



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 01:26 PM UTC

PDB ID : 4NCO / pdb_00004nco
Title : Crystal Structure of the BG505 SOSIP gp140 HIV-1 Env trimer in Complex with the Broadly Neutralizing Fab PGT122
Authors : Julien, J.-P.; Stanfield, R.L.; Lyumkis, D.; Ward, A.B.; Wilson, I.A.
Deposited on : 2013-10-24
Resolution : 4.70 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 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
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

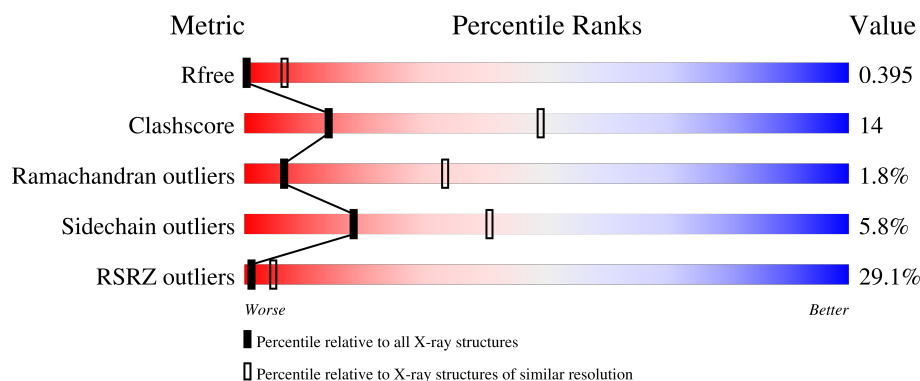
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.70 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.



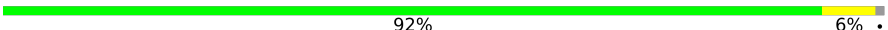
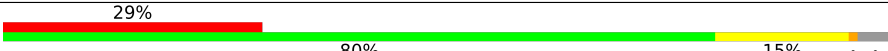
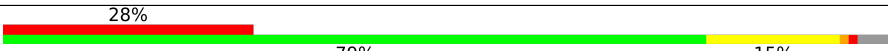
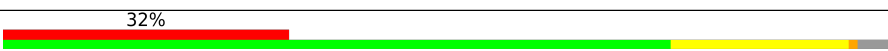
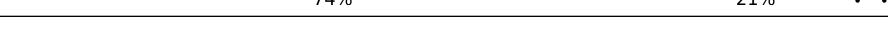
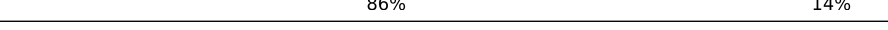
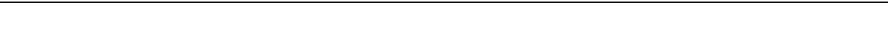
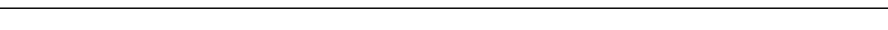
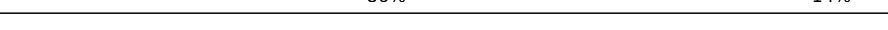
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1020 (5.48-3.92)
Clashscore	190562	1005 (5.40-3.96)
Ramachandran outliers	187476	1095 (5.50-3.90)
Sidechain outliers	187428	1077 (5.50-3.90)
RSRZ outliers	180081	1015 (5.48-3.92)

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	475	25% 48% 33% 7% 12%
1	E	475	24% 48% 34% 6% 12%
1	I	475	27% 48% 34% 6% 12%
2	B	78	95% ..
2	F	78	95% ..

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Mol	Chain	Length	Quality of chain
2	J	78	
3	C	211	
3	G	211	
3	K	211	
4	D	235	
4	H	235	
4	L	235	
5	M	7	
5	N	7	
5	O	7	
5	P	7	
5	S	7	
5	T	7	
5	U	7	
5	V	7	
5	Y	7	
5	Z	7	
5	a	7	
5	b	7	
6	Q	9	
6	W	9	
6	c	9	
7	R	2	
7	X	2	
7	d	2	

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
5	NAG	O	1	-	-	-	X
7	MAN	R	1	-	-	-	X
7	MAN	R	2	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 22014 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 BG505 SOSIP gp120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	420	Total	C	N	O	S	0	0	0
			3096	1937	544	590	25			
1	E	420	Total	C	N	O	S	0	0	0
			3096	1937	544	590	25			
1	I	420	Total	C	N	O	S	0	0	0
			3096	1937	544	590	25			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	332	ASN	THR	engineered mutation	UNP Q2N0S6
A	501	CYS	ALA	engineered mutation	UNP Q2N0S6
E	332	ASN	THR	engineered mutation	UNP Q2N0S6
E	501	CYS	ALA	engineered mutation	UNP Q2N0S6
I	332	ASN	THR	engineered mutation	UNP Q2N0S6
I	501	CYS	ALA	engineered mutation	UNP Q2N0S6

- Molecule 2 is a protein called BG505 SOSIP gp41.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	77	Total	C	N	O	0	0	0
			385	231	77	77			
2	F	77	Total	C	N	O	0	0	0
			385	231	77	77			
2	J	77	Total	C	N	O	0	0	0
			385	231	77	77			

- Molecule 3 is a protein called PGT122 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	202	Total	C	N	O	S	0	0	0
			1530	964	255	307	4			

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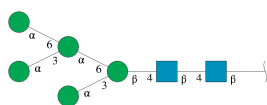
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	202	Total	C	N	O	S	0	0	0
			1530	964	255	307	4			
3	K	202	Total	C	N	O	S	0	0	0
			1530	964	255	307	4			

- Molecule 4 is a protein called PGT122 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	226	Total	C	N	O	S	0	0	0
			1728	1100	293	330	5			
4	H	226	Total	C	N	O	S	0	0	0
			1728	1100	293	330	5			
4	L	226	Total	C	N	O	S	0	0	0
			1728	1100	293	330	5			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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.



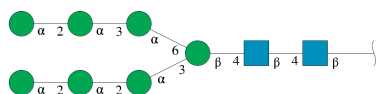
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	M	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	N	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	O	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	P	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	S	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	T	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	U	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	V	7	Total	C	N	O	0	0	0
			83	46	2	35			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	Y	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	Z	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	a	7	Total	C	N	O	0	0	0
			83	46	2	35			
5	b	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-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
6	Q	9	Total	C	N	O	0	0	0
			105	58	2	45			
6	W	9	Total	C	N	O	0	0	0
			105	58	2	45			
6	c	9	Total	C	N	O	0	0	0
			105	58	2	45			

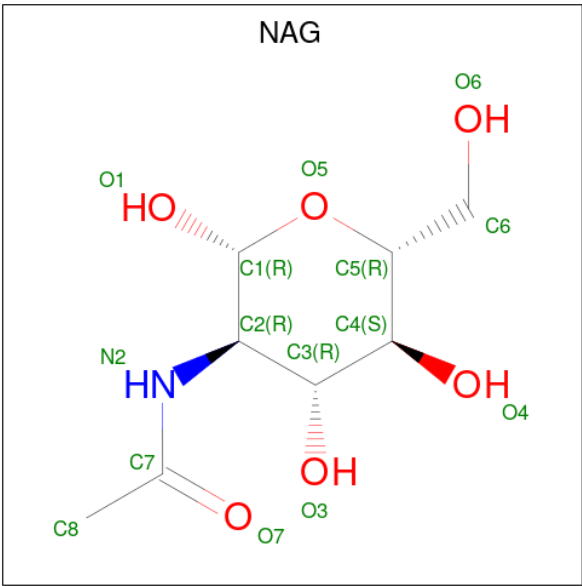
- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
7	R	2	Total	C	O	0	0	0
			22	12	10			
7	X	2	Total	C	O	0	0	0
			22	12	10			
7	d	2	Total	C	O	0	0	0
			22	12	10			

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula:

C₈H₁₅NO₆).



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

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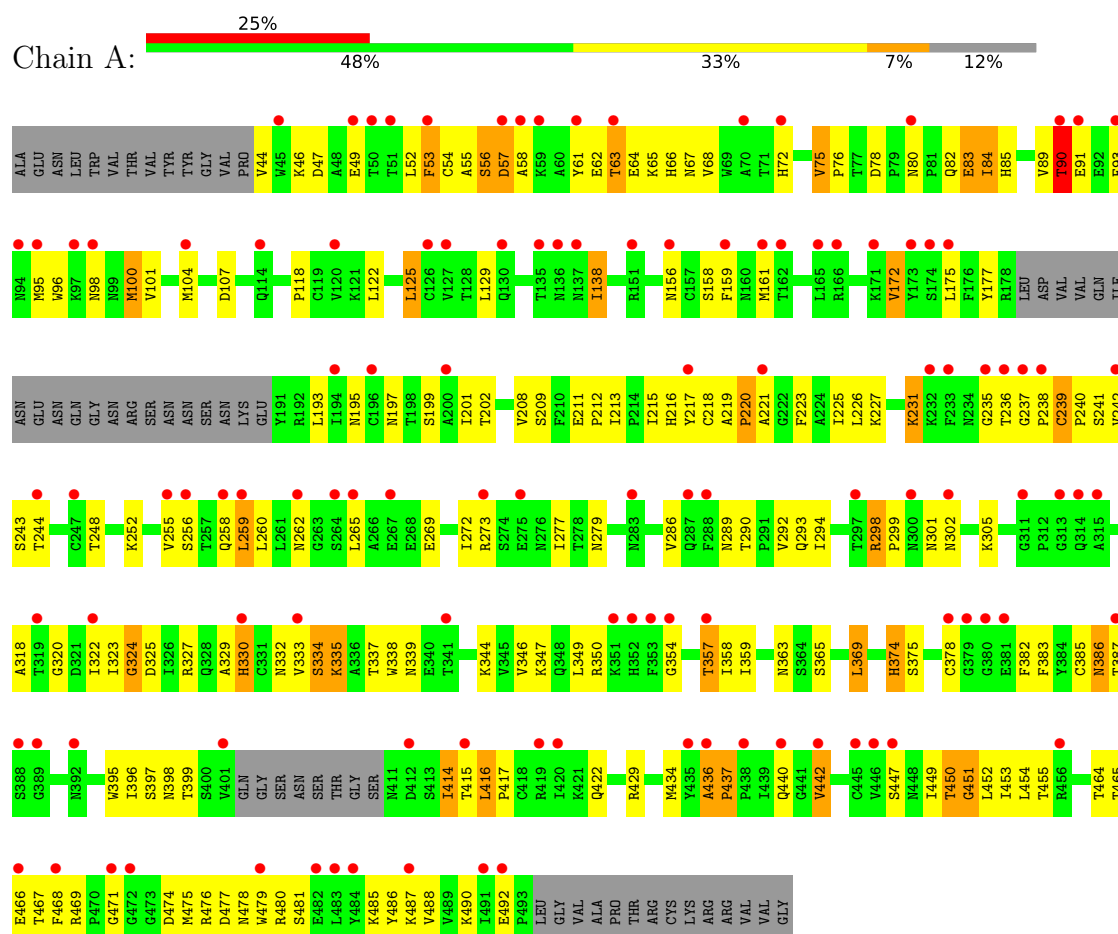
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	E	1	Total	C	N	O	0	0
			14	8	1	5		
8	E	1	Total	C	N	O	0	0
			14	8	1	5		
8	E	1	Total	C	N	O	0	0
			14	8	1	5		
8	E	1	Total	C	N	O	0	0
			14	8	1	5		
8	E	1	Total	C	N	O	0	0
			14	8	1	5		
8	E	1	Total	C	N	O	0	0
	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

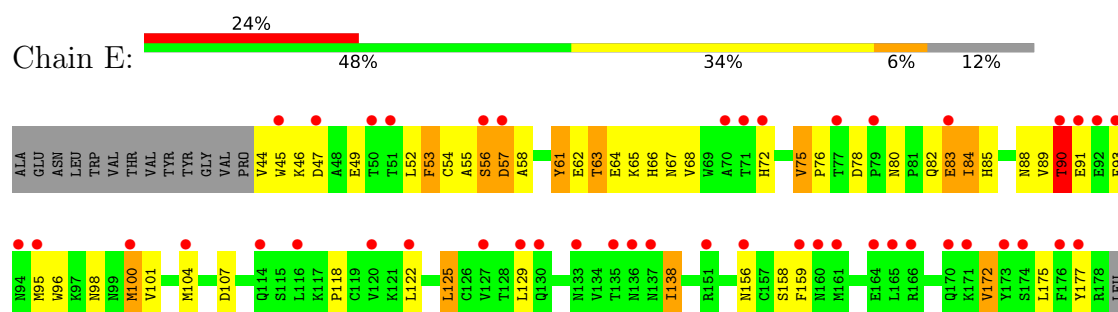
3 Residue-property plots [i](#)

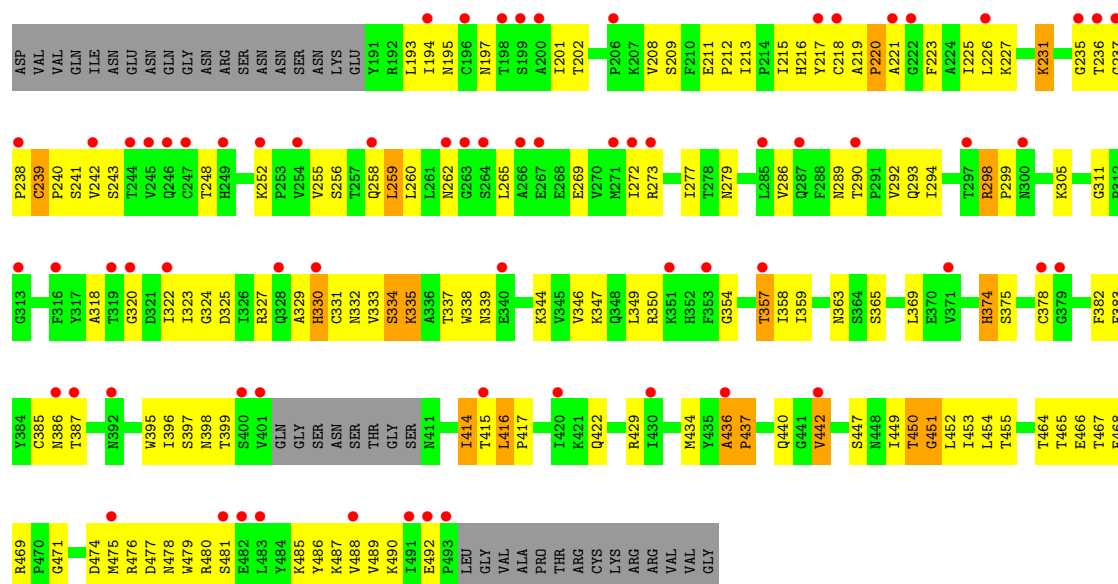
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: BG505 SOSIP gp120

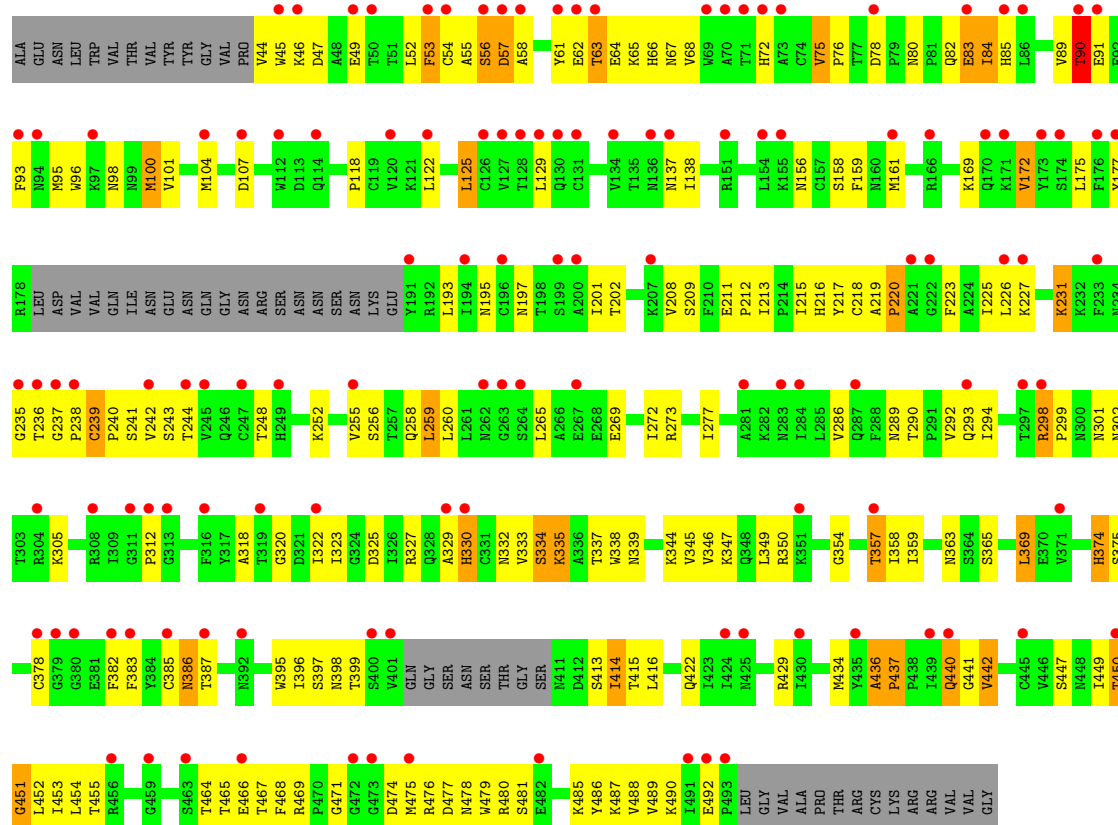


• Molecule 1: BG505 SOSIP gp120



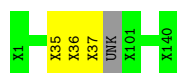


• Molecule 1: BG505 SOSIP gp120



• Molecule 2: BG505 SOSIP gp41





- Molecule 2: BG505 SOSIP gp41

Chain F: 95%



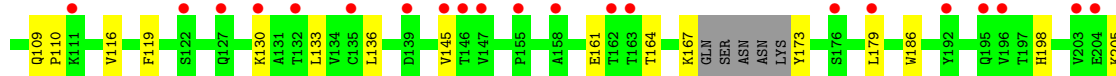
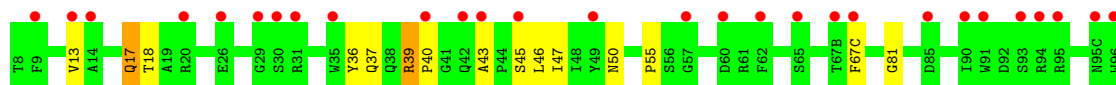
- Molecule 2: BG505 SOSIP gp41

Chain J: 92%



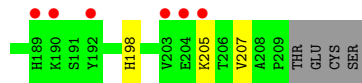
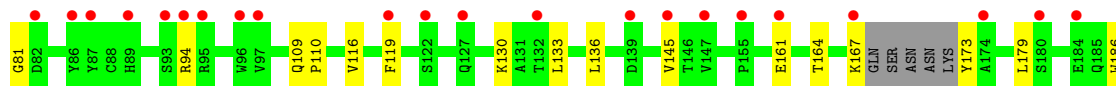
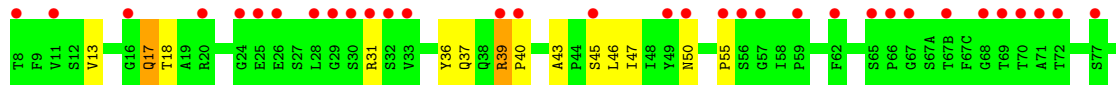
- Molecule 3: PGT122 light chain

Chain C: 24% 81% 14%



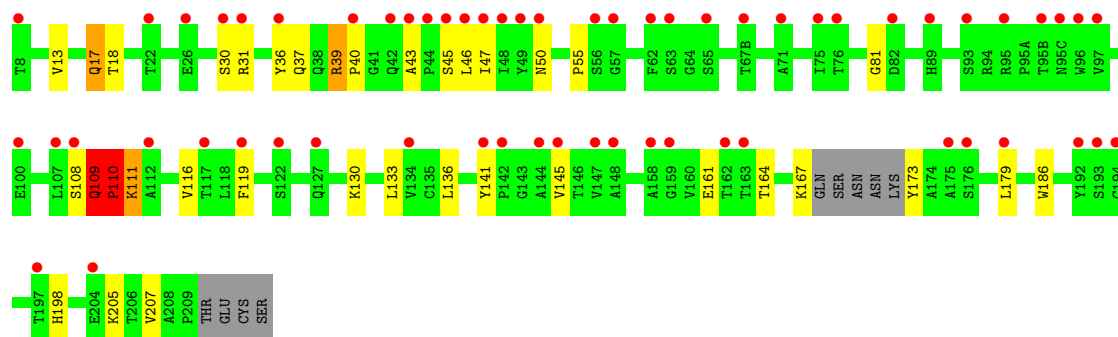
- Molecule 3: PGT122 light chain

Chain G: 29% 80% 15%

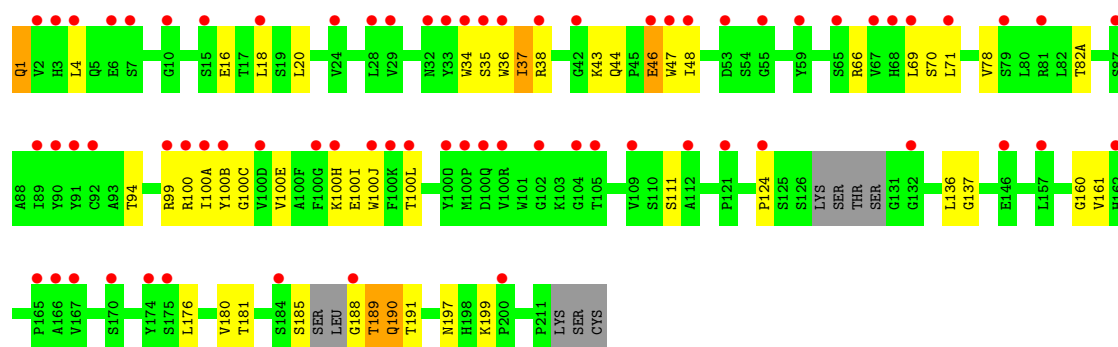
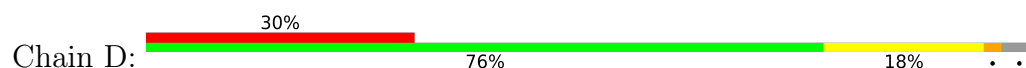


- Molecule 3: PGT122 light chain

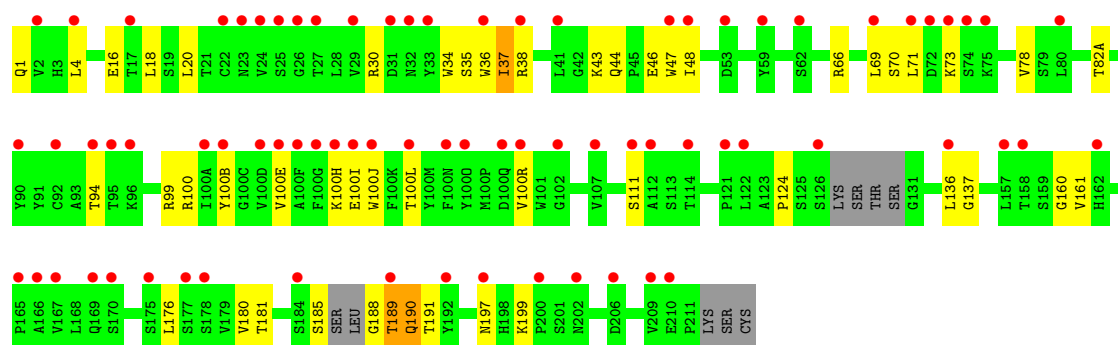
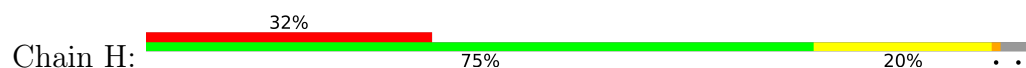
Chain K: 28% 79% 15%



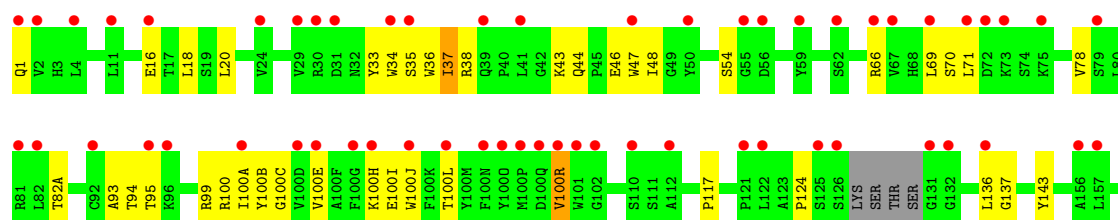
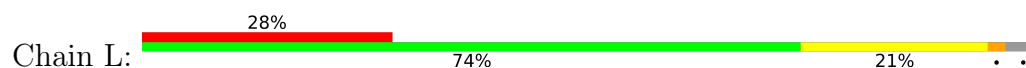
• Molecule 4: PGT122 heavy chain



• Molecule 4: PGT122 heavy chain



• Molecule 4: PGT122 heavy chain





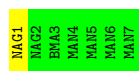
- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M: 57% 29% 14%



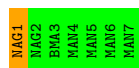
- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N: 86% 14%



- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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 O: 86% 14%



- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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 P: 86% 14%



- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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 S: 71% 14% 14%

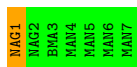


- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-

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

Chain T:  86% 14%

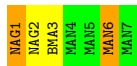
- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain U:  86% 14%

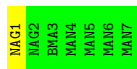
- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain V:  86% 14%

- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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 Y:  43% 29% 29%

- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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:  86% 14%

- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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 a:  86% 14%

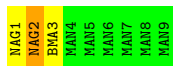
- Molecule 5: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain b:  86% 14%



- Molecule 6: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)- α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain Q:  67% 22% 11%



- Molecule 6: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)- α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain W:  78% 22%



- Molecule 6: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)- α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain c:  56% 22% 22%

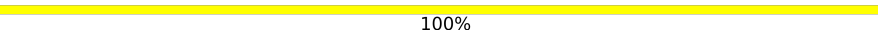


- Molecule 7: α -D-mannopyranose-(1-2)- α -D-mannopyranose

Chain R:  50% 50%



- Molecule 7: α -D-mannopyranose-(1-2)- α -D-mannopyranose

Chain X:  100%

MAN1
MAN2

- Molecule 7: α -D-mannopyranose-(1-2)- α -D-mannopyranose

Chain d:

50%

50%

MAN1
MAN2

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	152.20Å 260.72Å 283.18Å 90.00° 99.56° 90.00°	Depositor
Resolution (Å)	39.89 – 4.70 39.89 – 4.70	Depositor EDS
% Data completeness (in resolution range)	89.0 (39.89-4.70) 88.9 (39.89-4.70)	Depositor EDS
R_{merge}	0.30	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 4.63Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.375 , 0.389 0.377 , 0.395	Depositor DCC
R_{free} test set	2621 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	136.6	Xtriage
Anisotropy	1.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 597.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.31$, $\langle L^2 \rangle = 0.15$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.76	EDS
Total number of atoms	22014	wwPDB-VP
Average B, all atoms (Å ²)	262.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.45	0/3161	1.24	33/4306 (0.8%)
1	E	0.47	0/3161	1.25	36/4306 (0.8%)
1	I	0.45	0/3161	1.23	31/4306 (0.7%)
3	C	0.35	0/1571	0.84	4/2151 (0.2%)
3	G	0.36	0/1571	0.84	4/2151 (0.2%)
3	K	0.36	0/1571	0.91	9/2151 (0.4%)
4	D	0.39	1/1774 (0.1%)	0.84	6/2421 (0.2%)
4	H	0.46	1/1774 (0.1%)	0.83	6/2421 (0.2%)
4	L	0.67	1/1774 (0.1%)	0.86	9/2421 (0.4%)
All	All	0.45	3/19518 (0.0%)	1.06	138/26634 (0.5%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	E	0	3
1	I	0	3
3	C	0	1
3	G	0	1
3	K	0	1
All	All	0	12

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	100(R)	VAL	C-N	22.47	1.66	1.33
4	H	100(R)	VAL	C-N	10.38	1.48	1.33
4	D	46	GLU	C-N	-5.00	1.26	1.33

The worst 5 of 138 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	335	LYS	N-CA-C	13.96	127.77	111.71
1	A	335	LYS	N-CA-C	13.68	127.69	111.82
1	I	335	LYS	N-CA-C	13.50	127.48	111.82
1	I	239	CYS	CA-C-N	13.47	133.47	119.19
1	I	239	CYS	C-N-CA	13.47	133.47	119.19

There are no chirality outliers.

5 of 12 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	236	THR	Peptide
1	A	56	SER	Peptide
1	A	80	ASN	Peptide
3	C	110	PRO	Peptide
1	E	56	SER	Peptide

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	3096	0	2821	146	0
1	E	3096	0	2821	145	0
1	I	3096	0	2821	148	0
2	B	385	0	82	2	0
2	F	385	0	82	2	0
2	J	385	0	82	3	0
3	C	1530	0	1472	21	0
3	G	1530	0	1472	19	0
3	K	1530	0	1472	23	0
4	D	1728	0	1699	37	0
4	H	1728	0	1699	33	0
4	L	1728	0	1699	46	0
5	M	83	0	70	1	0
5	N	83	0	70	0	0
5	O	83	0	70	6	0
5	P	83	0	70	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	S	83	0	70	1	0
5	T	83	0	70	0	0
5	U	83	0	70	6	0
5	V	83	0	70	0	0
5	Y	83	0	70	3	0
5	Z	83	0	70	0	0
5	a	83	0	70	4	0
5	b	83	0	70	0	0
6	Q	105	0	88	2	0
6	W	105	0	88	0	0
6	c	105	0	88	4	0
7	R	22	0	19	3	0
7	X	22	0	19	1	0
7	d	22	0	19	2	0
8	A	140	0	130	3	0
8	E	140	0	130	3	0
8	I	140	0	130	2	0
All	All	22014	0	19773	603	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 14.

The worst 5 of 603 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:I:293:GLN:HB2	1:I:334:SER:HB3	1.33	1.11
1:E:293:GLN:HB2	1:E:334:SER:HB3	1.33	1.09
1:A:293:GLN:HB2	1:A:334:SER:HB3	1.33	1.08
1:I:91:GLU:HA	1:I:239:CYS:O	1.74	0.88
1:E:91:GLU:HA	1:E:239:CYS:O	1.74	0.86

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	414/475 (87%)	349 (84%)	51 (12%)	14 (3%)	3	20
1	E	414/475 (87%)	348 (84%)	51 (12%)	15 (4%)	2	20
1	I	414/475 (87%)	348 (84%)	52 (13%)	14 (3%)	3	20
3	C	198/211 (94%)	191 (96%)	7 (4%)	0	100	100
3	G	198/211 (94%)	191 (96%)	7 (4%)	0	100	100
3	K	198/211 (94%)	192 (97%)	6 (3%)	0	100	100
4	D	220/235 (94%)	213 (97%)	6 (3%)	1 (0%)	24	63
4	H	220/235 (94%)	213 (97%)	6 (3%)	1 (0%)	24	63
4	L	220/235 (94%)	213 (97%)	6 (3%)	1 (0%)	24	63
All	All	2496/2763 (90%)	2258 (90%)	192 (8%)	46 (2%)	6	33

5 of 46 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	57	ASP
1	A	83	GLU
1	A	323	ILE
1	A	429	ARG
1	E	57	ASP

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	317/422 (75%)	294 (93%)	23 (7%)	13	34
1	E	317/422 (75%)	295 (93%)	22 (7%)	14	37
1	I	317/422 (75%)	293 (92%)	24 (8%)	12	33
3	C	171/180 (95%)	163 (95%)	8 (5%)	23	45
3	G	171/180 (95%)	163 (95%)	8 (5%)	23	45

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	K	171/180 (95%)	163 (95%)	8 (5%)	23	45
4	D	196/205 (96%)	187 (95%)	9 (5%)	24	46
4	H	196/205 (96%)	187 (95%)	9 (5%)	24	46
4	L	196/205 (96%)	187 (95%)	9 (5%)	24	46
All	All	2052/2421 (85%)	1932 (94%)	120 (6%)	18	40

5 of 120 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	382	PHE
3	K	164	THR
4	H	16	GLU
3	K	161	GLU
4	L	190	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
3	K	198	HIS
3	K	109	GLN
1	E	398	ASN
3	K	17	GLN
1	E	374	HIS

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 ⓘ

117 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	M	1	5,1	14,14,15	0.57	0	17,19,21	1.35	1 (5%)
5	NAG	M	2	5	14,14,15	0.54	0	17,19,21	0.82	0
5	BMA	M	3	5	11,11,12	0.68	0	15,15,17	0.77	1 (6%)
5	MAN	M	4	5	11,11,12	0.69	0	15,15,17	0.76	0
5	MAN	M	5	5	11,11,12	0.62	0	15,15,17	0.56	0
5	MAN	M	6	5	11,11,12	0.64	0	15,15,17	1.02	1 (6%)
5	MAN	M	7	5	11,11,12	0.58	0	15,15,17	0.56	0
5	NAG	N	1	5,1	14,14,15	0.59	0	17,19,21	1.42	3 (17%)
5	NAG	N	2	5	14,14,15	0.55	0	17,19,21	0.88	0
5	BMA	N	3	5	11,11,12	0.75	0	15,15,17	0.77	0
5	MAN	N	4	5	11,11,12	0.71	0	15,15,17	0.99	0
5	MAN	N	5	5	11,11,12	0.62	0	15,15,17	0.56	0
5	MAN	N	6	5	11,11,12	0.69	0	15,15,17	0.69	0
5	MAN	N	7	5	11,11,12	0.57	0	15,15,17	0.60	0
5	NAG	O	1	5,1	14,14,15	0.74	0	17,19,21	1.45	3 (17%)
5	NAG	O	2	5	14,14,15	0.55	0	17,19,21	0.74	0
5	BMA	O	3	5	11,11,12	0.73	0	15,15,17	0.65	0
5	MAN	O	4	5	11,11,12	0.68	0	15,15,17	0.81	0
5	MAN	O	5	5	11,11,12	0.62	0	15,15,17	0.54	0
5	MAN	O	6	5	11,11,12	0.69	0	15,15,17	0.69	0
5	MAN	O	7	5	11,11,12	0.58	0	15,15,17	0.58	0
5	NAG	P	1	5,1	14,14,15	0.60	0	17,19,21	1.16	1 (5%)
5	NAG	P	2	5	14,14,15	0.57	0	17,19,21	0.77	0
5	BMA	P	3	5	11,11,12	0.71	0	15,15,17	0.71	0
5	MAN	P	4	5	11,11,12	0.69	0	15,15,17	0.81	0
5	MAN	P	5	5	11,11,12	0.61	0	15,15,17	0.54	0
5	MAN	P	6	5	11,11,12	0.67	0	15,15,17	0.70	0
5	MAN	P	7	5	11,11,12	0.57	0	15,15,17	0.56	0
6	NAG	Q	1	1,6	14,14,15	0.61	0	17,19,21	1.50	3 (17%)
6	NAG	Q	2	6	14,14,15	0.52	0	17,19,21	1.09	1 (5%)
6	BMA	Q	3	6	11,11,12	0.69	0	15,15,17	1.24	1 (6%)
6	MAN	Q	4	6	11,11,12	0.52	0	15,15,17	0.73	0
6	MAN	Q	5	6	11,11,12	0.59	0	15,15,17	0.73	0
6	MAN	Q	6	6	11,11,12	0.64	0	15,15,17	0.56	0
6	MAN	Q	7	6	11,11,12	0.65	0	15,15,17	0.95	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	MAN	Q	8	6	11,11,12	0.60	0	15,15,17	0.57	0
6	MAN	Q	9	6	11,11,12	0.61	0	15,15,17	0.50	0
7	MAN	R	1	7	11,11,12	0.70	0	15,15,17	1.08	1 (6%)
7	MAN	R	2	7	11,11,12	0.68	0	15,15,17	0.58	0
5	NAG	S	1	5,1	14,14,15	0.57	0	17,19,21	1.35	1 (5%)
5	NAG	S	2	5	14,14,15	0.53	0	17,19,21	0.83	0
5	BMA	S	3	5	11,11,12	0.68	0	15,15,17	0.77	1 (6%)
5	MAN	S	4	5	11,11,12	0.69	0	15,15,17	0.75	0
5	MAN	S	5	5	11,11,12	0.61	0	15,15,17	0.56	0
5	MAN	S	6	5	11,11,12	0.67	0	15,15,17	0.73	0
5	MAN	S	7	5	11,11,12	0.59	0	15,15,17	0.55	0
5	NAG	T	1	5,1	14,14,15	0.59	0	17,19,21	1.41	3 (17%)
5	NAG	T	2	5	14,14,15	0.54	0	17,19,21	0.89	0
5	BMA	T	3	5	11,11,12	0.76	0	15,15,17	0.78	0
5	MAN	T	4	5	11,11,12	0.72	0	15,15,17	0.98	0
5	MAN	T	5	5	11,11,12	0.61	0	15,15,17	0.56	0
5	MAN	T	6	5	11,11,12	0.69	0	15,15,17	0.69	0
5	MAN	T	7	5	11,11,12	0.57	0	15,15,17	0.60	0
5	NAG	U	1	5,1	14,14,15	0.73	0	17,19,21	1.43	3 (17%)
5	NAG	U	2	5	14,14,15	0.54	0	17,19,21	0.72	0
5	BMA	U	3	5	11,11,12	0.73	0	15,15,17	0.66	0
5	MAN	U	4	5	11,11,12	0.68	0	15,15,17	0.80	0
5	MAN	U	5	5	11,11,12	0.62	0	15,15,17	0.54	0
5	MAN	U	6	5	11,11,12	0.68	0	15,15,17	0.70	0
5	MAN	U	7	5	11,11,12	0.58	0	15,15,17	0.59	0
5	NAG	V	1	5,1	14,14,15	0.60	0	17,19,21	1.10	1 (5%)
5	NAG	V	2	5	14,14,15	0.56	0	17,19,21	0.77	0
5	BMA	V	3	5	11,11,12	0.70	0	15,15,17	0.73	0
5	MAN	V	4	5	11,11,12	0.69	0	15,15,17	0.81	0
5	MAN	V	5	5	11,11,12	0.63	0	15,15,17	0.54	0
5	MAN	V	6	5	11,11,12	0.68	0	15,15,17	0.69	0
5	MAN	V	7	5	11,11,12	0.59	0	15,15,17	0.56	0
6	NAG	W	1	1,6	14,14,15	0.62	0	17,19,21	1.54	3 (17%)
6	NAG	W	2	6	14,14,15	0.56	0	17,19,21	0.76	0
6	BMA	W	3	6	11,11,12	0.74	0	15,15,17	0.89	1 (6%)
6	MAN	W	4	6	11,11,12	0.52	0	15,15,17	0.70	0
6	MAN	W	5	6	11,11,12	0.58	0	15,15,17	0.74	0
6	MAN	W	6	6	11,11,12	0.65	0	15,15,17	0.56	0
6	MAN	W	7	6	11,11,12	0.65	0	15,15,17	0.95	0
6	MAN	W	8	6	11,11,12	0.61	0	15,15,17	0.56	0
6	MAN	W	9	6	11,11,12	0.62	0	15,15,17	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	MAN	X	1	7	11,11,12	0.64	0	15,15,17	0.93	1 (6%)
7	MAN	X	2	7	11,11,12	0.64	0	15,15,17	0.71	0
5	NAG	Y	1	5,1	14,14,15	0.57	0	17,19,21	1.42	1 (5%)
5	NAG	Y	2	5	14,14,15	0.51	0	17,19,21	0.96	1 (5%)
5	BMA	Y	3	5	11,11,12	0.66	0	15,15,17	0.82	1 (6%)
5	MAN	Y	4	5	11,11,12	0.68	0	15,15,17	0.73	0
5	MAN	Y	5	5	11,11,12	0.63	0	15,15,17	0.54	0
5	MAN	Y	6	5	11,11,12	0.67	0	15,15,17	1.04	1 (6%)
5	MAN	Y	7	5	11,11,12	0.58	0	15,15,17	0.55	0
5	NAG	Z	1	5,1	14,14,15	0.61	0	17,19,21	1.46	2 (11%)
5	NAG	Z	2	5	14,14,15	0.57	0	17,19,21	0.89	0
5	BMA	Z	3	5	11,11,12	0.75	0	15,15,17	0.78	0
5	MAN	Z	4	5	11,11,12	0.72	0	15,15,17	1.00	0
5	MAN	Z	5	5	11,11,12	0.62	0	15,15,17	0.56	0
5	MAN	Z	6	5	11,11,12	0.67	0	15,15,17	0.69	0
5	MAN	Z	7	5	11,11,12	0.55	0	15,15,17	0.61	0
5	NAG	a	1	5,1	14,14,15	0.71	0	17,19,21	1.47	3 (17%)
5	NAG	a	2	5	14,14,15	0.55	0	17,19,21	0.74	0
5	BMA	a	3	5	11,11,12	0.73	0	15,15,17	0.66	0
5	MAN	a	4	5	11,11,12	0.68	0	15,15,17	0.81	0
5	MAN	a	5	5	11,11,12	0.62	0	15,15,17	0.54	0
5	MAN	a	6	5	11,11,12	0.69	0	15,15,17	0.70	0
5	MAN	a	7	5	11,11,12	0.58	0	15,15,17	0.58	0
5	NAG	b	1	5,1	14,14,15	0.67	0	17,19,21	1.15	1 (5%)
5	NAG	b	2	5	14,14,15	0.57	0	17,19,21	0.79	0
5	BMA	b	3	5	11,11,12	0.72	0	15,15,17	0.71	0
5	MAN	b	4	5	11,11,12	0.69	0	15,15,17	0.81	0
5	MAN	b	5	5	11,11,12	0.62	0	15,15,17	0.54	0
5	MAN	b	6	5	11,11,12	0.69	0	15,15,17	0.69	0
5	MAN	b	7	5	11,11,12	0.58	0	15,15,17	0.57	0
6	NAG	c	1	1,6	14,14,15	0.55	0	17,19,21	1.56	3 (17%)
6	NAG	c	2	6	14,14,15	0.57	0	17,19,21	1.13	1 (5%)
6	BMA	c	3	6	11,11,12	0.70	0	15,15,17	1.30	1 (6%)
6	MAN	c	4	6	11,11,12	0.51	0	15,15,17	0.74	0
6	MAN	c	5	6	11,11,12	0.58	0	15,15,17	0.74	0
6	MAN	c	6	6	11,11,12	0.66	0	15,15,17	0.56	0
6	MAN	c	7	6	11,11,12	0.65	0	15,15,17	1.03	1 (6%)
6	MAN	c	8	6	11,11,12	0.61	0	15,15,17	0.56	0
6	MAN	c	9	6	11,11,12	0.62	0	15,15,17	0.50	0
7	MAN	d	1	7	11,11,12	0.71	0	15,15,17	1.03	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	MAN	d	2	7	11,11,12	0.73	0	15,15,17	0.68	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	M	1	5,1	-	3/6/23/26	0/1/1/1
5	NAG	M	2	5	-	0/6/23/26	0/1/1/1
5	BMA	M	3	5	-	0/2/19/22	0/1/1/1
5	MAN	M	4	5	-	0/2/19/22	0/1/1/1
5	MAN	M	5	5	-	0/2/19/22	0/1/1/1
5	MAN	M	6	5	-	0/2/19/22	0/1/1/1
5	MAN	M	7	5	-	0/2/19/22	0/1/1/1
5	NAG	N	1	5,1	-	0/6/23/26	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	-	1/2/19/22	0/1/1/1
5	MAN	N	5	5	-	0/2/19/22	0/1/1/1
5	MAN	N	6	5	-	0/2/19/22	0/1/1/1
5	MAN	N	7	5	-	0/2/19/22	0/1/1/1
5	NAG	O	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	O	2	5	-	0/6/23/26	0/1/1/1
5	BMA	O	3	5	-	0/2/19/22	0/1/1/1
5	MAN	O	4	5	-	2/2/19/22	0/1/1/1
5	MAN	O	5	5	-	0/2/19/22	0/1/1/1
5	MAN	O	6	5	-	0/2/19/22	0/1/1/1
5	MAN	O	7	5	-	0/2/19/22	0/1/1/1
5	NAG	P	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	P	2	5	-	0/6/23/26	0/1/1/1
5	BMA	P	3	5	-	0/2/19/22	0/1/1/1
5	MAN	P	4	5	-	2/2/19/22	0/1/1/1
5	MAN	P	5	5	-	0/2/19/22	0/1/1/1
5	MAN	P	6	5	-	0/2/19/22	0/1/1/1
5	MAN	P	7	5	-	0/2/19/22	0/1/1/1
6	NAG	Q	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	Q	2	6	-	0/6/23/26	0/1/1/1
6	BMA	Q	3	6	-	0/2/19/22	0/1/1/1
6	MAN	Q	4	6	-	0/2/19/22	0/1/1/1
6	MAN	Q	5	6	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	MAN	Q	6	6	-	0/2/19/22	0/1/1/1
6	MAN	Q	7	6	-	2/2/19/22	0/1/1/1
6	MAN	Q	8	6	-	0/2/19/22	0/1/1/1
6	MAN	Q	9	6	-	0/2/19/22	0/1/1/1
7	MAN	R	1	7	-	2/2/19/22	0/1/1/1
7	MAN	R	2	7	-	0/2/19/22	0/1/1/1
5	NAG	S	1	5,1	-	3/6/23/26	0/1/1/1
5	NAG	S	2	5	-	0/6/23/26	0/1/1/1
5	BMA	S	3	5	-	0/2/19/22	0/1/1/1
5	MAN	S	4	5	-	0/2/19/22	0/1/1/1
5	MAN	S	5	5	-	0/2/19/22	0/1/1/1
5	MAN	S	6	5	-	0/2/19/22	0/1/1/1
5	MAN	S	7	5	-	0/2/19/22	0/1/1/1
5	NAG	T	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	T	2	5	-	0/6/23/26	0/1/1/1
5	BMA	T	3	5	-	0/2/19/22	0/1/1/1
5	MAN	T	4	5	-	1/2/19/22	0/1/1/1
5	MAN	T	5	5	-	0/2/19/22	0/1/1/1
5	MAN	T	6	5	-	0/2/19/22	0/1/1/1
5	MAN	T	7	5	-	0/2/19/22	0/1/1/1
5	NAG	U	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	U	2	5	-	0/6/23/26	0/1/1/1
5	BMA	U	3	5	-	0/2/19/22	0/1/1/1
5	MAN	U	4	5	-	2/2/19/22	0/1/1/1
5	MAN	U	5	5	-	0/2/19/22	0/1/1/1
5	MAN	U	6	5	-	0/2/19/22	0/1/1/1
5	MAN	U	7	5	-	0/2/19/22	0/1/1/1
5	NAG	V	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	V	2	5	-	0/6/23/26	0/1/1/1
5	BMA	V	3	5	-	0/2/19/22	0/1/1/1
5	MAN	V	4	5	-	2/2/19/22	0/1/1/1
5	MAN	V	5	5	-	0/2/19/22	0/1/1/1
5	MAN	V	6	5	-	0/2/19/22	0/1/1/1
5	MAN	V	7	5	-	0/2/19/22	0/1/1/1
6	NAG	W	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	W	2	6	-	1/6/23/26	0/1/1/1
6	BMA	W	3	6	-	0/2/19/22	0/1/1/1
6	MAN	W	4	6	-	0/2/19/22	0/1/1/1
6	MAN	W	5	6	-	2/2/19/22	0/1/1/1
6	MAN	W	6	6	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	MAN	W	7	6	-	2/2/19/22	0/1/1/1
6	MAN	W	8	6	-	0/2/19/22	0/1/1/1
6	MAN	W	9	6	-	0/2/19/22	0/1/1/1
7	MAN	X	1	7	-	2/2/19/22	1/1/1/1
7	MAN	X	2	7	-	0/2/19/22	0/1/1/1
5	NAG	Y	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	Y	2	5	-	0/6/23/26	0/1/1/1
5	BMA	Y	3	5	-	0/2/19/22	0/1/1/1
5	MAN	Y	4	5	-	0/2/19/22	0/1/1/1
5	MAN	Y	5	5	-	0/2/19/22	0/1/1/1
5	MAN	Y	6	5	-	2/2/19/22	0/1/1/1
5	MAN	Y	7	5	-	0/2/19/22	0/1/1/1
5	NAG	Z	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	Z	2	5	-	0/6/23/26	0/1/1/1
5	BMA	Z	3	5	-	0/2/19/22	0/1/1/1
5	MAN	Z	4	5	-	1/2/19/22	0/1/1/1
5	MAN	Z	5	5	-	0/2/19/22	0/1/1/1
5	MAN	Z	6	5	-	0/2/19/22	0/1/1/1
5	MAN	Z	7	5	-	0/2/19/22	0/1/1/1
5	NAG	a	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	a	2	5	-	0/6/23/26	0/1/1/1
5	BMA	a	3	5	-	0/2/19/22	0/1/1/1
5	MAN	a	4	5	-	2/2/19/22	0/1/1/1
5	MAN	a	5	5	-	0/2/19/22	0/1/1/1
5	MAN	a	6	5	-	0/2/19/22	0/1/1/1
5	MAN	a	7	5	-	0/2/19/22	0/1/1/1
5	NAG	b	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	b	2	5	-	0/6/23/26	0/1/1/1
5	BMA	b	3	5	-	0/2/19/22	0/1/1/1
5	MAN	b	4	5	-	2/2/19/22	0/1/1/1
5	MAN	b	5	5	-	0/2/19/22	0/1/1/1
5	MAN	b	6	5	-	0/2/19/22	0/1/1/1
5	MAN	b	7	5	-	0/2/19/22	0/1/1/1
6	NAG	c	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	c	2	6	-	2/6/23/26	0/1/1/1
6	BMA	c	3	6	-	0/2/19/22	0/1/1/1
6	MAN	c	4	6	-	0/2/19/22	0/1/1/1
6	MAN	c	5	6	-	2/2/19/22	0/1/1/1
6	MAN	c	6	6	-	0/2/19/22	0/1/1/1
6	MAN	c	7	6	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	MAN	c	8	6	-	0/2/19/22	0/1/1/1
6	MAN	c	9	6	-	0/2/19/22	0/1/1/1
7	MAN	d	1	7	-	2/2/19/22	0/1/1/1
7	MAN	d	2	7	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 47 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	Y	1	NAG	C4-C3-C2	4.11	117.05	111.02
5	S	1	NAG	C4-C3-C2	3.87	116.69	111.02
5	M	1	NAG	C4-C3-C2	3.84	116.65	111.02
6	c	3	BMA	C1-C2-C3	3.69	115.01	109.64
6	c	1	NAG	O5-C1-C2	-3.53	105.83	111.29

There are no chirality outliers.

5 of 55 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	M	1	NAG	C1-C2-N2-C7
5	O	1	NAG	C8-C7-N2-C2
5	O	1	NAG	O7-C7-N2-C2
5	S	1	NAG	C1-C2-N2-C7
5	U	1	NAG	C8-C7-N2-C2

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	X	1	MAN	C1-C2-C3-C4-C5-O5

15 monomers are involved in 33 short contacts:

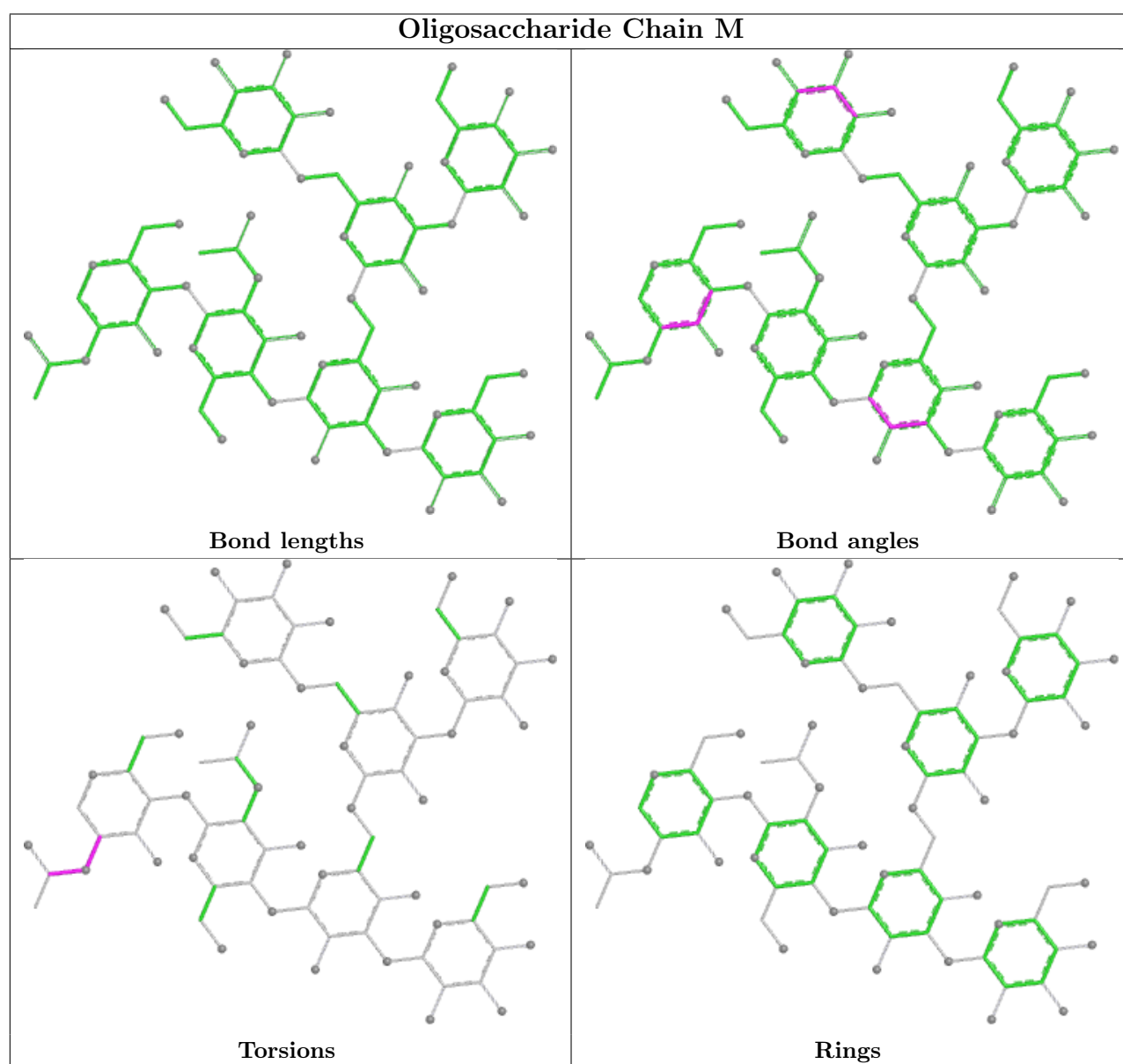
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	U	1	NAG	6	0
5	O	1	NAG	6	0
5	Y	1	NAG	1	0
5	a	1	NAG	4	0
5	S	1	NAG	1	0
6	c	2	NAG	3	0
7	R	1	MAN	2	0
6	Q	2	NAG	2	0

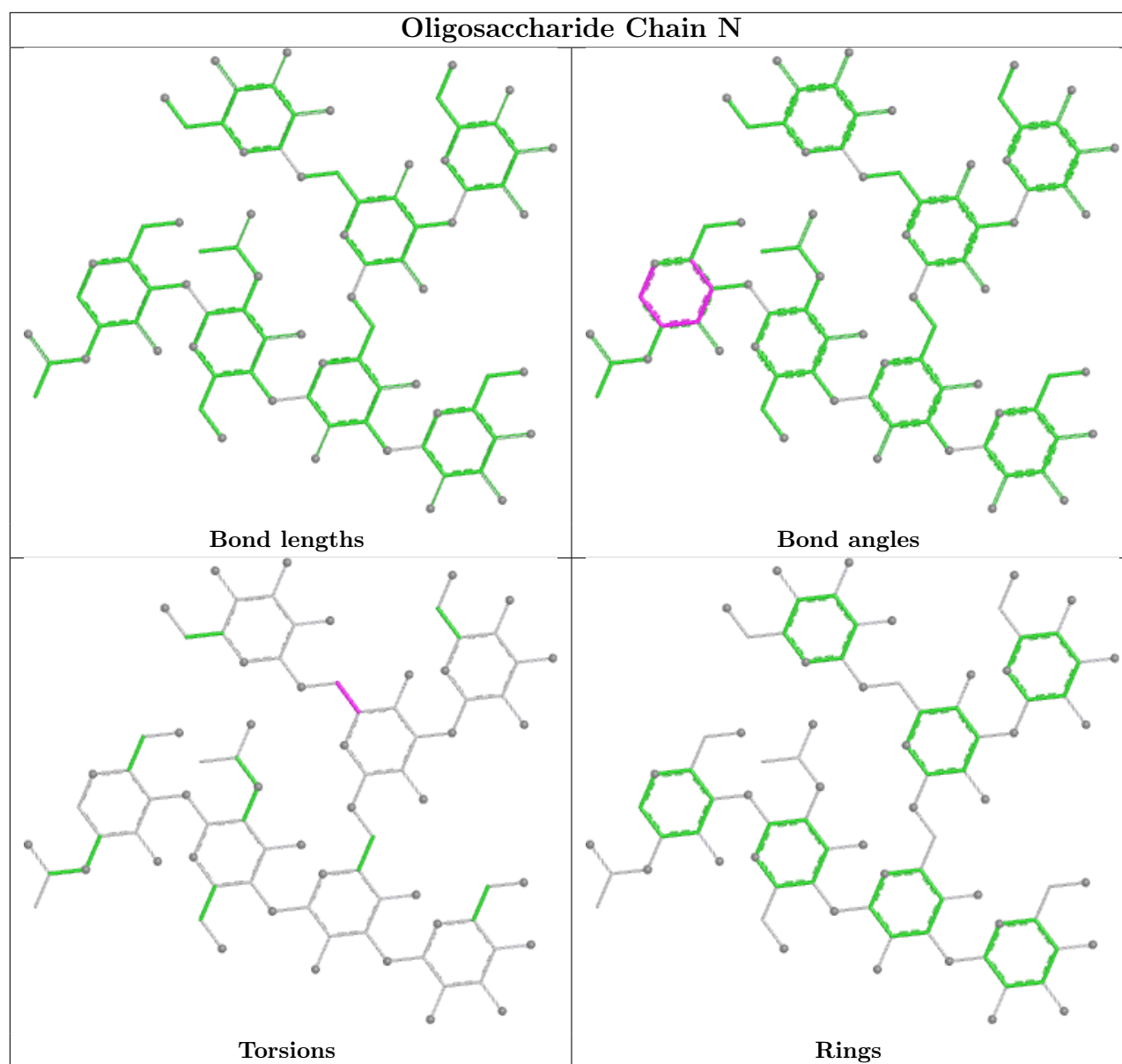
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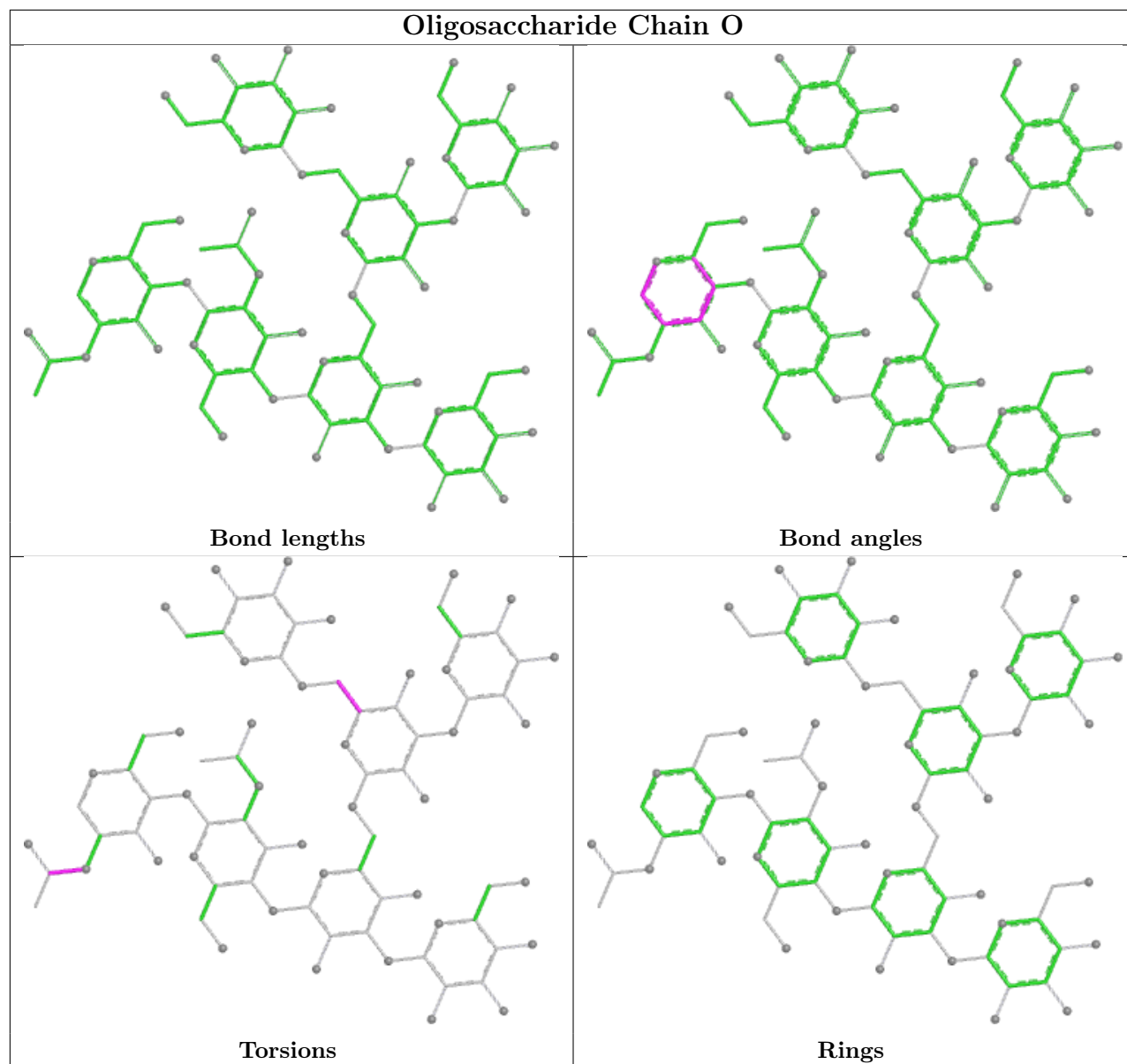
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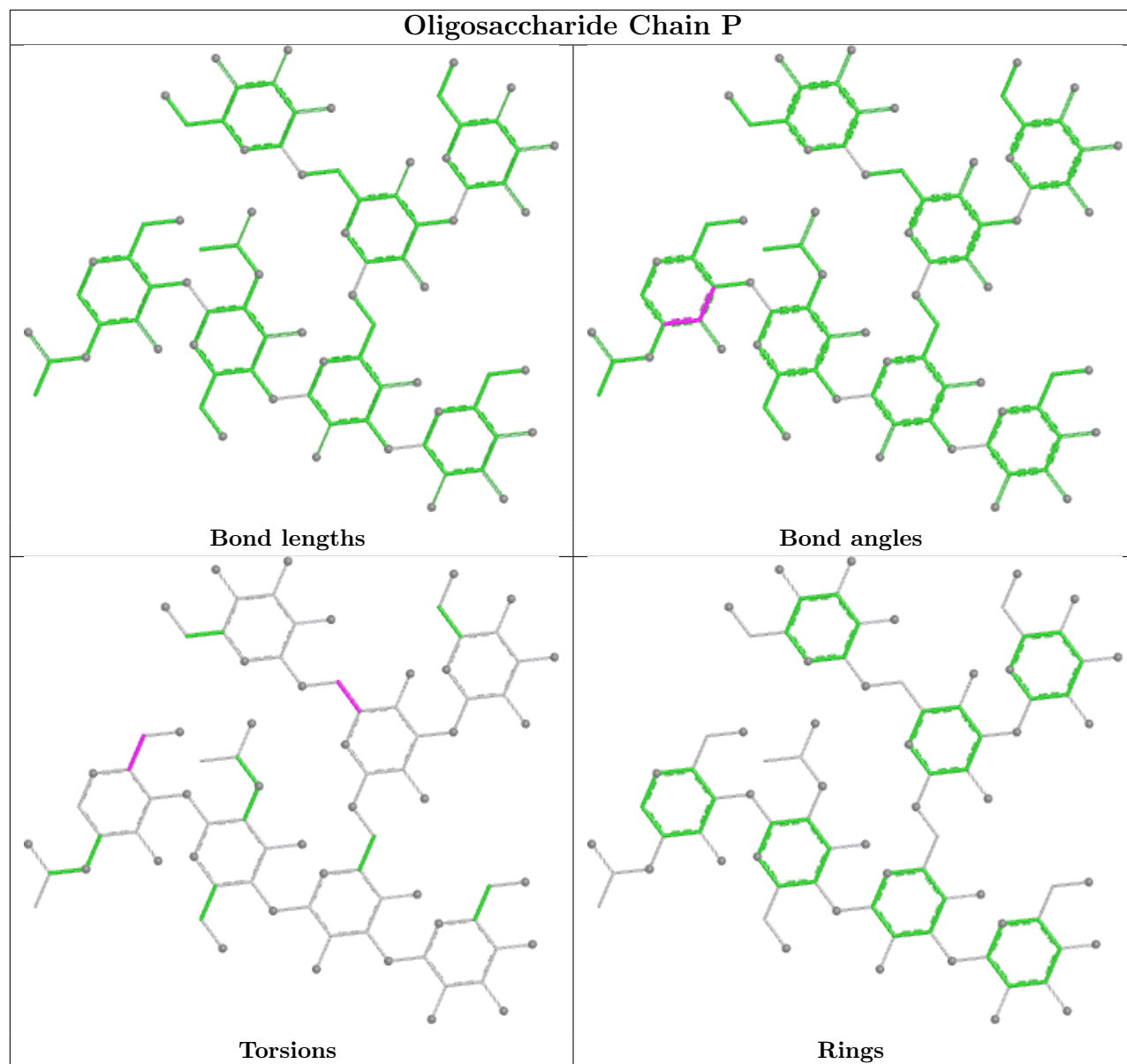
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	R	2	MAN	1	0
5	Y	6	MAN	2	0
7	d	2	MAN	1	0
6	c	1	NAG	1	0
5	M	1	NAG	1	0
7	d	1	MAN	1	0
7	X	2	MAN	1	0

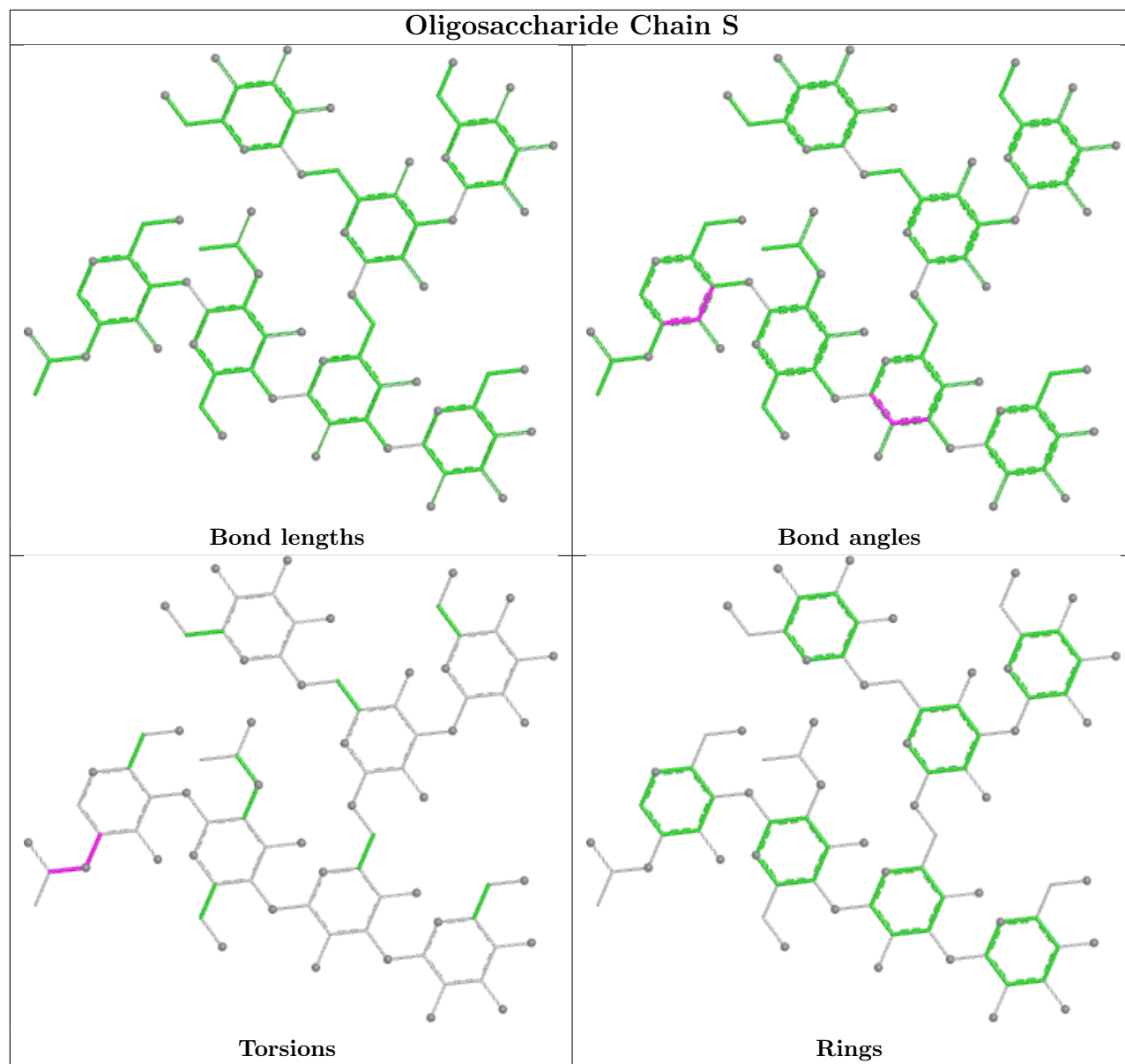
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

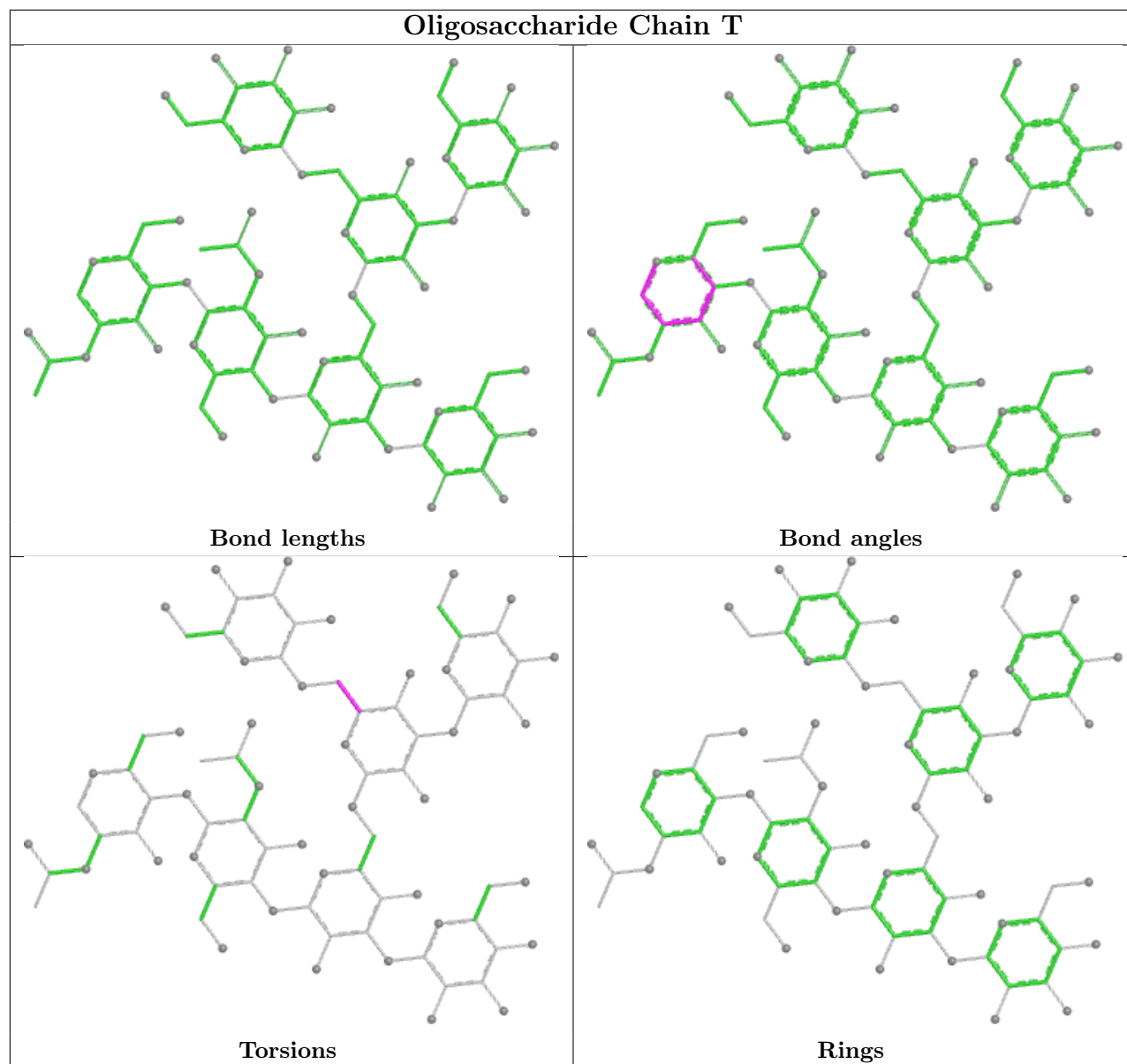


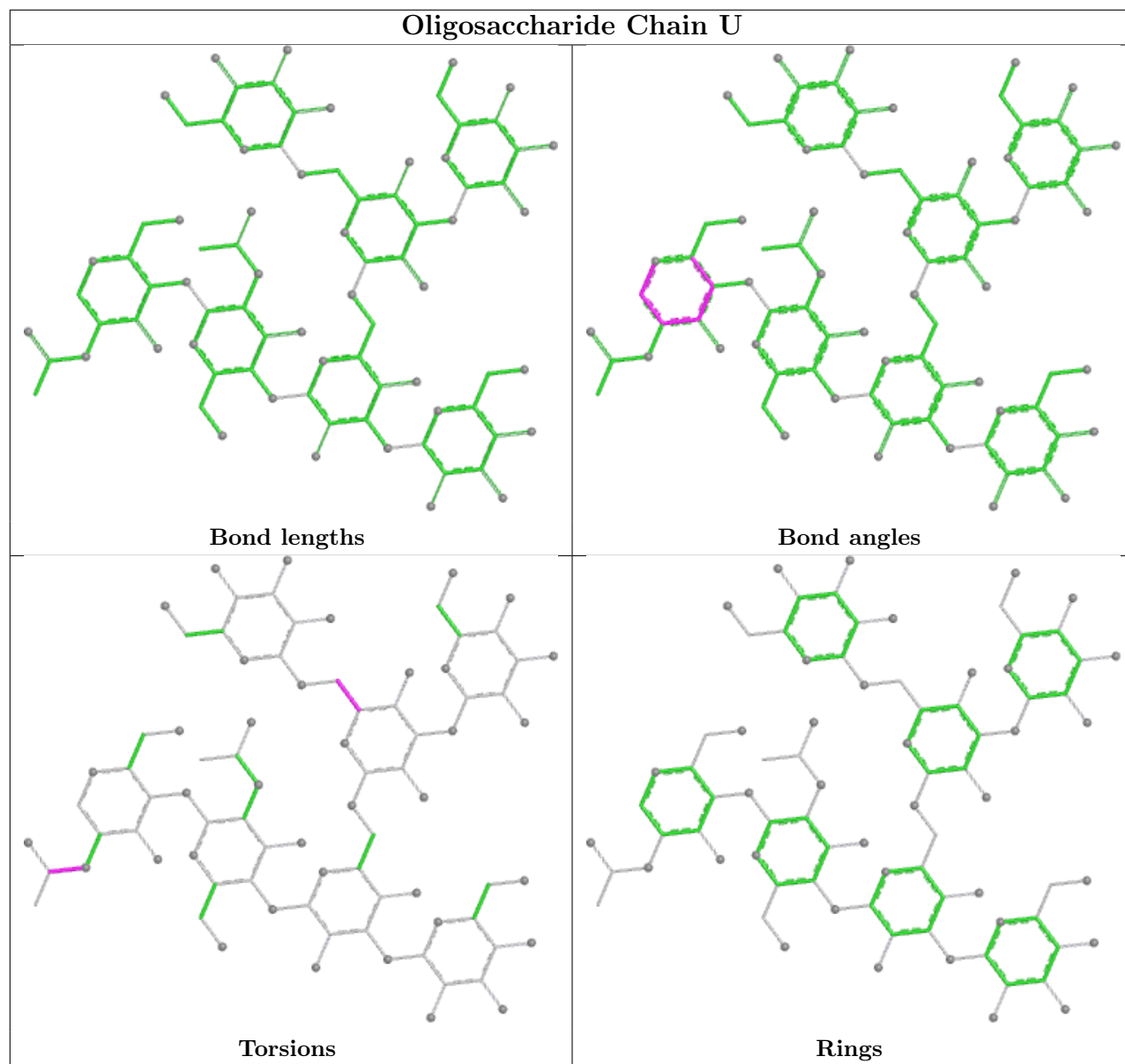


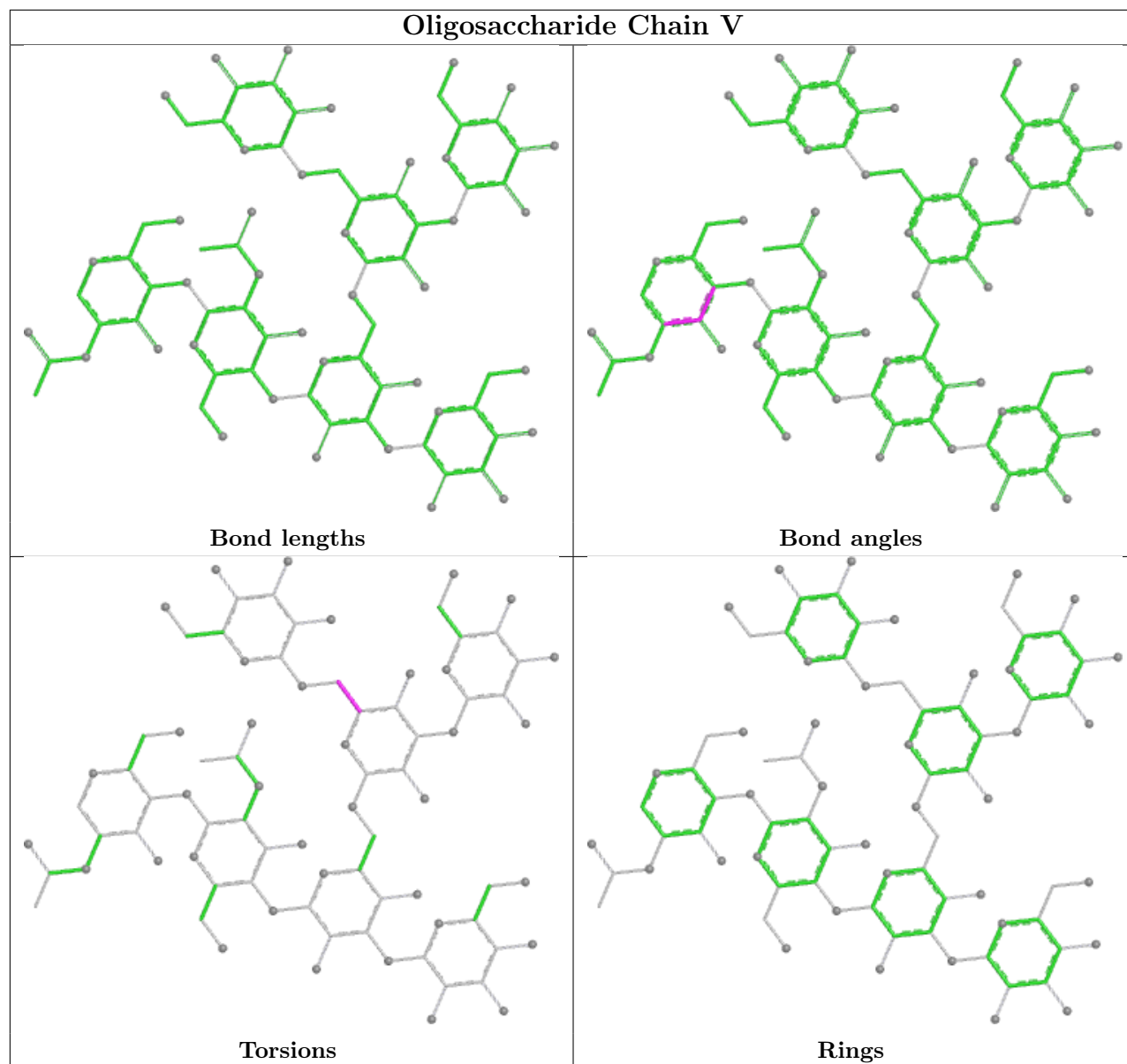


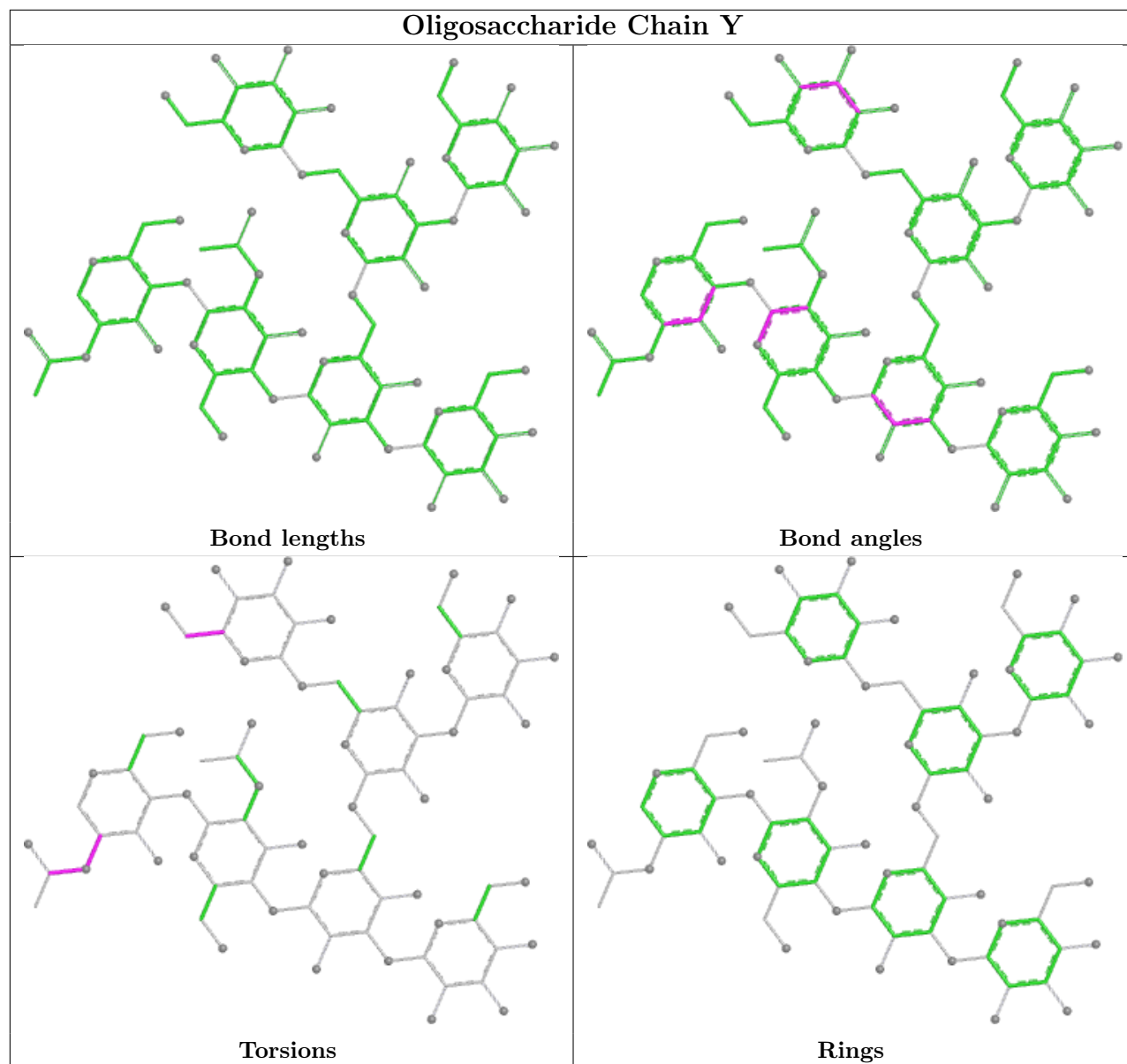


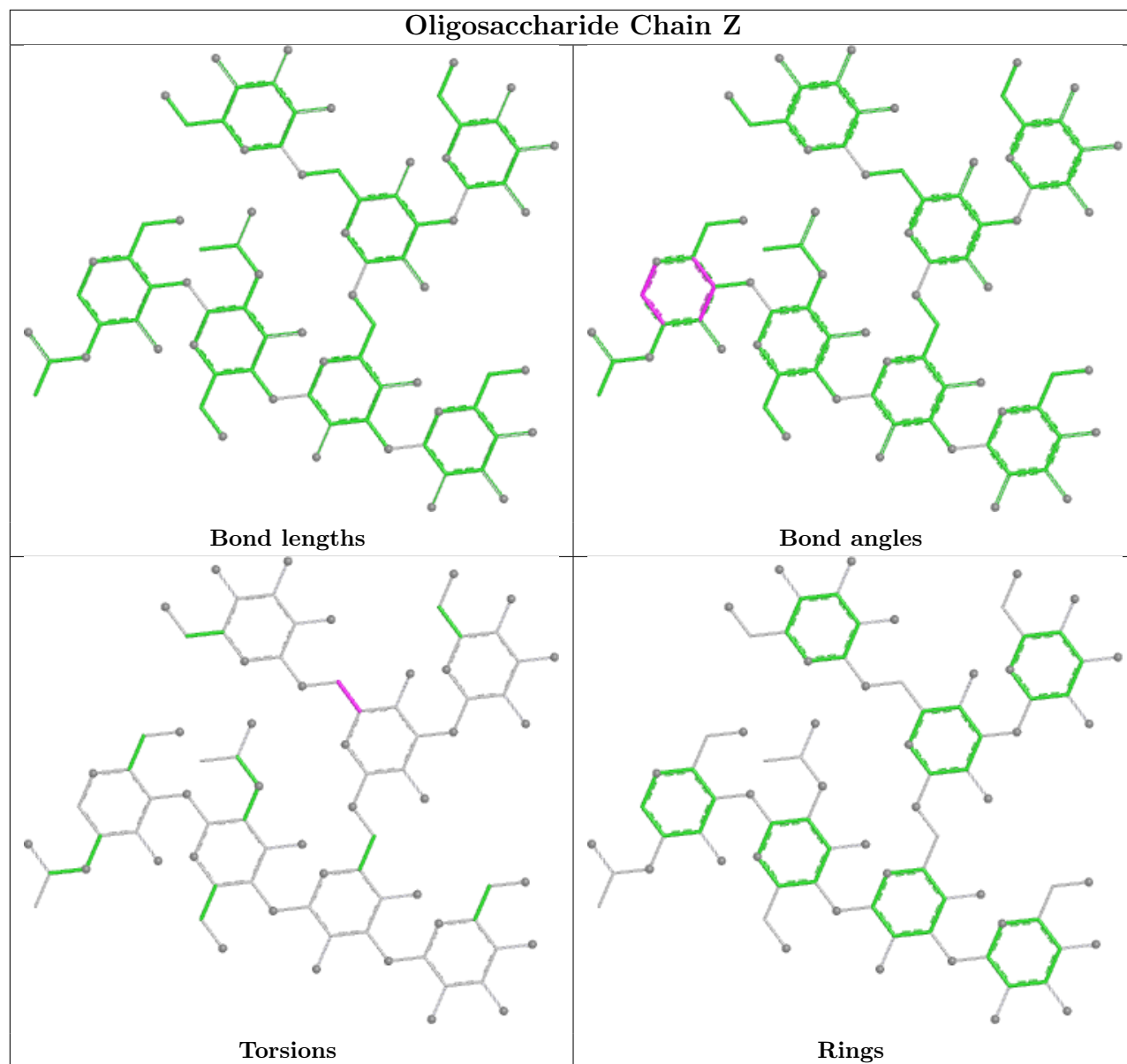


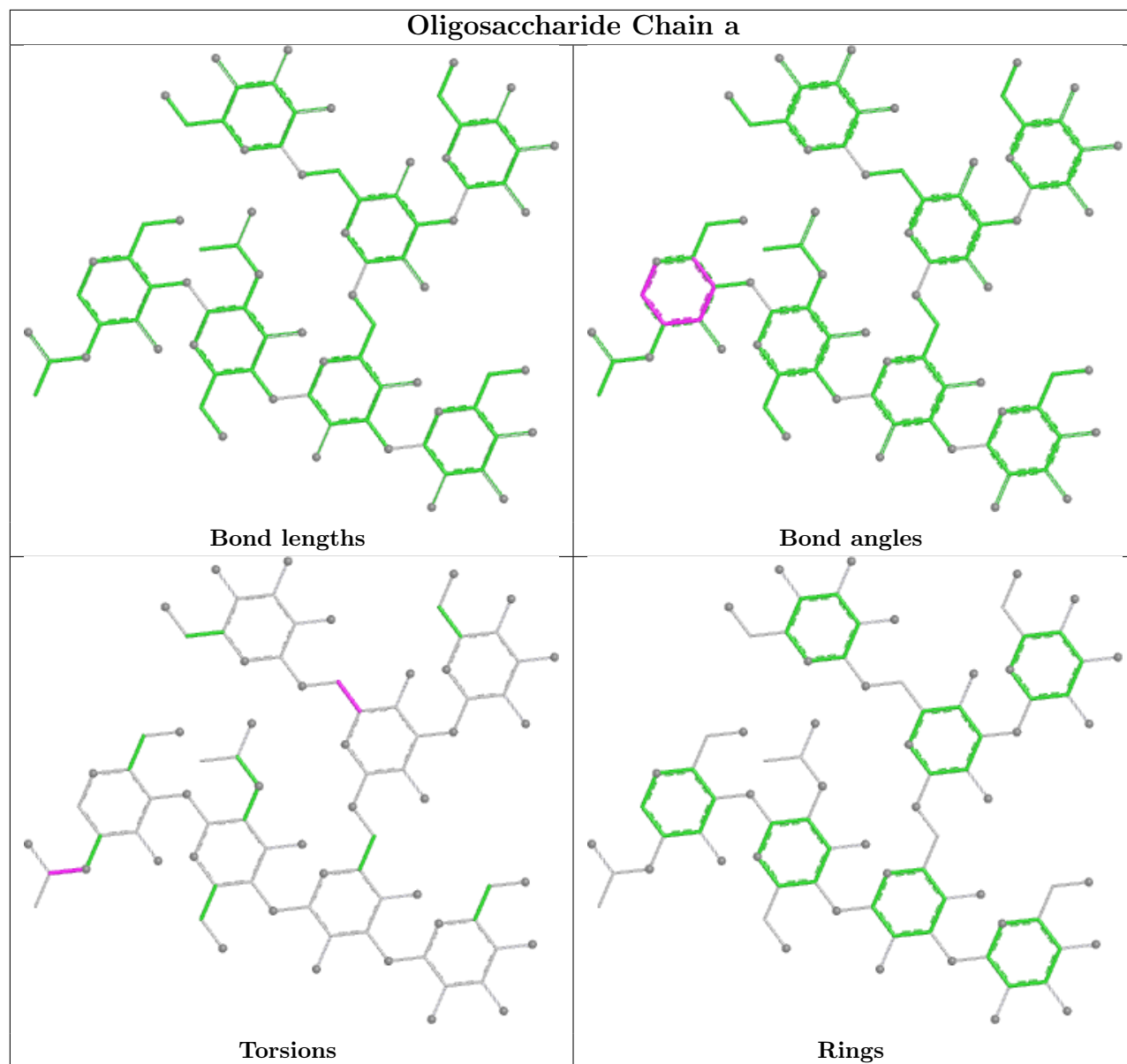


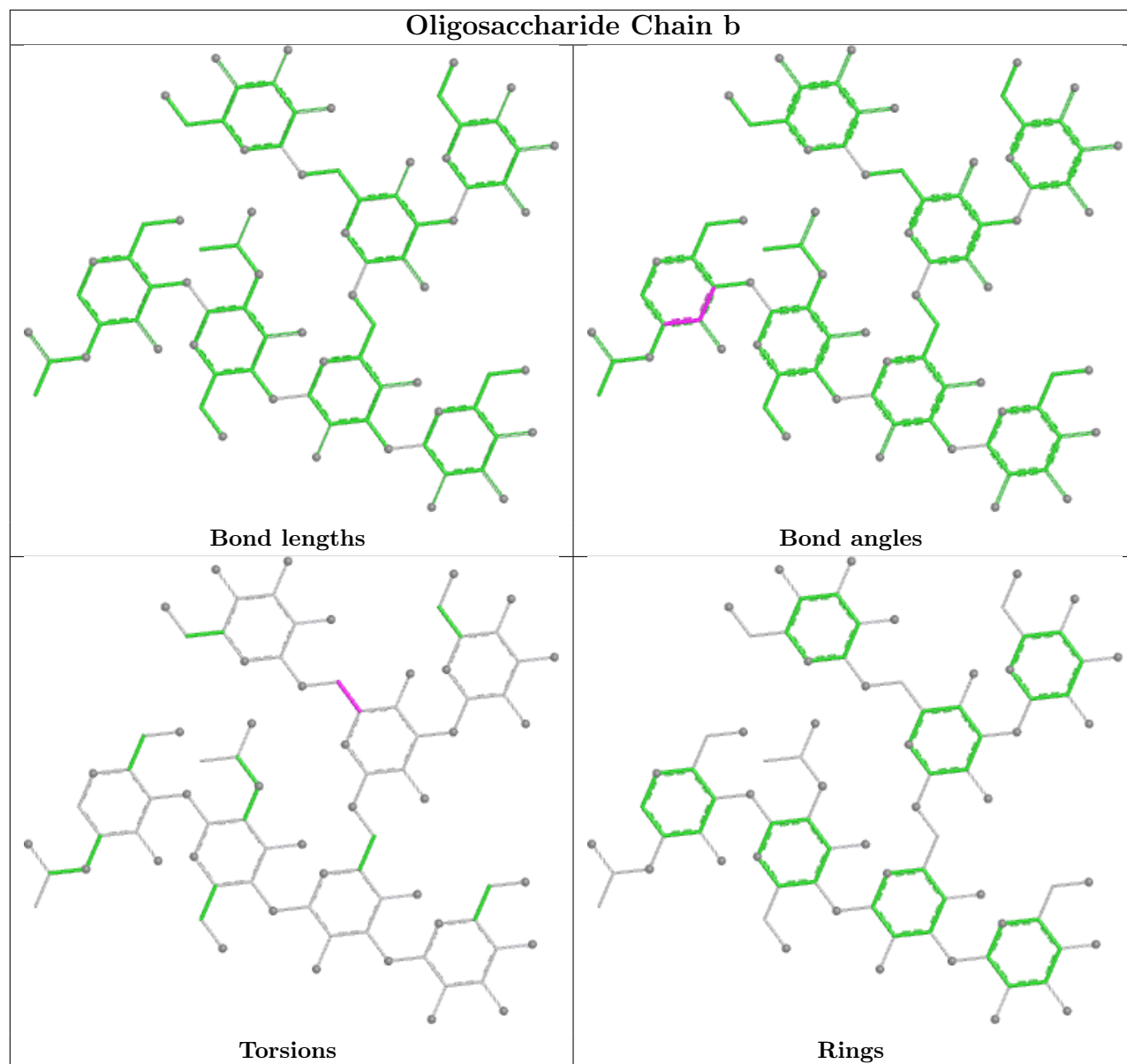


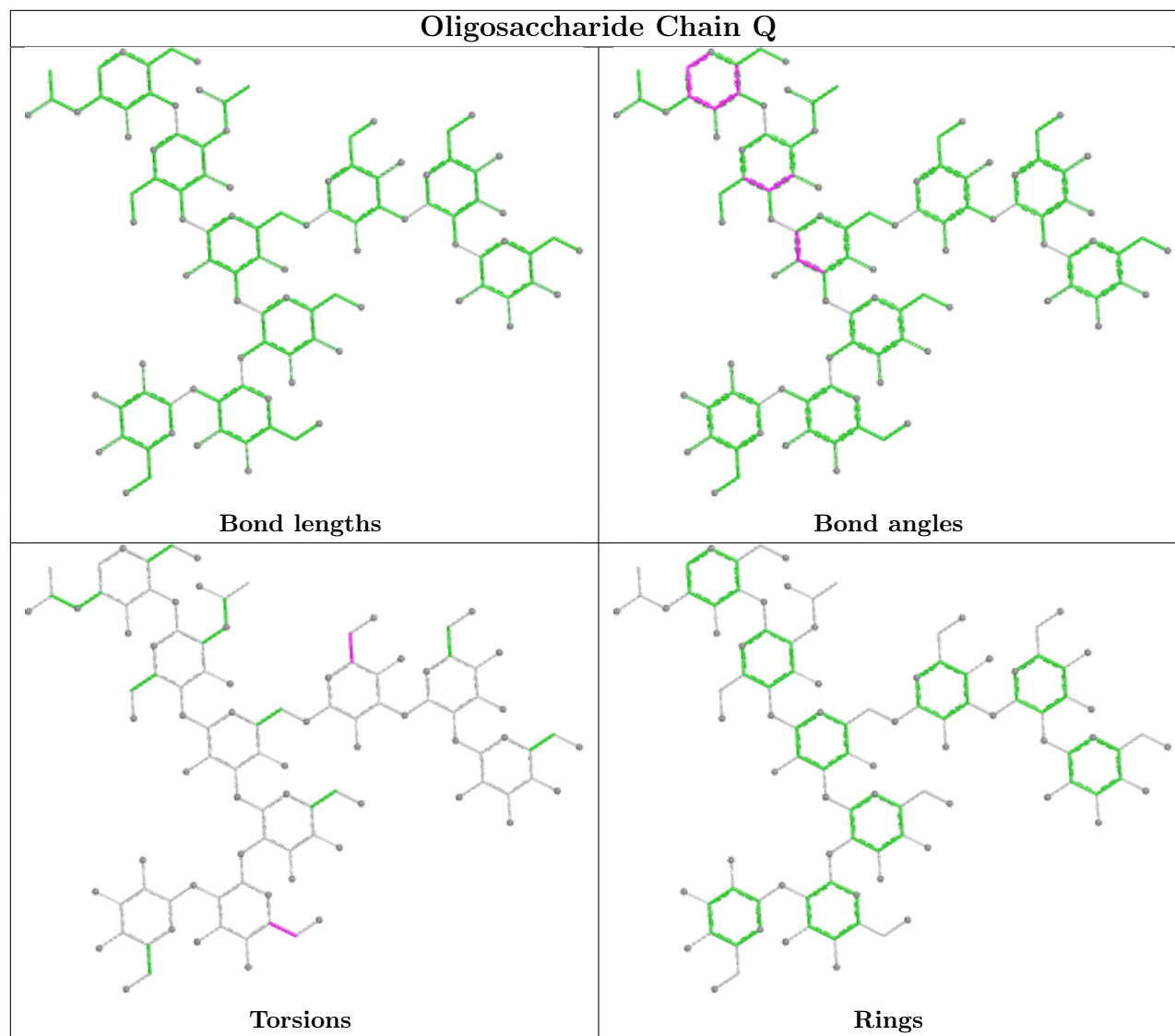


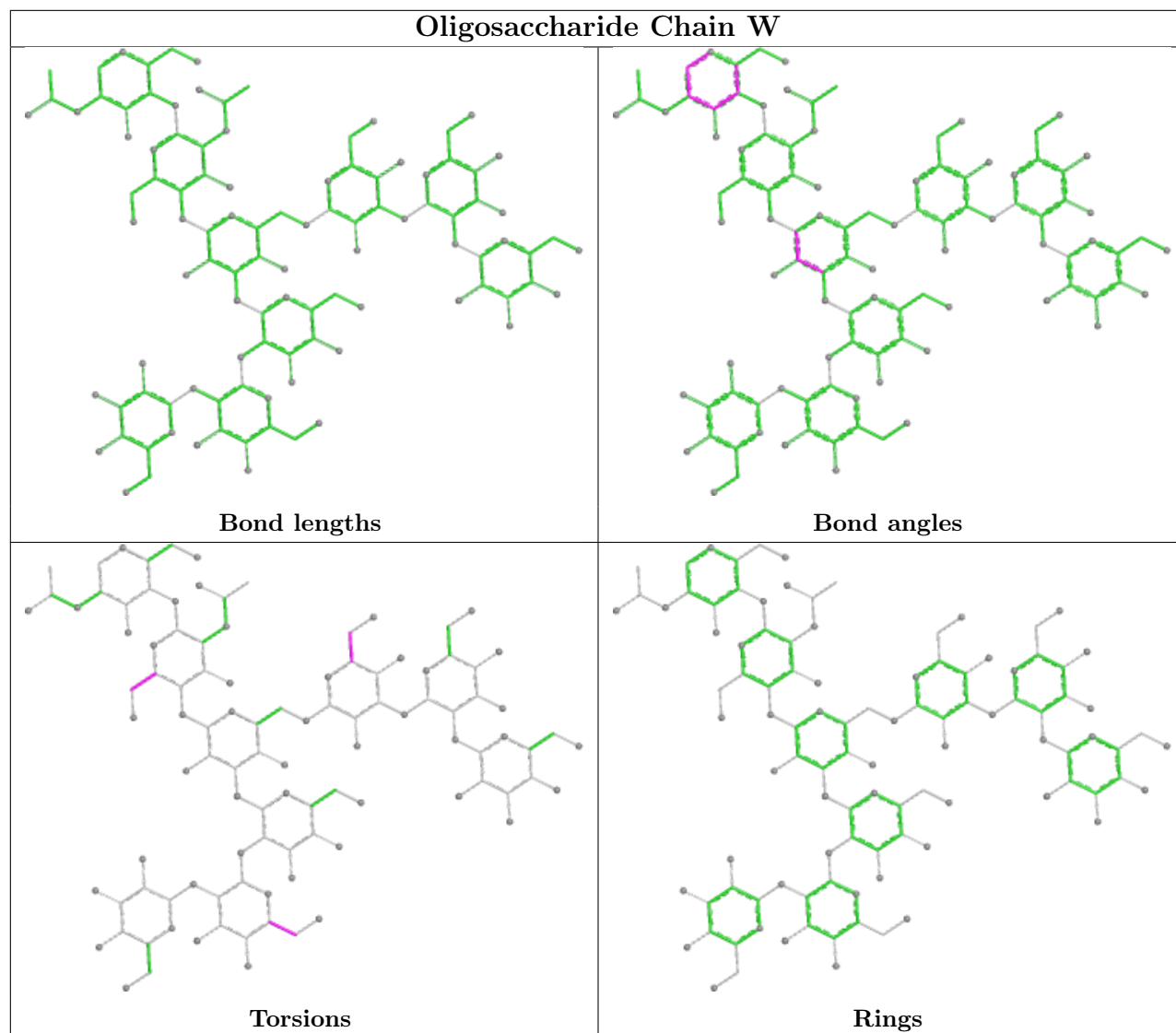


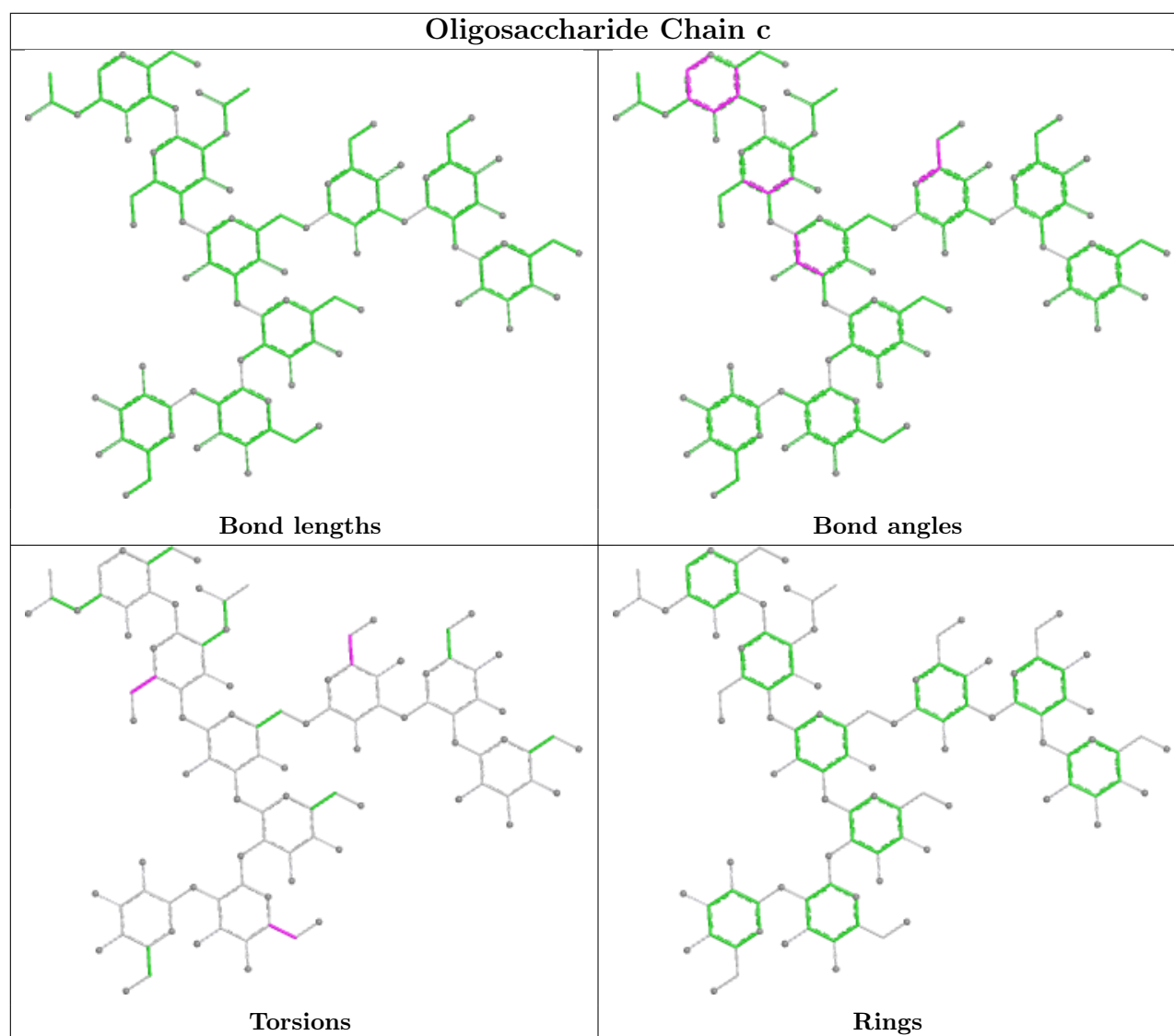


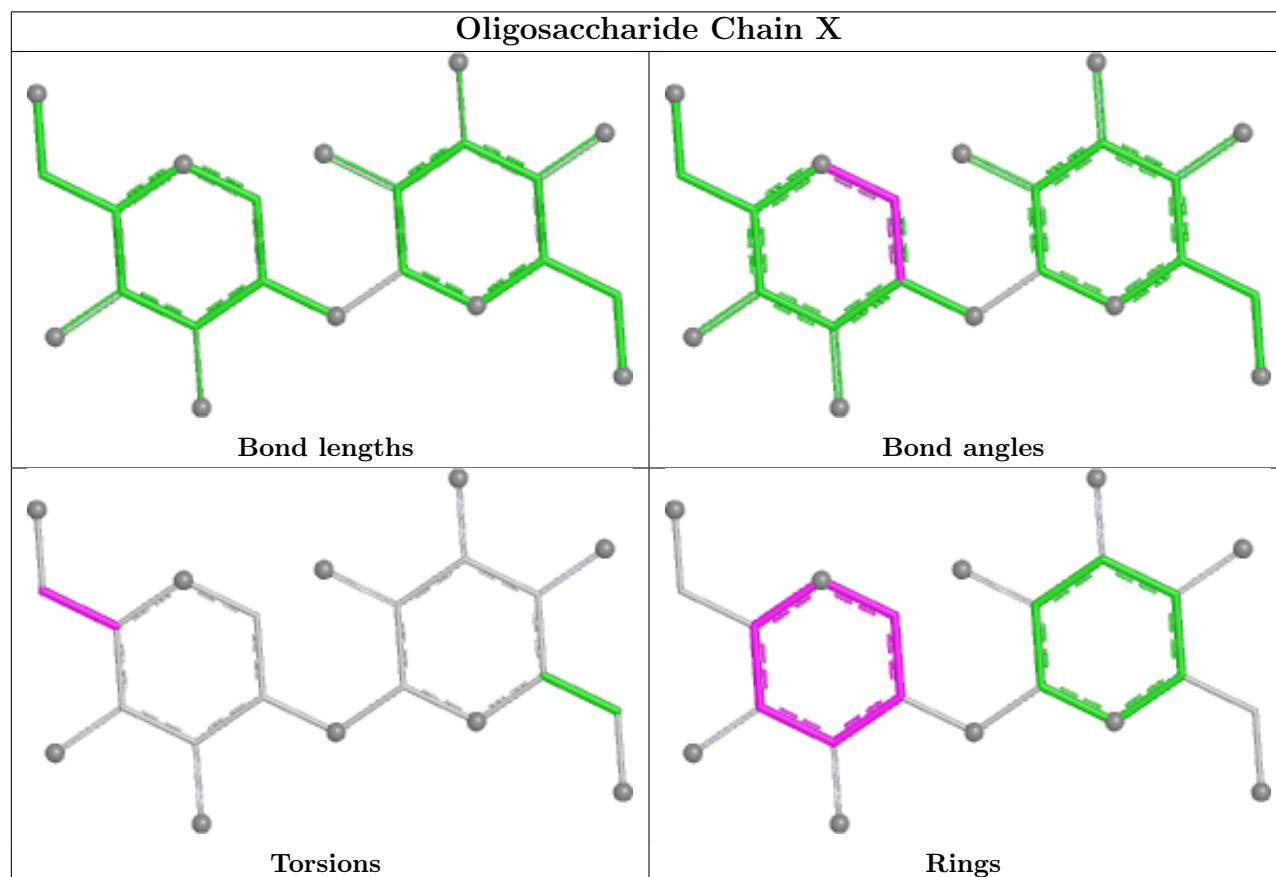
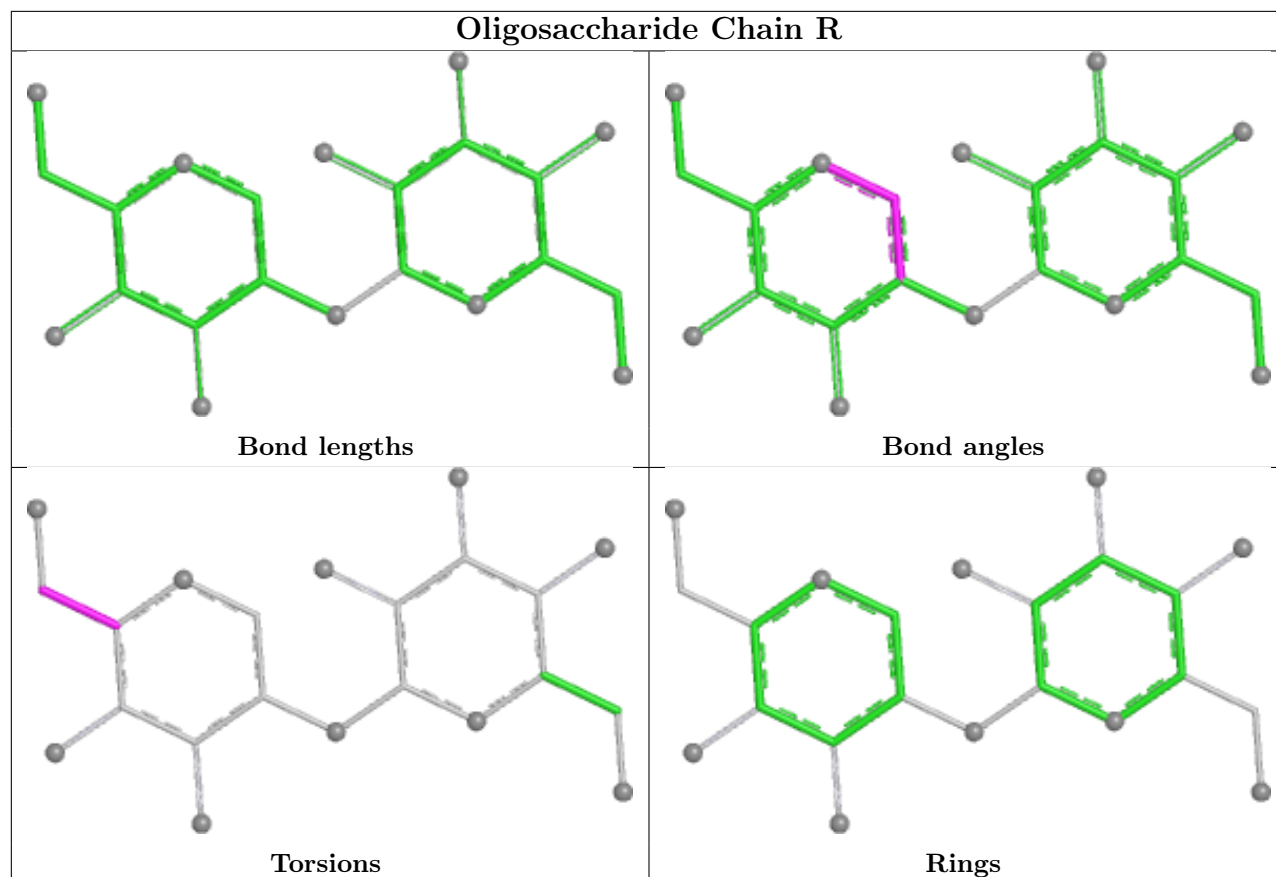


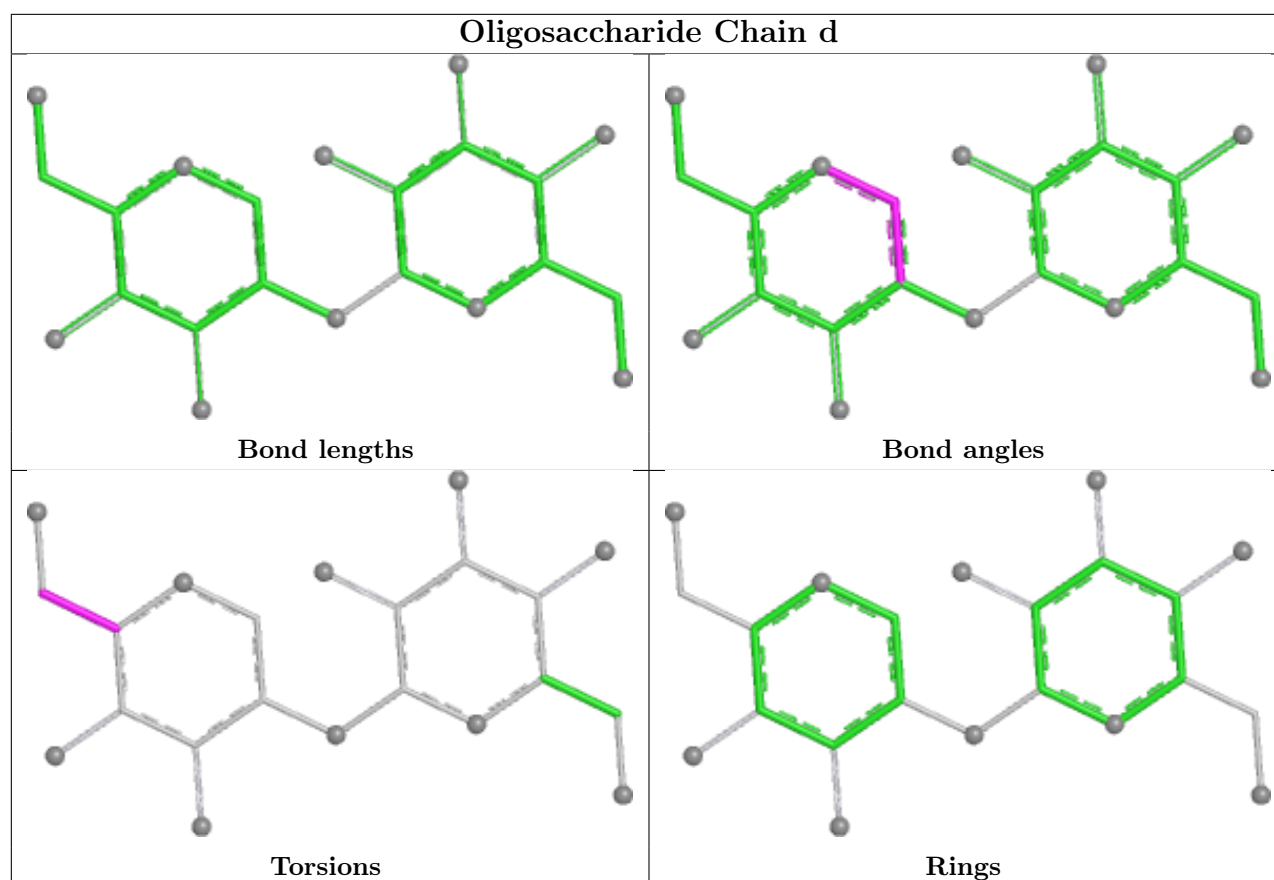












5.6 Ligand geometry [i](#)

30 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	A	1392	1	14,14,15	0.44	0	17,19,21	0.70	0
8	NAG	I	1276	1	14,14,15	0.45	0	17,19,21	1.03	1 (5%)
8	NAG	E	1088	1	14,14,15	0.55	0	17,19,21	0.73	0
8	NAG	E	1295	1	14,14,15	0.50	0	17,19,21	0.75	0
8	NAG	A	1197	1	14,14,15	0.45	0	17,19,21	0.88	1 (5%)
8	NAG	E	1197	1	14,14,15	0.45	0	17,19,21	0.80	0
8	NAG	I	1448	1	14,14,15	0.45	0	17,19,21	1.58	3 (17%)
8	NAG	E	1234	1	14,14,15	0.42	0	17,19,21	1.15	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	E	1392	1	14,14,15	0.47	0	17,19,21	0.67	0
8	NAG	A	1448	1	14,14,15	0.49	0	17,19,21	1.50	3 (17%)
8	NAG	I	1160	1	14,14,15	0.41	0	17,19,21	0.87	1 (5%)
8	NAG	E	1448	1	14,14,15	0.48	0	17,19,21	1.56	3 (17%)
8	NAG	E	1355	1	14,14,15	0.49	0	17,19,21	0.71	0
8	NAG	I	1392	1	14,14,15	0.42	0	17,19,21	0.75	0
8	NAG	I	1088	1	14,14,15	0.50	0	17,19,21	0.86	1 (5%)
8	NAG	A	1160	1	14,14,15	0.48	0	17,19,21	0.75	0
8	NAG	A	1088	1	14,14,15	0.49	0	17,19,21	0.87	1 (5%)
8	NAG	A	1295	1	14,14,15	0.53	0	17,19,21	0.71	0
8	NAG	I	1355	1	14,14,15	0.46	0	17,19,21	0.85	1 (5%)
8	NAG	A	1276	1	14,14,15	0.40	0	17,19,21	1.18	2 (11%)
8	NAG	E	1160	1	14,14,15	0.49	0	17,19,21	0.69	0
8	NAG	A	1234	1	14,14,15	0.43	0	17,19,21	1.19	2 (11%)
8	NAG	A	1355	1	14,14,15	0.49	0	17,19,21	0.74	0
8	NAG	E	1386	1	14,14,15	0.58	0	17,19,21	0.72	0
8	NAG	E	1276	1	14,14,15	0.44	0	17,19,21	1.25	3 (17%)
8	NAG	I	1234	1	14,14,15	0.45	0	17,19,21	1.18	2 (11%)
8	NAG	I	1295	1	14,14,15	0.51	0	17,19,21	0.82	1 (5%)
8	NAG	I	1386	1	14,14,15	0.54	0	17,19,21	0.79	1 (5%)
8	NAG	I	1197	1	14,14,15	0.49	0	17,19,21	0.80	0
8	NAG	A	1386	1	14,14,15	0.58	0	17,19,21	0.75	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	A	1392	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1276	1	-	0/6/23/26	0/1/1/1
8	NAG	E	1088	1	-	0/6/23/26	0/1/1/1
8	NAG	E	1295	1	-	0/6/23/26	0/1/1/1
8	NAG	A	1197	1	-	0/6/23/26	0/1/1/1
8	NAG	E	1197	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1448	1	-	4/6/23/26	0/1/1/1
8	NAG	E	1234	1	-	0/6/23/26	0/1/1/1
8	NAG	E	1392	1	-	0/6/23/26	0/1/1/1
8	NAG	A	1448	1	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	I	1160	1	-	2/6/23/26	0/1/1/1
8	NAG	E	1448	1	-	4/6/23/26	0/1/1/1
8	NAG	E	1355	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1392	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1088	1	-	0/6/23/26	0/1/1/1
8	NAG	A	1160	1	-	2/6/23/26	0/1/1/1
8	NAG	A	1088	1	-	0/6/23/26	0/1/1/1
8	NAG	A	1295	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1355	1	-	0/6/23/26	0/1/1/1
8	NAG	A	1276	1	-	0/6/23/26	0/1/1/1
8	NAG	E	1160	1	-	2/6/23/26	0/1/1/1
8	NAG	A	1234	1	-	0/6/23/26	0/1/1/1
8	NAG	A	1355	1	-	0/6/23/26	0/1/1/1
8	NAG	E	1386	1	-	1/6/23/26	0/1/1/1
8	NAG	E	1276	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1234	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1295	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1386	1	-	0/6/23/26	0/1/1/1
8	NAG	I	1197	1	-	0/6/23/26	0/1/1/1
8	NAG	A	1386	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	I	1448	NAG	C1-O5-C5	3.37	116.71	112.19
8	E	1448	NAG	C1-O5-C5	3.35	116.68	112.19
8	A	1448	NAG	C1-O5-C5	3.13	116.38	112.19
8	A	1276	NAG	C1-O5-C5	3.03	116.25	112.19
8	E	1276	NAG	C1-O5-C5	3.01	116.21	112.19

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	E	1160	NAG	C8-C7-N2-C2
8	A	1448	NAG	O5-C5-C6-O6
8	E	1448	NAG	O5-C5-C6-O6
8	A	1160	NAG	C8-C7-N2-C2
8	E	1160	NAG	O7-C7-N2-C2

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	E	1295	NAG	2	0
8	A	1295	NAG	2	0
8	A	1276	NAG	1	0
8	E	1276	NAG	1	0
8	I	1295	NAG	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
4	L	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	L	100(R):VAL	C	101:TRP	N	1.66

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	420/475 (88%)	1.31	117 (27%) 1 5	7, 227, 434, 596	0
1	E	420/475 (88%)	1.43	114 (27%) 1 5	16, 221, 423, 559	0
1	I	420/475 (88%)	1.50	126 (30%) 1 5	25, 216, 434, 550	0
2	B	0/78	-	-	-	-
2	F	0/78	-	-	-	-
2	J	0/78	-	-	-	-
3	C	202/211 (95%)	1.25	50 (24%) 2 5	74, 234, 422, 563	0
3	G	202/211 (95%)	1.52	61 (30%) 1 4	43, 229, 392, 572	0
3	K	202/211 (95%)	1.48	60 (29%) 1 5	47, 291, 454, 533	0
4	D	226/235 (96%)	1.48	70 (30%) 1 4	67, 232, 488, 582	0
4	H	226/235 (96%)	1.56	76 (33%) 1 4	56, 232, 421, 563	0
4	L	226/235 (96%)	1.49	66 (29%) 1 5	71, 287, 466, 582	0
All	All	2544/2997 (84%)	1.44	740 (29%) 1 5	7, 236, 440, 596	0

The worst 5 of 740 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	100(A)	ILE	13.8
4	H	100(A)	ILE	9.1
1	I	236	THR	8.7
1	E	313	GLY	8.6
4	H	72	ASP	7.9

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	MAN	X	1	11/12	-0.42	0.34	427,427,427,427	0
5	MAN	Y	4	11/12	-0.34	0.21	550,550,550,550	0
7	MAN	R	1	11/12	-0.24	0.43	550,550,550,550	0
5	NAG	Y	1	14/15	-0.15	0.29	530,530,530,530	0
5	NAG	Y	2	14/15	-0.09	0.23	460,460,460,460	0
5	MAN	S	4	11/12	-0.05	0.15	550,550,550,550	0
5	MAN	M	4	11/12	-0.03	0.16	517,517,517,517	0
5	MAN	S	6	11/12	0.06	0.19	497,497,497,497	0
7	MAN	R	2	11/12	0.09	0.42	489,489,489,489	0
5	MAN	P	6	11/12	0.09	0.16	470,470,470,470	0
6	MAN	Q	9	11/12	0.13	0.17	303,303,303,303	0
5	MAN	U	6	11/12	0.15	0.17	515,515,515,515	0
5	MAN	Y	6	11/12	0.16	0.17	494,494,494,494	0
5	MAN	Z	7	11/12	0.16	0.34	550,550,550,550	0
5	MAN	Z	6	11/12	0.17	0.18	428,428,428,428	0
5	MAN	P	4	11/12	0.18	0.10	483,483,483,483	0
5	MAN	V	6	11/12	0.19	0.14	454,454,454,454	0
5	BMA	V	3	11/12	0.26	0.15	492,492,492,492	0
5	MAN	V	4	11/12	0.27	0.13	471,471,471,471	0
5	MAN	N	7	11/12	0.28	0.37	498,498,498,498	0
5	MAN	S	5	11/12	0.29	0.16	550,550,550,550	0
5	MAN	O	6	11/12	0.30	0.17	455,455,455,455	0
5	MAN	U	4	11/12	0.31	0.23	550,550,550,550	0
5	NAG	M	1	14/15	0.31	0.22	439,439,439,439	0
6	MAN	W	7	11/12	0.31	0.15	418,418,418,418	0
5	MAN	M	7	11/12	0.32	0.15	451,451,451,451	0
5	MAN	M	5	11/12	0.32	0.15	481,481,481,481	0
6	MAN	Q	7	11/12	0.33	0.12	309,309,309,309	0
5	NAG	S	1	14/15	0.33	0.28	550,550,550,550	0
5	BMA	S	3	11/12	0.34	0.16	500,500,500,500	0
6	MAN	Q	5	11/12	0.35	0.27	302,302,302,302	0
5	MAN	N	6	11/12	0.37	0.16	464,464,464,464	0
6	BMA	W	3	11/12	0.38	0.13	400,400,400,400	0
6	MAN	W	5	11/12	0.38	0.27	330,330,330,330	0
5	MAN	Y	5	11/12	0.39	0.19	477,477,477,477	0
5	MAN	Y	7	11/12	0.40	0.26	520,520,520,520	0
5	MAN	N	4	11/12	0.40	0.21	550,550,550,550	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MAN	Z	4	11/12	0.41	0.14	550,550,550,550	0
6	MAN	W	6	11/12	0.41	0.31	452,452,452,452	0
5	MAN	S	7	11/12	0.43	0.19	518,518,518,518	0
5	MAN	M	6	11/12	0.46	0.16	434,434,434,434	0
5	BMA	Y	3	11/12	0.47	0.13	402,402,402,402	0
5	MAN	O	4	11/12	0.47	0.20	550,550,550,550	0
5	MAN	P	7	11/12	0.48	0.29	447,447,447,447	0
5	MAN	T	5	11/12	0.49	0.19	422,422,422,422	0
7	MAN	X	2	11/12	0.49	0.40	545,545,545,545	0
5	MAN	Z	5	11/12	0.50	0.18	410,410,410,410	0
5	BMA	O	3	11/12	0.50	0.15	506,506,506,506	0
5	MAN	T	7	11/12	0.52	0.32	542,542,542,542	0
5	MAN	T	6	11/12	0.53	0.16	446,446,446,446	0
6	MAN	W	9	11/12	0.53	0.09	335,335,335,335	0
5	BMA	P	3	11/12	0.55	0.13	455,455,455,455	0
5	NAG	U	1	14/15	0.55	0.38	408,408,408,408	0
5	MAN	U	5	11/12	0.57	0.19	550,550,550,550	0
5	MAN	T	4	11/12	0.57	0.15	550,550,550,550	0
5	BMA	U	3	11/12	0.58	0.12	340,340,340,340	0
5	MAN	O	5	11/12	0.58	0.18	468,468,468,468	0
5	MAN	V	5	11/12	0.59	0.21	495,495,495,495	0
5	BMA	Z	3	11/12	0.59	0.13	491,491,491,491	0
5	NAG	P	2	14/15	0.60	0.16	307,307,307,307	0
5	BMA	M	3	11/12	0.62	0.11	356,356,356,356	0
6	NAG	Q	1	14/15	0.65	0.32	366,366,366,366	0
5	BMA	N	3	11/12	0.66	0.10	483,483,483,483	0
6	MAN	W	4	11/12	0.66	0.34	550,550,550,550	0
5	MAN	N	5	11/12	0.67	0.14	329,329,329,329	0
5	NAG	O	1	14/15	0.70	0.40	423,423,423,423	0
5	BMA	T	3	11/12	0.71	0.10	466,466,466,466	0
6	MAN	Q	8	11/12	0.72	0.22	500,500,500,500	0
6	MAN	Q	4	11/12	0.72	0.30	550,550,550,550	0
5	NAG	V	2	14/15	0.73	0.16	335,335,335,335	0
5	NAG	a	1	14/15	-	-	437,437,437,437	0
5	NAG	a	2	14/15	-	-	231,231,231,231	0
5	BMA	a	3	11/12	-	-	297,297,297,297	0
5	MAN	a	4	11/12	-	-	545,545,545,545	0
5	MAN	a	5	11/12	-	-	507,507,507,507	0
5	MAN	a	6	11/12	-	-	532,532,532,532	0
5	MAN	a	7	11/12	-	-	197,197,197,197	0
5	NAG	b	1	14/15	-	-	276,276,276,276	0
5	NAG	b	2	14/15	-	-	382,382,382,382	0

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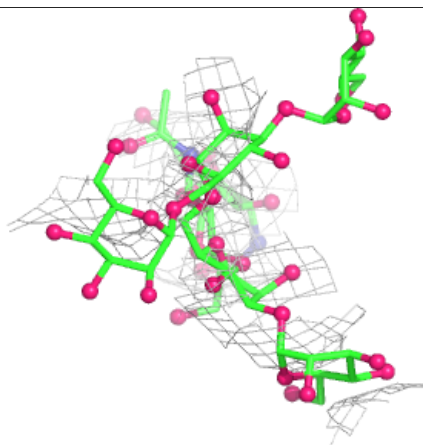
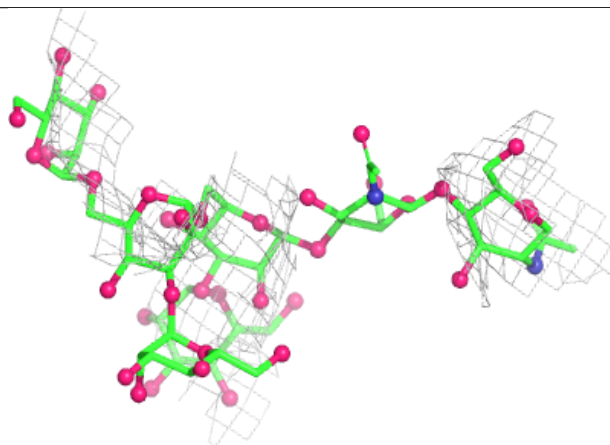
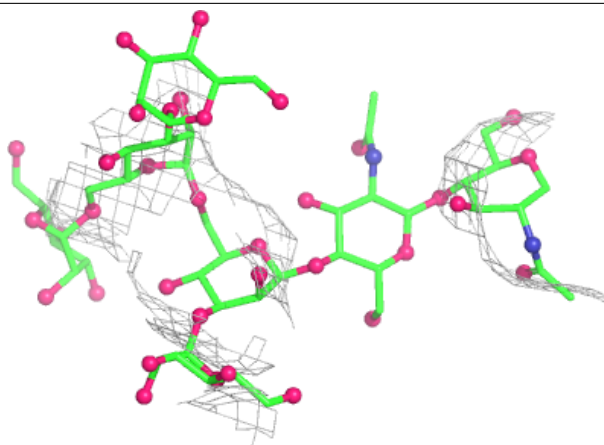
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	BMA	b	3	11/12	-	-	474,474,474,474	0
5	MAN	b	4	11/12	-	-	467,467,467,467	0
5	MAN	b	5	11/12	-	-	482,482,482,482	0
5	MAN	b	6	11/12	-	-	474,474,474,474	0
5	MAN	b	7	11/12	-	-	330,330,330,330	0
6	NAG	W	1	14/15	0.73	0.24	302,302,302,302	0
6	NAG	W	2	14/15	0.74	0.24	388,388,388,388	0
5	NAG	M	2	14/15	0.75	0.18	519,519,519,519	0
6	MAN	Q	6	11/12	0.75	0.21	422,422,422,422	0
6	BMA	Q	3	11/12	0.75	0.10	292,292,292,292	0
5	NAG	P	1	14/15	0.75	0.21	202,202,202,202	0
6	MAN	W	8	11/12	0.75	0.23	491,491,491,491	0
5	MAN	V	7	11/12	0.76	0.18	377,377,377,377	0
5	NAG	U	2	14/15	0.77	0.29	310,310,310,310	0
5	MAN	P	5	11/12	0.78	0.18	418,418,418,418	0
5	NAG	S	2	14/15	0.79	0.15	488,488,488,488	0
5	NAG	T	1	14/15	0.82	0.22	186,186,186,186	0
6	NAG	Q	2	14/15	0.83	0.13	258,258,258,258	0
5	MAN	O	7	11/12	0.84	0.10	229,229,229,229	0
5	MAN	U	7	11/12	0.84	0.10	195,195,195,195	0
5	NAG	N	1	14/15	0.84	0.28	356,356,356,356	0
5	NAG	V	1	14/15	0.85	0.25	242,242,242,242	0
5	NAG	O	2	14/15	0.89	0.27	410,410,410,410	0
6	NAG	c	1	14/15	-	-	292,292,292,292	0
6	NAG	c	2	14/15	-	-	213,213,213,213	0
6	BMA	c	3	11/12	-	-	345,345,345,345	0
6	MAN	c	4	11/12	-	-	550,550,550,550	0
6	MAN	c	5	11/12	-	-	277,277,277,277	0
6	MAN	c	6	11/12	-	-	436,436,436,436	0
6	MAN	c	7	11/12	-	-	382,382,382,382	0
6	MAN	c	8	11/12	-	-	507,507,507,507	0
6	MAN	c	9	11/12	-	-	362,362,362,362	0
5	NAG	Z	2	14/15	0.90	0.14	375,375,375,375	0
5	NAG	Z	1	14/15	0.90	0.22	337,337,337,337	0
5	NAG	N	2	14/15	0.91	0.14	310,310,310,310	0
5	NAG	T	2	14/15	0.97	0.13	301,301,301,301	0
7	MAN	d	1	11/12	-	-	408,408,408,408	0
7	MAN	d	2	11/12	-	-	550,550,550,550	0

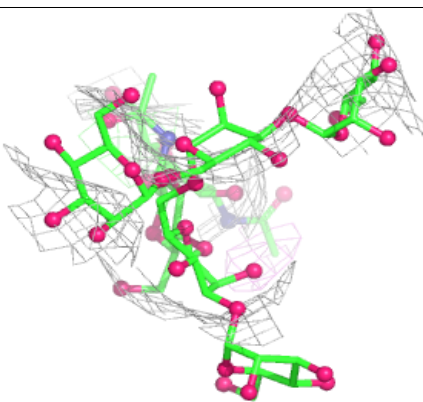
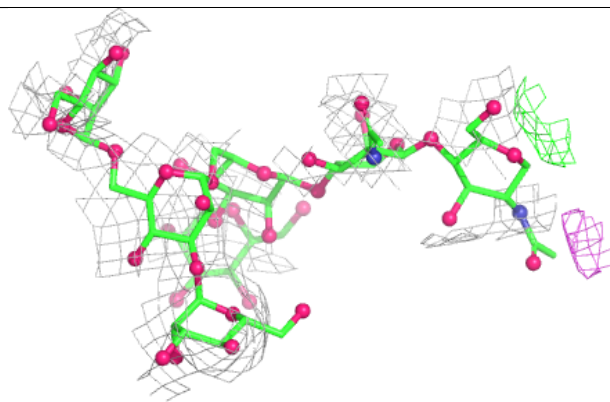
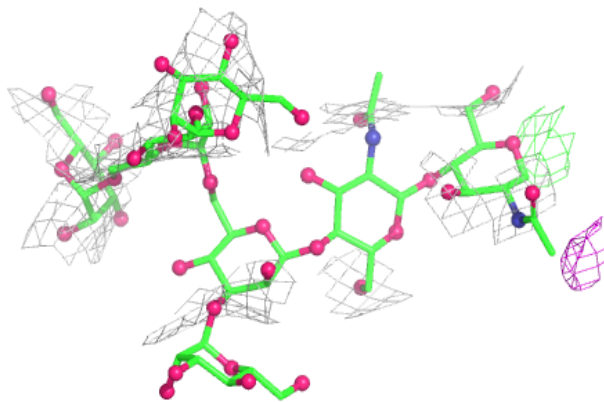
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain M:

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

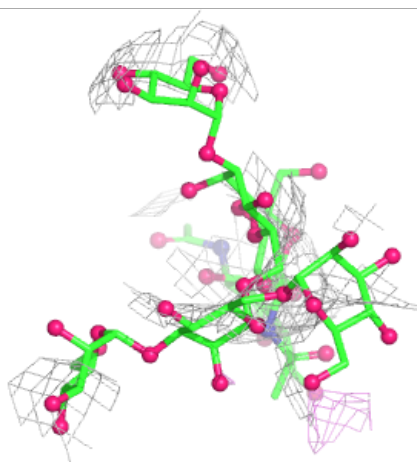
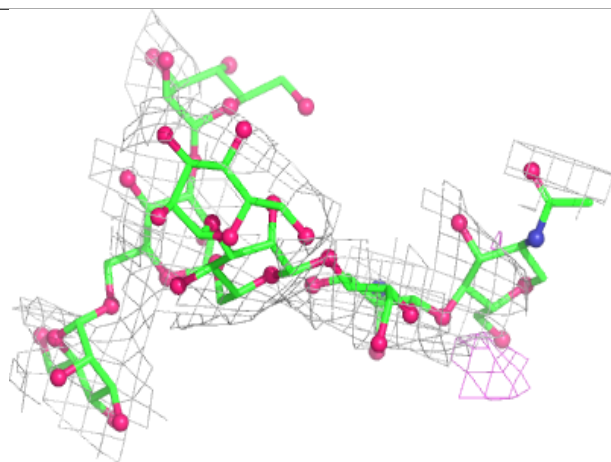
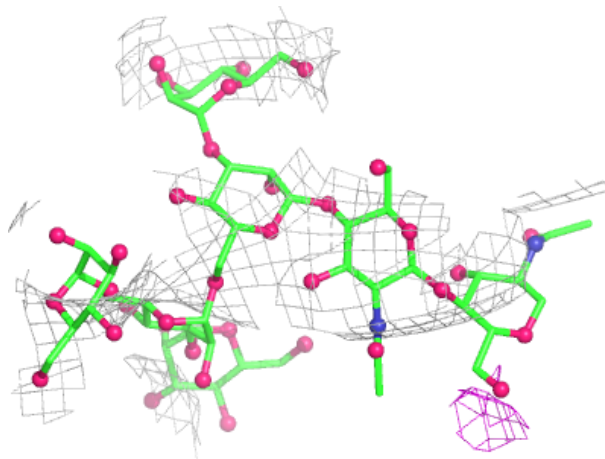
**Electron density around Chain N:**

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



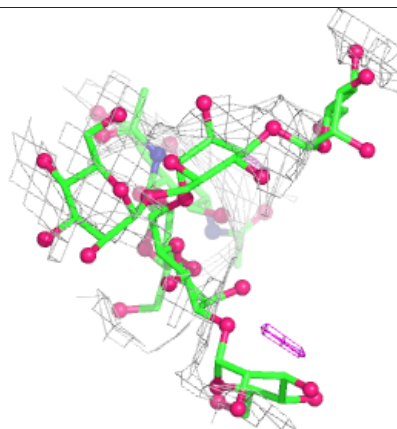
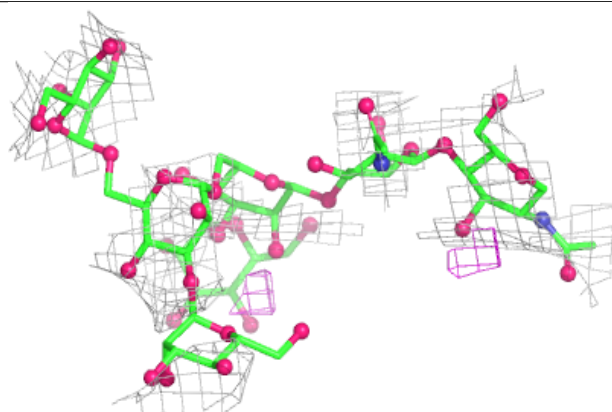
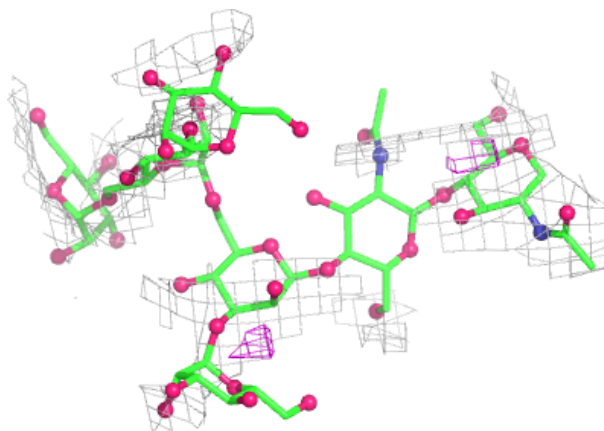
Electron density around Chain O:

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



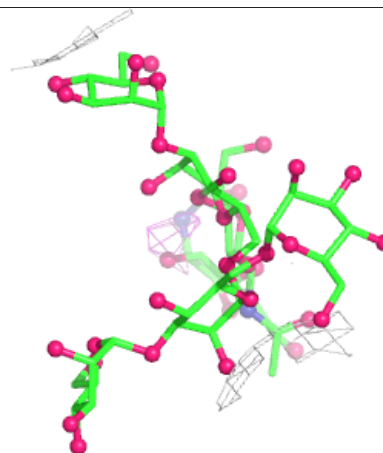
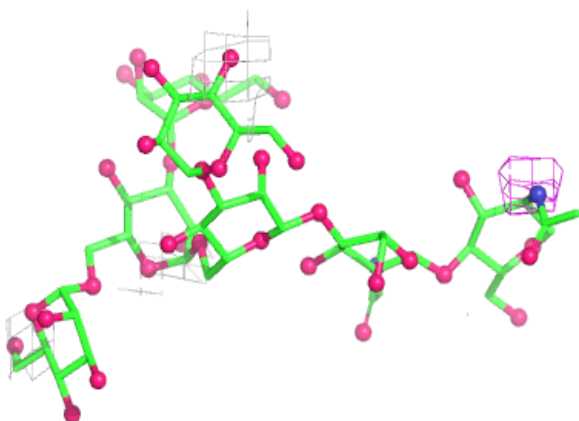
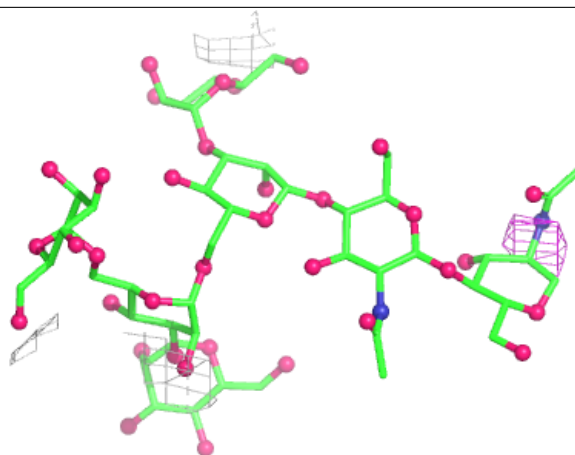
Electron density around Chain P:

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

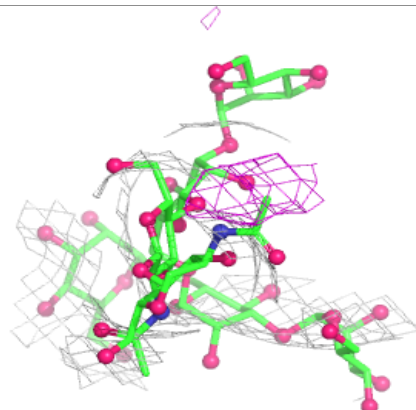
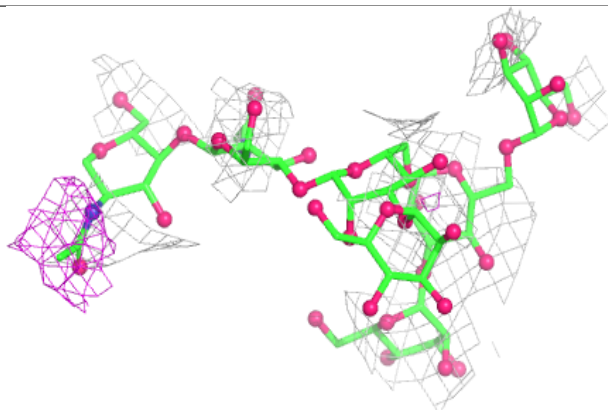
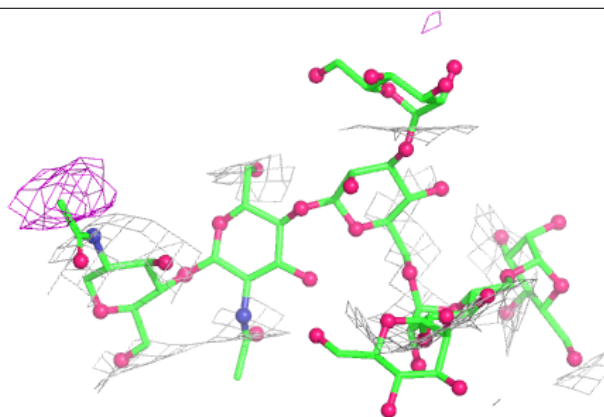


Electron density around Chain S:

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

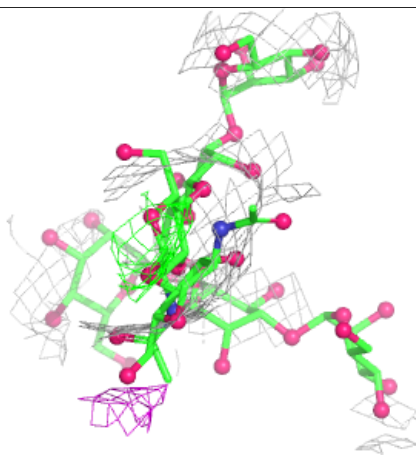
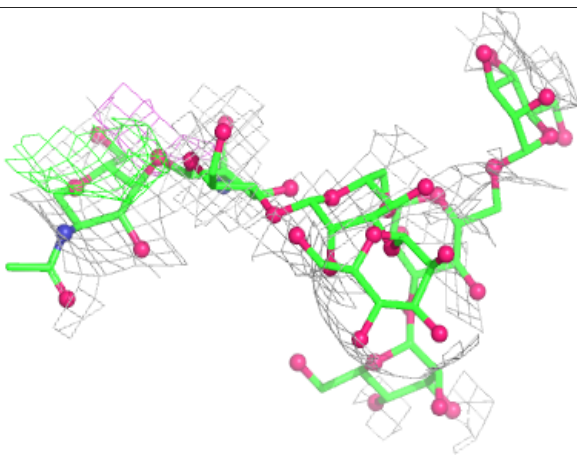
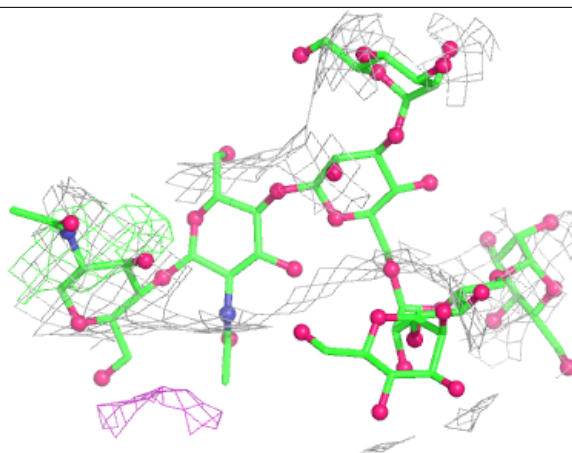
**Electron density around Chain T:**

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

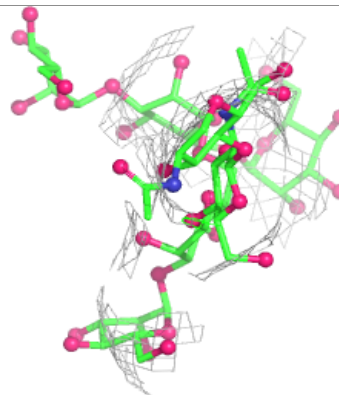
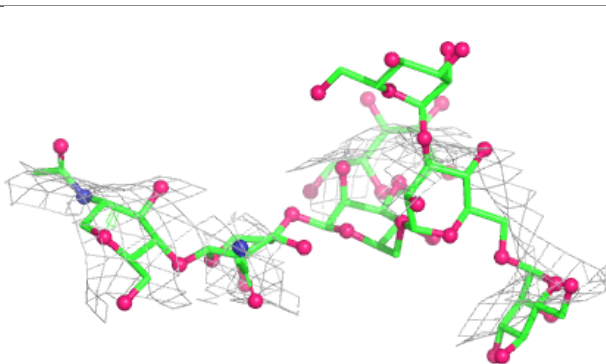
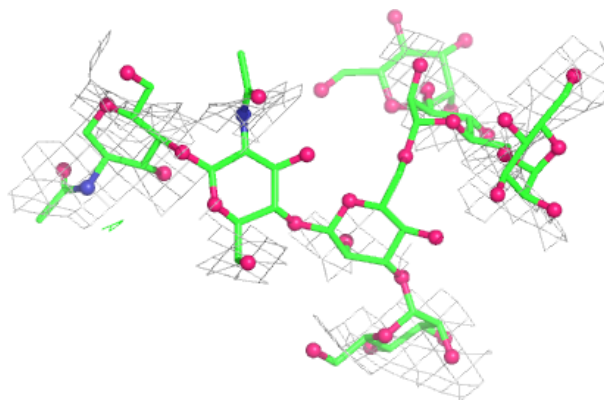


Electron density around Chain U:

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

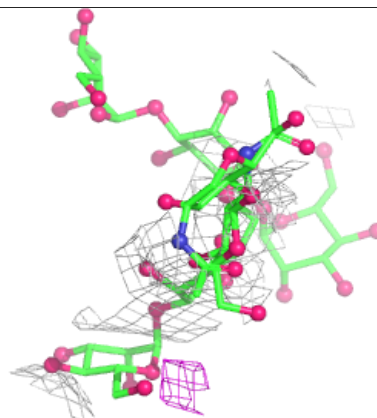
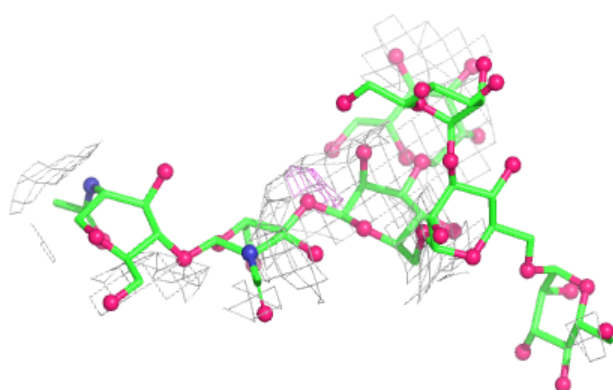
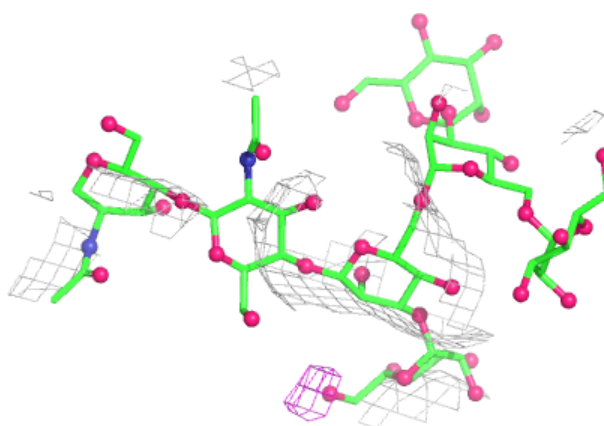
**Electron density around Chain V:**

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

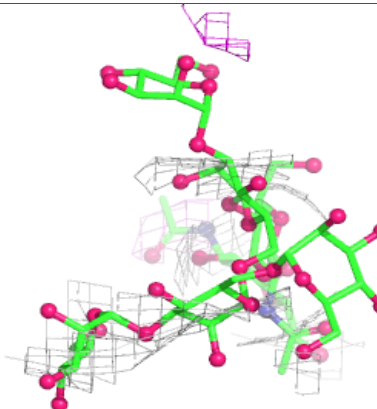
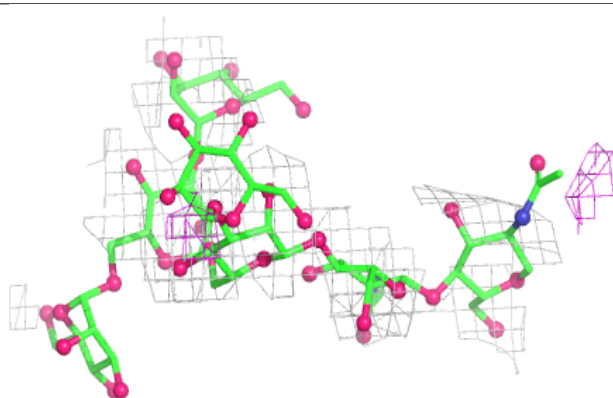
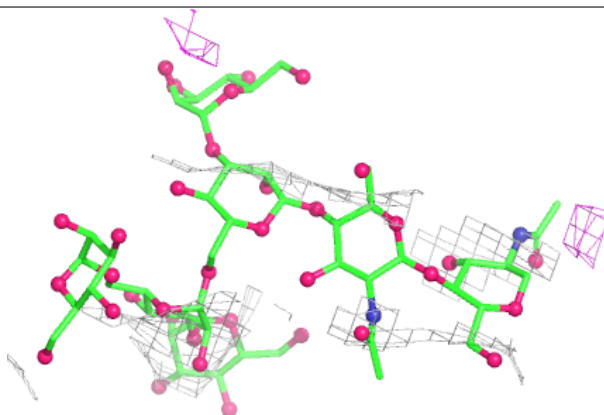


Electron density around Chain Y:

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

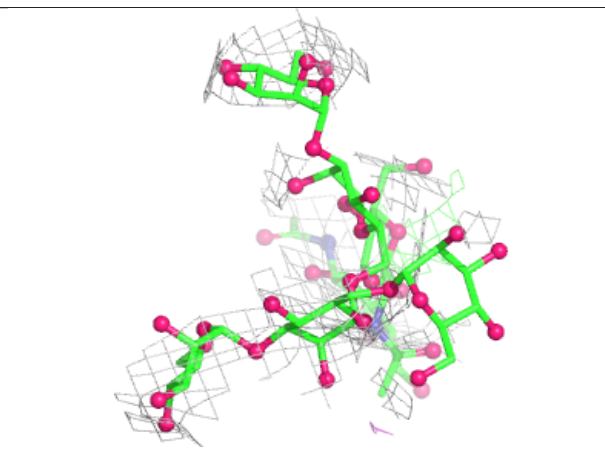
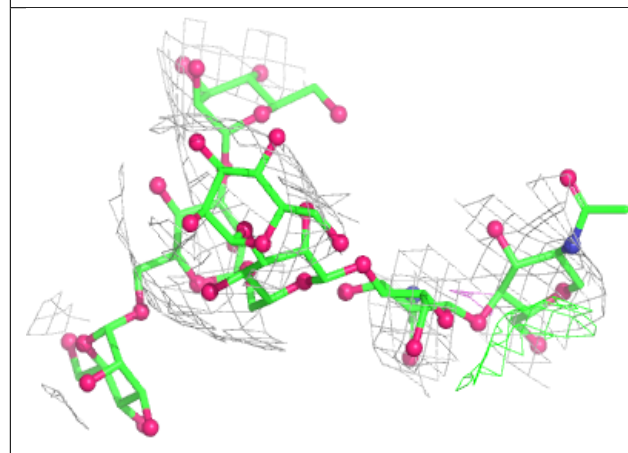
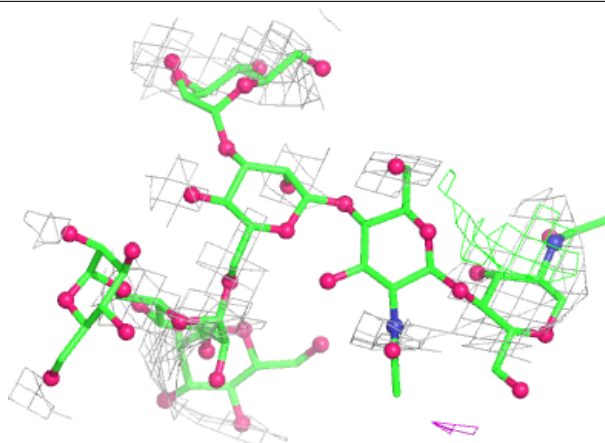
**Electron density around Chain Z:**

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

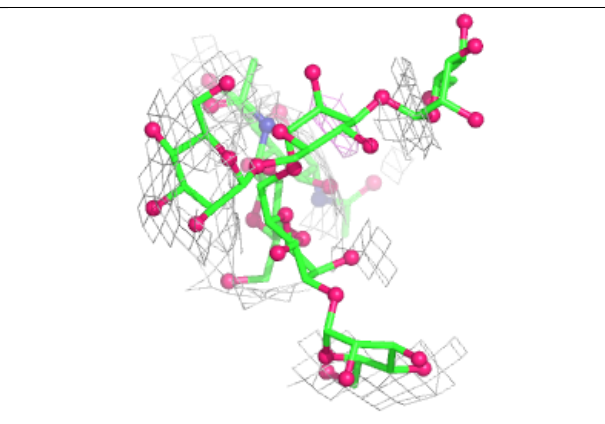
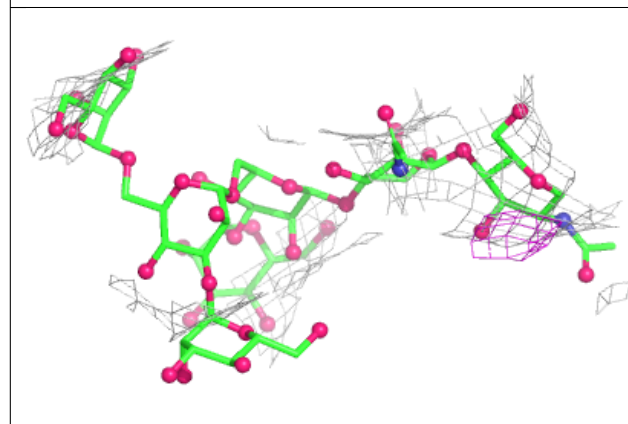
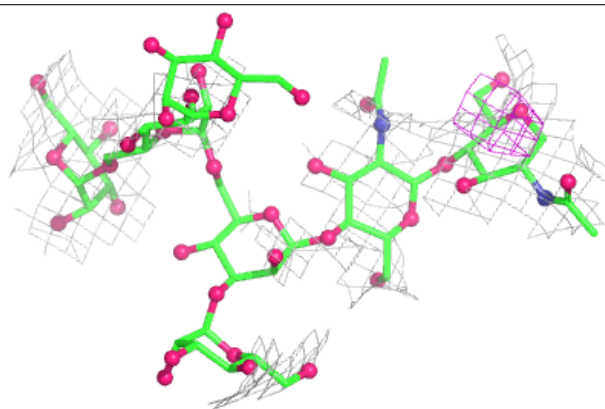


Electron density around Chain a:

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

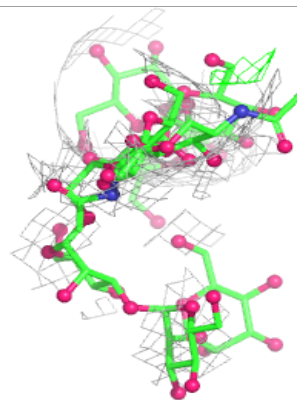
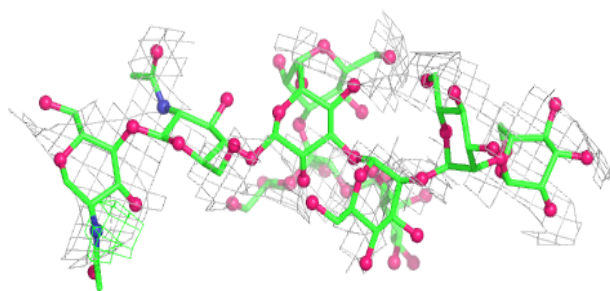
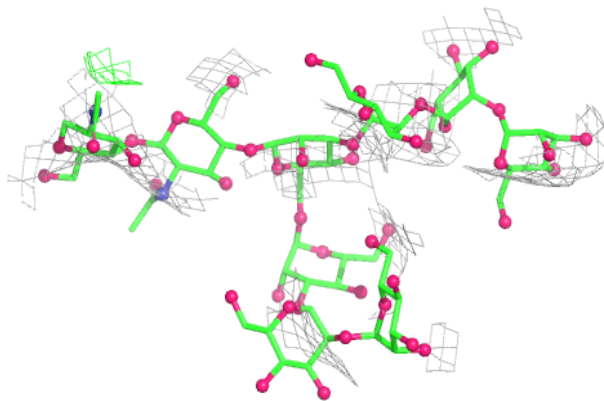
**Electron density around Chain b:**

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

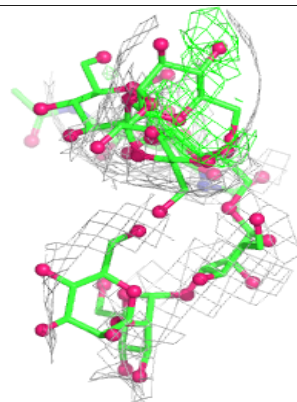
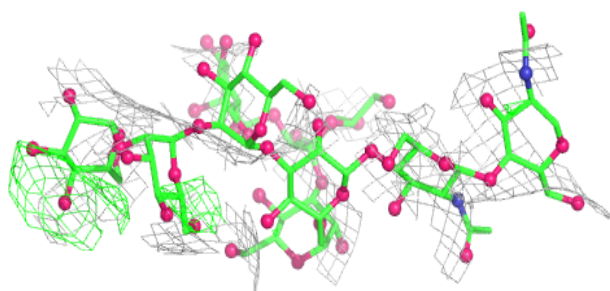
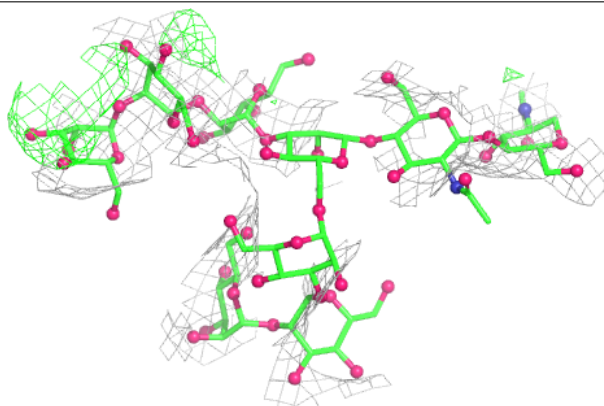


Electron density around Chain Q:

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

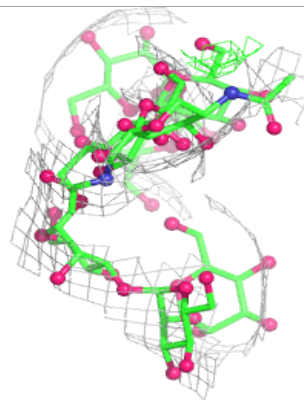
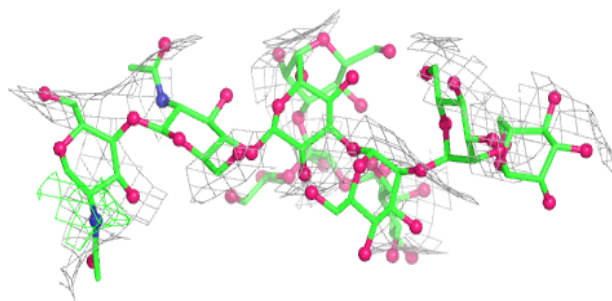
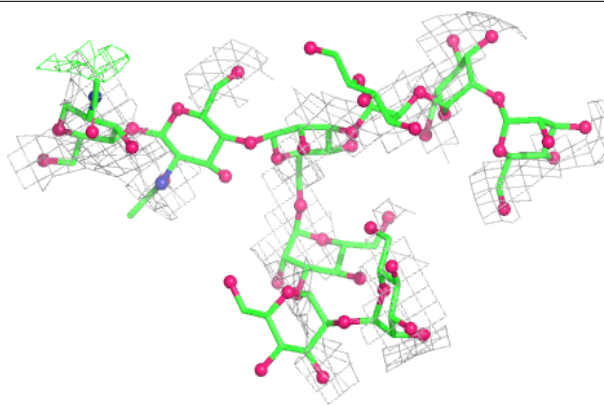
**Electron density around Chain W:**

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



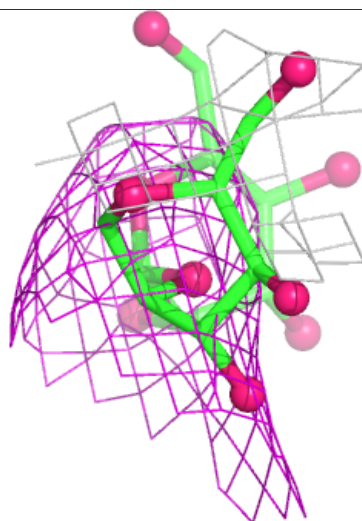
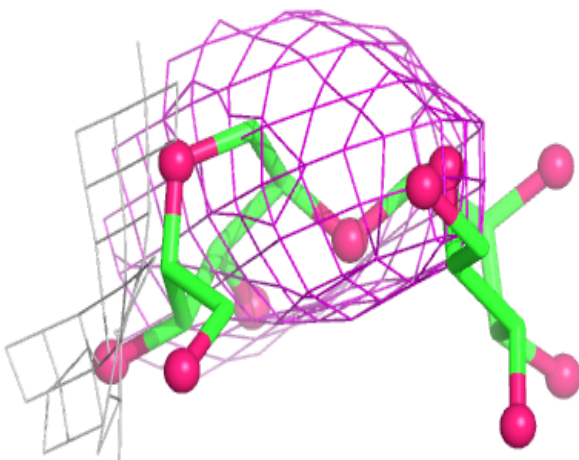
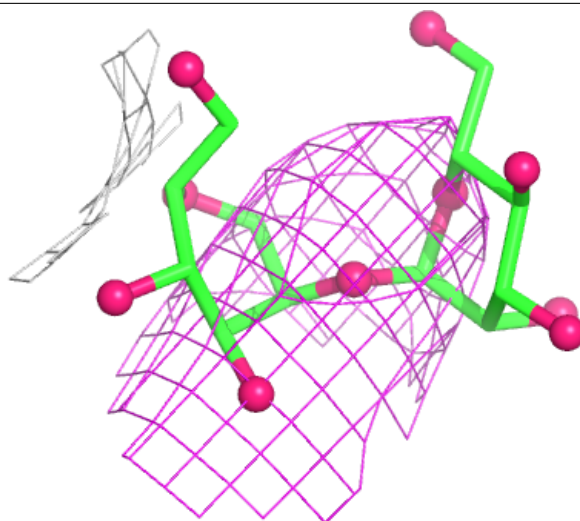
Electron density around Chain c:

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



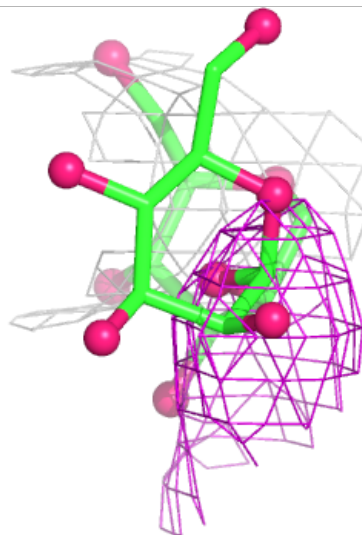
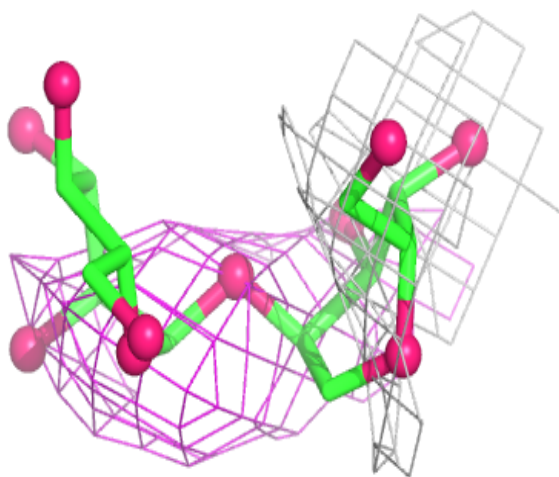
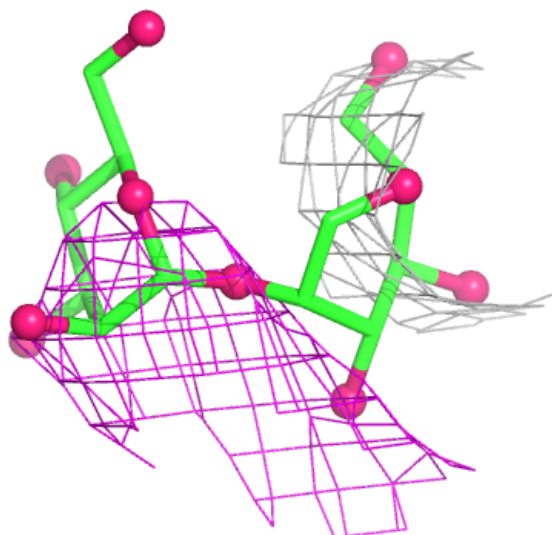
Electron density around Chain R:

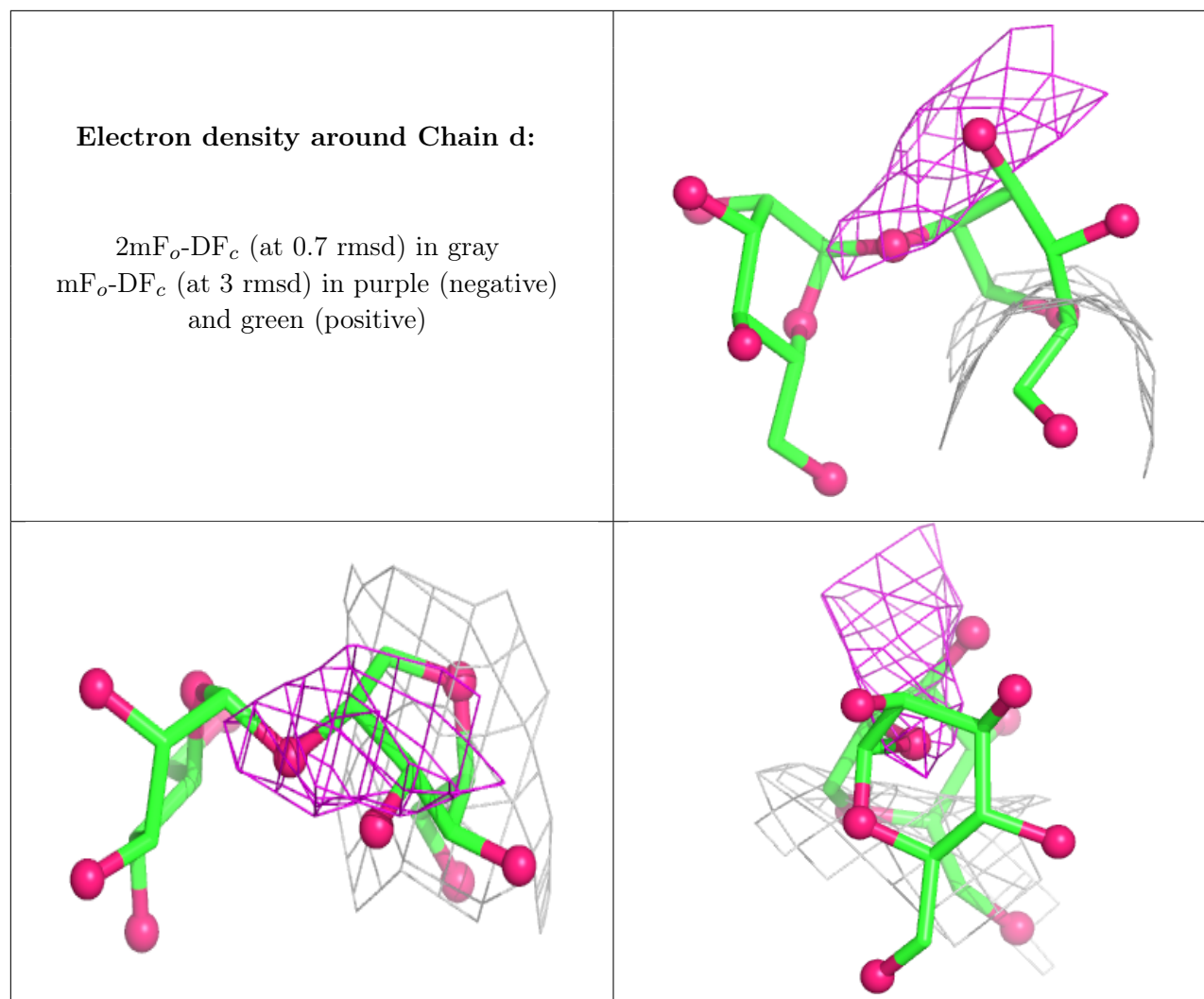
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain X:

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





6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	NAG	E	1355	14/15	0.15	0.21	411,411,411,411	0
8	NAG	E	1392	14/15	0.23	0.20	305,305,305,305	0
8	NAG	I	1355	14/15	0.23	0.13	266,266,266,266	0
8	NAG	A	1392	14/15	0.25	0.32	386,386,386,386	0
8	NAG	A	1355	14/15	0.37	0.13	238,238,238,238	0
8	NAG	A	1448	14/15	0.39	0.33	279,279,279,279	0
8	NAG	I	1160	14/15	0.45	0.22	362,362,362,362	0
8	NAG	E	1448	14/15	0.47	0.37	361,361,361,361	0
8	NAG	E	1088	14/15	0.49	0.14	486,486,486,486	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	NAG	I	1234	14/15	0.54	0.25	345,345,345,345	0
8	NAG	I	1392	14/15	0.55	0.22	345,345,345,345	0
8	NAG	I	1197	14/15	0.59	0.24	304,304,304,304	0
8	NAG	A	1088	14/15	0.60	0.15	474,474,474,474	0
8	NAG	E	1160	14/15	0.63	0.17	245,245,245,245	0
8	NAG	I	1088	14/15	0.64	0.20	503,503,503,503	0
8	NAG	E	1234	14/15	0.64	0.26	405,405,405,405	0
8	NAG	A	1160	14/15	0.66	0.15	267,267,267,267	0
8	NAG	I	1448	14/15	0.66	0.29	287,287,287,287	0
8	NAG	E	1276	14/15	0.67	0.17	313,313,313,313	0
8	NAG	A	1234	14/15	0.68	0.27	431,431,431,431	0
8	NAG	E	1197	14/15	0.72	0.26	288,288,288,288	0
8	NAG	E	1295	14/15	0.73	0.34	356,356,356,356	0
8	NAG	A	1386	14/15	0.75	0.31	356,356,356,356	0
8	NAG	A	1276	14/15	0.78	0.14	288,288,288,288	0
8	NAG	I	1276	14/15	0.80	0.14	281,281,281,281	0
8	NAG	I	1386	14/15	0.80	0.27	331,331,331,331	0
8	NAG	A	1295	14/15	0.82	0.32	330,330,330,330	0
8	NAG	I	1295	14/15	0.84	0.36	323,323,323,323	0
8	NAG	E	1386	14/15	0.85	0.25	427,427,427,427	0
8	NAG	A	1197	14/15	0.86	0.20	302,302,302,302	0

6.5 Other polymers

There are no such residues in this entry.