



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 8, 2020 – 10:02 AM BST

PDB ID : 4UO5
Title : Structure of the A_Canine_Colorado_17864_06 H3 haemagglutinin in complex with 3SLN
Authors : Vachieri, S.G.; Collins, P.J.; Haire, L.F.; Ogradowicz, R.W.; Martin, S.R.; Walker, P.A.; Xiong, X.; Gamblin, S.J.; Skehel, J.J.
Deposited on : 2014-05-31
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

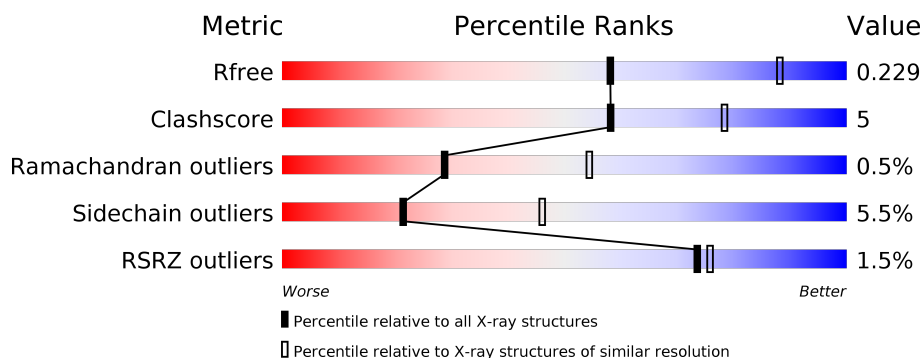
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 on 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	328	<div> <div>2%</div> <div>84%</div> <div>12%</div> <div>..</div> </div>
1	C	328	<div> <div>2%</div> <div>82%</div> <div>13%</div> <div>..</div> </div>
1	E	328	<div> <div>2%</div> <div>80%</div> <div>16%</div> <div>..</div> </div>
2	B	175	<div> <div>2%</div> <div>87%</div> <div>11%</div> <div>..</div> </div>
2	D	175	<div> <div>2%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>
2	F	175	<div> <div>2%</div> <div>83%</div> <div>15%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
3	G	2	 100%
3	L	2	 100%
3	Q	2	 100%
3	R	2	 50% 50%
3	T	2	 100%
4	H	3	 100%
4	J	3	 100%
4	M	3	 33% 67%
5	I	5	 40% 60%
5	N	5	 40% 60%
5	S	5	 20% 80%
6	K	3	 33% 67%
6	P	3	 100%
6	U	3	 33% 67%
7	O	5	 40% 40% 20%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	L	2	-	-	-	X
3	NAG	R	2	-	-	-	X
3	NAG	T	2	-	-	-	X
4	NAG	H	2	-	-	-	X
4	BMA	H	3	-	-	-	X
4	NAG	M	1	X	-	-	-
4	NAG	M	2	-	-	-	X
5	MAN	S	4	-	-	-	X
7	NAG	O	5	-	-	-	X

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 12482 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 H3 HAEMAGGLUTININ HA1 CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	0	0
			2474	1544	437	477	16			
1	C	319	Total	C	N	O	S	0	0	0
			2475	1544	438	477	16			
1	E	319	Total	C	N	O	S	0	0	0
			2476	1545	438	477	16			

- Molecule 2 is a protein called H3 HAEMAGGLUTININ HA2 CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	175	Total	C	N	O	S	0	0	0
			1425	886	250	283	6			
2	D	175	Total	C	N	O	S	0	0	0
			1419	883	247	283	6			
2	F	175	Total	C	N	O	S	0	0	0
			1406	876	243	281	6			

There are 12 discrepancies between the modelled and reference sequences:

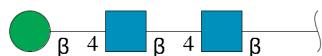
Chain	Residue	Modelled	Actual	Comment	Reference
B	173	SER	-	expression tag	UNP E0UVR5
B	174	GLY	-	expression tag	UNP E0UVR5
B	175	ARG	-	expression tag	UNP E0UVR5
B	131	GLU	ASP	conflict	UNP E0UVR5
D	173	SER	-	expression tag	UNP E0UVR5
D	174	GLY	-	expression tag	UNP E0UVR5
D	175	ARG	-	expression tag	UNP E0UVR5
D	131	GLU	ASP	conflict	UNP E0UVR5
F	173	SER	-	expression tag	UNP E0UVR5
F	174	GLY	-	expression tag	UNP E0UVR5
F	175	ARG	-	expression tag	UNP E0UVR5
F	131	GLU	ASP	conflict	UNP E0UVR5

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



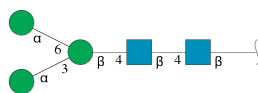
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	L	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	Q	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	R	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	T	2	Total	C	N	O	0	0	0
			28	16	2	10			

- 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	H	3	Total	C	N	O	0	0	0
			39	22	2	15			
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			

- 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	I	5	Total	C	N	O	0	0	0
			61	34	2	25			
5	N	5	Total	C	N	O	0	0	0
			61	34	2	25			
5	S	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 6 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	K	3	Total	C	N	O	0	0	0
			46	25	2	19			
6	P	3	Total	C	N	O	0	0	0
			46	25	2	19			
6	U	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 7 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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.



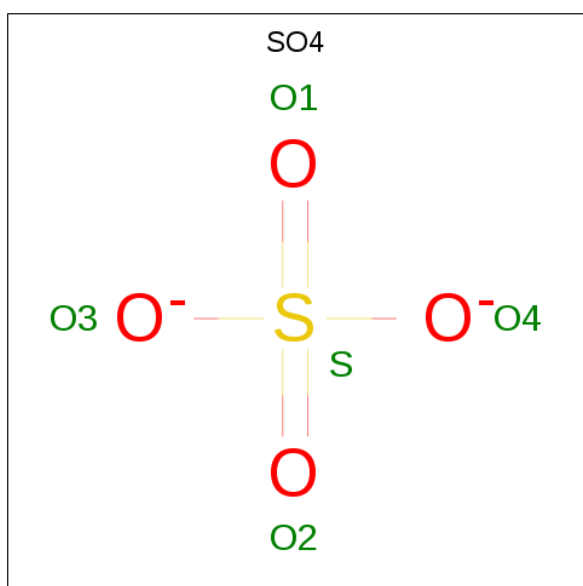
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	O	5	Total	C	N	O	0	0	0
			64	36	3	25			

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 9 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	O	S	0	0
			5	4	1		
9	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	F	1	Total	O	S	0	0
			5	4	1		

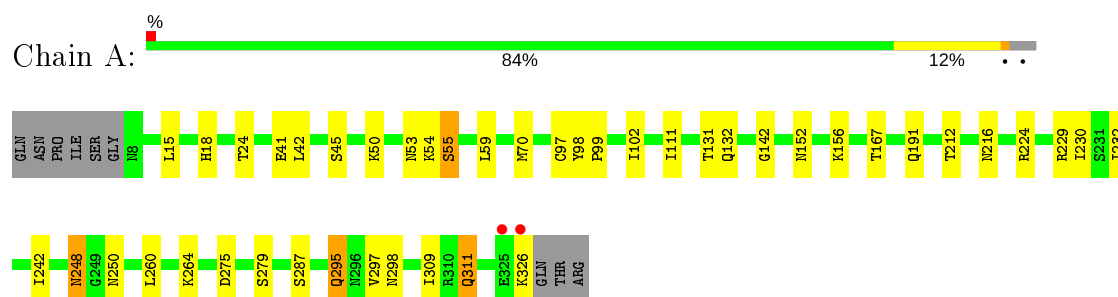
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	21	Total	O	0	0
			21	21		
10	B	13	Total	O	0	0
			13	13		
10	C	36	Total	O	0	0
			36	36		
10	D	11	Total	O	0	0
			11	11		
10	E	35	Total	O	0	0
			35	35		
10	F	6	Total	O	0	0
			6	6		

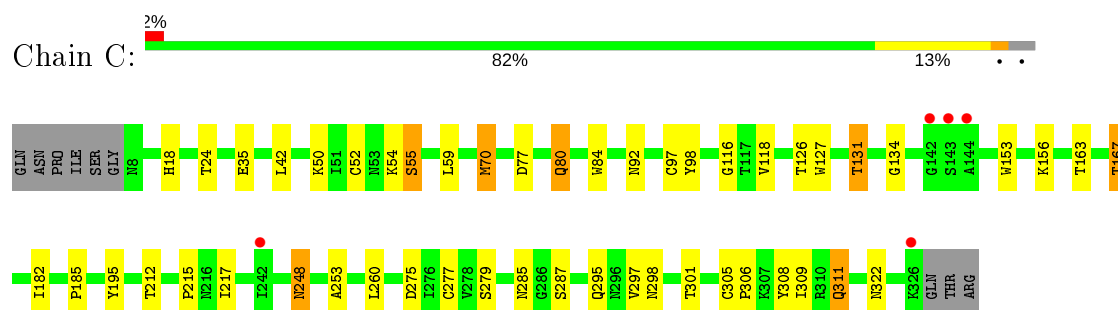
3 Residue-property plots [i](#)

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: H3 HAEMAGGLUTININ HA1 CHAIN



- Molecule 1: H3 HAEMAGGLUTININ HA1 CHAIN




- Molecule 2: H3 HAEMAGGLUTININ HA2 CHAIN

Chain D: 

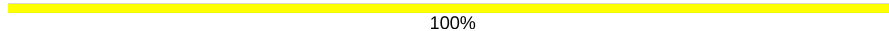


- Molecule 2: H3 HAEMAGGLUTININ HA2 CHAIN

Chain F: 

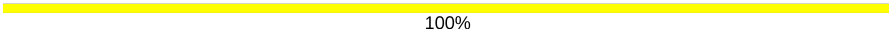


- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G: 

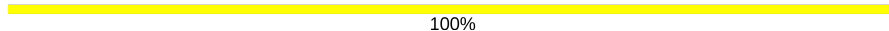


- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L: 



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

Chain Q: 



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

Chain R: 



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

Chain T: 

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 H:  100%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 J:  100%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:  33% 67%MAG1
MAG2
BMA3

- 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 I:  40% 60%MAG1
MAG2
BMA3
MAN4
MAN5

- 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 N:  40% 60%MAG1
MAG2
BMA3
MAN4
MAN5

- 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 S:  20% 80%



- Molecule 6: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  33% 67%



- Molecule 6: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:  100%



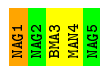
- Molecule 6: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain U:  33% 67%



- Molecule 7: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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 O:  40% 40% 20%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	167.17Å 349.15Å 96.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	174.58 – 2.70 47.80 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.5 (174.58-2.70) 96.5 (47.80-2.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.197 , 0.229 0.197 , 0.229	Depositor DCC
R_{free} test set	3789 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	67.7	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.1	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	12482	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, SIA, GAL, SO4, MAN

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	0.31	0/2525	0.52	0/3425
1	C	0.33	0/2522	0.54	0/3413
1	E	0.32	0/2525	0.53	0/3421
2	B	0.32	0/1450	0.52	1/1949 (0.1%)
2	D	0.31	0/1444	0.53	1/1942 (0.1%)
2	F	0.33	0/1431	0.54	1/1927 (0.1%)
All	All	0.32	0/11897	0.53	3/16077 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	110	LEU	CA-CB-CG	7.36	132.23	115.30
2	D	110	LEU	CA-CB-CG	6.86	131.08	115.30
2	B	110	LEU	CA-CB-CG	6.27	129.72	115.30

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	2474	0	2424	21	0
1	C	2475	0	2422	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2476	0	2430	27	0
2	B	1425	0	1344	15	0
2	D	1419	0	1333	16	0
2	F	1406	0	1309	22	0
3	G	28	0	25	0	0
3	L	28	0	25	0	0
3	Q	28	0	25	0	0
3	R	28	0	25	0	0
3	T	28	0	25	0	0
4	H	39	0	34	0	0
4	J	39	0	34	0	0
4	M	39	0	34	0	0
5	I	61	0	52	3	0
5	N	61	0	52	0	0
5	S	61	0	52	0	0
6	K	46	0	40	0	0
6	P	46	0	40	0	0
6	U	46	0	40	0	0
7	O	64	0	55	2	0
8	A	14	0	13	0	0
8	C	14	0	13	1	0
9	B	5	0	0	0	0
9	D	5	0	0	0	0
9	F	5	0	0	0	0
10	A	21	0	0	0	0
10	B	13	0	0	1	0
10	C	36	0	0	1	0
10	D	11	0	0	0	0
10	E	35	0	0	0	0
10	F	6	0	0	0	0
All	All	12482	0	11846	117	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

The worst 5 of 117 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)
5:I:3:BMA:H5	5:I:5:MAN:H2	1.58	0.84
1:C:285:ASN:HD21	7:O:1:NAG:C1	1.93	0.81
2:B:134:GLY:HA2	2:D:124:ARG:HD3	1.65	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:25:ARG:HE	2:B:34:GLN:HE21	1.33	0.75
1:C:295:GLN:NE2	1:C:298:ASN:H	1.87	0.72

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	317/328 (97%)	301 (95%)	14 (4%)	2 (1%)	25	50
1	C	315/328 (96%)	301 (96%)	12 (4%)	2 (1%)	25	50
1	E	316/328 (96%)	300 (95%)	13 (4%)	3 (1%)	17	40
2	B	173/175 (99%)	164 (95%)	9 (5%)	0	100	100
2	D	173/175 (99%)	163 (94%)	10 (6%)	0	100	100
2	F	173/175 (99%)	163 (94%)	10 (6%)	0	100	100
All	All	1467/1509 (97%)	1392 (95%)	68 (5%)	7 (0%)	29	54

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	55	SER
1	E	196	ILE
1	A	55	SER
1	C	55	SER
1	E	158	GLY

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	280/289 (97%)	263 (94%)	17 (6%)	18	41
1	C	279/289 (96%)	265 (95%)	14 (5%)	24	51
1	E	280/289 (97%)	260 (93%)	20 (7%)	14	34
2	B	148/148 (100%)	143 (97%)	5 (3%)	37	66
2	D	147/148 (99%)	141 (96%)	6 (4%)	30	59
2	F	144/148 (97%)	136 (94%)	8 (6%)	21	45
All	All	1278/1311 (98%)	1208 (94%)	70 (6%)	21	46

5 of 70 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	167	THR
2	D	110	LEU
2	F	86	ASP
1	C	195	TYR
1	C	311	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	250	ASN
2	D	27	GLN
2	F	116	ASN
1	C	311	GLN
2	D	34	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

48 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	G	1	1,3	14,14,15	0.51	0	17,19,21	0.95	1 (5%)
3	NAG	G	2	3	14,14,15	0.63	0	17,19,21	1.39	3 (17%)
4	NAG	H	1	1,4	14,14,15	0.58	0	17,19,21	0.86	1 (5%)
4	NAG	H	2	4	14,14,15	0.66	0	17,19,21	1.24	2 (11%)
4	BMA	H	3	4	11,11,12	0.47	0	15,15,17	0.94	1 (6%)
5	NAG	I	1	1,5	14,14,15	0.59	0	17,19,21	1.10	1 (5%)
5	NAG	I	2	5	14,14,15	0.53	0	17,19,21	1.32	2 (11%)
5	BMA	I	3	5	11,11,12	0.74	0	15,15,17	1.89	5 (33%)
5	MAN	I	4	5	11,11,12	0.62	0	15,15,17	1.74	3 (20%)
5	MAN	I	5	5	11,11,12	0.56	0	15,15,17	1.15	1 (6%)
4	NAG	J	1	4	14,14,15	0.57	0	17,19,21	1.70	3 (17%)
4	NAG	J	2	4	14,14,15	0.61	0	17,19,21	1.29	1 (5%)
4	BMA	J	3	4	11,11,12	0.47	0	15,15,17	1.50	2 (13%)
6	NAG	K	1	6	15,15,15	0.53	0	21,21,21	1.57	3 (14%)
6	GAL	K	2	6	11,11,12	0.60	0	15,15,17	0.92	0
6	SIA	K	3	6	17,20,21	0.39	0	21,28,31	0.91	1 (4%)
3	NAG	L	1	1,3	14,14,15	0.66	0	17,19,21	1.21	2 (11%)
3	NAG	L	2	3	14,14,15	0.60	0	17,19,21	1.08	2 (11%)
4	NAG	M	1	1,4	14,14,15	0.44	0	17,19,21	1.39	3 (17%)
4	NAG	M	2	4	14,14,15	0.58	0	17,19,21	1.29	2 (11%)
4	BMA	M	3	4	11,11,12	0.44	0	15,15,17	0.76	0
5	NAG	N	1	1,5	14,14,15	0.60	0	17,19,21	1.17	3 (17%)
5	NAG	N	2	5	14,14,15	0.59	0	17,19,21	0.67	0
5	BMA	N	3	5	11,11,12	0.43	0	15,15,17	1.39	1 (6%)
5	MAN	N	4	5	11,11,12	0.60	0	15,15,17	1.34	2 (13%)
5	MAN	N	5	5	11,11,12	0.66	0	15,15,17	0.86	0
7	NAG	O	1	7	14,14,15	0.63	0	17,19,21	1.56	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NAG	O	2	7	14,14,15	0.52	0	17,19,21	1.06	0
7	BMA	O	3	7	11,11,12	0.48	0	15,15,17	1.19	2 (13%)
7	MAN	O	4	7	11,11,12	0.58	0	15,15,17	0.89	1 (6%)
7	NAG	O	5	7	14,14,15	0.48	0	17,19,21	0.92	0
6	NAG	P	1	6	15,15,15	0.53	0	21,21,21	1.17	1 (4%)
6	GAL	P	2	6	11,11,12	0.64	0	15,15,17	1.30	2 (13%)
6	SIA	P	3	6	17,20,21	0.45	0	21,28,31	0.82	1 (4%)
3	NAG	Q	1	1,3	14,14,15	0.58	0	17,19,21	0.90	1 (5%)
3	NAG	Q	2	3	14,14,15	0.46	0	17,19,21	1.01	1 (5%)
3	NAG	R	1	1,3	14,14,15	0.56	0	17,19,21	0.98	0
3	NAG	R	2	3	14,14,15	0.65	0	17,19,21	1.22	1 (5%)
5	NAG	S	1	1,5	14,14,15	0.60	0	17,19,21	1.08	1 (5%)
5	NAG	S	2	5	14,14,15	0.50	0	17,19,21	0.93	1 (5%)
5	BMA	S	3	5	11,11,12	0.48	0	15,15,17	2.19	2 (13%)
5	MAN	S	4	5	11,11,12	0.71	0	15,15,17	1.61	3 (20%)
5	MAN	S	5	5	11,11,12	0.68	0	15,15,17	0.89	0
3	NAG	T	1	3	14,14,15	0.54	0	17,19,21	1.12	2 (11%)
3	NAG	T	2	3	14,14,15	0.57	0	17,19,21	1.24	2 (11%)
6	NAG	U	1	6	15,15,15	0.49	0	21,21,21	1.03	2 (9%)
6	GAL	U	2	6	11,11,12	0.68	0	15,15,17	0.88	0
6	SIA	U	3	6	17,20,21	0.44	0	21,28,31	0.92	1 (4%)

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	G	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
4	NAG	H	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1
4	BMA	H	3	4	-	2/2/19/22	0/1/1/1
5	NAG	I	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	I	2	5	-	2/6/23/26	0/1/1/1
5	BMA	I	3	5	-	2/2/19/22	1/1/1/1
5	MAN	I	4	5	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	MAN	I	5	5	-	0/2/19/22	1/1/1/1
4	NAG	J	1	4	-	2/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1
4	BMA	J	3	4	-	0/2/19/22	0/1/1/1
6	NAG	K	1	6	-	0/6/26/26	0/1/1/1
6	GAL	K	2	6	-	2/2/19/22	0/1/1/1
6	SIA	K	3	6	-	0/14/34/38	0/1/1/1
3	NAG	L	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	L	2	3	-	2/6/23/26	0/1/1/1
4	NAG	M	1	1,4	1/1/5/7	0/6/23/26	0/1/1/1
4	NAG	M	2	4	-	4/6/23/26	0/1/1/1
4	BMA	M	3	4	-	0/2/19/22	0/1/1/1
5	NAG	N	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	N	2	5	-	0/6/23/26	0/1/1/1
5	BMA	N	3	5	-	0/2/19/22	0/1/1/1
5	MAN	N	4	5	-	2/2/19/22	0/1/1/1
5	MAN	N	5	5	-	0/2/19/22	0/1/1/1
7	NAG	O	1	7	-	0/6/23/26	0/1/1/1
7	NAG	O	2	7	-	2/6/23/26	0/1/1/1
7	BMA	O	3	7	-	2/2/19/22	0/1/1/1
7	MAN	O	4	7	-	1/2/19/22	0/1/1/1
7	NAG	O	5	7	-	2/6/23/26	0/1/1/1
6	NAG	P	1	6	-	0/6/26/26	0/1/1/1
6	GAL	P	2	6	-	2/2/19/22	0/1/1/1
6	SIA	P	3	6	-	0/14/34/38	0/1/1/1
3	NAG	Q	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	Q	2	3	-	0/6/23/26	0/1/1/1
3	NAG	R	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	R	2	3	-	2/6/23/26	0/1/1/1
5	NAG	S	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	S	2	5	-	2/6/23/26	0/1/1/1
5	BMA	S	3	5	-	2/2/19/22	0/1/1/1
5	MAN	S	4	5	-	0/2/19/22	0/1/1/1
5	MAN	S	5	5	-	2/2/19/22	0/1/1/1
3	NAG	T	1	3	-	0/6/23/26	0/1/1/1
3	NAG	T	2	3	-	2/6/23/26	0/1/1/1
6	NAG	U	1	6	-	2/6/26/26	0/1/1/1
6	GAL	U	2	6	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	SIA	U	3	6	-	1/14/34/38	0/1/1/1

There are no bond length outliers.

The worst 5 of 72 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	S	3	BMA	C1-O5-C5	7.46	122.30	112.19
4	J	1	NAG	C3-C4-C5	4.53	118.31	110.24
5	I	4	MAN	C1-O5-C5	4.50	118.30	112.19
5	S	4	MAN	C1-C2-C3	4.30	114.95	109.67
5	I	4	MAN	C1-C2-C3	4.05	114.64	109.67

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	M	1	NAG	C1

5 of 58 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	U	1	NAG	C1-C2-N2-C7
5	S	5	MAN	O5-C5-C6-O6
3	T	2	NAG	O5-C5-C6-O6
3	R	2	NAG	C4-C5-C6-O6
7	O	2	NAG	O5-C5-C6-O6

All (2) ring outliers are listed below:

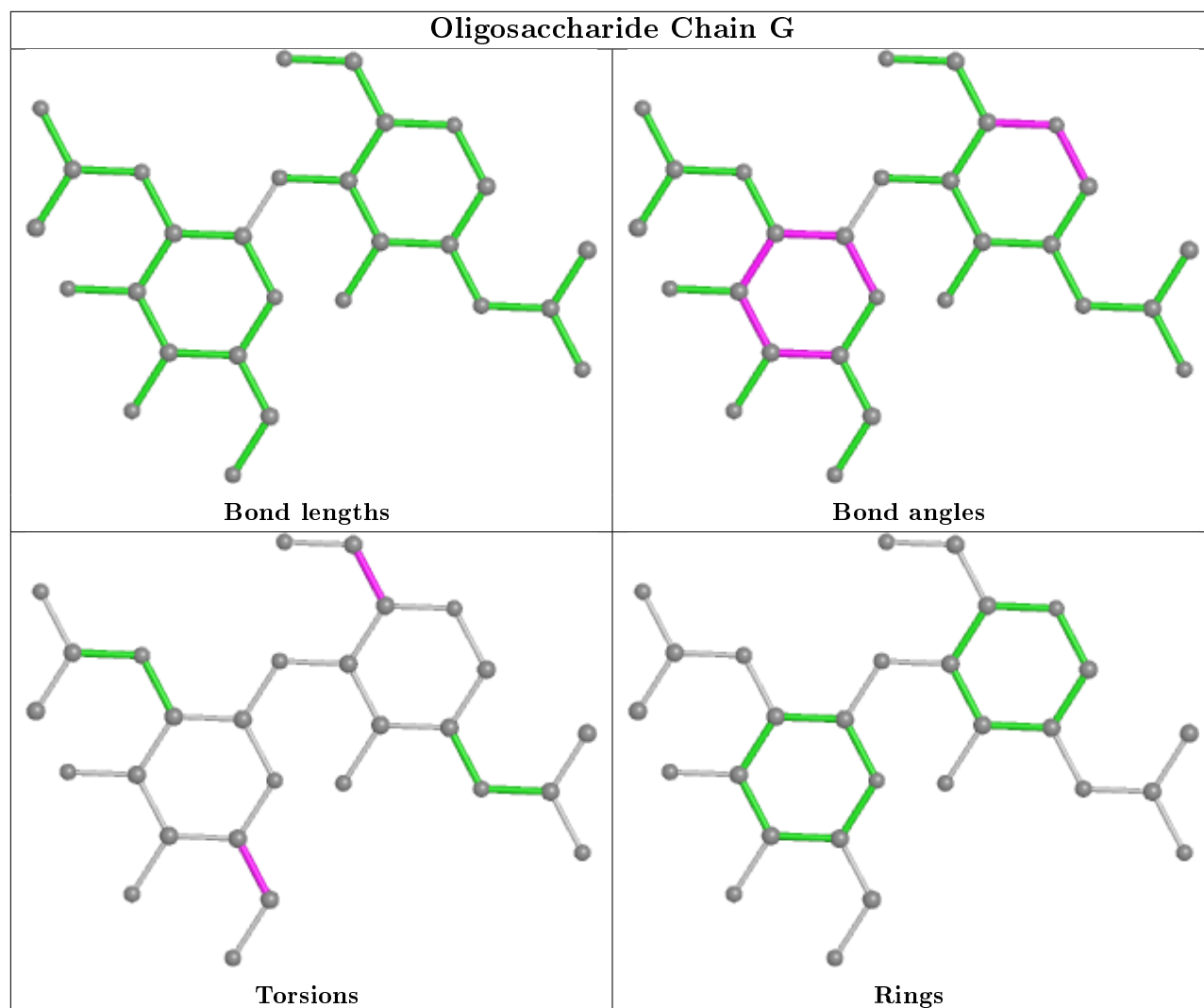
Mol	Chain	Res	Type	Atoms
5	I	3	BMA	C1-C2-C3-C4-C5-O5
5	I	5	MAN	C1-C2-C3-C4-C5-O5

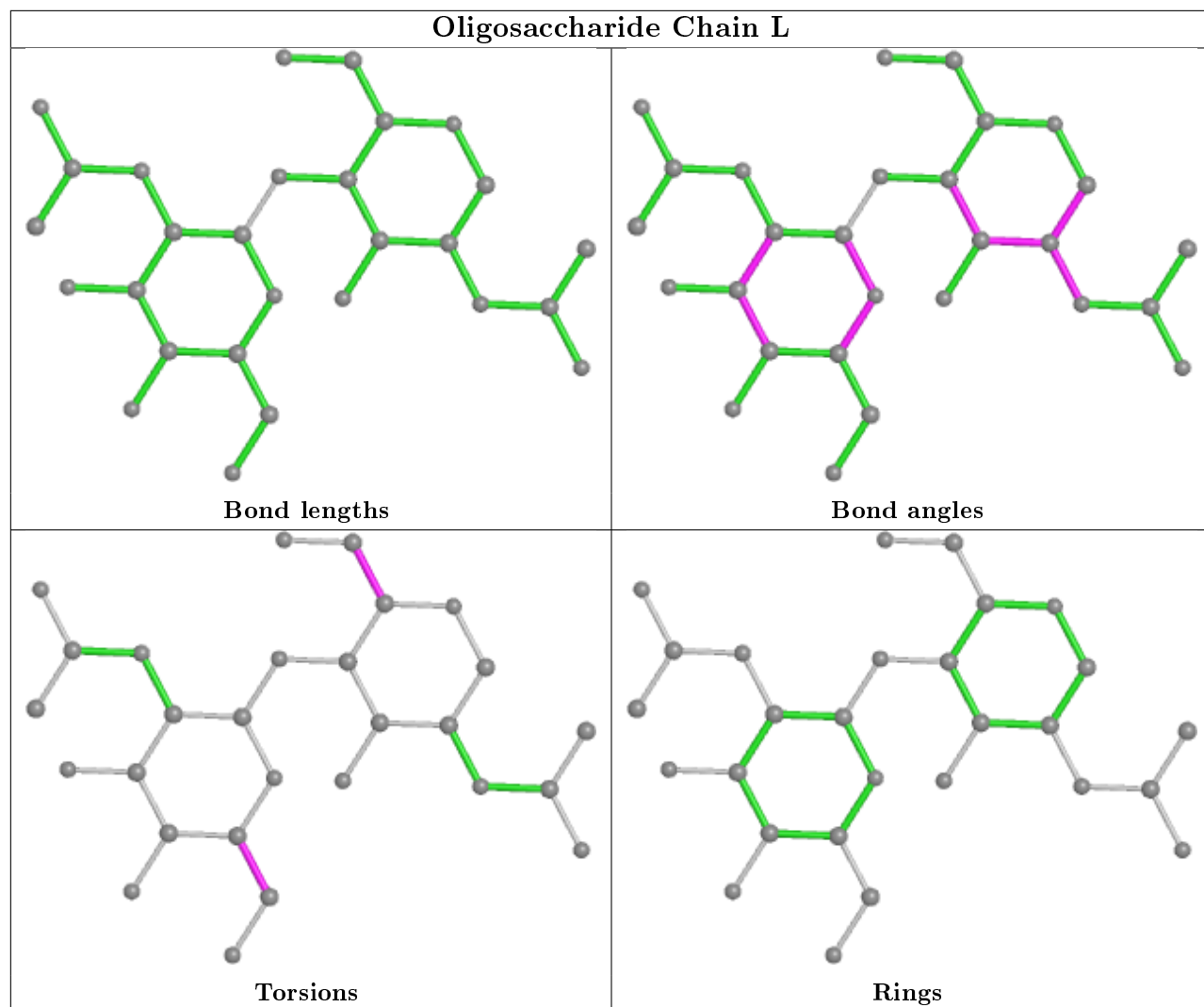
4 monomers are involved in 5 short contacts:

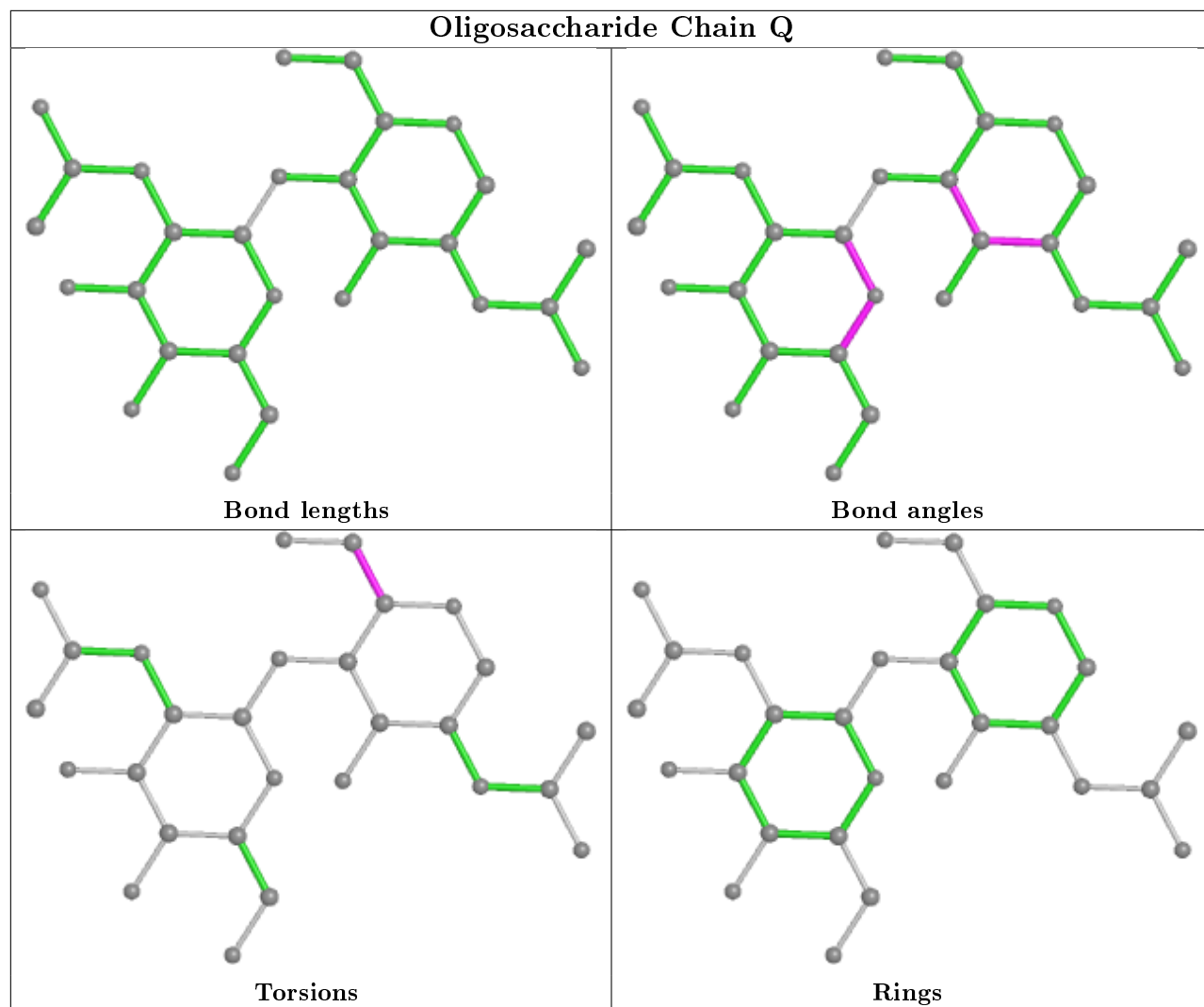
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	I	5	MAN	1	0
5	I	3	BMA	3	0
5	I	4	MAN	1	0
7	O	1	NAG	2	0

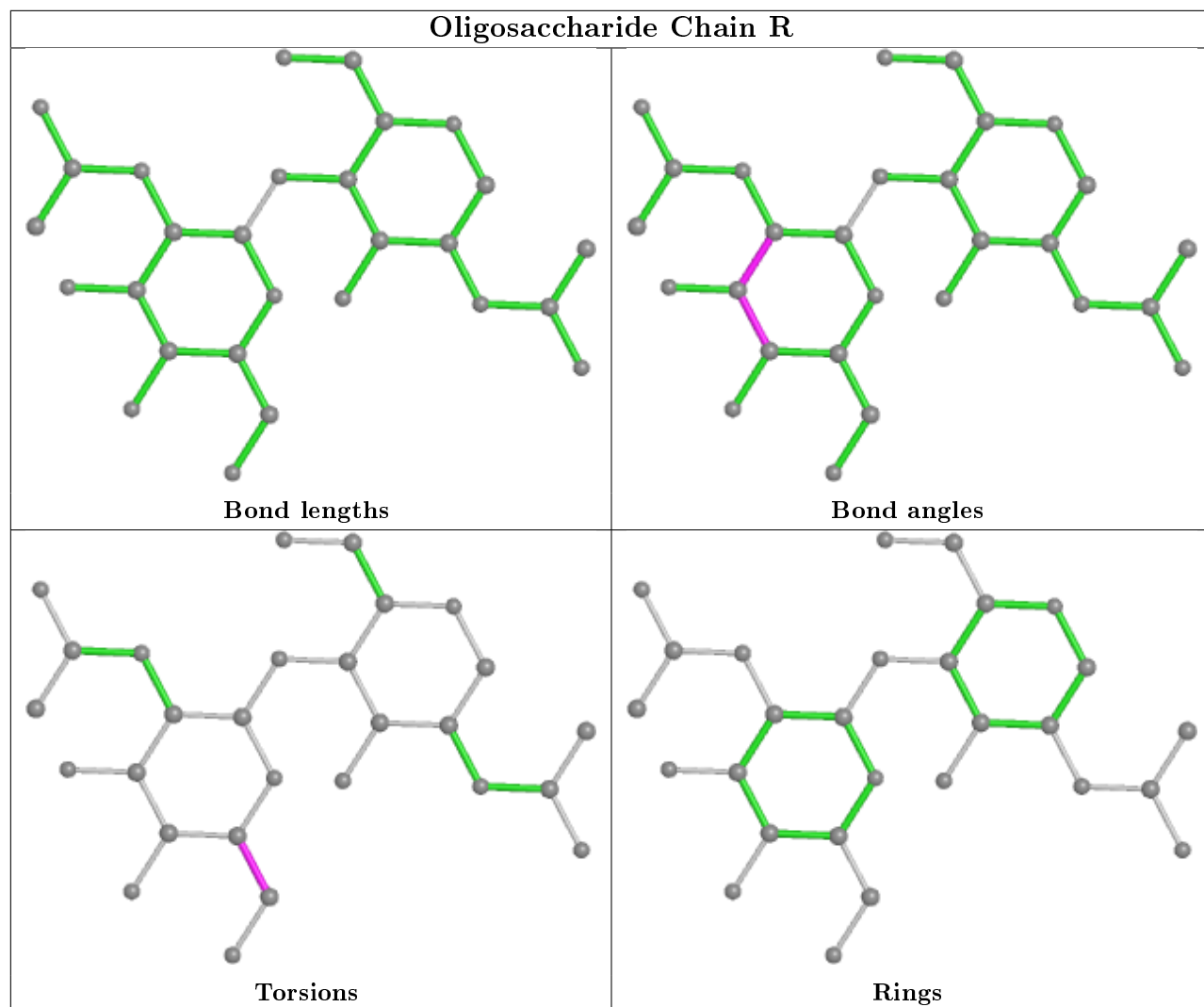
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

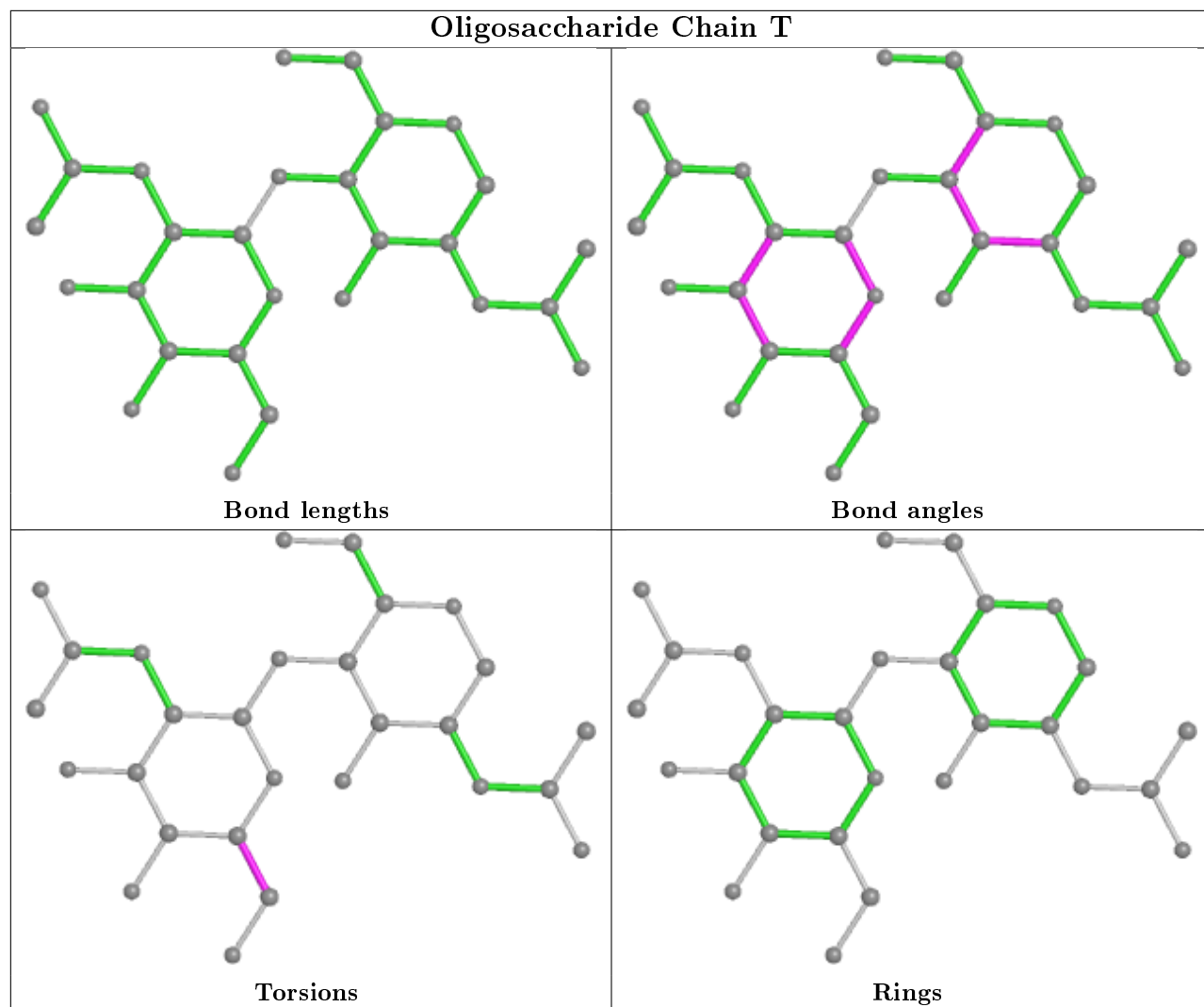
bond angles, torsion angles, and ring geometry for oligosaccharide.

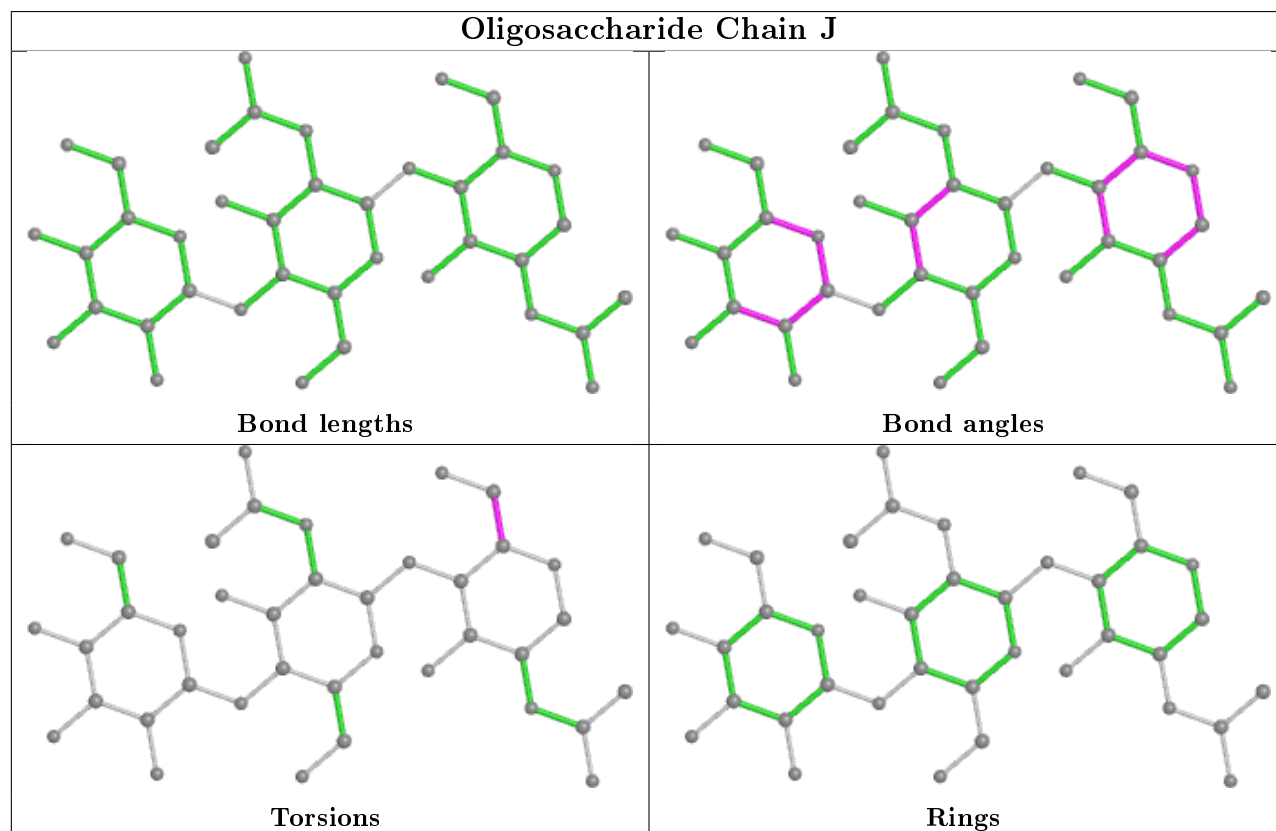
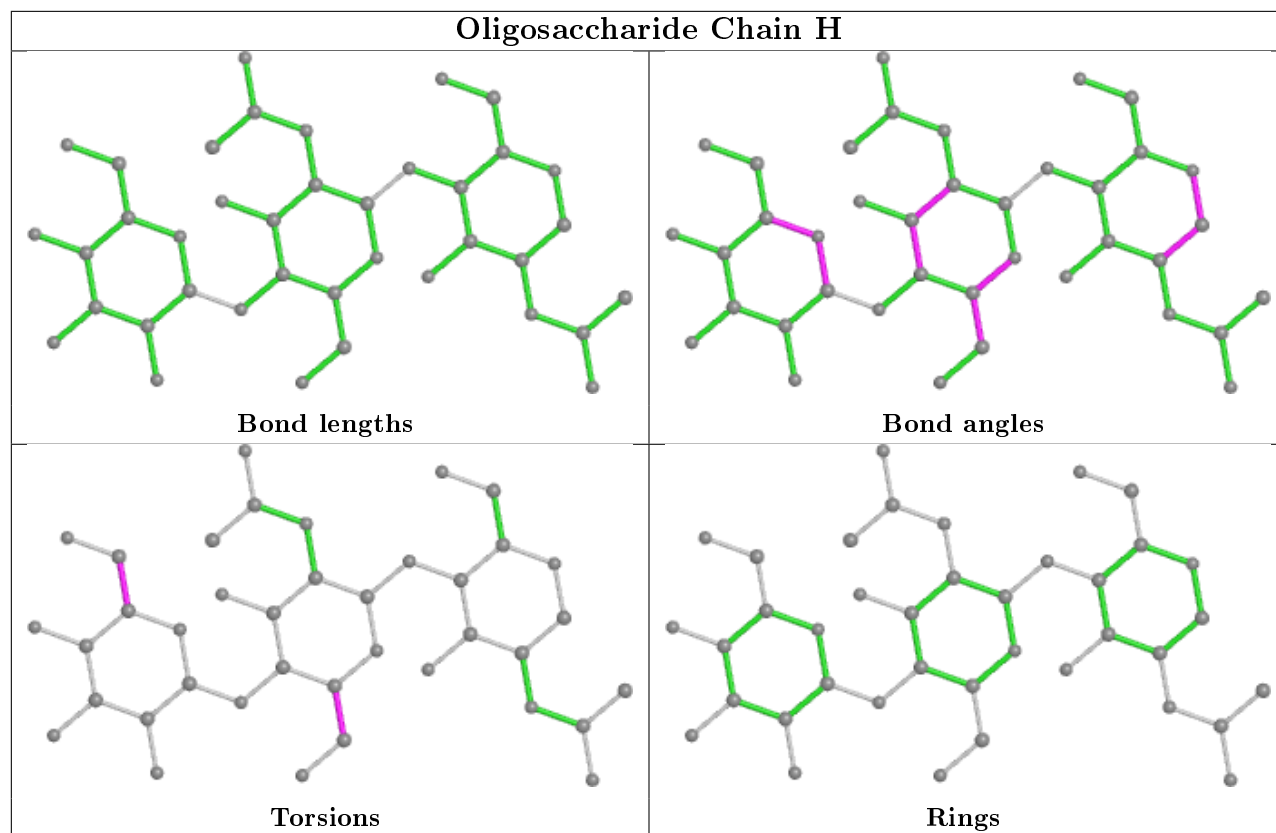


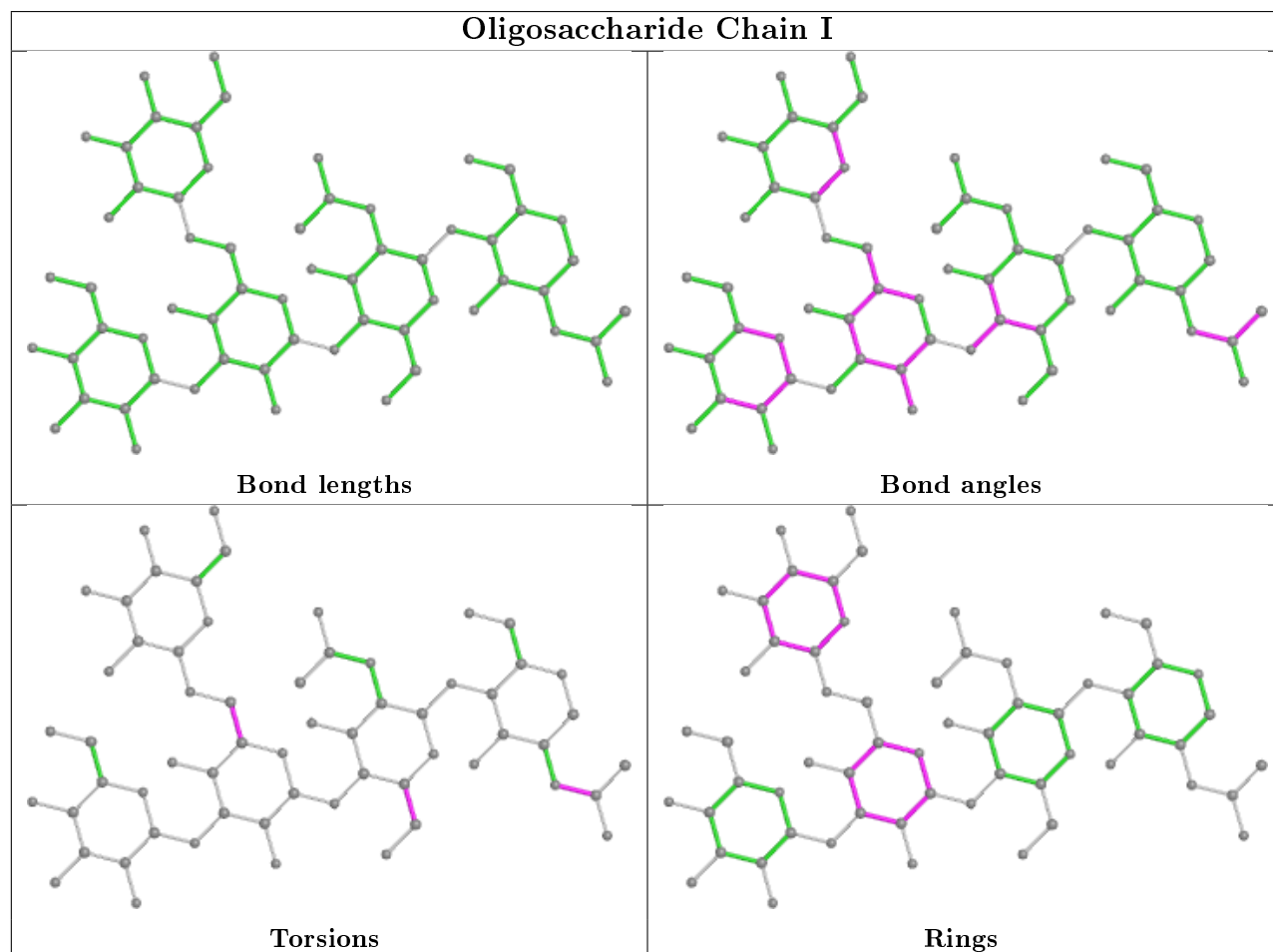
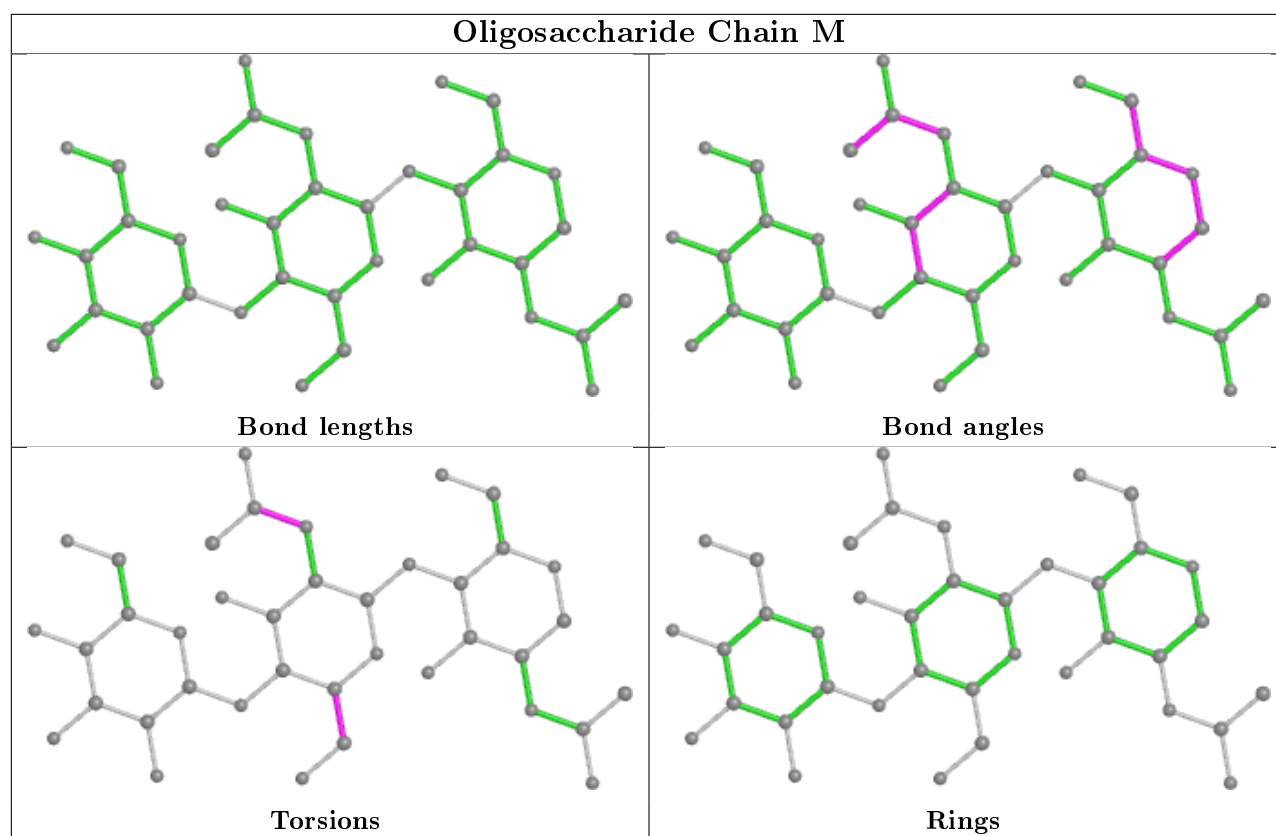


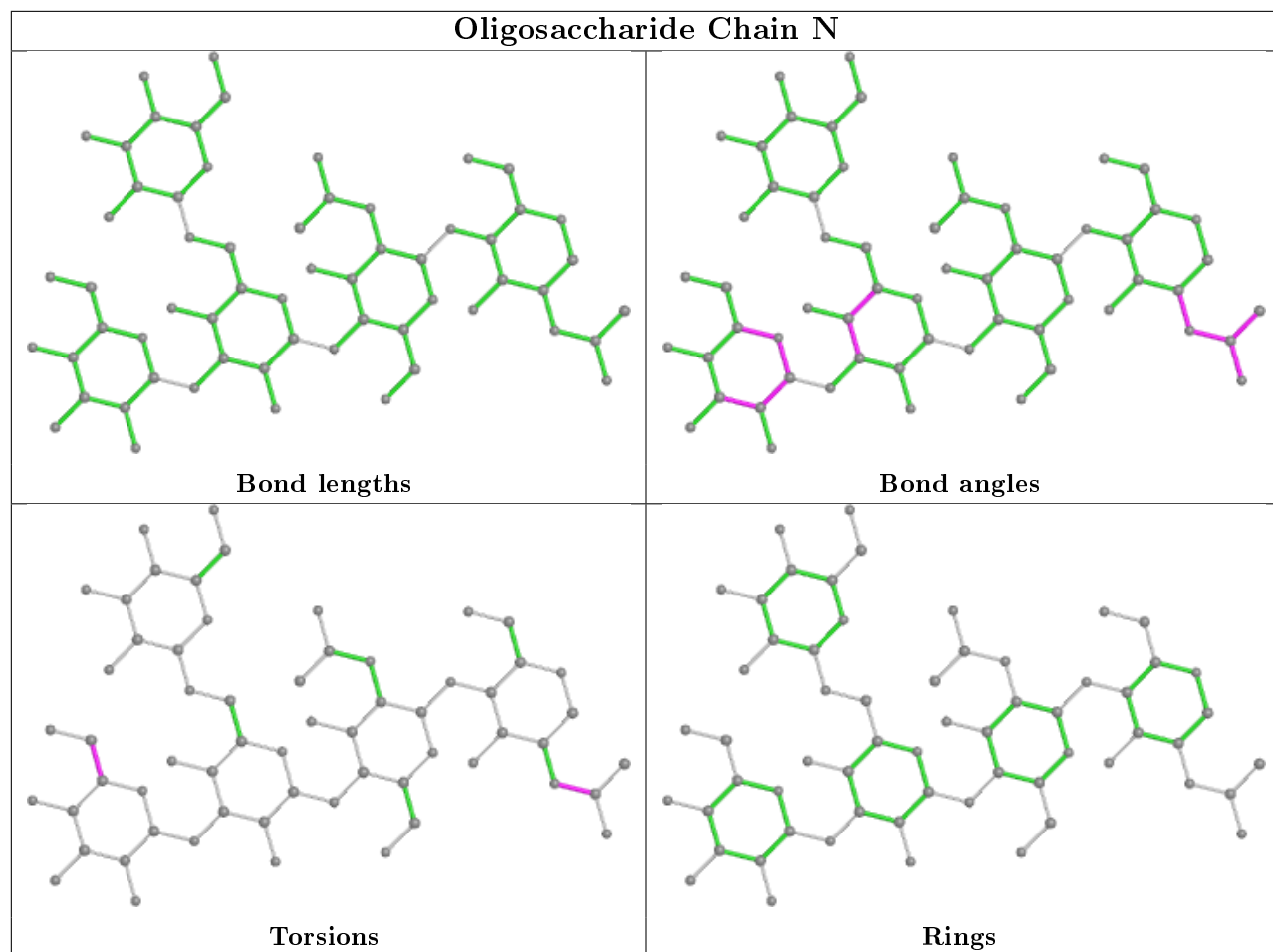


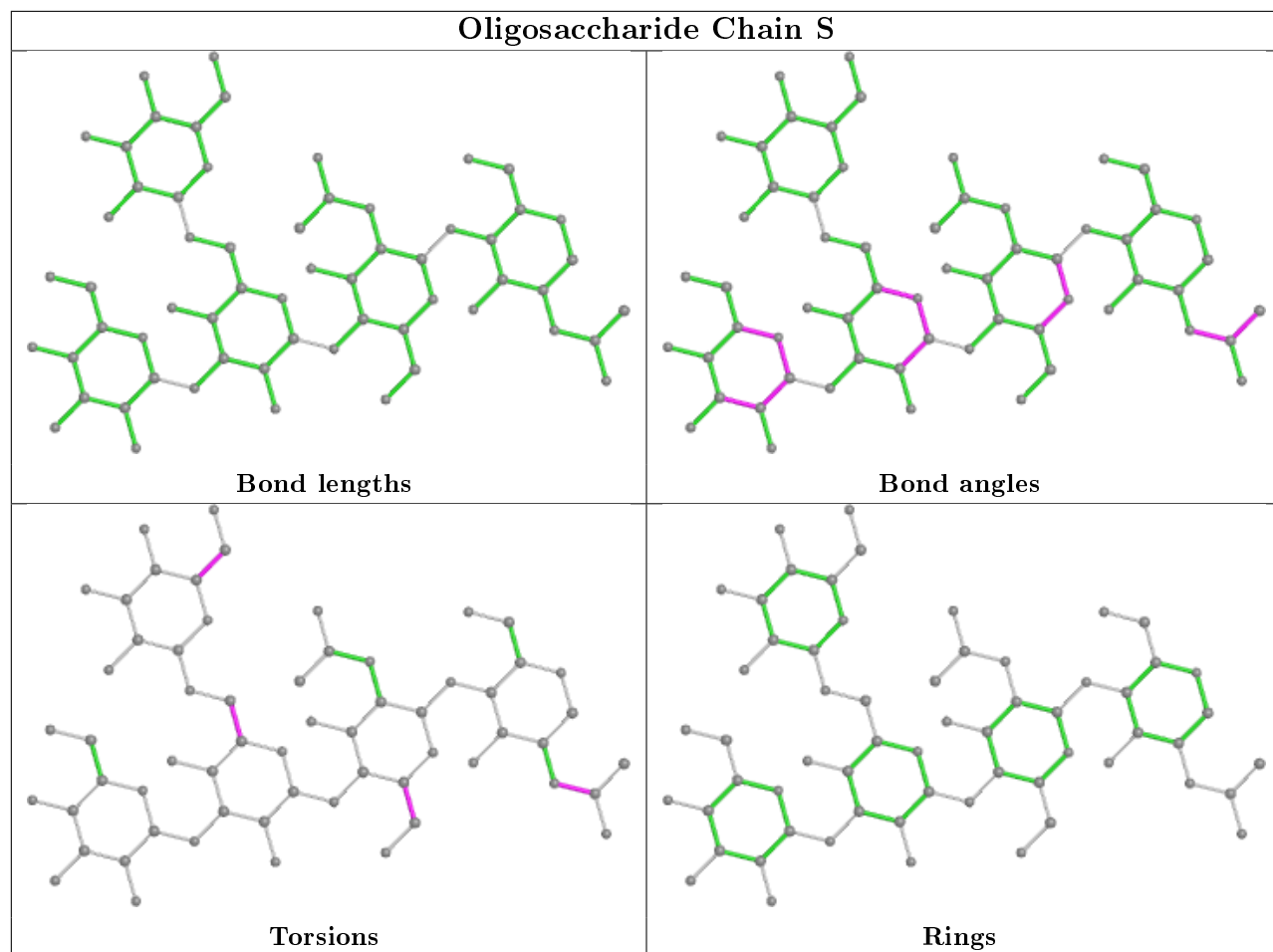


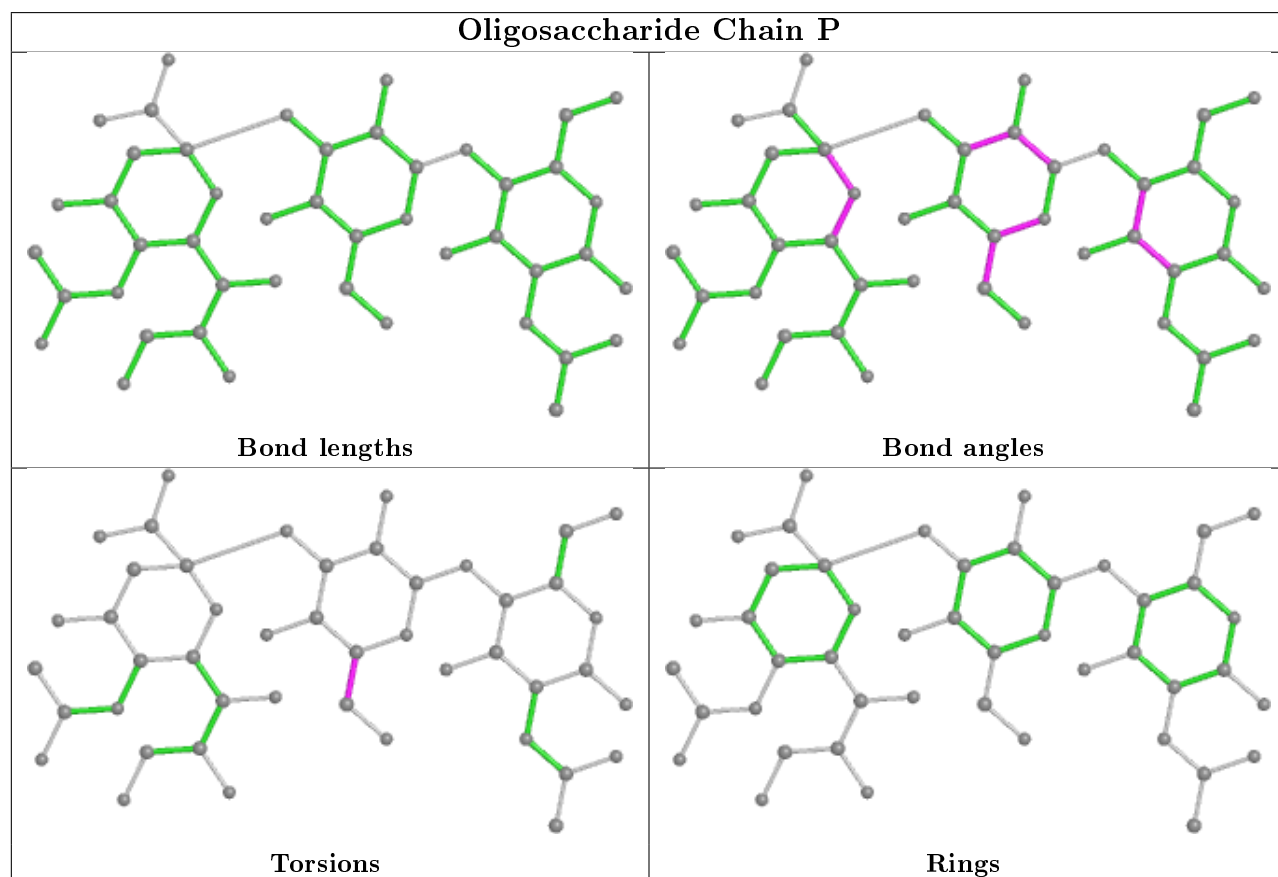
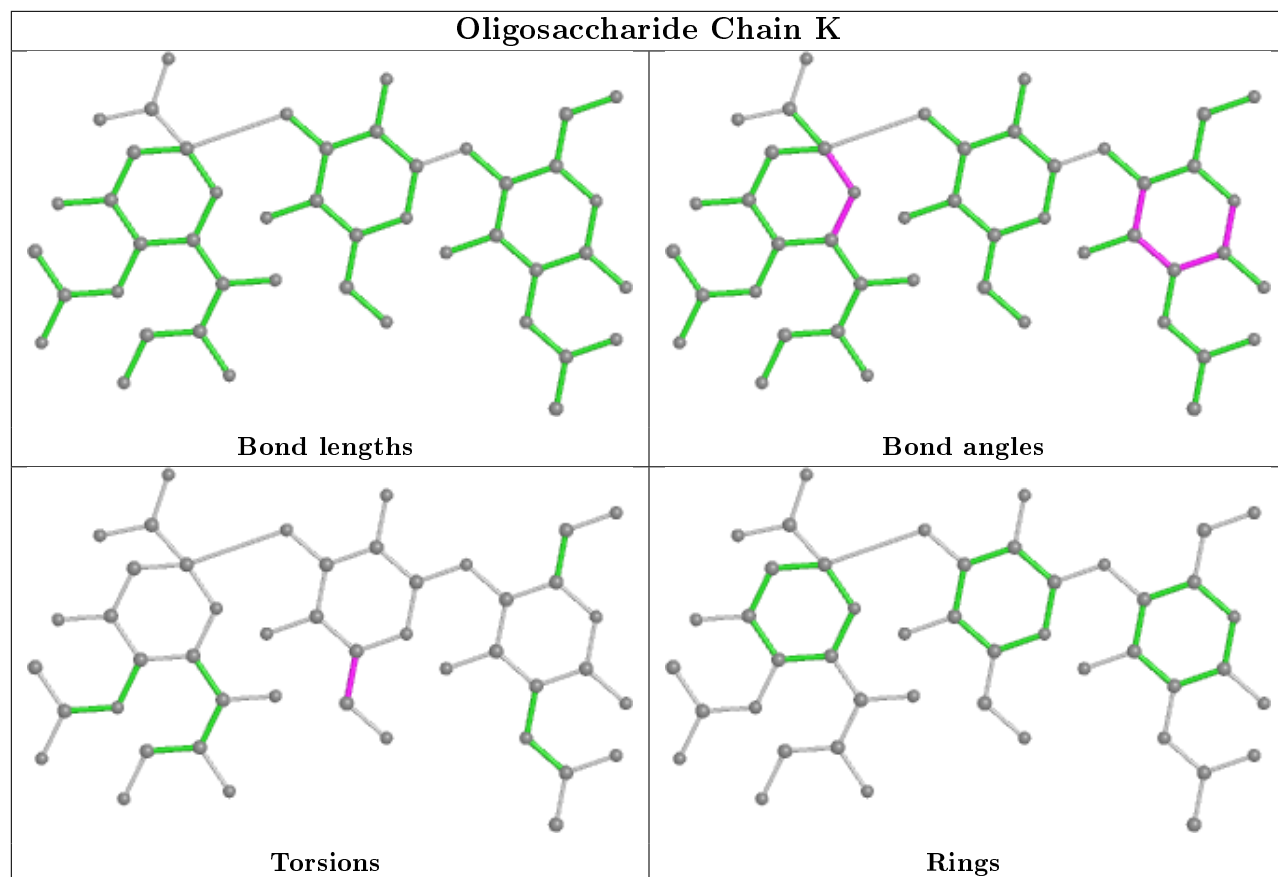


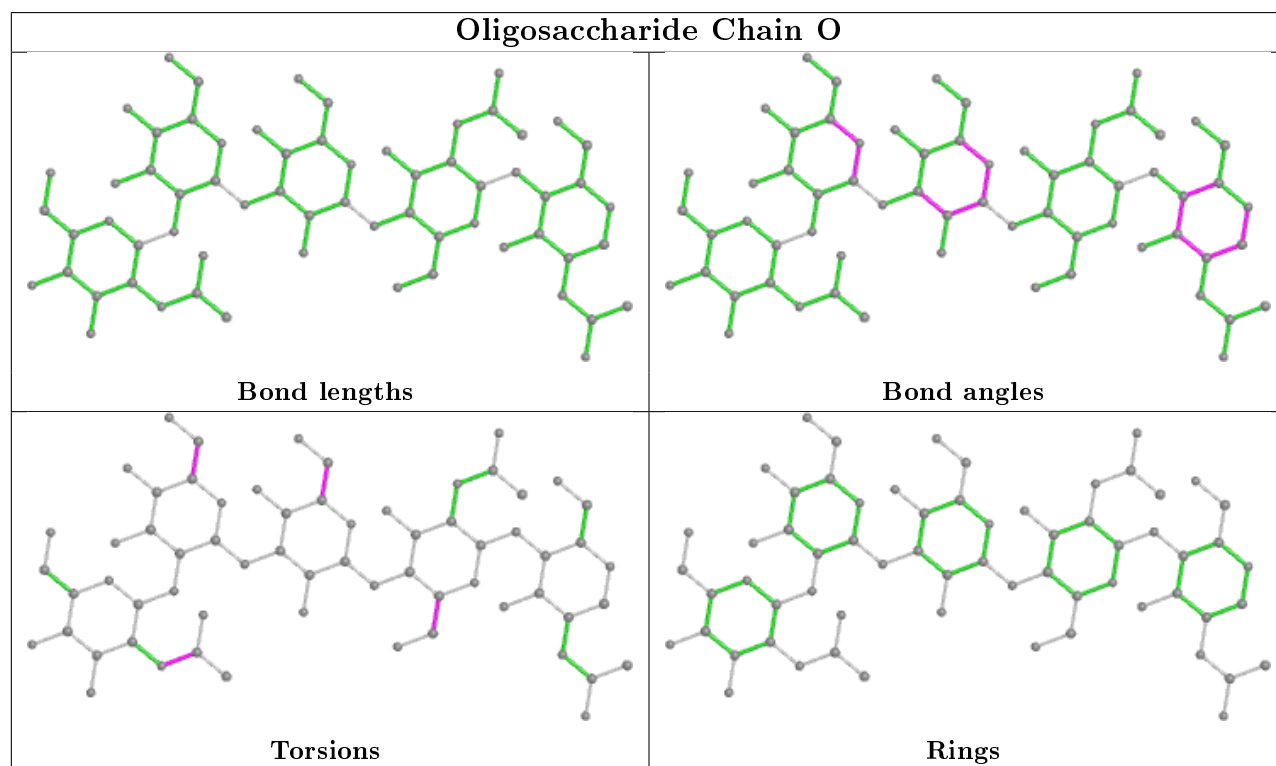
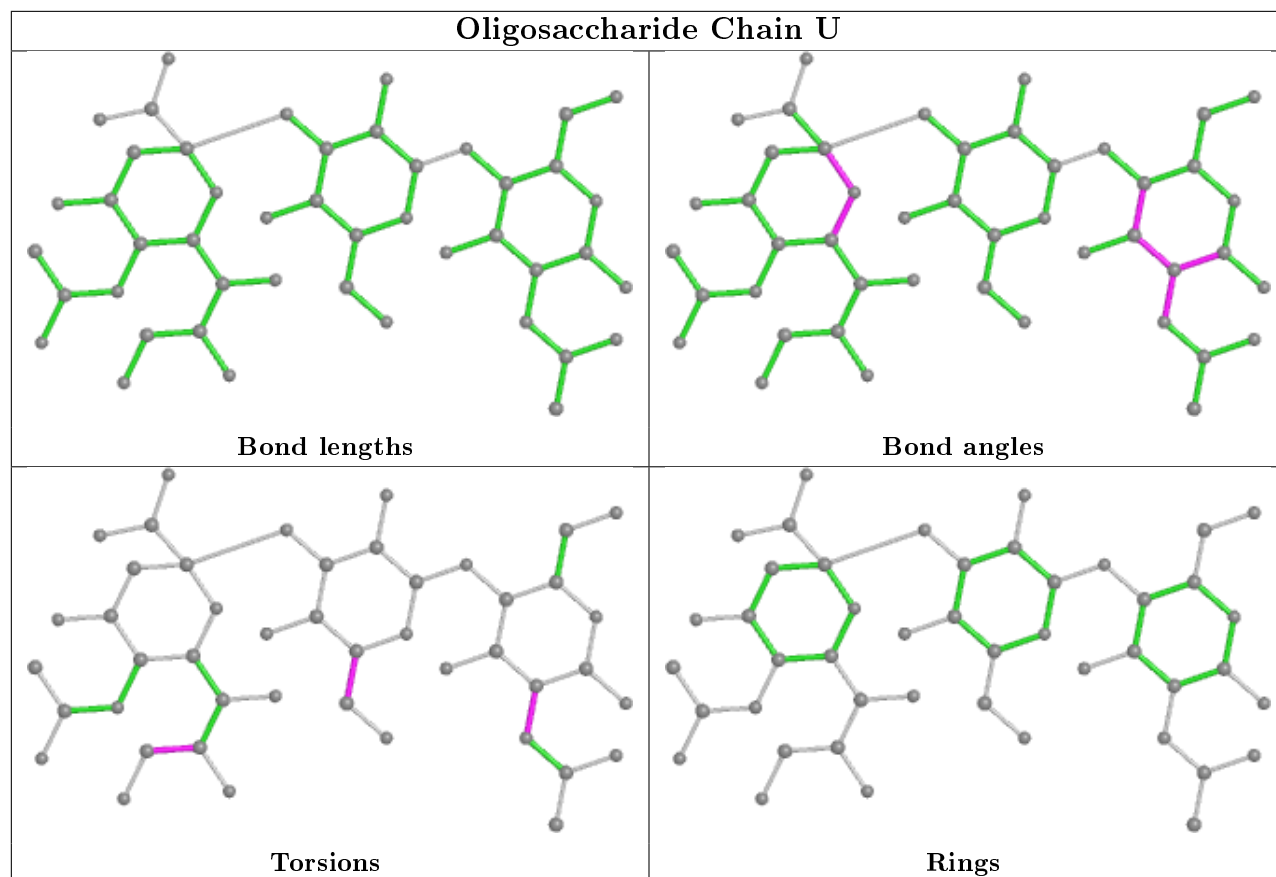












5.6 Ligand geometry

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	$\# Z > 2$
8	NAG	C	601	1	14,14,15	0.64	0	17,19,21	2.94	5 (29%)
9	SO4	F	1176	-	4,4,4	0.34	0	6,6,6	0.08	0
9	SO4	B	1176	-	4,4,4	0.35	0	6,6,6	0.08	0
9	SO4	D	1176	-	4,4,4	0.35	0	6,6,6	0.14	0
8	NAG	A	601	1	14,14,15	0.61	0	17,19,21	1.58	4 (23%)

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
8	NAG	C	601	1	-	5/6/23/26	0/1/1/1
8	NAG	A	601	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	601	NAG	C1-O5-C5	8.20	123.30	112.19
8	C	601	NAG	C2-N2-C7	6.19	131.71	122.90
8	C	601	NAG	O5-C1-C2	4.13	117.81	111.29
8	A	601	NAG	C1-O5-C5	4.02	117.64	112.19
8	C	601	NAG	C8-C7-N2	3.41	121.87	116.10

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	C	601	NAG	C8-C7-N2-C2
8	C	601	NAG	O7-C7-N2-C2
8	A	601	NAG	C8-C7-N2-C2
8	A	601	NAG	O7-C7-N2-C2
8	C	601	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	C	601	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 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	319/328 (97%)	-0.00	2 (0%) 89 91	46, 66, 93, 130	0
1	C	319/328 (97%)	-0.04	5 (1%) 72 74	45, 64, 96, 132	0
1	E	319/328 (97%)	0.02	5 (1%) 72 74	46, 64, 95, 127	0
2	B	175/175 (100%)	0.22	4 (2%) 60 62	44, 75, 97, 152	0
2	D	175/175 (100%)	0.15	3 (1%) 70 72	43, 73, 98, 137	0
2	F	175/175 (100%)	0.10	3 (1%) 70 72	42, 71, 95, 128	0
All	All	1482/1509 (98%)	0.05	22 (1%) 73 76	42, 68, 96, 152	0

The worst 5 of 22 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	174	GLY	8.9
2	B	175	ARG	8.1
2	D	173	SER	6.8
2	D	175	ARG	5.1
2	D	174	GLY	4.9

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.

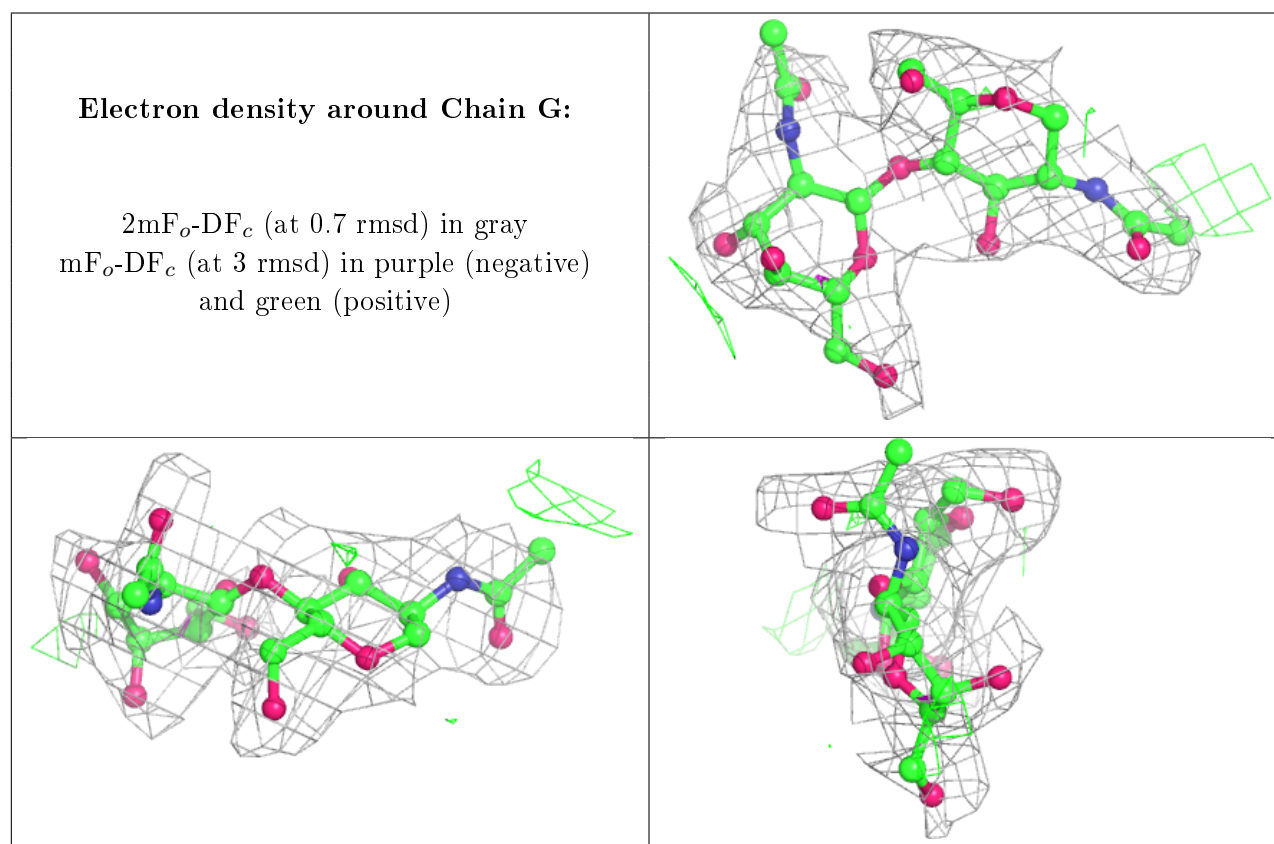
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	T	2	14/15	0.52	0.41	125,138,144,146	0
7	NAG	O	5	14/15	0.66	0.42	130,150,156,158	0
4	NAG	M	2	14/15	0.67	0.42	141,148,155,156	0
4	BMA	M	3	11/12	0.69	0.36	149,158,160,160	0
5	MAN	S	5	11/12	0.70	0.27	142,148,153,154	0
4	BMA	H	3	11/12	0.72	0.44	147,153,155,155	0
5	MAN	I	5	11/12	0.74	0.33	154,157,160,161	0
4	NAG	H	2	14/15	0.74	0.52	135,148,153,154	0
5	MAN	S	4	11/12	0.75	0.49	145,150,157,160	0
5	MAN	N	5	11/12	0.75	0.28	135,141,150,152	0
5	MAN	I	4	11/12	0.76	0.30	146,150,154,156	0
5	BMA	I	3	11/12	0.76	0.29	124,139,146,155	0
3	NAG	L	2	14/15	0.78	0.50	120,124,134,136	0
3	NAG	R	2	14/15	0.78	0.45	130,135,138,139	0
5	BMA	S	3	11/12	0.79	0.27	120,131,141,142	0
3	NAG	T	1	14/15	0.79	0.33	109,125,128,131	0
5	MAN	N	4	11/12	0.79	0.35	132,135,141,141	0
4	BMA	J	3	11/12	0.80	0.33	137,143,146,146	0
6	NAG	U	1	15/15	0.81	0.35	102,110,121,123	0
7	NAG	O	2	14/15	0.81	0.39	127,134,141,143	0
3	NAG	Q	2	14/15	0.83	0.42	119,125,129,132	0
7	BMA	O	3	11/12	0.83	0.32	136,143,147,147	0
3	NAG	G	2	14/15	0.84	0.25	116,122,128,129	0
7	MAN	O	4	11/12	0.84	0.44	144,147,149,149	0
4	NAG	M	1	14/15	0.84	0.29	101,108,119,132	0
4	NAG	J	1	14/15	0.84	0.20	103,107,114,116	0
7	NAG	O	1	14/15	0.84	0.31	111,124,129,132	0
6	NAG	P	1	15/15	0.85	0.52	111,120,135,136	0
5	BMA	N	3	11/12	0.85	0.22	115,125,132,137	0
4	NAG	H	1	14/15	0.85	0.35	104,113,122,133	0
3	NAG	L	1	14/15	0.87	0.24	91,98,106,115	0
3	NAG	R	1	14/15	0.88	0.30	99,108,113,122	0
6	NAG	K	1	15/15	0.88	0.36	100,107,118,118	0
4	NAG	J	2	14/15	0.88	0.32	116,124,132,138	0
3	NAG	Q	1	14/15	0.92	0.25	92,98,106,113	0
6	SIA	U	3	20/21	0.94	0.16	71,73,82,85	0
3	NAG	G	1	14/15	0.94	0.14	84,89,98,109	0
6	GAL	U	2	11/12	0.95	0.13	87,93,97,97	0
6	GAL	K	2	11/12	0.95	0.22	84,90,95,95	0
6	GAL	P	2	11/12	0.95	0.18	92,99,105,105	0
5	NAG	N	2	14/15	0.96	0.21	76,89,98,106	0
5	NAG	S	2	14/15	0.96	0.14	77,93,100,109	0
5	NAG	I	2	14/15	0.96	0.18	87,95,107,114	0

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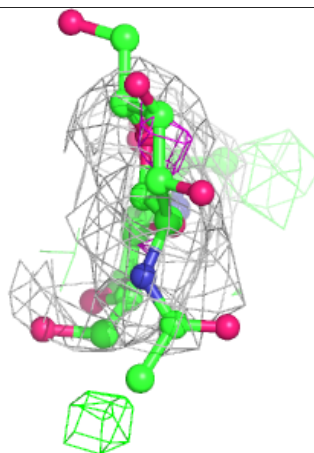
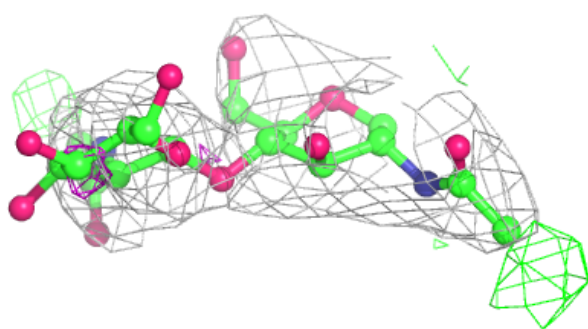
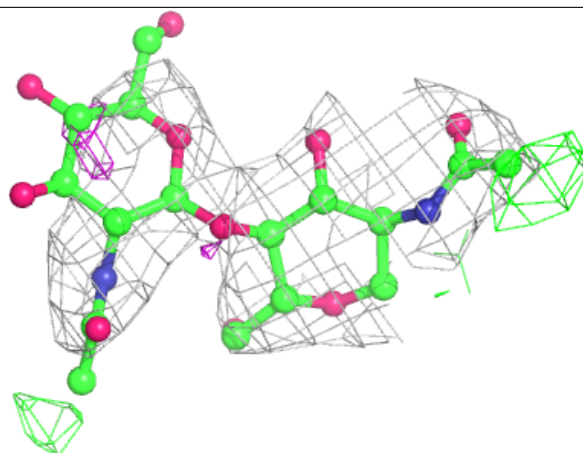
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	NAG	I	1	14/15	0.96	0.15	78,82,85,86	0
6	SIA	P	3	20/21	0.96	0.23	72,79,86,86	0
5	NAG	S	1	14/15	0.97	0.13	82,86,92,93	0
6	SIA	K	3	20/21	0.97	0.20	70,75,81,82	0
5	NAG	N	1	14/15	0.97	0.13	75,79,81,82	0

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.

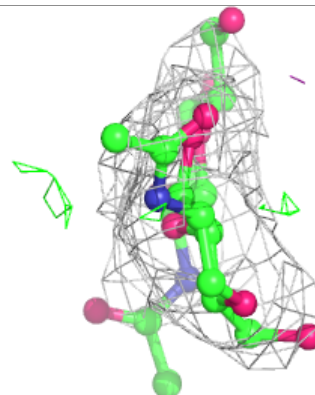
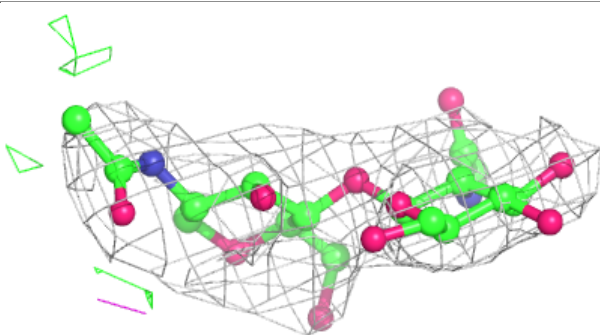
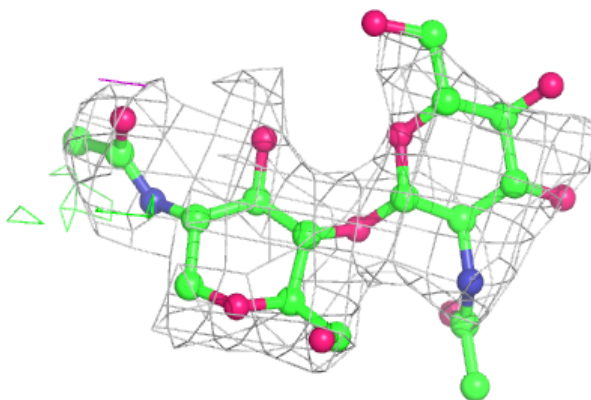


Electron density around Chain L:

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

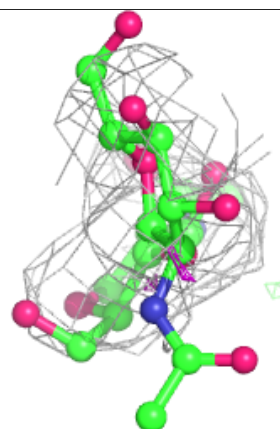
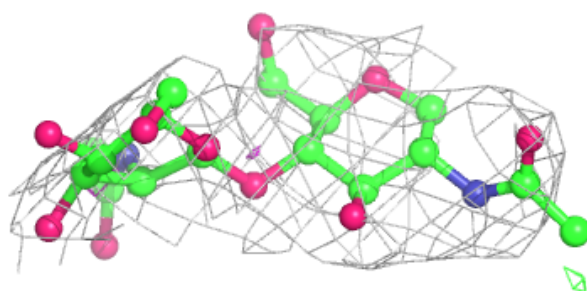
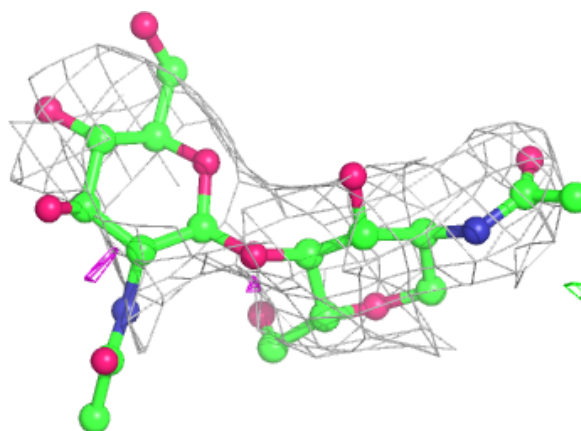
**Electron density around Chain Q:**

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



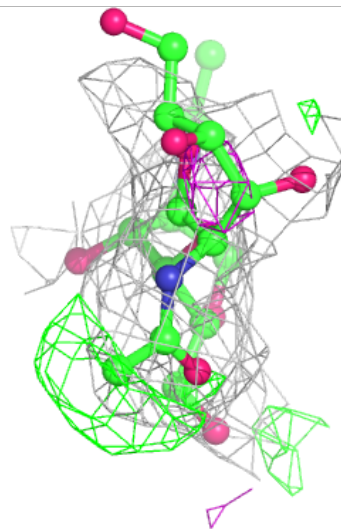
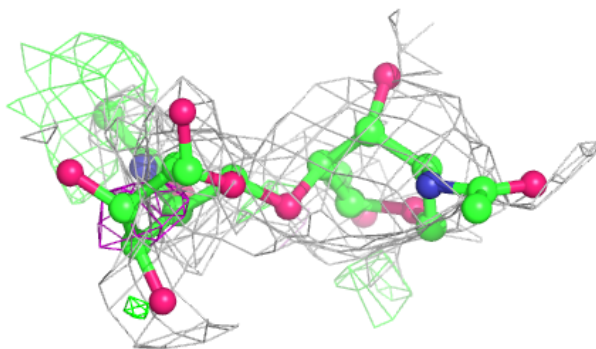
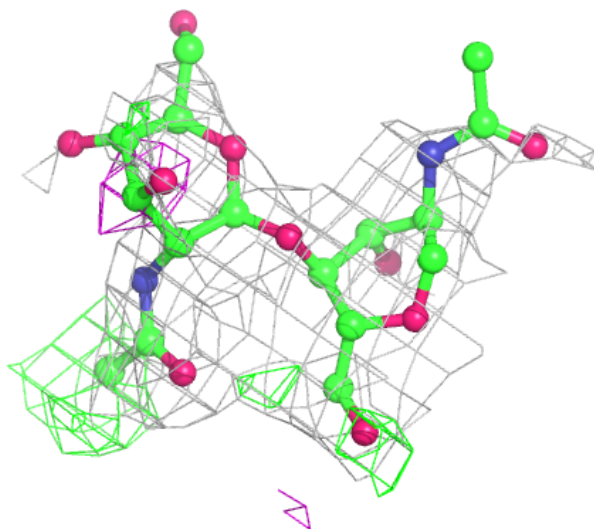
Electron density around Chain R:

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



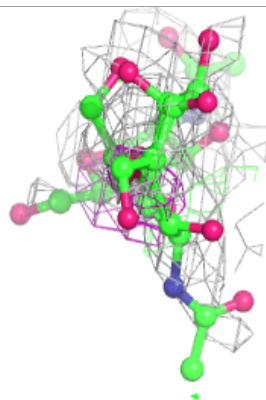
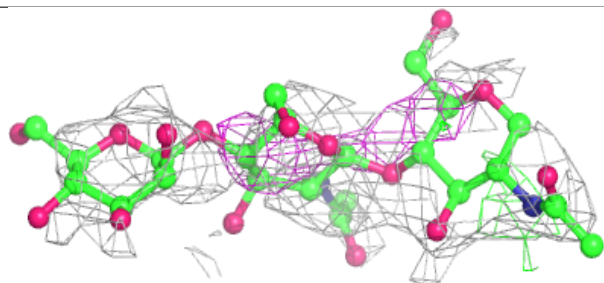
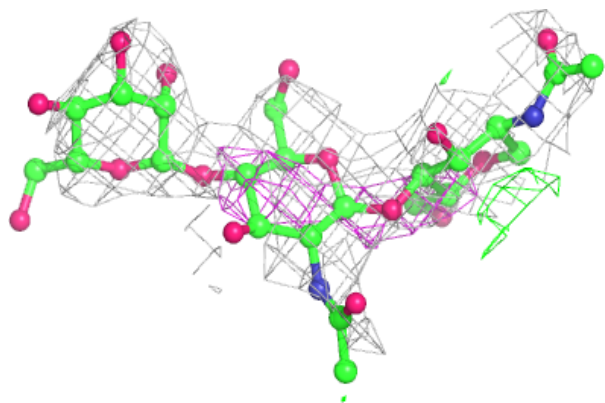
Electron density around Chain T:

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

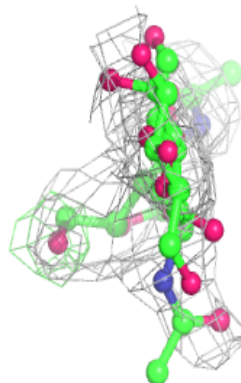
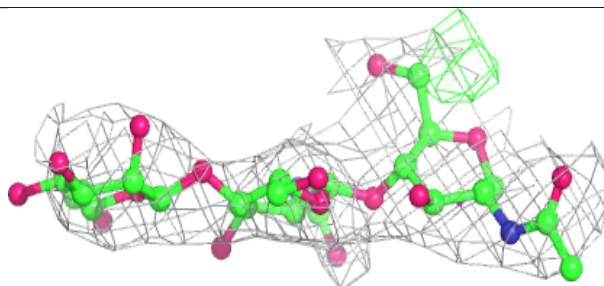
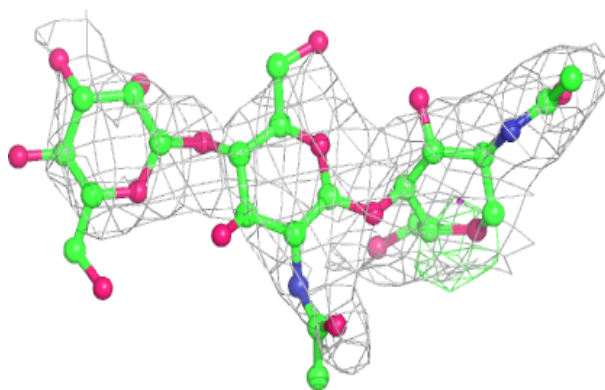


Electron density around Chain H:

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

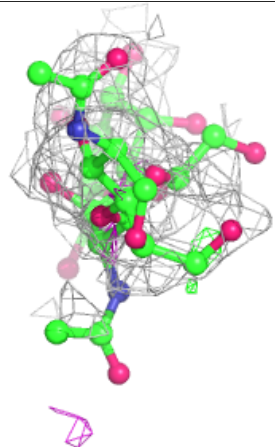
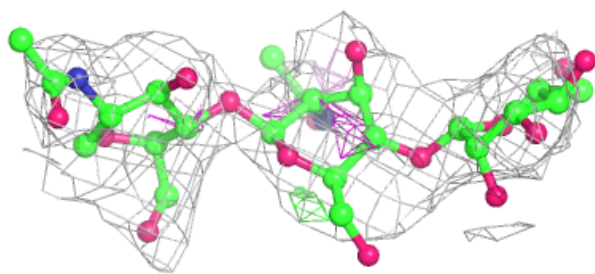
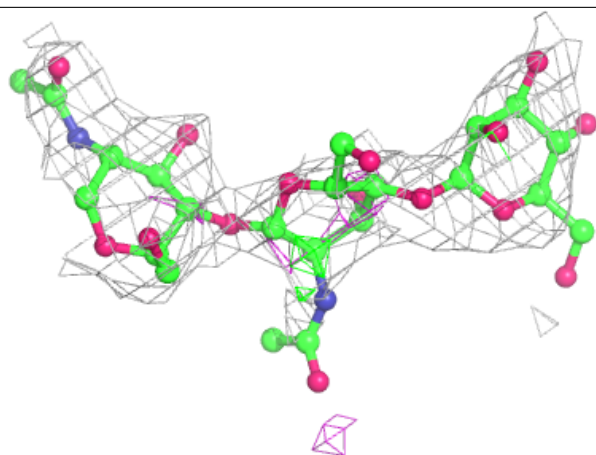
**Electron density around Chain J:**

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

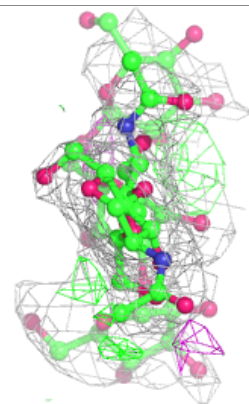
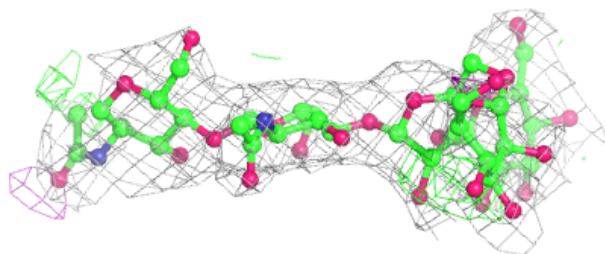
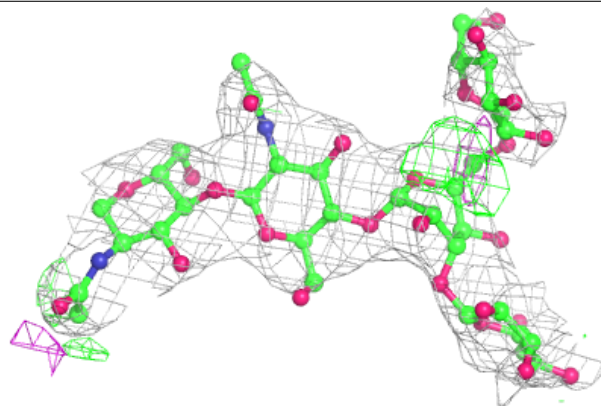


Electron density around Chain M:

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

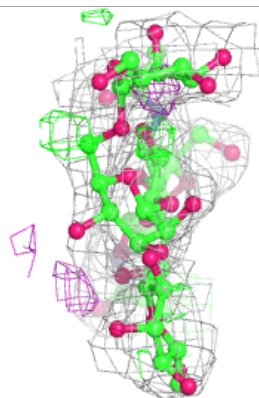
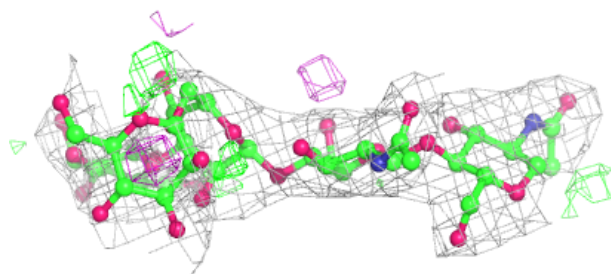
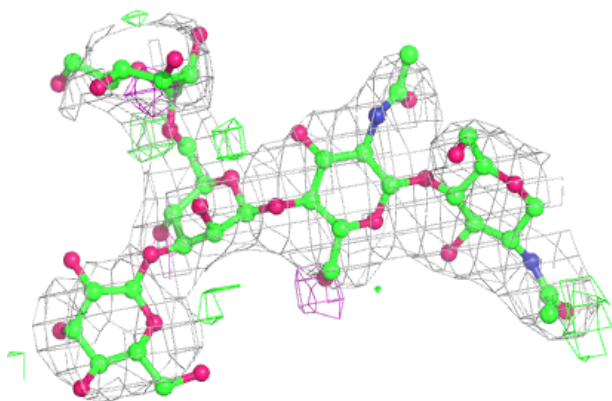
**Electron density around Chain I:**

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

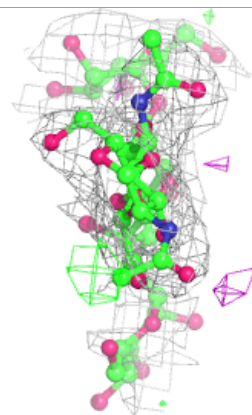
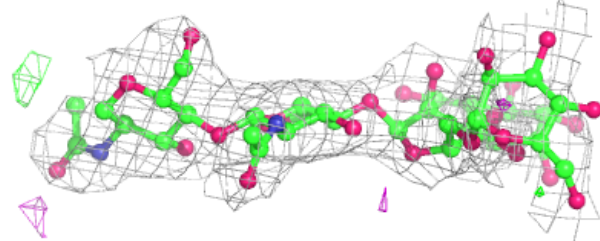
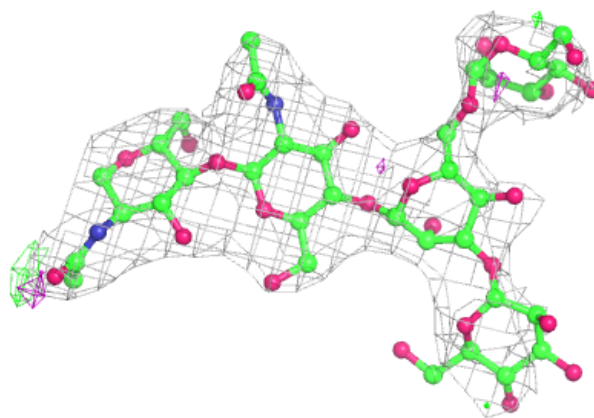


Electron density around Chain N:

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

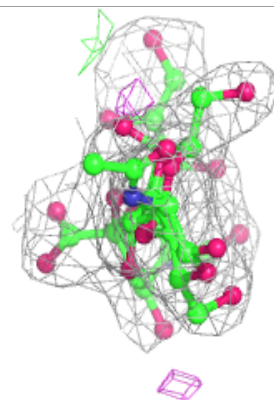
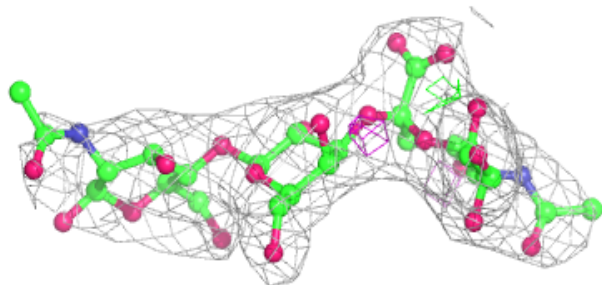
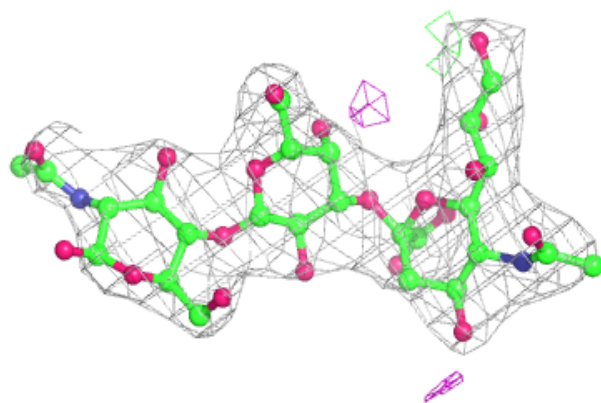
**Electron density around Chain S:**

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

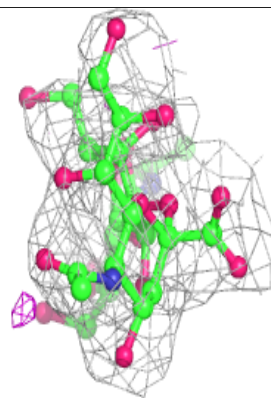
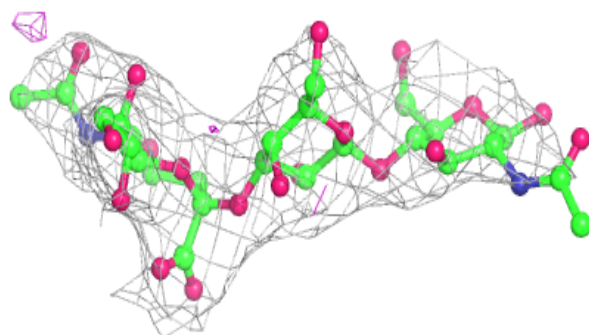
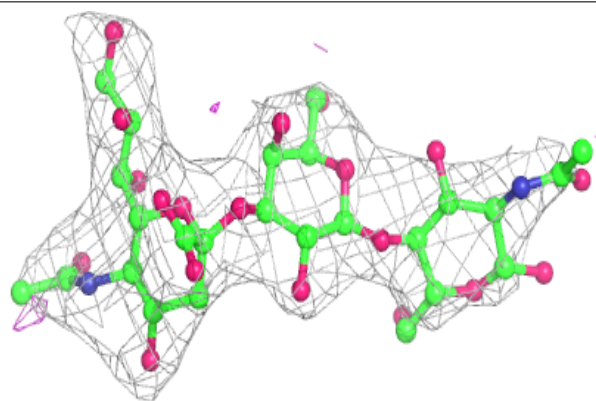


Electron density around Chain K:

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

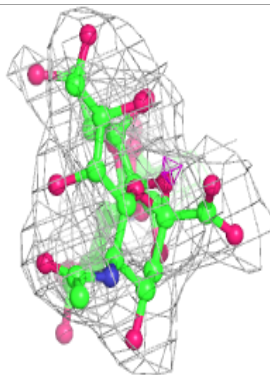
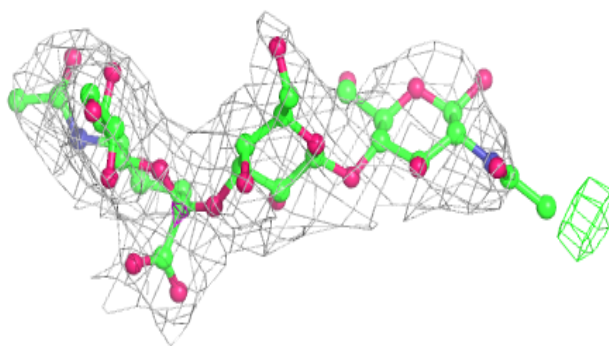
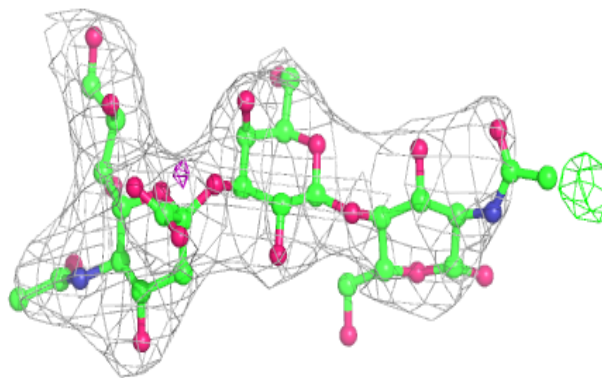
**Electron density around Chain P:**

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

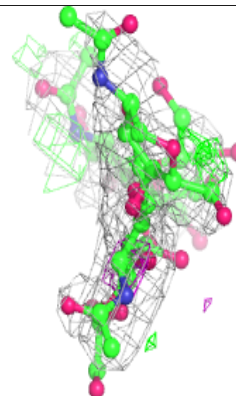
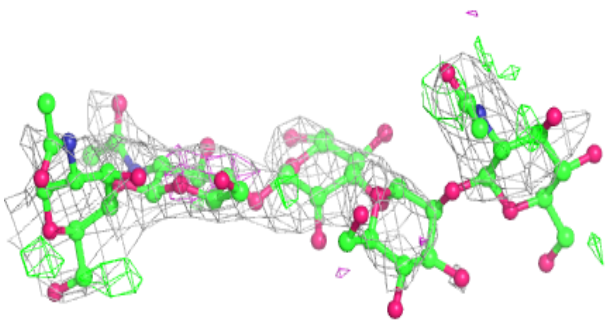
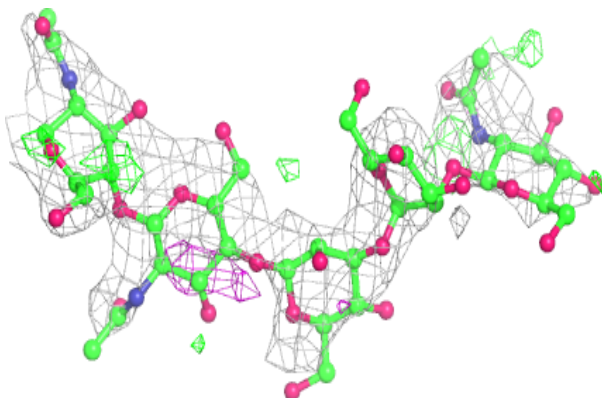


Electron density around Chain U:

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

**Electron density around Chain O:**

$2mF_o-DF_c$ (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(\AA^2)	Q<0.9
8	NAG	C	601	14/15	0.78	0.40	107,113,124,126	0
8	NAG	A	601	14/15	0.78	0.34	114,121,125,126	0
9	SO4	D	1176	5/5	0.93	0.13	92,94,96,96	0
9	SO4	B	1176	5/5	0.95	0.20	93,99,99,104	0
9	SO4	F	1176	5/5	0.96	0.17	91,91,93,94	0

6.5 Other polymers [i](#)

There are no such residues in this entry.