	0.41
1:G:259:LYS:HE3	1:G:259:LYS:HB3	1.84	0.41
2:H:54:ARG:HB2	2:H:54:ARG:NH1	2.36	0.41
2:F:115:MET:HE3	2:F:115:MET:HA	2.03	0.40
1:A:84:TRP:CE2	1:A:116:GLY:HA2	2.56	0.40
1:A:316:LEU:HD23	2:B:52:LEU:HD13	2.03	0.40
11:B:202:PG4:H52	11:B:202:PG4:H72	1.69	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	315/323 (98%)	304 (96%)	11 (4%)	0	100	100
1	C	315/323 (98%)	304 (96%)	11 (4%)	0	100	100
1	E	315/323 (98%)	305 (97%)	9 (3%)	1 (0%)	36	28
1	G	315/323 (98%)	304 (96%)	11 (4%)	0	100	100
2	B	170/174 (98%)	160 (94%)	10 (6%)	0	100	100
2	D	170/174 (98%)	161 (95%)	9 (5%)	0	100	100
2	F	170/174 (98%)	161 (95%)	9 (5%)	0	100	100
2	H	170/174 (98%)	160 (94%)	10 (6%)	0	100	100
All	All	1940/1988 (98%)	1859 (96%)	80 (4%)	1 (0%)	48	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	196	VAL

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	278/283 (98%)	277 (100%)	1 (0%)	84	85
1	C	278/283 (98%)	276 (99%)	2 (1%)	76	76
1	E	278/283 (98%)	276 (99%)	2 (1%)	76	76
1	G	278/283 (98%)	276 (99%)	2 (1%)	76	76
2	B	145/148 (98%)	145 (100%)	0	100	100
2	D	145/148 (98%)	144 (99%)	1 (1%)	76	76
2	F	146/148 (99%)	145 (99%)	1 (1%)	76	76
2	H	146/148 (99%)	146 (100%)	0	100	100
All	All	1694/1724 (98%)	1685 (100%)	9 (0%)	81	82

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

Mol	Chain	Res	Type
1	A	325	GLU
1	C	24	THR
1	C	137	TYR
2	D	59	THR
1	E	193	ASN
1	E	321	ARG
2	F	59	THR
1	G	321	ARG
1	G	325	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	132	GLN

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Mol	Chain	Res	Type
1	A	193	ASN
1	A	296	ASN
1	C	31	ASN
1	C	132	GLN
1	C	296	ASN
2	D	105	GLN
1	E	31	ASN
1	E	75	HIS
1	E	193	ASN
1	G	31	ASN
1	G	296	ASN
2	H	27	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

66 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	I	1	1,3	14,14,15	0.18	0	17,19,21	0.73	0
3	NAG	I	2	3	14,14,15	0.51	0	17,19,21	0.46	0
4	NAG	J	1	1,4	14,14,15	0.89	1 (7%)	17,19,21	0.47	0
4	NAG	J	2	4	14,14,15	0.53	0	17,19,21	0.61	0
4	BMA	J	3	4	11,11,12	0.86	0	15,15,17	0.59	0
5	NAG	K	1	1,5	14,14,15	0.30	0	17,19,21	0.60	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	K	2	5	14,14,15	0.16	0	17,19,21	0.48	0
5	BMA	K	3	5	11,11,12	0.81	0	15,15,17	1.00	0
5	MAN	K	4	5	11,11,12	0.81	0	15,15,17	0.94	1 (6%)
5	MAN	K	5	5	11,11,12	0.90	0	15,15,17	1.04	2 (13%)
3	NAG	L	1	1,3	14,14,15	0.46	0	17,19,21	0.56	0
3	NAG	L	2	3	14,14,15	1.40	2 (14%)	17,19,21	1.08	1 (5%)
4	NAG	M	1	1,4	14,14,15	0.24	0	17,19,21	0.79	1 (5%)
4	NAG	M	2	4	14,14,15	0.51	0	17,19,21	0.59	0
4	BMA	M	3	4	11,11,12	0.96	0	15,15,17	0.72	0
4	NAG	N	1	1,4	14,14,15	0.61	0	17,19,21	0.46	0
4	NAG	N	2	4	14,14,15	0.43	0	17,19,21	0.51	0
4	BMA	N	3	4	11,11,12	0.91	0	15,15,17	0.76	0
3	NAG	O	1	1,3	14,14,15	0.30	0	17,19,21	0.58	0
3	NAG	O	2	3	14,14,15	0.20	0	17,19,21	0.84	0
6	NAG	P	1	1,6	14,14,15	0.42	0	17,19,21	0.74	1 (5%)
6	FUC	P	2	6	10,10,11	0.84	0	14,14,16	0.82	0
6	NAG	P	3	6	14,14,15	0.34	0	17,19,21	0.52	0
6	FUC	P	4	6	10,10,11	0.87	0	14,14,16	0.68	0
5	NAG	Q	1	1,5	14,14,15	0.28	0	17,19,21	0.48	0
5	NAG	Q	2	5	14,14,15	0.20	0	17,19,21	0.47	0
5	BMA	Q	3	5	11,11,12	0.51	0	15,15,17	0.90	1 (6%)
5	MAN	Q	4	5	11,11,12	0.83	0	15,15,17	0.93	2 (13%)
5	MAN	Q	5	5	11,11,12	0.97	0	15,15,17	0.83	1 (6%)
5	NAG	R	1	1,5	14,14,15	0.27	0	17,19,21	0.61	0
5	NAG	R	2	5	14,14,15	0.25	0	17,19,21	0.48	0
5	BMA	R	3	5	11,11,12	0.79	0	15,15,17	0.75	0
5	MAN	R	4	5	11,11,12	1.40	2 (18%)	15,15,17	1.10	1 (6%)
5	MAN	R	5	5	11,11,12	0.81	0	15,15,17	1.03	2 (13%)
3	NAG	S	1	1,3	14,14,15	0.22	0	17,19,21	0.71	0
3	NAG	S	2	3	14,14,15	0.96	1 (7%)	17,19,21	1.32	1 (5%)
3	NAG	T	1	1,3	14,14,15	0.72	1 (7%)	17,19,21	0.46	0
3	NAG	T	2	3	14,14,15	0.47	0	17,19,21	0.40	0
3	NAG	U	1	1,3	14,14,15	0.42	0	17,19,21	0.46	0
3	NAG	U	2	3	14,14,15	0.55	0	17,19,21	0.51	0
3	NAG	V	1	1,3	14,14,15	0.28	0	17,19,21	0.76	1 (5%)
3	NAG	V	2	3	14,14,15	0.33	0	17,19,21	0.59	0
3	NAG	W	1	1,3	14,14,15	0.20	0	17,19,21	0.73	1 (5%)
3	NAG	W	2	3	14,14,15	0.46	0	17,19,21	0.49	0
3	NAG	X	1	1,3	14,14,15	0.64	1 (7%)	17,19,21	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	X	2	3	14,14,15	0.71	1 (7%)	17,19,21	0.69	0
6	NAG	Y	1	1,6	14,14,15	0.38	0	17,19,21	0.66	1 (5%)
6	FUC	Y	2	6	10,10,11	0.92	0	14,14,16	0.84	0
6	NAG	Y	3	6	14,14,15	0.49	0	17,19,21	0.53	0
6	FUC	Y	4	6	10,10,11	1.00	1 (10%)	14,14,16	0.97	0
7	NAG	Z	1	1,7	14,14,15	0.37	0	17,19,21	0.44	0
7	NAG	Z	2	7	14,14,15	0.16	0	17,19,21	0.57	0
7	BMA	Z	3	7	11,11,12	0.58	0	15,15,17	0.79	0
7	MAN	Z	4	7	11,11,12	1.23	1 (9%)	15,15,17	1.11	2 (13%)
7	MAN	Z	5	7	11,11,12	0.69	0	15,15,17	1.03	2 (13%)
7	MAN	Z	6	7	11,11,12	0.82	0	15,15,17	0.99	2 (13%)
8	NAG	a	1	8	15,15,15	1.50	3 (20%)	21,21,21	1.48	3 (14%)
8	GAL	a	2	8	11,11,12	0.79	0	15,15,17	1.00	0
8	SIA	a	3	8	20,20,21	1.71	2 (10%)	21,28,31	1.40	2 (9%)
9	GAL	b	1	9	12,12,12	0.54	0	17,17,17	0.81	0
9	SIA	b	2	9	20,20,21	1.76	3 (15%)	21,28,31	1.51	4 (19%)
8	NAG	c	1	8	15,15,15	1.36	1 (6%)	21,21,21	1.25	3 (14%)
8	GAL	c	2	8	11,11,12	0.81	0	15,15,17	1.09	1 (6%)
8	SIA	c	3	8	20,20,21	1.64	2 (10%)	21,28,31	1.23	2 (9%)
9	GAL	d	1	9	12,12,12	0.65	0	17,17,17	0.82	0
9	SIA	d	2	9	20,20,21	1.66	3 (15%)	21,28,31	1.42	4 (19%)

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

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	L	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	L	2	3	-	2/6/23/26	0/1/1/1
4	NAG	M	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	M	2	4	-	2/6/23/26	0/1/1/1
4	BMA	M	3	4	-	0/2/19/22	0/1/1/1
4	NAG	N	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	N	2	4	-	0/6/23/26	0/1/1/1
4	BMA	N	3	4	-	0/2/19/22	0/1/1/1
3	NAG	O	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	O	2	3	-	4/6/23/26	0/1/1/1
6	NAG	P	1	1,6	-	0/6/23/26	0/1/1/1
6	FUC	P	2	6	-	-	0/1/1/1
6	NAG	P	3	6	-	0/6/23/26	0/1/1/1
6	FUC	P	4	6	-	-	0/1/1/1
5	NAG	Q	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	Q	2	5	-	0/6/23/26	0/1/1/1
5	BMA	Q	3	5	-	2/2/19/22	0/1/1/1
5	MAN	Q	4	5	-	1/2/19/22	0/1/1/1
5	MAN	Q	5	5	-	0/2/19/22	0/1/1/1
5	NAG	R	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	R	2	5	-	0/6/23/26	0/1/1/1
5	BMA	R	3	5	-	2/2/19/22	0/1/1/1
5	MAN	R	4	5	-	0/2/19/22	0/1/1/1
5	MAN	R	5	5	-	2/2/19/22	0/1/1/1
3	NAG	S	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	S	2	3	-	2/6/23/26	0/1/1/1
3	NAG	T	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	T	2	3	-	2/6/23/26	0/1/1/1
3	NAG	U	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	U	2	3	-	2/6/23/26	0/1/1/1
3	NAG	V	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	V	2	3	-	2/6/23/26	0/1/1/1
3	NAG	W	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	W	2	3	-	0/6/23/26	0/1/1/1
3	NAG	X	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	X	2	3	-	2/6/23/26	0/1/1/1
6	NAG	Y	1	1,6	-	1/6/23/26	0/1/1/1
6	FUC	Y	2	6	-	-	0/1/1/1
6	NAG	Y	3	6	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	FUC	Y	4	6	-	-	0/1/1/1
7	NAG	Z	1	1,7	-	0/6/23/26	0/1/1/1
7	NAG	Z	2	7	-	0/6/23/26	0/1/1/1
7	BMA	Z	3	7	-	2/2/19/22	0/1/1/1
7	MAN	Z	4	7	-	0/2/19/22	0/1/1/1
7	MAN	Z	5	7	-	0/2/19/22	0/1/1/1
7	MAN	Z	6	7	-	0/2/19/22	0/1/1/1
8	NAG	a	1	8	-	2/6/26/26	0/1/1/1
8	GAL	a	2	8	-	0/2/19/22	0/1/1/1
8	SIA	a	3	8	-	0/18/34/38	0/1/1/1
9	GAL	b	1	9	-	0/2/22/22	0/1/1/1
9	SIA	b	2	9	-	1/18/34/38	0/1/1/1
8	NAG	c	1	8	-	1/6/26/26	0/1/1/1
8	GAL	c	2	8	-	1/2/19/22	0/1/1/1
8	SIA	c	3	8	-	0/18/34/38	0/1/1/1
9	GAL	d	1	9	-	0/2/22/22	0/1/1/1
9	SIA	d	2	9	-	0/18/34/38	0/1/1/1

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	b	2	SIA	C2-C1	6.40	1.60	1.52
9	d	2	SIA	C2-C1	5.95	1.59	1.52
8	a	3	SIA	C2-C1	5.85	1.59	1.52
8	c	3	SIA	C2-C1	5.71	1.59	1.52
8	c	1	NAG	C1-C2	4.83	1.58	1.52
8	a	1	NAG	C1-C2	4.64	1.58	1.52
3	L	2	NAG	O5-C1	-3.47	1.37	1.43
3	L	2	NAG	C1-C2	3.37	1.56	1.52
3	S	2	NAG	O5-C1	3.33	1.49	1.43
5	R	4	MAN	O5-C1	-2.99	1.38	1.43
4	J	1	NAG	O5-C1	-2.93	1.38	1.43
7	Z	4	MAN	C2-C3	2.84	1.56	1.52
8	a	3	SIA	C7-C6	2.61	1.56	1.52
8	c	3	SIA	C7-C6	2.57	1.56	1.52
3	X	2	NAG	C1-C2	2.43	1.55	1.52
8	a	1	NAG	O5-C1	-2.36	1.37	1.42
9	d	2	SIA	C7-C6	2.28	1.55	1.52
8	a	1	NAG	C3-C2	2.21	1.57	1.53
5	R	4	MAN	C4-C5	2.20	1.57	1.53
9	b	2	SIA	C7-C6	2.20	1.55	1.52
3	X	1	NAG	C1-C2	2.10	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	T	1	NAG	O5-C1	-2.08	1.40	1.43
9	d	2	SIA	O6-C2	2.07	1.47	1.43
6	Y	4	FUC	O5-C1	-2.03	1.40	1.43
9	b	2	SIA	O6-C2	2.02	1.47	1.43

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	S	2	NAG	C1-O5-C5	5.11	119.03	112.19
8	a	1	NAG	C4-C3-C2	4.46	116.90	110.40
8	a	3	SIA	O1A-C1-C2	-4.03	114.14	122.85
8	a	1	NAG	C1-C2-C3	3.75	115.66	110.54
9	b	2	SIA	O1A-C1-C2	-3.71	114.84	122.85
8	c	1	NAG	C1-C2-C3	3.48	115.29	110.54
3	L	2	NAG	C4-C3-C2	3.39	115.98	111.02
9	d	2	SIA	O1A-C1-C2	-3.36	115.60	122.85
8	c	3	SIA	O1A-C1-C2	-3.31	115.70	122.85
8	c	1	NAG	C4-C3-C2	3.21	115.07	110.40
7	Z	4	MAN	O3-C3-C2	2.79	115.75	110.05
8	c	3	SIA	O1B-C1-O1A	2.78	130.39	124.08
5	K	5	MAN	C1-O5-C5	2.78	115.91	112.19
8	a	3	SIA	O1B-C1-O1A	2.76	130.34	124.08
7	Z	5	MAN	C1-O5-C5	2.67	115.77	112.19
5	R	5	MAN	C1-O5-C5	2.65	115.73	112.19
8	a	1	NAG	O5-C5-C4	-2.65	104.93	109.70
3	V	1	NAG	C1-O5-C5	2.56	115.62	112.19
9	b	2	SIA	O6-C2-C3	-2.51	107.18	110.56
4	M	1	NAG	C1-O5-C5	2.44	115.46	112.19
9	d	2	SIA	O1B-C1-O1A	2.42	129.57	124.08
7	Z	6	MAN	C1-O5-C5	2.41	115.42	112.19
7	Z	4	MAN	C1-O5-C5	2.40	115.40	112.19
5	R	4	MAN	C3-C4-C5	2.35	114.50	110.23
7	Z	6	MAN	O2-C2-C3	-2.30	105.38	110.15
9	d	2	SIA	O6-C2-C3	-2.28	107.49	110.56
8	c	1	NAG	O5-C5-C4	-2.22	105.69	109.70
6	P	1	NAG	C1-O5-C5	2.22	115.17	112.19
8	c	2	GAL	C1-O5-C5	2.22	115.17	112.19
5	K	4	MAN	C1-O5-C5	2.20	115.14	112.19
5	Q	3	BMA	C1-O5-C5	2.19	115.12	112.19
7	Z	5	MAN	O2-C2-C3	-2.19	105.62	110.15
9	b	2	SIA	O1B-C1-O1A	2.18	129.03	124.08
5	Q	5	MAN	O2-C2-C3	-2.13	105.73	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	Q	4	MAN	O2-C2-C3	-2.10	105.80	110.15
9	d	2	SIA	O6-C2-C1	2.07	111.62	107.72
9	b	2	SIA	O6-C2-C1	2.06	111.60	107.72
3	W	1	NAG	C1-O5-C5	2.05	114.93	112.19
5	R	5	MAN	O2-C2-C3	-2.04	105.92	110.15
5	Q	4	MAN	C1-O5-C5	2.04	114.92	112.19
6	Y	1	NAG	C1-O5-C5	2.02	114.90	112.19
5	K	5	MAN	O2-C2-C3	-2.00	106.00	110.15

There are no chirality outliers.

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	I	1	NAG	C1-C2-N2-C7
3	S	1	NAG	C1-C2-N2-C7
3	X	2	NAG	O5-C5-C6-O6
3	O	2	NAG	C4-C5-C6-O6
7	Z	3	BMA	C4-C5-C6-O6
5	Q	3	BMA	C4-C5-C6-O6
3	U	1	NAG	O5-C5-C6-O6
3	S	2	NAG	O5-C5-C6-O6
3	T	2	NAG	O5-C5-C6-O6
3	L	2	NAG	O5-C5-C6-O6
5	R	3	BMA	O5-C5-C6-O6
7	Z	3	BMA	O5-C5-C6-O6
4	J	2	NAG	O5-C5-C6-O6
3	X	2	NAG	C4-C5-C6-O6
3	O	2	NAG	O5-C5-C6-O6
5	Q	3	BMA	O5-C5-C6-O6
5	R	3	BMA	C4-C5-C6-O6
3	T	2	NAG	C4-C5-C6-O6
3	U	1	NAG	C4-C5-C6-O6
4	J	2	NAG	C4-C5-C6-O6
3	L	2	NAG	C4-C5-C6-O6
3	I	1	NAG	O5-C5-C6-O6
5	K	4	MAN	C4-C5-C6-O6
3	I	1	NAG	C4-C5-C6-O6
5	R	5	MAN	O5-C5-C6-O6
5	K	2	NAG	C4-C5-C6-O6
8	a	1	NAG	C4-C5-C6-O6
5	K	4	MAN	O5-C5-C6-O6
5	K	3	BMA	C4-C5-C6-O6

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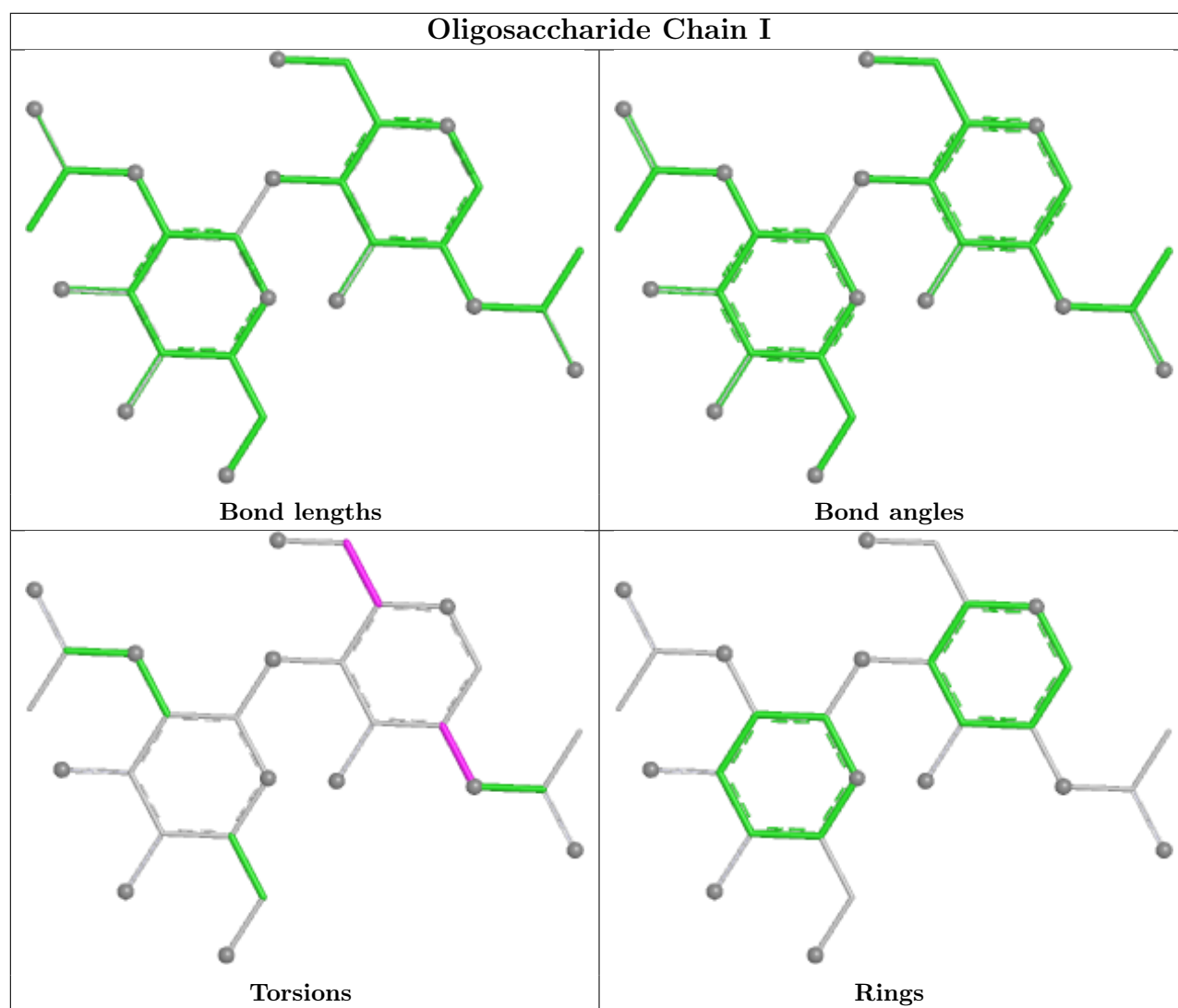
Mol	Chain	Res	Type	Atoms
5	K	2	NAG	O5-C5-C6-O6
3	S	2	NAG	C4-C5-C6-O6
8	c	2	GAL	O5-C5-C6-O6
5	Q	4	MAN	O5-C5-C6-O6
3	U	2	NAG	C4-C5-C6-O6
5	K	3	BMA	O5-C5-C6-O6
8	c	1	NAG	C4-C5-C6-O6
8	a	1	NAG	O5-C5-C6-O6
3	O	2	NAG	C3-C2-N2-C7
3	V	2	NAG	C3-C2-N2-C7
3	S	1	NAG	C4-C5-C6-O6
5	R	5	MAN	C4-C5-C6-O6
3	U	2	NAG	O5-C5-C6-O6
3	O	2	NAG	C1-C2-N2-C7
3	V	2	NAG	C1-C2-N2-C7
6	Y	1	NAG	C1-C2-N2-C7
6	Y	3	NAG	C1-C2-N2-C7
3	I	1	NAG	C3-C2-N2-C7
4	M	2	NAG	O5-C5-C6-O6
4	M	2	NAG	C4-C5-C6-O6
9	b	2	SIA	C6-C7-C8-O8
3	S	1	NAG	O5-C5-C6-O6

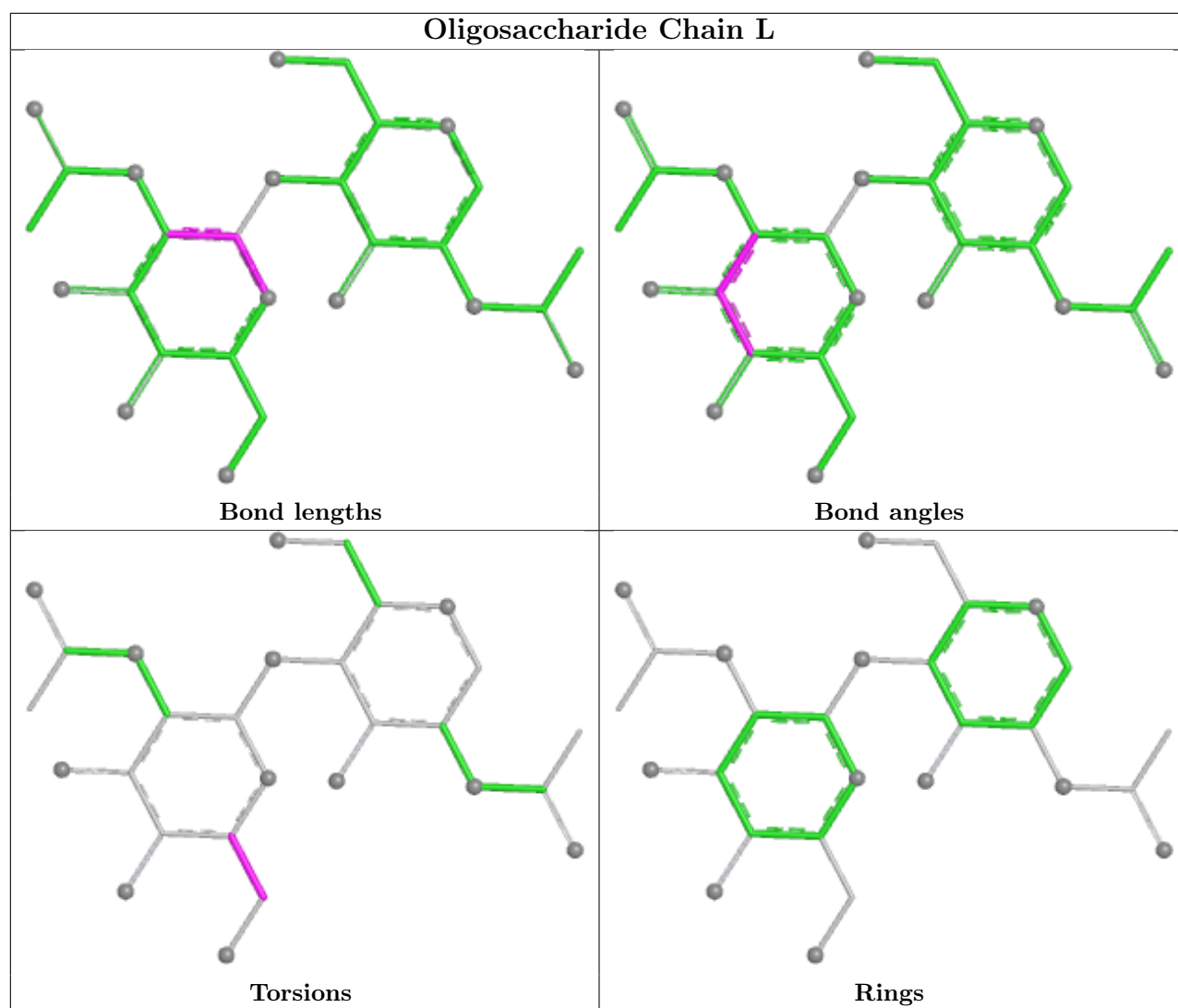
There are no ring outliers.

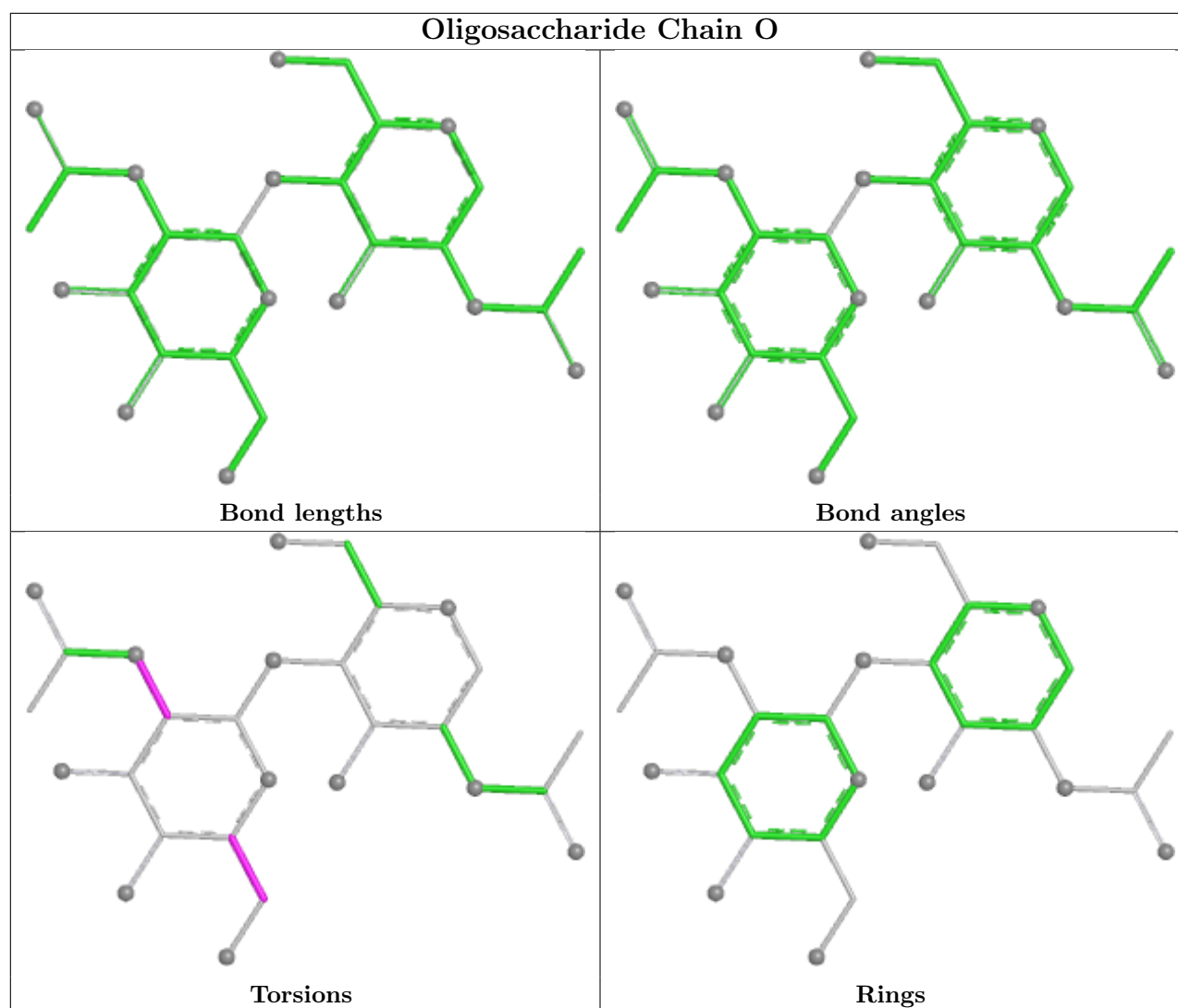
8 monomers are involved in 5 short contacts:

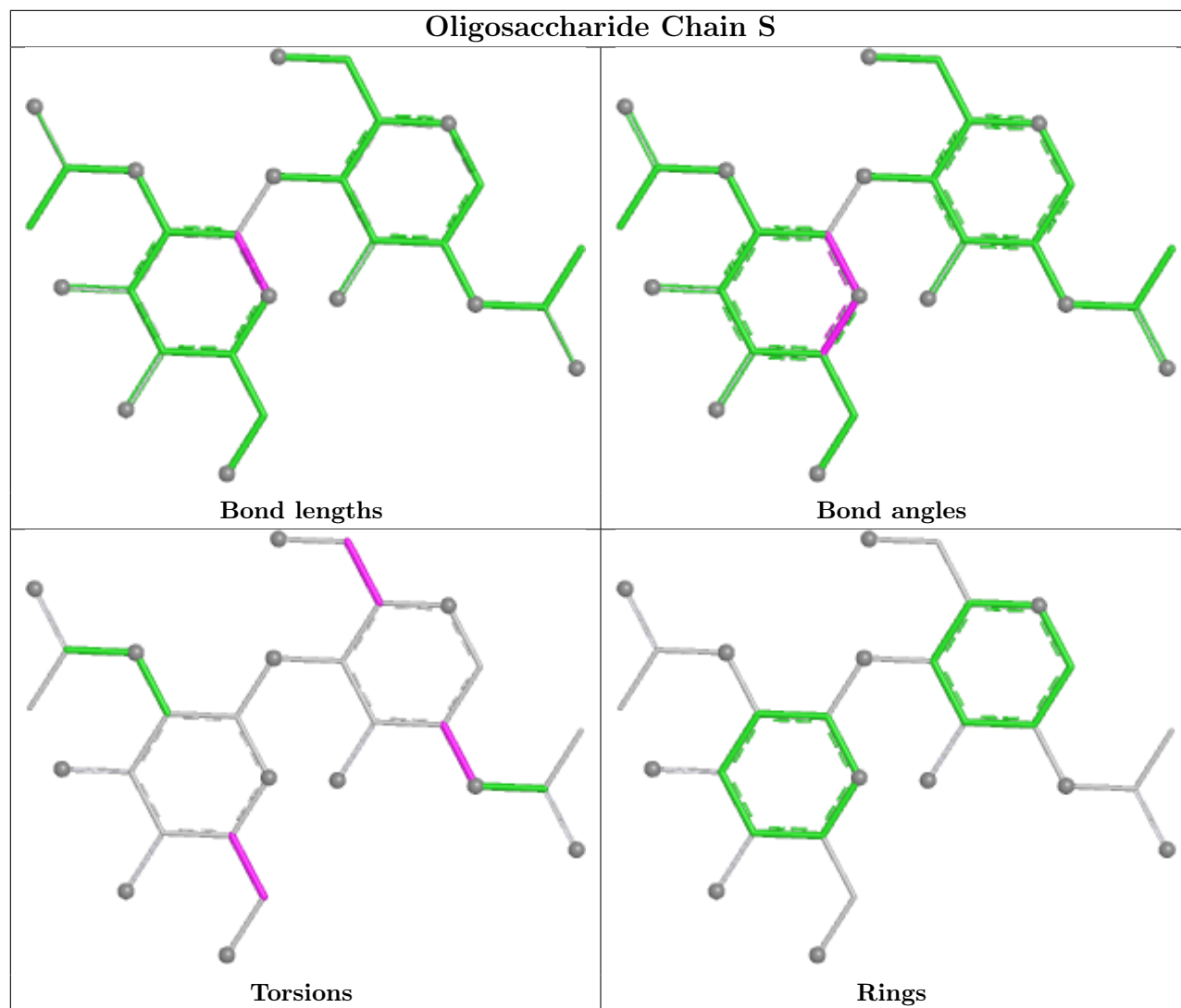
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	Q	3	BMA	1	0
6	Y	3	NAG	1	0
7	Z	4	MAN	1	0
3	W	1	NAG	1	0
6	Y	2	FUC	2	0
5	Q	5	MAN	1	0
7	Z	3	BMA	1	0
3	S	1	NAG	1	0

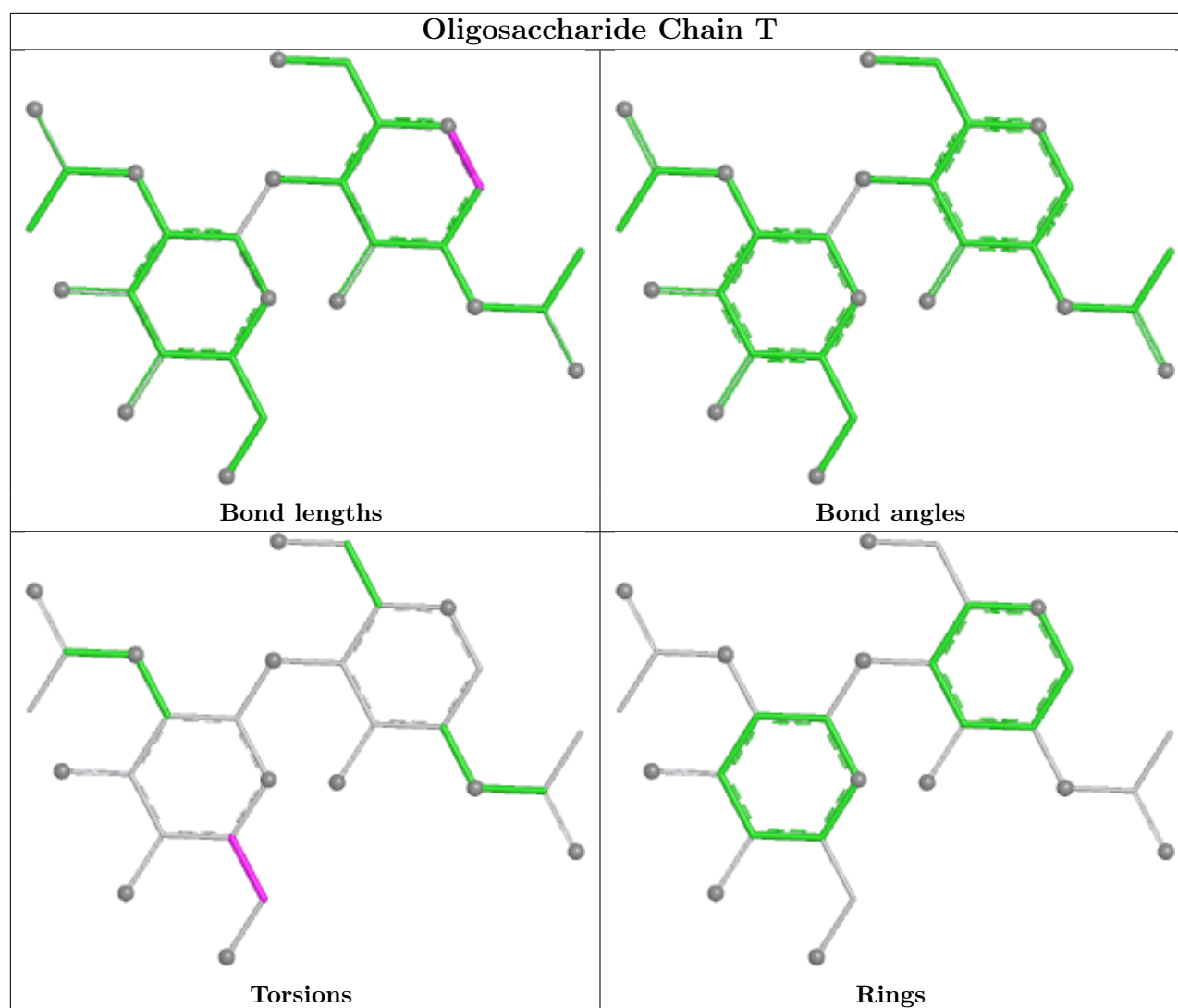
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

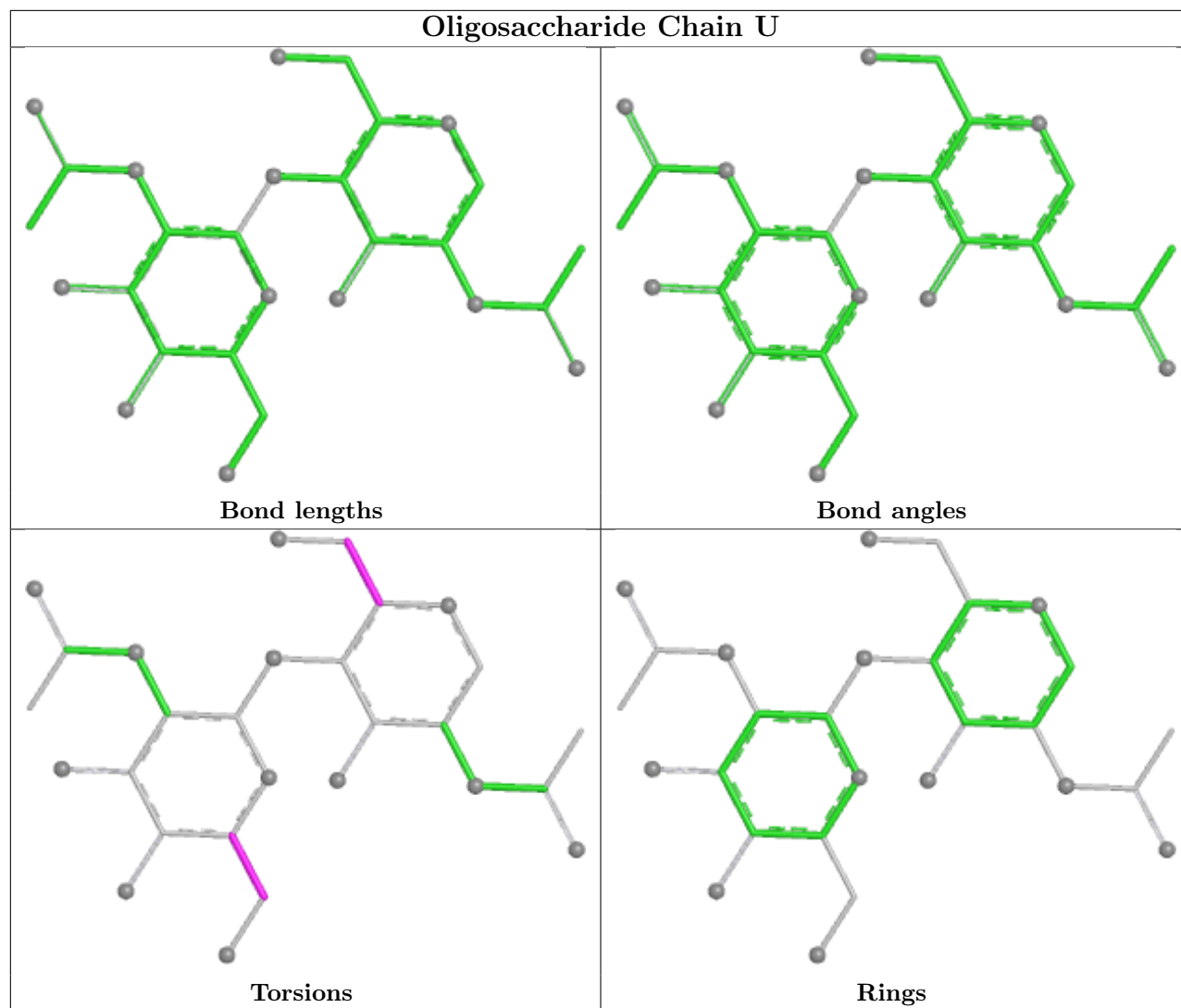


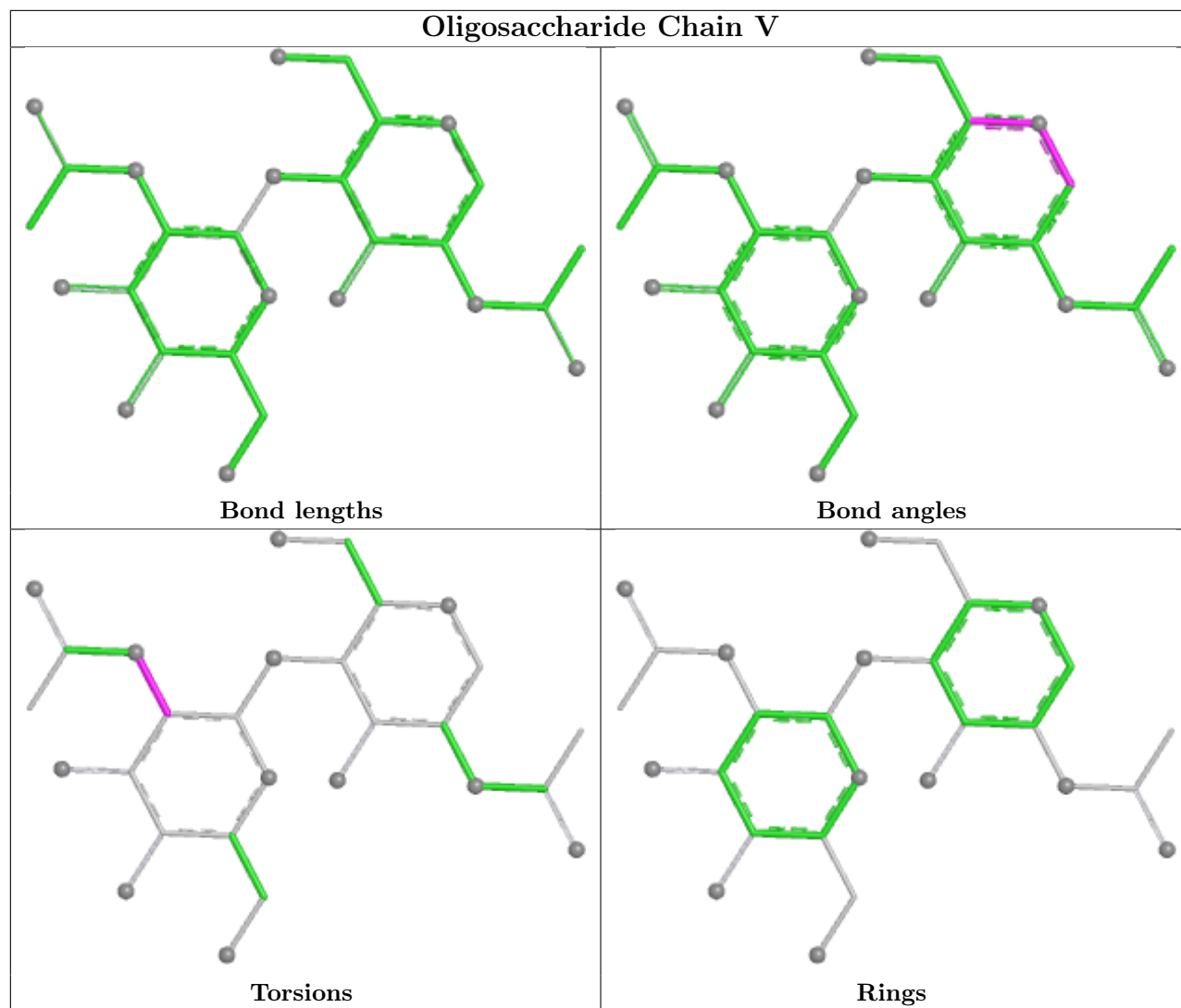


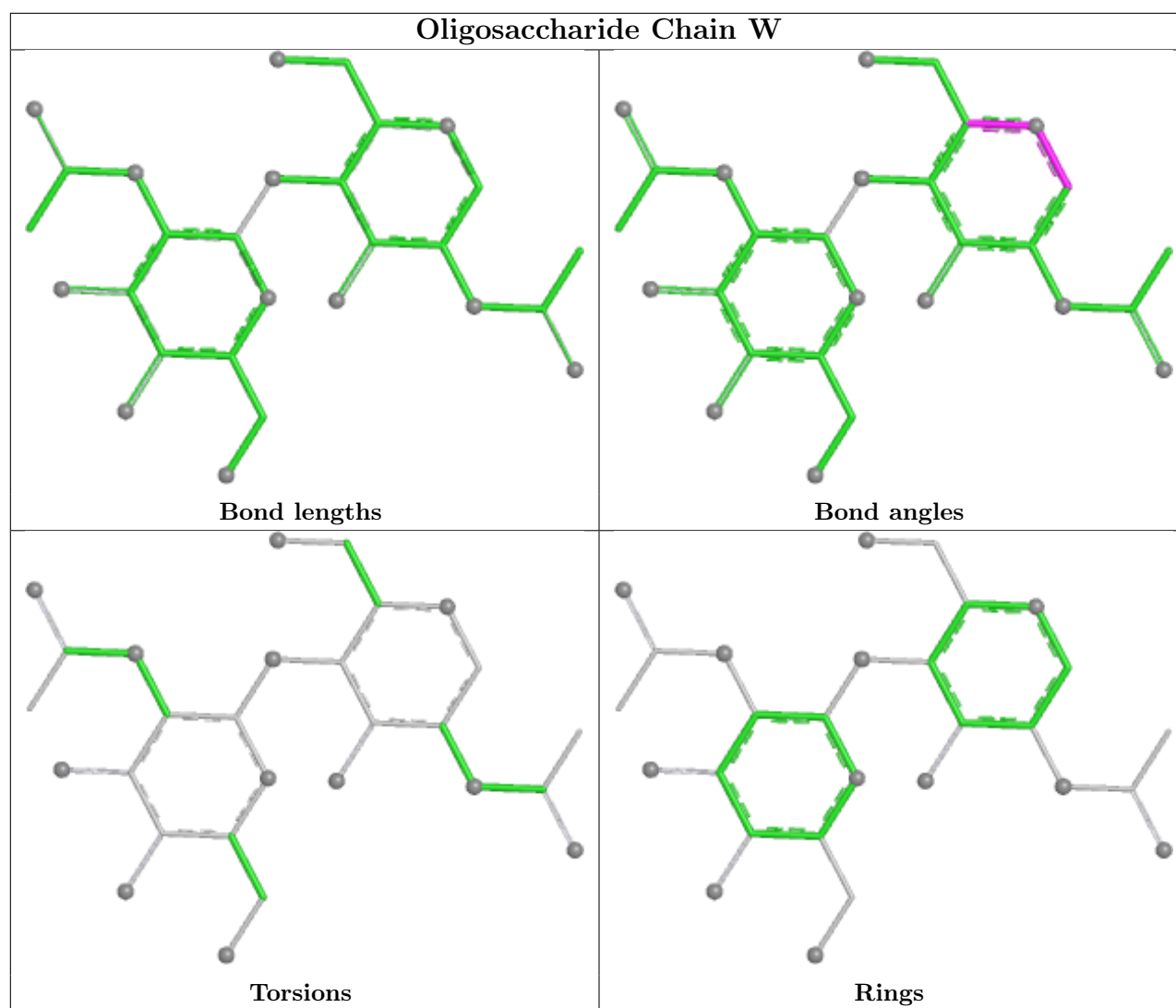


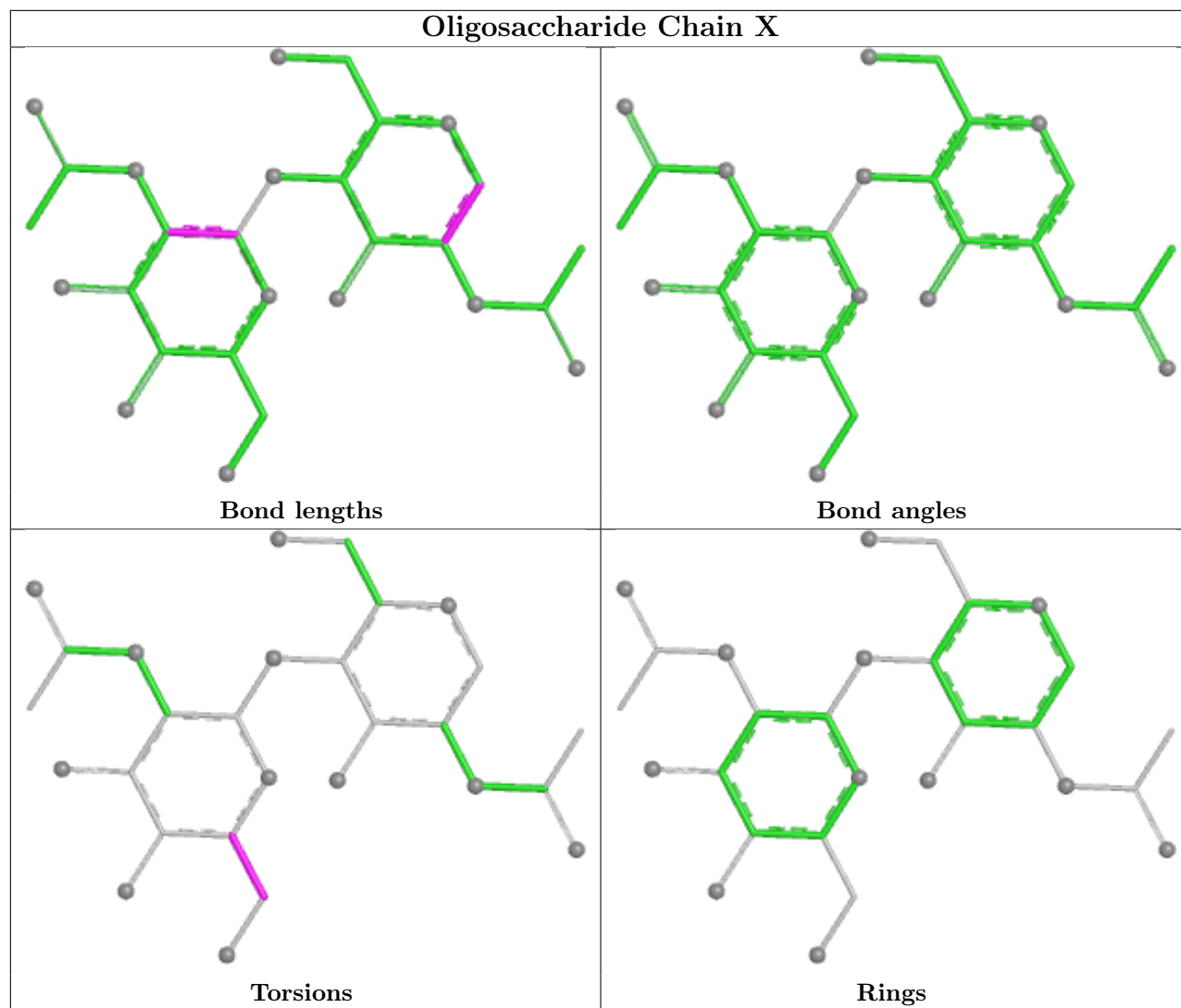


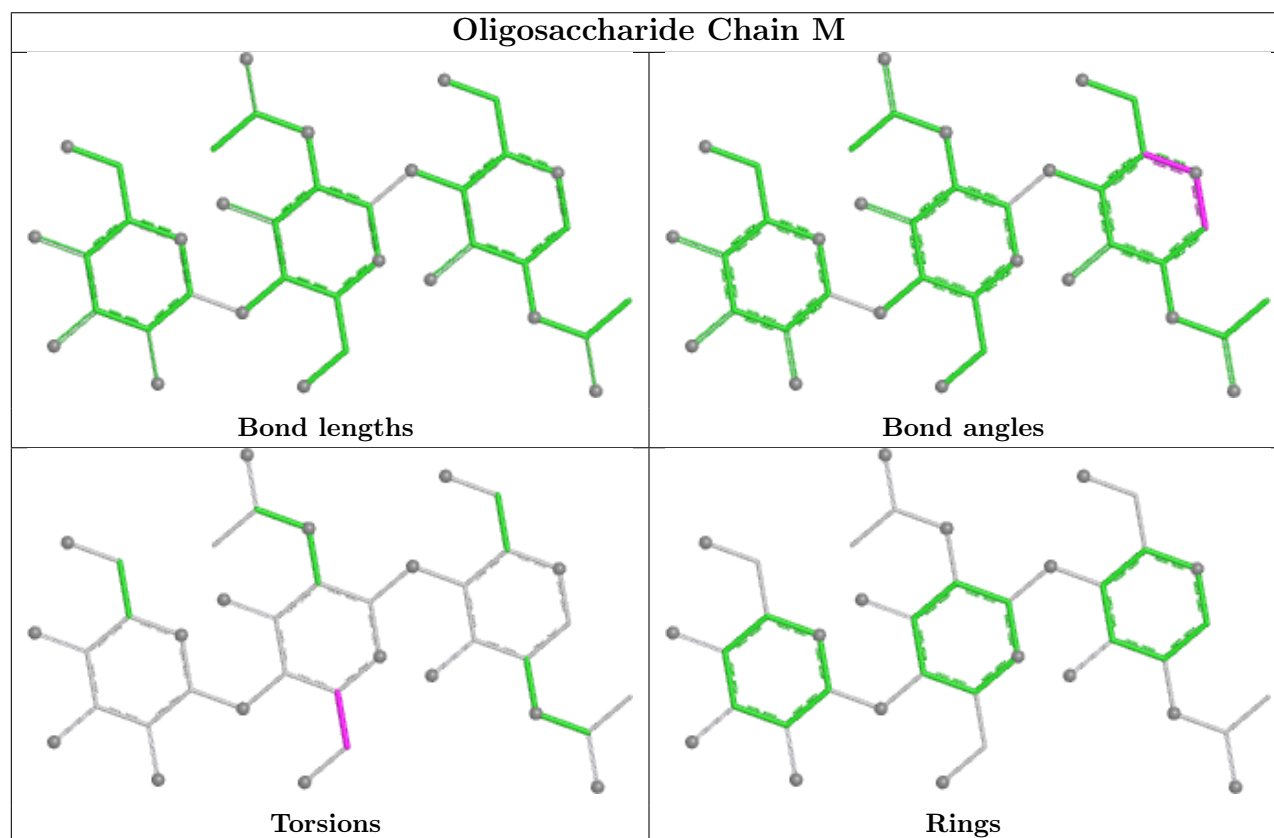
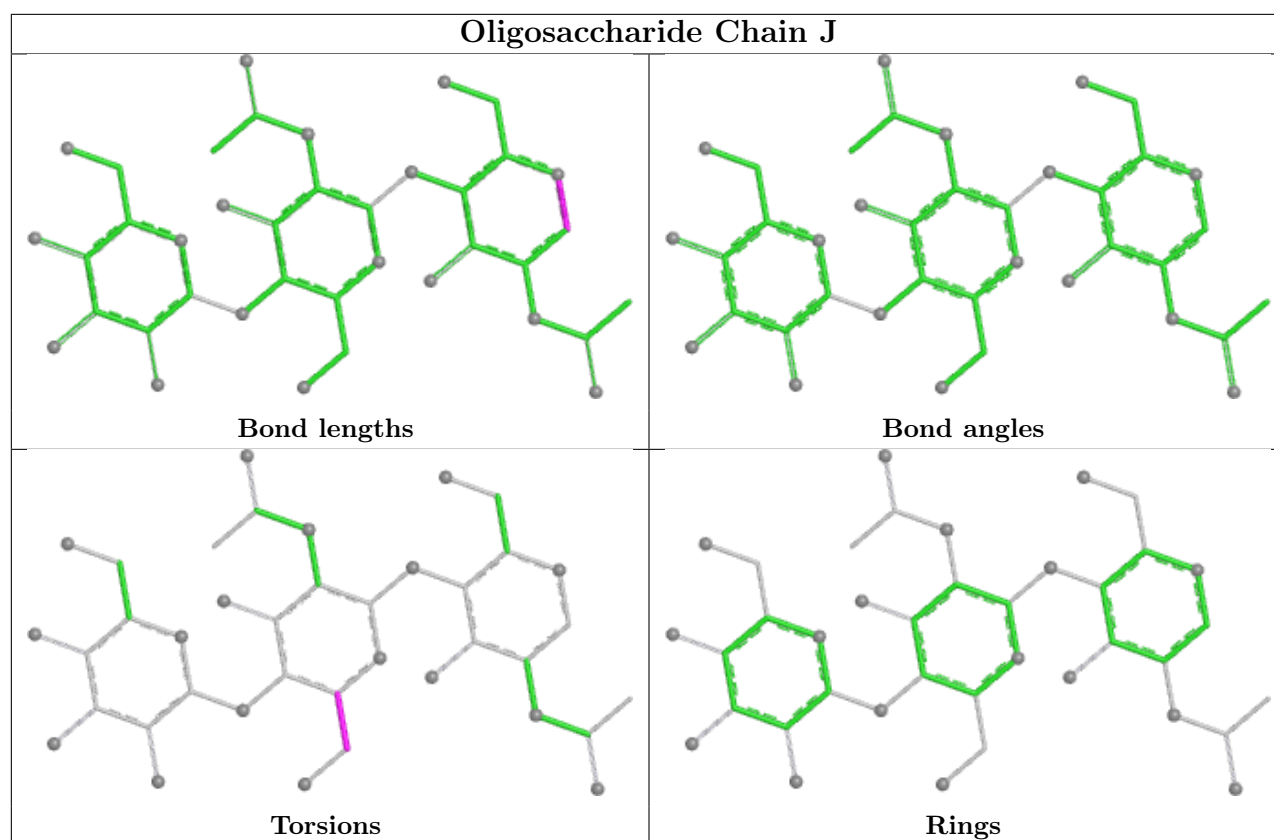


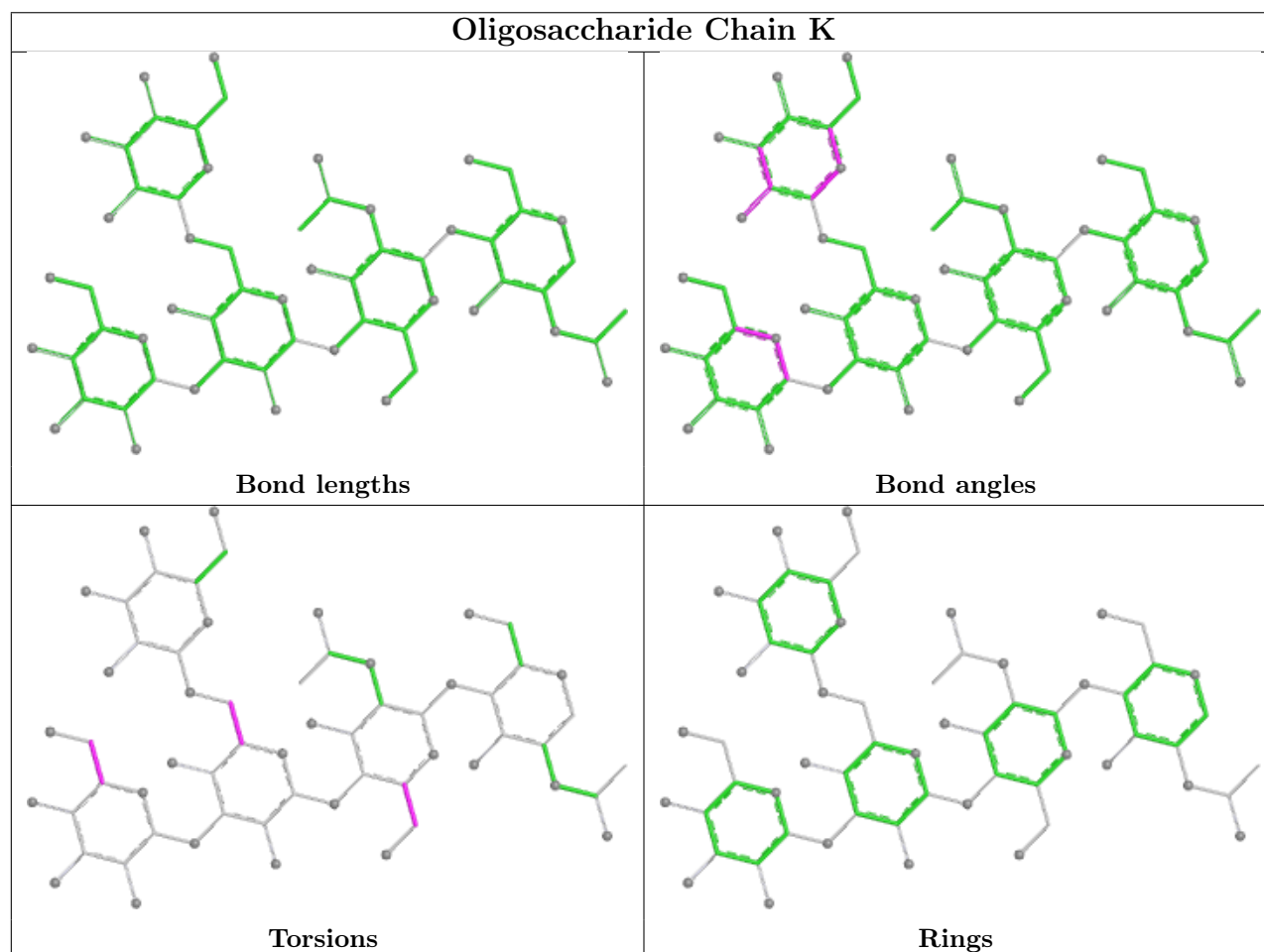
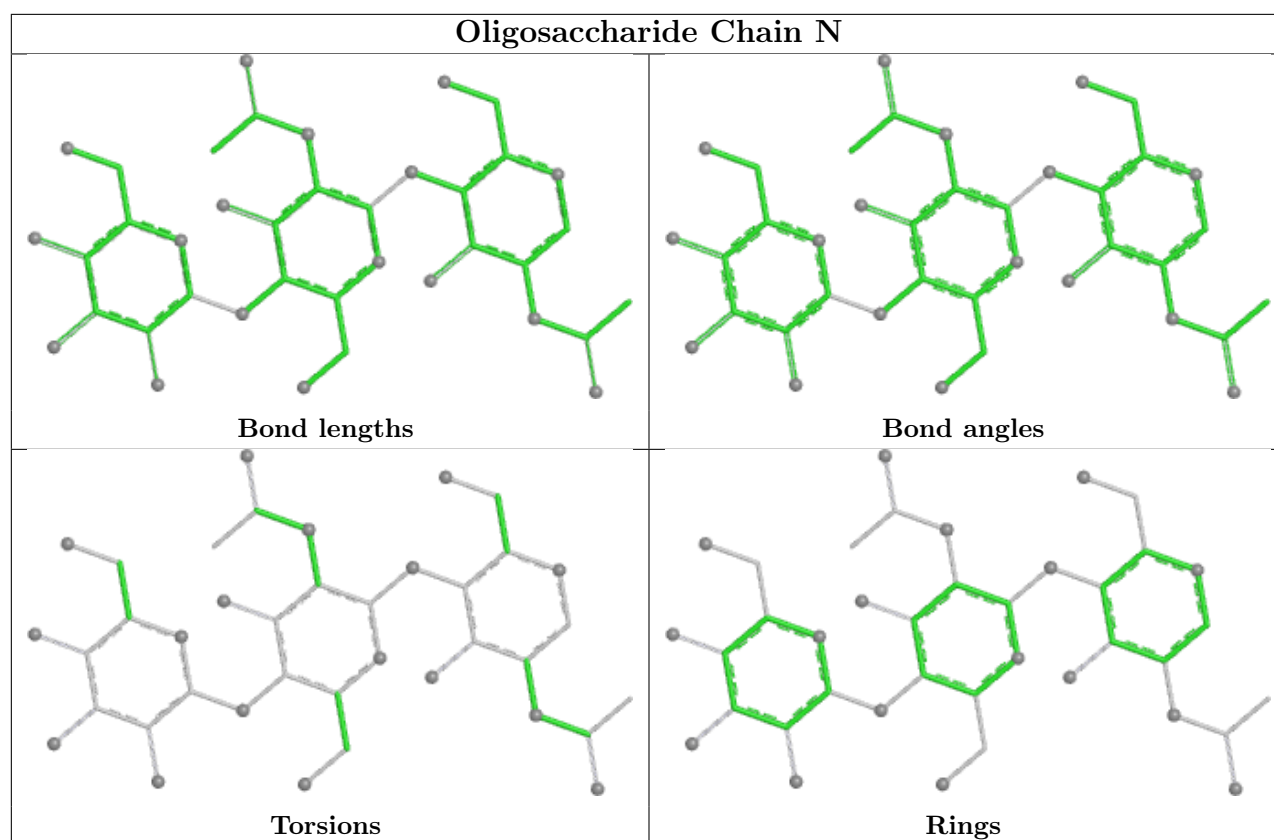


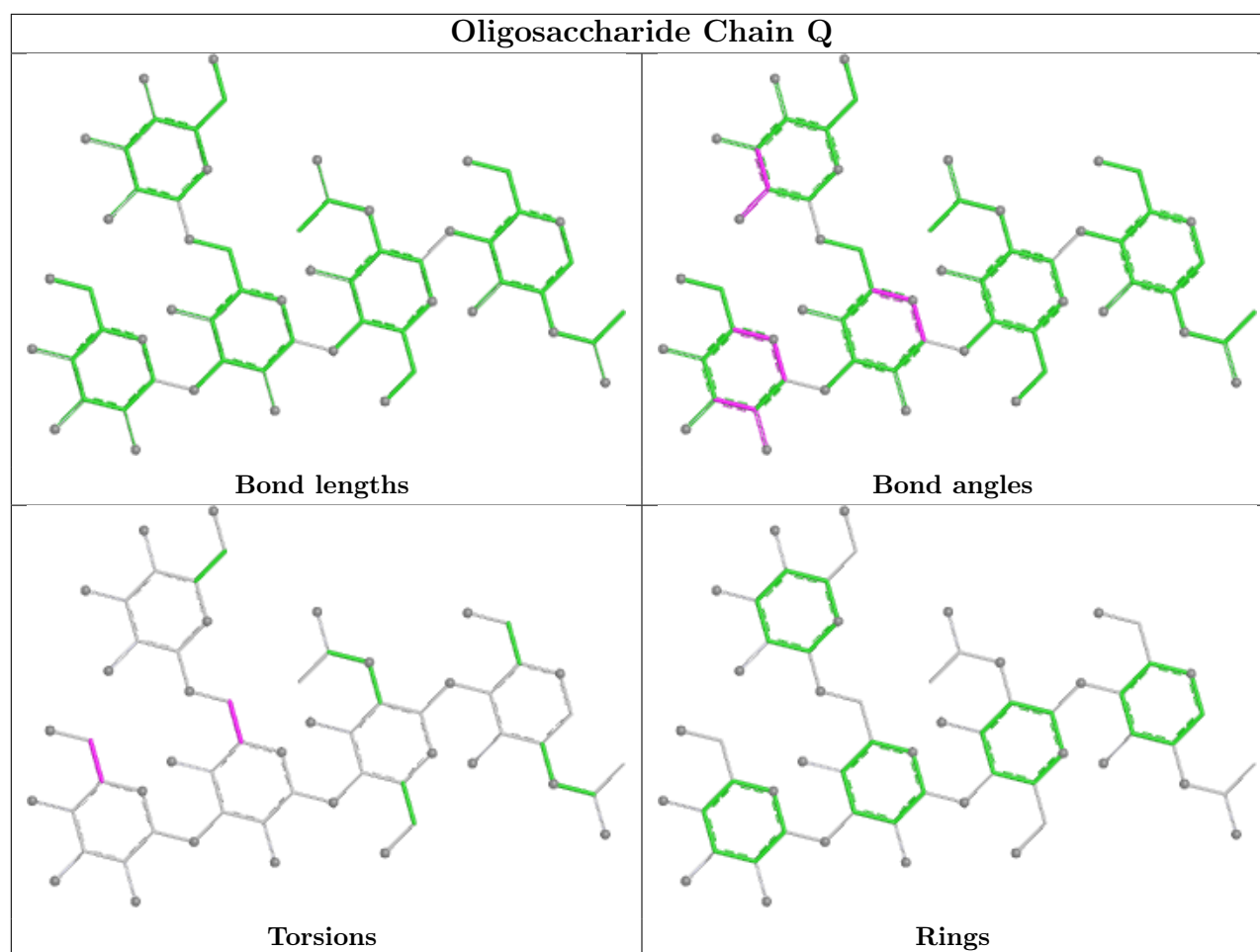


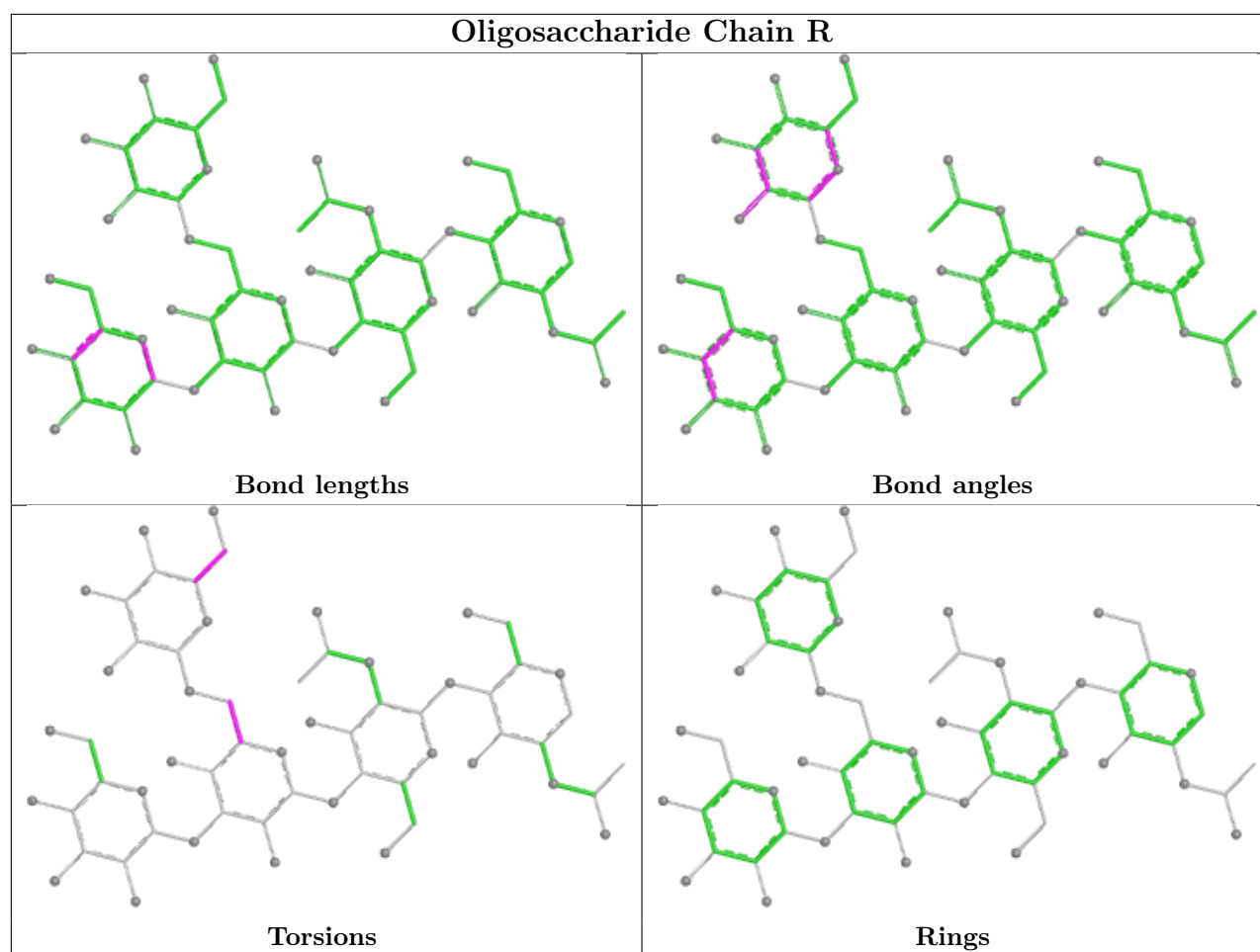


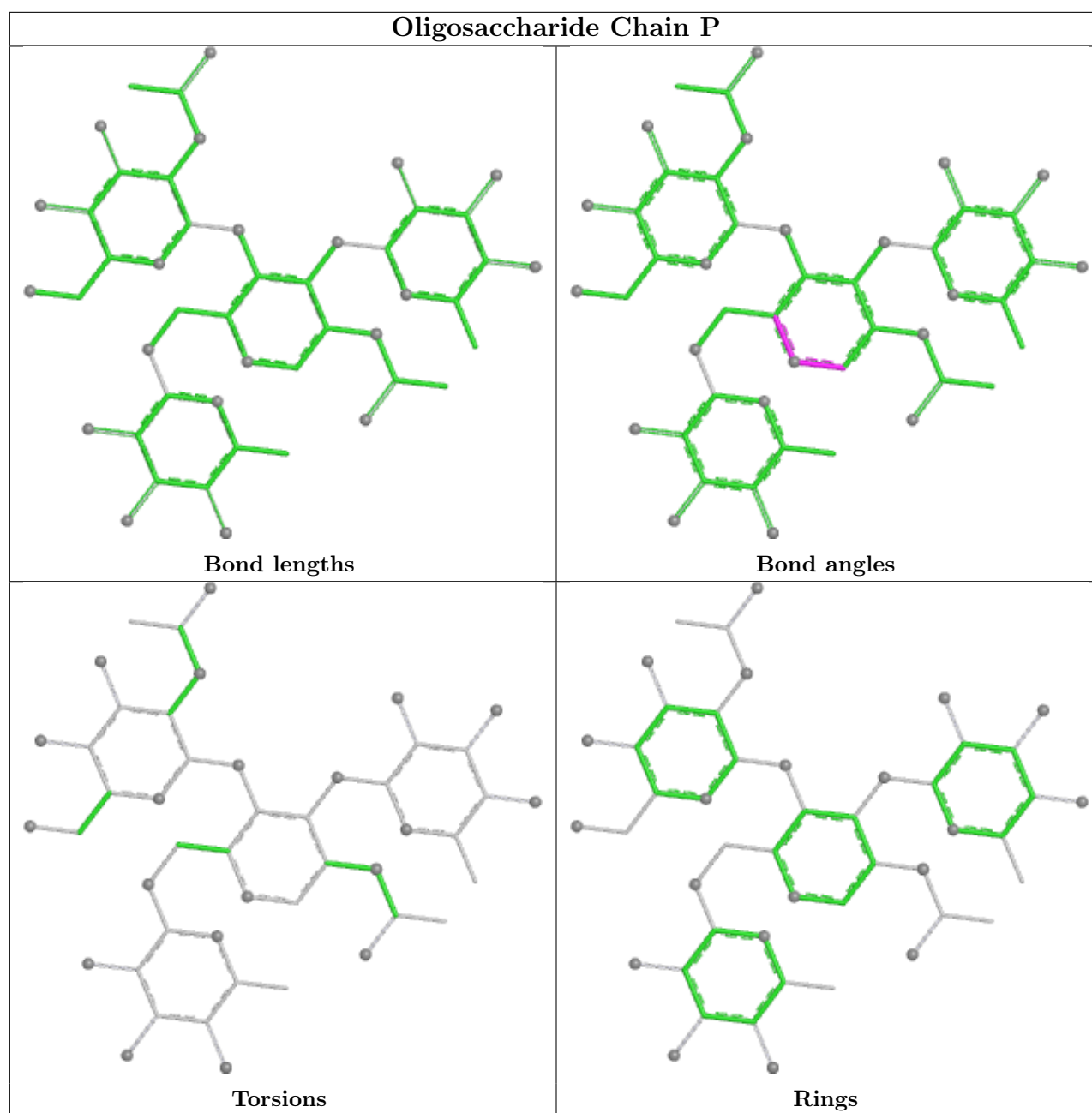


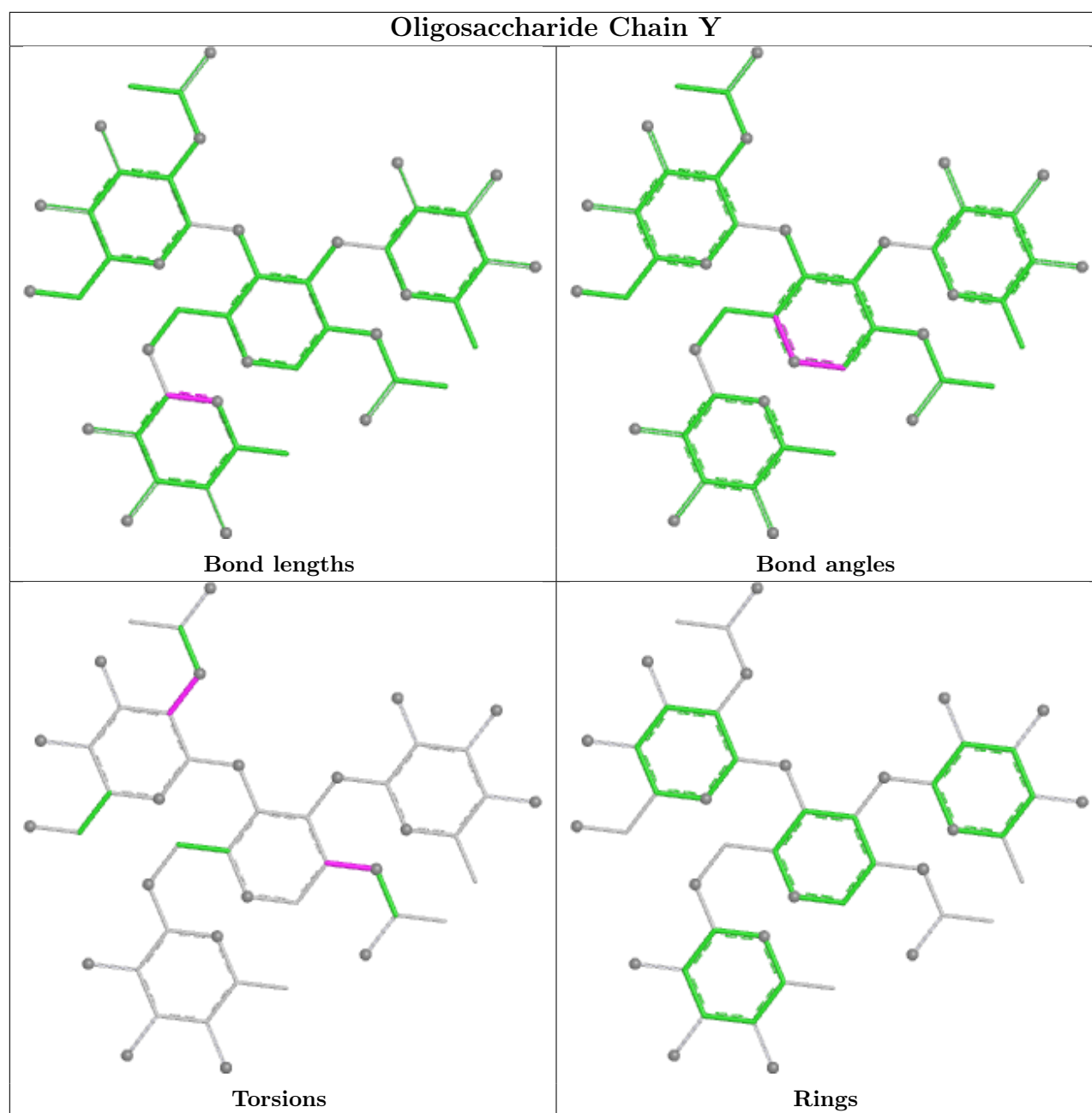


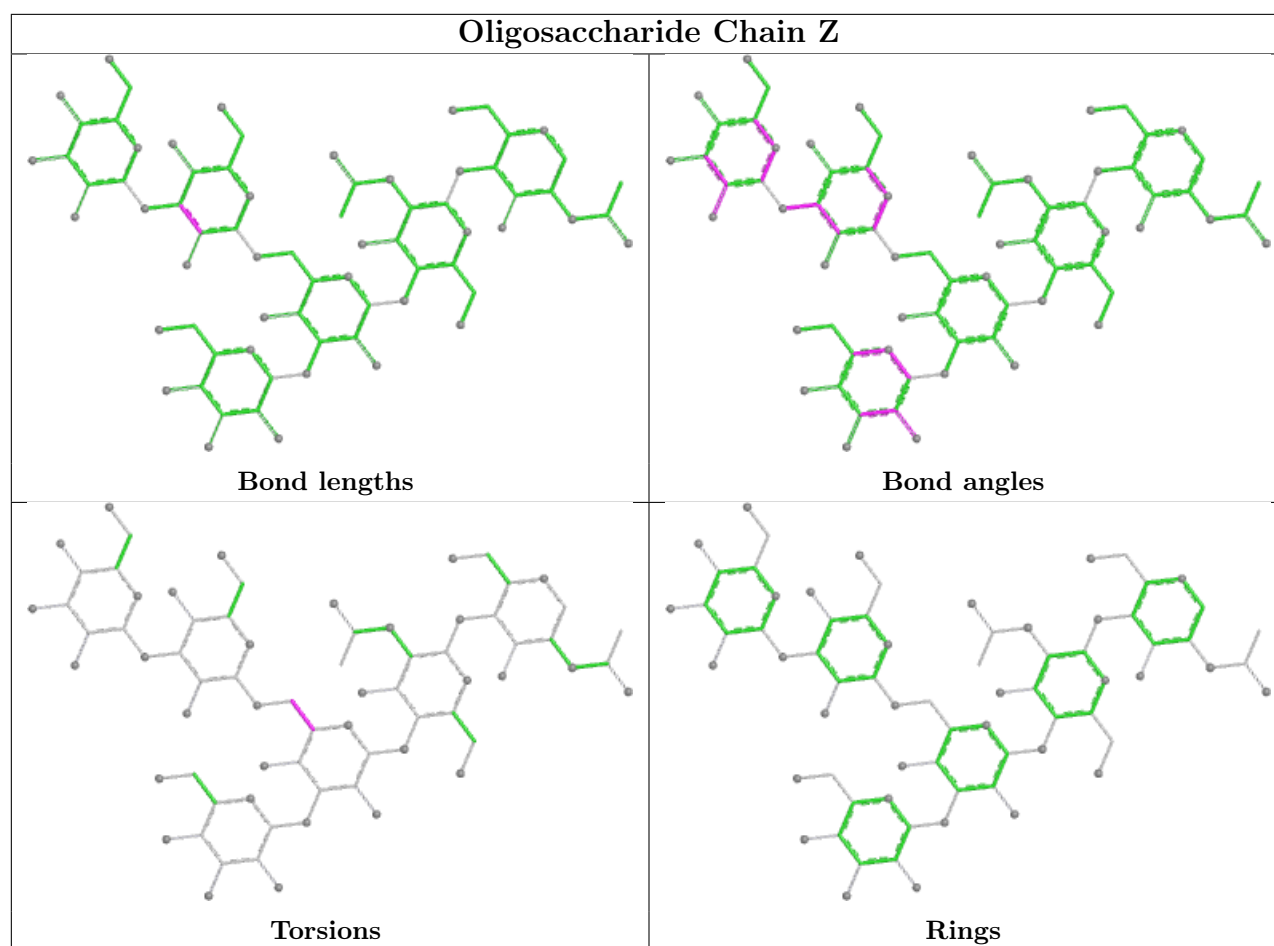


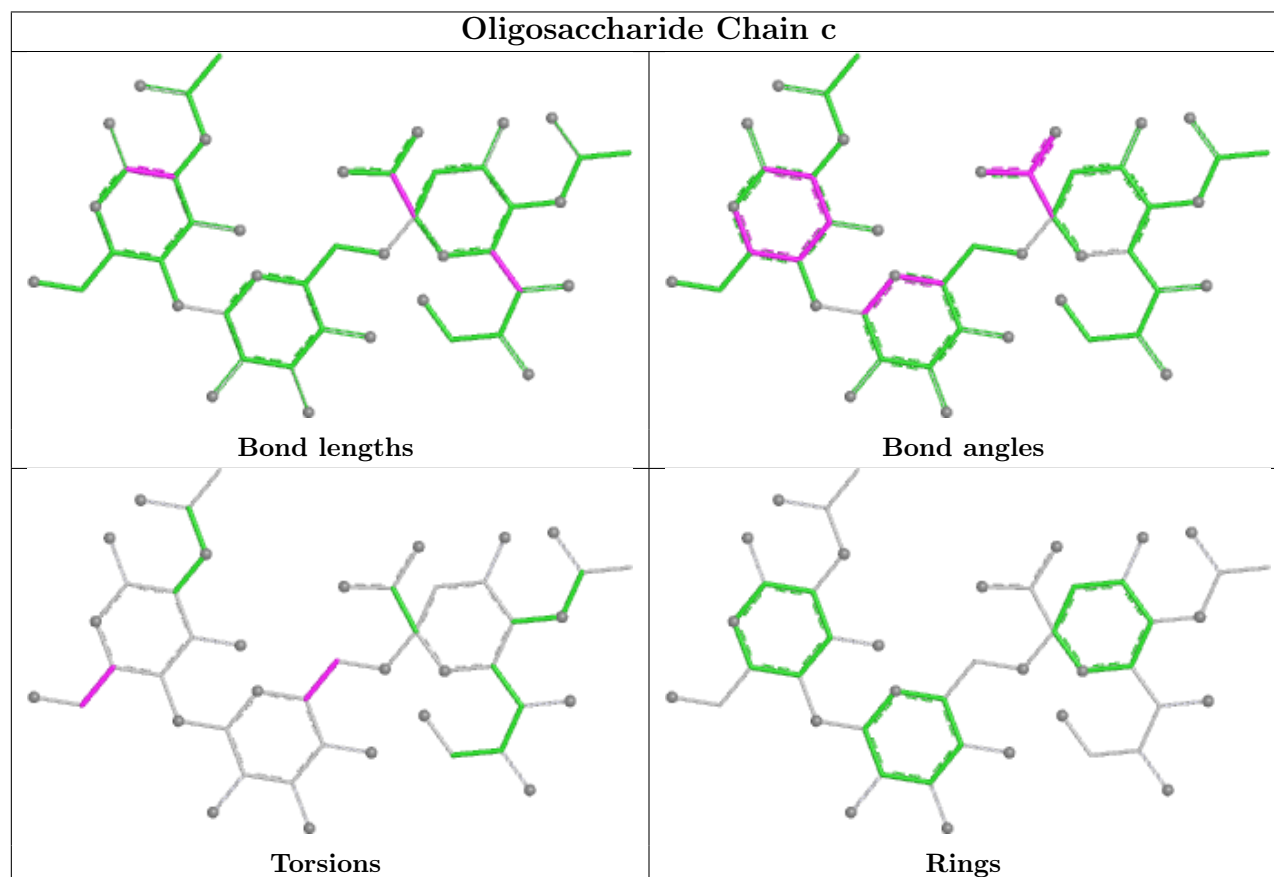
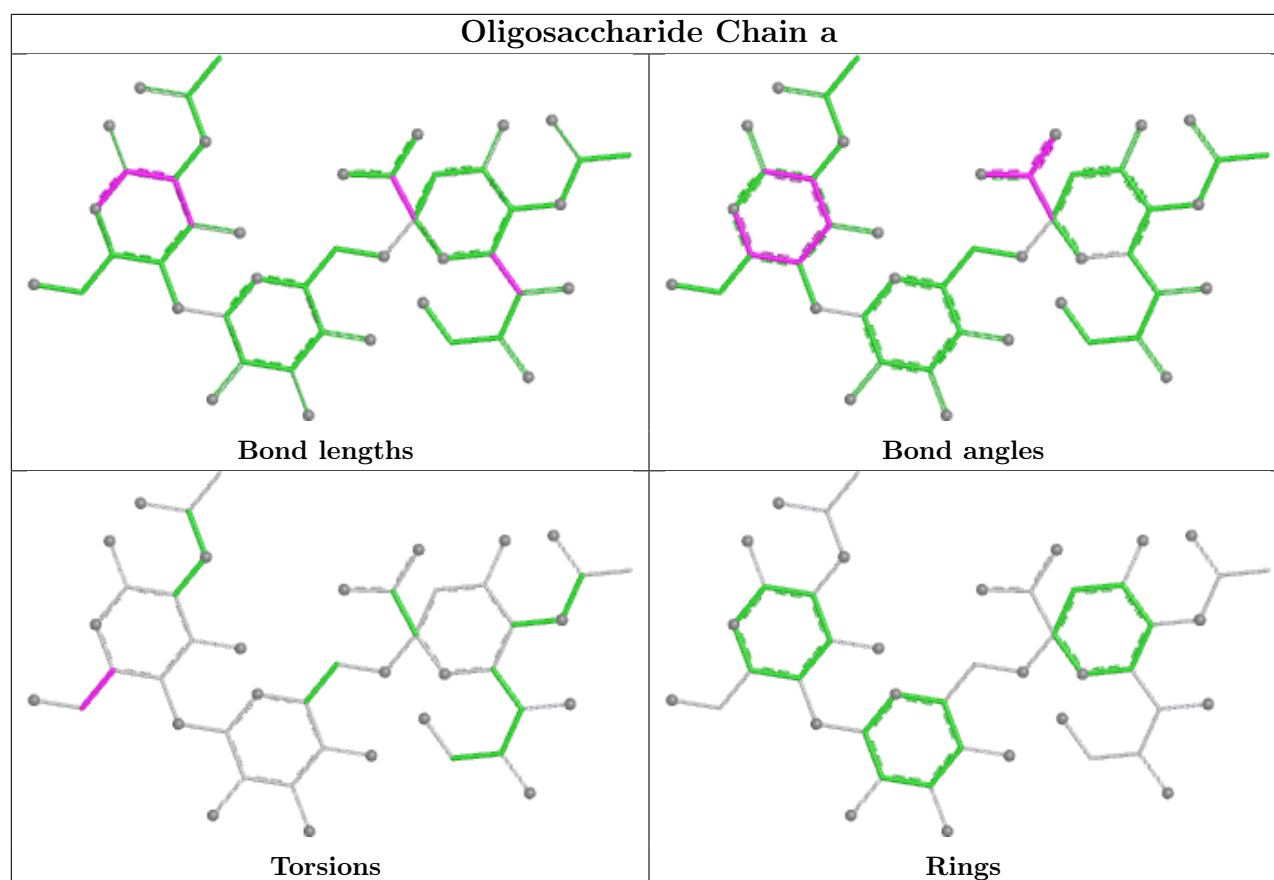




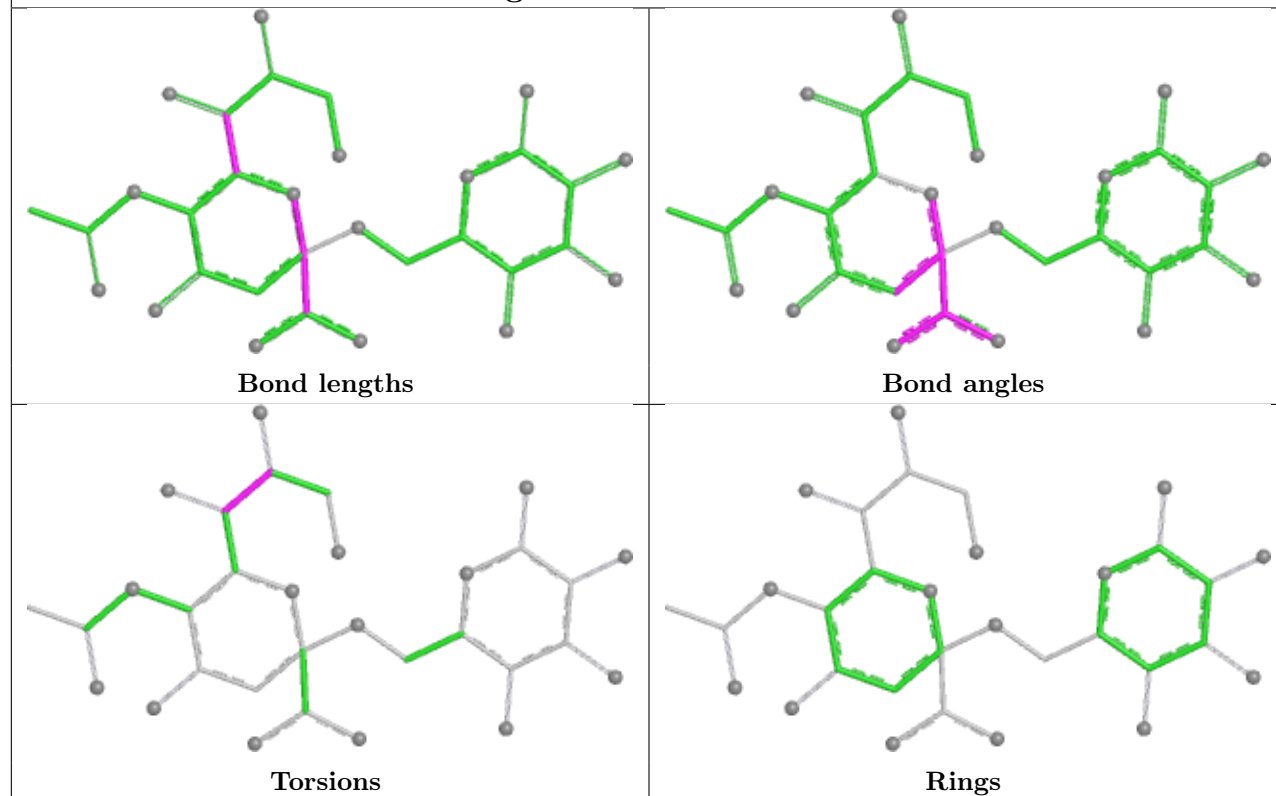




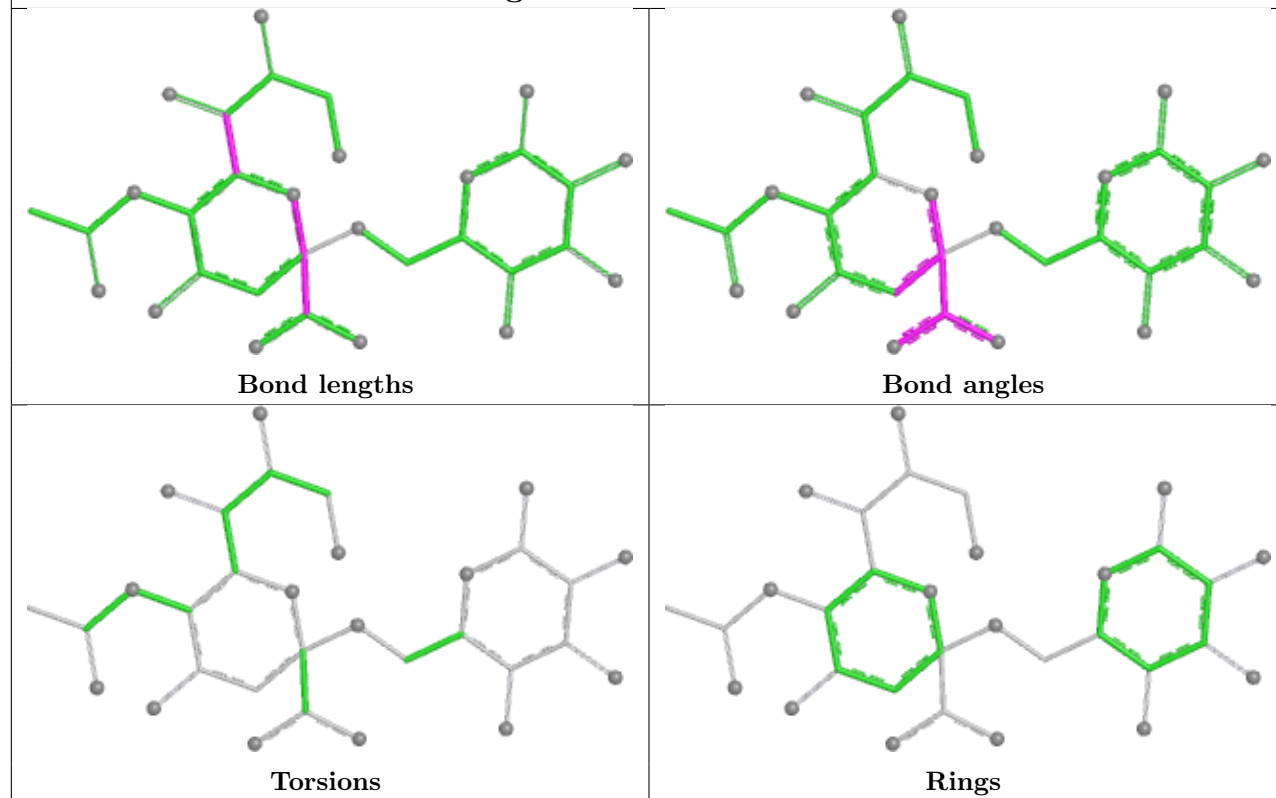




Oligosaccharide Chain b



Oligosaccharide Chain d



5.6 Ligand geometry

14 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
10	NAG	D	201	2	14,14,15	0.42	0	17,19,21	1.06	2 (11%)
10	NAG	H	201	2	14,14,15	0.80	1 (7%)	17,19,21	0.62	0
10	NAG	A	401	1	14,14,15	0.52	0	17,19,21	0.74	0
10	NAG	G	401	1	14,14,15	0.70	0	17,19,21	0.52	0
12	PG0	H	202	-	7,7,7	0.48	0	6,6,6	0.34	0
12	PG0	B	203	-	7,7,7	0.48	0	6,6,6	0.23	0
13	PEG	F	204	-	6,6,6	0.19	0	5,5,5	0.06	0
11	PG4	B	202	-	12,12,12	0.25	0	11,11,11	0.65	0
10	NAG	F	201	2	14,14,15	0.67	1 (7%)	17,19,21	0.51	0
13	PEG	F	203	-	6,6,6	0.18	0	5,5,5	0.16	0
13	PEG	E	402	-	6,6,6	0.21	0	5,5,5	0.16	0
10	NAG	E	401	1	14,14,15	0.64	0	17,19,21	0.50	0
10	NAG	B	201	2	14,14,15	0.65	1 (7%)	17,19,21	0.58	0
11	PG4	F	202	-	12,12,12	0.21	0	11,11,11	0.67	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
10	NAG	D	201	2	-	1/6/23/26	0/1/1/1
10	NAG	H	201	2	-	0/6/23/26	0/1/1/1
10	NAG	A	401	1	-	2/6/23/26	0/1/1/1
10	NAG	G	401	1	-	0/6/23/26	0/1/1/1
12	PG0	H	202	-	-	2/5/5/5	-
12	PG0	B	203	-	-	5/5/5/5	-
13	PEG	F	204	-	-	1/4/4/4	-
11	PG4	B	202	-	-	7/10/10/10	-
10	NAG	F	201	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
13	PEG	F	203	-	-	2/4/4/4	-
13	PEG	E	402	-	-	2/4/4/4	-
10	NAG	E	401	1	-	0/6/23/26	0/1/1/1
10	NAG	B	201	2	-	0/6/23/26	0/1/1/1
11	PG4	F	202	-	-	7/10/10/10	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	H	201	NAG	C1-C2	2.61	1.55	1.52
10	F	201	NAG	C1-C2	2.39	1.55	1.52
10	B	201	NAG	C1-C2	2.26	1.55	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	D	201	NAG	C1-O5-C5	3.36	116.69	112.19
10	D	201	NAG	C2-N2-C7	2.14	125.77	122.90

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
13	F	204	PEG	C1-C2-O2-C3
13	F	203	PEG	C1-C2-O2-C3
12	B	203	PG0	C1-C2-O1-C3
11	F	202	PG4	C4-C3-O2-C2
13	E	402	PEG	C1-C2-O2-C3
11	B	202	PG4	C4-C3-O2-C2
11	F	202	PG4	O1-C1-C2-O2
12	H	202	PG0	O1-C3-C4-O2
11	B	202	PG4	C5-C6-O4-C7
12	B	203	PG0	O1-C3-C4-O2
11	B	202	PG4	O1-C1-C2-O2
13	E	402	PEG	O2-C3-C4-O4
10	D	201	NAG	C3-C2-N2-C7
12	H	202	PG0	OTT-C1-C2-O1
11	F	202	PG4	O3-C5-C6-O4
10	A	401	NAG	C1-C2-N2-C7
12	B	203	PG0	C3-C4-O2-C5

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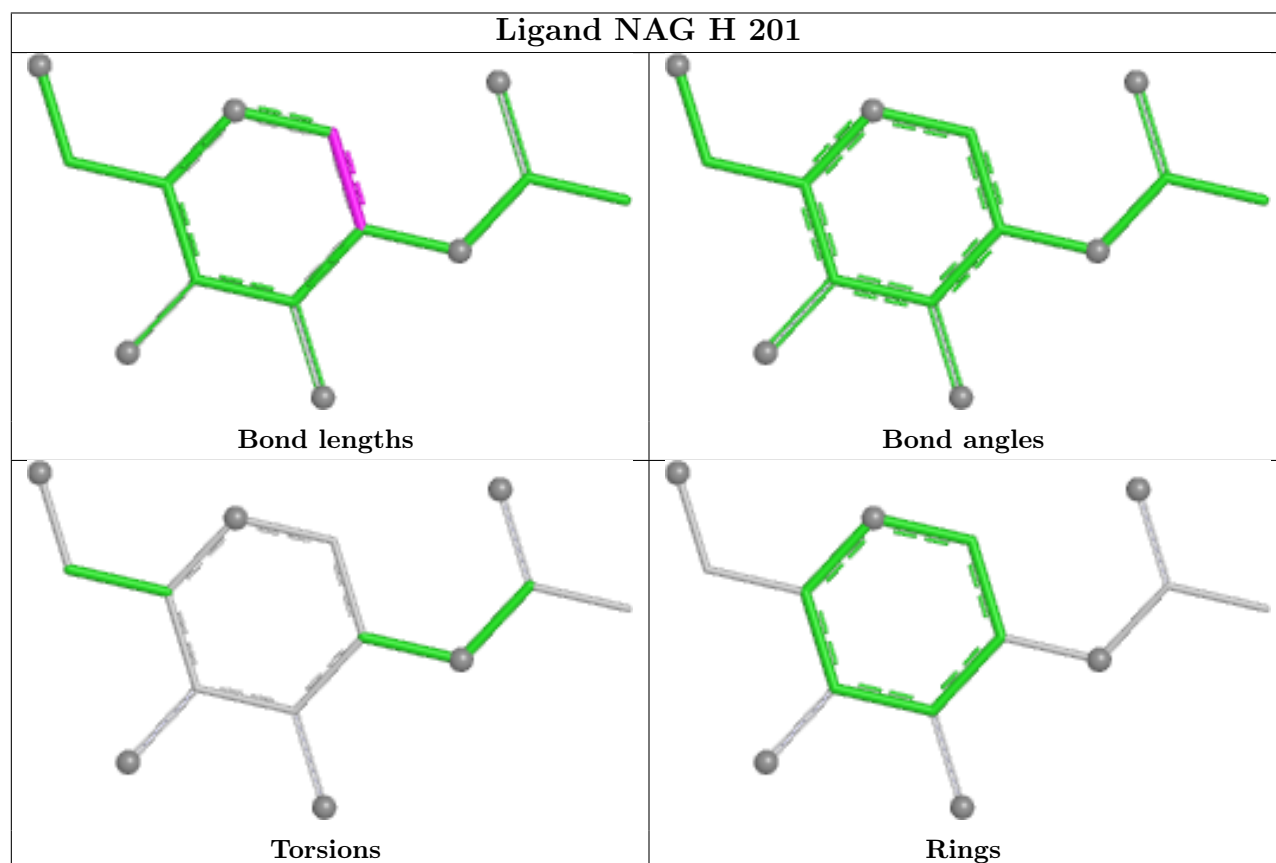
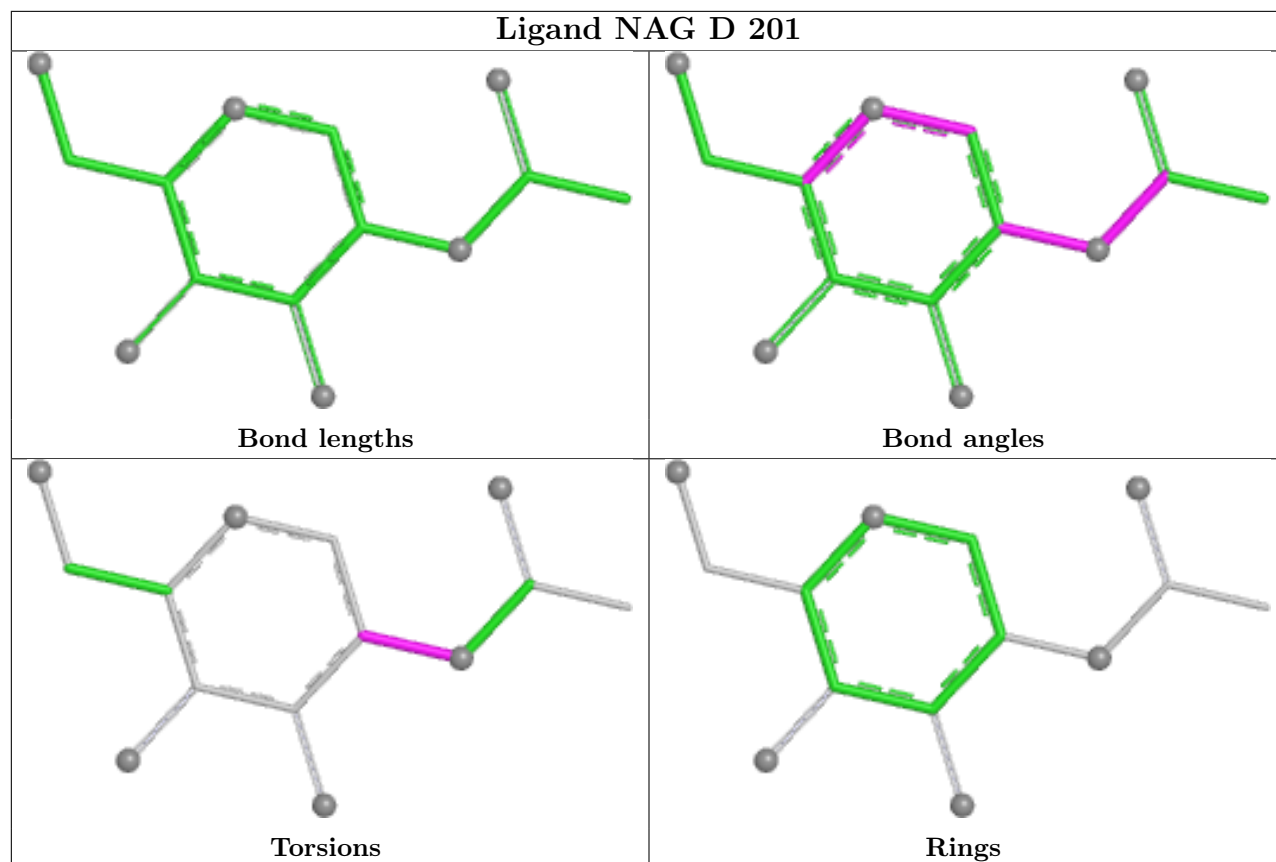
Mol	Chain	Res	Type	Atoms
11	B	202	PG4	C1-C2-O2-C3
11	B	202	PG4	O4-C7-C8-O5
13	F	203	PEG	C4-C3-O2-C2
12	B	203	PG0	OTT-C1-C2-O1
11	F	202	PG4	C8-C7-O4-C6
11	F	202	PG4	C5-C6-O4-C7
11	F	202	PG4	C1-C2-O2-C3
10	F	201	NAG	C4-C5-C6-O6
10	F	201	NAG	O5-C5-C6-O6
11	B	202	PG4	O3-C5-C6-O4
11	B	202	PG4	O2-C3-C4-O3
10	A	401	NAG	C3-C2-N2-C7
11	F	202	PG4	O2-C3-C4-O3
12	B	203	PG0	C4-C3-O1-C2

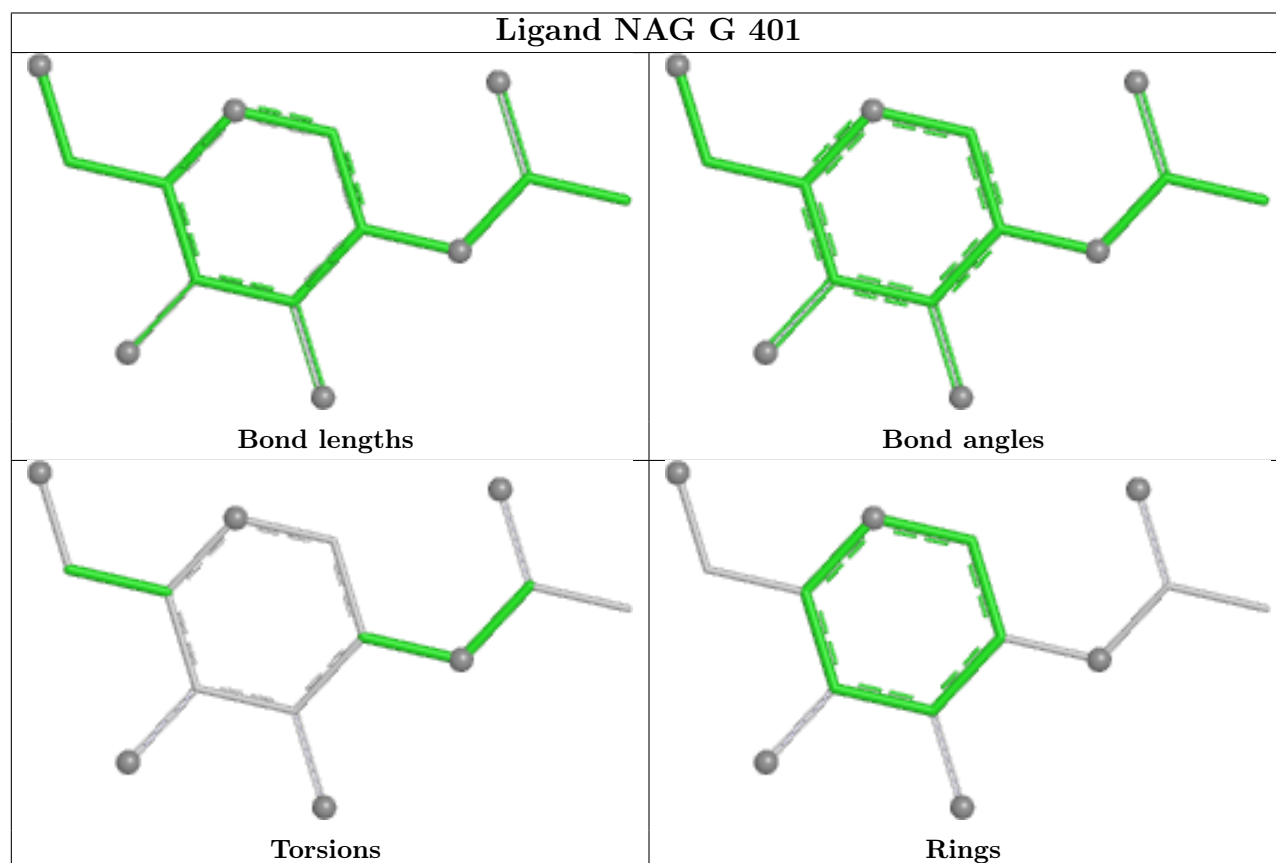
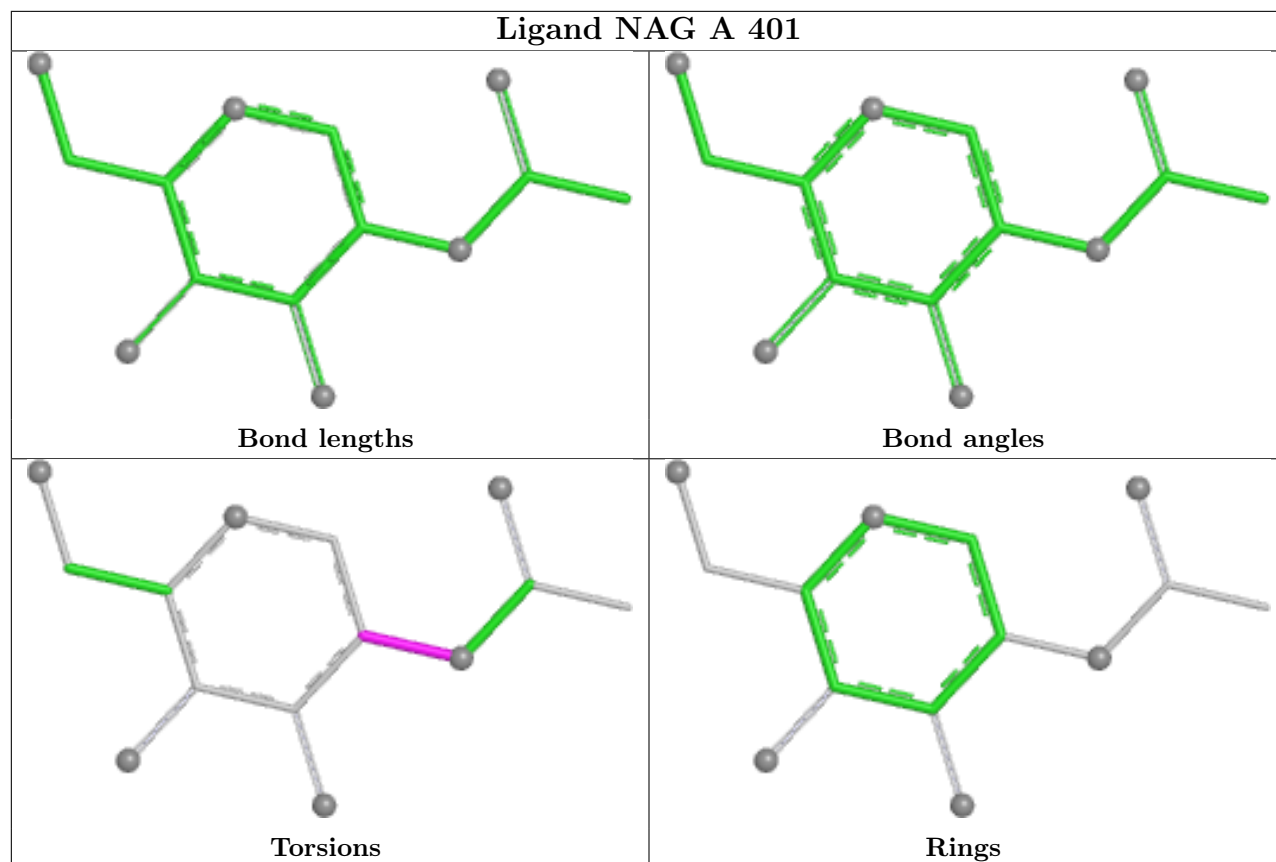
There are no ring outliers.

6 monomers are involved in 8 short contacts:

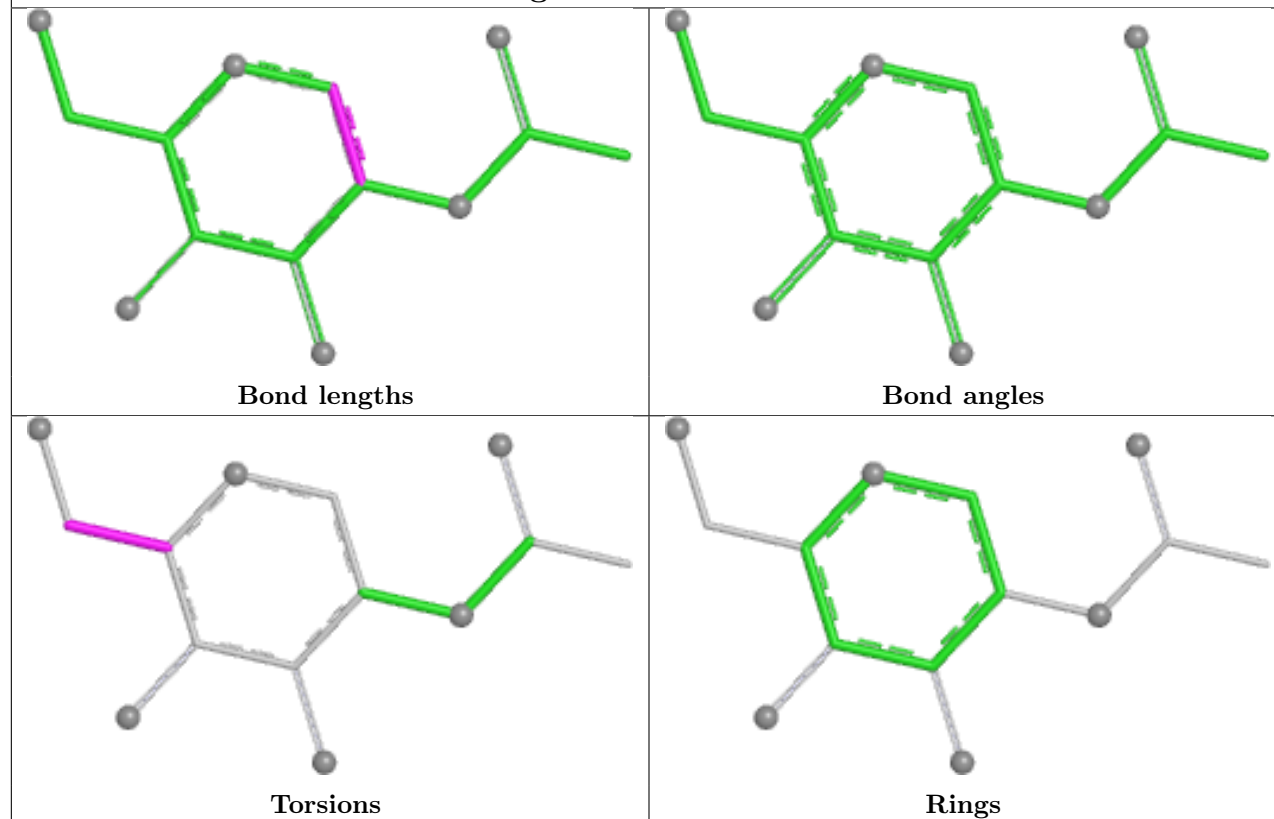
Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	B	203	PG0	1	0
13	F	204	PEG	1	0
11	B	202	PG4	2	0
13	F	203	PEG	2	0
13	E	402	PEG	2	0
11	F	202	PG4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

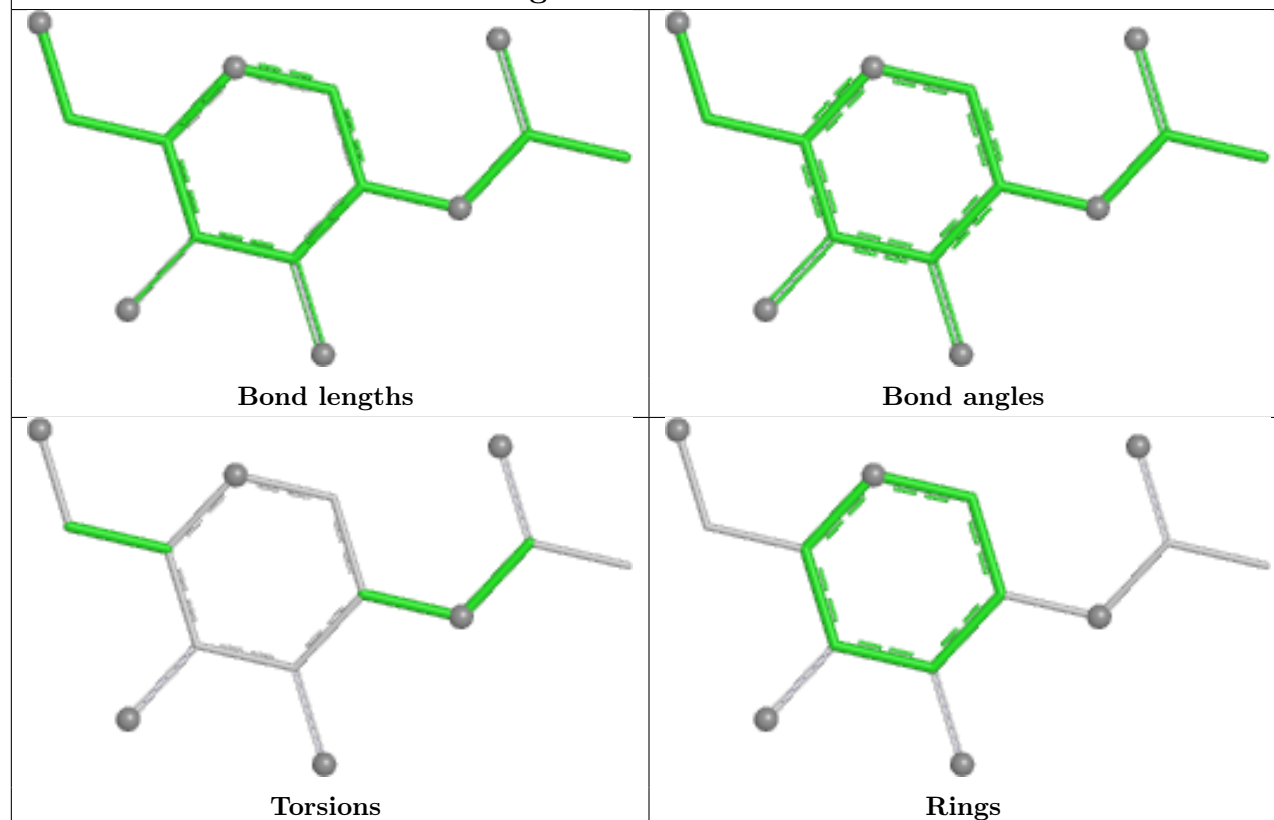


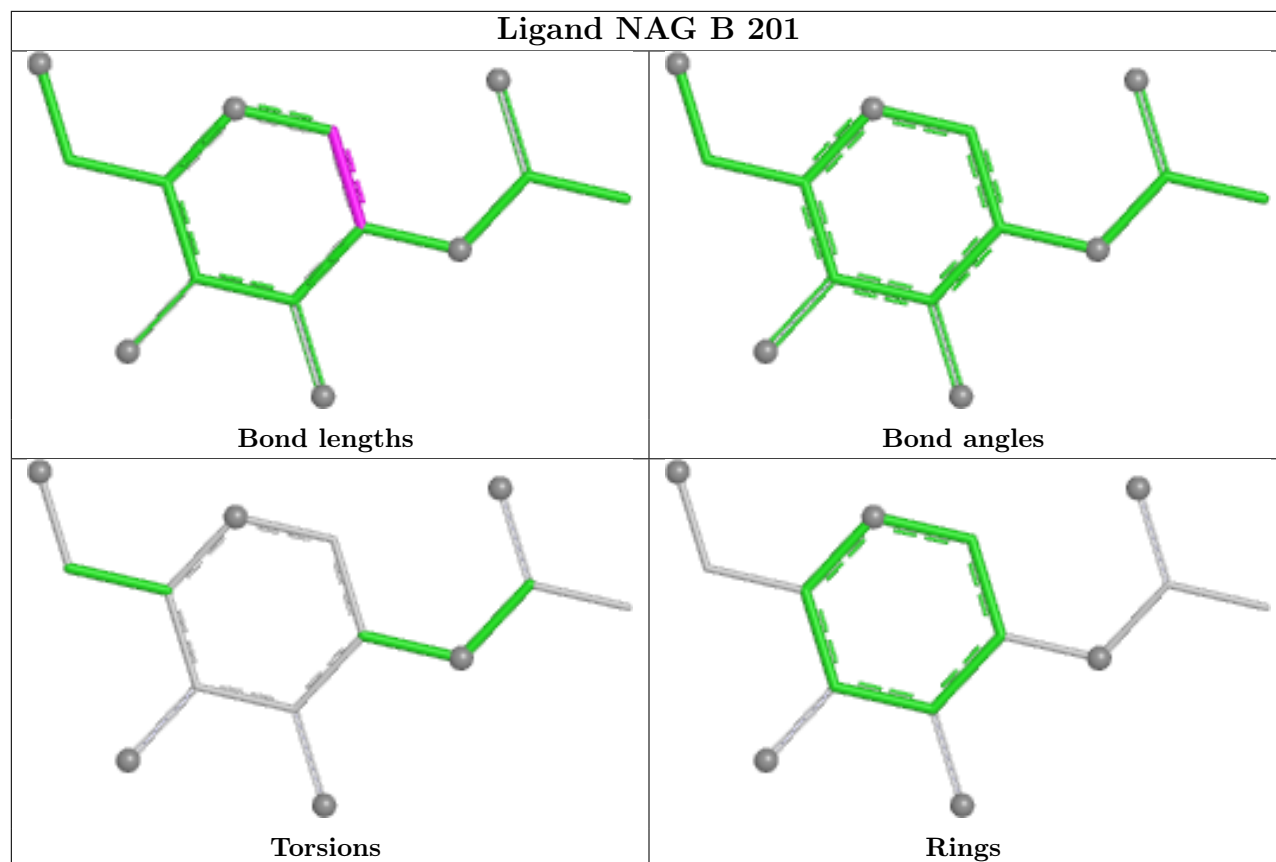


Ligand NAG F 201



Ligand NAG E 401





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	317/323 (98%)	-1.20	0 100 100	18, 36, 58, 76	0
1	C	317/323 (98%)	-1.19	0 100 100	17, 34, 61, 82	0
1	E	317/323 (98%)	-1.18	0 100 100	17, 36, 60, 75	0
1	G	317/323 (98%)	-1.20	0 100 100	17, 35, 61, 83	0
2	B	172/174 (98%)	-1.35	0 100 100	17, 26, 42, 69	0
2	D	172/174 (98%)	-1.33	0 100 100	16, 26, 46, 80	0
2	F	172/174 (98%)	-1.35	0 100 100	18, 26, 43, 68	0
2	H	172/174 (98%)	-1.32	0 100 100	16, 26, 45, 78	0
All	All	1956/1988 (98%)	-1.24	0 100 100	16, 31, 58, 83	0

There are no RSRZ outliers to report.

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	BMA	M	3	11/12	0.93	0.08	92,102,104,105	0
5	MAN	K	4	11/12	0.94	0.06	83,95,99,99	0
4	BMA	N	3	11/12	0.95	0.06	92,102,110,110	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MAN	K	5	11/12	0.95	0.07	98,103,107,108	0
5	MAN	Q	5	11/12	0.95	0.06	76,82,93,94	0
5	MAN	R	4	11/12	0.95	0.07	97,102,104,106	0
6	FUC	Y	4	10/11	0.95	0.07	88,93,99,100	0
7	MAN	Z	6	11/12	0.95	0.06	82,87,89,90	0
3	NAG	O	2	14/15	0.96	0.06	87,96,102,104	0
3	NAG	U	2	14/15	0.96	0.05	85,97,103,105	0
3	NAG	V	2	14/15	0.96	0.06	87,94,99,100	0
4	BMA	J	3	11/12	0.96	0.05	92,100,108,108	0
4	NAG	M	2	14/15	0.96	0.08	61,78,94,102	0
5	MAN	R	5	11/12	0.96	0.05	91,100,104,106	0
6	FUC	P	4	10/11	0.96	0.06	77,85,90,90	0
3	NAG	L	2	14/15	0.96	0.06	89,98,107,107	0
4	NAG	N	2	14/15	0.96	0.06	58,80,92,101	0
5	MAN	Q	4	11/12	0.97	0.06	84,88,92,94	0
3	NAG	T	2	14/15	0.97	0.05	61,71,79,82	0
3	NAG	I	1	14/15	0.97	0.05	46,62,69,72	0
3	NAG	I	2	14/15	0.97	0.06	66,80,92,92	0
6	FUC	P	2	10/11	0.97	0.04	59,64,67,68	0
6	NAG	P	3	14/15	0.97	0.05	74,81,86,87	0
3	NAG	W	2	14/15	0.97	0.06	62,72,86,90	0
6	FUC	Y	2	10/11	0.97	0.05	58,68,71,71	0
6	NAG	Y	3	14/15	0.97	0.05	75,83,87,87	0
3	NAG	X	2	14/15	0.97	0.06	59,77,83,87	0
7	MAN	Z	4	11/12	0.97	0.04	82,85,94,97	0
7	MAN	Z	5	11/12	0.97	0.05	82,93,98,99	0
3	NAG	S	2	14/15	0.97	0.05	54,76,86,90	0
4	NAG	J	2	14/15	0.98	0.04	58,72,93,98	0
6	NAG	P	1	14/15	0.98	0.04	43,61,74,80	0
5	NAG	K	2	14/15	0.98	0.05	41,55,72,76	0
5	BMA	K	3	11/12	0.98	0.04	71,78,88,90	0
3	NAG	S	1	14/15	0.98	0.05	46,62,67,71	0
6	NAG	Y	1	14/15	0.98	0.04	45,62,77,84	0
3	NAG	U	1	14/15	0.98	0.04	53,65,80,89	0
5	BMA	Q	3	11/12	0.98	0.04	63,69,77,87	0
3	NAG	X	1	14/15	0.98	0.04	30,42,54,60	0
7	BMA	Z	3	11/12	0.98	0.04	62,70,79,85	0
4	NAG	N	1	14/15	0.98	0.04	32,39,52,65	0
5	BMA	R	3	11/12	0.98	0.04	77,82,91,101	0
3	NAG	T	1	14/15	0.98	0.05	30,41,56,58	0
5	NAG	Q	1	14/15	0.99	0.04	47,54,65,66	0
5	NAG	Q	2	14/15	0.99	0.04	44,55,60,62	0

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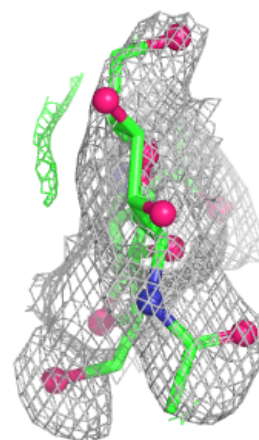
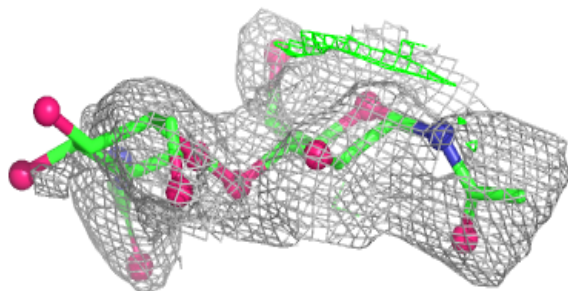
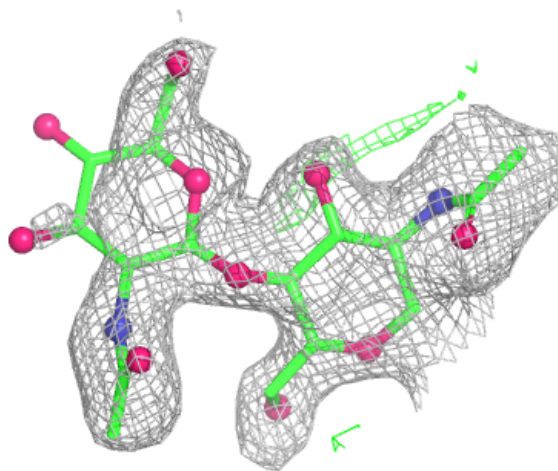
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	W	1	14/15	0.99	0.03	45,50,60,61	0
3	NAG	L	1	14/15	0.99	0.04	55,66,81,89	0
5	NAG	K	1	14/15	0.99	0.04	45,51,57,59	0
5	NAG	R	1	14/15	0.99	0.04	45,51,58,59	0
5	NAG	R	2	14/15	0.99	0.04	41,60,75,79	0
7	NAG	Z	1	14/15	0.99	0.03	51,55,62,62	0
7	NAG	Z	2	14/15	0.99	0.03	47,58,63,63	0
4	NAG	M	1	14/15	0.99	0.04	44,52,60,60	0
3	NAG	V	1	14/15	0.99	0.04	55,62,72,84	0
3	NAG	O	1	14/15	0.99	0.03	54,61,71,86	0
4	NAG	J	1	14/15	0.99	0.03	27,39,52,56	0
8	NAG	a	1	15/15	-	-	73,96,107,111	0
8	GAL	a	2	11/12	-	-	45,72,85,90	0
8	SIA	a	3	20/21	-	-	37,45,52,55	0
8	NAG	c	1	15/15	-	-	71,97,103,103	0
8	GAL	c	2	11/12	-	-	48,72,89,95	0
8	SIA	c	3	20/21	-	-	39,46,53,53	0
9	GAL	b	1	12/12	-	-	49,72,79,87	0
9	SIA	b	2	20/21	-	-	36,45,51,59	0
9	GAL	d	1	12/12	-	-	43,65,79,79	0
9	SIA	d	2	20/21	-	-	36,44,50,55	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.

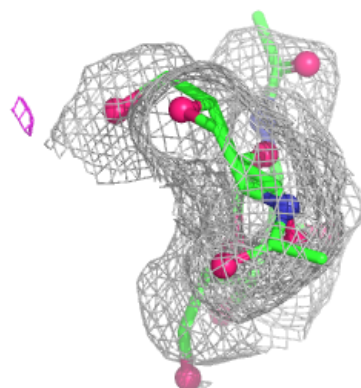
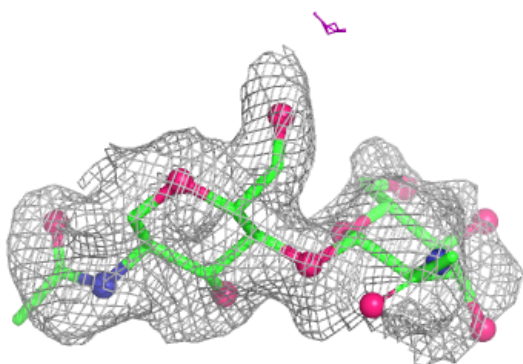
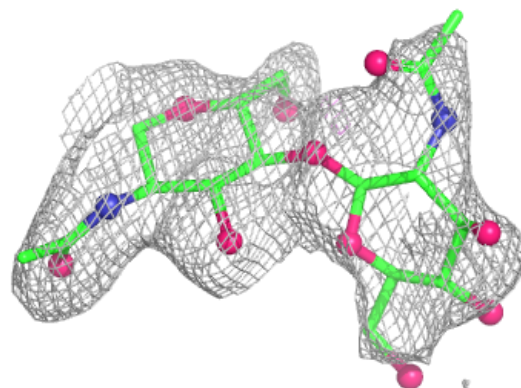
Electron density around Chain I:

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

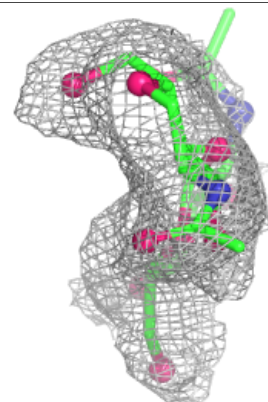
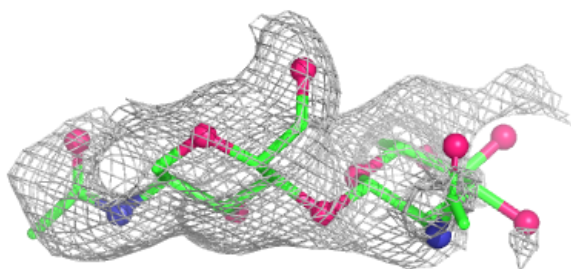
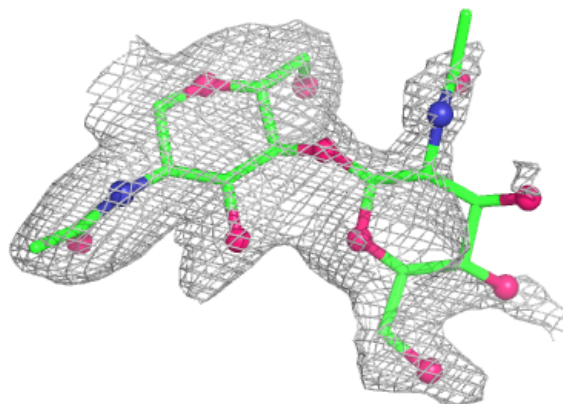


Electron density around Chain L:

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

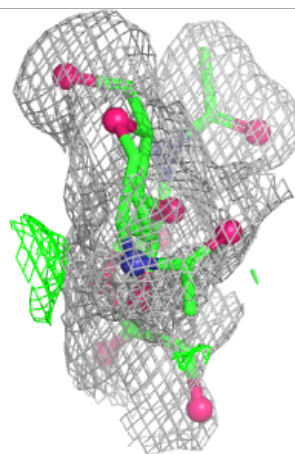
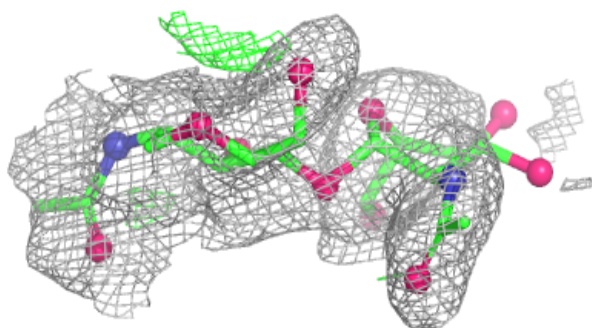
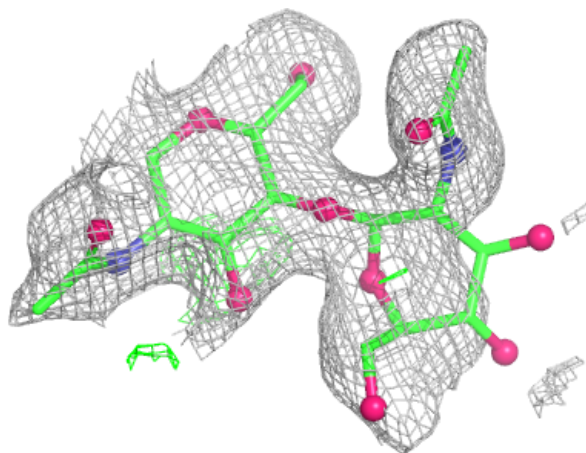
**Electron density around Chain O:**

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



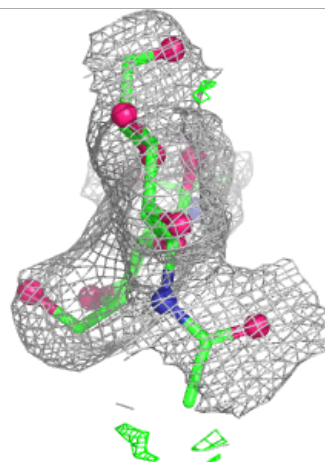
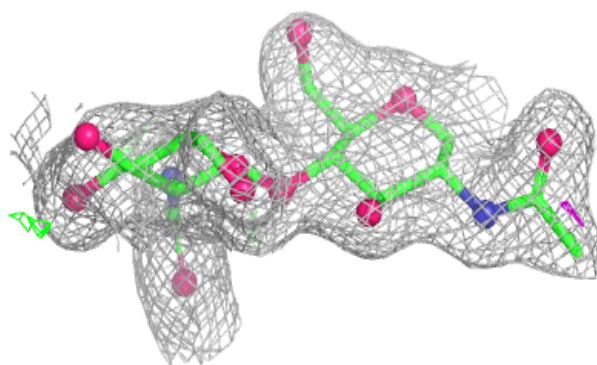
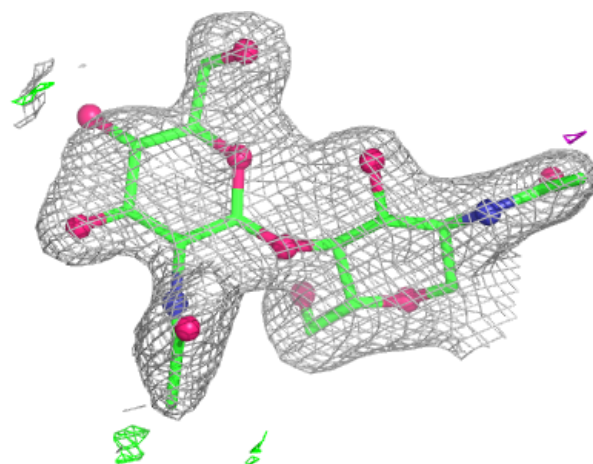
Electron density around Chain S:

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



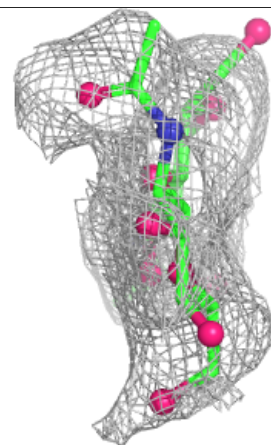
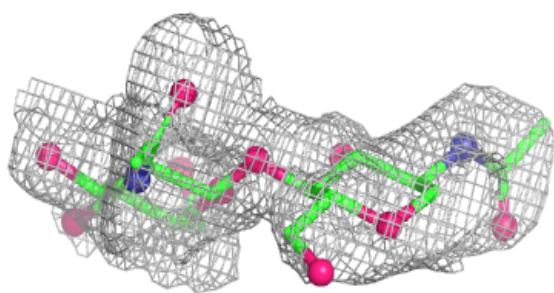
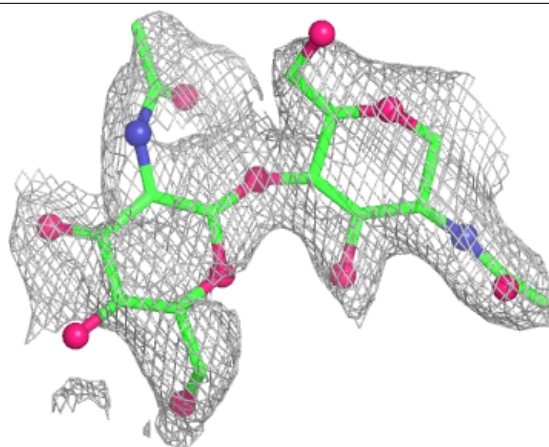
Electron density around Chain T:

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



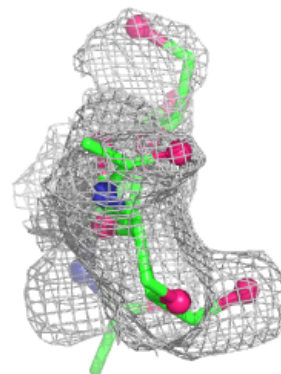
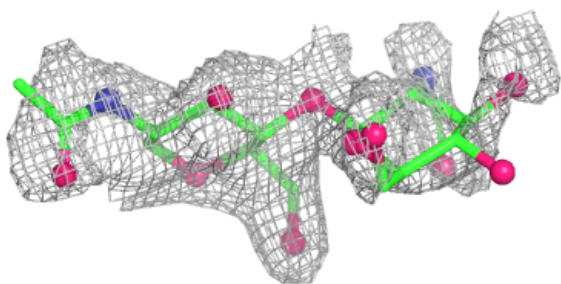
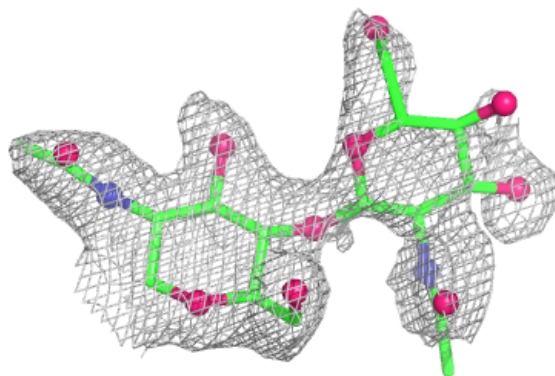
Electron density around Chain U:

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

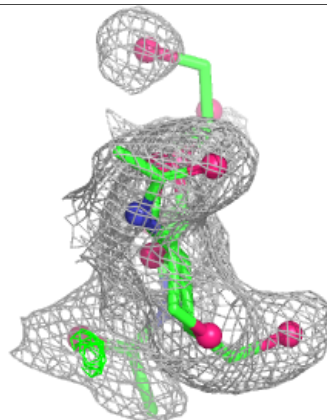
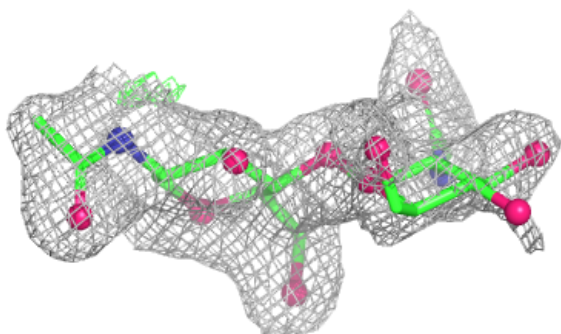
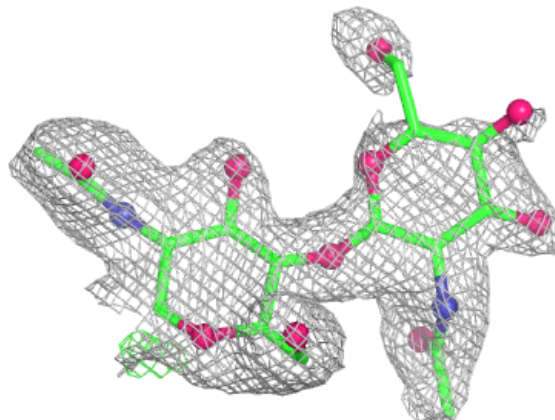


Electron density around Chain V:

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

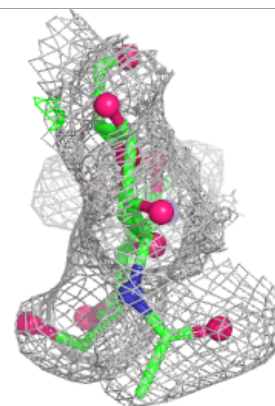
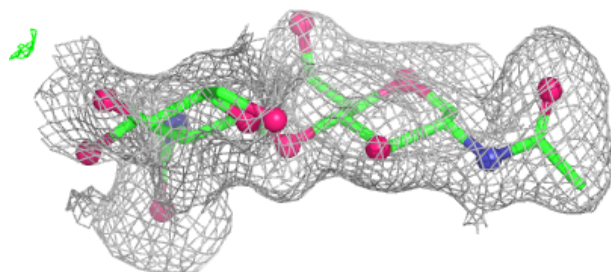
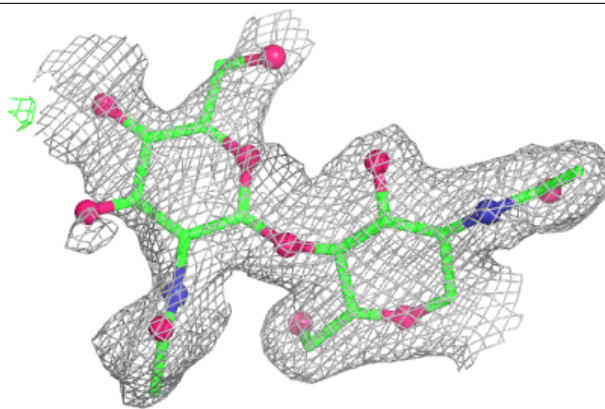
**Electron density around Chain W:**

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

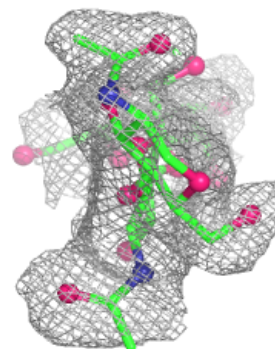
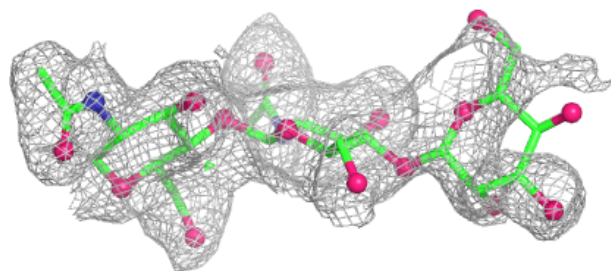
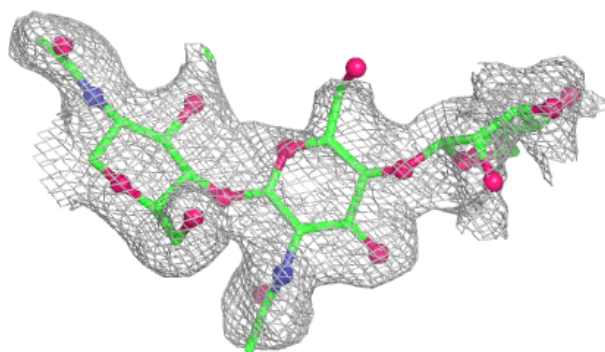


Electron density around Chain X:

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

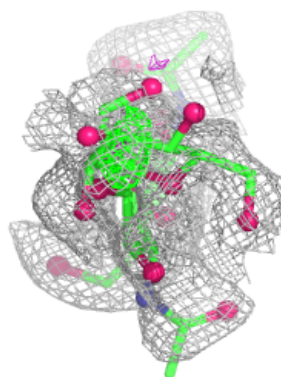
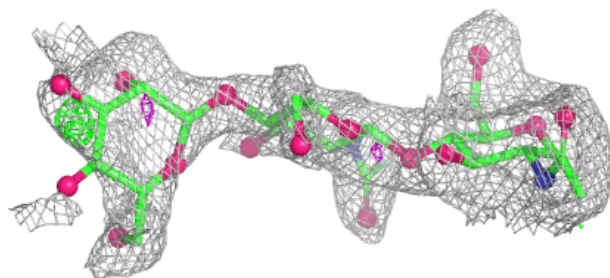
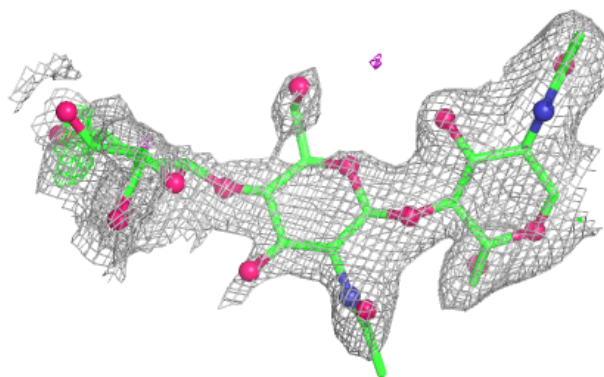
**Electron density around Chain J:**

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

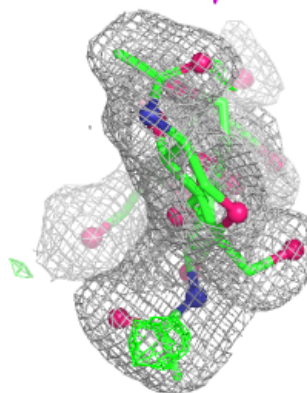
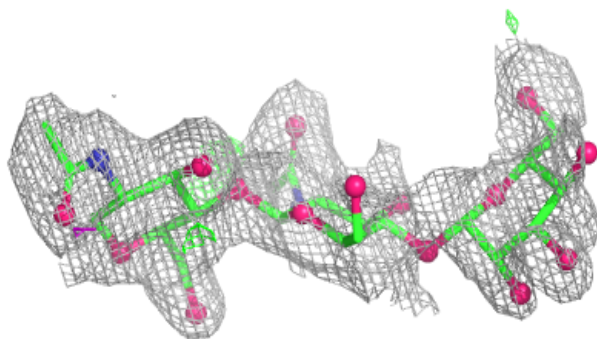
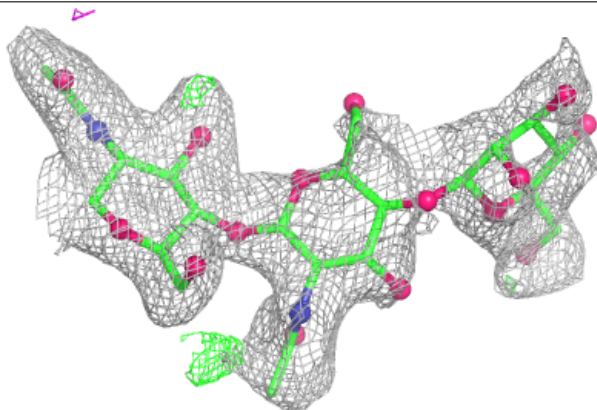


Electron density around Chain M:

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

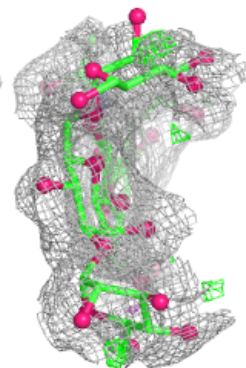
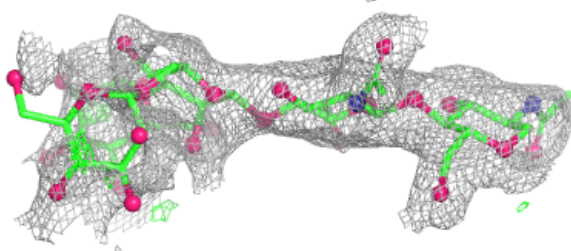
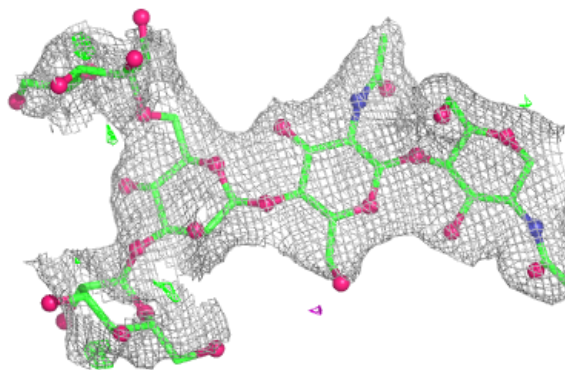
**Electron density around Chain N:**

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

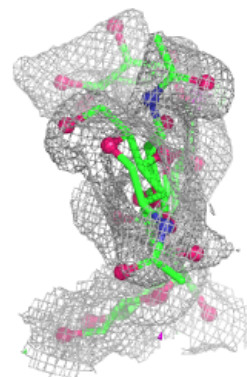
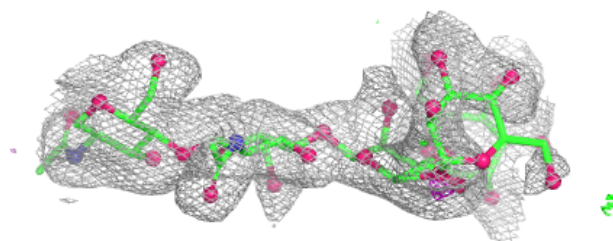
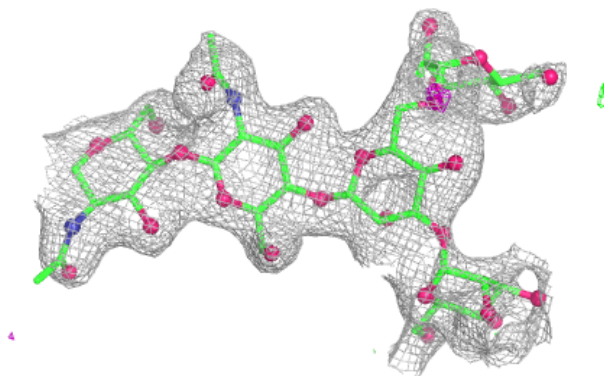


Electron density around Chain K:

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

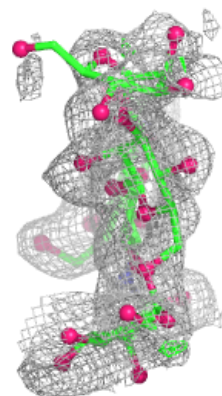
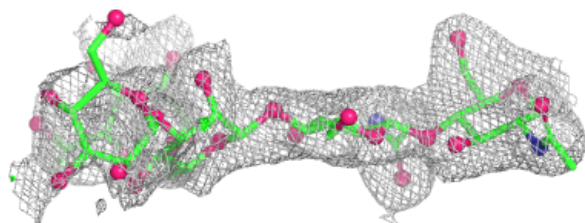
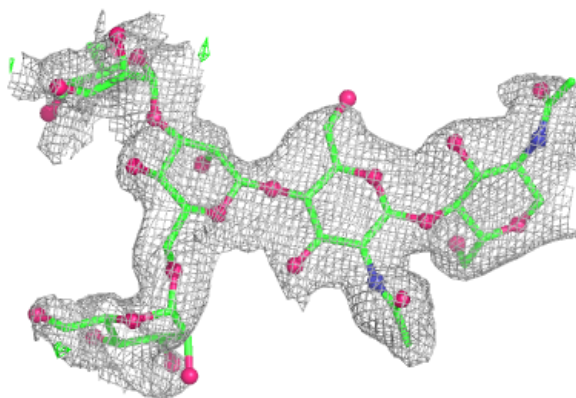
**Electron density around Chain Q:**

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



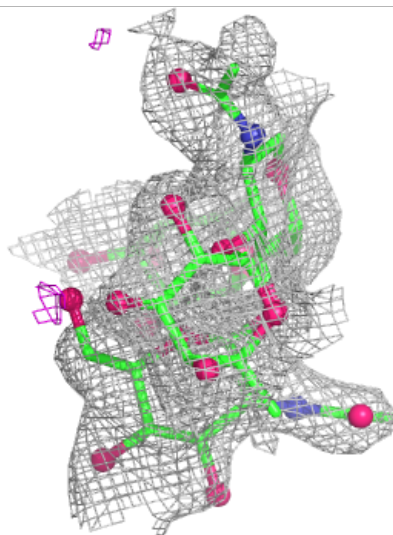
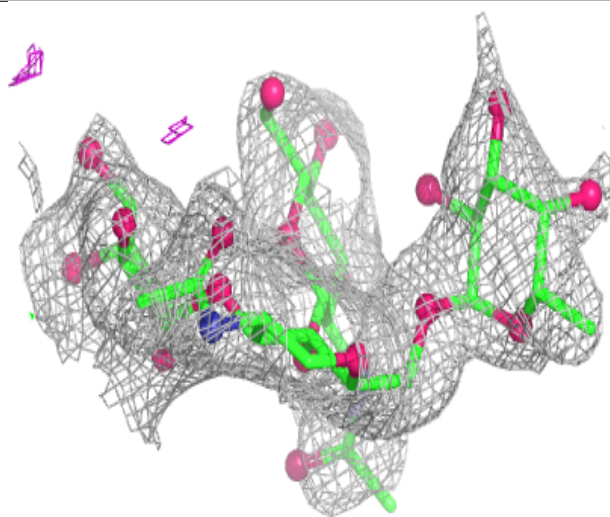
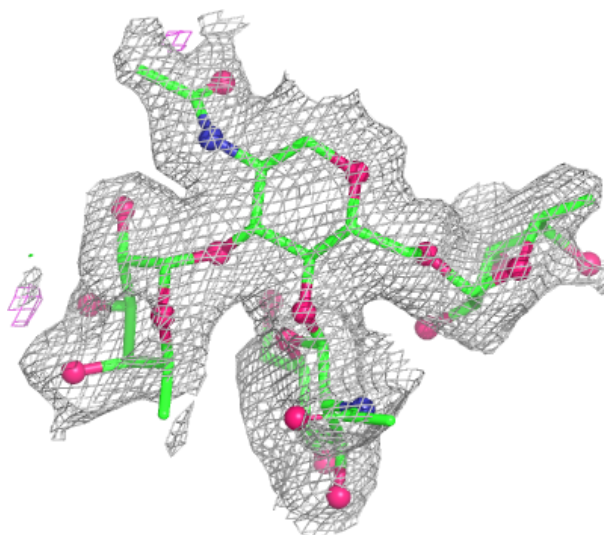
Electron density around Chain R:

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



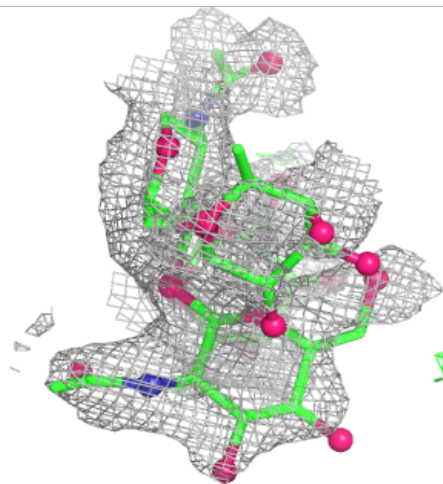
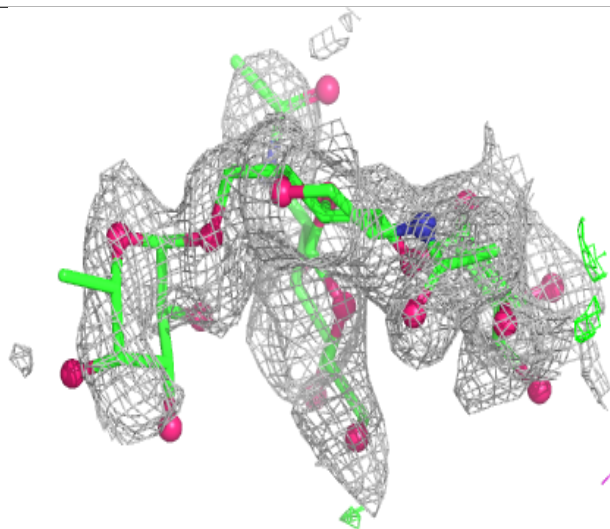
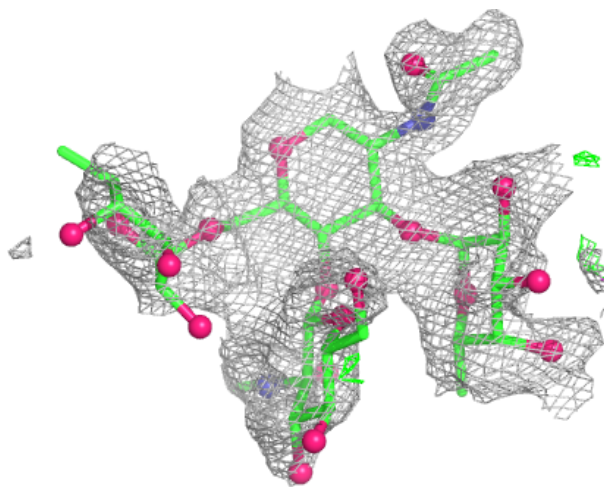
Electron density around Chain P:

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



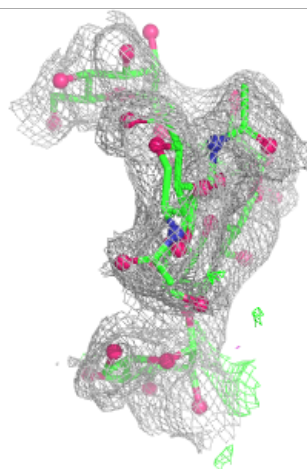
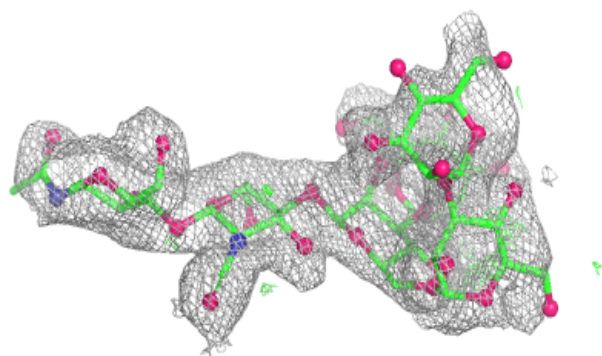
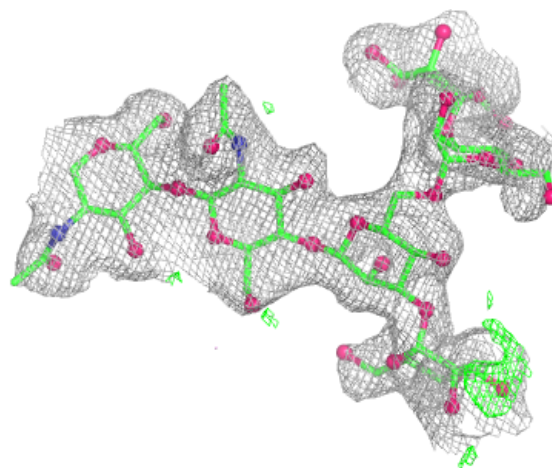
Electron density around Chain Y:

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



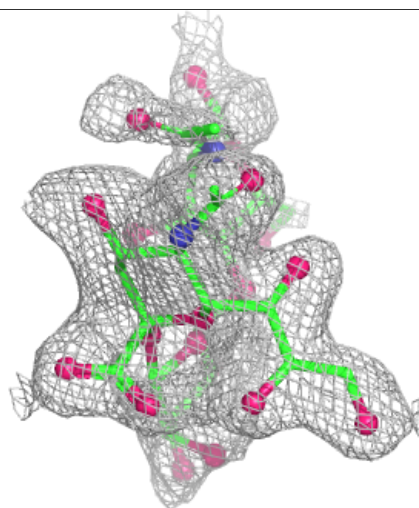
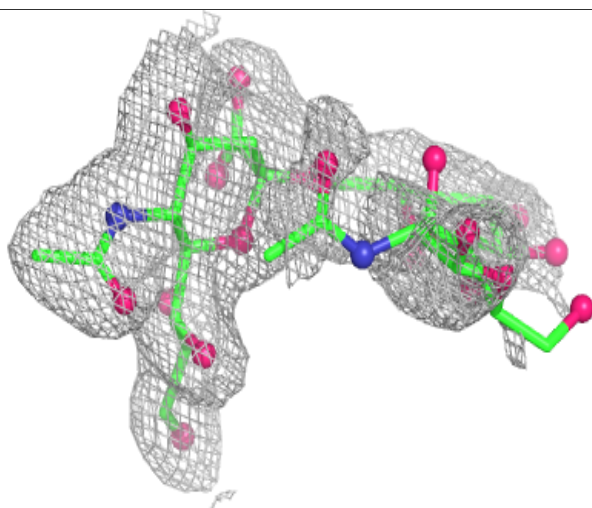
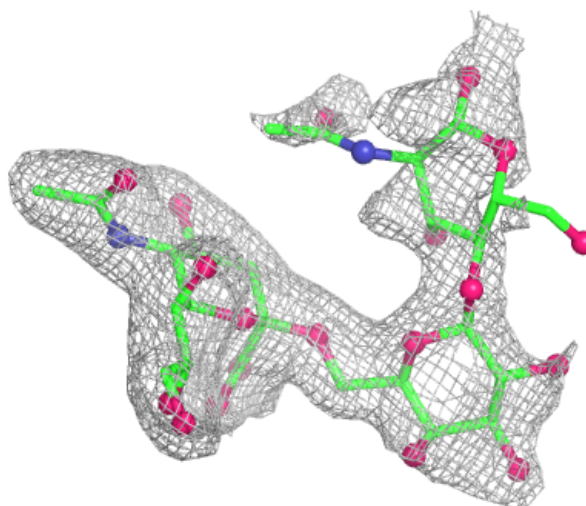
Electron density around Chain Z:

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



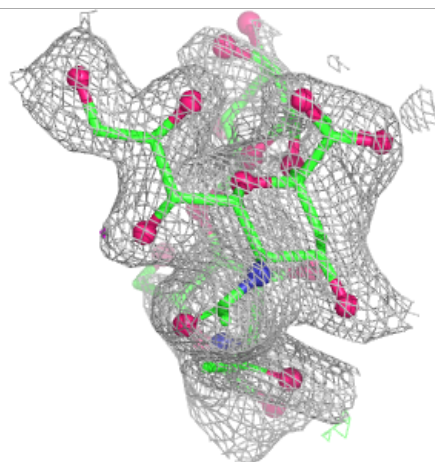
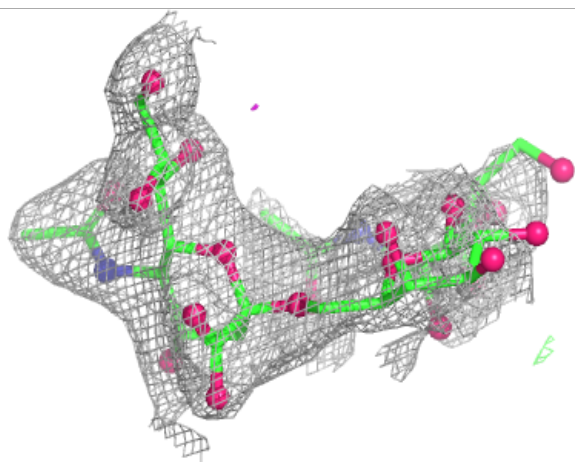
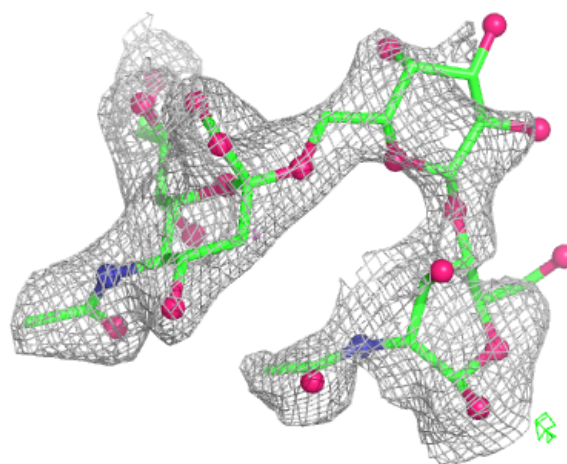
Electron density around Chain a:

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



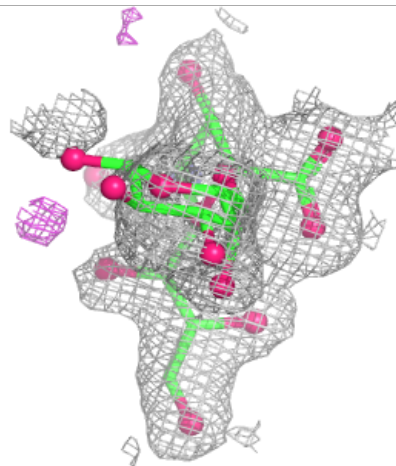
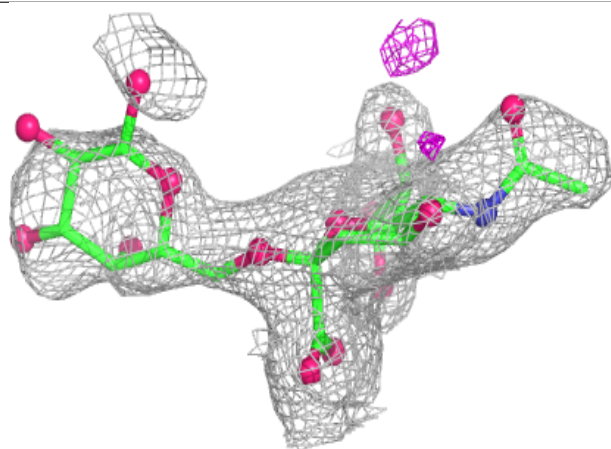
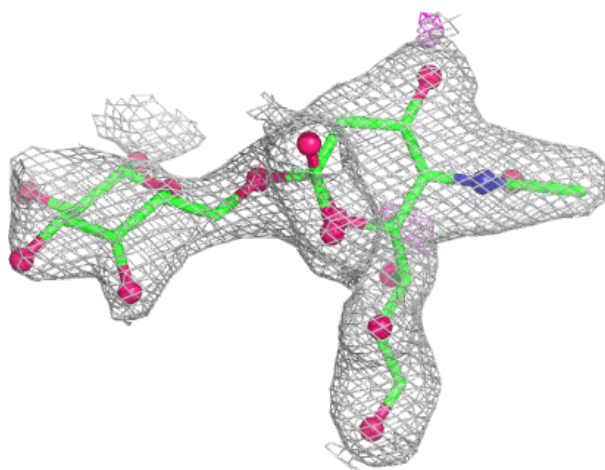
Electron density around Chain c:

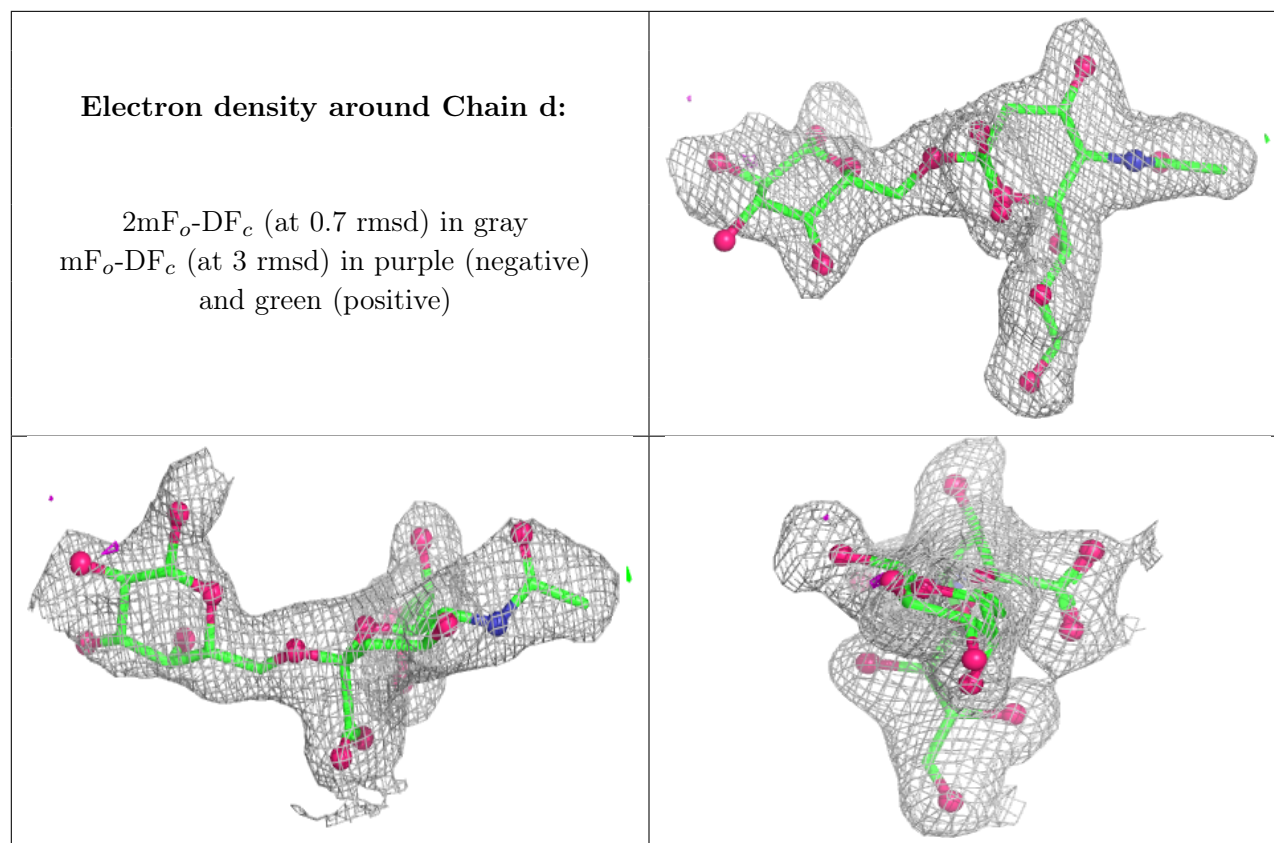
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain b:

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



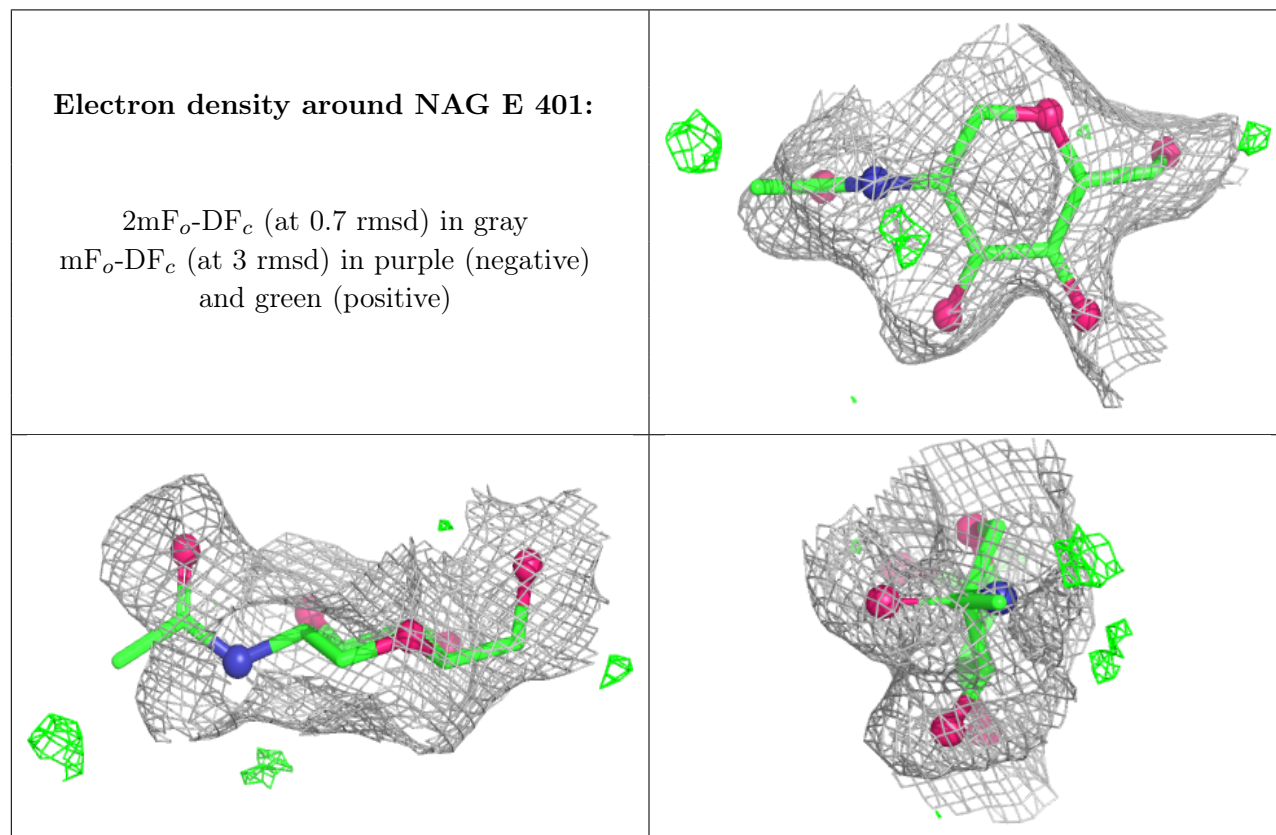


6.4 Ligands [i](#)

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.

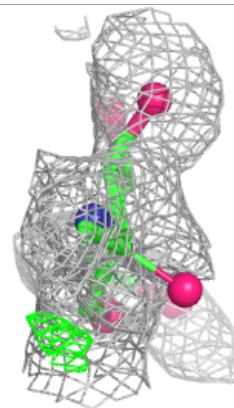
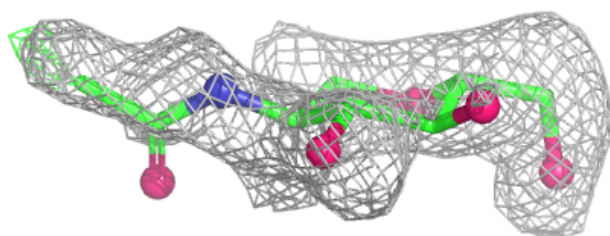
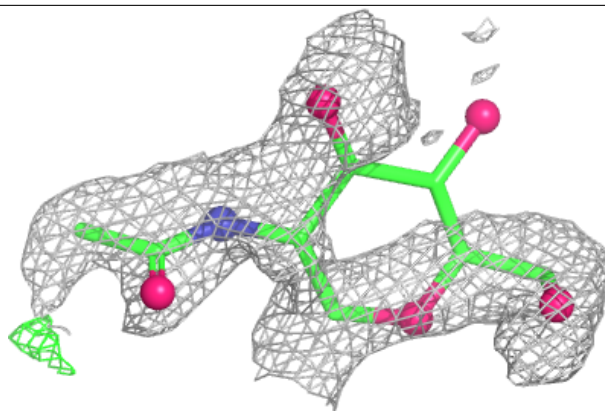
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
10	NAG	E	401	14/15	0.96	0.05	71,75,80,82	0
10	NAG	H	201	14/15	0.96	0.06	45,74,81,82	0
10	NAG	F	201	14/15	0.97	0.05	47,58,70,71	0
10	NAG	G	401	14/15	0.97	0.04	70,76,79,80	0
10	NAG	A	401	14/15	0.97	0.06	67,79,87,92	0
11	PG4	B	202	13/13	0.97	0.06	44,50,60,61	0
12	PG0	B	203	8/8	0.97	0.07	42,54,61,66	0
12	PG0	H	202	8/8	0.97	0.06	42,48,50,59	0
11	PG4	F	202	13/13	0.98	0.06	41,51,59,61	0
10	NAG	D	201	14/15	0.98	0.04	48,70,74,77	0
10	NAG	B	201	14/15	0.98	0.05	46,57,70,73	0
13	PEG	E	402	7/7	0.98	0.06	37,42,50,53	0
13	PEG	F	203	7/7	0.98	0.05	47,52,58,59	0
13	PEG	F	204	7/7	0.98	0.05	46,51,55,63	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

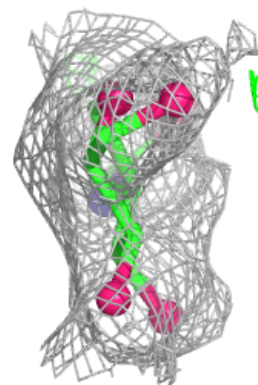
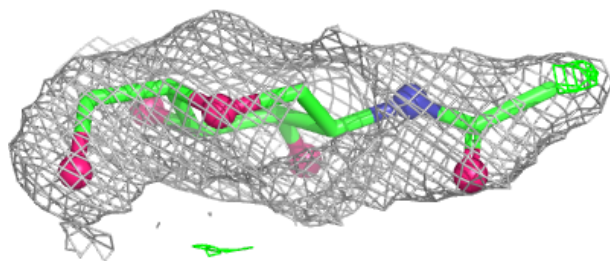
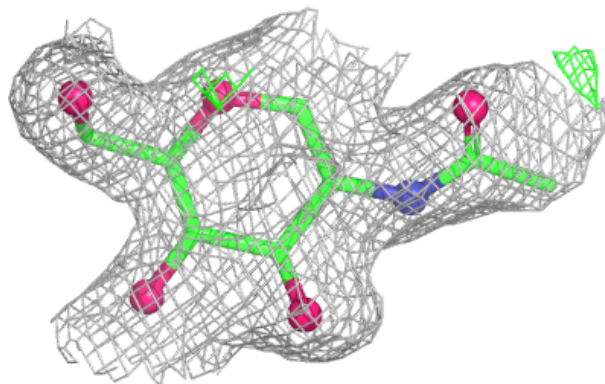


Electron density around NAG H 201:

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

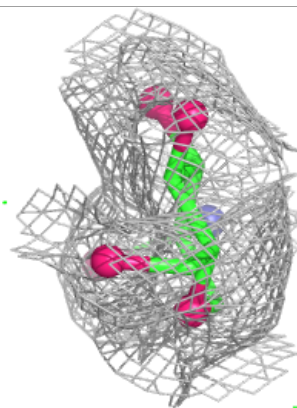
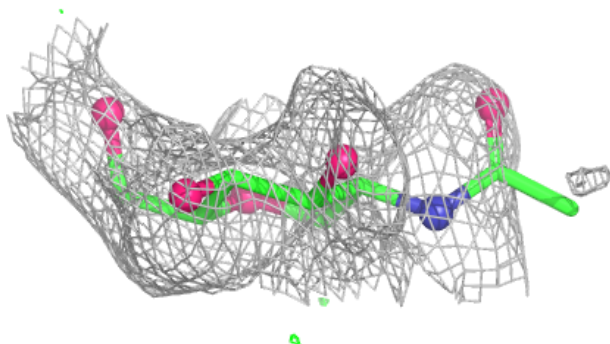
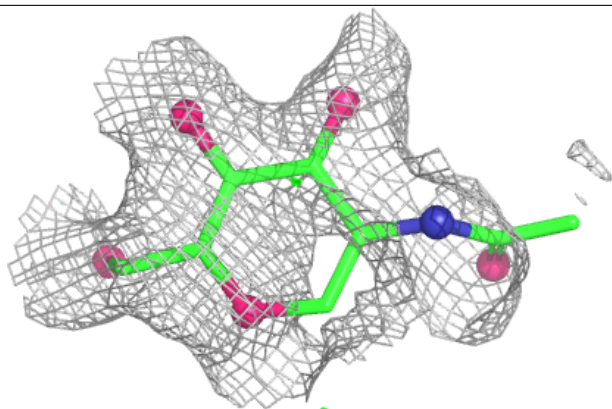
**Electron density around NAG F 201:**

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

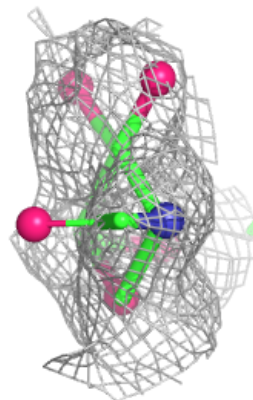
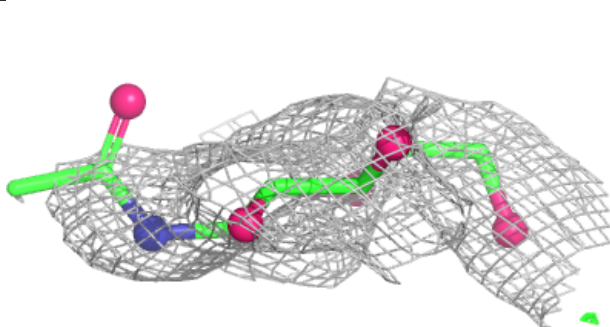
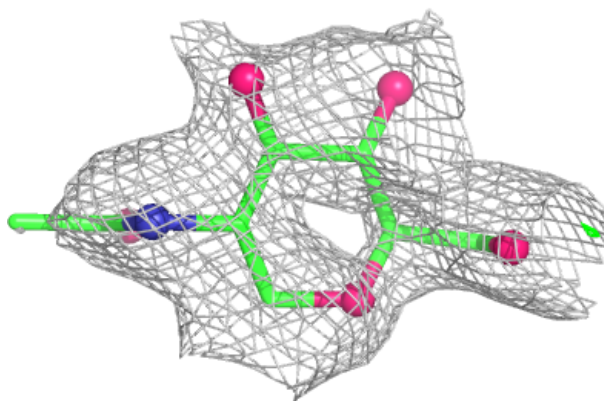


Electron density around NAG G 401:

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

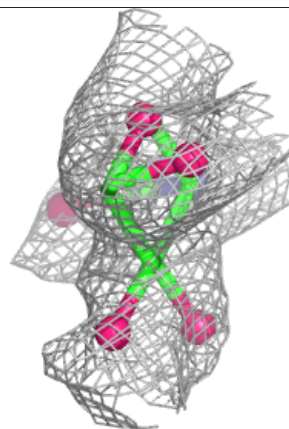
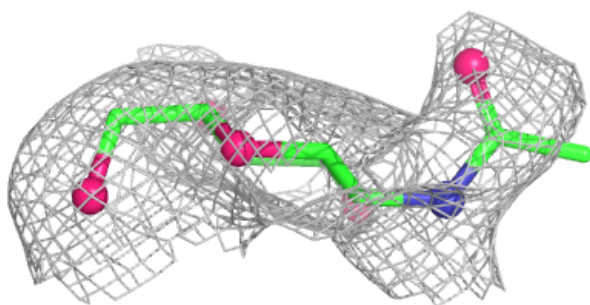
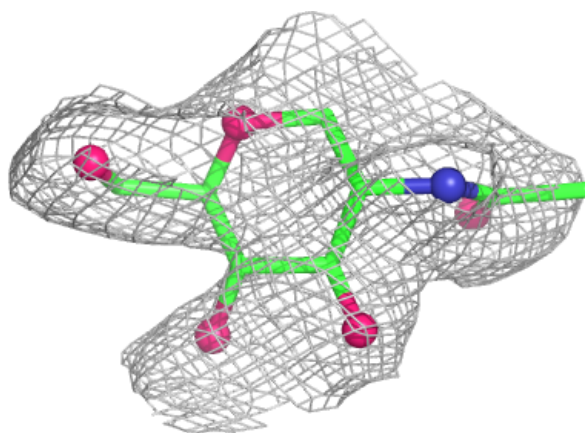
**Electron density around NAG A 401:**

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

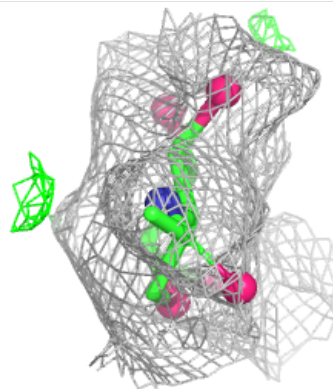
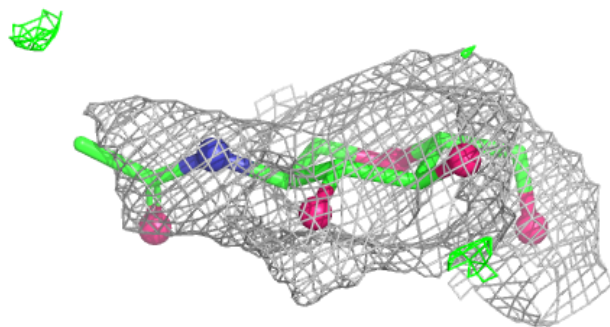
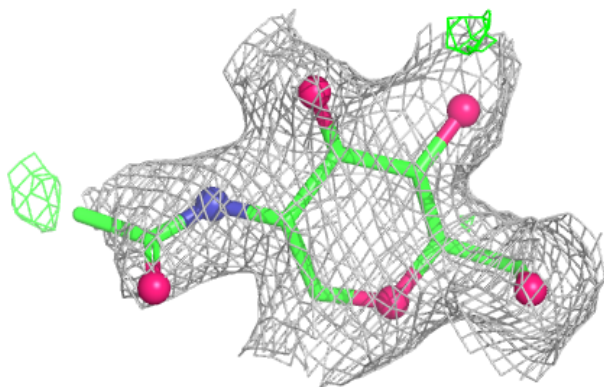


Electron density around NAG D 201:

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

**Electron density around NAG B 201:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.