



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:06 PM UTC

PDB ID : 9S80 / pdb_00009s80
Title : Crystal structure of the BRL3 ectodomain from Arabidopsis thaliana in complex with brassinolide.
Authors : Hohmann, U.; Hothorn, M.
Deposited on : 2025-08-05
Resolution : 2.45 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

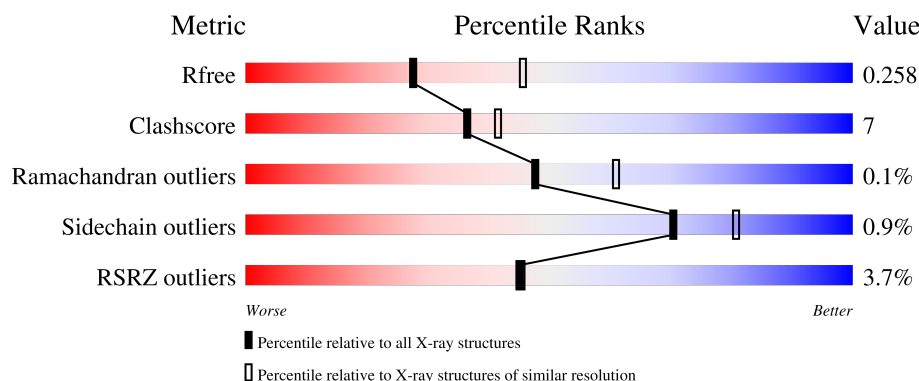
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



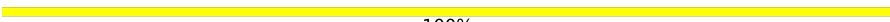
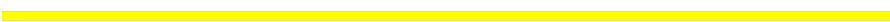
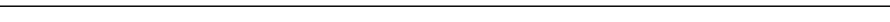
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	770	4% 80% 13% 6%
1	B	770	3% 82% 12% 6%
2	C	2	50% 50%
2	D	2	50% 50%
2	E	2	50% 50%

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Mol	Chain	Length	Quality of chain
2	H	2	 100%
2	J	2	 100%
2	K	2	 50% 50%
2	L	2	 100%
2	M	2	 100%
2	P	2	 50% 50%
2	R	2	 100%
2	S	2	 100%
2	T	2	 100%
3	F	11	 9% 73% 18%
3	N	11	 18% 82%
4	G	3	 33% 67%
4	O	3	 33% 67%
5	I	3	 33% 67%
5	Q	3	 33% 67%

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
12	ACT	B	808	-	-	X	-

2 Entry composition [i](#)

There are 13 unique types of molecules in this entry. The entry contains 23739 atoms, of which 11699 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 Receptor-like protein kinase BRI1-like 3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	721	Total	C	H	N	O	S	0	4	0
			10775	3396	5367	909	1075	28			
1	B	727	Total	C	H	N	O	S	0	7	0
			10908	3436	5436	924	1083	29			

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



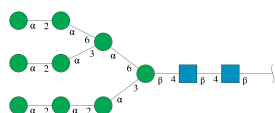
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	D	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	E	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	H	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	J	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	K	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	L	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	M	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	P	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	R	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			

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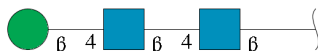
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	S	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	T	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			

- Molecule 3 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-(1-2)-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.



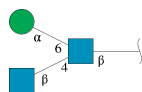
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	11	Total	C	H	N	O	0	0	0
			233	70	106	2	55			
3	N	11	Total	C	H	N	O	0	0	0
			233	70	106	2	55			

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



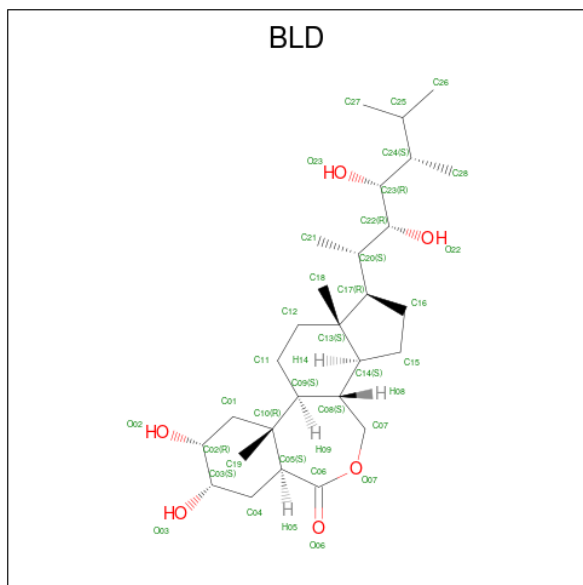
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	3	Total	C	H	N	O	0	0	0
			73	22	34	2	15			
4	O	3	Total	C	H	N	O	0	0	0
			73	22	34	2	15			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-mannopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	I	3	Total	C	H	N	O	0	0	0
			73	22	34	2	15			
5	Q	3	Total	C	H	N	O	0	0	0
			73	22	34	2	15			

- Molecule 6 is Brassinolide (CCD ID: BLD) (formula: $C_{28}H_{48}O_6$).



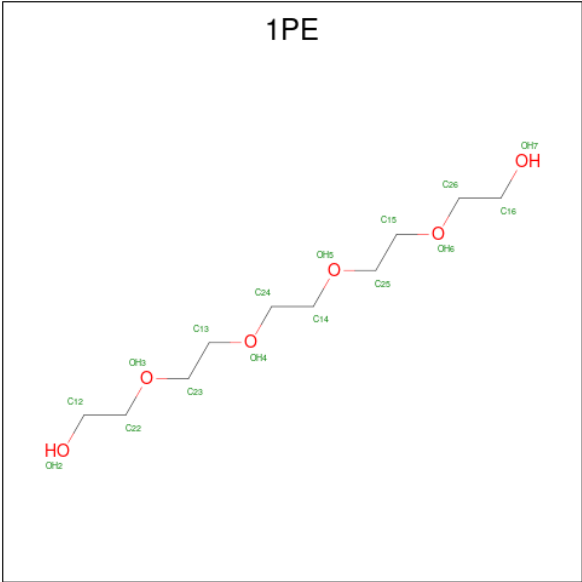
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	H	O		0	0
			82	28	48	6			
6	B	1	Total	C	H	O		0	0
			82	28	48	6			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



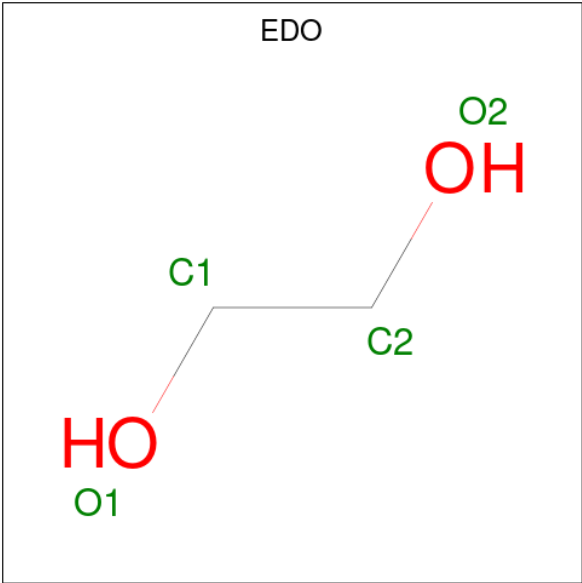
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
7	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
7	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
7	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
7	B	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
7	B	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
7	B	1	Total	C	H	N	O	0	0
			27	8	13	1	5		

- Molecule 8 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



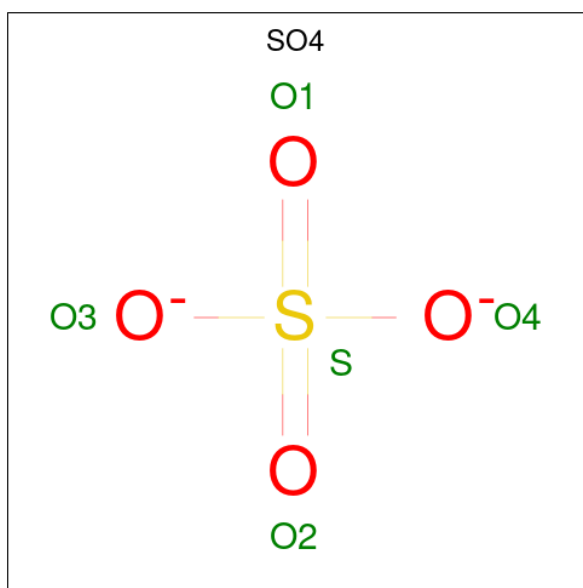
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	H	O	0	0
			38	10	22	6		
8	B	1	Total	C	H	O	0	0
			38	10	22	6		

- Molecule 9 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



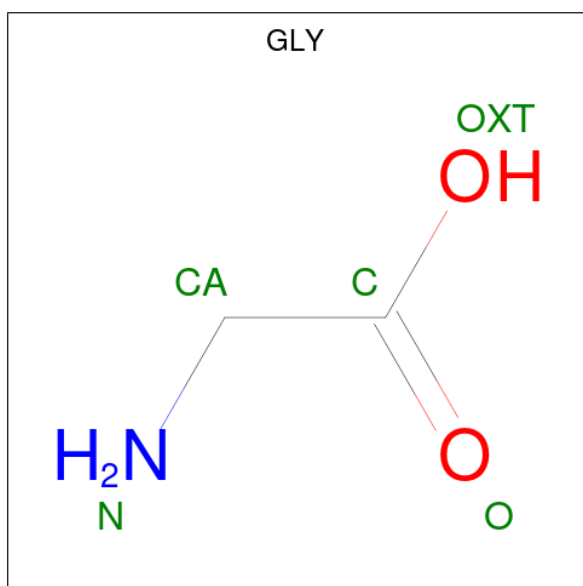
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	H	O	0	0
			10	2	6	2		
9	B	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 10 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



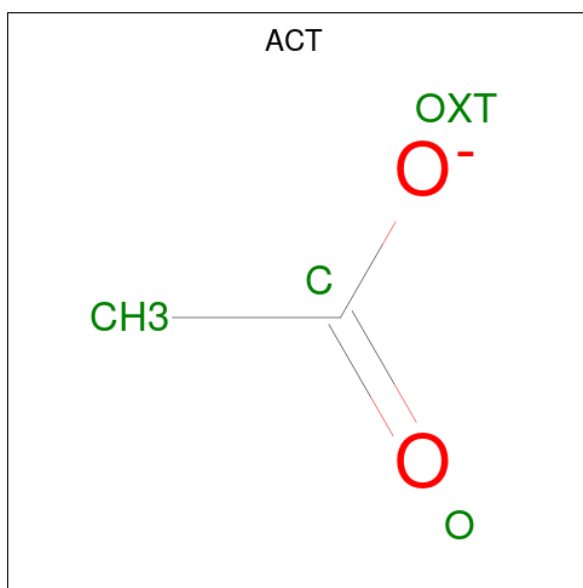
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	O	S	0	0
			5	4	1		
10	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 11 is GLYCINE (CCD ID: GLY) (formula: C₂H₅NO₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	B	1	Total	C	H	N	O	0	0
			6	2	2	1	1		

- Molecule 12 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	B	1	Total	C	H	O	0	0
			7	2	3	2		

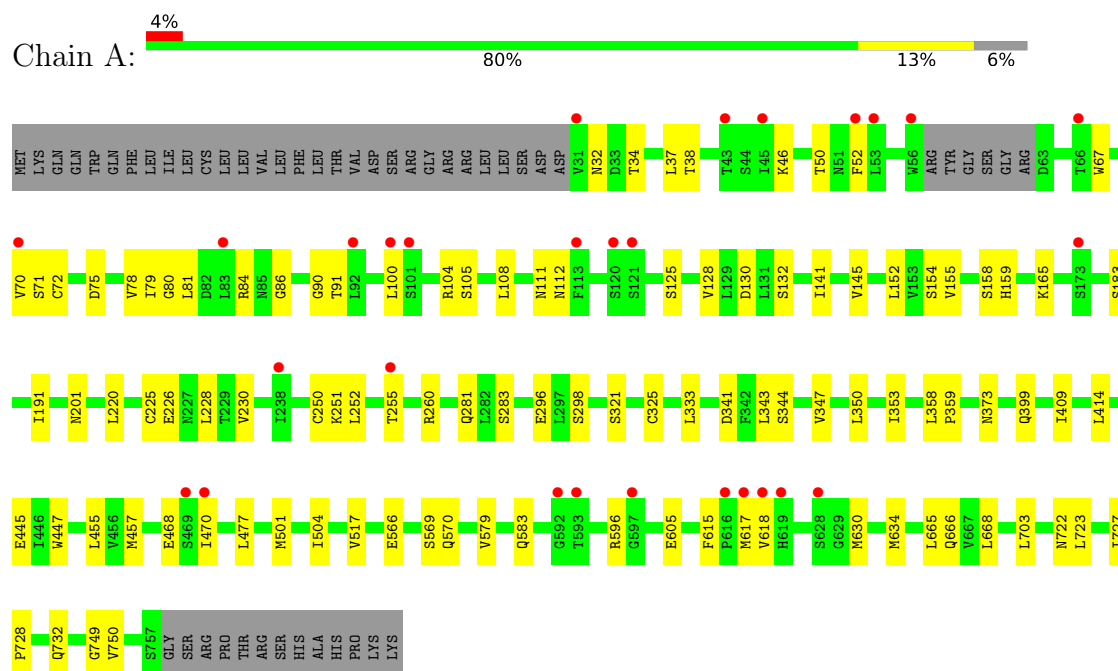
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	94	Total	O	0	0
			94	94		
13	B	96	Total	O	0	0
			96	96		

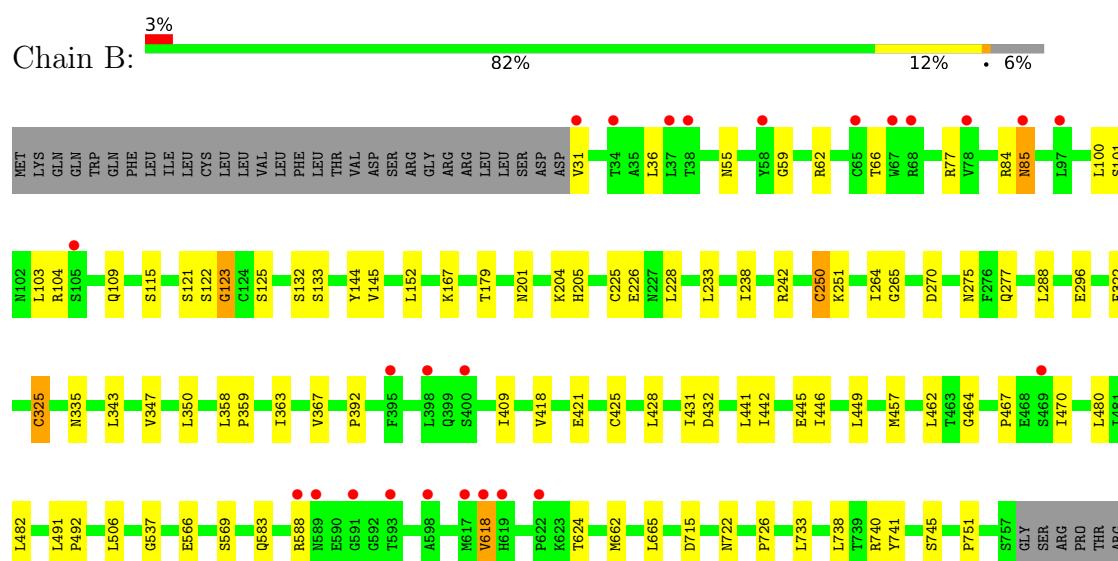
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Receptor-like protein kinase BRI1-like 3



• Molecule 1: Receptor-like protein kinase BRI1-like 3



SER
HIS
ALA
HIS
PRO
LYS
LYS

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

Chain C:  50% 50%

NAG1
NAG2

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

Chain D:  50% 50%

NAG1
NAG2

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

Chain E:  50% 50%

NAG1
NAG2

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

Chain H:  100%

NAG1
NAG2

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

Chain J:  100%

NAG1
NAG2

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

Chain K:  50% 50%

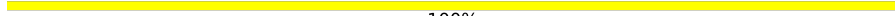
NAG1
NAG2

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

Chain L:  100%

MAG1
MAG2

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

Chain M:  100%

MAG1
MAG2

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

Chain P:  50% 50%

MAG1
MAG2

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

Chain R:  100%

MAG1
MAG2

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

Chain S:  100%

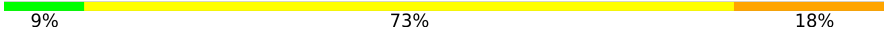
MAG1
MAG2

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

Chain T:  100%

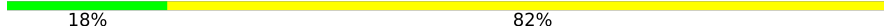
MAG1
MAG2

- Molecule 3: 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-(1-2)-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 F:  9% 73% 18%

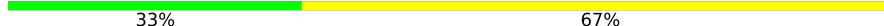


- Molecule 3: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-6)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy-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

Chain N:  18% 82%



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

Chain G:  33% 67%



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

Chain O:  33% 67%



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

Chain I:  33% 67%



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

Chain Q:  33% 67%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.46Å 81.76Å 122.19Å 107.24° 91.94° 112.39°	Depositor
Resolution (Å)	48.66 – 2.45 48.66 – 2.45	Depositor EDS
% Data completeness (in resolution range)	98.0 (48.66-2.45) 98.0 (48.66-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.222 , 0.258 0.222 , 0.258	Depositor DCC
R_{free} test set	3668 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	50.8	Xtriage
Anisotropy	0.336	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 38.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.008 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	23739	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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: EDO, ACT, NAG, SO4, 1PE, BMA, MAN, BLD

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	0.16	0/5516	0.35	0/7495
1	B	0.16	0/5588	0.35	0/7592
All	All	0.16	0/11104	0.35	0/15087

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5408	5367	5355	84	0
1	B	5472	5436	5432	79	0
2	C	28	25	25	1	0
2	D	28	25	25	1	0
2	E	28	25	25	0	0
2	H	28	25	25	2	0
2	J	28	25	25	0	0
2	K	28	25	25	1	0
2	L	28	25	25	0	0
2	M	28	25	25	0	0
2	P	28	25	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	R	28	25	25	1	0
2	S	28	25	25	0	0
2	T	28	25	25	0	0
3	F	127	106	106	2	0
3	N	127	106	106	0	0
4	G	39	34	34	0	0
4	O	39	34	34	0	0
5	I	39	34	34	1	0
5	Q	39	34	34	0	0
6	A	34	48	48	0	0
6	B	34	48	48	1	0
7	A	56	52	52	2	0
7	B	42	39	39	3	0
8	A	16	22	22	1	0
8	B	16	22	22	0	0
9	A	4	6	6	0	0
9	B	4	6	6	0	0
10	A	5	0	0	0	0
10	B	5	0	0	0	0
11	B	4	2	2	0	0
12	B	4	3	3	2	0
13	A	94	0	0	8	0
13	B	96	0	0	8	0
All	All	12040	11699	11683	170	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 7.

All (170) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:ASN:OD1	1:A:34:THR:OG1	1.81	0.97
1:B:343:LEU:HD11	1:B:358:LEU:HD13	1.60	0.82
1:A:296:GLU:N	1:A:296:GLU:OE1	2.13	0.81
1:B:77:ARG:NH1	1:B:101:SER:OG	2.13	0.81
1:B:715:ASP:OD1	1:B:740:ARG:NH2	2.14	0.81
1:B:335[A]:ASN:OD1	13:B:901:HOH:O	2.01	0.79
1:A:605:GLU:OE1	13:A:901:HOH:O	1.99	0.79
1:B:467:PRO:O	1:B:470:ILE:HD11	1.84	0.77
1:A:37:LEU:HD21	1:A:100:LEU:HD11	1.66	0.76
1:B:277:GLN:NE2	13:B:904:HOH:O	2.18	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:806:1PE:OH7	13:A:902:HOH:O	2.05	0.75
1:B:343:LEU:HD11	1:B:358:LEU:CD1	2.17	0.75
1:B:418:VAL:HG21	1:B:457:MET:HE1	1.70	0.74
1:B:421:GLU:N	1:B:421:GLU:OE1	2.20	0.73
1:B:491[B]:LEU:HD21	1:B:506:LEU:CD1	2.19	0.71
1:B:432:ASP:OD2	13:B:902:HOH:O	2.10	0.70
1:A:130:ASP:OD1	1:A:132:SER:OG	2.10	0.69
1:B:726:PRO:O	13:B:903:HOH:O	2.12	0.68
1:A:583:GLN:HA	1:A:583:GLN:HE21	1.59	0.68
1:B:233:LEU:HD13	1:B:238:ILE:HD13	1.79	0.65
1:A:566:GLU:O	1:A:569:SER:OG	2.14	0.65
1:A:732:GLN:N	1:A:732:GLN:OE1	2.30	0.65
1:A:50:THR:OG1	1:A:52:PHE:HB3	1.97	0.65
1:A:201:ASN:ND2	1:A:226:GLU:OE1	2.29	0.64
1:B:442:ILE:HD11	1:B:462:LEU:HD13	1.79	0.64
1:A:255[A]:THR:HG22	1:A:281:GLN:HG3	1.80	0.63
1:B:343:LEU:HD21	1:B:363:ILE:HD13	1.79	0.63
1:A:141:ILE:O	1:A:145:VAL:HG23	1.98	0.63
1:B:85:ASN:N	1:B:85:ASN:OD1	2.33	0.61
1:B:491[B]:LEU:HD21	1:B:506:LEU:HD13	1.81	0.61
1:A:341:ASP:OD2	1:A:344:SER:HB3	2.01	0.61
1:B:55:ASN:ND2	1:B:66:THR:O	2.34	0.61
1:B:733:LEU:O	1:B:741:TYR:OH	2.12	0.60
1:B:467:PRO:C	1:B:470:ILE:HD11	2.27	0.60
1:A:70:VAL:HG13	1:A:78:VAL:HG13	1.83	0.59
1:A:722:ASN:HD21	7:A:805:NAG:H83	1.67	0.59
1:A:70:VAL:HG22	1:A:81:LEU:HD13	1.84	0.59
1:A:191:ILE:HB	1:A:220:LEU:HD11	1.84	0.59
1:B:425:CYS:HB3	1:B:428:LEU:HG	1.84	0.58
1:A:298:SER:HB3	1:A:321:SER:OG	2.03	0.58
2:C:2:NAG:H3	2:C:2:NAG:H83	1.85	0.57
1:B:84:ARG:NH1	7:B:803:NAG:O6	2.36	0.57
1:A:132:SER:O	1:A:158:SER:O	2.22	0.57
1:A:165:LYS:HD2	1:A:165:LYS:H	1.69	0.56
2:H:1:NAG:O6	2:H:2:NAG:H82	2.05	0.56
1:B:228:LEU:HB3	1:B:250[A]:CYS:SG	2.46	0.56
1:A:373:ASN:ND2	13:A:911:HOH:O	2.38	0.56
1:B:133:SER:HB2	7:B:803:NAG:H81	1.88	0.56
1:A:732:GLN:N	1:A:732:GLN:CD	2.63	0.56
13:A:993:HOH:O	2:K:1:NAG:O6	2.17	0.56
1:B:347:VAL:HA	1:B:350:LEU:HD12	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:LEU:HB2	1:A:501:MET:HE1	1.88	0.55
1:B:264:ILE:HG22	1:B:265:GLY:N	2.21	0.55
1:A:46:LYS:HG2	1:A:90:GLY:HA2	1.89	0.55
1:A:409:ILE:HG21	1:A:414:LEU:HD11	1.89	0.55
1:A:343:LEU:HD11	1:A:358:LEU:CD1	2.38	0.54
1:A:230:VAL:HG13	1:A:255[A]:THR:OG1	2.08	0.54
1:A:152:LEU:HD21	1:A:155:VAL:HB	1.90	0.54
1:B:588:ARG:NH2	1:B:715:ASP:OD2	2.41	0.54
1:B:343:LEU:CD2	1:B:363:ILE:HD13	2.38	0.54
1:B:201:ASN:ND2	1:B:226:GLU:OE1	2.33	0.54
1:B:428:LEU:HD13	1:B:431:ILE:HD11	1.90	0.53
1:B:467:PRO:O	1:B:492:PRO:HG2	2.09	0.53
1:B:233:LEU:HD13	1:B:238:ILE:CD1	2.39	0.53
1:B:537:GLY:N	13:B:911:HOH:O	2.29	0.53
1:B:264:ILE:HG12	1:B:288:LEU:HD12	1.91	0.53
1:B:445:GLU:O	1:B:449:LEU:HD22	2.09	0.53
1:A:501:MET:HG2	1:A:504:ILE:HD11	1.92	0.52
1:A:583:GLN:HA	1:A:583:GLN:NE2	2.24	0.52
1:B:583:GLN:NE2	13:B:916:HOH:O	2.40	0.52
1:A:50:THR:HG21	1:A:86:GLY:O	2.10	0.52
1:A:666:GLN:HG2	13:A:936:HOH:O	2.09	0.52
1:A:358:LEU:N	1:A:359:PRO:CD	2.73	0.52
1:A:341:ASP:O	1:A:341:ASP:CG	2.52	0.51
1:B:740:ARG:NH1	12:B:808:ACT:H2	2.25	0.51
1:A:46:LYS:HE2	1:A:91:THR:HG22	1.91	0.51
1:B:740:ARG:HH11	12:B:808:ACT:H2	1.75	0.51
1:B:179:THR:HG22	1:B:205[B]:HIS:CD2	2.46	0.51
1:A:70:VAL:CG2	1:A:81:LEU:HD13	2.41	0.51
1:B:167:LYS:O	13:B:905:HOH:O	2.19	0.51
1:B:121:SER:HA	1:B:144:TYR:HE1	1.77	0.50
1:A:333:LEU:HB2	1:A:358:LEU:HD23	1.94	0.50
1:B:358:LEU:N	1:B:359:PRO:CD	2.75	0.50
1:A:501:MET:HG2	1:A:504:ILE:CD1	2.42	0.50
5:I:2:NAG:O7	5:I:3:MAN:O6	2.19	0.49
1:B:59:GLY:N	1:B:62:ARG:HG2	2.27	0.49
1:B:745:SER:O	1:B:745:SER:OG	2.25	0.49
1:A:617:MET:C	1:A:617:MET:SD	2.95	0.48
1:A:165:LYS:HD2	1:A:165:LYS:N	2.28	0.48
1:A:665:LEU:HD21	1:A:668:LEU:HB2	1.94	0.48
1:A:634:MET:HE2	13:A:951:HOH:O	2.12	0.48
1:B:59:GLY:O	1:B:62:ARG:HG3	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:617:MET:SD	1:A:618:VAL:N	2.86	0.48
1:B:133:SER:CB	7:B:803:NAG:H81	2.43	0.48
1:B:145:VAL:CG1	1:B:152:LEU:HD11	2.43	0.48
1:B:31:VAL:O	1:B:31:VAL:HG13	2.13	0.48
1:B:618:VAL:O	1:B:624:THR:HG21	2.14	0.48
1:B:264:ILE:HG22	1:B:265:GLY:H	1.79	0.48
1:A:75:ASP:OD2	1:A:75:ASP:N	2.46	0.48
1:A:347:VAL:HA	1:A:350:LEU:CD1	2.44	0.47
1:A:615:PHE:CD2	1:A:617:MET:HG3	2.50	0.47
1:B:367:VAL:HG11	1:B:392:PRO:HG3	1.97	0.47
1:A:108:LEU:O	1:A:111:ASN:OD1	2.33	0.47
1:B:251:LYS:NZ	1:B:275:ASN:OD1	2.46	0.47
1:A:255[A]:THR:HG22	1:A:281:GLN:CG	2.45	0.47
1:B:100:LEU:HD12	1:B:103:LEU:HD22	1.96	0.47
1:B:145:VAL:HG13	1:B:152:LEU:HD11	1.96	0.46
2:R:1:NAG:O3	2:R:2:NAG:H83	2.15	0.46
1:B:322:PHE:O	1:B:325[A]:CYS:SG	2.74	0.46
1:A:251:LYS:C	1:A:252:LEU:HD23	2.41	0.46
1:A:445:GLU:N	1:A:445:GLU:OE1	2.46	0.45
1:B:109:GLN:HB2	1:B:132:SER:OG	2.16	0.45
1:A:298:SER:CB	1:A:321:SER:OG	2.65	0.45
1:A:570:GLN:HB2	13:A:959:HOH:O	2.16	0.45
1:A:80:GLY:HA2	1:A:105:SER:O	2.16	0.45
1:B:446:ILE:HA	1:B:449:LEU:CD2	2.47	0.45
1:A:37:LEU:CD2	1:A:100:LEU:HD11	2.40	0.44
1:A:260:ARG:HD2	3:F:2:NAG:H81	1.99	0.44
1:B:358:LEU:N	1:B:359:PRO:HD3	2.32	0.44
1:A:225[B]:CYS:SG	1:A:228:LEU:HD23	2.58	0.44
1:A:517:VAL:O	13:A:903:HOH:O	2.21	0.44
1:B:347:VAL:HA	1:B:350:LEU:CD1	2.47	0.44
1:A:343:LEU:HD11	1:A:358:LEU:HD13	2.00	0.44
1:A:455:LEU:HD11	1:A:457:MET:HE2	1.98	0.44
1:B:104:ARG:NH1	1:B:125:SER:HB2	2.33	0.44
1:B:491[B]:LEU:HD21	1:B:506:LEU:HD12	1.99	0.43
1:B:204:LYS:HA	1:B:228:LEU:HA	2.00	0.43
1:A:112:ASN:HD22	2:D:1:NAG:H83	1.83	0.43
1:A:749:GLY:O	1:A:750:VAL:C	2.61	0.43
1:B:264:ILE:CD1	1:B:288:LEU:HD12	2.48	0.43
1:B:122:SER:O	1:B:123:GLY:C	2.61	0.43
1:B:233:LEU:HD22	1:B:238:ILE:CD1	2.48	0.43
1:B:242:ARG:HH11	1:B:242:ARG:CG	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:SER:HB2	3:F:1:NAG:H82	2.00	0.43
1:B:270:ASP:O	1:B:296:GLU:HB3	2.19	0.43
1:B:583:GLN:HG3	13:B:983:HOH:O	2.19	0.43
1:A:596:ARG:HB2	2:H:1:NAG:H81	2.01	0.42
1:B:738:LEU:HD12	1:B:751:PRO:HB2	2.00	0.42
1:A:630:MET:HE2	1:A:630:MET:HB3	1.93	0.42
1:B:480:LEU:HD11	1:B:482:LEU:CD1	2.49	0.42
1:A:84:ARG:NH1	7:A:802:NAG:O6	2.49	0.42
1:A:158:SER:HA	1:A:183:SER:O	2.20	0.42
1:A:37:LEU:HD13	1:A:70:VAL:HG21	2.00	0.42
1:A:468:GLU:O	1:A:468:GLU:HG3	2.18	0.42
1:B:566:GLU:OE1	1:B:566:GLU:N	2.53	0.42
1:A:225[A]:CYS:SG	1:A:250:CYS:N	2.92	0.42
1:B:264:ILE:HD11	1:B:288:LEU:HD12	2.01	0.42
1:B:662:MET:HE3	1:B:665:LEU:HD22	2.02	0.42
6:B:802:BLD:H112	6:B:802:BLD:H221	2.00	0.42
1:A:38:THR:HG23	1:A:67:TRP:CH2	2.55	0.42
1:A:447:TRP:CE3	1:A:470:ILE:HG22	2.55	0.42
1:A:566:GLU:C	1:A:569:SER:HG	2.24	0.42
1:A:617:MET:HE2	1:A:617:MET:O	2.19	0.42
1:A:583:GLN:HE21	1:A:583:GLN:CA	2.23	0.41
1:A:71:SER:O	1:A:79:ILE:HG13	2.20	0.41
1:A:722:ASN:O	1:A:722:ASN:OD1	2.38	0.41
1:A:128:VAL:HG13	1:A:154:SER:OG	2.21	0.41
1:B:457:MET:HE2	1:B:462:LEU:HD11	2.01	0.41
1:B:441:LEU:HD22	1:B:464:GLY:HA3	2.03	0.41
1:B:445:GLU:O	1:B:449:LEU:CD2	2.68	0.41
1:B:446:ILE:O	1:B:449:LEU:HD23	2.21	0.41
1:B:566:GLU:O	1:B:569:SER:OG	2.34	0.41
1:A:579:VAL:O	1:A:579:VAL:HG22	2.20	0.40
1:A:703:LEU:HD21	1:A:723:LEU:HD13	2.03	0.40
1:A:727:ILE:O	1:A:728:PRO:C	2.64	0.40
1:A:37:LEU:HD21	1:A:100:LEU:CD1	2.45	0.40
1:A:104:ARG:HG2	1:A:125:SER:O	2.22	0.40
1:A:158:SER:O	1:A:159:HIS:C	2.65	0.40
1:B:31:VAL:HG12	1:B:36:LEU:HG	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	721/770 (94%)	663 (92%)	58 (8%)	0	100	100
1	B	732/770 (95%)	661 (90%)	70 (10%)	1 (0%)	48	61
All	All	1453/1540 (94%)	1324 (91%)	128 (9%)	1 (0%)	48	61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	123	GLY

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	633/673 (94%)	630 (100%)	3 (0%)	81	87
1	B	639/673 (95%)	629 (98%)	10 (2%)	55	69
All	All	1272/1346 (94%)	1259 (99%)	13 (1%)	70	77

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	CYS
1	A	325	CYS
1	A	353	ILE
1	B	85	ASN
1	B	115	SER

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Mol	Chain	Res	Type
1	B	225	CYS
1	B	250[A]	CYS
1	B	250[B]	CYS
1	B	325[A]	CYS
1	B	325[B]	CYS
1	B	409	ILE
1	B	618	VAL
1	B	722	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	ASN
1	A	111	ASN
1	A	317	GLN
1	A	538	ASN
1	A	583	GLN
1	B	278	ASN
1	B	320	GLN
1	B	589	ASN
1	B	614	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 ⓘ

58 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	1,2	14,14,15	0.65	0	17,19,21	0.82	0
2	NAG	C	2	2	14,14,15	0.73	0	17,19,21	1.54	1 (5%)
2	NAG	D	1	1,2	14,14,15	0.71	0	17,19,21	1.16	1 (5%)
2	NAG	D	2	2	14,14,15	0.65	0	17,19,21	0.94	0
2	NAG	E	1	1,2	14,14,15	0.65	0	17,19,21	1.03	0
2	NAG	E	2	2	14,14,15	0.65	0	17,19,21	1.19	1 (5%)
3	NAG	F	1	3,1	14,14,15	0.63	0	17,19,21	1.47	4 (23%)
3	MAN	F	10	3	11,11,12	0.81	0	15,15,17	0.94	1 (6%)
3	MAN	F	11	3	11,11,12	0.75	0	15,15,17	1.11	1 (6%)
3	NAG	F	2	3	14,14,15	0.78	1 (7%)	17,19,21	1.26	2 (11%)
3	BMA	F	3	3	11,11,12	0.74	0	15,15,17	2.35	5 (33%)
3	MAN	F	4	3	11,11,12	0.79	0	15,15,17	0.86	1 (6%)
3	MAN	F	5	3	11,11,12	0.95	1 (9%)	15,15,17	1.04	1 (6%)
3	MAN	F	6	3	11,11,12	0.74	0	15,15,17	0.93	0
3	MAN	F	7	3	11,11,12	0.81	1 (9%)	15,15,17	1.06	1 (6%)
3	MAN	F	8	3	11,11,12	0.95	1 (9%)	15,15,17	0.87	0
3	MAN	F	9	3	11,11,12	0.82	0	15,15,17	0.97	1 (6%)
4	NAG	G	1	4,1	14,14,15	0.65	0	17,19,21	0.94	0
4	NAG	G	2	4	14,14,15	0.70	0	17,19,21	1.04	1 (5%)
4	BMA	G	3	4	11,11,12	0.70	0	15,15,17	2.59	5 (33%)
2	NAG	H	1	1,2	14,14,15	0.74	0	17,19,21	1.08	1 (5%)
2	NAG	H	2	2	14,14,15	0.68	0	17,19,21	1.12	1 (5%)
5	NAG	I	1	1,5	14,14,15	0.78	0	17,19,21	1.05	0
5	NAG	I	2	5	14,14,15	0.81	0	17,19,21	1.01	1 (5%)
5	MAN	I	3	5	11,11,12	0.77	0	15,15,17	0.93	1 (6%)
2	NAG	J	1	1,2	14,14,15	0.63	0	17,19,21	1.63	4 (23%)
2	NAG	J	2	2	14,14,15	0.72	0	17,19,21	1.18	2 (11%)
2	NAG	K	1	1,2	14,14,15	0.64	0	17,19,21	1.36	2 (11%)
2	NAG	K	2	2	14,14,15	0.70	0	17,19,21	1.22	3 (17%)
2	NAG	L	1	1,2	14,14,15	0.65	0	17,19,21	1.21	1 (5%)
2	NAG	L	2	2	14,14,15	0.65	0	17,19,21	1.13	1 (5%)
2	NAG	M	1	1,2	14,14,15	0.68	0	17,19,21	1.07	1 (5%)
2	NAG	M	2	2	14,14,15	0.67	0	17,19,21	1.13	1 (5%)
3	NAG	N	1	3,1	14,14,15	0.68	0	17,19,21	1.36	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MAN	N	10	3	11,11,12	0.73	0	15,15,17	1.38	1 (6%)
3	MAN	N	11	3	11,11,12	0.78	0	15,15,17	0.87	0
3	NAG	N	2	3	14,14,15	0.63	0	17,19,21	1.24	2 (11%)
3	BMA	N	3	3	11,11,12	0.84	0	15,15,17	2.03	4 (26%)
3	MAN	N	4	3	11,11,12	0.81	1 (9%)	15,15,17	1.18	1 (6%)
3	MAN	N	5	3	11,11,12	0.82	0	15,15,17	1.07	1 (6%)
3	MAN	N	6	3	11,11,12	0.69	0	15,15,17	0.98	0
3	MAN	N	7	3	11,11,12	0.87	0	15,15,17	1.06	2 (13%)
3	MAN	N	8	3	11,11,12	0.77	0	15,15,17	1.13	1 (6%)
3	MAN	N	9	3	11,11,12	0.83	1 (9%)	15,15,17	0.84	0
4	NAG	O	1	4,1	14,14,15	0.68	0	17,19,21	1.02	1 (5%)
4	NAG	O	2	4	14,14,15	0.71	0	17,19,21	0.97	0
4	BMA	O	3	4	11,11,12	0.69	0	15,15,17	2.65	6 (40%)
2	NAG	P	1	1,2	14,14,15	0.66	0	17,19,21	1.00	0
2	NAG	P	2	2	14,14,15	0.64	0	17,19,21	1.06	1 (5%)
5	NAG	Q	1	1,5	14,14,15	0.78	0	17,19,21	0.92	1 (5%)
5	NAG	Q	2	5	14,14,15	0.78	0	17,19,21	1.14	1 (5%)
5	MAN	Q	3	5	11,11,12	0.78	0	15,15,17	0.80	0
2	NAG	R	1	1,2	14,14,15	0.82	1 (7%)	17,19,21	1.25	2 (11%)
2	NAG	R	2	2	14,14,15	0.70	0	17,19,21	1.36	1 (5%)
2	NAG	S	1	1,2	14,14,15	0.86	1 (7%)	17,19,21	1.91	5 (29%)
2	NAG	S	2	2	14,14,15	0.73	0	17,19,21	1.13	1 (5%)
2	NAG	T	1	1,2	14,14,15	0.62	0	17,19,21	1.17	1 (5%)
2	NAG	T	2	2	14,14,15	0.74	0	17,19,21	1.34	2 (11%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	4/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	1/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	F	1	3,1	-	0/6/23/26	0/1/1/1
3	MAN	F	10	3	-	0/2/19/22	0/1/1/1
3	MAN	F	11	3	-	0/2/19/22	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	0/2/19/22	0/1/1/1
3	MAN	F	4	3	-	0/2/19/22	0/1/1/1
3	MAN	F	5	3	-	0/2/19/22	0/1/1/1
3	MAN	F	6	3	-	0/2/19/22	0/1/1/1
3	MAN	F	7	3	-	0/2/19/22	0/1/1/1
3	MAN	F	8	3	-	2/2/19/22	0/1/1/1
3	MAN	F	9	3	-	0/2/19/22	0/1/1/1
4	NAG	G	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	1/6/23/26	0/1/1/1
4	BMA	G	3	4	-	0/2/19/22	0/1/1/1
2	NAG	H	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	H	2	2	-	1/6/23/26	0/1/1/1
5	NAG	I	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	3/6/23/26	0/1/1/1
5	MAN	I	3	5	-	0/2/19/22	0/1/1/1
2	NAG	J	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	NAG	K	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	K	2	2	-	2/6/23/26	0/1/1/1
2	NAG	L	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	L	2	2	-	0/6/23/26	0/1/1/1
2	NAG	M	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	M	2	2	-	2/6/23/26	0/1/1/1
3	NAG	N	1	3,1	-	2/6/23/26	0/1/1/1
3	MAN	N	10	3	-	2/2/19/22	0/1/1/1
3	MAN	N	11	3	-	0/2/19/22	0/1/1/1
3	NAG	N	2	3	-	0/6/23/26	0/1/1/1
3	BMA	N	3	3	-	0/2/19/22	0/1/1/1
3	MAN	N	4	3	-	2/2/19/22	0/1/1/1
3	MAN	N	5	3	-	0/2/19/22	0/1/1/1
3	MAN	N	6	3	-	0/2/19/22	0/1/1/1
3	MAN	N	7	3	-	0/2/19/22	0/1/1/1
3	MAN	N	8	3	-	2/2/19/22	0/1/1/1
3	MAN	N	9	3	-	2/2/19/22	0/1/1/1
4	NAG	O	1	4,1	-	0/6/23/26	0/1/1/1

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

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	5	MAN	O5-C1	-2.47	1.39	1.43
3	F	8	MAN	O5-C1	-2.42	1.39	1.43
3	F	2	NAG	O5-C1	-2.20	1.40	1.43
2	R	1	NAG	O5-C1	-2.13	1.40	1.43
3	F	7	MAN	O5-C1	-2.07	1.40	1.43
3	N	9	MAN	O5-C1	-2.06	1.40	1.43
2	S	1	NAG	O5-C1	-2.04	1.40	1.43
3	N	4	MAN	O5-C1	-2.03	1.40	1.43

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	O	3	BMA	C1-O5-C5	8.01	122.93	112.19
4	G	3	BMA	C1-O5-C5	7.64	122.42	112.19
3	F	3	BMA	C1-O5-C5	6.75	121.23	112.19
3	N	3	BMA	C1-O5-C5	5.36	119.37	112.19
2	C	2	NAG	C2-N2-C7	5.09	129.72	122.90
2	S	1	NAG	C1-O5-C5	4.36	118.03	112.19
2	R	2	NAG	C2-N2-C7	4.23	128.56	122.90
4	O	3	BMA	C3-C4-C5	3.94	117.37	110.23
2	S	1	NAG	O5-C1-C2	-3.90	105.25	111.29
4	G	3	BMA	C3-C4-C5	3.83	117.17	110.23
2	J	1	NAG	O4-C4-C3	-3.70	101.65	110.38
2	J	1	NAG	C2-N2-C7	3.68	127.84	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	2	NAG	C1-O5-C5	3.67	117.10	112.19
2	D	1	NAG	C2-N2-C7	3.55	127.66	122.90
2	E	2	NAG	C2-N2-C7	3.51	127.60	122.90
2	L	1	NAG	C2-N2-C7	3.43	127.50	122.90
3	N	10	MAN	C1-O5-C5	3.40	116.74	112.19
2	M	2	NAG	C2-N2-C7	3.39	127.45	122.90
2	S	2	NAG	O5-C1-C2	-3.26	106.25	111.29
3	N	1	NAG	C1-O5-C5	3.25	116.55	112.19
2	J	2	NAG	C2-N2-C7	3.20	127.19	122.90
2	R	1	NAG	C2-N2-C7	3.12	127.08	122.90
2	S	1	NAG	O5-C5-C6	-2.97	101.88	107.66
5	Q	2	NAG	C2-N2-C7	2.89	126.78	122.90
2	K	1	NAG	C1-O5-C5	2.85	116.01	112.19
3	N	3	BMA	O4-C4-C3	-2.84	103.69	110.38
4	G	2	NAG	C1-O5-C5	2.81	115.96	112.19
4	G	3	BMA	C2-C3-C4	2.80	115.78	110.86
3	N	2	NAG	C1-O5-C5	2.78	115.92	112.19
3	F	3	BMA	O6-C6-C5	-2.78	101.88	111.33
2	T	1	NAG	C1-O5-C5	2.76	115.89	112.19
2	H	1	NAG	O5-C1-C2	-2.75	107.04	111.29
3	F	7	MAN	C1-O5-C5	2.73	115.85	112.19
3	F	11	MAN	C1-O5-C5	2.69	115.79	112.19
3	F	3	BMA	C1-C2-C3	2.64	113.49	109.64
3	N	4	MAN	C1-O5-C5	2.57	115.63	112.19
3	F	3	BMA	O4-C4-C3	-2.55	104.36	110.38
2	H	2	NAG	O5-C1-C2	-2.55	107.35	111.29
3	F	1	NAG	O4-C4-C3	-2.51	104.45	110.38
3	F	1	NAG	O5-C1-C2	-2.51	107.41	111.29
2	T	2	NAG	C4-C3-C2	-2.44	107.44	111.02
3	N	7	MAN	O2-C2-C3	2.40	115.12	110.15
3	N	1	NAG	O5-C1-C2	-2.38	107.60	111.29
2	S	1	NAG	O4-C4-C3	-2.38	104.77	110.38
2	K	2	NAG	O5-C1-C2	-2.38	107.61	111.29
4	O	3	BMA	C2-C3-C4	2.36	115.02	110.86
3	N	3	BMA	O6-C6-C5	-2.35	103.32	111.33
4	O	3	BMA	O4-C4-C3	-2.35	104.84	110.38
3	N	2	NAG	O4-C4-C3	-2.35	104.84	110.38
2	J	2	NAG	O5-C1-C2	-2.33	107.68	111.29
3	N	5	MAN	C1-O5-C5	2.32	115.30	112.19
5	I	2	NAG	O5-C1-C2	-2.32	107.70	111.29
3	F	4	MAN	C1-O5-C5	2.30	115.27	112.19
3	F	2	NAG	C1-O5-C5	2.29	115.26	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	S	1	NAG	C1-C2-N2	2.29	114.05	110.43
3	N	8	MAN	C1-O5-C5	2.28	115.24	112.19
4	G	3	BMA	O4-C4-C3	-2.27	105.02	110.38
2	K	1	NAG	O5-C1-C2	-2.27	107.78	111.29
4	O	3	BMA	O5-C5-C4	2.26	116.32	110.83
2	K	2	NAG	C1-O5-C5	2.26	115.21	112.19
2	J	1	NAG	O5-C1-C2	-2.25	107.81	111.29
2	K	2	NAG	C1-C2-N2	2.22	113.93	110.43
4	O	1	NAG	O5-C1-C2	-2.21	107.87	111.29
2	R	1	NAG	O4-C4-C3	-2.19	105.21	110.38
2	J	1	NAG	C4-C3-C2	2.18	114.22	111.02
3	F	9	MAN	C1-O5-C5	2.17	115.09	112.19
5	Q	1	NAG	O5-C1-C2	-2.13	108.00	111.29
2	M	1	NAG	C1-O5-C5	2.13	115.04	112.19
2	P	2	NAG	C1-O5-C5	2.13	115.03	112.19
3	N	7	MAN	C1-O5-C5	2.11	115.01	112.19
4	O	3	BMA	O3-C3-C2	-2.09	105.79	110.05
2	L	2	NAG	C1-O5-C5	2.08	114.97	112.19
3	F	10	MAN	C1-O5-C5	2.07	114.95	112.19
5	I	3	MAN	C1-O5-C5	2.06	114.95	112.19
3	F	1	NAG	C4-C3-C2	2.05	114.02	111.02
3	N	3	BMA	O2-C2-C3	2.04	114.38	110.15
4	G	3	BMA	O3-C3-C2	-2.03	105.91	110.05
3	F	2	NAG	O4-C4-C3	-2.03	105.59	110.38
3	F	5	MAN	C1-C2-C3	-2.03	106.69	109.64
3	F	1	NAG	C8-C7-N2	-2.02	112.77	116.12
3	F	3	BMA	C3-C4-C5	2.01	113.87	110.23

There are no chirality outliers.

All (59) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	N	4	MAN	O5-C5-C6-O6
2	R	1	NAG	O5-C5-C6-O6
3	N	10	MAN	O5-C5-C6-O6
3	N	1	NAG	C4-C5-C6-O6
2	R	1	NAG	C4-C5-C6-O6
3	N	1	NAG	O5-C5-C6-O6
3	N	4	MAN	C4-C5-C6-O6
3	N	10	MAN	C4-C5-C6-O6
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	D	1	NAG	C8-C7-N2-C2
2	D	1	NAG	O7-C7-N2-C2
2	E	2	NAG	C8-C7-N2-C2
2	E	2	NAG	O7-C7-N2-C2
2	J	1	NAG	C8-C7-N2-C2
2	J	1	NAG	O7-C7-N2-C2
2	J	2	NAG	C8-C7-N2-C2
2	J	2	NAG	O7-C7-N2-C2
2	L	1	NAG	C8-C7-N2-C2
2	L	1	NAG	O7-C7-N2-C2
2	M	2	NAG	C8-C7-N2-C2
2	M	2	NAG	O7-C7-N2-C2
2	R	1	NAG	C8-C7-N2-C2
2	R	1	NAG	O7-C7-N2-C2
2	R	2	NAG	C8-C7-N2-C2
2	R	2	NAG	O7-C7-N2-C2
4	O	2	NAG	O5-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
2	H	1	NAG	O5-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
5	Q	1	NAG	C4-C5-C6-O6
2	K	2	NAG	C4-C5-C6-O6
3	N	8	MAN	C4-C5-C6-O6
3	N	9	MAN	C4-C5-C6-O6
5	I	2	NAG	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
2	P	2	NAG	O5-C5-C6-O6
2	P	1	NAG	O5-C5-C6-O6
5	Q	2	NAG	O5-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
5	Q	1	NAG	O5-C5-C6-O6
5	Q	2	NAG	C3-C2-N2-C7
2	E	1	NAG	C4-C5-C6-O6
3	N	8	MAN	O5-C5-C6-O6
3	N	9	MAN	O5-C5-C6-O6
4	O	2	NAG	C4-C5-C6-O6
5	Q	3	MAN	C4-C5-C6-O6
2	T	2	NAG	C4-C5-C6-O6
2	C	2	NAG	C3-C2-N2-C7
5	I	2	NAG	C3-C2-N2-C7
2	E	1	NAG	O5-C5-C6-O6
3	F	8	MAN	O5-C5-C6-O6

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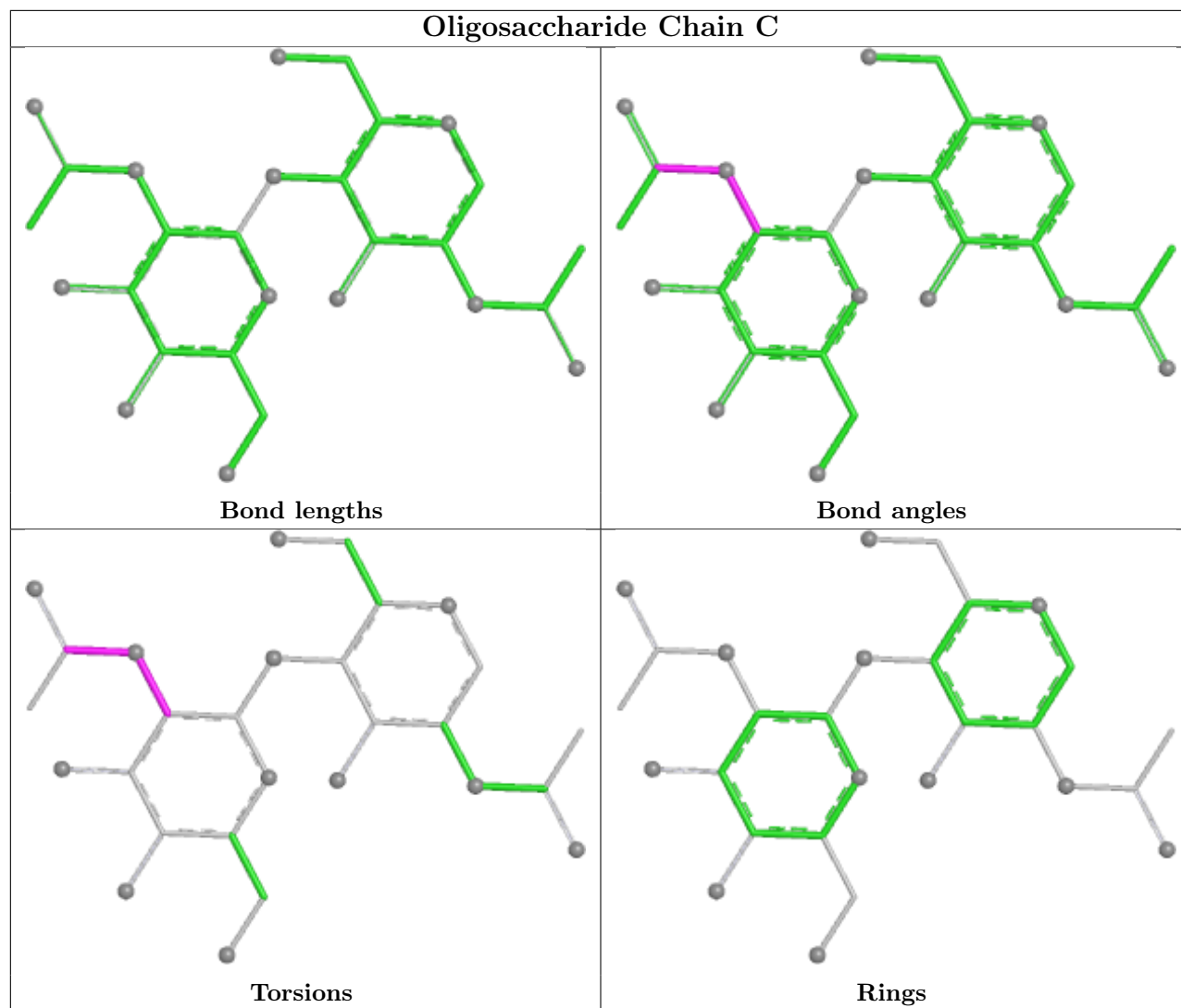
Mol	Chain	Res	Type	Atoms
3	F	8	MAN	C4-C5-C6-O6
2	C	2	NAG	C1-C2-N2-C7
2	D	2	NAG	C1-C2-N2-C7
2	T	2	NAG	C1-C2-N2-C7
5	I	2	NAG	C1-C2-N2-C7
2	S	1	NAG	O5-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6

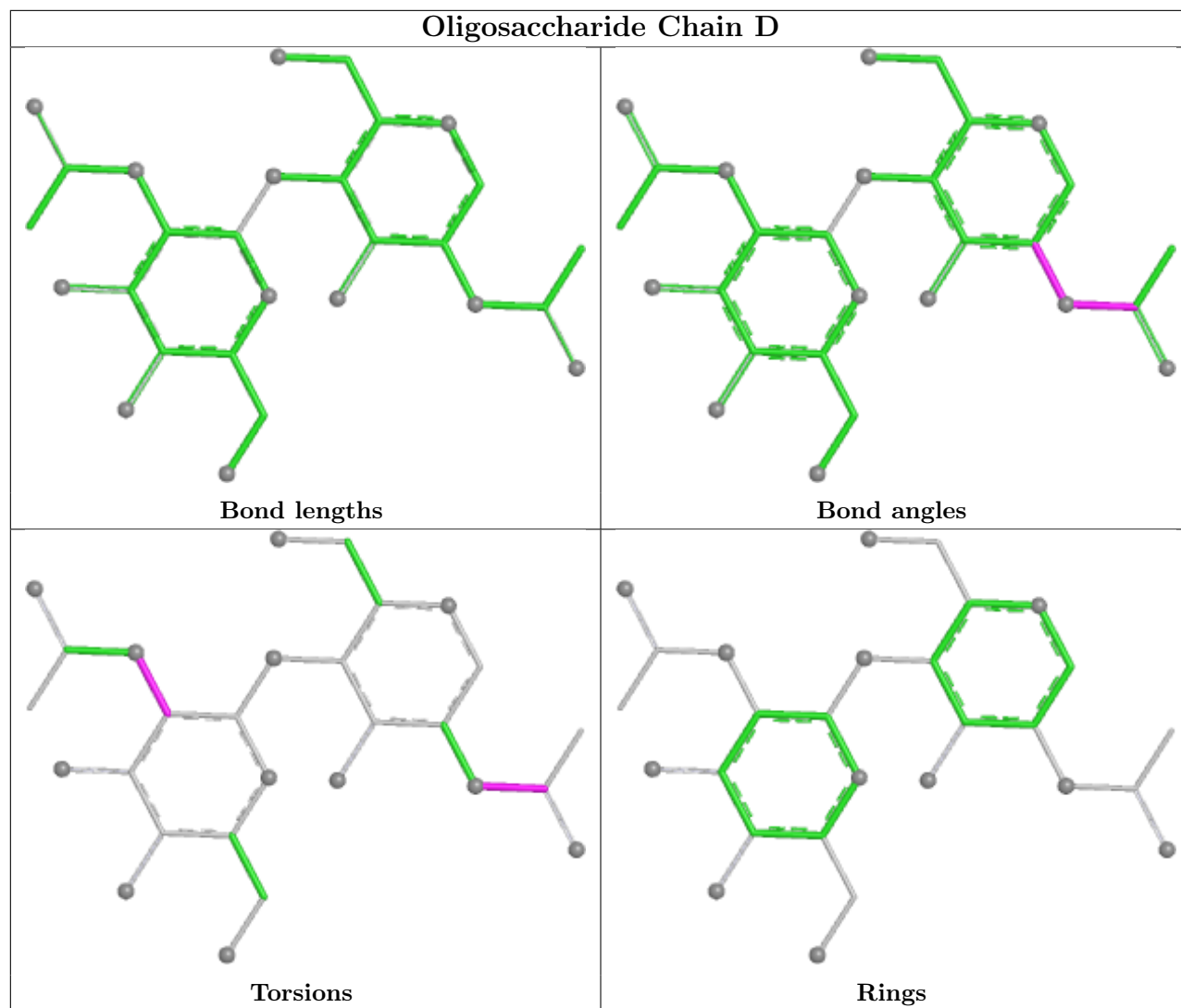
There are no ring outliers.

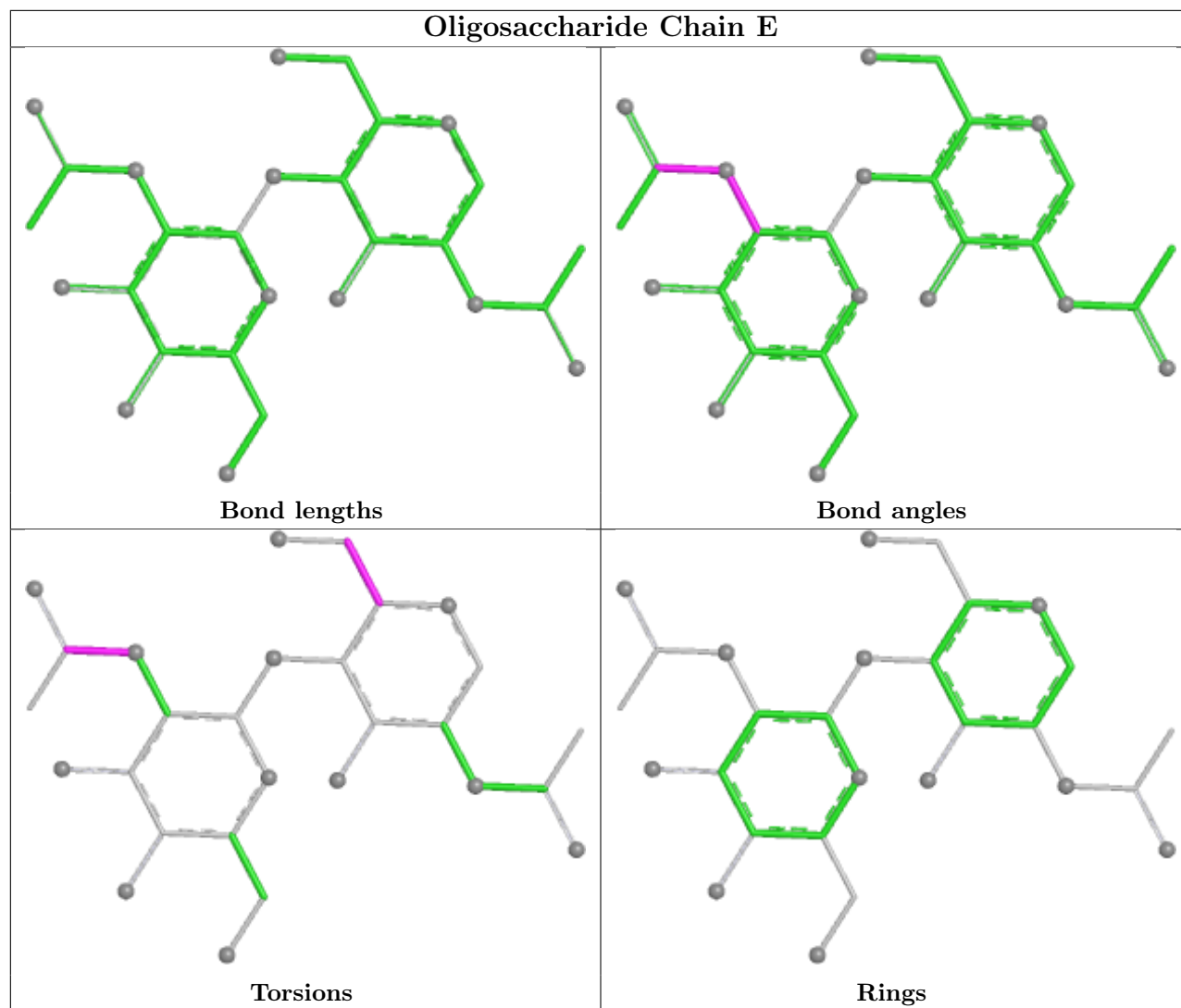
11 monomers are involved in 9 short contacts:

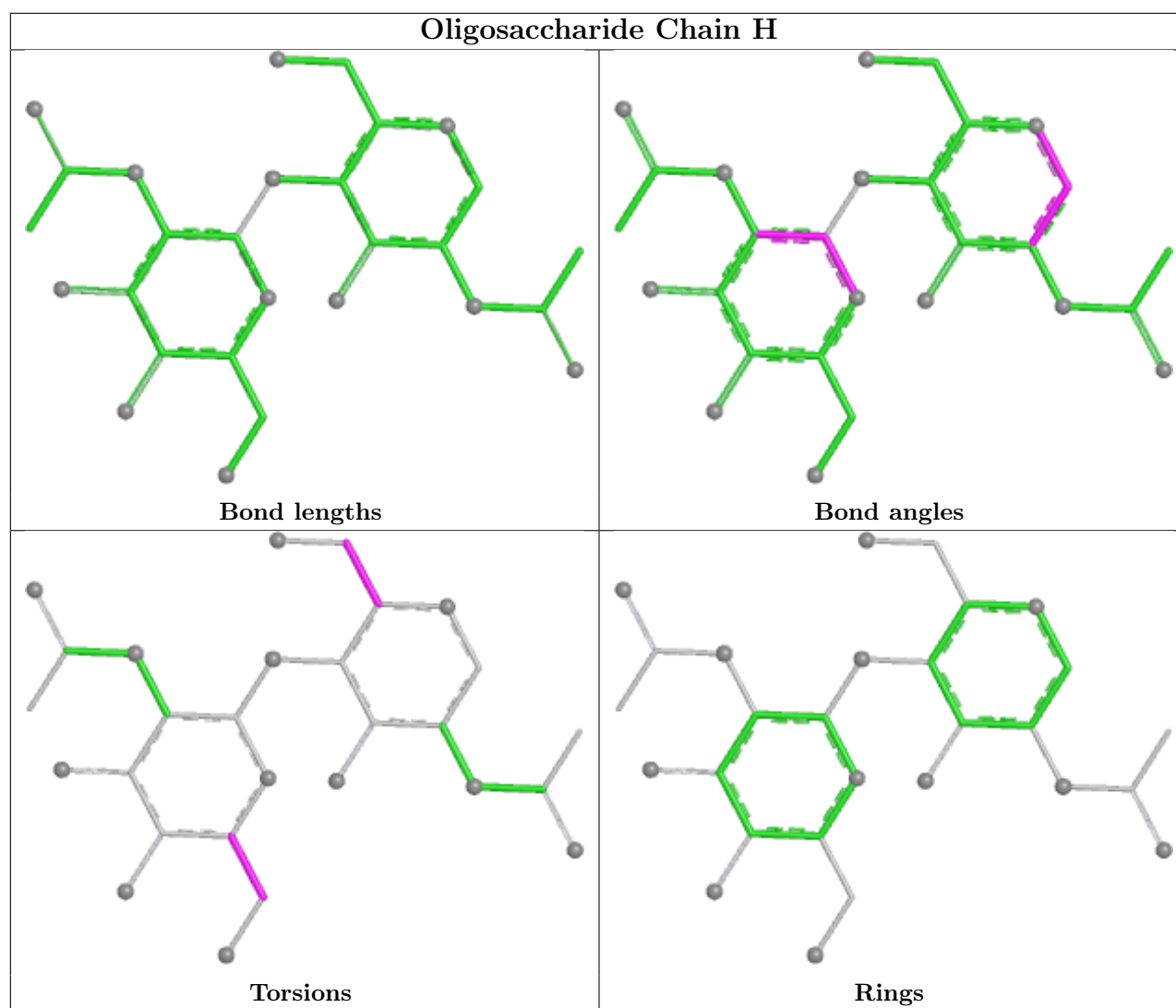
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	2	NAG	1	0
2	R	1	NAG	1	0
2	H	1	NAG	2	0
3	F	1	NAG	1	0
2	D	1	NAG	1	0
2	K	1	NAG	1	0
2	R	2	NAG	1	0
2	C	2	NAG	1	0
2	H	2	NAG	1	0
5	I	2	NAG	1	0
5	I	3	MAN	1	0

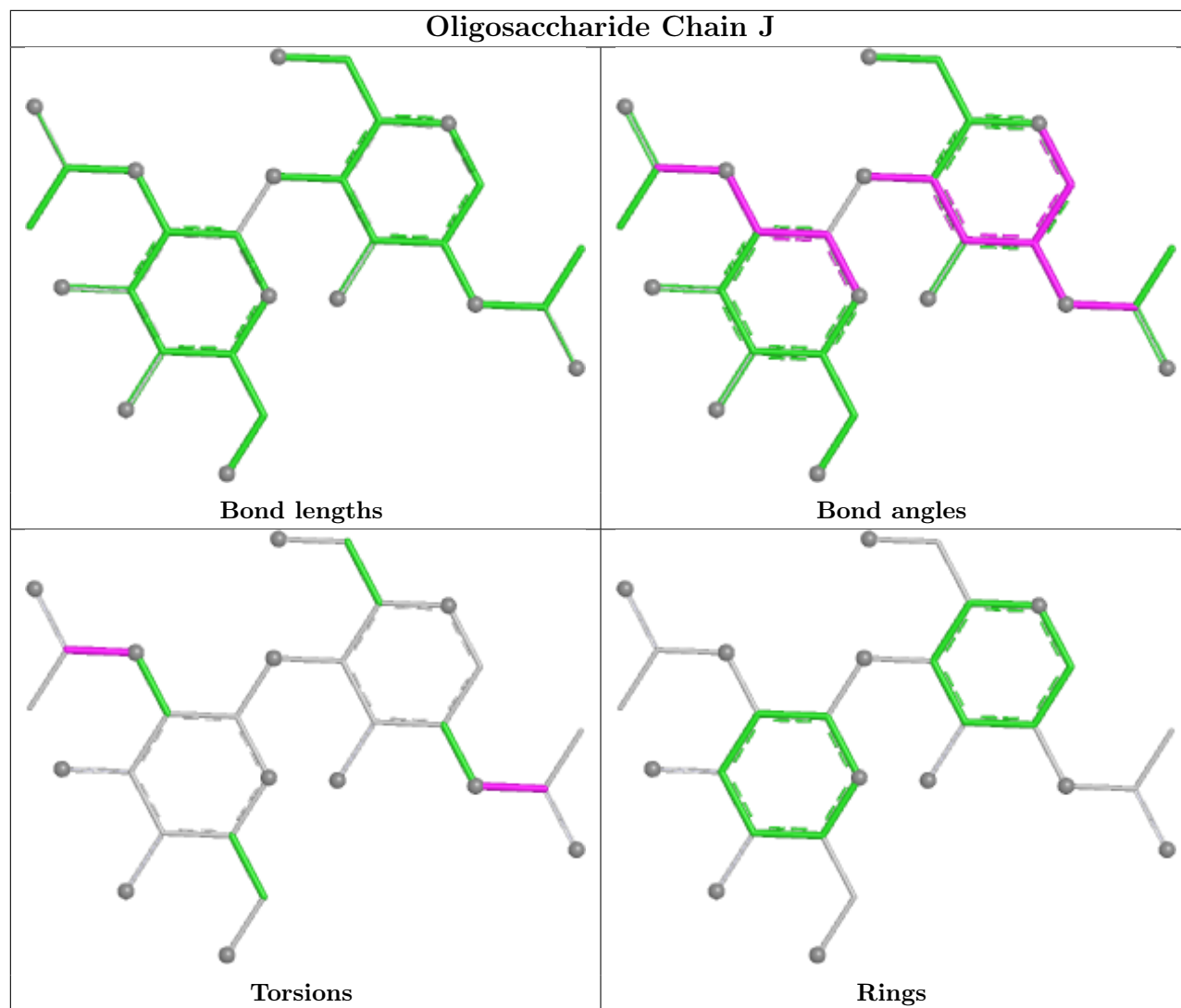
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

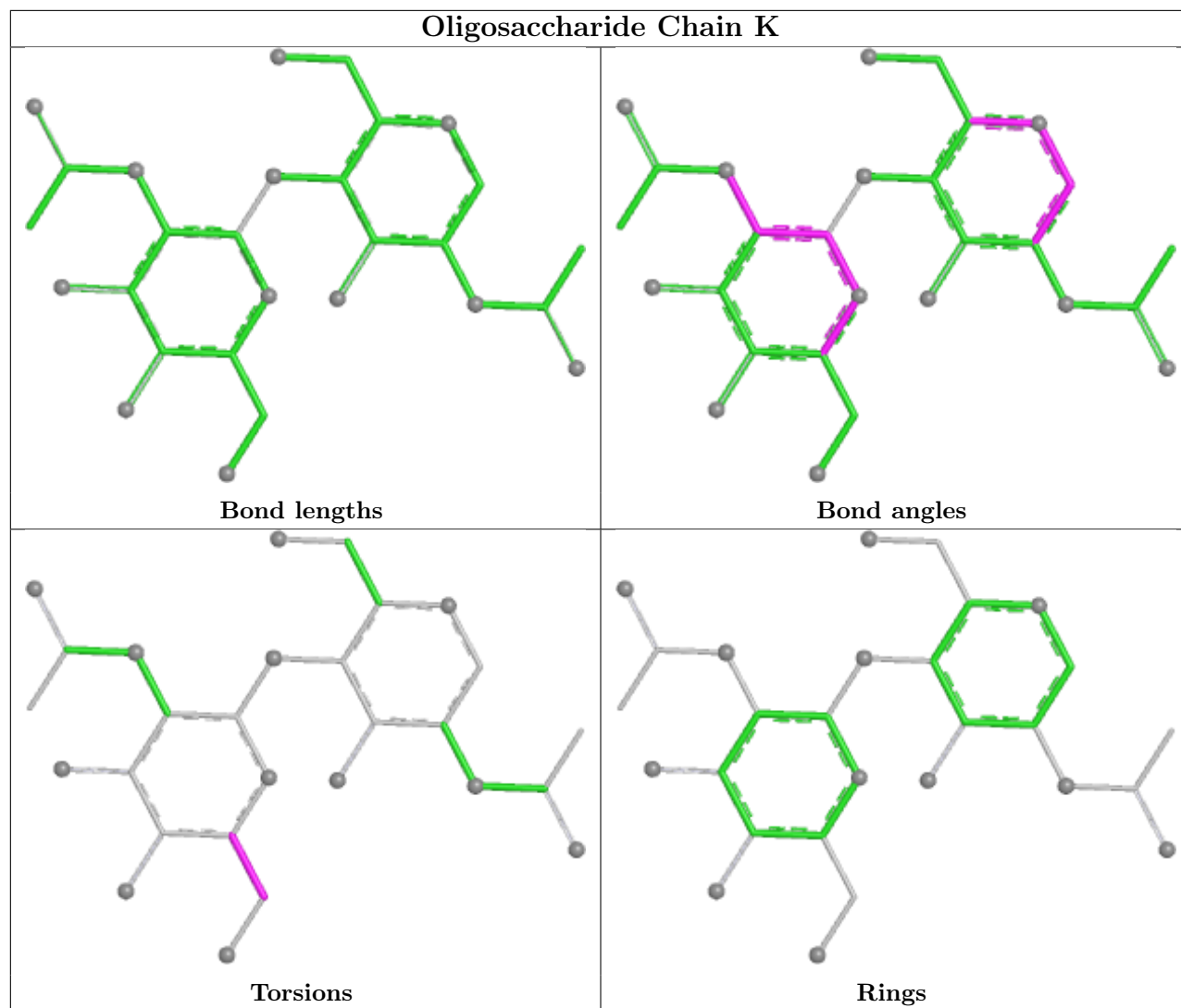


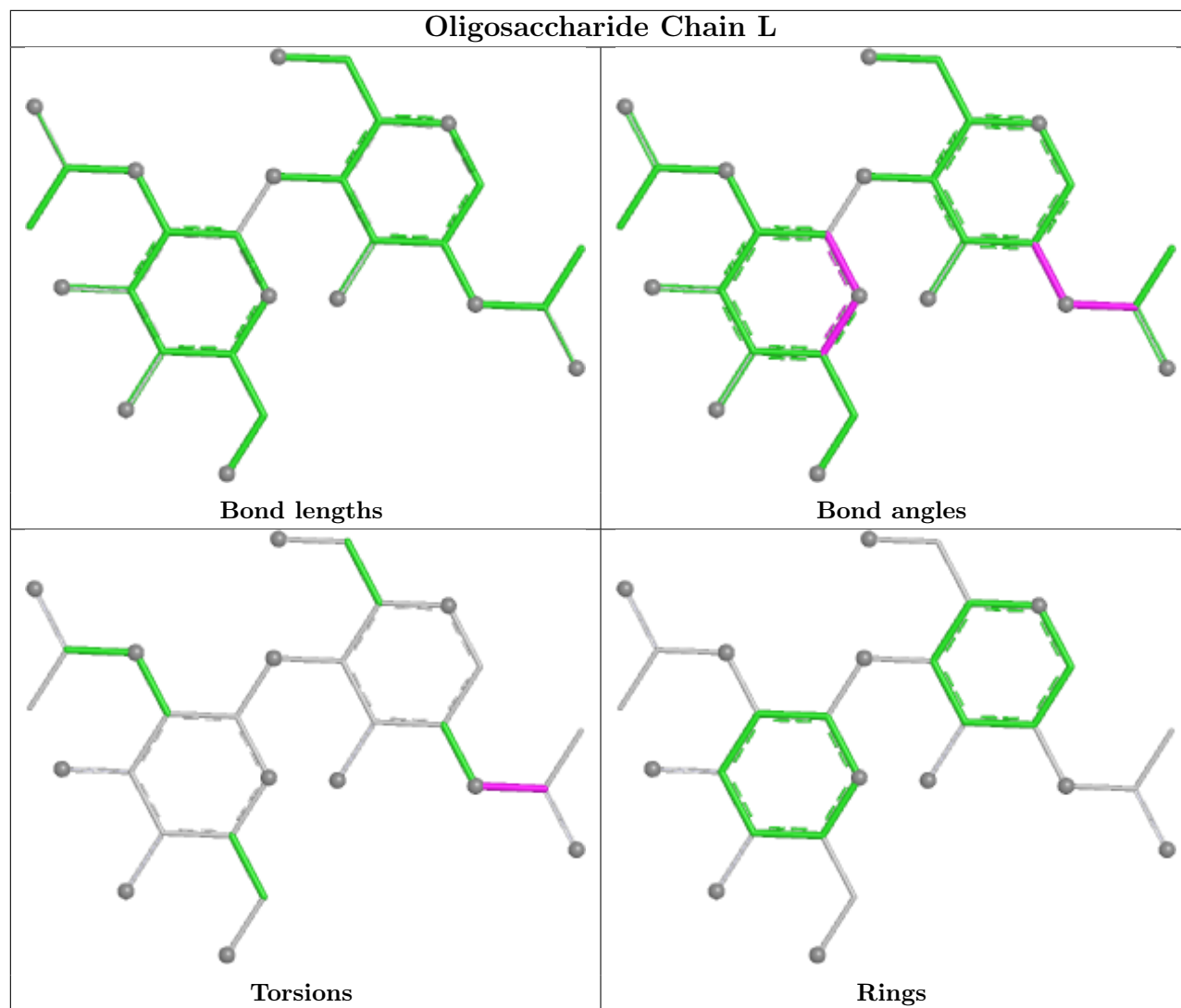


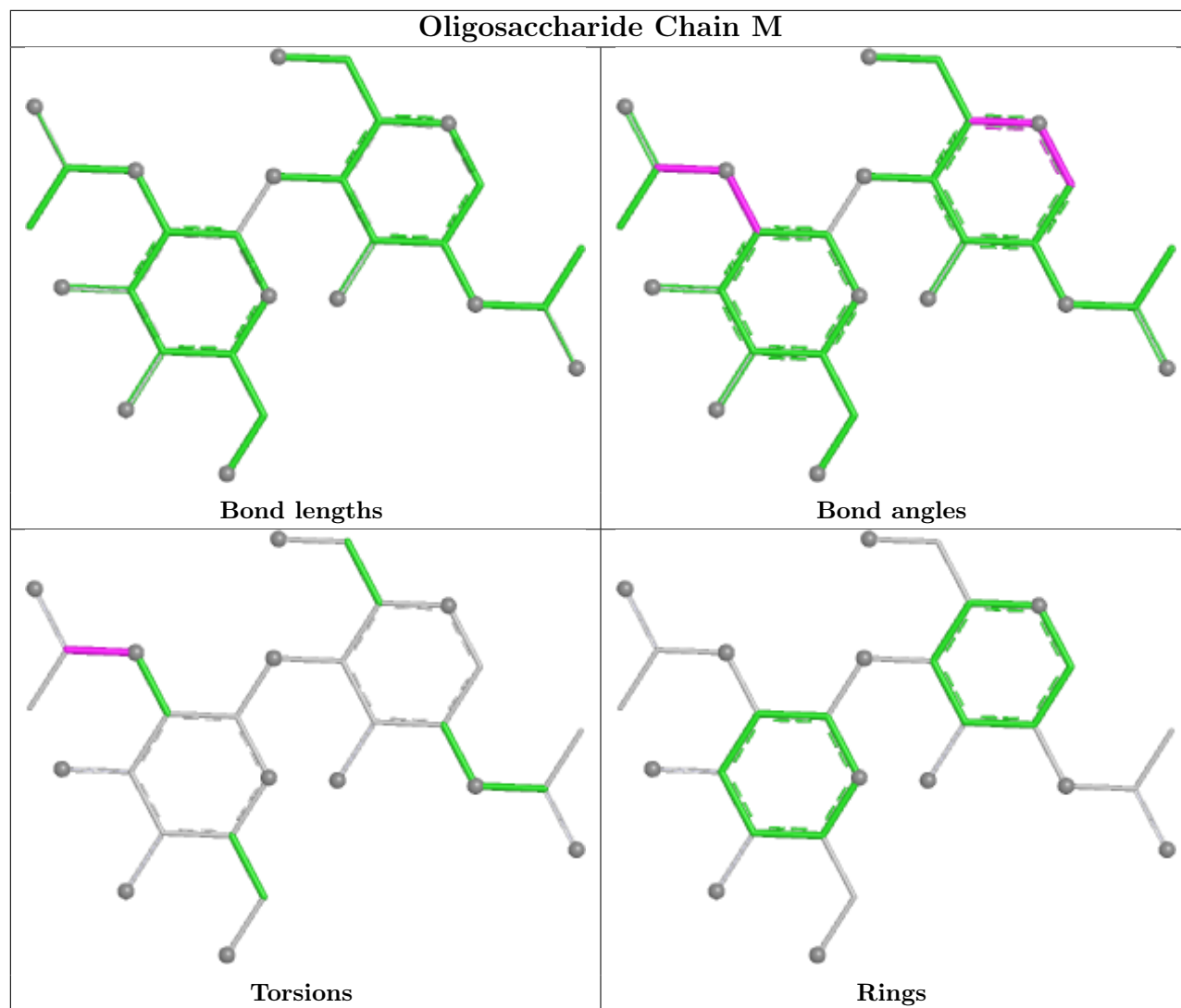


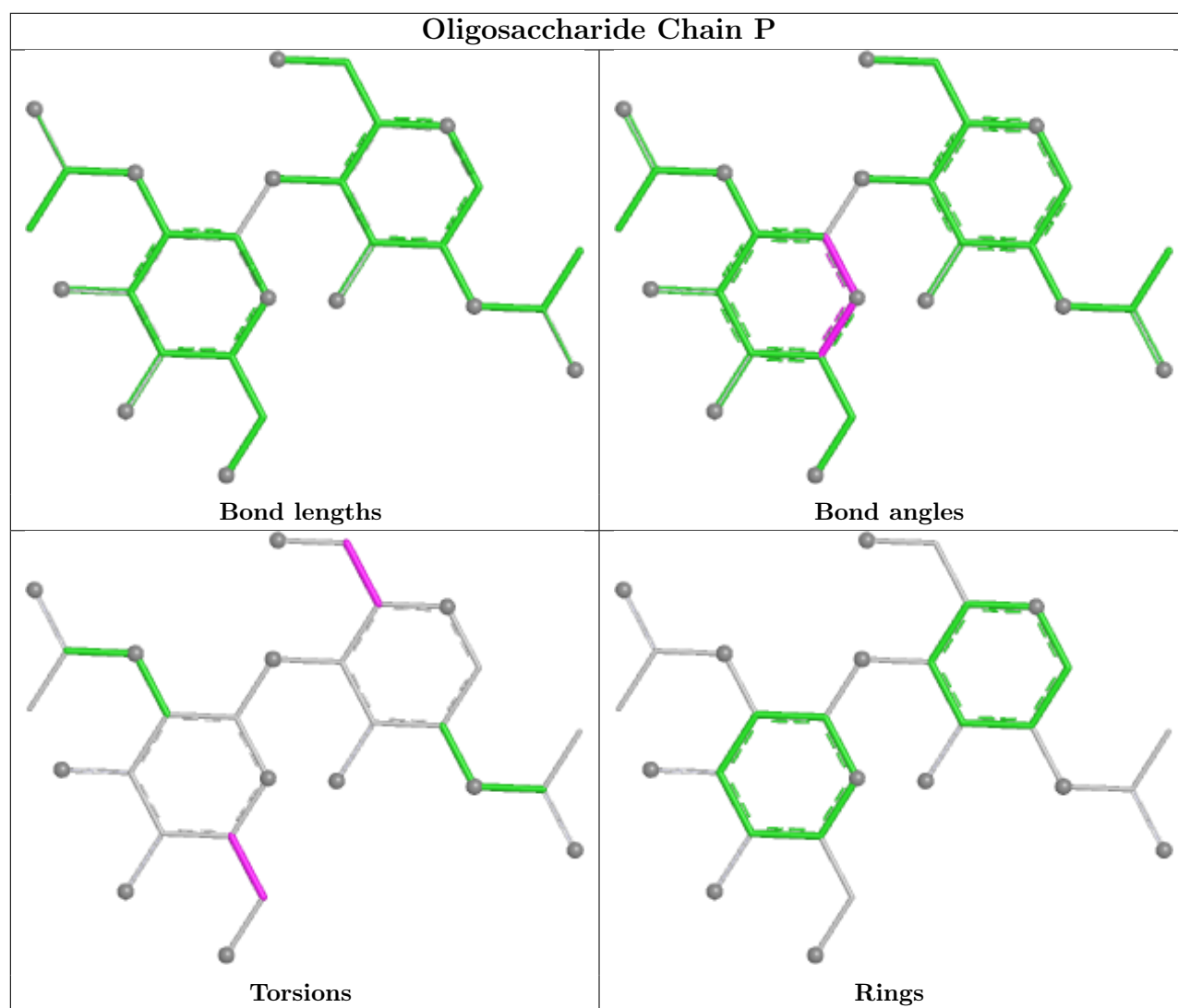


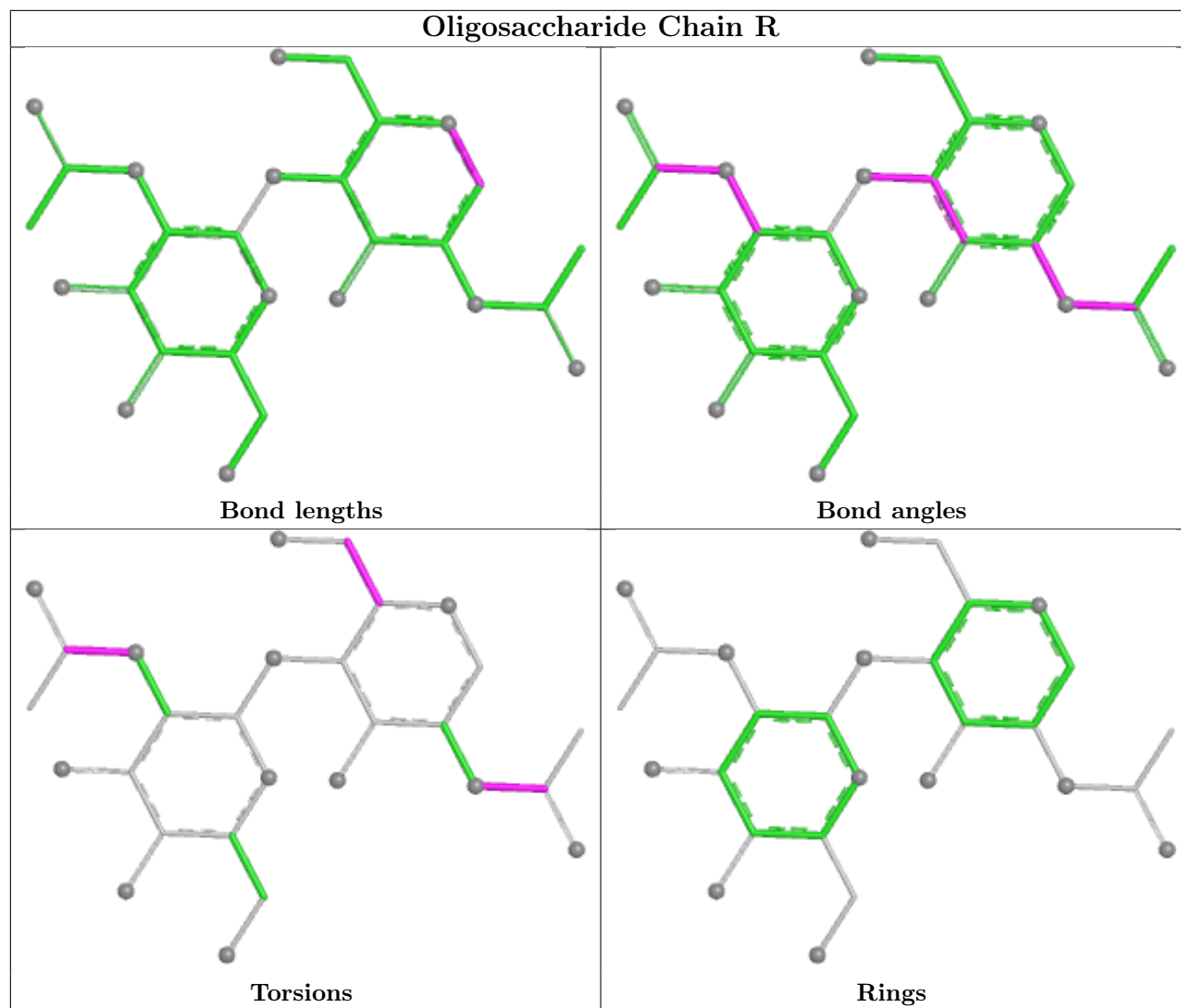


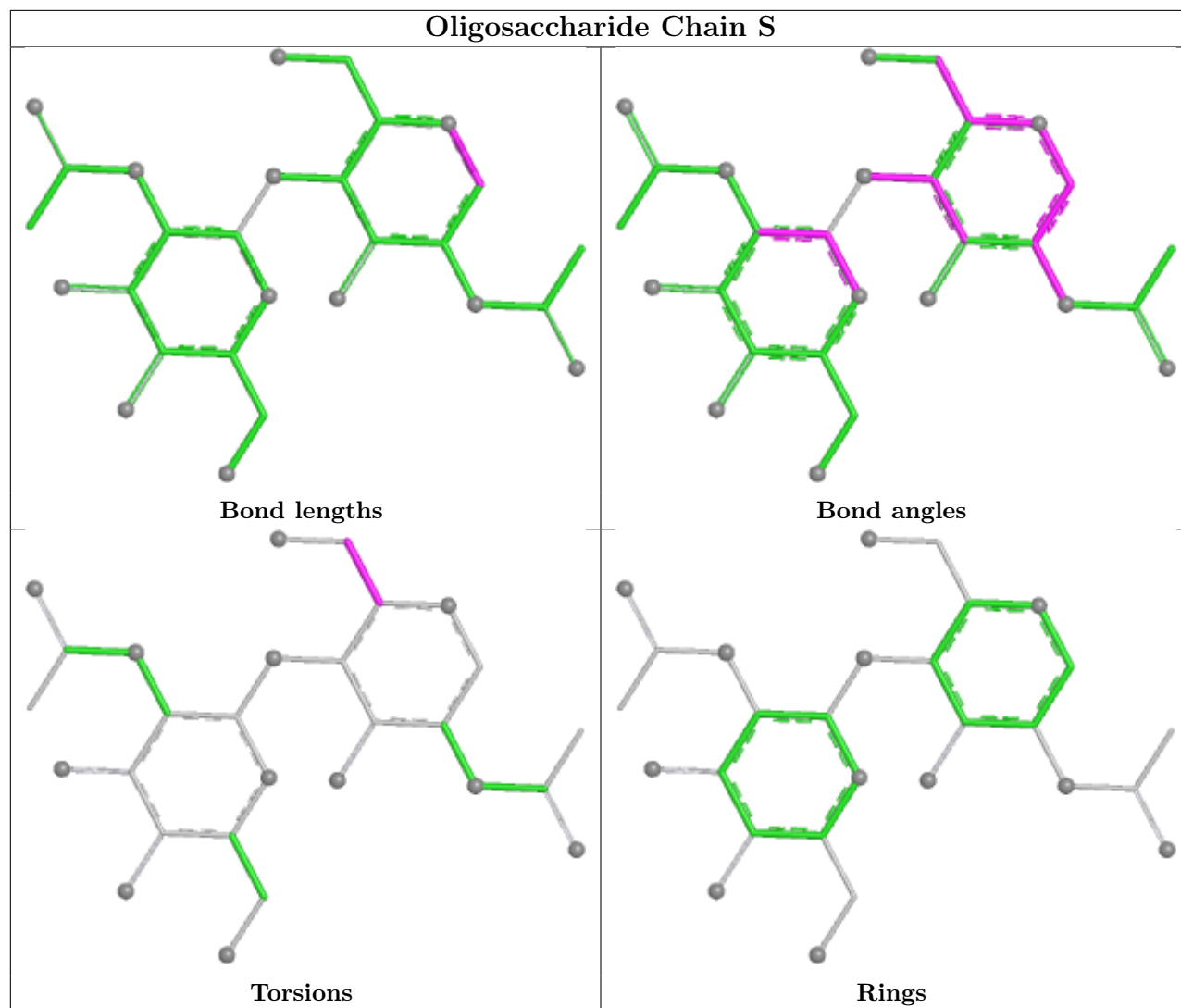


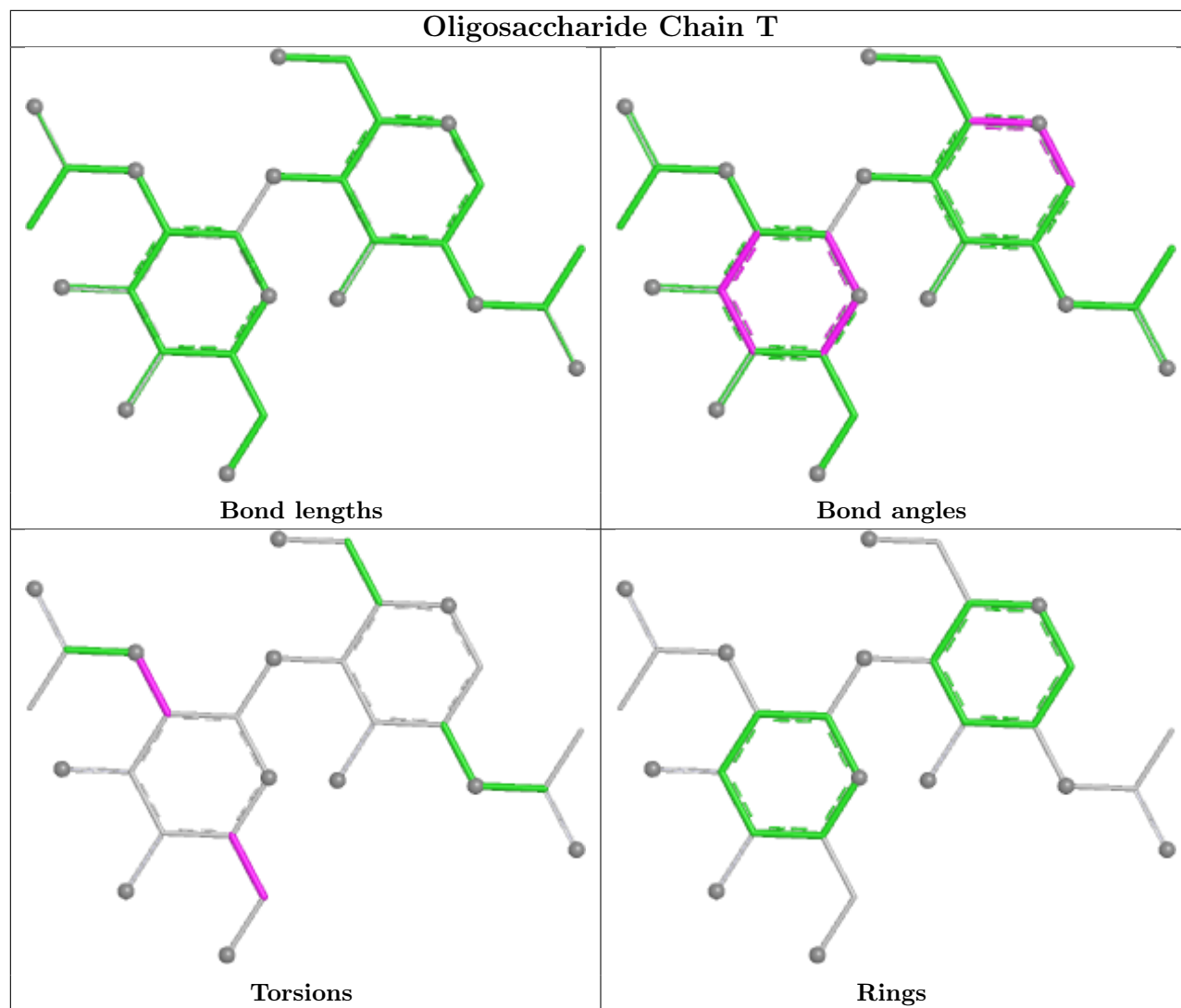




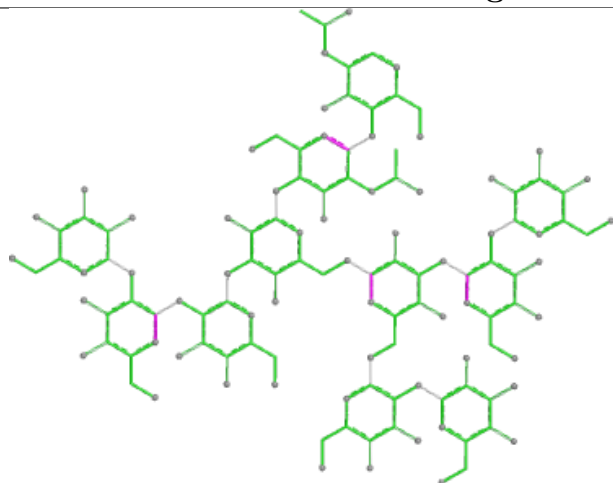




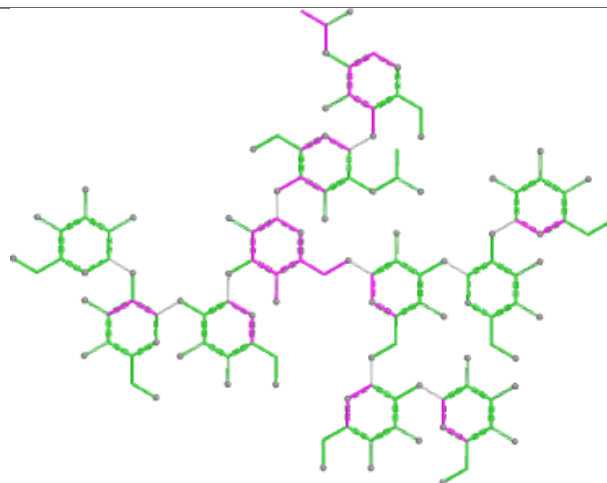




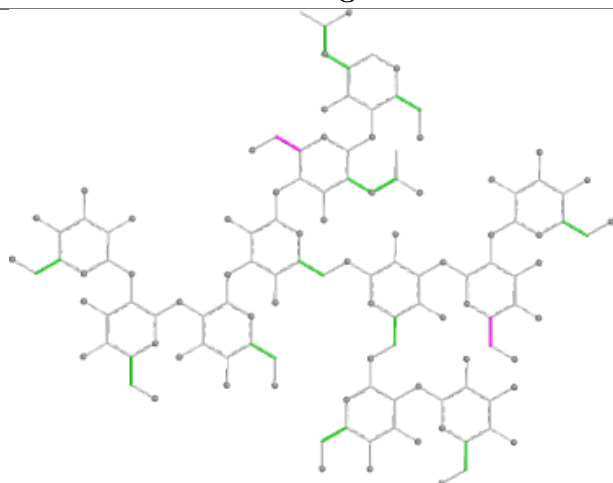
Oligosaccharide Chain F



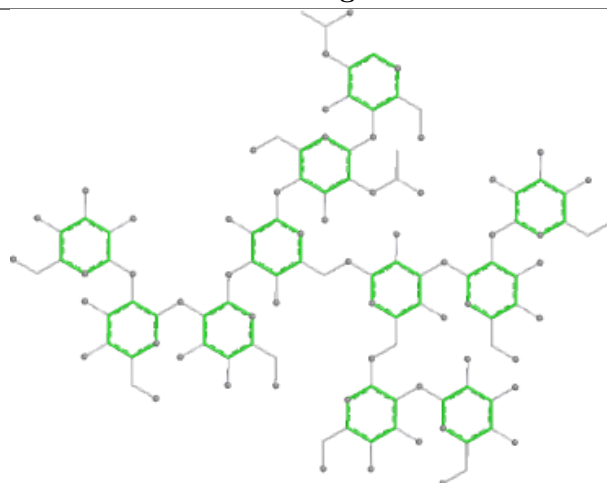
Bond lengths



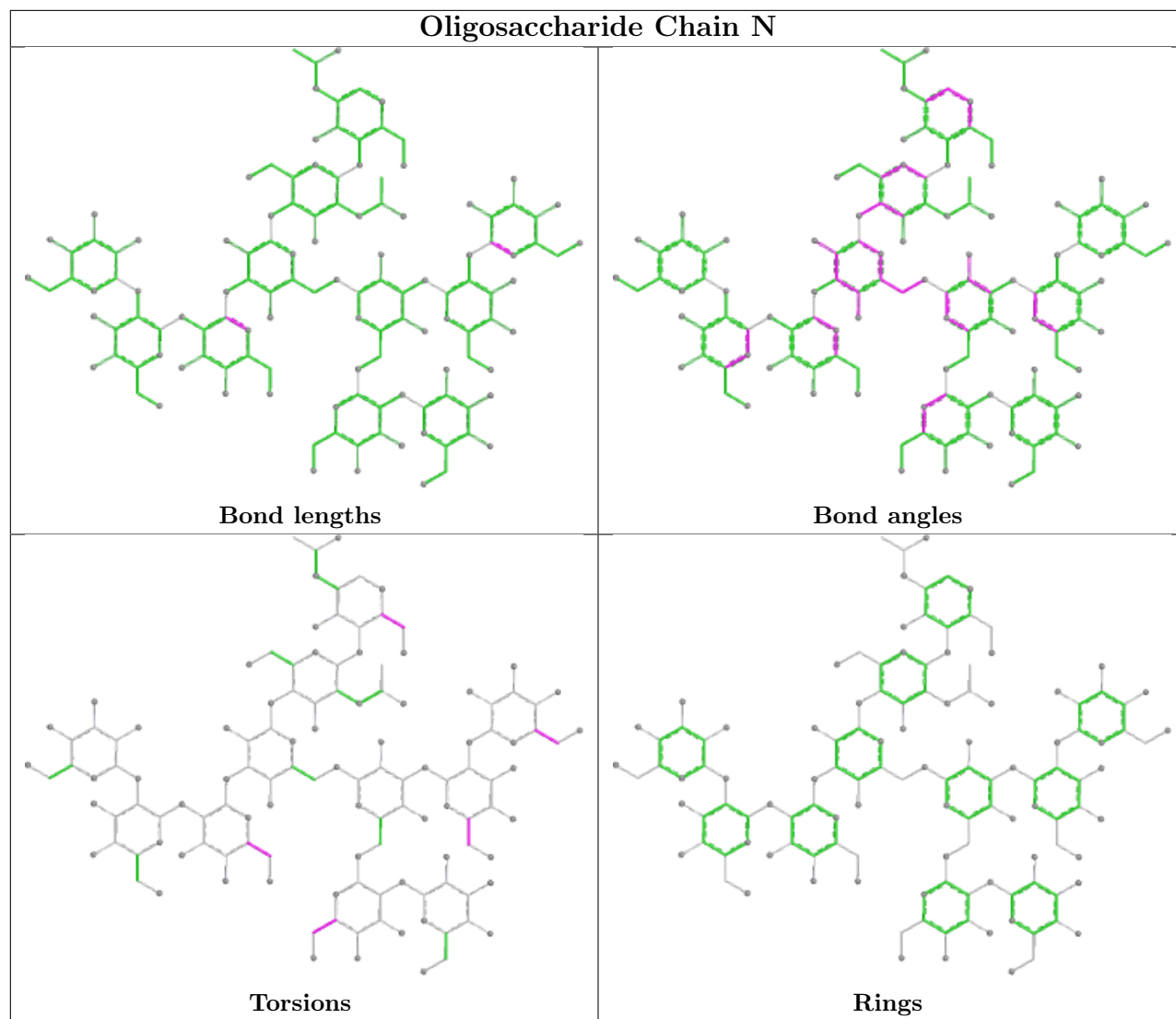
Bond angles

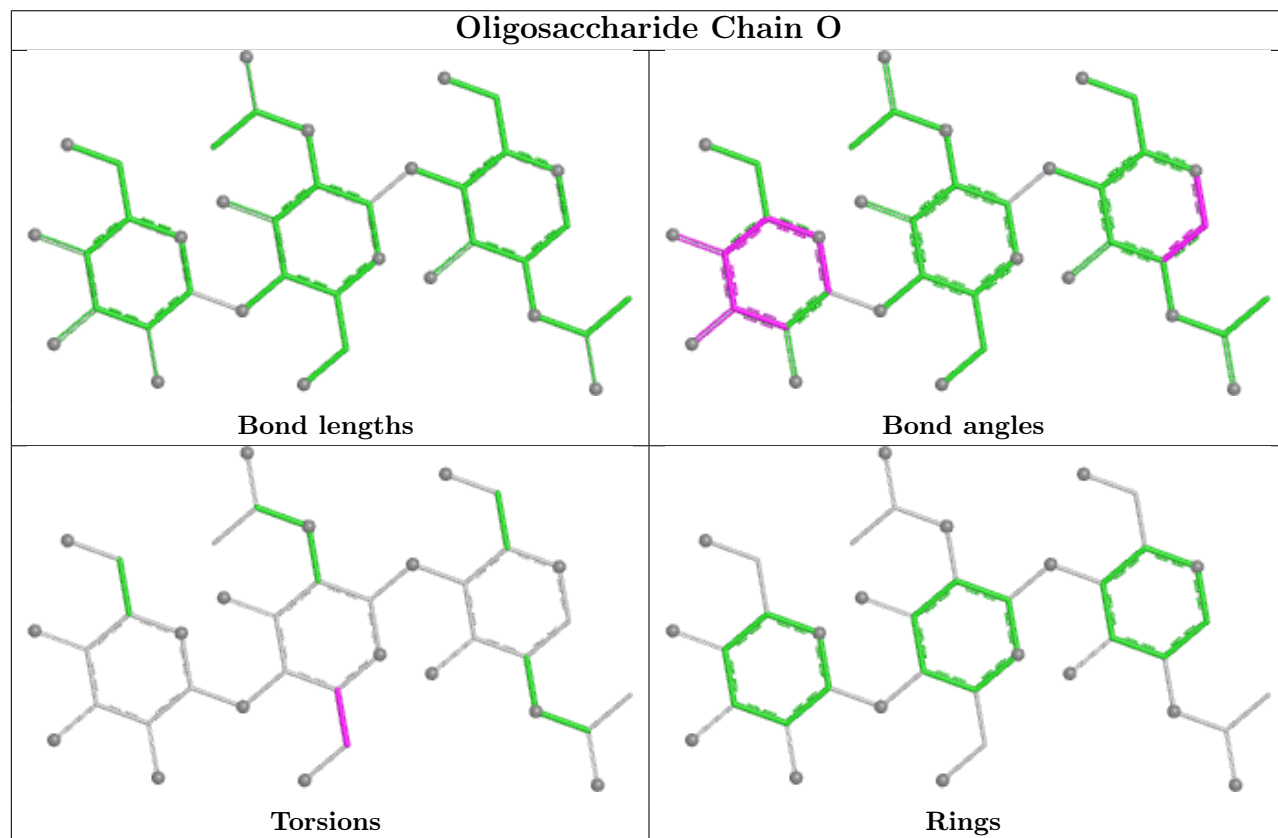
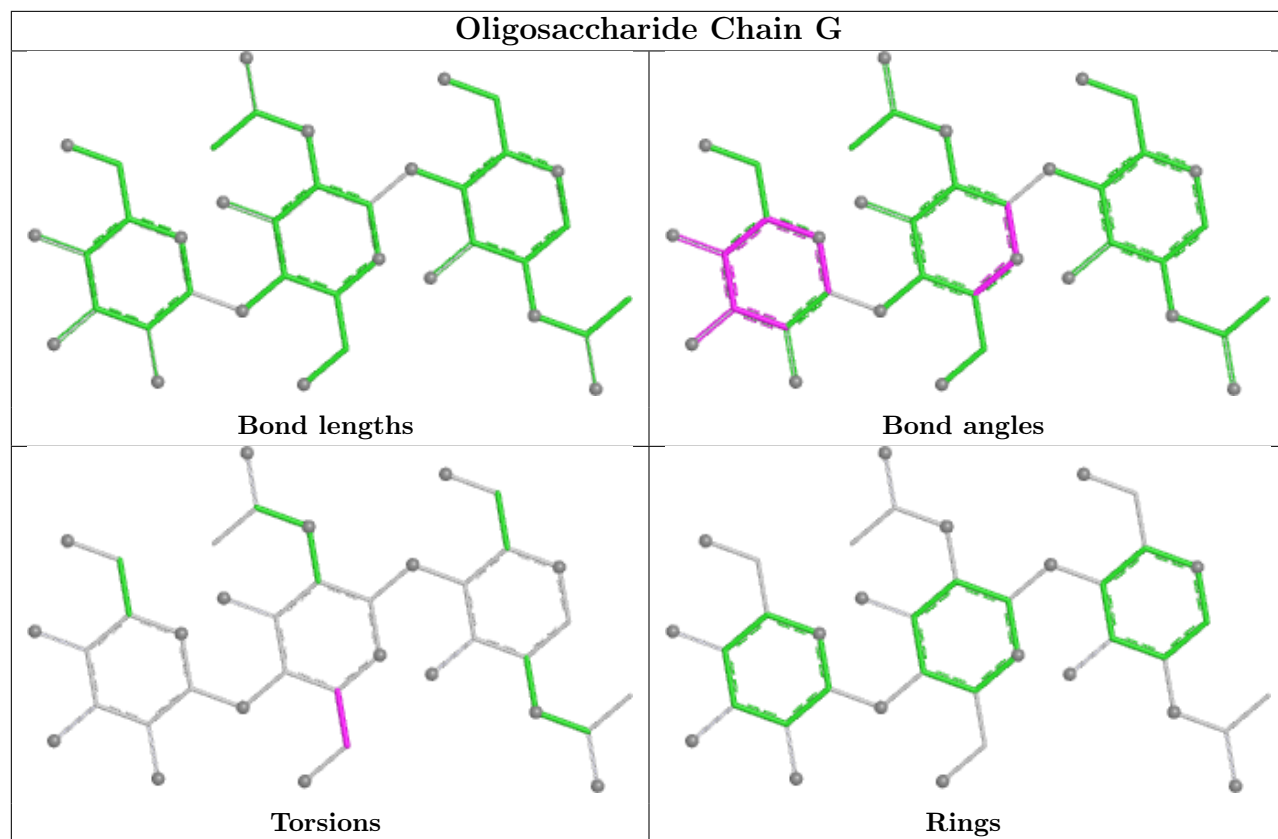


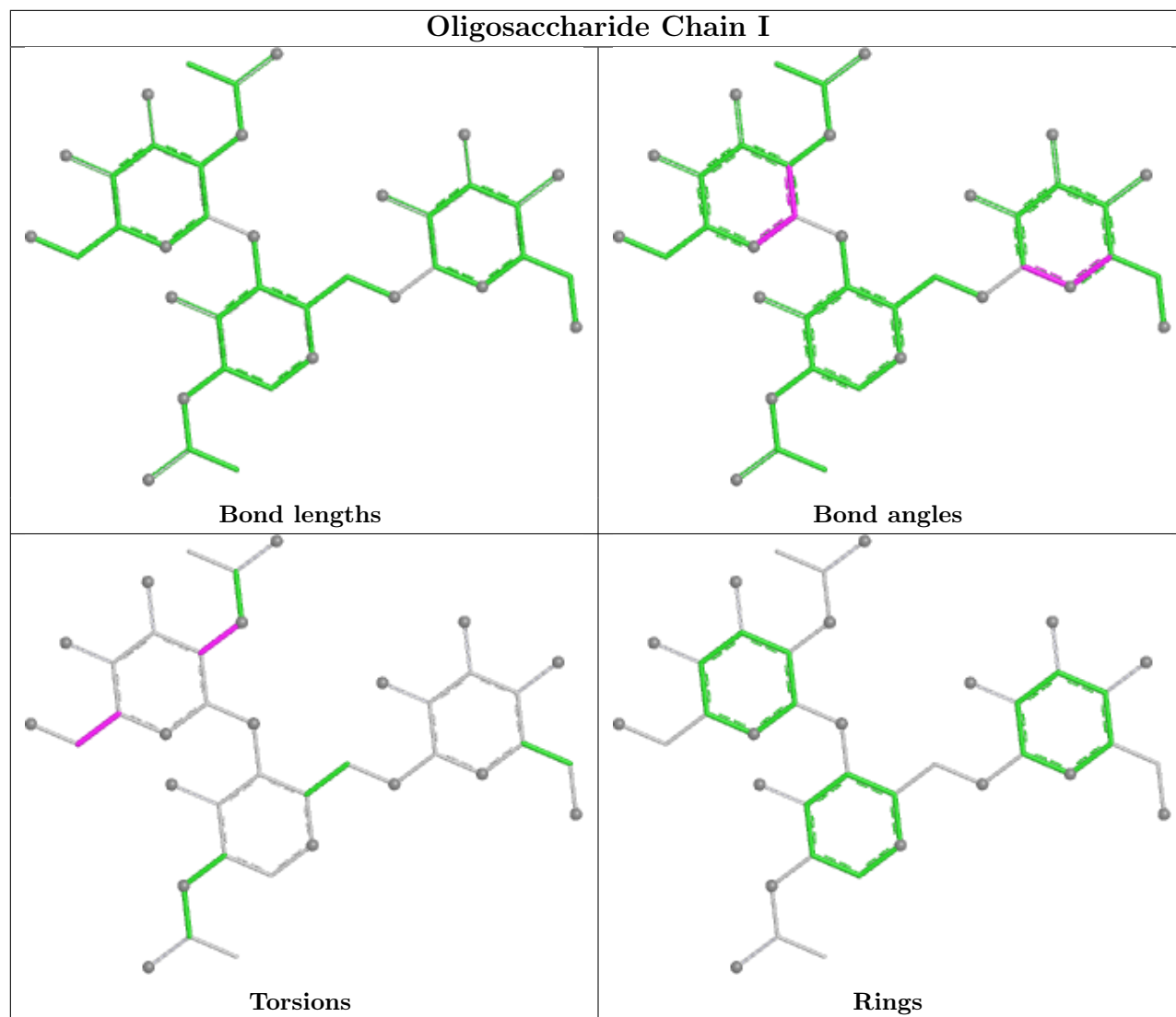
Torsions

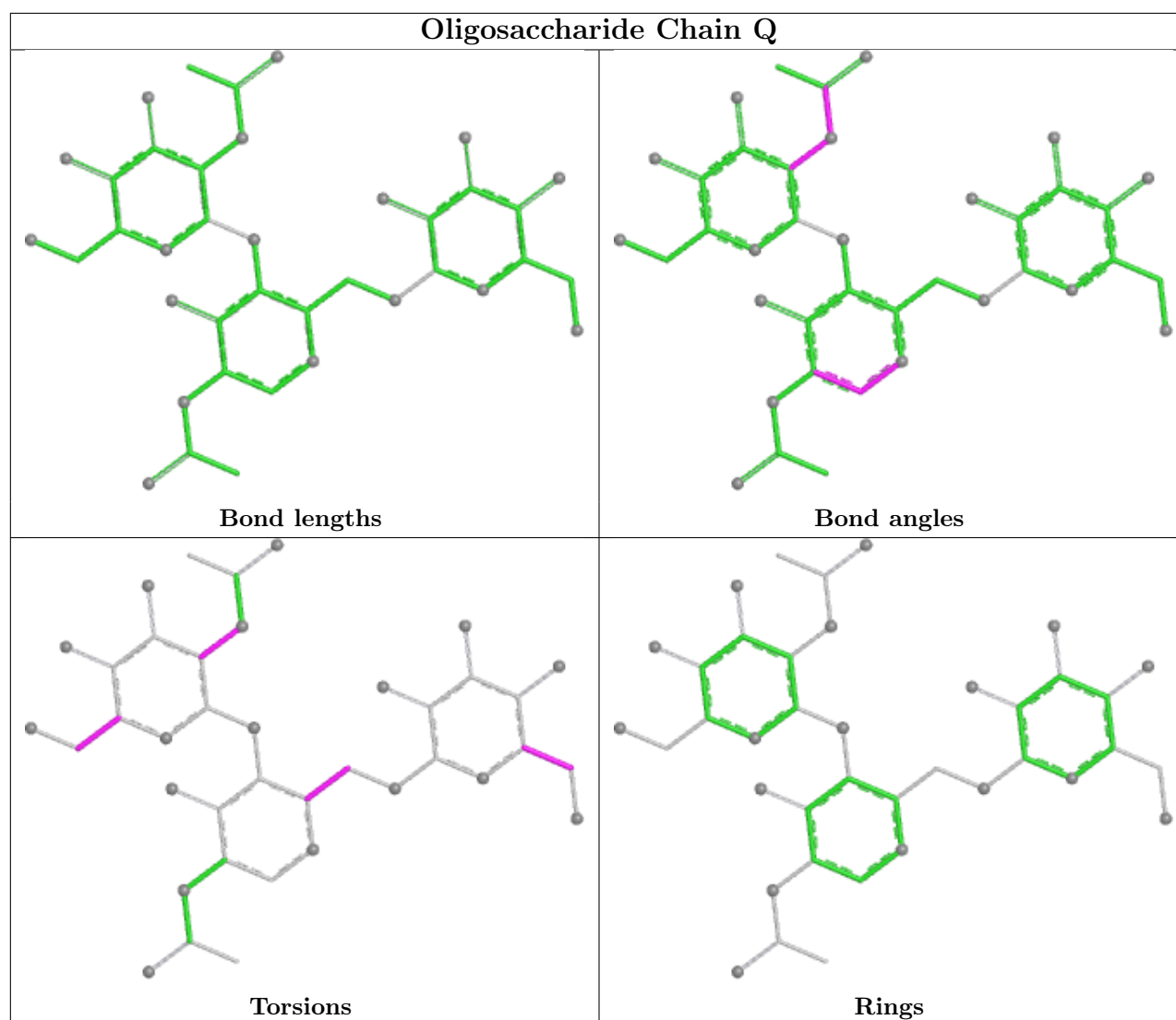


Rings









5.6 Ligand geometry [i](#)

17 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
10	SO4	B	809	-	4,4,4	0.70	0	6,6,6	0.17	0
7	NAG	B	805	1	14,14,15	0.65	0	17,19,21	1.06	1 (5%)
12	ACT	B	808	-	3,3,3	1.38	0	3,3,3	1.04	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	EDO	A	807	-	3,3,3	0.26	0	2,2,2	0.35	0
7	NAG	B	804	1	14,14,15	0.77	0	17,19,21	0.97	0
7	NAG	A	802	1	14,14,15	0.73	0	17,19,21	1.25	1 (5%)
7	NAG	B	803	1	14,14,15	0.70	0	17,19,21	1.15	1 (5%)
7	NAG	A	805	1	14,14,15	0.60	0	17,19,21	1.70	2 (11%)
8	1PE	A	806	-	15,15,15	0.29	0	14,14,14	0.20	0
7	NAG	A	804	1	14,14,15	0.64	0	17,19,21	1.02	1 (5%)
8	1PE	B	806	-	15,15,15	0.30	0	14,14,14	0.19	0
7	NAG	A	803	1	14,14,15	0.68	0	17,19,21	0.87	0
9	EDO	B	807	-	3,3,3	0.25	0	2,2,2	0.32	0
11	GLY	B	801	-	3,3,4	0.57	0	1,2,4	0.39	0
10	SO4	A	808	-	4,4,4	0.69	0	6,6,6	0.14	0
6	BLD	A	801	-	36,37,37	0.90	2 (5%)	49,59,59	1.01	3 (6%)
6	BLD	B	802	-	36,37,37	0.80	1 (2%)	49,59,59	0.94	1 (2%)

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	B	805	1	-	1/6/23/26	0/1/1/1
9	EDO	A	807	-	-	0/1/1/1	-
7	NAG	B	804	1	-	0/6/23/26	0/1/1/1
7	NAG	A	802	1	-	2/6/23/26	0/1/1/1
7	NAG	B	803	1	-	2/6/23/26	0/1/1/1
7	NAG	A	805	1	-	5/6/23/26	0/1/1/1
8	1PE	A	806	-	-	8/13/13/13	-
7	NAG	A	804	1	-	1/6/23/26	0/1/1/1
8	1PE	B	806	-	-	8/13/13/13	-
7	NAG	A	803	1	-	0/6/23/26	0/1/1/1
9	EDO	B	807	-	-	0/1/1/1	-
11	GLY	B	801	-	-	0/0/1/2	-
6	BLD	A	801	-	-	0/20/85/85	0/4/4/4
6	BLD	B	802	-	-	2/20/85/85	0/4/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	801	BLD	O07-C06	-2.70	1.31	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	802	BLD	O07-C07	-2.45	1.41	1.45
6	A	801	BLD	O07-C07	-2.44	1.41	1.45

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	805	NAG	C2-N2-C7	4.97	129.56	122.90
6	A	801	BLD	C11-C09-C08	-4.08	106.09	111.78
7	A	802	NAG	C2-N2-C7	3.59	127.71	122.90
7	A	805	NAG	C1-O5-C5	3.43	116.78	112.19
7	B	803	NAG	C2-N2-C7	3.09	127.04	122.90
6	B	802	BLD	O23-C23-C24	3.03	116.28	109.56
7	A	804	NAG	C1-O5-C5	2.71	115.81	112.19
7	B	805	NAG	C1-O5-C5	2.58	115.65	112.19
6	A	801	BLD	C04-C03-C02	-2.49	107.30	110.38
6	A	801	BLD	C10-C01-C02	2.31	117.97	114.17

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	802	BLD	C23-C24-C25-C26
7	A	805	NAG	O5-C5-C6-O6
7	A	805	NAG	C4-C5-C6-O6
7	A	802	NAG	C8-C7-N2-C2
7	A	802	NAG	O7-C7-N2-C2
7	A	805	NAG	C8-C7-N2-C2
7	A	805	NAG	O7-C7-N2-C2
7	B	803	NAG	C8-C7-N2-C2
7	B	803	NAG	O7-C7-N2-C2
8	A	806	1PE	OH2-C12-C22-OH3
8	B	806	1PE	OH5-C14-C24-OH4
8	B	806	1PE	OH4-C13-C23-OH3
7	A	804	NAG	O5-C5-C6-O6
8	A	806	1PE	OH4-C13-C23-OH3
8	B	806	1PE	OH7-C16-C26-OH6
7	B	805	NAG	O5-C5-C6-O6
8	A	806	1PE	OH5-C14-C24-OH4
8	B	806	1PE	OH6-C15-C25-OH5
8	A	806	1PE	C12-C22-OH3-C23
8	A	806	1PE	C15-C25-OH5-C14
8	A	806	1PE	C23-C13-OH4-C24

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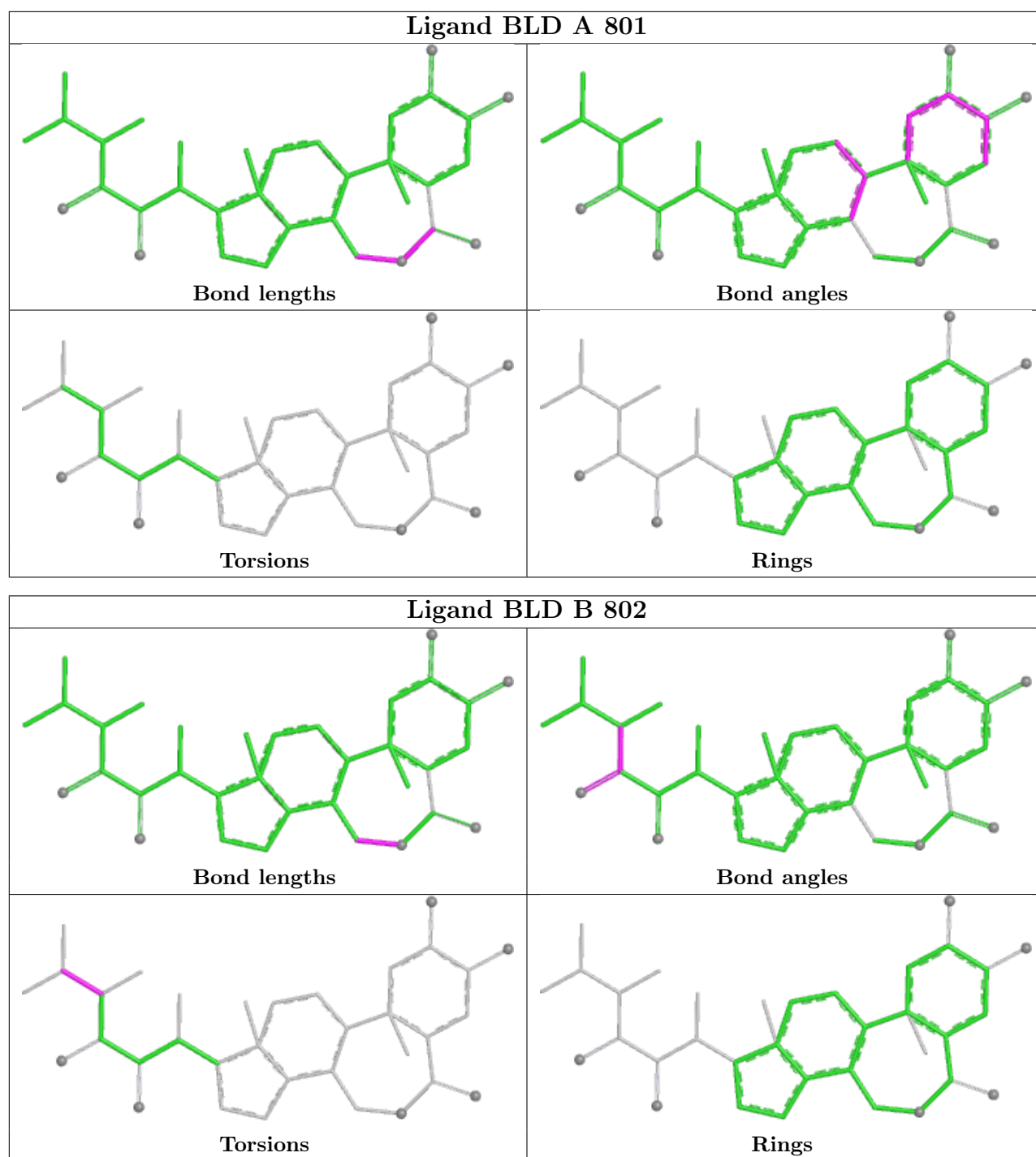
Mol	Chain	Res	Type	Atoms
8	B	806	1PE	C23-C13-OH4-C24
8	B	806	1PE	C13-C23-OH3-C22
8	A	806	1PE	C24-C14-OH5-C25
7	A	805	NAG	C1-C2-N2-C7
6	B	802	BLD	C28-C24-C25-C26
8	A	806	1PE	C25-C15-OH6-C26
8	B	806	1PE	C15-C25-OH5-C14
8	B	806	1PE	C24-C14-OH5-C25

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	B	808	ACT	2	0
7	A	802	NAG	1	0
7	B	803	NAG	3	0
7	A	805	NAG	1	0
8	A	806	1PE	1	0
6	B	802	BLD	1	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	721/770 (93%)	0.38	28 (3%)	43 43	38, 67, 130, 160	3 (0%)
1	B	727/770 (94%)	0.36	25 (3%)	48 48	29, 65, 124, 160	6 (0%)
All	All	1448/1540 (94%)	0.37	53 (3%)	45 45	29, 66, 127, 160	9 (0%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	53	LEU	4.0
1	B	618	VAL	4.0
1	A	617	MET	3.9
1	A	52	PHE	3.8
1	B	58	TYR	3.2
1	A	593	THR	3.1
1	A	45	ILE	2.9
1	B	617	MET	2.9
1	B	37	LEU	2.8
1	A	56	TRP	2.8
1	A	618	VAL	2.8
1	B	97	LEU	2.7
1	A	120	SER	2.7
1	B	619	HIS	2.7
1	B	591	GLY	2.7
1	B	589	ASN	2.6
1	B	395	PHE	2.6
1	B	398	LEU	2.6
1	B	31	VAL	2.6
1	B	38	THR	2.6
1	A	113	PHE	2.6
1	B	400	SER	2.6
1	A	66	THR	2.5
1	A	255[A]	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	469	SER	2.4
1	A	70	VAL	2.4
1	B	78	VAL	2.4
1	B	588	ARG	2.3
1	B	34	THR	2.3
1	A	173	SER	2.3
1	A	592	GLY	2.3
1	A	597	GLY	2.3
1	A	100	LEU	2.3
1	B	65	CYS	2.2
1	A	83	LEU	2.2
1	A	92	LEU	2.2
1	B	469	SER	2.2
1	B	598	ALA	2.2
1	A	238	ILE	2.1
1	A	628	SER	2.1
1	A	31	VAL	2.1
1	A	619	HIS	2.1
1	A	101	SER	2.1
1	A	616	PRO	2.1
1	B	85	ASN	2.1
1	B	68	ARG	2.1
1	B	593	THR	2.1
1	A	470	ILE	2.1
1	B	105	SER	2.1
1	A	43	THR	2.0
1	A	121	SER	2.0
1	B	622	PRO	2.0
1	B	67	TRP	2.0

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

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.

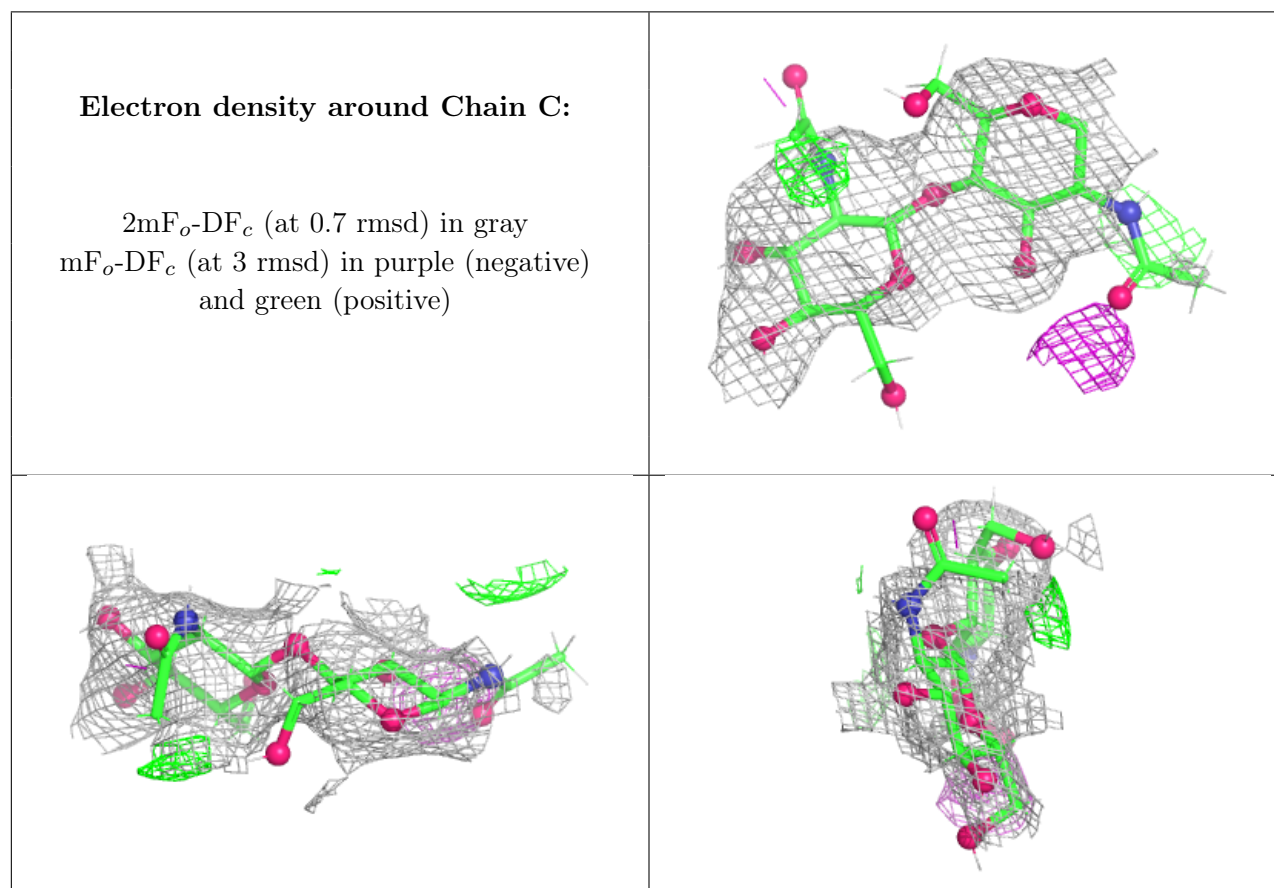
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	J	2	14/15	0.31	0.14	117,129,154,156	0
2	NAG	C	2	14/15	0.38	0.14	127,134,157,162	0
4	BMA	O	3	11/12	0.39	0.14	120,127,152,154	0
2	NAG	R	2	14/15	0.41	0.15	127,134,160,164	0
2	NAG	P	2	14/15	0.44	0.18	122,135,160,162	0
2	NAG	S	1	14/15	0.45	0.20	87,121,147,149	0
5	MAN	Q	3	11/12	0.45	0.17	120,129,153,154	0
2	NAG	C	1	14/15	0.50	0.17	113,129,153,157	0
4	BMA	G	3	11/12	0.52	0.14	99,116,139,142	0
2	NAG	S	2	14/15	0.53	0.17	110,134,161,164	0
2	NAG	L	2	14/15	0.55	0.12	113,123,148,148	0
5	MAN	I	3	11/12	0.57	0.13	106,118,142,147	0
3	MAN	F	6	11/12	0.59	0.22	87,96,113,115	0
3	MAN	N	5	11/12	0.59	0.12	94,106,125,128	0
5	NAG	Q	2	14/15	0.59	0.13	115,125,149,151	0
2	NAG	H	2	14/15	0.59	0.15	109,123,148,166	0
3	MAN	N	6	11/12	0.63	0.18	101,113,133,137	0
2	NAG	D	2	14/15	0.63	0.12	116,123,148,151	0
2	NAG	K	2	14/15	0.68	0.13	94,105,125,126	0
2	NAG	M	2	14/15	0.69	0.14	83,91,109,112	0
4	NAG	O	2	14/15	0.70	0.12	86,108,142,149	0
5	NAG	I	2	14/15	0.71	0.12	111,120,139,145	0
2	NAG	R	1	14/15	0.72	0.14	102,122,146,147	0
2	NAG	E	2	14/15	0.73	0.11	104,112,134,136	0
2	NAG	D	1	14/15	0.74	0.10	107,115,137,138	0
2	NAG	T	2	14/15	0.74	0.14	96,105,125,126	0
2	NAG	J	1	14/15	0.74	0.12	82,107,129,139	0
2	NAG	P	1	14/15	0.77	0.14	83,106,130,149	0
3	MAN	F	5	11/12	0.77	0.10	76,83,99,104	0
2	NAG	H	1	14/15	0.77	0.14	74,93,116,127	0
2	NAG	L	1	14/15	0.77	0.12	99,116,135,138	0
2	NAG	E	1	14/15	0.79	0.12	88,107,120,128	0
3	MAN	N	4	11/12	0.81	0.13	77,86,103,104	0
5	NAG	Q	1	14/15	0.82	0.14	99,119,139,143	0
4	NAG	G	2	14/15	0.82	0.12	92,105,125,129	0
5	NAG	I	1	14/15	0.82	0.14	84,101,124,131	0
2	NAG	T	1	14/15	0.84	0.12	63,83,105,105	0
3	MAN	F	11	11/12	0.85	0.10	54,67,77,82	0
2	NAG	K	1	14/15	0.85	0.12	62,88,109,119	0
3	NAG	F	2	14/15	0.86	0.12	46,57,68,73	0
3	BMA	F	3	11/12	0.88	0.11	51,61,73,82	0
3	MAN	F	10	11/12	0.88	0.11	50,59,71,74	0
3	MAN	F	4	11/12	0.89	0.08	64,74,89,90	0

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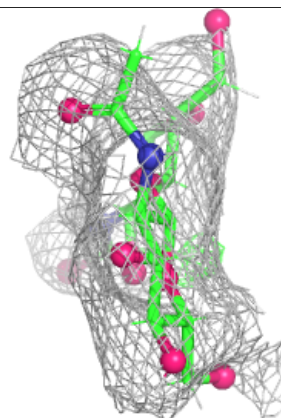
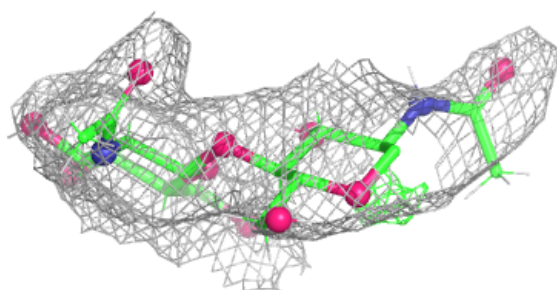
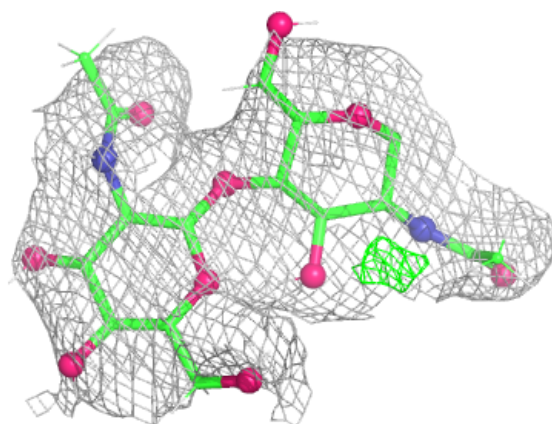
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	N	2	14/15	0.89	0.09	38,53,64,80	0
2	NAG	M	1	14/15	0.89	0.08	67,80,92,98	0
4	NAG	O	1	14/15	0.90	0.10	67,81,95,100	0
3	MAN	N	10	11/12	0.90	0.10	54,63,76,83	0
3	MAN	N	11	11/12	0.90	0.09	53,62,71,77	0
4	NAG	G	1	14/15	0.91	0.10	65,81,96,97	0
3	BMA	N	3	11/12	0.91	0.09	44,58,73,78	0
3	NAG	F	1	14/15	0.92	0.10	51,63,80,80	0
3	MAN	F	9	11/12	0.93	0.08	40,49,56,61	0
3	MAN	N	9	11/12	0.94	0.07	36,43,50,54	0
3	NAG	N	1	14/15	0.94	0.08	41,51,67,67	0
3	MAN	F	7	11/12	0.94	0.07	48,53,64,64	0
3	MAN	N	7	11/12	0.95	0.06	40,47,57,58	0
3	MAN	F	8	11/12	0.96	0.06	40,46,54,55	0
3	MAN	N	8	11/12	0.97	0.07	40,47,56,58	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.

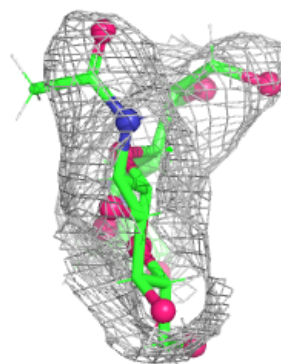
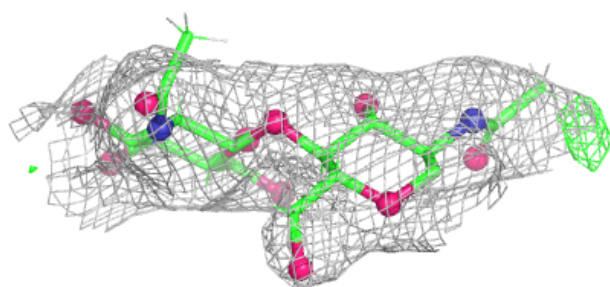
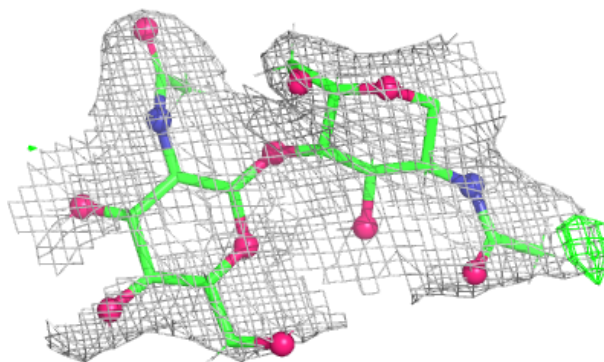


Electron density around Chain D:

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

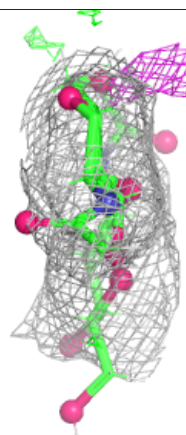
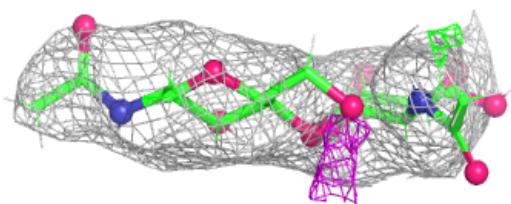
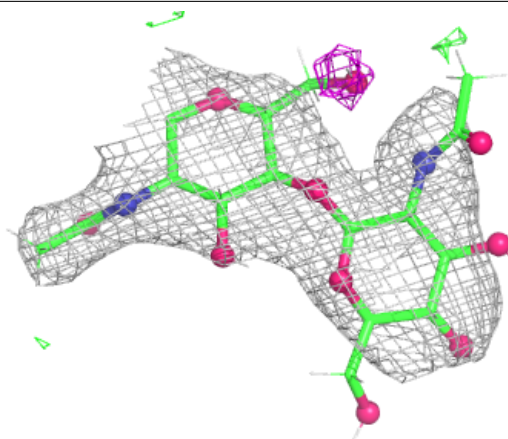
**Electron density around Chain E:**

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

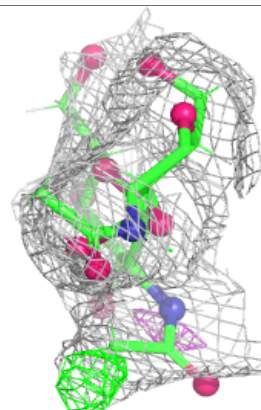
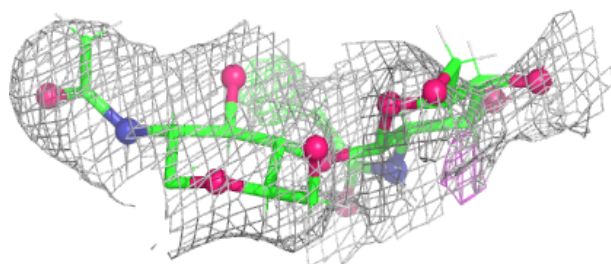
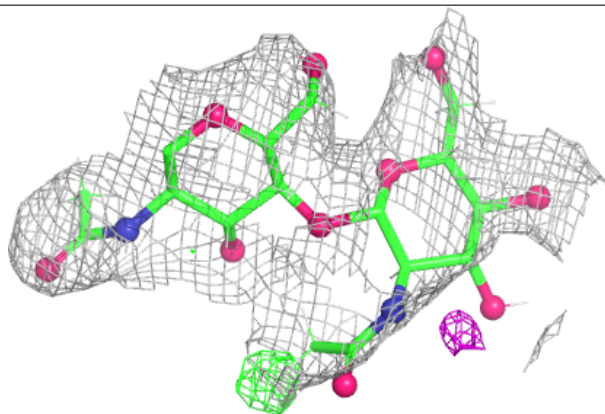


Electron density around Chain H:

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

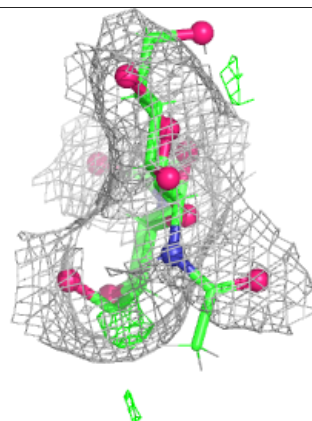
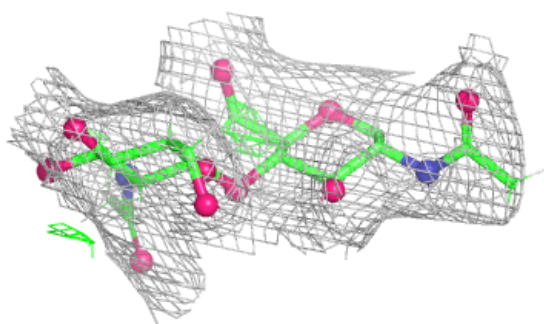
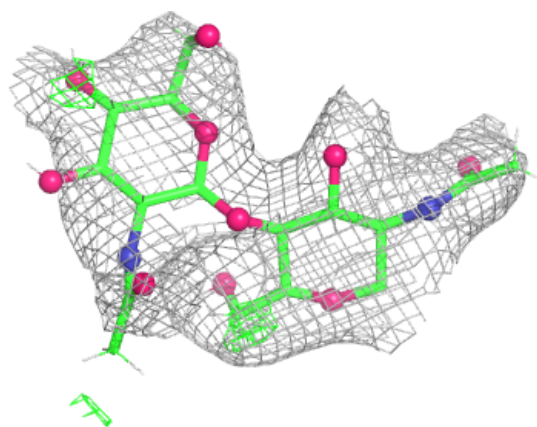
**Electron density around Chain J:**

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



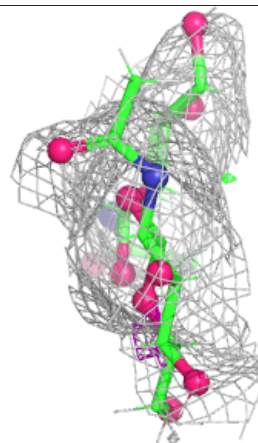
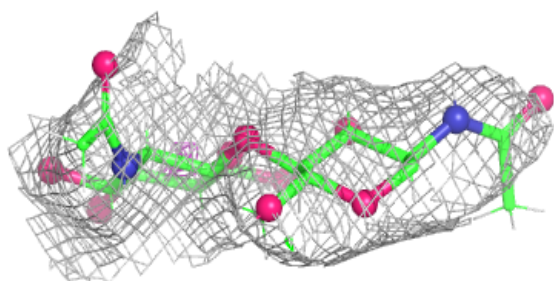
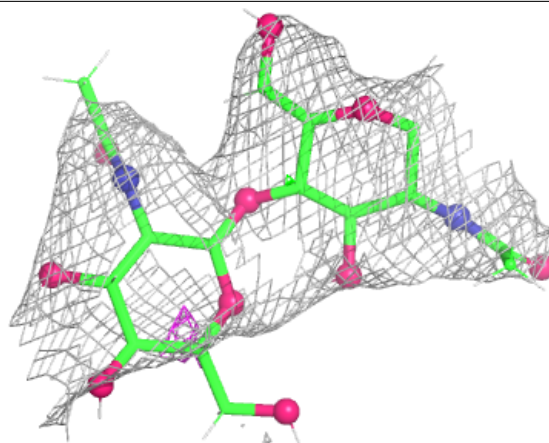
Electron density around Chain K:

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



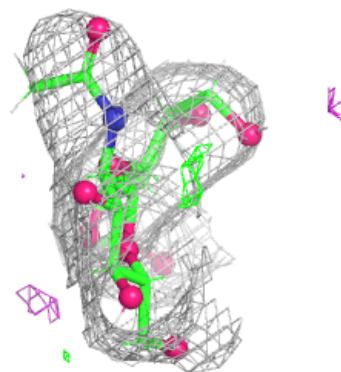
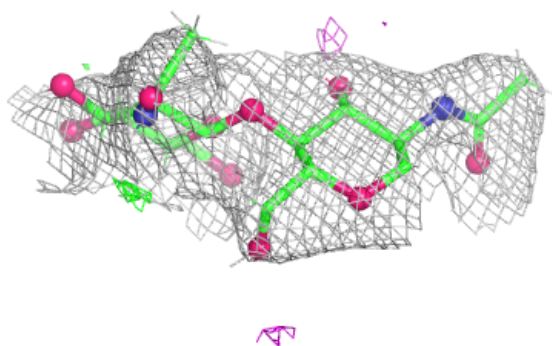
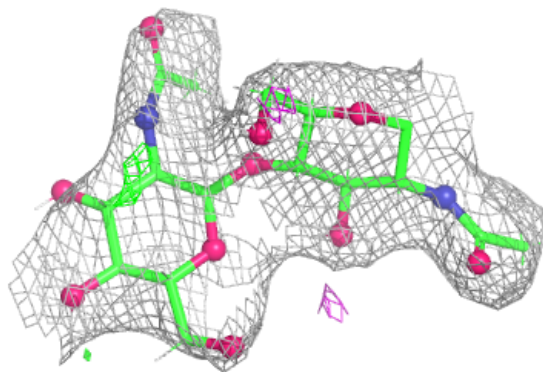
Electron density around Chain L:

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

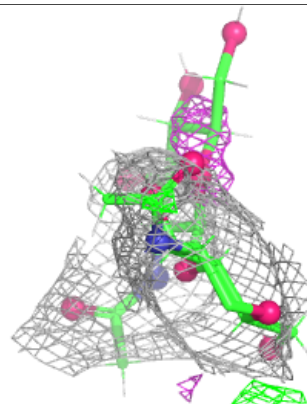
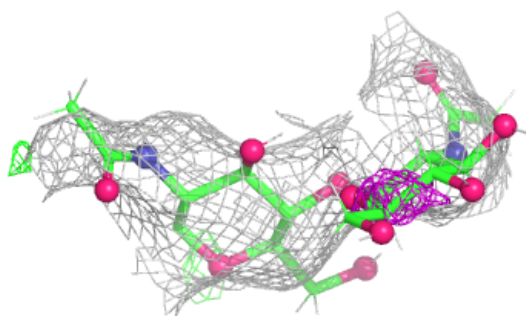
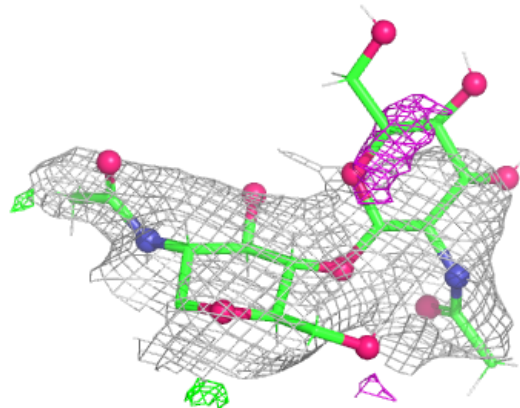


Electron density around Chain M:

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

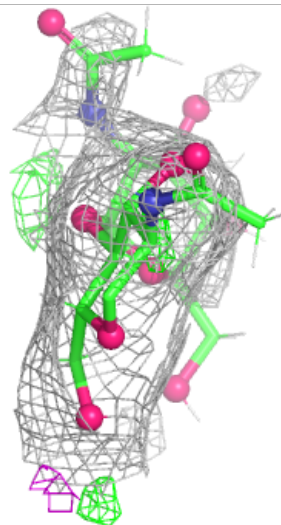
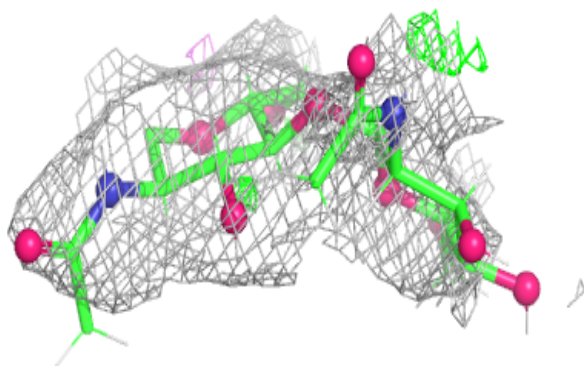
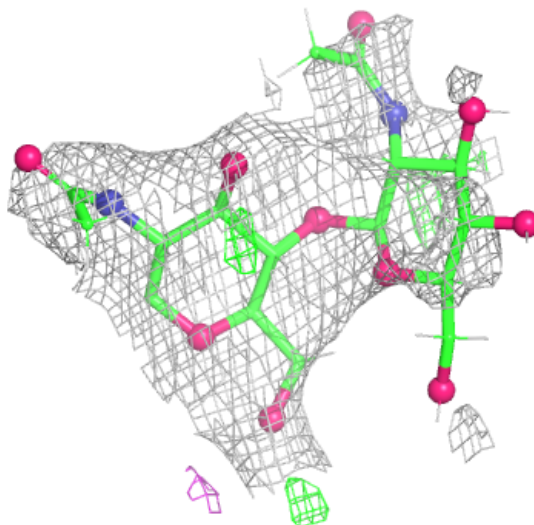
**Electron density around Chain P:**

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



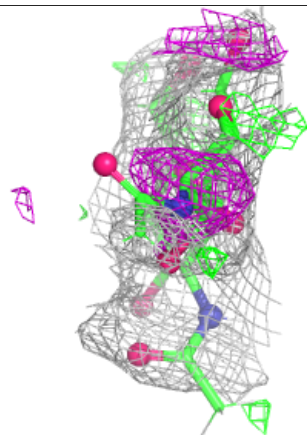
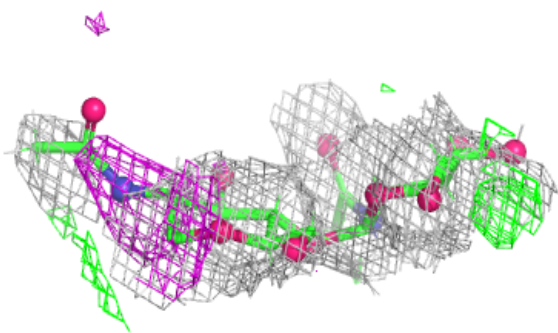
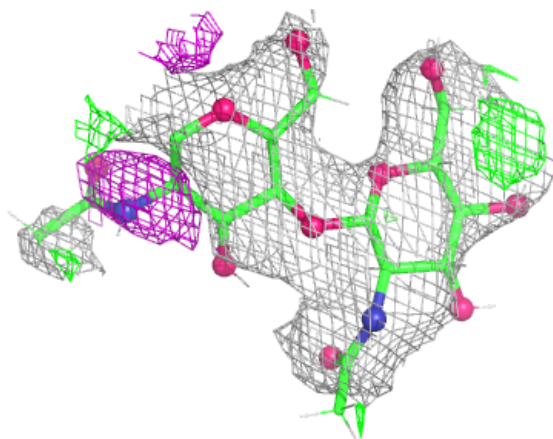
Electron density around Chain R:

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



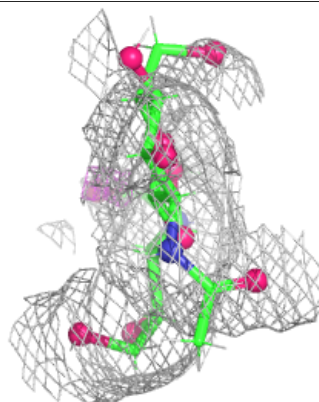
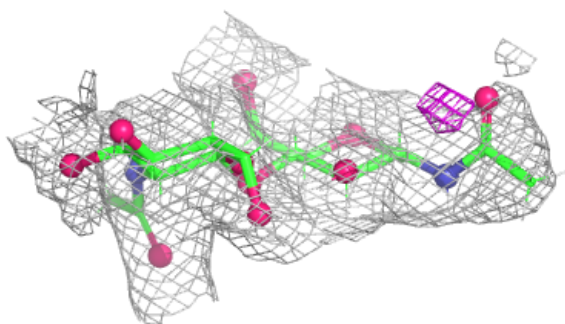
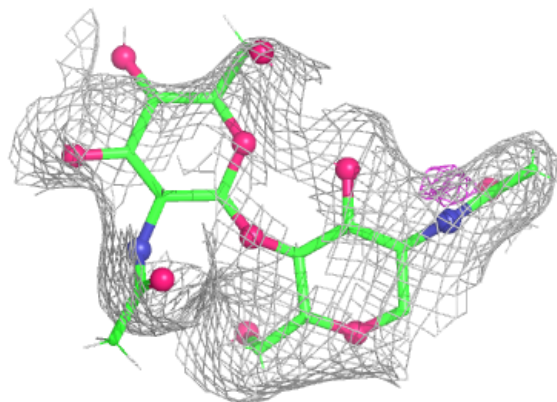
Electron density around Chain S:

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



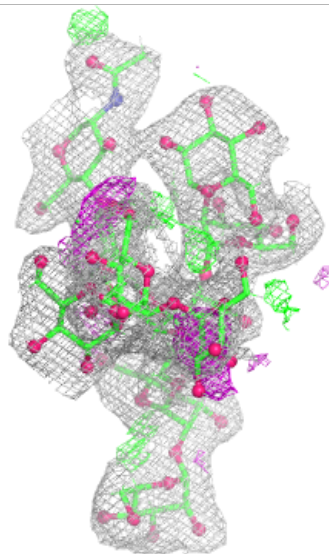
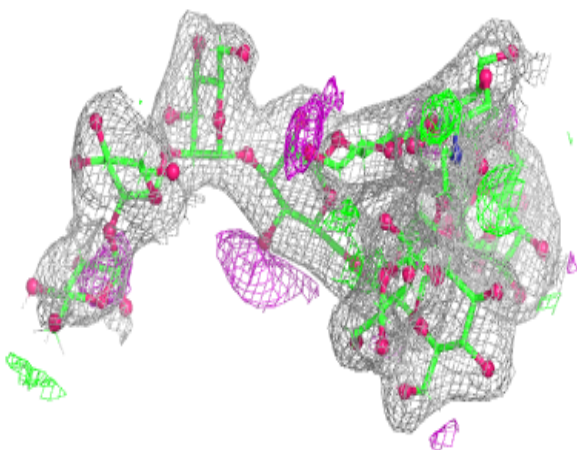
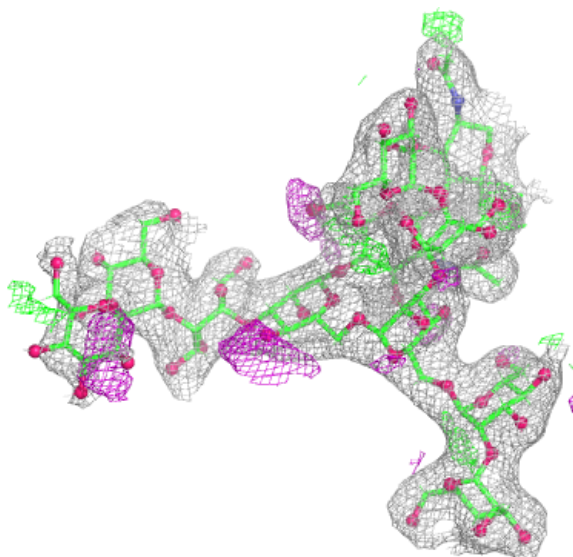
Electron density around Chain T:

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



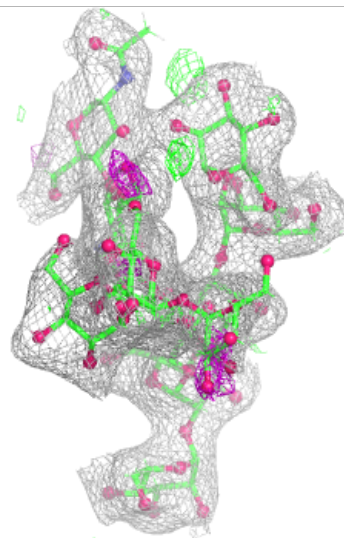
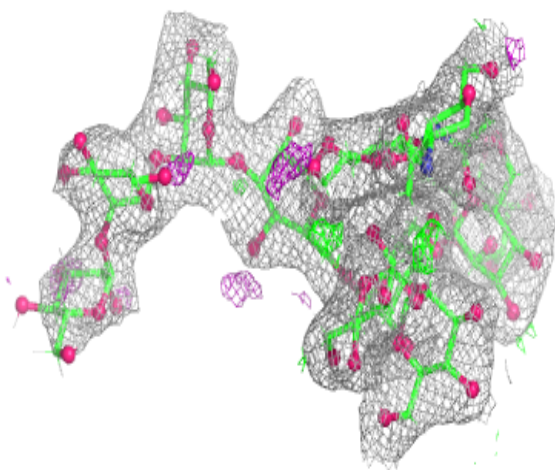
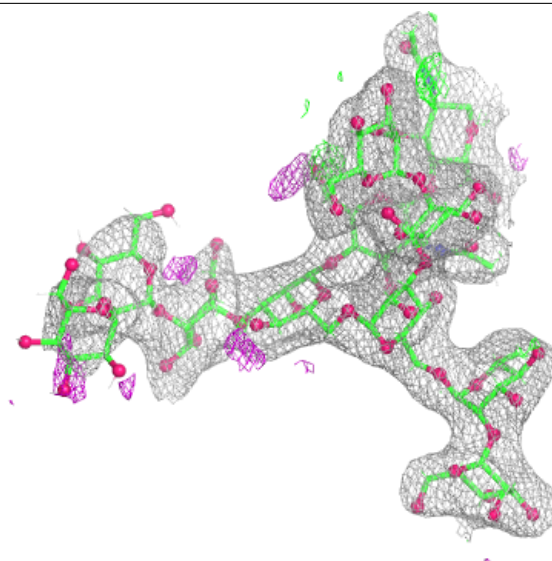
Electron density around Chain F:

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



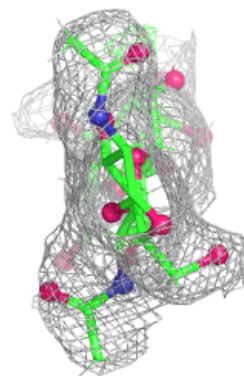
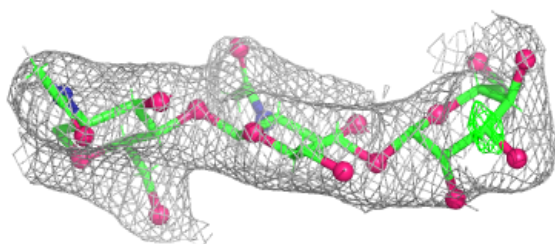
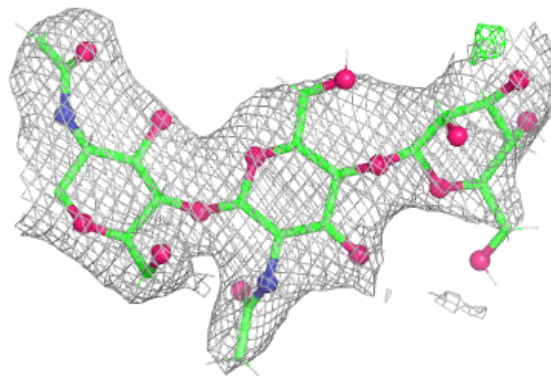
Electron density around Chain N:

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

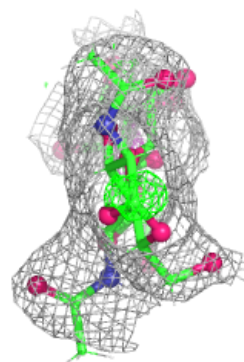
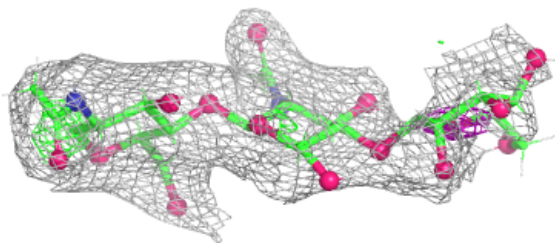
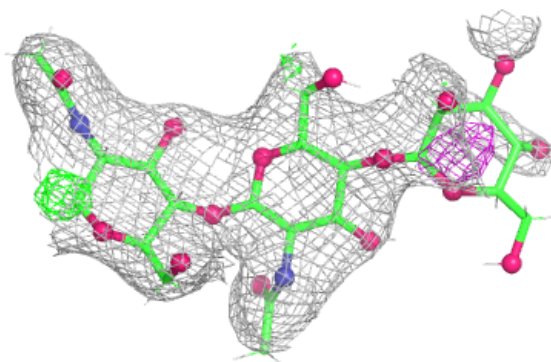


Electron density around Chain G:

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

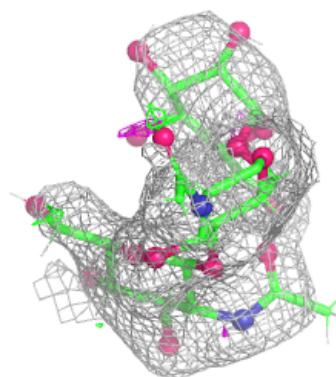
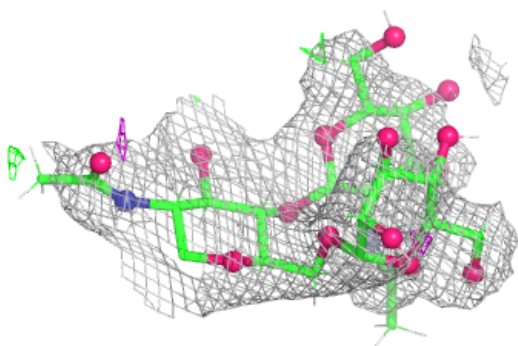
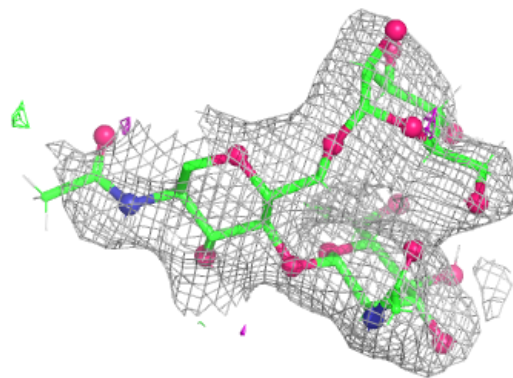
**Electron density around Chain O:**

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

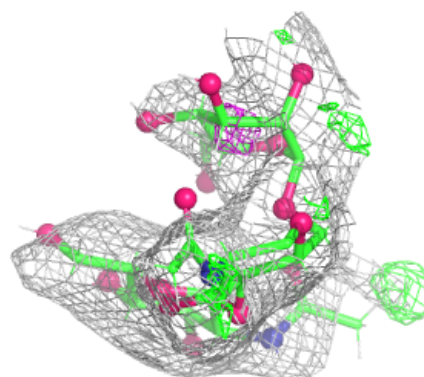
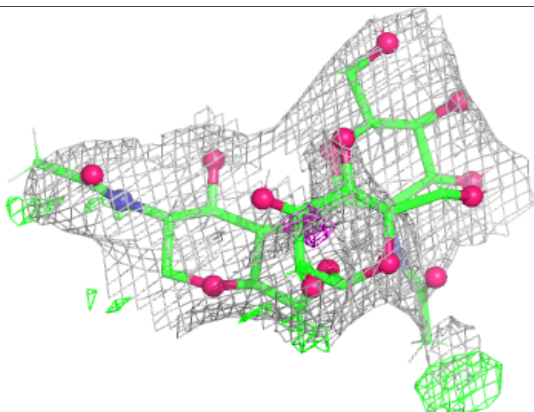
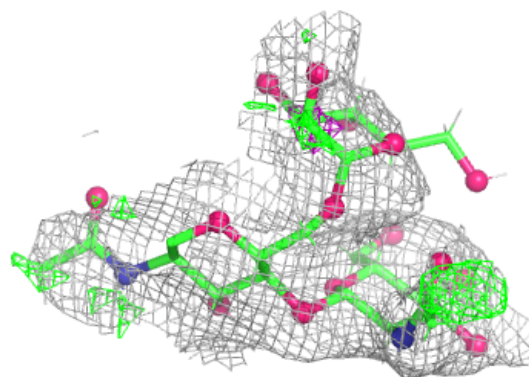


Electron density around Chain I:

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

**Electron density around Chain Q:**

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



6.4 Ligands ⓘ

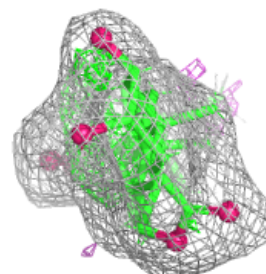
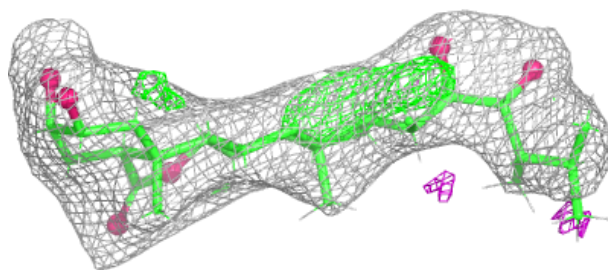
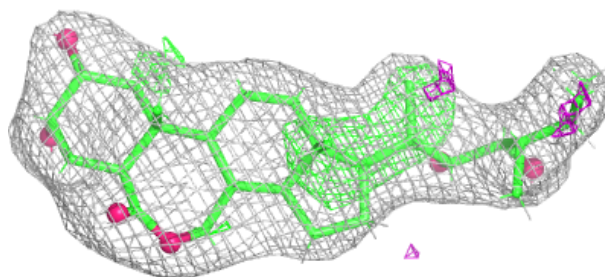
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	NAG	A	805	14/15	0.44	0.14	100,110,132,136	0
11	GLY	B	801	4/5	0.46	0.22	67,71,100,100	0
7	NAG	A	802	14/15	0.49	0.15	107,129,154,165	0
7	NAG	B	803	14/15	0.57	0.15	90,108,129,131	0
7	NAG	B	805	14/15	0.66	0.14	75,87,104,112	0
12	ACT	B	808	4/4	0.69	0.21	69,71,83,83	0
7	NAG	B	804	14/15	0.72	0.12	84,97,116,121	0
8	1PE	A	806	16/16	0.74	0.19	59,77,93,95	0
7	NAG	A	804	14/15	0.75	0.12	70,86,105,120	0
7	NAG	A	803	14/15	0.76	0.13	88,100,120,127	0
10	SO4	A	808	5/5	0.82	0.11	79,84,99,116	0
8	1PE	B	806	16/16	0.82	0.16	57,72,87,89	0
9	EDO	A	807	4/4	0.82	0.16	70,84,96,97	0
10	SO4	B	809	5/5	0.83	0.22	62,65,74,82	0
9	EDO	B	807	4/4	0.85	0.16	50,60,70,71	0
6	BLD	B	802	34/34	0.90	0.15	49,61,68,73	0
6	BLD	A	801	34/34	0.92	0.12	44,61,75,76	0

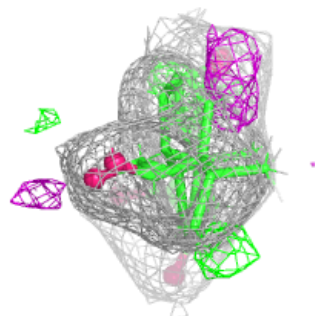
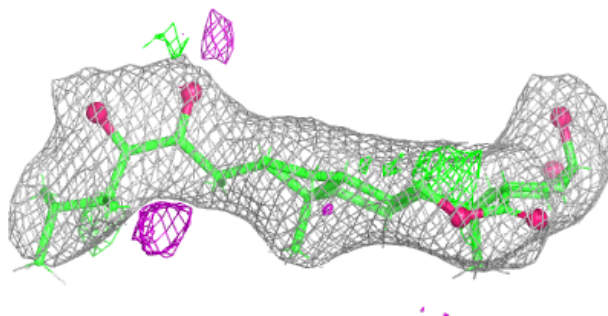
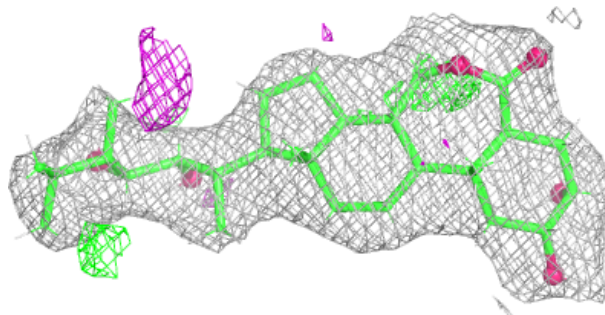
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around BLD B 802:

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

**Electron density around BLD A 801:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.