



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 08:02 PM UTC

PDB ID : 6MCO / pdb_00006mco
Title : Crystal structure of the B41 SOSIP.664 Env trimer with PGT124 and 35O22 Fabs, in P23 space group
Authors : Kumar, S.; Sarkar, A.; Wilson, I.A.
Deposited on : 2018-08-31
Resolution : 3.53 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
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

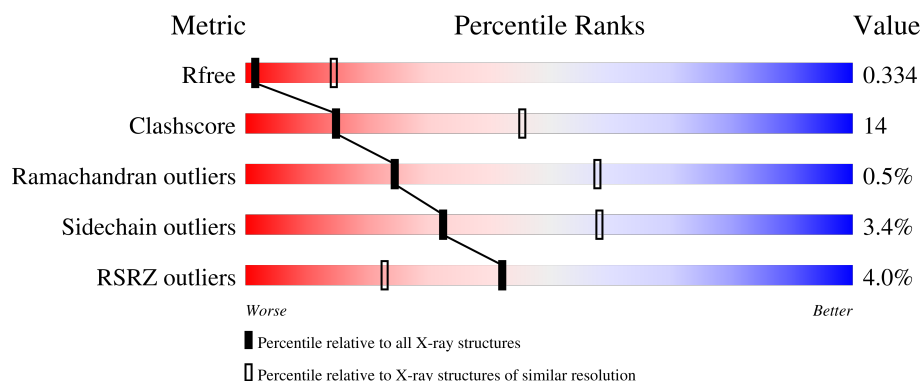
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.53 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	1025 (3.56-3.48)
Clashscore	190562	1079 (3.56-3.48)
Ramachandran outliers	187476	1052 (3.56-3.48)
Sidechain outliers	187428	1053 (3.56-3.48)
RSRZ outliers	180081	1024 (3.56-3.48)

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	B	153	10% 65% 18% • 13%
2	D	243	78% 20% •
3	E	216	% 84% 13% ••
4	G	482	3% 69% 23% • 5%
5	H	236	6% 67% 30% ••

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Mol	Chain	Length	Quality of chain
6	L	214	
7	A	6	
8	C	2	
8	I	2	
8	Q	2	
9	F	7	
10	J	4	
10	M	4	
11	K	3	
11	N	3	
11	O	3	
11	U	3	
12	P	8	
13	R	4	
14	S	5	
15	T	3	

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
8	NAG	I	2	-	-	-	X

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 12268 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 Transmembrane protein gp41.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	133	Total	C	N	O	S	0	0	0
			1058	675	178	198	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	605	CYS	THR	conflict	UNP B3UF08

- Molecule 2 is a protein called 35O22 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	240	Total	C	N	O	S	0	0	0
			1813	1150	303	352	8			

- Molecule 3 is a protein called 35O22 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	213	Total	C	N	O	S	0	0	0
			1615	1012	267	328	8			

- Molecule 4 is a protein called Surface protein gp120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	456	Total	C	N	O	S	0	0	0
			3576	2244	632	674	26			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	501	CYS	ALA	conflict	UNP B3UES2

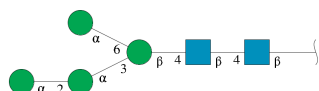
- Molecule 5 is a protein called PGT124 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	H	231	Total	C	N	O	S	0	0	0
			1754	1111	293	345	5			

- Molecule 6 is a protein called PGT124 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	L	211	Total	C	N	O	S	0	0	0
			1601	1008	271	317	5			

- Molecule 7 is an oligosaccharide called 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.



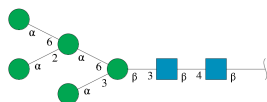
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	A	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 8 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
8	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
8	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
8	Q	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 9 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



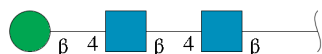
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	F	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 10 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



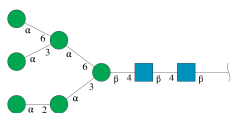
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
10	J	4	Total	C	N	O	0	0	0
			50	28	2	20			
10	M	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 11 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
11	K	3	Total	C	N	O	0	0	0
			39	22	2	15			
11	N	3	Total	C	N	O	0	0	0
			39	22	2	15			
11	O	3	Total	C	N	O	0	0	0
			39	22	2	15			
11	U	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 12 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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.



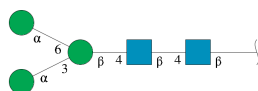
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
12	P	8	Total	C	N	O	0	0	0
			94	52	2	40			

- Molecule 13 is an oligosaccharide called 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.



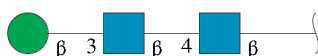
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
13	R	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 14 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.



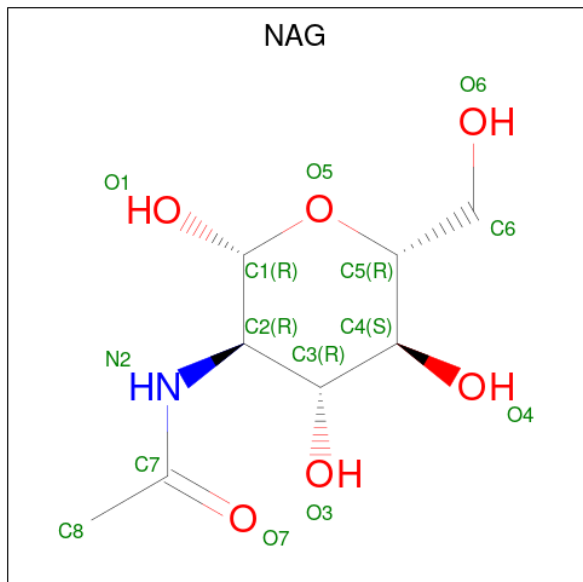
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
14	S	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 15 is an oligosaccharide called beta-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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

- Molecule 16 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

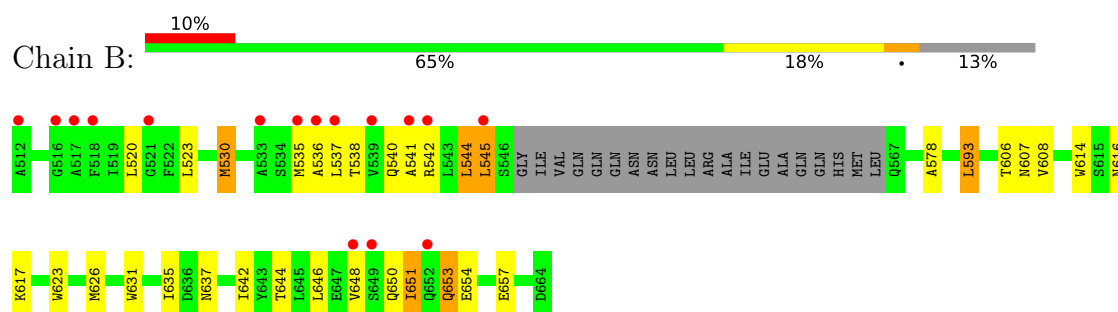


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
16	B	1	Total	C	N	O	0	0
			14	8	1	5		
16	B	1	Total	C	N	O	0	0
			14	8	1	5		
16	G	1	Total	C	N	O	0	0
			14	8	1	5		
16	G	1	Total	C	N	O	0	0
			14	8	1	5		
16	G	1	Total	C	N	O	0	0
			14	8	1	5		
16	G	1	Total	C	N	O	0	0
			14	8	1	5		
16	G	1	Total	C	N	O	0	0
			14	8	1	5		

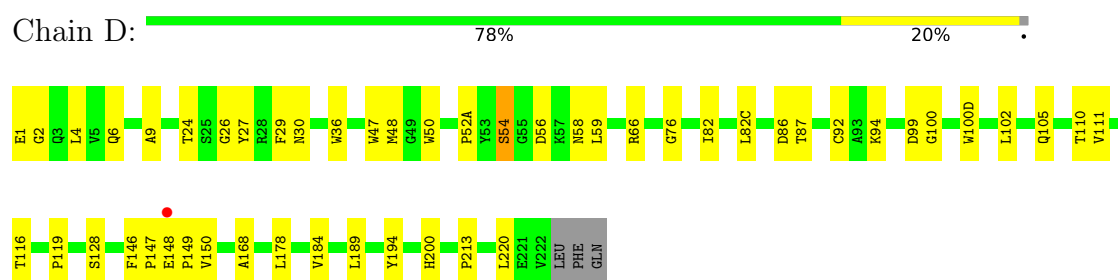
3 Residue-property plots [i](#)

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.

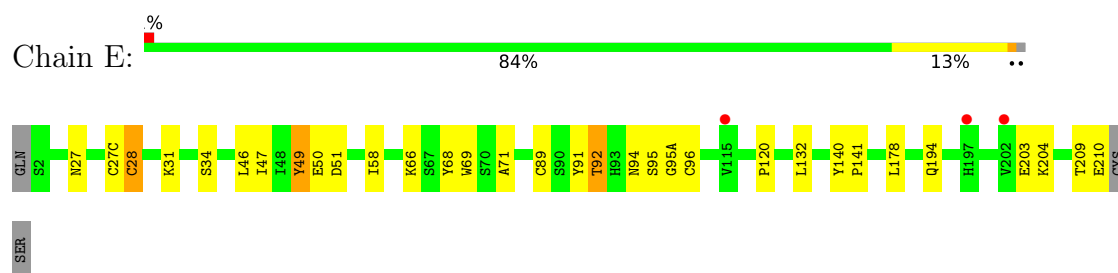
- Molecule 1: Transmembrane protein gp41



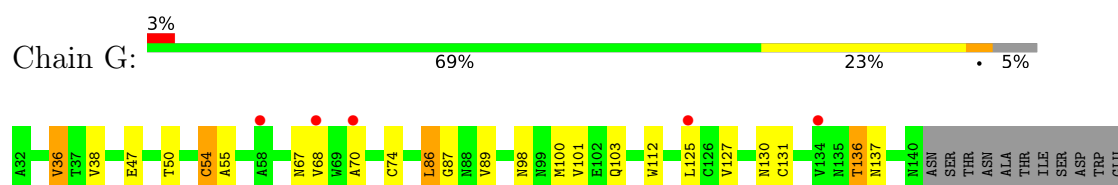
- Molecule 2: 35O22 Fab heavy chain

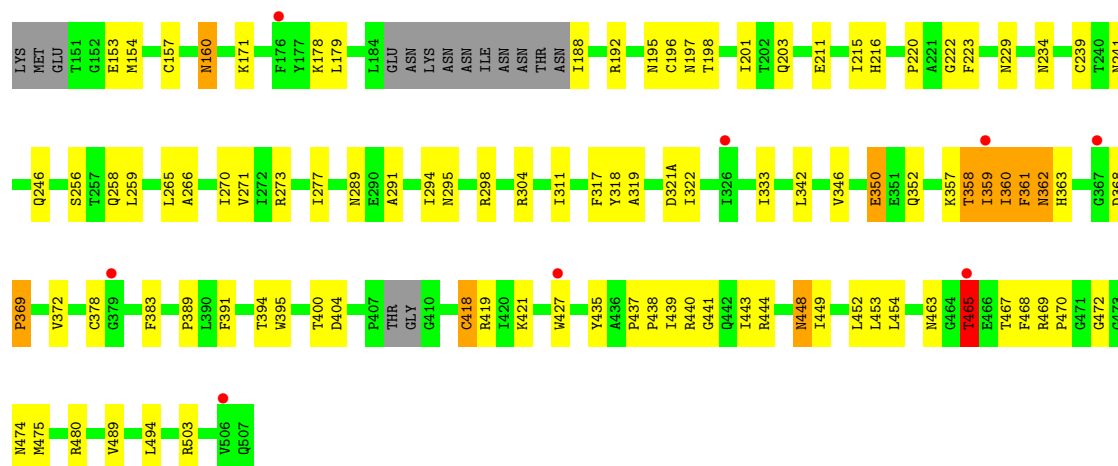


- Molecule 3: 35O22 Fab light chain

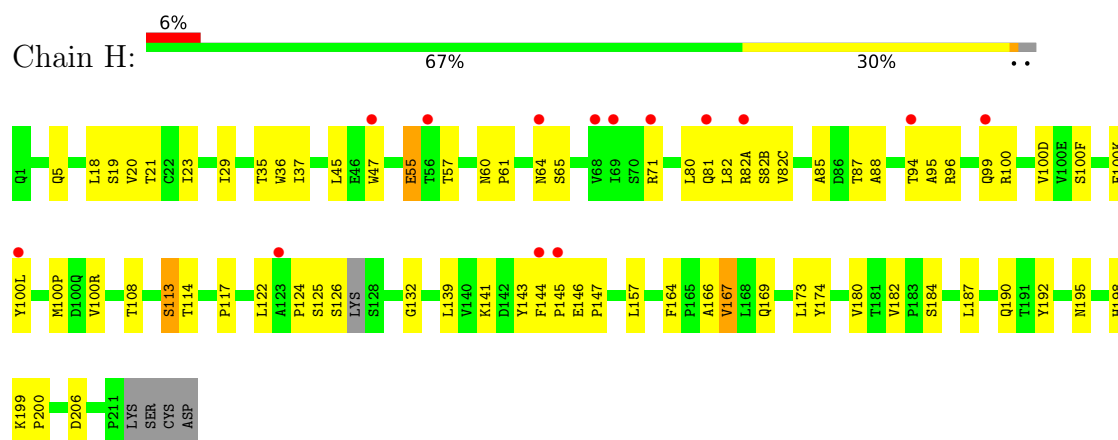


- Molecule 4: Surface protein gp120

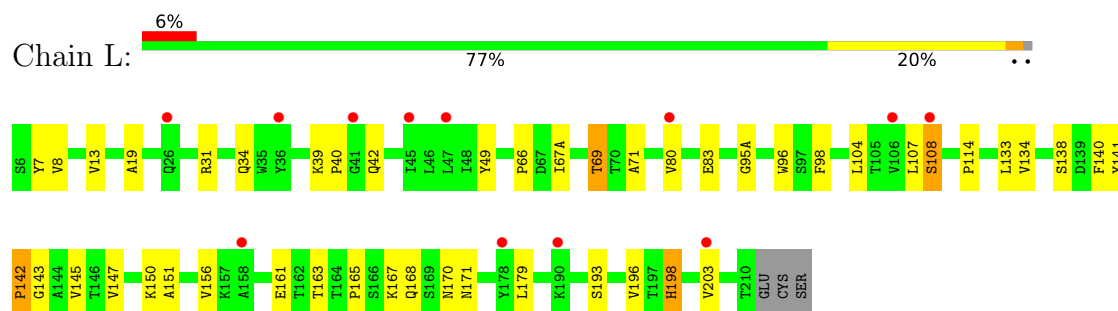




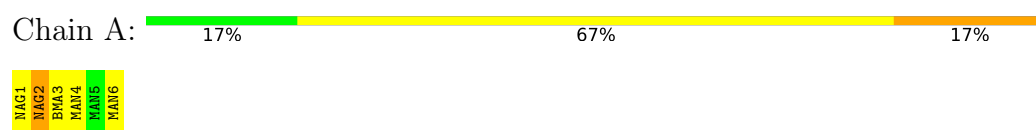
• Molecule 5: PGT124 Fab heavy chain



• Molecule 6: PGT124 Fab light chain



• Molecule 7: 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

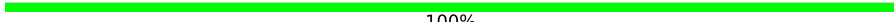


• Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  50% 50%



- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  100%

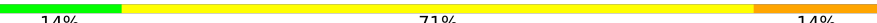


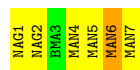
- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  50% 50%



- Molecule 9: alpha-D-mannopyranose-(1-2)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  14% 71% 14%



- Molecule 10: 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 J:  50% 50%



- Molecule 10: 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 M:  50% 50%



- Molecule 11: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  33% 33% 33%



- Molecule 11: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N: 33% 33% 33%



- Molecule 11: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O: 33% 67%



- Molecule 11: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain U: 100%



- Molecule 12: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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 P: 12% 62% 25%



- Molecule 13: 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 R: 25% 75%



- Molecule 14: 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: 40% 60%



- Molecule 15: beta-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:  67% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 3	Depositor
Cell constants a, b, c, α , β , γ	212.33Å 212.33Å 212.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.05 – 3.53 50.05 – 3.53	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.05-3.53) 99.7 (50.05-3.53)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 3.48Å)	Xtriage
Refinement program	PHENIX (1.12 _2829)	Depositor
R, R_{free}	0.306 , 0.322 0.320 , 0.334	Depositor DCC
R_{free} test set	1960 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	106.1	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 71.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.023 for l,-k,h	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	12268	wwPDB-VP
Average B, all atoms (Å ²)	124.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.96% 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: NAG, MAN, 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	B	0.26	1/1077 (0.1%)	0.50	2/1461 (0.1%)
2	D	0.14	0/1860	0.33	0/2533
3	E	0.18	0/1659	0.40	1/2269 (0.0%)
4	G	0.15	0/3651	0.38	0/4965
5	H	0.18	0/1797	0.56	5/2453 (0.2%)
6	L	0.20	0/1644	0.51	3/2246 (0.1%)
All	All	0.18	1/11688 (0.0%)	0.44	11/15927 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	545	LEU	CA-C	6.57	1.62	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	545	LEU	CA-C-O	7.59	127.85	121.02
6	L	69	THR	N-CA-C	-7.17	101.06	110.53
3	E	49	TYR	N-CA-C	7.00	119.00	111.36
6	L	151	ALA	N-CA-C	-6.86	98.06	108.96
5	H	55	GLU	CB-CA-C	-6.12	101.03	112.30
5	H	57	THR	N-CA-CB	-5.83	101.47	110.56
5	H	96	ARG	N-CA-C	5.66	118.13	108.90
6	L	71	ALA	N-CA-CB	-5.61	100.40	109.60
1	B	544	LEU	CB-CA-C	-5.60	104.09	111.82
5	H	88	ALA	N-CA-C	5.26	116.84	108.79
5	H	113	SER	N-CA-C	5.03	117.02	110.53

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1058	0	1042	38	0
2	D	1813	0	1784	34	0
3	E	1615	0	1544	23	0
4	G	3576	0	3493	123	0
5	H	1754	0	1719	85	0
6	L	1601	0	1546	47	0
7	A	72	0	61	3	0
8	C	28	0	25	0	0
8	I	28	0	25	0	0
8	Q	28	0	25	2	0
9	F	83	0	69	1	0
10	J	50	0	43	0	0
10	M	50	0	43	0	0
11	K	39	0	34	5	0
11	N	39	0	34	2	0
11	O	39	0	34	2	0
11	U	39	0	34	0	0
12	P	94	0	79	7	0
13	R	50	0	43	0	0
14	S	61	0	52	1	0
15	T	39	0	34	0	0
16	B	28	0	26	1	0
16	G	84	0	78	5	0
All	All	12268	0	11867	333	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.

All (333) 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)
4:G:360:ILE:CG1	4:G:395:TRP:H	1.26	1.47
4:G:360:ILE:HG13	4:G:395:TRP:N	1.31	1.44
1:B:545:LEU:O	1:B:545:LEU:HD23	1.43	1.16
5:H:94:THR:HG22	5:H:100(R):VAL:HB	1.35	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:29:ILE:HB	5:H:71:ARG:HD2	1.11	1.07
4:G:361:PHE:HB3	4:G:468:PHE:H	1.15	1.07
4:G:361:PHE:HD2	4:G:468:PHE:HB2	1.14	1.05
4:G:361:PHE:CD2	4:G:468:PHE:HB2	1.91	1.04
5:H:94:THR:CG2	5:H:100(R):VAL:HB	1.88	1.03
5:H:99:GLN:NE2	5:H:100(L):TYR:HE1	1.57	1.03
4:G:360:ILE:HD11	4:G:395:TRP:O	1.61	0.99
5:H:29:ILE:HB	5:H:71:ARG:CD	2.01	0.91
5:H:99:GLN:HG2	5:H:100(L):TYR:CE1	2.06	0.89
3:E:51:ASP:OD1	3:E:66:LYS:CE	2.20	0.89
5:H:100(D):VAL:HG22	12:P:1:NAG:H61	1.57	0.86
5:H:35:THR:HG21	5:H:100(P):MET:HE2	1.57	0.86
6:L:39:LYS:HG2	6:L:42:GLN:NE2	1.91	0.85
4:G:361:PHE:HB3	4:G:468:PHE:N	1.90	0.85
6:L:140:PHE:HZ	6:L:165:PRO:HB3	1.42	0.85
4:G:360:ILE:CG1	4:G:395:TRP:N	2.09	0.84
6:L:141:TYR:HB3	6:L:142:PRO:HD3	1.58	0.84
4:G:360:ILE:CG1	4:G:394:THR:HA	2.08	0.83
4:G:360:ILE:HB	4:G:394:THR:HA	1.60	0.83
5:H:29:ILE:HD13	5:H:71:ARG:HE	1.42	0.83
5:H:99:GLN:HG2	5:H:100(L):TYR:CD1	2.15	0.82
4:G:360:ILE:HG13	4:G:395:TRP:H	0.65	0.81
3:E:51:ASP:OD1	3:E:66:LYS:NZ	2.14	0.79
5:H:29:ILE:CB	5:H:71:ARG:HD2	2.04	0.79
4:G:360:ILE:CB	4:G:394:THR:HA	2.13	0.79
4:G:360:ILE:HG13	4:G:394:THR:C	2.08	0.79
5:H:99:GLN:HE21	5:H:100(L):TYR:HE1	0.81	0.78
5:H:99:GLN:NE2	5:H:100(L):TYR:CE1	2.40	0.78
5:H:29:ILE:CD1	5:H:71:ARG:HE	1.97	0.77
5:H:35:THR:HG21	5:H:100(P):MET:CE	2.14	0.77
4:G:277:ILE:HD13	4:G:352:GLN:HG2	1.67	0.76
4:G:361:PHE:CB	4:G:468:PHE:H	1.95	0.76
4:G:154:MET:HG2	4:G:322:ILE:HD11	1.65	0.76
5:H:94:THR:HG22	5:H:100(R):VAL:CB	2.13	0.74
5:H:61:PRO:HA	5:H:64:ASN:HD22	1.53	0.74
4:G:400:THR:HG21	16:G:604:NAG:H3	1.70	0.74
1:B:520:LEU:HD12	1:B:523:LEU:HB2	1.71	0.73
1:B:545:LEU:O	1:B:545:LEU:CD2	2.30	0.73
5:H:20:VAL:HG12	5:H:21:THR:N	2.02	0.72
5:H:169:GLN:HG3	6:L:161:GLU:HG3	1.70	0.72
5:H:100(D):VAL:HG22	12:P:1:NAG:C6	2.19	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:51:ASP:OD1	3:E:66:LYS:HE2	1.89	0.72
6:L:141:TYR:CB	6:L:142:PRO:HD3	2.17	0.72
6:L:66:PRO:HB2	12:P:5:MAN:O3	1.90	0.71
5:H:94:THR:HG23	5:H:94:THR:O	1.90	0.71
1:B:545:LEU:HD22	4:G:246:GLN:CD	2.14	0.71
3:E:34:SER:OG	3:E:89:CYS:HB3	1.92	0.70
6:L:142:PRO:CD	6:L:198:HIS:NE2	2.54	0.70
2:D:150:VAL:HG11	2:D:178:LEU:HD21	1.73	0.70
4:G:360:ILE:HG21	4:G:395:TRP:HD1	1.56	0.70
4:G:67:ASN:ND2	4:G:211:GLU:O	2.25	0.70
4:G:400:THR:HB	16:G:604:NAG:H5	1.73	0.69
6:L:107:LEU:O	6:L:108:SER:HB3	1.90	0.69
5:H:139:LEU:HD11	6:L:134:VAL:HG21	1.74	0.69
5:H:166:ALA:HB1	5:H:174:TYR:HB3	1.74	0.69
4:G:360:ILE:CD1	4:G:395:TRP:H	2.03	0.68
4:G:295:ASN:HD22	11:N:1:NAG:H83	1.58	0.68
5:H:195:ASN:ND2	5:H:206:ASP:OD2	2.28	0.67
2:D:48:MET:HE1	2:D:82:ILE:HD11	1.77	0.67
4:G:192:ARG:HH21	4:G:196:CYS:HB2	1.60	0.67
1:B:545:LEU:HD22	4:G:246:GLN:NE2	2.10	0.66
6:L:150:LYS:HB2	6:L:193:SER:HB2	1.76	0.66
2:D:47:TRP:HH2	2:D:58:ASN:HD22	1.41	0.66
6:L:142:PRO:HD3	6:L:198:HIS:NE2	2.09	0.66
5:H:99:GLN:CG	5:H:100(L):TYR:CE1	2.78	0.66
5:H:29:ILE:HD13	5:H:71:ARG:NE	2.11	0.66
5:H:94:THR:CG2	5:H:100(R):VAL:CB	2.70	0.66
4:G:361:PHE:HD2	4:G:468:PHE:CB	1.99	0.65
4:G:361:PHE:HA	4:G:467:THR:HB	1.78	0.65
6:L:114:PRO:HB3	6:L:140:PHE:HB3	1.78	0.65
4:G:454:LEU:HA	4:G:470:PRO:HA	1.79	0.64
4:G:101:VAL:HG11	4:G:480:ARG:HG3	1.79	0.64
4:G:294:ILE:HG13	4:G:333:ILE:HG12	1.79	0.64
4:G:360:ILE:HD11	4:G:395:TRP:C	2.23	0.64
5:H:35:THR:OG1	5:H:47:TRP:NE1	2.30	0.64
5:H:5:GLN:HB3	5:H:23:ILE:CG1	2.27	0.64
4:G:361:PHE:HD1	4:G:361:PHE:O	1.81	0.64
4:G:427:TRP:HD1	4:G:475:MET:HG2	1.62	0.64
2:D:27:TYR:HB3	7:A:2:NAG:H83	1.79	0.63
5:H:61:PRO:HA	5:H:64:ASN:ND2	2.14	0.62
4:G:360:ILE:HG21	4:G:395:TRP:CD1	2.35	0.62
3:E:27:ASN:HA	3:E:27(C):CYS:HB2	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:132:LEU:HD12	3:E:178:LEU:HD23	1.81	0.61
1:B:617:LYS:HE3	1:B:626:MET:HE1	1.82	0.61
4:G:389:PRO:HB3	14:S:1:NAG:H62	1.82	0.61
2:D:4:LEU:HG	2:D:24:THR:HG22	1.81	0.61
4:G:54:CYS:SG	4:G:55:ALA:N	2.73	0.61
6:L:39:LYS:CG	6:L:42:GLN:NE2	2.64	0.61
6:L:142:PRO:HD2	6:L:198:HIS:NE2	2.15	0.61
5:H:20:VAL:CG1	5:H:21:THR:N	2.64	0.61
4:G:360:ILE:HD13	4:G:360:ILE:N	2.17	0.60
6:L:141:TYR:CD2	6:L:142:PRO:HG3	2.37	0.60
4:G:360:ILE:CG1	4:G:394:THR:CA	2.80	0.60
4:G:444:ARG:HH12	12:P:1:NAG:H81	1.67	0.60
5:H:94:THR:HG21	5:H:100(R):VAL:HB	1.78	0.60
4:G:137:ASN:HB3	6:L:95(A):GLY:HA2	1.84	0.60
1:B:607:ASN:H	1:B:650:GLN:CD	2.10	0.59
5:H:20:VAL:HG12	5:H:21:THR:H	1.65	0.59
4:G:362:ASN:HA	4:G:391:PHE:O	2.03	0.59
5:H:124:PRO:C	5:H:126:SER:H	2.11	0.59
5:H:47:TRP:HB2	6:L:98:PHE:HE1	1.67	0.59
5:H:55:GLU:HG3	5:H:71:ARG:NH2	2.17	0.59
1:B:530:MET:HE1	1:B:631:TRP:HB2	1.83	0.59
4:G:463:ASN:OD1	16:G:606:NAG:N2	2.36	0.59
4:G:304:ARG:HH12	4:G:440:ARG:HD3	1.66	0.59
4:G:383:PHE:HB3	4:G:418:CYS:SG	2.42	0.59
1:B:616:ASN:HD22	16:B:701:NAG:C7	2.15	0.58
2:D:102:LEU:HD11	7:A:2:NAG:H2	1.84	0.58
4:G:358:THR:OG1	4:G:359:ILE:N	2.37	0.58
5:H:55:GLU:CG	5:H:71:ARG:NH2	2.67	0.58
5:H:20:VAL:CG1	5:H:21:THR:H	2.16	0.58
4:G:203:GLN:HE22	4:G:318:TYR:HD2	1.52	0.58
1:B:540:GLN:HG3	1:B:544:LEU:HD13	1.85	0.58
6:L:141:TYR:HD2	6:L:142:PRO:HG3	1.69	0.58
1:B:578:ALA:HB1	4:G:220:PRO:HB3	1.86	0.57
2:D:94:LYS:HB3	2:D:102:LEU:HB2	1.85	0.57
5:H:94:THR:CG2	5:H:94:THR:O	2.52	0.57
5:H:19:SER:O	5:H:20:VAL:HG23	2.04	0.57
2:D:54:SER:HB2	2:D:56:ASP:OD2	2.04	0.57
1:B:606:THR:OG1	4:G:36:VAL:O	2.22	0.57
4:G:201:ILE:HD11	4:G:435:TYR:HB2	1.86	0.57
4:G:229:ASN:ND2	11:K:1:NAG:H62	2.19	0.57
6:L:80:VAL:HG11	6:L:171:ASN:HB3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:131:CYS:HA	4:G:157:CYS:HA	1.87	0.56
1:B:608:VAL:N	1:B:650:GLN:OE1	2.39	0.56
4:G:55:ALA:HB3	4:G:216:HIS:HB2	1.86	0.56
1:B:535:MET:HG3	1:B:536:ALA:H	1.71	0.56
4:G:229:ASN:HD22	11:K:1:NAG:H62	1.70	0.56
4:G:360:ILE:HG13	4:G:394:THR:CA	2.35	0.56
1:B:614:TRP:HB3	1:B:635:ILE:HD11	1.87	0.56
5:H:85:ALA:C	5:H:87:THR:N	2.64	0.55
4:G:259:LEU:HD23	4:G:449:ILE:HG21	1.88	0.55
5:H:100:ARG:HH12	12:P:5:MAN:H5	1.71	0.55
4:G:67:ASN:OD1	4:G:70:ALA:N	2.37	0.54
5:H:117:PRO:HB3	5:H:143:TYR:HB3	1.87	0.54
6:L:143:GLY:O	6:L:165:PRO:HG2	2.07	0.54
3:E:94:ASN:HB2	9:F:6:MAN:H61	1.88	0.54
6:L:142:PRO:HD2	6:L:198:HIS:CD2	2.43	0.54
1:B:535:MET:HE2	4:G:87:GLY:H	1.73	0.54
5:H:5:GLN:HB3	5:H:23:ILE:HD11	1.89	0.54
5:H:146:GLU:N	5:H:147:PRO:HD2	2.23	0.54
4:G:160:ASN:HA	4:G:171:LYS:HA	1.90	0.53
4:G:362:ASN:OD1	4:G:362:ASN:N	2.40	0.53
5:H:37:ILE:HG21	5:H:45:LEU:HD23	1.90	0.53
2:D:9:ALA:HB3	2:D:149:PRO:HG3	1.90	0.53
4:G:363:HIS:HA	4:G:469:ARG:HG2	1.89	0.53
4:G:256:SER:OG	4:G:259:LEU:O	2.25	0.53
4:G:360:ILE:CD1	4:G:395:TRP:O	2.48	0.53
5:H:132:GLY:O	5:H:184:SER:N	2.35	0.53
6:L:13:VAL:HG11	6:L:19:ALA:HB2	1.90	0.53
1:B:545:LEU:HD21	4:G:246:GLN:HG2	1.91	0.53
1:B:545:LEU:HD13	4:G:246:GLN:OE1	2.09	0.52
3:E:27(C):CYS:SG	3:E:28:CYS:N	2.83	0.52
5:H:94:THR:HG21	5:H:100(R):VAL:CG2	2.40	0.52
6:L:107:LEU:O	6:L:108:SER:CB	2.58	0.52
4:G:474:ASN:OD1	4:G:475:MET:N	2.40	0.52
3:E:31:LYS:HD3	3:E:92:THR:HG23	1.92	0.52
3:E:47:ILE:HA	3:E:58:ILE:HG13	1.92	0.52
6:L:34:GLN:HG3	6:L:49:TYR:HA	1.90	0.52
3:E:49:TYR:O	3:E:91:TYR:OH	2.26	0.52
2:D:168:ALA:HA	2:D:178:LEU:HB3	1.90	0.52
11:K:1:NAG:H61	11:K:2:NAG:HN2	1.75	0.52
1:B:545:LEU:CD2	4:G:246:GLN:CD	2.84	0.51
4:G:448:ASN:OD1	4:G:448:ASN:N	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:469:ARG:HG3	8:Q:1:NAG:H81	1.92	0.51
6:L:156:VAL:HG11	6:L:179:LEU:HD11	1.93	0.51
5:H:85:ALA:C	5:H:87:THR:H	2.16	0.51
4:G:125:LEU:HD11	4:G:311:ILE:HG21	1.92	0.51
5:H:19:SER:O	5:H:20:VAL:CG2	2.59	0.51
1:B:535:MET:N	1:B:538:THR:OG1	2.44	0.51
4:G:350:GLU:HG2	4:G:357:LYS:HG2	1.93	0.51
5:H:47:TRP:HB2	6:L:98:PHE:CE1	2.45	0.51
6:L:147:VAL:HG22	6:L:196:VAL:HG22	1.93	0.51
4:G:197:ASN:OD1	4:G:198:THR:HG23	2.12	0.50
5:H:5:GLN:HB3	5:H:23:ILE:HG13	1.92	0.50
5:H:18:LEU:O	5:H:81:GLN:HA	2.12	0.50
4:G:360:ILE:O	4:G:467:THR:HG22	2.12	0.50
4:G:453:LEU:HD23	4:G:472:GLY:HA2	1.92	0.50
5:H:122:LEU:HD21	5:H:139:LEU:HB2	1.93	0.50
1:B:542:ARG:HD3	4:G:222:GLY:HA2	1.92	0.50
3:E:46:LEU:HD21	3:E:49:TYR:HB3	1.94	0.50
1:B:607:ASN:ND2	1:B:654:GLU:OE2	2.45	0.50
3:E:49:TYR:CD2	3:E:50:GLU:HG3	2.47	0.50
4:G:195:ASN:HD21	4:G:201:ILE:HD12	1.77	0.50
5:H:184:SER:HA	5:H:187:LEU:HB2	1.92	0.50
1:B:646:LEU:HD11	4:G:36:VAL:HG21	1.93	0.49
2:D:87:THR:HG23	2:D:110:THR:HA	1.93	0.49
4:G:197:ASN:OD1	4:G:198:THR:N	2.45	0.49
1:B:606:THR:HG22	4:G:503:ARG:NH2	2.27	0.49
4:G:360:ILE:HG13	4:G:394:THR:HA	1.87	0.49
5:H:145:PRO:HG2	5:H:200:PRO:HG2	1.94	0.49
6:L:140:PHE:CD1	6:L:142:PRO:O	2.66	0.49
1:B:650:GLN:O	1:B:654:GLU:N	2.31	0.48
2:D:66:ARG:NH2	2:D:86:ASP:OD2	2.40	0.48
3:E:140:TYR:N	3:E:141:PRO:HD2	2.29	0.48
4:G:265:LEU:HD21	4:G:291:ALA:HB2	1.94	0.48
1:B:545:LEU:HD11	4:G:246:GLN:HB2	1.96	0.48
6:L:140:PHE:CZ	6:L:165:PRO:HB3	2.34	0.48
4:G:298:ARG:NH2	4:G:441:GLY:O	2.47	0.48
5:H:64:ASN:OD1	5:H:65:SER:N	2.47	0.48
4:G:317:PHE:CE2	4:G:319:ALA:HB2	2.49	0.48
2:D:2:GLY:HA2	2:D:26:GLY:HA3	1.96	0.47
2:D:189:LEU:HD22	2:D:213:PRO:HB2	1.95	0.47
4:G:112:TRP:CD2	4:G:427:TRP:HH2	2.32	0.47
5:H:100(P):MET:SD	5:H:100(P):MET:N	2.87	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:180:VAL:HG12	5:H:182:VAL:HG13	1.96	0.47
5:H:5:GLN:HB3	5:H:23:ILE:CD1	2.44	0.47
5:H:55:GLU:HG3	5:H:71:ARG:HH22	1.78	0.47
2:D:128:SER:HB2	2:D:220:LEU:HB2	1.96	0.47
6:L:39:LYS:HD2	6:L:40:PRO:HD2	1.95	0.47
11:K:2:NAG:O4	11:K:2:NAG:O6	2.32	0.47
2:D:168:ALA:HB2	2:D:178:LEU:HD23	1.97	0.47
4:G:361:PHE:N	4:G:361:PHE:CD1	2.73	0.47
5:H:124:PRO:O	5:H:126:SER:N	2.44	0.47
5:H:143:TYR:CE2	5:H:174:TYR:HB2	2.49	0.47
5:H:157:LEU:HD21	5:H:180:VAL:HG11	1.96	0.47
6:L:67(A):ILE:HG23	12:P:5:MAN:O3	2.15	0.47
6:L:168:GLN:HG3	6:L:170:ASN:H	1.79	0.47
11:N:1:NAG:H61	11:N:2:NAG:C7	2.45	0.47
2:D:47:TRP:HZ2	2:D:50:TRP:CD1	2.32	0.47
6:L:198:HIS:ND1	6:L:198:HIS:C	2.73	0.47
6:L:34:GLN:NE2	6:L:49:TYR:O	2.48	0.46
4:G:179:LEU:HB2	4:G:421:LYS:HG2	1.97	0.46
4:G:359:ILE:HD13	4:G:465:THR:HG21	1.97	0.46
5:H:167:VAL:HB	6:L:163:THR:OG1	2.15	0.46
2:D:6:GLN:H	2:D:105:GLN:HE22	1.62	0.46
4:G:87:GLY:O	4:G:89:VAL:HG23	2.16	0.46
4:G:427:TRP:HE1	4:G:475:MET:HE2	1.81	0.46
6:L:133:LEU:HD12	6:L:179:LEU:HD23	1.98	0.46
2:D:100(D):TRP:HZ2	3:E:91:TYR:HB2	1.81	0.46
4:G:360:ILE:HG12	4:G:394:THR:HG23	1.98	0.46
4:G:360:ILE:HG12	4:G:395:TRP:N	2.22	0.46
2:D:1:GLU:HB3	7:A:2:NAG:H3	1.97	0.45
4:G:160:ASN:HD22	16:G:601:NAG:H83	1.80	0.45
4:G:321(A):ASP:OD1	4:G:321(A):ASP:N	2.44	0.45
1:B:530:MET:HG2	1:B:623:TRP:HA	1.98	0.45
5:H:61:PRO:O	5:H:64:ASN:ND2	2.49	0.45
4:G:258:GLN:HG3	4:G:470:PRO:HB2	1.97	0.45
1:B:642:ILE:HG22	1:B:646:LEU:HD13	1.97	0.45
4:G:86:LEU:HB3	4:G:89:VAL:HG21	1.98	0.45
4:G:368:ASP:HB2	4:G:369:PRO:HD3	1.98	0.45
2:D:29:PHE:CG	2:D:76:GLY:HA3	2.52	0.45
2:D:47:TRP:HH2	2:D:58:ASN:ND2	2.09	0.45
6:L:140:PHE:HE1	6:L:143:GLY:HA2	1.81	0.45
5:H:99:GLN:CD	5:H:100(L):TYR:HE1	2.23	0.45
8:Q:1:NAG:HO6	8:Q:2:NAG:HN2	1.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:94:THR:HG21	5:H:100(R):VAL:CB	2.44	0.44
5:H:199:LYS:HB2	5:H:200:PRO:HD3	1.98	0.44
1:B:545:LEU:CD1	4:G:246:GLN:HB2	2.47	0.44
3:E:51:ASP:OD1	3:E:66:LYS:CD	2.65	0.44
4:G:361:PHE:HD2	4:G:468:PHE:CD2	2.35	0.44
2:D:116:THR:HG23	2:D:147:PRO:HG2	1.99	0.44
3:E:194:GLN:HG2	3:E:203:GLU:HG3	2.00	0.44
4:G:153:GLU:HA	4:G:178:LYS:HD2	2.00	0.43
4:G:220:PRO:HG2	4:G:223:PHE:CD1	2.53	0.43
4:G:266:ALA:O	4:G:289:ASN:HA	2.18	0.43
4:G:197:ASN:ND2	16:G:602:NAG:C7	2.81	0.43
4:G:391:PHE:CD2	4:G:470:PRO:HG3	2.53	0.43
5:H:19:SER:C	5:H:20:VAL:HG23	2.43	0.43
5:H:87:THR:HG23	5:H:108:THR:HA	2.01	0.43
5:H:182:VAL:HG11	5:H:192:TYR:CE2	2.54	0.43
6:L:80:VAL:HG13	6:L:167:LYS:NZ	2.33	0.43
6:L:7:TYR:HB3	6:L:8:VAL:H	1.66	0.43
4:G:271:VAL:HG11	4:G:273:ARG:CZ	2.49	0.43
6:L:83:GLU:HG3	6:L:104:LEU:O	2.19	0.43
1:B:644:THR:O	1:B:648:VAL:HG12	2.19	0.43
1:B:606:THR:HG22	4:G:503:ARG:HH21	1.84	0.43
2:D:146:PHE:HB3	2:D:147:PRO:HD3	2.01	0.43
4:G:319:ALA:C	4:G:438:PRO:HG2	2.44	0.43
4:G:289:ASN:OD1	4:G:289:ASN:N	2.52	0.42
11:O:1:NAG:H62	11:O:2:NAG:C7	2.49	0.42
1:B:653:GLN:O	1:B:657:GLU:HG2	2.19	0.42
5:H:36:TRP:CD2	5:H:80:LEU:HD22	2.55	0.42
6:L:150:LYS:HB3	6:L:150:LYS:HE2	1.83	0.42
2:D:59:LEU:HD23	2:D:59:LEU:HA	1.80	0.42
5:H:122:LEU:HD21	5:H:139:LEU:HD12	2.00	0.42
1:B:651:ILE:HD12	1:B:651:ILE:HA	1.77	0.42
4:G:178:LYS:HD3	4:G:419:ARG:HH11	1.85	0.42
2:D:99:ASP:OD1	2:D:100:GLY:N	2.52	0.42
4:G:437:PRO:HA	4:G:438:PRO:HD3	1.95	0.42
4:G:195:ASN:ND2	4:G:201:ILE:HB	2.35	0.42
4:G:239:CYS:SG	4:G:241:ASN:O	2.77	0.42
4:G:494:LEU:HA	4:G:494:LEU:HD23	1.72	0.42
5:H:100:ARG:HG2	5:H:100(K):PHE:CZ	2.55	0.41
5:H:145:PRO:HD2	5:H:198:HIS:HE2	1.84	0.41
2:D:119:PRO:HD3	2:D:200:HIS:ND1	2.35	0.41
4:G:361:PHE:CD2	4:G:468:PHE:HD2	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:100(D):VAL:HG22	12:P:1:NAG:C5	2.50	0.41
1:B:607:ASN:N	1:B:650:GLN:OE1	2.53	0.41
2:D:148:GLU:N	2:D:149:PRO:HD2	2.35	0.41
5:H:190:GLN:HG2	5:H:192:TYR:H	1.86	0.41
2:D:36:TRP:CZ3	2:D:92:CYS:HB3	2.55	0.41
3:E:95:SER:OG	3:E:95(A):GLY:N	2.54	0.41
11:K:1:NAG:O7	11:K:1:NAG:O3	2.28	0.41
5:H:164:PHE:CZ	6:L:138:SER:HB2	2.56	0.41
6:L:31:ARG:NH2	6:L:69:THR:OG1	2.54	0.41
3:E:68:TYR:CD1	3:E:69:TRP:HD1	2.38	0.41
5:H:95:ALA:HB2	5:H:100(P):MET:HG3	2.02	0.41
5:H:113:SER:OG	5:H:114:THR:N	2.54	0.41
2:D:29:PHE:CE2	2:D:52(A):PRO:HB3	2.56	0.41
6:L:167:LYS:HE3	6:L:167:LYS:HB2	1.95	0.41
1:B:537:LEU:O	1:B:541:ALA:N	2.50	0.41
2:D:147:PRO:HD2	2:D:200:HIS:CE1	2.56	0.41
5:H:55:GLU:HG2	5:H:71:ARG:NH2	2.36	0.41
6:L:141:TYR:HB3	6:L:198:HIS:NE2	2.36	0.41
6:L:198:HIS:ND1	6:L:198:HIS:O	2.53	0.41
1:B:593:LEU:HD11	4:G:494:LEU:HD21	2.03	0.41
1:B:635:ILE:HD12	1:B:635:ILE:HA	1.80	0.41
2:D:184:VAL:HG11	2:D:194:TYR:CE1	2.56	0.41
3:E:120:PRO:HD3	3:E:132:LEU:HD23	2.03	0.41
4:G:50:THR:O	4:G:103:GLN:NE2	2.54	0.41
4:G:361:PHE:CD1	4:G:361:PHE:C	2.99	0.41
4:G:427:TRP:CD1	4:G:475:MET:HG2	2.50	0.41
5:H:100(D):VAL:HG12	5:H:100(F):SER:H	1.85	0.41
2:D:30:ASN:HA	2:D:52(A):PRO:HB2	2.02	0.40
3:E:66:LYS:HA	3:E:71:ALA:HA	2.03	0.40
3:E:204:LYS:HD3	3:E:204:LYS:HA	1.91	0.40
3:E:209:THR:HG23	3:E:210:GLU:HG3	2.03	0.40
4:G:98:ASN:HD21	4:G:100:MET:HB2	1.86	0.40
5:H:144:PHE:HA	5:H:173:LEU:HG	2.03	0.40
2:D:82(C):LEU:HD22	2:D:111:VAL:HG21	2.03	0.40
4:G:47:GLU:HA	4:G:489:VAL:HG12	2.03	0.40
4:G:439:ILE:HB	4:G:443:ILE:HD11	2.02	0.40
1:B:606:THR:HG22	1:B:650:GLN:HE22	1.86	0.40
4:G:361:PHE:CA	4:G:467:THR:HB	2.47	0.40
11:O:1:NAG:H62	11:O:2:NAG:N2	2.37	0.40
4:G:136:THR:HG22	4:G:137:ASN:H	1.86	0.40
4:G:346:VAL:O	4:G:350:GLU:HB2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:60:ASN:ND2	6:L:96:TRP:HB3	2.37	0.40

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	B	129/153 (84%)	116 (90%)	12 (9%)	1 (1%)	16	50
2	D	238/243 (98%)	226 (95%)	12 (5%)	0	100	100
3	E	211/216 (98%)	198 (94%)	13 (6%)	0	100	100
4	G	448/482 (93%)	413 (92%)	31 (7%)	4 (1%)	14	48
5	H	227/236 (96%)	211 (93%)	15 (7%)	1 (0%)	30	62
6	L	209/214 (98%)	194 (93%)	13 (6%)	2 (1%)	12	45
All	All	1462/1544 (95%)	1358 (93%)	96 (7%)	8 (0%)	24	58

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	G	358	THR
4	G	369	PRO
5	H	125	SER
6	L	142	PRO
6	L	108	SER
1	B	530	MET
4	G	234	ASN
4	G	465	THR

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	B	113/130 (87%)	109 (96%)	4 (4%)	32	57
2	D	203/206 (98%)	202 (100%)	1 (0%)	81	81
3	E	186/189 (98%)	183 (98%)	3 (2%)	55	70
4	G	403/427 (94%)	377 (94%)	26 (6%)	15	42
5	H	199/204 (98%)	193 (97%)	6 (3%)	36	60
6	L	177/180 (98%)	174 (98%)	3 (2%)	53	70
All	All	1281/1336 (96%)	1238 (97%)	43 (3%)	32	58

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	593	LEU
1	B	637	ASN
1	B	651	ILE
1	B	653	GLN
2	D	54	SER
3	E	28	CYS
3	E	92	THR
3	E	96	CYS
4	G	36	VAL
4	G	38	VAL
4	G	54	CYS
4	G	68	VAL
4	G	74	CYS
4	G	86	LEU
4	G	127	VAL
4	G	130	ASN
4	G	136	THR
4	G	160	ASN
4	G	188	ILE
4	G	215	ILE
4	G	270	ILE
4	G	342	LEU

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Mol	Chain	Res	Type
4	G	350	GLU
4	G	359	ILE
4	G	360	ILE
4	G	361	PHE
4	G	362	ASN
4	G	372	VAL
4	G	378	CYS
4	G	404	ASP
4	G	418	CYS
4	G	448	ASN
4	G	452	LEU
4	G	465	THR
5	H	82	LEU
5	H	82(A)	ARG
5	H	82(B)	SER
5	H	82(C)	VAL
5	H	141	LYS
5	H	167	VAL
6	L	145	VAL
6	L	198	HIS
6	L	203	VAL

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

Mol	Chain	Res	Type
1	B	607	ASN
3	E	170	ASN
4	G	132	ASN
4	G	190	ASN
4	G	195	ASN
4	G	203	GLN
4	G	229	ASN
4	G	328	GLN
4	G	330	HIS
4	G	344	GLN
4	G	422	GLN
6	L	42	GLN
6	L	189	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 ⓘ

59 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	$\# Z > 2$
7	NAG	A	1	1,7	14,14,15	0.33	0	17,19,21	1.06	2 (11%)
7	NAG	A	2	7	14,14,15	0.29	0	17,19,21	1.32	1 (5%)
7	BMA	A	3	7	11,11,12	0.49	0	15,15,17	1.73	4 (26%)
7	MAN	A	4	7	11,11,12	1.22	1 (9%)	15,15,17	0.87	1 (6%)
7	MAN	A	5	7	11,11,12	1.21	0	15,15,17	0.80	0
7	MAN	A	6	7	11,11,12	0.78	0	15,15,17	1.16	2 (13%)
8	NAG	C	1	1,8	14,14,15	0.42	0	17,19,21	1.41	2 (11%)
8	NAG	C	2	8	14,14,15	0.36	0	17,19,21	0.54	0
9	NAG	F	1	4,9	14,14,15	0.37	0	17,19,21	1.04	1 (5%)
9	NAG	F	2	9	14,14,15	0.66	0	17,19,21	1.09	2 (11%)
9	BMA	F	3	9	11,11,12	0.43	0	15,15,17	0.98	0
9	MAN	F	4	9	11,11,12	2.08	4 (36%)	15,15,17	1.58	4 (26%)
9	MAN	F	5	9	11,11,12	0.77	1 (9%)	15,15,17	1.20	2 (13%)
9	MAN	F	6	9	11,11,12	0.55	0	15,15,17	1.23	2 (13%)
9	MAN	F	7	9	11,11,12	0.68	0	15,15,17	0.89	1 (6%)
8	NAG	I	1	4,8	14,14,15	0.30	0	17,19,21	0.70	0
8	NAG	I	2	8	14,14,15	0.30	0	17,19,21	0.61	0
10	NAG	J	1	4,10	14,14,15	0.34	0	17,19,21	0.84	0
10	NAG	J	2	10	14,14,15	0.43	0	17,19,21	1.70	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	BMA	J	3	10	11,11,12	0.29	0	15,15,17	0.63	0
10	MAN	J	4	10	11,11,12	0.71	0	15,15,17	0.92	2 (13%)
11	NAG	K	1	4,11	14,14,15	0.35	0	17,19,21	0.85	0
11	NAG	K	2	11	14,14,15	0.47	0	17,19,21	1.48	3 (17%)
11	BMA	K	3	11	11,11,12	0.22	0	15,15,17	0.56	0
10	NAG	M	1	4,10	14,14,15	0.30	0	17,19,21	0.76	0
10	NAG	M	2	10	14,14,15	0.35	0	17,19,21	0.75	1 (5%)
10	BMA	M	3	10	11,11,12	0.27	0	15,15,17	0.51	0
10	MAN	M	4	10	11,11,12	0.81	0	15,15,17	1.13	1 (6%)
11	NAG	N	1	4,11	14,14,15	0.36	0	17,19,21	0.95	1 (5%)
11	NAG	N	2	11	14,14,15	0.32	0	17,19,21	0.69	0
11	BMA	N	3	11	11,11,12	0.26	0	15,15,17	0.65	0
11	NAG	O	1	4,11	14,14,15	0.28	0	17,19,21	0.56	0
11	NAG	O	2	11	14,14,15	0.28	0	17,19,21	0.84	0
11	BMA	O	3	11	11,11,12	0.30	0	15,15,17	0.69	0
12	NAG	P	1	4,12	14,14,15	0.29	0	17,19,21	1.23	2 (11%)
12	NAG	P	2	12	14,14,15	0.33	0	17,19,21	1.07	2 (11%)
12	BMA	P	3	12	11,11,12	0.62	0	15,15,17	1.58	3 (20%)
12	MAN	P	4	12	11,11,12	0.84	0	15,15,17	1.58	2 (13%)
12	MAN	P	5	12	11,11,12	0.98	0	15,15,17	1.40	2 (13%)
12	MAN	P	6	12	11,11,12	0.77	0	15,15,17	1.11	1 (6%)
12	MAN	P	7	12	11,11,12	0.77	0	15,15,17	0.87	1 (6%)
12	MAN	P	8	12	11,11,12	0.89	0	15,15,17	0.93	0
8	NAG	Q	1	4,8	14,14,15	0.36	0	17,19,21	1.06	2 (11%)
8	NAG	Q	2	8	14,14,15	0.30	0	17,19,21	0.91	0
13	NAG	R	1	13,4	14,14,15	0.30	0	17,19,21	1.02	1 (5%)
13	NAG	R	2	13	14,14,15	0.42	0	17,19,21	0.88	1 (5%)
13	BMA	R	3	13	11,11,12	0.42	0	15,15,17	0.80	0
13	MAN	R	4	13	11,11,12	0.85	0	15,15,17	0.94	1 (6%)
14	NAG	S	1	4,14	14,14,15	0.27	0	17,19,21	0.61	0
14	NAG	S	2	14	14,14,15	0.33	0	17,19,21	0.62	0
14	BMA	S	3	14	11,11,12	0.32	0	15,15,17	0.79	0
14	MAN	S	4	14	11,11,12	0.75	0	15,15,17	0.85	1 (6%)
14	MAN	S	5	14	11,11,12	0.62	0	15,15,17	1.06	2 (13%)
15	NAG	T	1	15,4	14,14,15	0.39	0	17,19,21	0.89	0
15	NAG	T	2	15	14,14,15	0.58	0	17,19,21	2.07	4 (23%)
15	BMA	T	3	15	11,11,12	0.29	0	15,15,17	0.70	0
11	NAG	U	1	4,11	14,14,15	0.31	0	17,19,21	0.74	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	NAG	U	2	11	14,14,15	0.29	0	17,19,21	0.67	0
11	BMA	U	3	11	11,11,12	0.27	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
7	NAG	A	1	1,7	-	2/6/23/26	0/1/1/1
7	NAG	A	2	7	-	2/6/23/26	0/1/1/1
7	BMA	A	3	7	-	1/2/19/22	0/1/1/1
7	MAN	A	4	7	-	2/2/19/22	0/1/1/1
7	MAN	A	5	7	-	2/2/19/22	0/1/1/1
7	MAN	A	6	7	-	0/2/19/22	0/1/1/1
8	NAG	C	1	1,8	-	3/6/23/26	0/1/1/1
8	NAG	C	2	8	-	0/6/23/26	0/1/1/1
9	NAG	F	1	4,9	-	4/6/23/26	0/1/1/1
9	NAG	F	2	9	-	2/6/23/26	0/1/1/1
9	BMA	F	3	9	-	1/2/19/22	0/1/1/1
9	MAN	F	4	9	-	2/2/19/22	0/1/1/1
9	MAN	F	5	9	-	0/2/19/22	0/1/1/1
9	MAN	F	6	9	-	1/2/19/22	0/1/1/1
9	MAN	F	7	9	-	2/2/19/22	0/1/1/1
8	NAG	I	1	4,8	-	0/6/23/26	0/1/1/1
8	NAG	I	2	8	-	1/6/23/26	0/1/1/1
10	NAG	J	1	4,10	-	0/6/23/26	0/1/1/1
10	NAG	J	2	10	-	4/6/23/26	0/1/1/1
10	BMA	J	3	10	-	2/2/19/22	0/1/1/1
10	MAN	J	4	10	-	2/2/19/22	0/1/1/1
11	NAG	K	1	4,11	-	3/6/23/26	0/1/1/1
11	NAG	K	2	11	-	6/6/23/26	0/1/1/1
11	BMA	K	3	11	-	0/2/19/22	0/1/1/1
10	NAG	M	1	4,10	-	0/6/23/26	0/1/1/1
10	NAG	M	2	10	-	0/6/23/26	0/1/1/1
10	BMA	M	3	10	-	2/2/19/22	0/1/1/1
10	MAN	M	4	10	-	1/2/19/22	0/1/1/1
11	NAG	N	1	4,11	-	4/6/23/26	0/1/1/1
11	NAG	N	2	11	-	3/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	BMA	N	3	11	-	0/2/19/22	0/1/1/1
11	NAG	O	1	4,11	-	0/6/23/26	0/1/1/1
11	NAG	O	2	11	-	4/6/23/26	0/1/1/1
11	BMA	O	3	11	-	0/2/19/22	0/1/1/1
12	NAG	P	1	4,12	-	3/6/23/26	0/1/1/1
12	NAG	P	2	12	-	1/6/23/26	0/1/1/1
12	BMA	P	3	12	-	0/2/19/22	0/1/1/1
12	MAN	P	4	12	-	2/2/19/22	0/1/1/1
12	MAN	P	5	12	-	0/2/19/22	0/1/1/1
12	MAN	P	6	12	-	2/2/19/22	0/1/1/1
12	MAN	P	7	12	-	2/2/19/22	0/1/1/1
12	MAN	P	8	12	-	0/2/19/22	0/1/1/1
8	NAG	Q	1	4,8	-	3/6/23/26	0/1/1/1
8	NAG	Q	2	8	-	3/6/23/26	0/1/1/1
13	NAG	R	1	13,4	-	3/6/23/26	0/1/1/1
13	NAG	R	2	13	-	2/6/23/26	0/1/1/1
13	BMA	R	3	13	-	1/2/19/22	0/1/1/1
13	MAN	R	4	13	-	1/2/19/22	0/1/1/1
14	NAG	S	1	4,14	-	2/6/23/26	0/1/1/1
14	NAG	S	2	14	-	2/6/23/26	0/1/1/1
14	BMA	S	3	14	-	2/2/19/22	0/1/1/1
14	MAN	S	4	14	-	0/2/19/22	0/1/1/1
14	MAN	S	5	14	-	0/2/19/22	0/1/1/1
15	NAG	T	1	15,4	-	2/6/23/26	0/1/1/1
15	NAG	T	2	15	-	5/6/23/26	0/1/1/1
15	BMA	T	3	15	-	2/2/19/22	0/1/1/1
11	NAG	U	1	4,11	-	1/6/23/26	0/1/1/1
11	NAG	U	2	11	-	0/6/23/26	0/1/1/1
11	BMA	U	3	11	-	0/2/19/22	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	F	4	MAN	O2-C2	4.29	1.52	1.43
9	F	4	MAN	O5-C1	-3.17	1.38	1.43
9	F	4	MAN	C2-C3	3.07	1.57	1.52
7	A	4	MAN	O5-C1	-2.87	1.38	1.43
9	F	4	MAN	C4-C3	2.47	1.58	1.52
9	F	5	MAN	C1-C2	2.46	1.58	1.52

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	T	2	NAG	C2-N2-C7	6.03	130.98	122.90
10	J	2	NAG	C2-N2-C7	5.40	130.14	122.90
12	P	4	MAN	C1-O5-C5	4.46	118.17	112.19
12	P	3	BMA	O3-C3-C2	-4.44	101.00	110.05
9	F	6	MAN	C1-O5-C5	3.81	117.29	112.19
15	T	2	NAG	C1-C2-N2	3.78	116.40	110.43
7	A	3	BMA	O3-C3-C4	-3.57	101.96	110.38
12	P	5	MAN	C1-C2-C3	-3.55	104.48	109.64
7	A	2	NAG	O4-C4-C3	-3.45	102.23	110.38
10	M	4	MAN	C1-O5-C5	3.45	116.81	112.19
7	A	6	MAN	C1-O5-C5	3.42	116.77	112.19
8	C	1	NAG	O4-C4-C3	-3.38	102.40	110.38
11	K	2	NAG	C4-C3-C2	3.33	115.90	111.02
12	P	2	NAG	O4-C4-C5	-3.18	101.50	109.32
8	C	1	NAG	O4-C4-C5	-3.17	101.51	109.32
14	S	5	MAN	C1-O5-C5	3.15	116.41	112.19
12	P	5	MAN	C1-O5-C5	3.11	116.35	112.19
12	P	6	MAN	C1-O5-C5	3.07	116.30	112.19
11	K	2	NAG	O3-C3-C2	-3.05	103.06	109.40
11	N	1	NAG	C2-N2-C7	2.95	126.85	122.90
9	F	4	MAN	O2-C2-C3	2.67	115.68	110.15
10	J	2	NAG	C8-C7-N2	2.65	120.52	116.12
12	P	4	MAN	C1-C2-C3	2.64	113.50	109.64
9	F	5	MAN	O2-C2-C3	-2.62	104.72	110.15
9	F	5	MAN	C1-O5-C5	2.62	115.69	112.19
7	A	1	NAG	C2-N2-C7	-2.53	119.50	122.90
13	R	2	NAG	O4-C4-C3	-2.52	104.43	110.38
7	A	3	BMA	C1-O5-C5	-2.49	108.85	112.19
7	A	4	MAN	O2-C2-C3	-2.47	105.03	110.15
13	R	4	MAN	C1-O5-C5	2.47	115.50	112.19
7	A	1	NAG	O4-C4-C3	-2.44	104.62	110.38
8	Q	1	NAG	O5-C1-C2	-2.38	107.61	111.29
15	T	2	NAG	C8-C7-N2	2.38	120.06	116.12
9	F	4	MAN	C1-C2-C3	-2.36	106.20	109.64
12	P	3	BMA	O5-C5-C4	-2.35	105.10	110.83
7	A	3	BMA	O2-C2-C3	-2.35	105.28	110.15
9	F	4	MAN	C2-C3-C4	2.35	114.99	110.86
9	F	7	MAN	O2-C2-C3	-2.32	105.34	110.15
9	F	1	NAG	O5-C1-C2	-2.31	107.72	111.29
8	Q	1	NAG	C1-O5-C5	2.29	115.26	112.19
7	A	3	BMA	O3-C3-C2	-2.26	105.45	110.05
12	P	1	NAG	O4-C4-C5	-2.26	103.77	109.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	P	3	BMA	C1-C2-C3	2.21	112.87	109.64
15	T	2	NAG	O5-C1-C2	-2.21	107.87	111.29
12	P	7	MAN	O2-C2-C3	-2.19	105.62	110.15
12	P	2	NAG	O5-C5-C6	2.18	111.91	107.66
7	A	6	MAN	O2-C2-C3	-2.16	105.67	110.15
12	P	1	NAG	O5-C1-C2	-2.15	107.96	111.29
14	S	4	MAN	O2-C2-C3	-2.15	105.70	110.15
13	R	1	NAG	C2-N2-C7	2.15	125.78	122.90
9	F	4	MAN	O3-C3-C2	2.14	114.41	110.05
11	K	2	NAG	C2-N2-C7	-2.09	120.09	122.90
10	J	4	MAN	C1-O5-C5	2.09	114.98	112.19
9	F	2	NAG	O3-C3-C2	2.08	113.71	109.40
10	J	4	MAN	O2-C2-C3	-2.07	105.86	110.15
14	S	5	MAN	O2-C2-C3	-2.03	105.95	110.15
10	M	2	NAG	O5-C1-C2	-2.02	108.17	111.29
9	F	2	NAG	O4-C4-C5	-2.01	104.37	109.32
9	F	6	MAN	O2-C2-C3	-2.01	105.99	110.15

There are no chirality outliers.

All (95) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	2	NAG	C1-C2-N2-C7
8	Q	1	NAG	O7-C7-N2-C2
8	Q	2	NAG	C3-C2-N2-C7
10	J	2	NAG	C3-C2-N2-C7
10	J	2	NAG	C8-C7-N2-C2
10	J	2	NAG	O7-C7-N2-C2
11	K	1	NAG	C3-C2-N2-C7
11	K	1	NAG	C8-C7-N2-C2
11	K	1	NAG	O7-C7-N2-C2
11	K	2	NAG	C8-C7-N2-C2
11	K	2	NAG	O7-C7-N2-C2
12	P	1	NAG	C1-C2-N2-C7
15	T	2	NAG	C1-C2-N2-C7
15	T	2	NAG	C8-C7-N2-C2
15	T	2	NAG	O7-C7-N2-C2
8	Q	1	NAG	C8-C7-N2-C2
7	A	1	NAG	O5-C5-C6-O6
10	M	3	BMA	C4-C5-C6-O6
12	P	4	MAN	O5-C5-C6-O6
13	R	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
12	P	7	MAN	O5-C5-C6-O6
13	R	2	NAG	O5-C5-C6-O6
15	T	3	BMA	O5-C5-C6-O6
7	A	1	NAG	C4-C5-C6-O6
9	F	4	MAN	C4-C5-C6-O6
9	F	7	MAN	O5-C5-C6-O6
8	C	1	NAG	C4-C5-C6-O6
14	S	3	BMA	O5-C5-C6-O6
9	F	1	NAG	C4-C5-C6-O6
14	S	3	BMA	C4-C5-C6-O6
8	C	1	NAG	O5-C5-C6-O6
15	T	2	NAG	O5-C5-C6-O6
9	F	4	MAN	O5-C5-C6-O6
10	M	3	BMA	O5-C5-C6-O6
7	A	4	MAN	O5-C5-C6-O6
8	Q	2	NAG	C8-C7-N2-C2
11	N	1	NAG	C8-C7-N2-C2
11	N	1	NAG	O7-C7-N2-C2
13	R	1	NAG	C4-C5-C6-O6
15	T	3	BMA	C4-C5-C6-O6
12	P	6	MAN	O5-C5-C6-O6
15	T	2	NAG	C4-C5-C6-O6
9	F	1	NAG	O5-C5-C6-O6
11	N	1	NAG	O5-C5-C6-O6
12	P	7	MAN	C4-C5-C6-O6
11	K	2	NAG	O5-C5-C6-O6
11	N	1	NAG	C4-C5-C6-O6
8	Q	2	NAG	O7-C7-N2-C2
9	F	1	NAG	C8-C7-N2-C2
11	N	2	NAG	C8-C7-N2-C2
11	N	2	NAG	O7-C7-N2-C2
11	O	2	NAG	C8-C7-N2-C2
11	O	2	NAG	O7-C7-N2-C2
14	S	1	NAG	C8-C7-N2-C2
14	S	1	NAG	O7-C7-N2-C2
14	S	2	NAG	C8-C7-N2-C2
15	T	1	NAG	C8-C7-N2-C2
15	T	1	NAG	O7-C7-N2-C2
10	J	4	MAN	O5-C5-C6-O6
13	R	2	NAG	C4-C5-C6-O6
9	F	7	MAN	C4-C5-C6-O6
11	O	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
9	F	1	NAG	O7-C7-N2-C2
14	S	2	NAG	O7-C7-N2-C2
12	P	4	MAN	C4-C5-C6-O6
10	J	3	BMA	O5-C5-C6-O6
12	P	2	NAG	O5-C5-C6-O6
10	J	2	NAG	O5-C5-C6-O6
11	U	1	NAG	O5-C5-C6-O6
12	P	6	MAN	C4-C5-C6-O6
7	A	3	BMA	O5-C5-C6-O6
7	A	2	NAG	O5-C5-C6-O6
11	O	2	NAG	O5-C5-C6-O6
13	R	3	BMA	O5-C5-C6-O6
11	N	2	NAG	O5-C5-C6-O6
11	K	2	NAG	C3-C2-N2-C7
8	Q	1	NAG	O5-C5-C6-O6
9	F	6	MAN	O5-C5-C6-O6
7	A	4	MAN	C4-C5-C6-O6
7	A	5	MAN	C4-C5-C6-O6
7	A	5	MAN	O5-C5-C6-O6
11	K	2	NAG	C1-C2-N2-C7
10	M	4	MAN	O5-C5-C6-O6
10	J	3	BMA	C4-C5-C6-O6
9	F	2	NAG	C8-C7-N2-C2
10	J	4	MAN	C4-C5-C6-O6
8	C	1	NAG	C3-C2-N2-C7
11	K	2	NAG	C4-C5-C6-O6
9	F	2	NAG	O7-C7-N2-C2
9	F	3	BMA	C4-C5-C6-O6
12	P	1	NAG	C8-C7-N2-C2
12	P	1	NAG	O7-C7-N2-C2
13	R	1	NAG	C3-C2-N2-C7
8	I	2	NAG	C4-C5-C6-O6
13	R	4	MAN	C4-C5-C6-O6

There are no ring outliers.

13 monomers are involved in 23 short contacts:

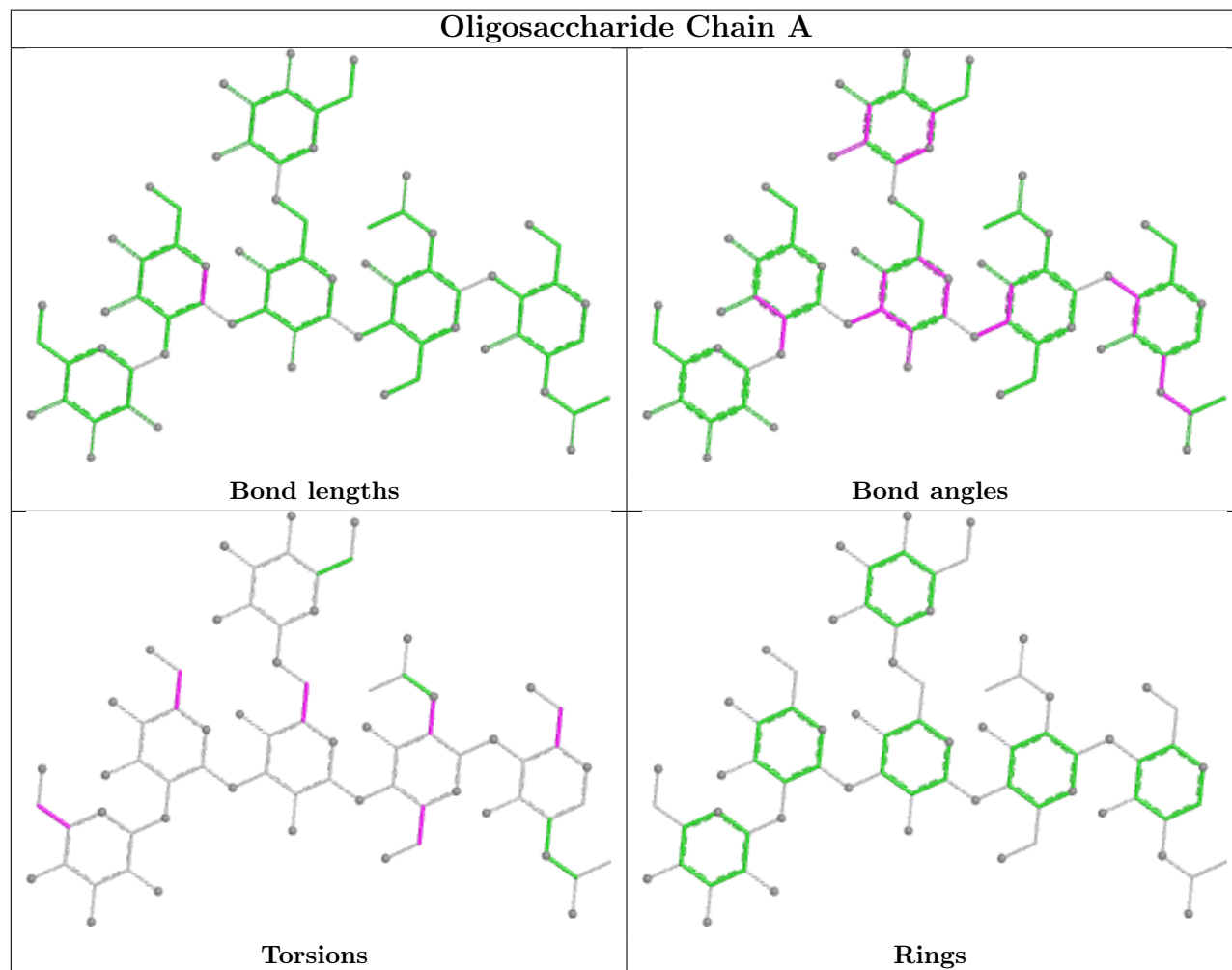
Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	K	2	NAG	2	0
12	P	1	NAG	4	0
11	N	2	NAG	1	0
8	Q	1	NAG	2	0

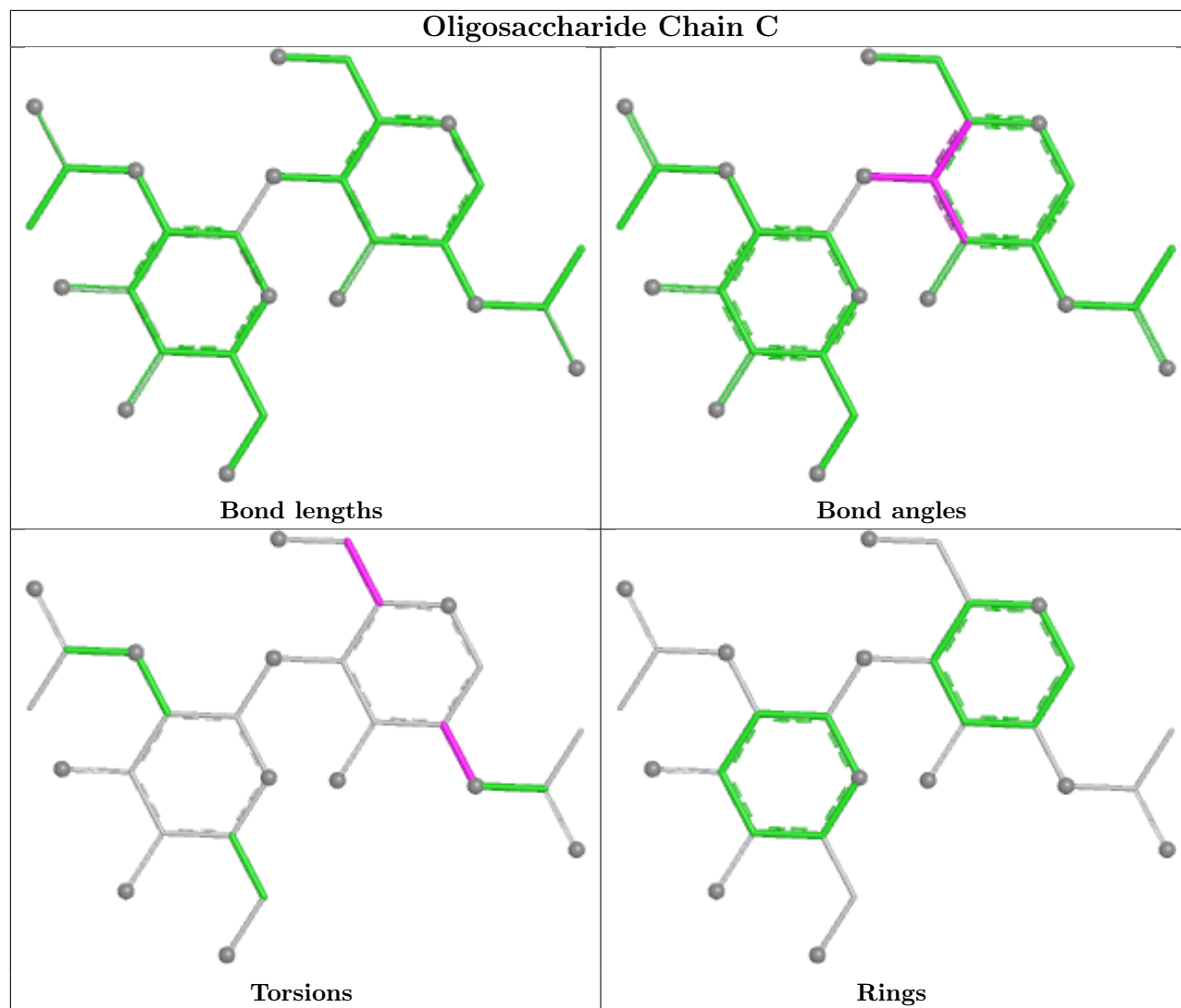
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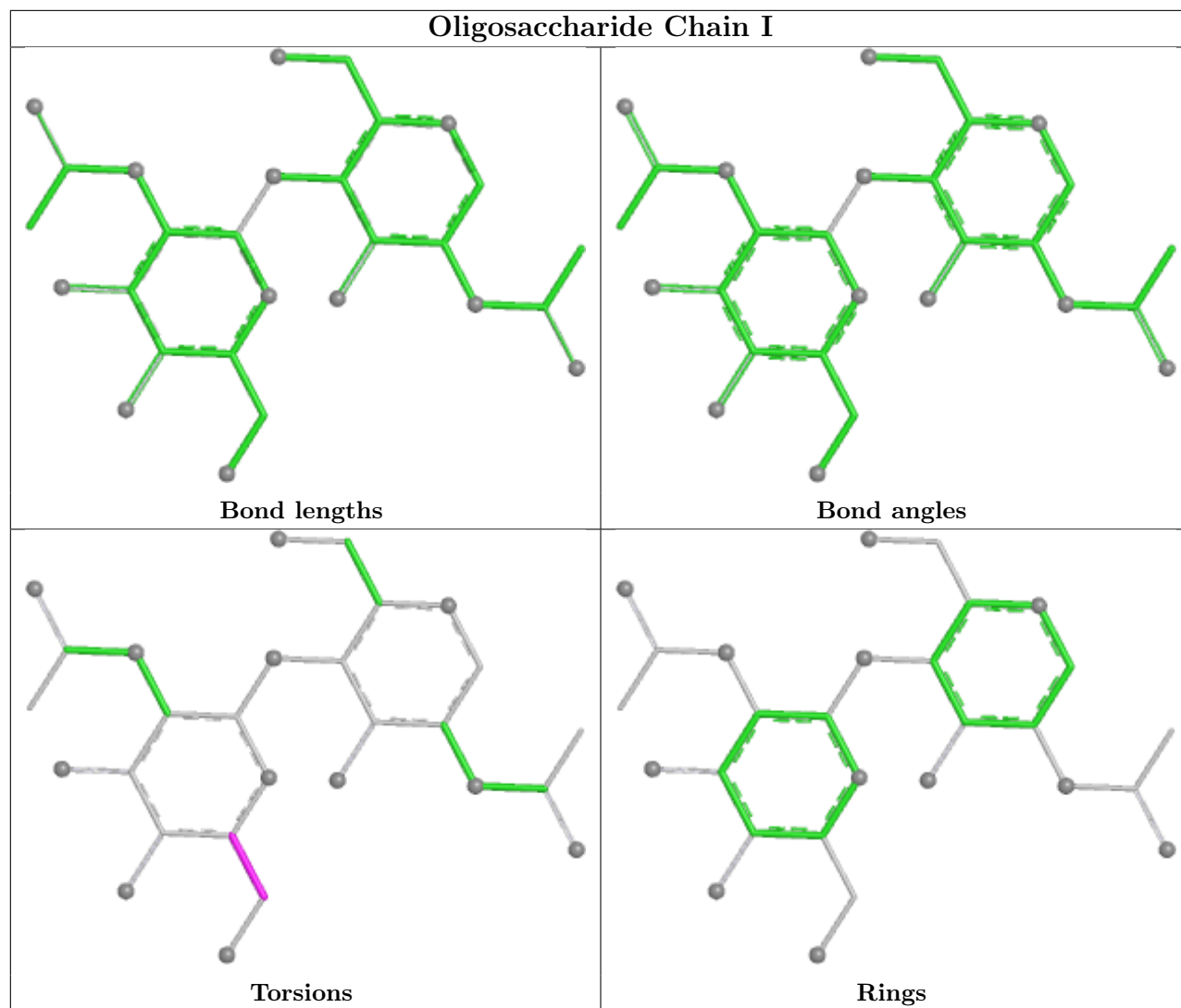
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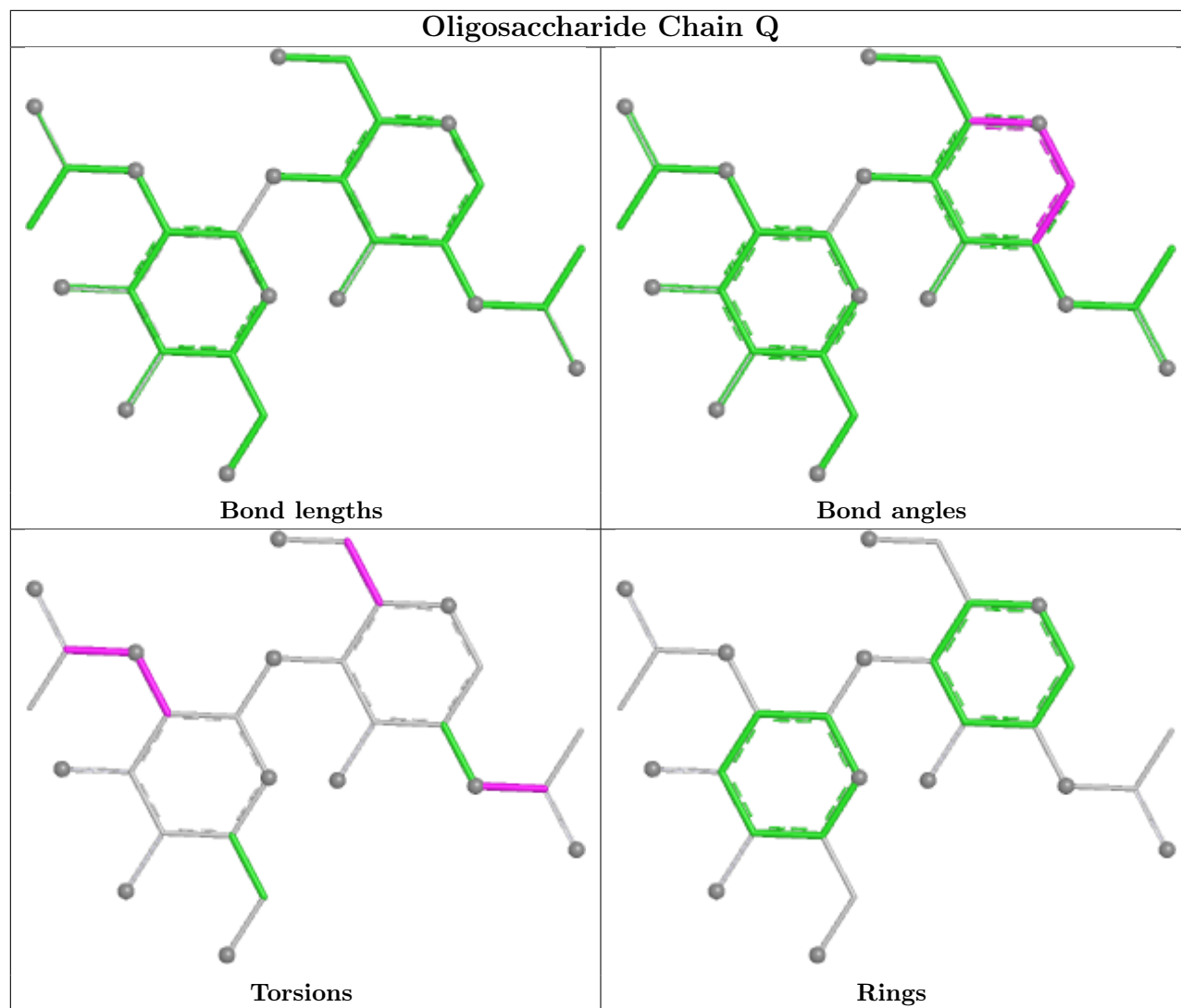
Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	S	1	NAG	1	0
11	O	1	NAG	2	0
9	F	6	MAN	1	0
8	Q	2	NAG	1	0
11	N	1	NAG	2	0
11	K	1	NAG	4	0
12	P	5	MAN	3	0
7	A	2	NAG	3	0
11	O	2	NAG	2	0

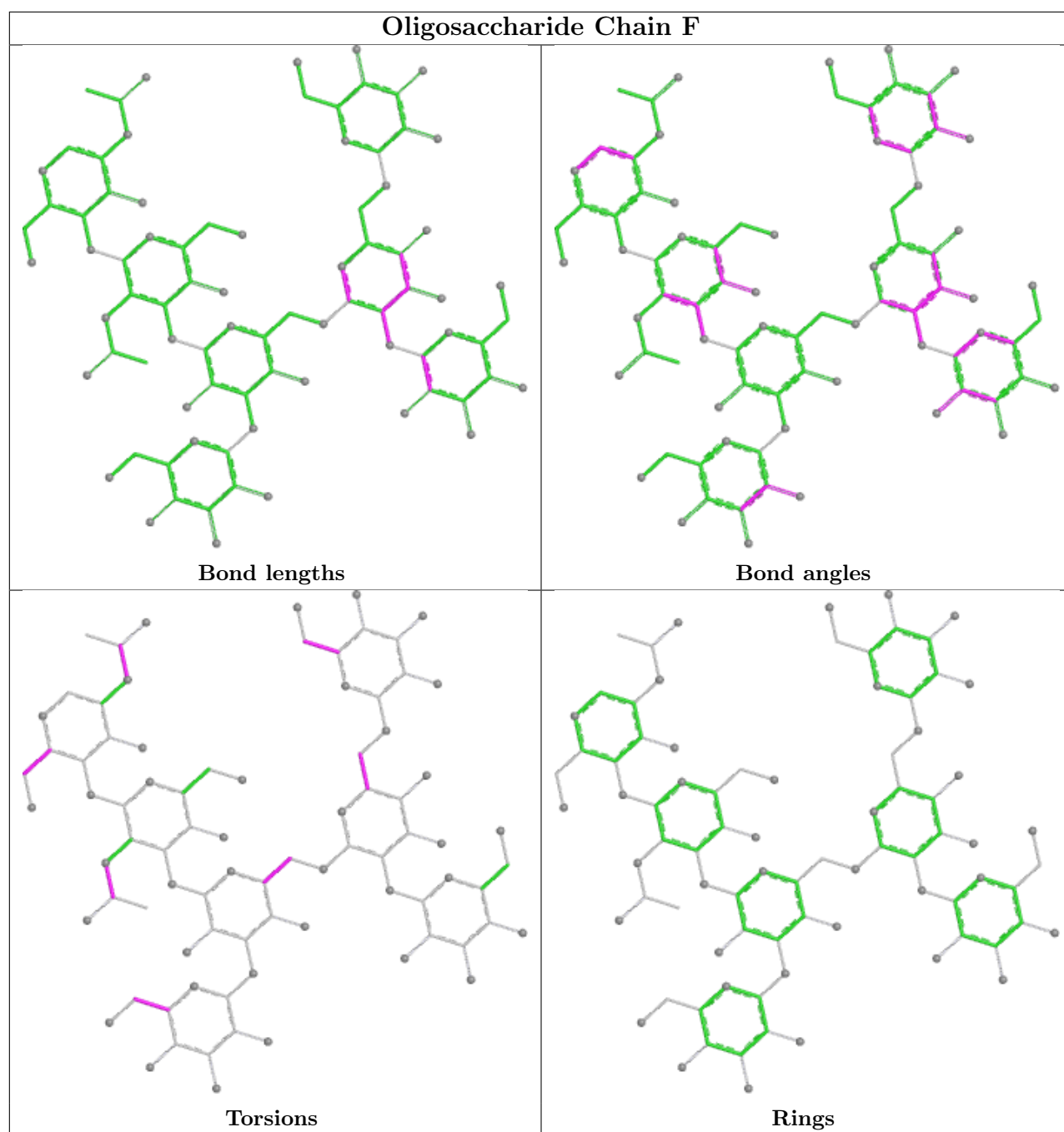
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

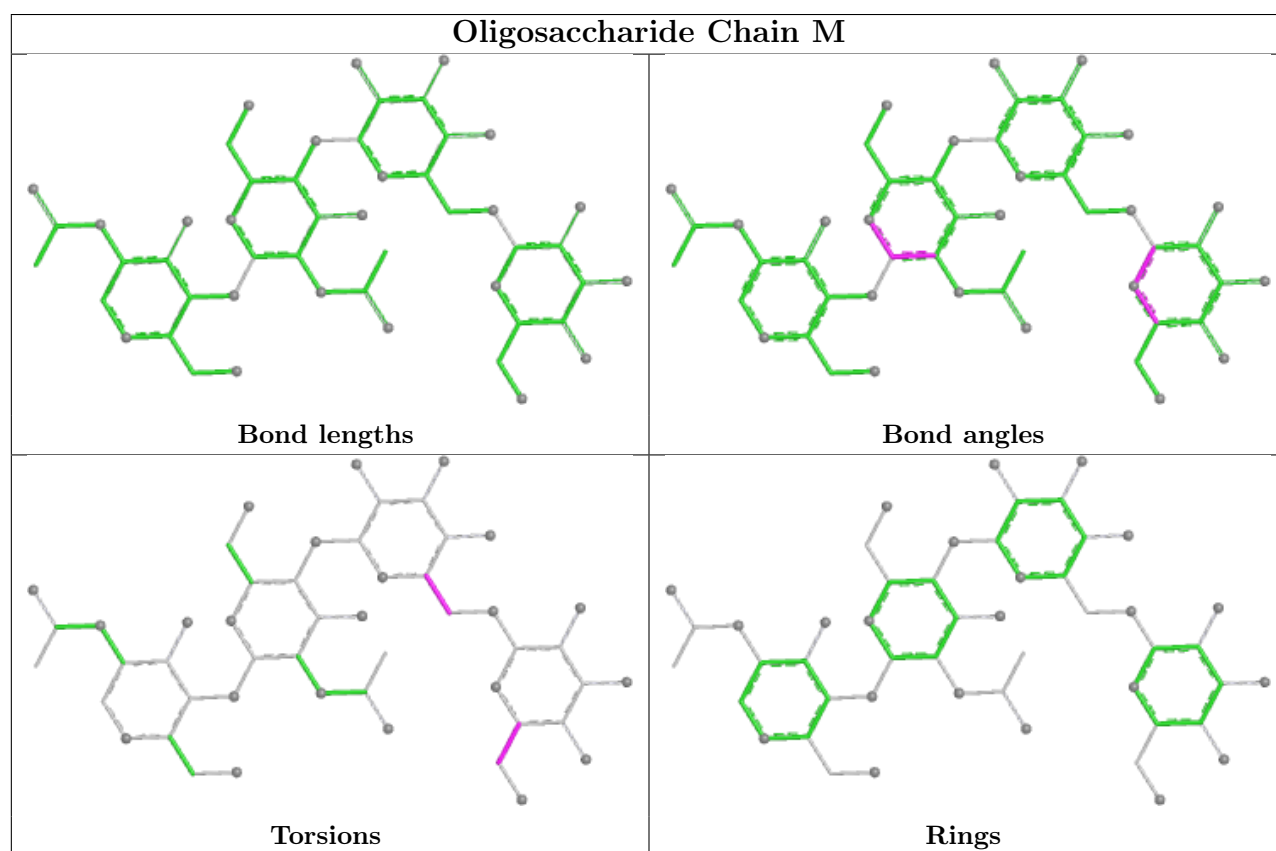
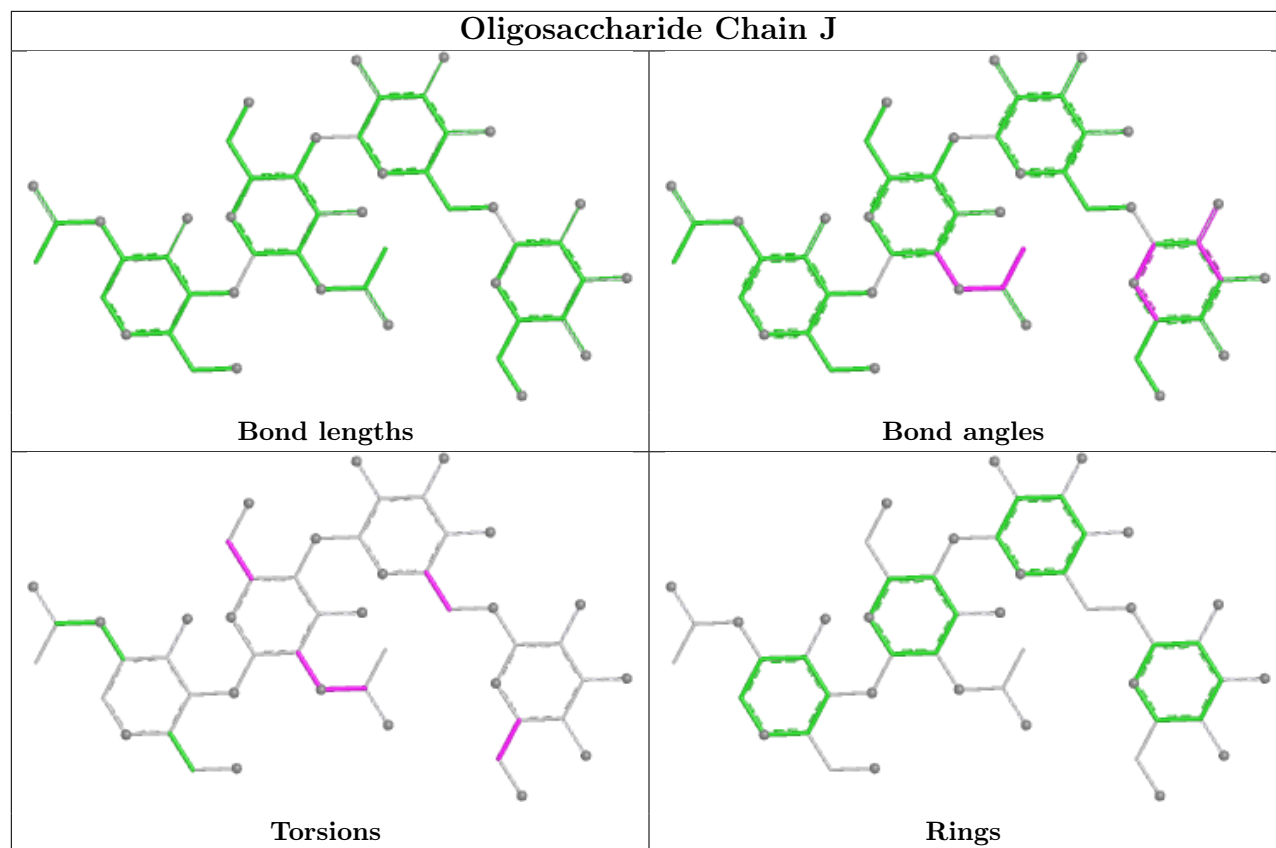


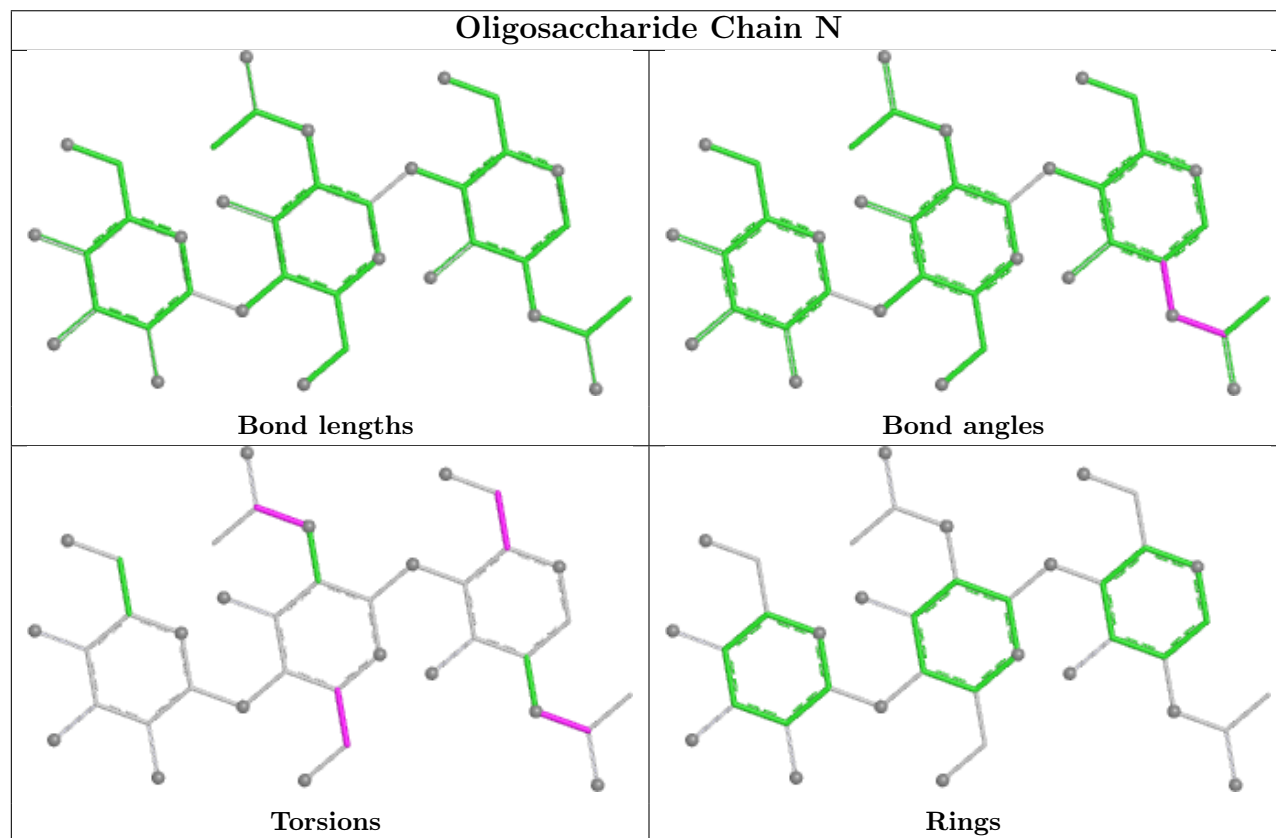
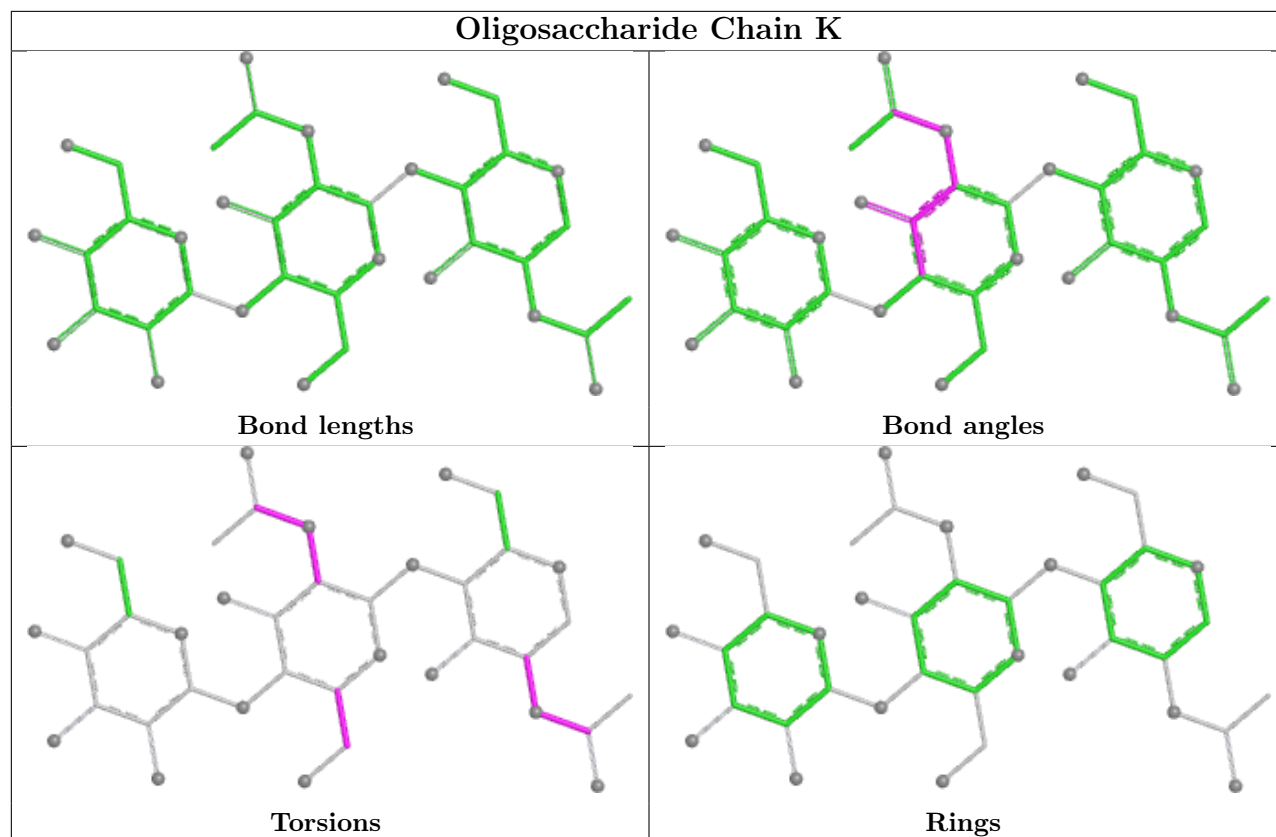


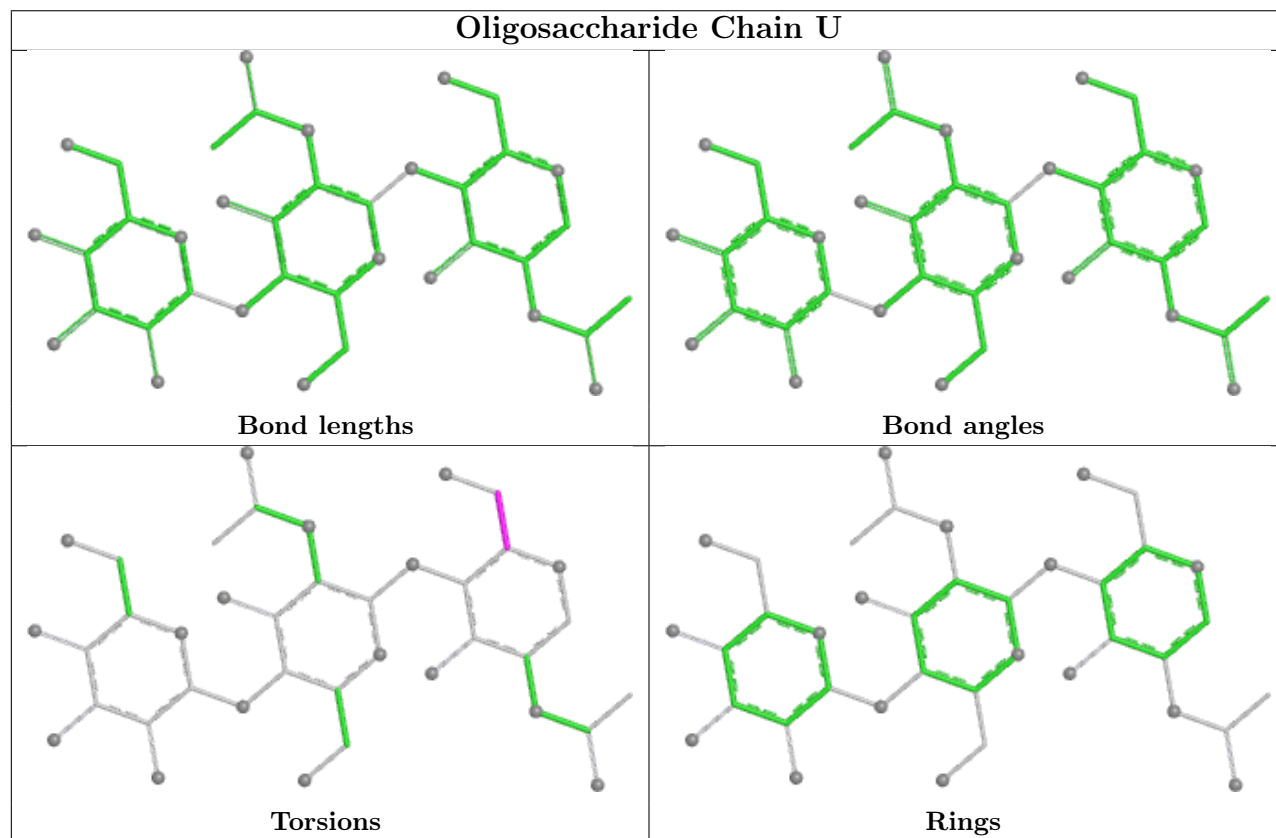
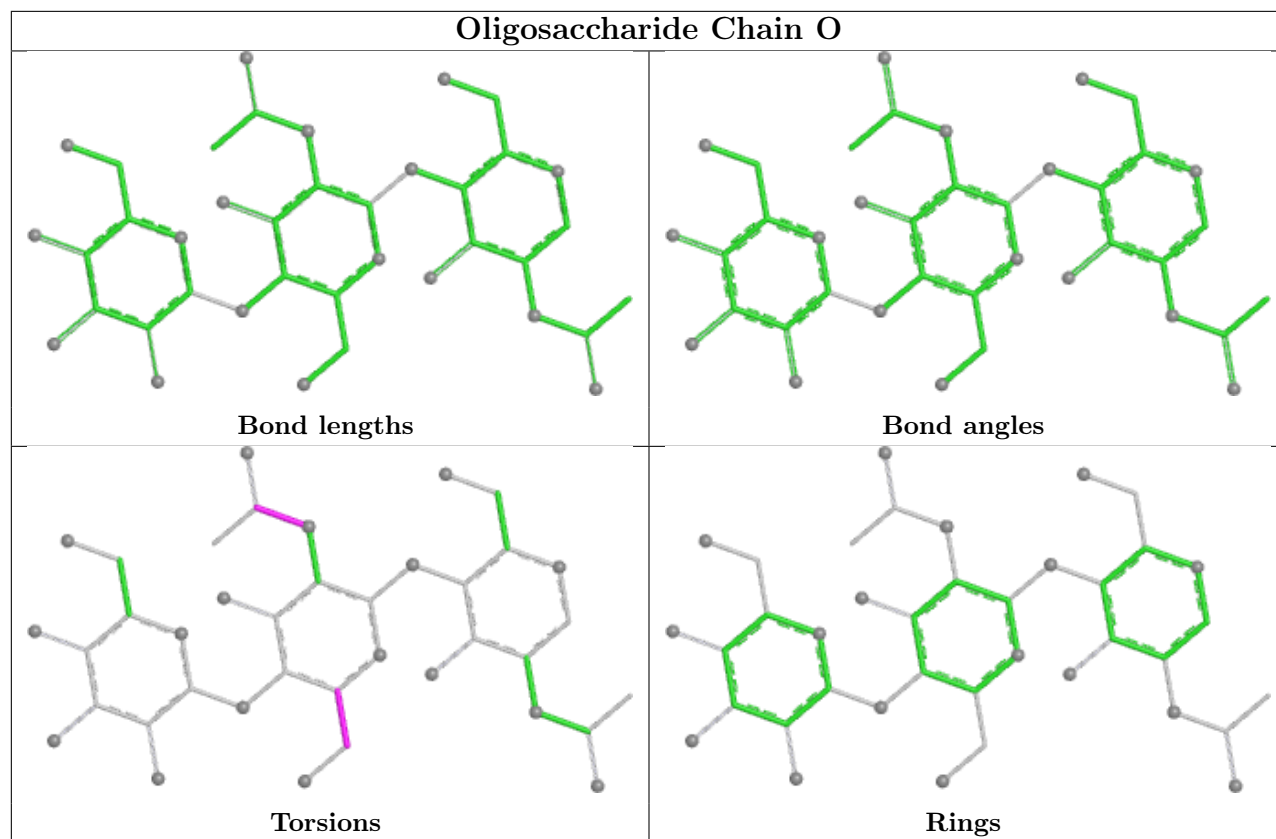


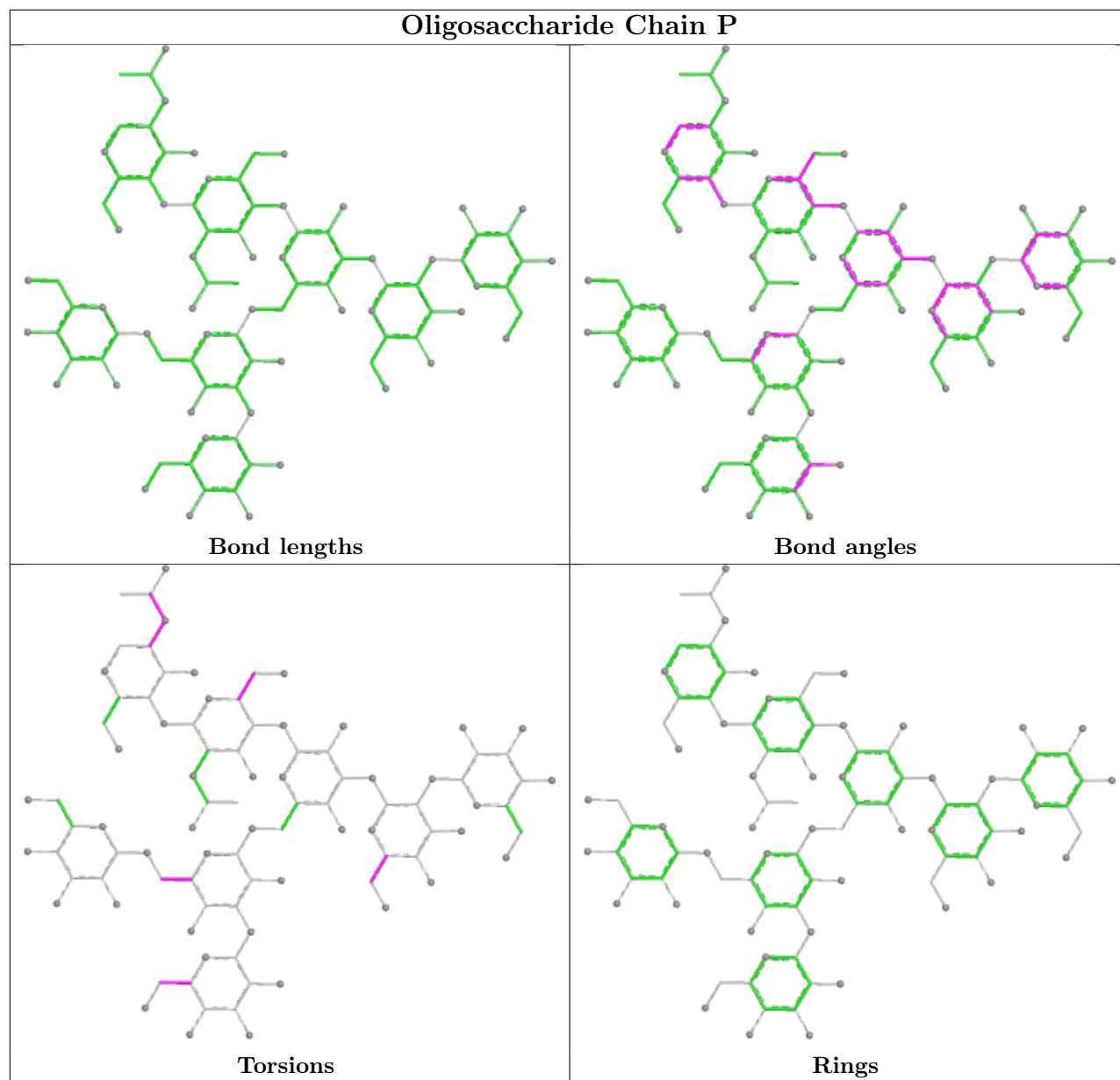


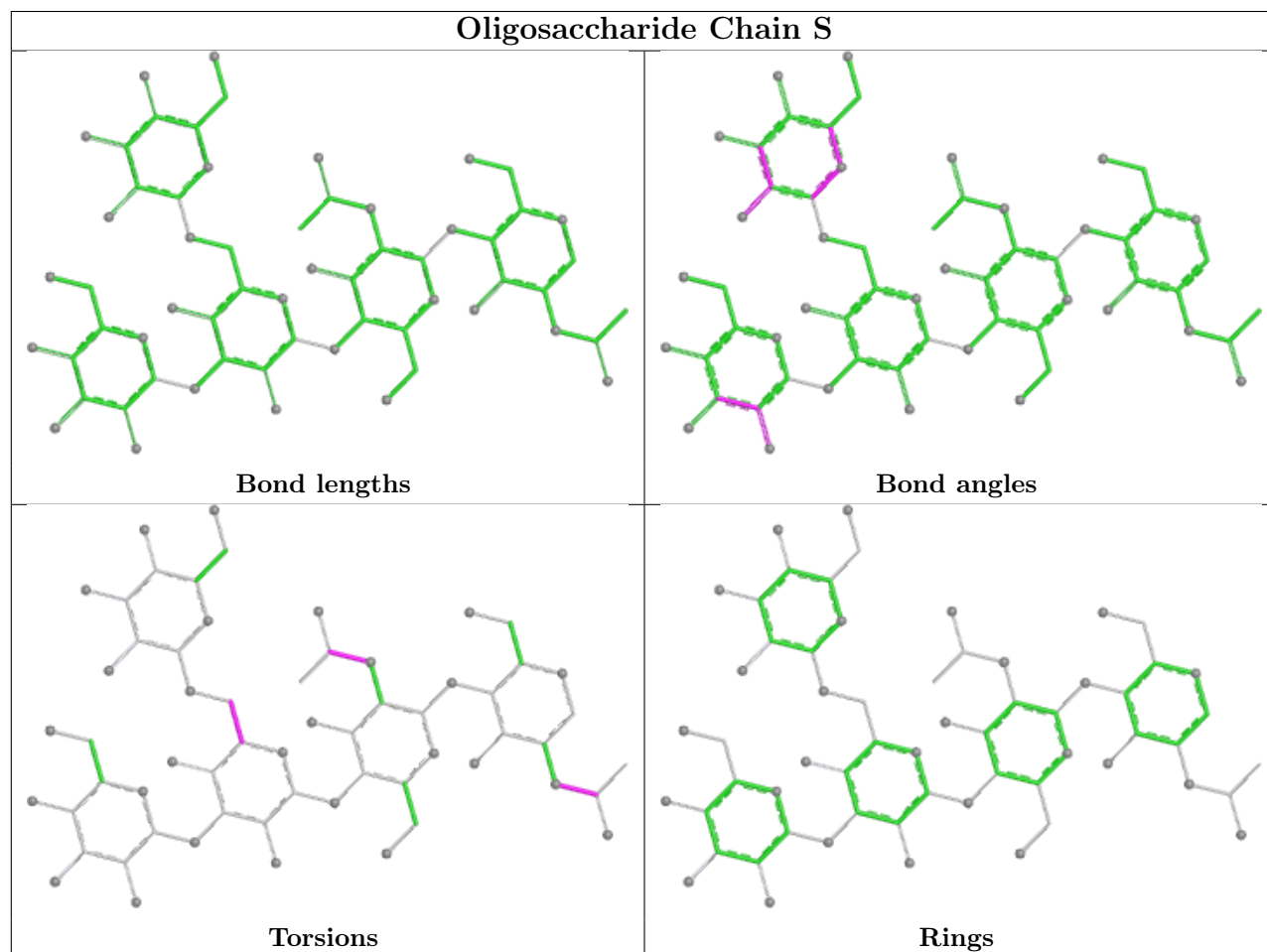
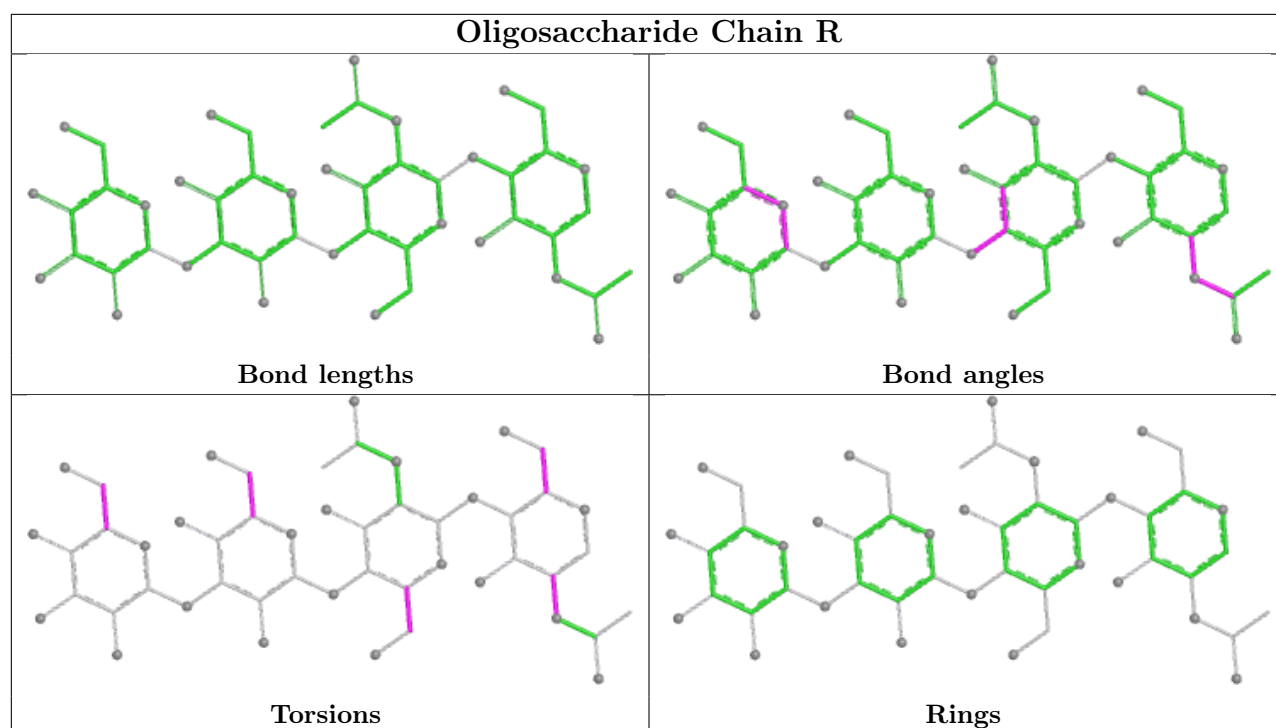


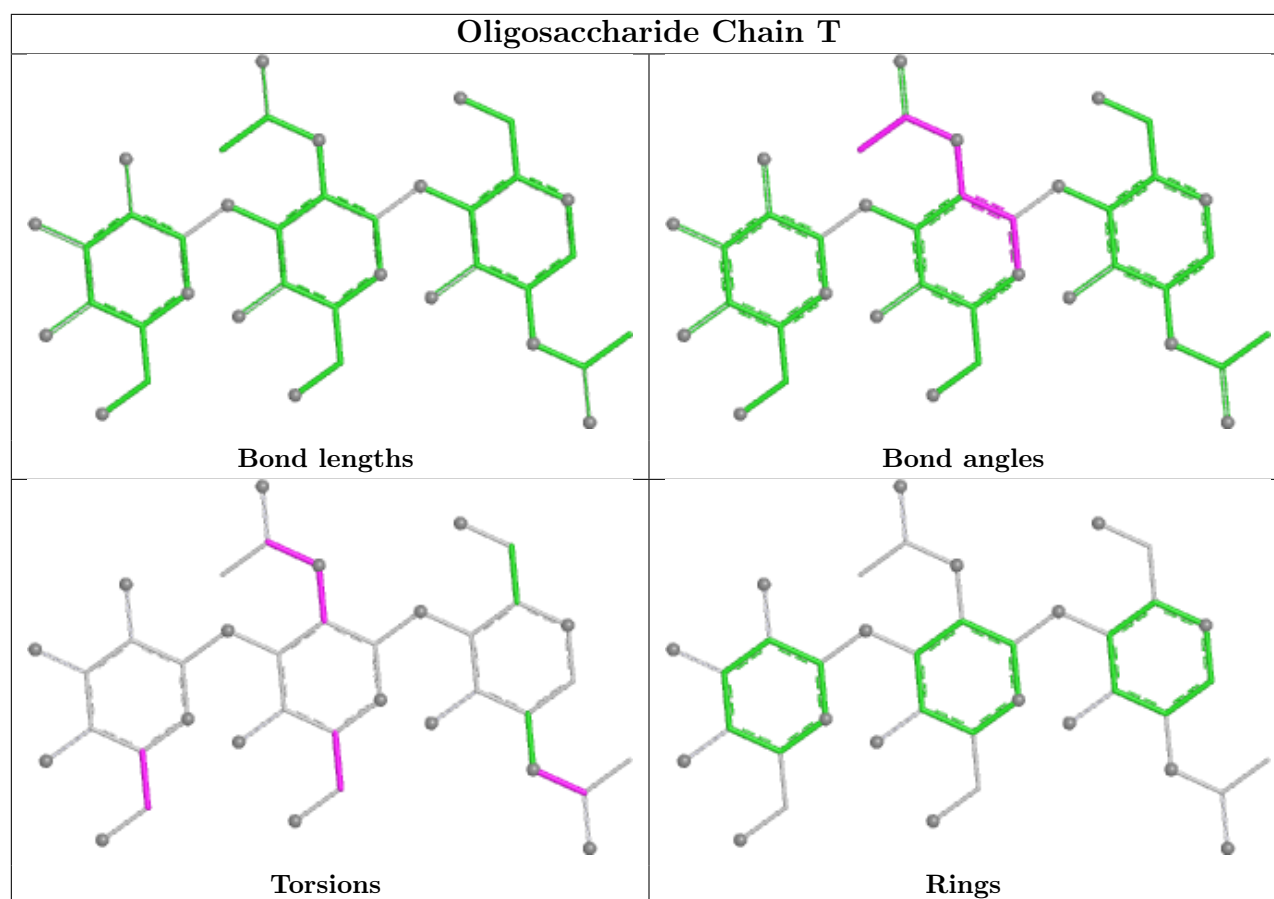












5.6 Ligand geometry [i](#)

8 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	$\# Z > 2$
16	NAG	G	603	4	14,14,15	0.27	0	17,19,21	0.51	0
16	NAG	G	605	4	14,14,15	0.28	0	17,19,21	0.52	0
16	NAG	G	601	4	14,14,15	0.29	0	17,19,21	0.67	0
16	NAG	B	701	1	14,14,15	0.27	0	17,19,21	0.76	0
16	NAG	G	606	4	14,14,15	0.52	0	17,19,21	2.02	3 (17%)
16	NAG	G	602	4	14,14,15	0.30	0	17,19,21	0.71	0
16	NAG	B	702	1	14,14,15	0.28	0	17,19,21	0.55	0
16	NAG	G	604	4	14,14,15	0.30	0	17,19,21	0.59	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
16	NAG	G	603	4	-	0/6/23/26	0/1/1/1
16	NAG	G	605	4	-	0/6/23/26	0/1/1/1
16	NAG	G	601	4	-	2/6/23/26	0/1/1/1
16	NAG	B	701	1	-	0/6/23/26	0/1/1/1
16	NAG	G	606	4	-	3/6/23/26	0/1/1/1
16	NAG	G	602	4	-	1/6/23/26	0/1/1/1
16	NAG	B	702	1	-	1/6/23/26	0/1/1/1
16	NAG	G	604	4	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	G	606	NAG	C2-N2-C7	6.57	131.70	122.90
16	G	606	NAG	C8-C7-N2	3.16	121.36	116.12
16	G	606	NAG	C1-O5-C5	2.24	115.19	112.19

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
16	G	601	NAG	C8-C7-N2-C2
16	G	606	NAG	C8-C7-N2-C2
16	G	606	NAG	O7-C7-N2-C2
16	G	601	NAG	O7-C7-N2-C2
16	B	702	NAG	O5-C5-C6-O6
16	G	606	NAG	C3-C2-N2-C7
16	G	602	NAG	C1-C2-N2-C7

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
16	G	601	NAG	1	0
16	B	701	NAG	1	0
16	G	606	NAG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
16	G	602	NAG	1	0
16	G	604	NAG	2	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

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	B	133/153 (86%)	0.68	16 (12%) 9 6	46, 76, 144, 167	0
2	D	240/243 (98%)	0.24	1 (0%) 88 70	46, 98, 243, 277	0
3	E	213/216 (98%)	0.28	3 (1%) 73 46	45, 113, 181, 199	0
4	G	456/482 (94%)	0.28	13 (2%) 53 30	44, 105, 174, 218	0
5	H	231/236 (97%)	0.77	14 (6%) 27 16	64, 161, 217, 255	0
6	L	211/214 (98%)	0.63	12 (5%) 29 17	62, 156, 210, 224	0
All	All	1484/1544 (96%)	0.44	59 (3%) 42 23	44, 119, 206, 277	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	536	ALA	7.1
1	B	517	ALA	5.6
1	B	652	GLN	4.2
5	H	99	GLN	4.2
5	H	81	GLN	4.0
4	G	176	PHE	3.5
5	H	47	TRP	3.5
5	H	64	ASN	3.5
1	B	537	LEU	3.5
4	G	506	VAL	3.4
6	L	108	SER	3.4
6	L	80	VAL	3.3
1	B	535	MET	3.3
3	E	115	VAL	3.2
1	B	542	ARG	3.2
1	B	516	GLY	3.1
5	H	123	ALA	3.1
4	G	359	ILE	3.0
1	B	539	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
6	L	47	LEU	2.9
5	H	82(A)	ARG	2.9
4	G	68	VAL	2.9
6	L	158	ALA	2.8
1	B	649	SER	2.8
3	E	202	VAL	2.8
5	H	145	PRO	2.8
5	H	144	PHE	2.7
5	H	100(L)	TYR	2.7
5	H	71	ARG	2.6
4	G	58	ALA	2.6
6	L	41	GLY	2.6
1	B	541	ALA	2.5
1	B	518	PHE	2.5
4	G	134	VAL	2.5
1	B	521	GLY	2.5
1	B	512	ALA	2.5
1	B	545	LEU	2.3
4	G	326	ILE	2.3
5	H	94	THR	2.3
6	L	36	TYR	2.3
5	H	69	ILE	2.2
4	G	379	GLY	2.2
4	G	427	TRP	2.2
4	G	465	THR	2.2
5	H	56	THR	2.2
1	B	648	VAL	2.2
6	L	45	ILE	2.2
4	G	70	ALA	2.2
5	H	68	VAL	2.2
1	B	533	ALA	2.1
6	L	203	VAL	2.1
6	L	106	VAL	2.1
6	L	26	GLN	2.1
2	D	148	GLU	2.1
6	L	190	LYS	2.0
3	E	197	HIS	2.0
6	L	178	TYR	2.0
4	G	367	GLY	2.0
4	G	125	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	NAG	A	1	14/15	-	-	179,179,179,179	0
7	NAG	A	2	14/15	-	-	179,179,179,179	0
7	BMA	A	3	11/12	-	-	179,179,179,179	0
7	MAN	A	4	11/12	-	-	138,138,138,138	0
7	MAN	A	5	11/12	-	-	90,90,90,90	0
7	MAN	A	6	11/12	-	-	138,138,138,138	0
8	NAG	C	1	14/15	-	-	138,138,138,138	0
8	NAG	C	2	14/15	-	-	138,138,138,138	0
15	BMA	T	3	11/12	0.21	0.14	179,179,179,179	0
14	MAN	S	5	11/12	0.27	0.12	179,179,179,179	0
11	BMA	N	3	11/12	0.32	0.11	179,179,179,179	0
8	NAG	Q	1	14/15	0.34	0.14	179,179,179,179	0
9	NAG	F	1	14/15	-	-	179,179,179,179	0
9	NAG	F	2	14/15	-	-	179,179,179,179	0
9	BMA	F	3	11/12	-	-	179,179,179,179	0
9	MAN	F	4	11/12	-	-	179,179,179,179	0
9	MAN	F	5	11/12	-	-	179,179,179,179	0
9	MAN	F	6	11/12	-	-	179,179,179,179	0
9	MAN	F	7	11/12	-	-	179,179,179,179	0
8	NAG	Q	2	14/15	0.35	0.18	179,179,179,179	0
11	BMA	O	3	11/12	0.36	0.12	138,138,138,138	0
10	BMA	J	3	11/12	-	-	138,138,138,138	0
10	MAN	J	4	11/12	-	-	138,138,138,138	0
11	BMA	K	3	11/12	0.37	0.14	41,41,41,41	0
11	NAG	U	2	14/15	0.38	0.12	179,179,179,179	0
13	MAN	R	4	11/12	0.44	0.12	138,138,138,138	0
14	MAN	S	4	11/12	0.48	0.10	179,179,179,179	0
14	BMA	S	3	11/12	0.49	0.10	179,179,179,179	0
10	BMA	M	3	11/12	0.50	0.10	179,179,179,179	0
11	BMA	U	3	11/12	0.52	0.10	179,179,179,179	0
12	MAN	P	6	11/12	0.54	0.12	138,138,138,138	0
12	MAN	P	5	11/12	0.54	0.15	138,138,138,138	0

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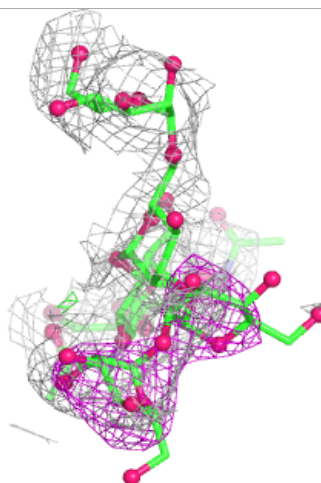
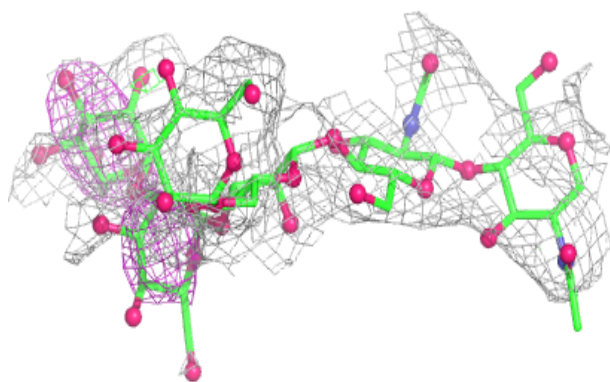
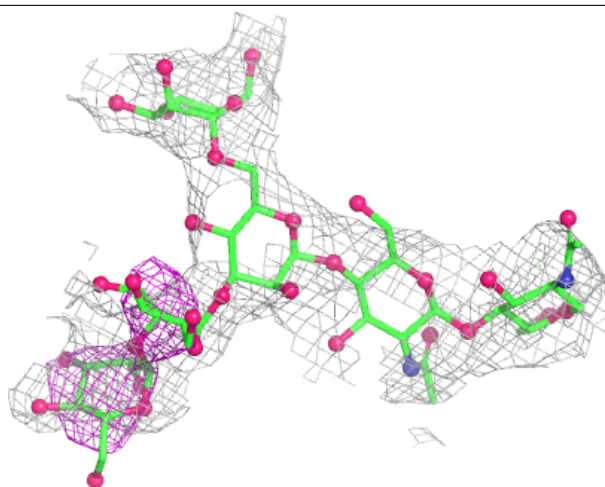
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
12	MAN	P	7	11/12	0.55	0.11	138,138,138,138	0
11	NAG	O	2	14/15	0.59	0.14	138,138,138,138	0
10	MAN	M	4	11/12	0.59	0.10	138,138,138,138	0
13	BMA	R	3	11/12	0.60	0.10	138,138,138,138	0
13	NAG	R	2	14/15	0.63	0.10	179,179,179,179	0
11	NAG	N	2	14/15	0.64	0.10	179,179,179,179	0
12	MAN	P	8	11/12	0.66	0.25	138,138,138,138	0
11	NAG	U	1	14/15	0.66	0.20	179,179,179,179	0
11	NAG	K	2	14/15	0.66	0.13	41,41,41,41	0
8	NAG	I	2	14/15	0.66	0.47	179,179,179,179	0
15	NAG	T	2	14/15	0.70	0.12	179,179,179,179	0
12	BMA	P	3	11/12	0.72	0.12	138,138,138,138	0
10	NAG	J	2	14/15	0.74	0.11	138,138,138,138	0
15	NAG	T	1	14/15	0.75	0.10	179,179,179,179	0
12	NAG	P	2	14/15	0.76	0.18	179,179,179,179	0
10	NAG	M	2	14/15	0.76	0.15	179,179,179,179	0
13	NAG	R	1	14/15	0.76	0.16	179,179,179,179	0
11	NAG	K	1	14/15	0.77	0.17	179,179,179,179	0
12	MAN	P	4	11/12	0.77	0.13	138,138,138,138	0
11	NAG	N	1	14/15	0.79	0.11	179,179,179,179	0
12	NAG	P	1	14/15	0.80	0.19	179,179,179,179	0
14	NAG	S	2	14/15	0.81	0.13	179,179,179,179	0
10	NAG	M	1	14/15	0.85	0.21	179,179,179,179	0
14	NAG	S	1	14/15	0.86	0.14	179,179,179,179	0
11	NAG	O	1	14/15	0.87	0.20	179,179,179,179	0
8	NAG	I	1	14/15	0.88	0.39	179,179,179,179	0
10	NAG	J	1	14/15	0.92	0.12	179,179,179,179	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.

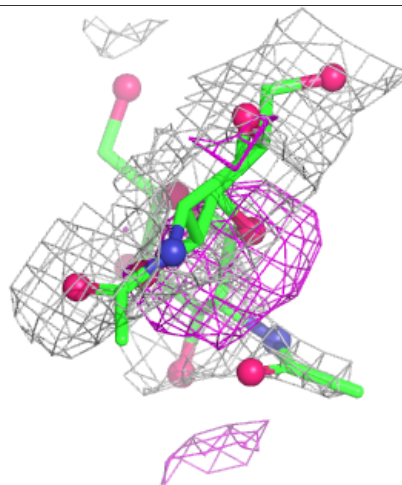
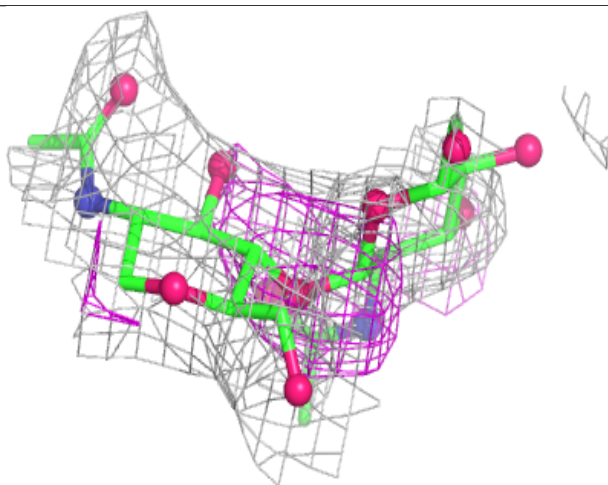
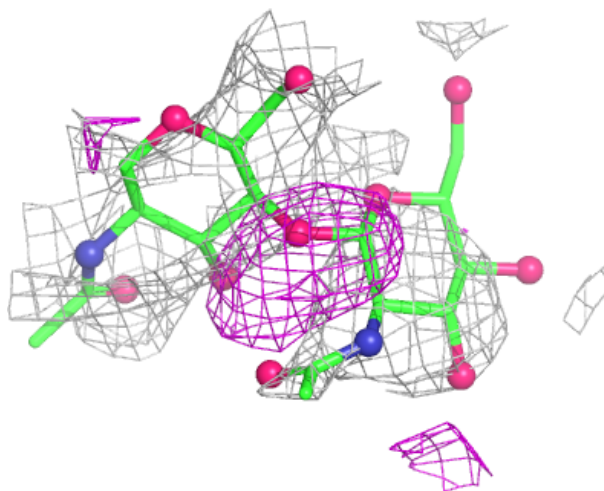
Electron density around Chain A:

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



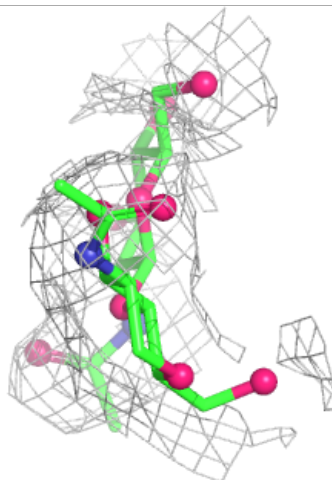
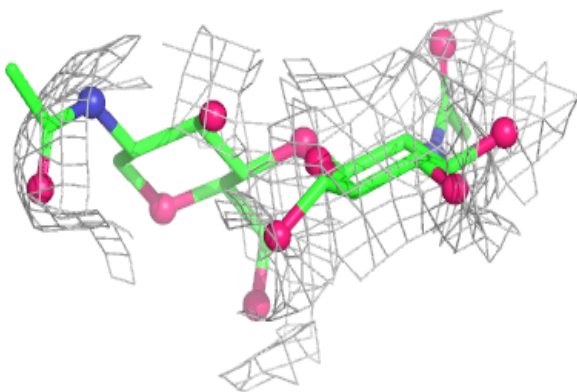
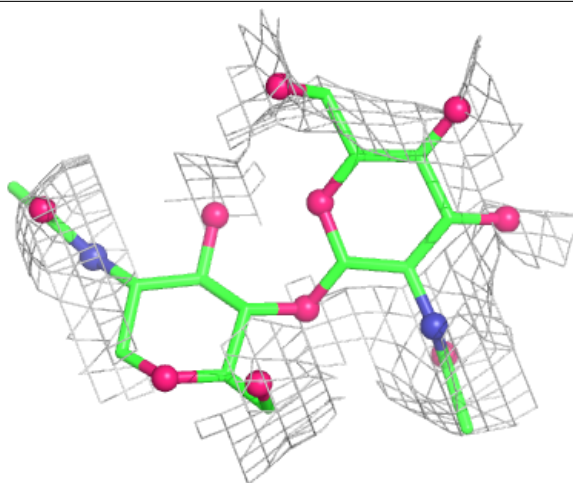
Electron density around Chain C:

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



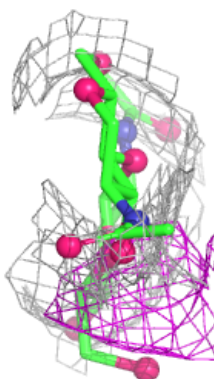
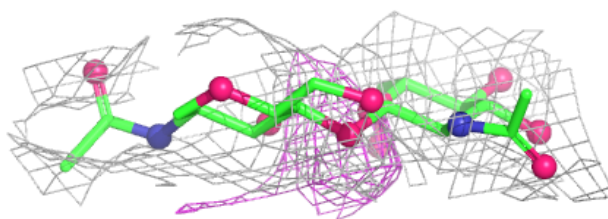
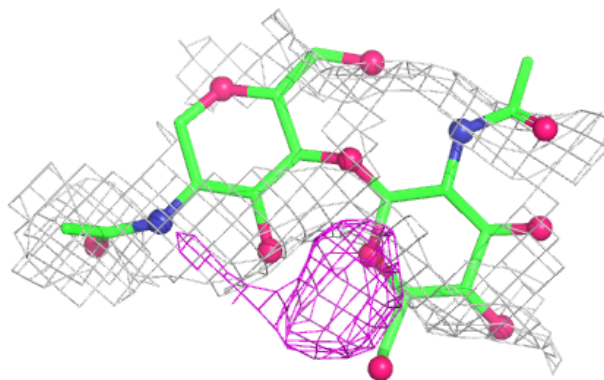
Electron density around Chain I:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



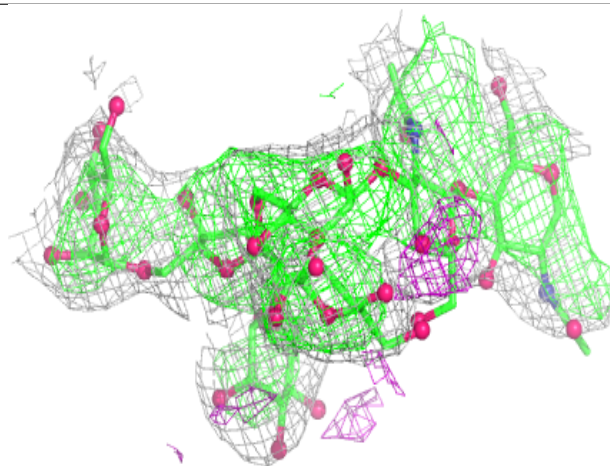
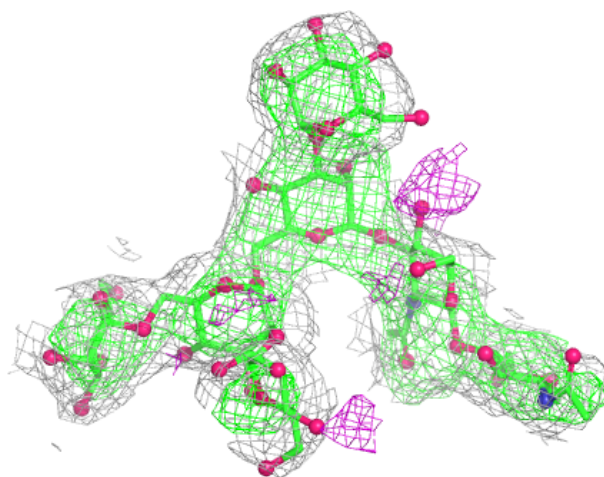
Electron density around Chain Q:

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



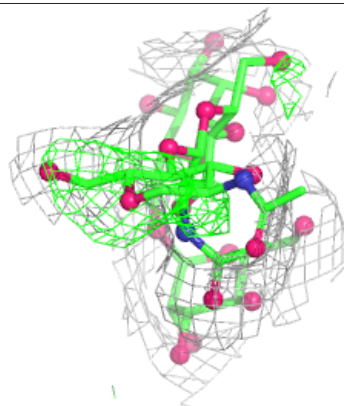
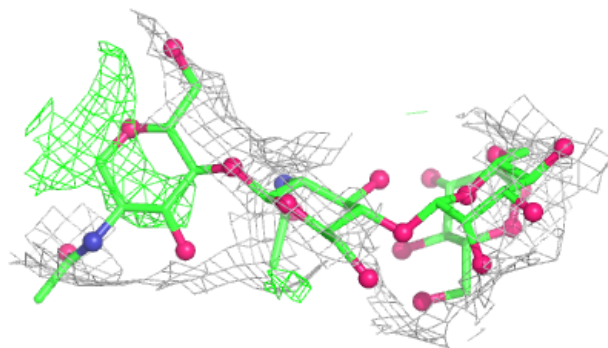
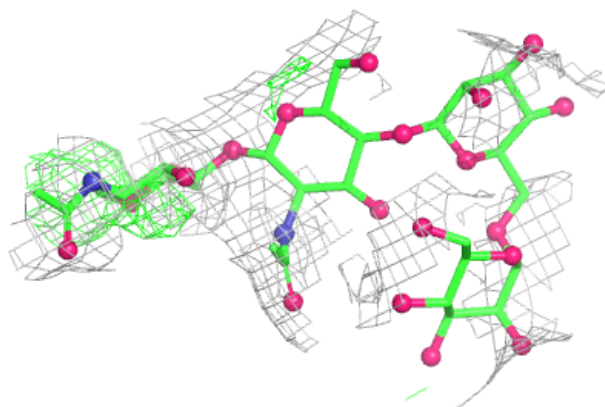
Electron density around Chain F:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

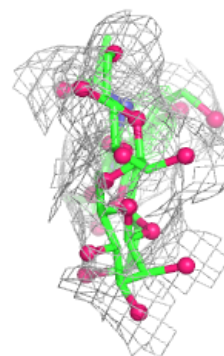
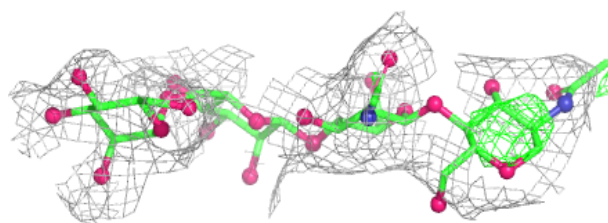
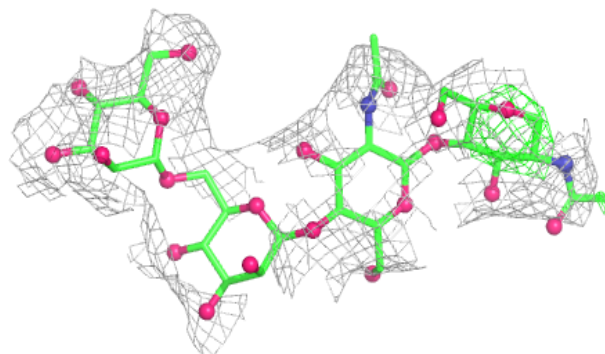


Electron density around Chain J:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

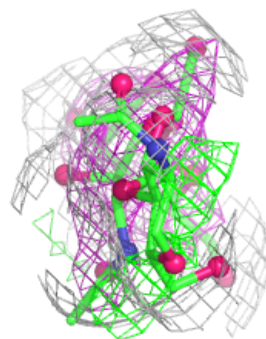
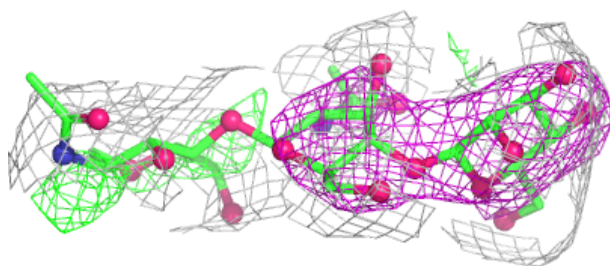
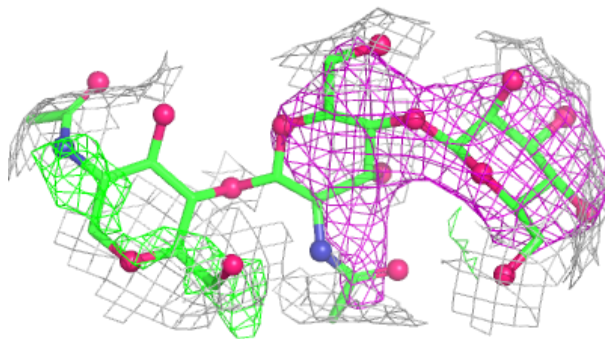
**Electron density around Chain M:**

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

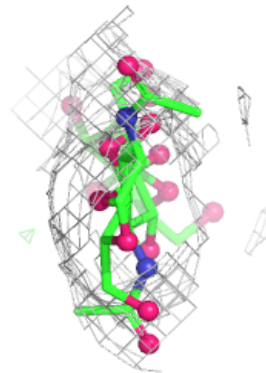
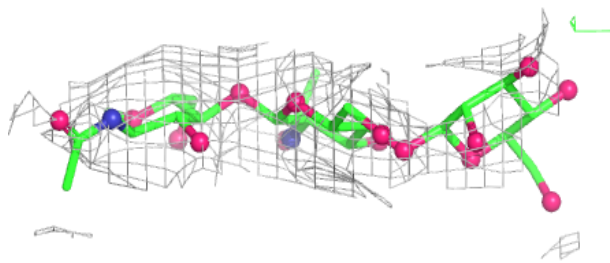
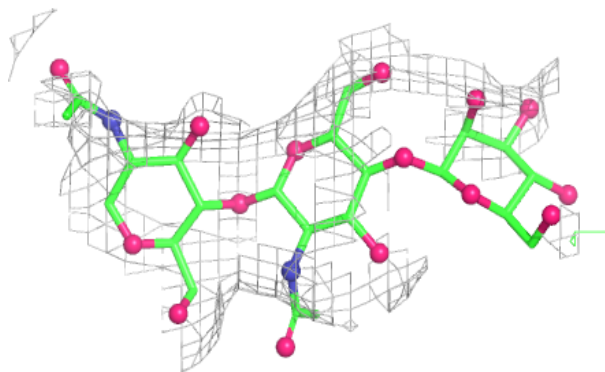


Electron density around Chain K:

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

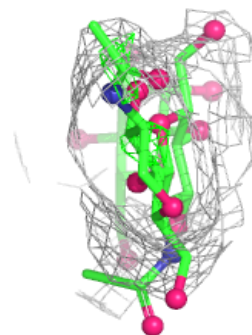
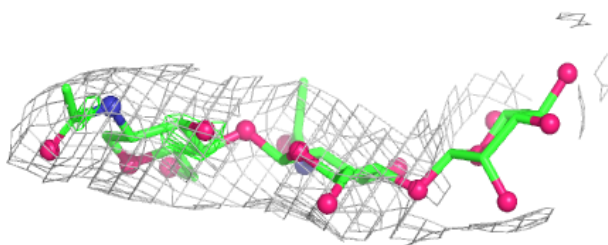
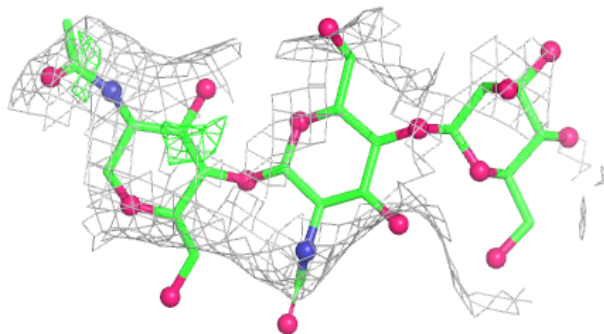
**Electron density around Chain N:**

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

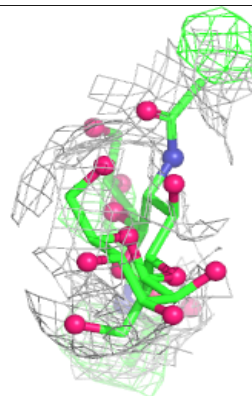
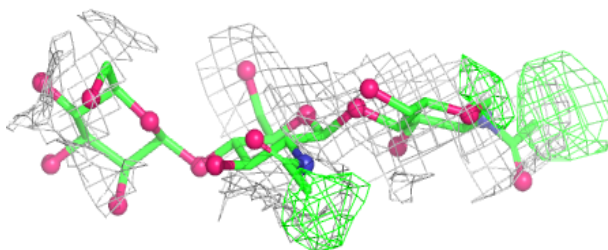
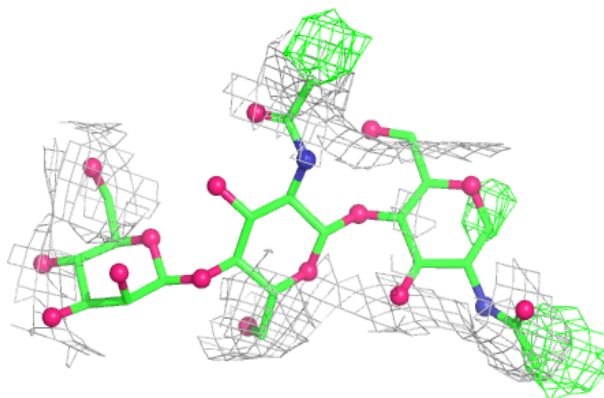


Electron density around Chain O:

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and green (positive)

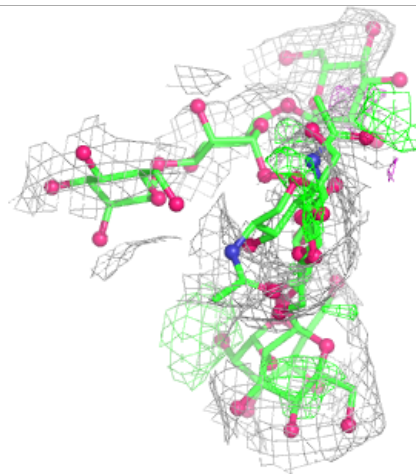
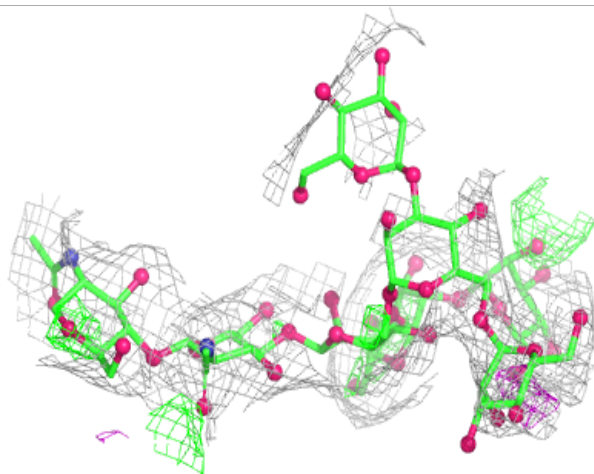
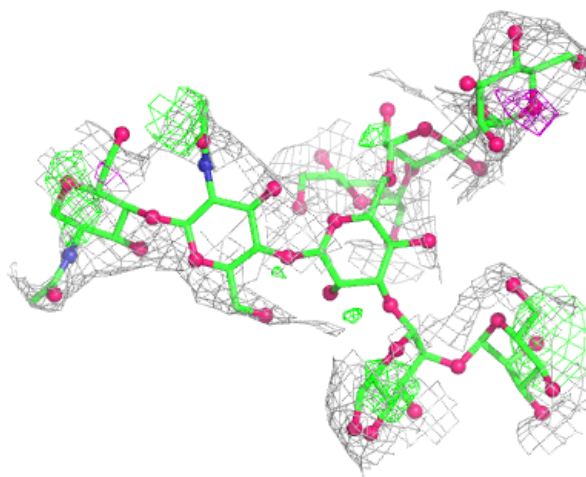
**Electron density around Chain U:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



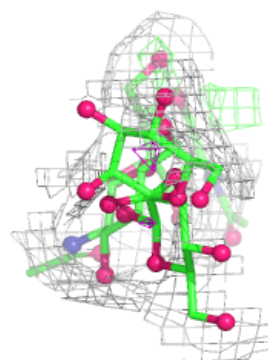
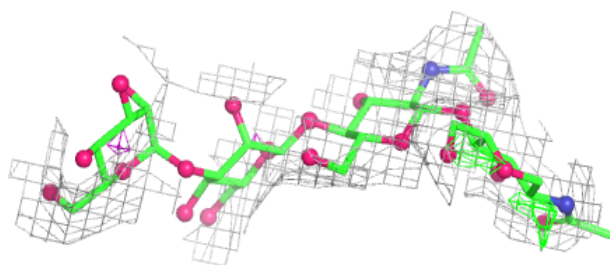
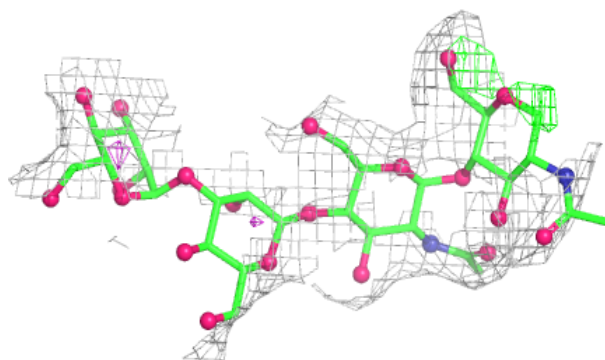
Electron density around Chain P:

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and green (positive)



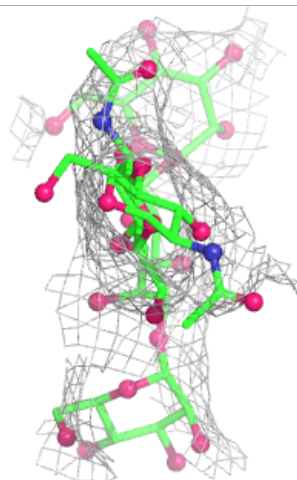
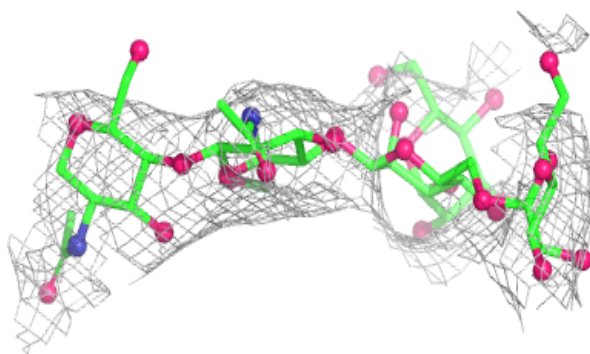
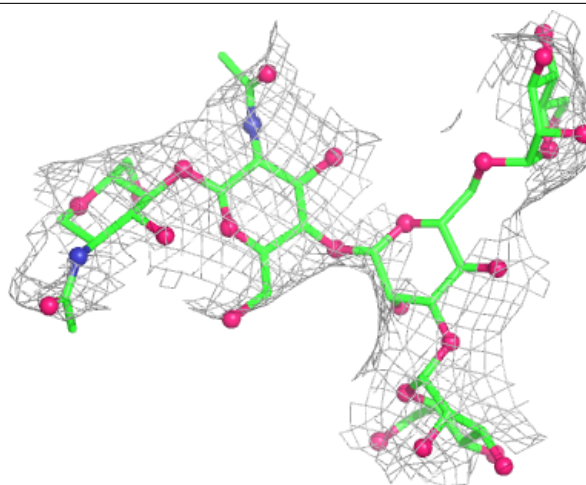
Electron density around Chain R:

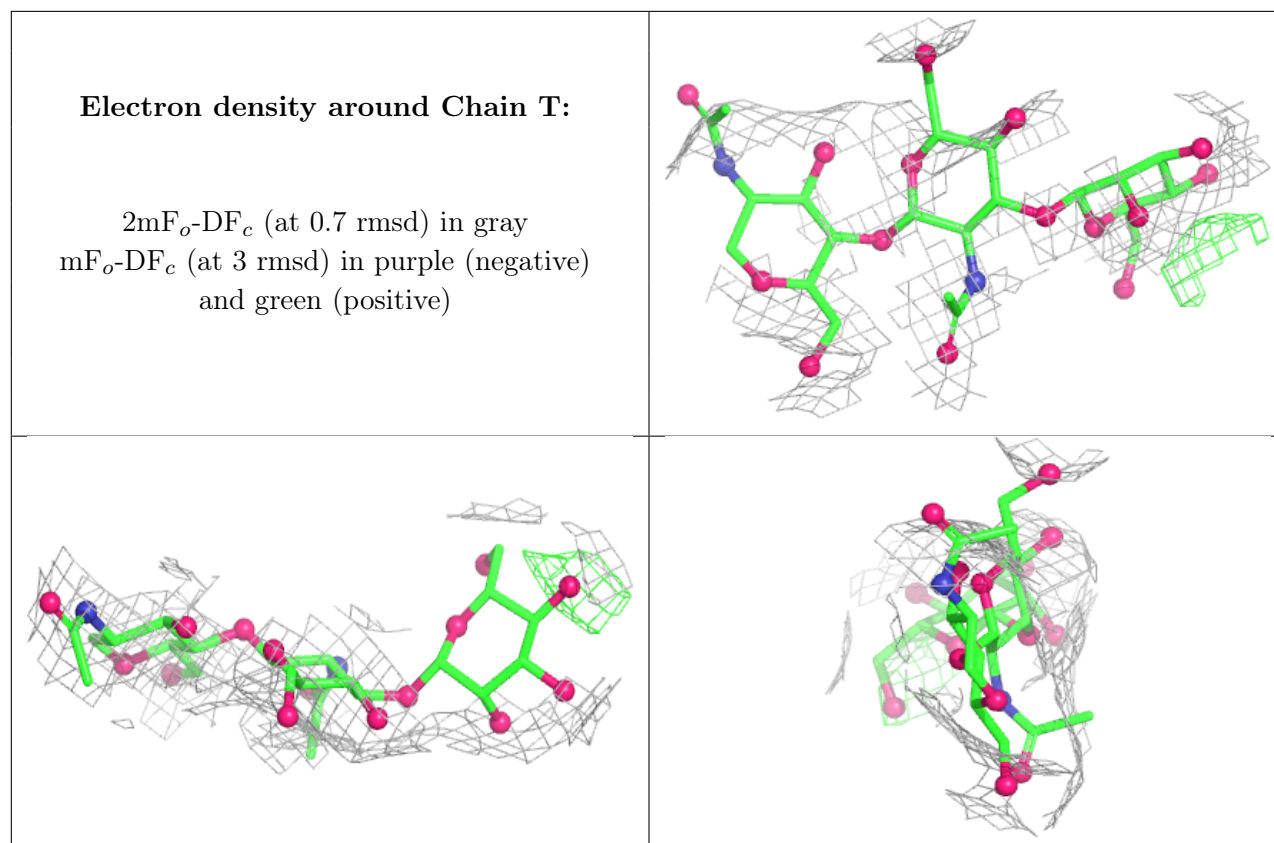
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain S:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
16	NAG	B	701	14/15	0.34	0.16	179,179,179,179	0
16	NAG	G	606	14/15	0.39	0.10	179,179,179,179	0
16	NAG	G	605	14/15	0.48	0.09	179,179,179,179	0
16	NAG	B	702	14/15	0.51	0.13	179,179,179,179	0
16	NAG	G	603	14/15	0.51	0.13	179,179,179,179	0
16	NAG	G	604	14/15	0.68	0.19	179,179,179,179	0
16	NAG	G	602	14/15	0.70	0.12	179,179,179,179	0
16	NAG	G	601	14/15	0.72	0.12	179,179,179,179	0

6.5 Other polymers [i](#)

There are no such residues in this entry.