



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 26, 2025 – 04:28 pm BST

PDB ID : 9S9A / pdb\_00009s9a  
Title : Crystal structure of the BRL3 ectodomain from Arabidopsis thaliana in complex with 6-deoxocastasterone.  
Authors : Caregnato, A.; Hothorn, M.  
Deposited on : 2025-08-06  
Resolution : 3.20 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](mailto: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>.

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

MolProbity : 4-5-2 with Phenix2.0rc1  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.45.1

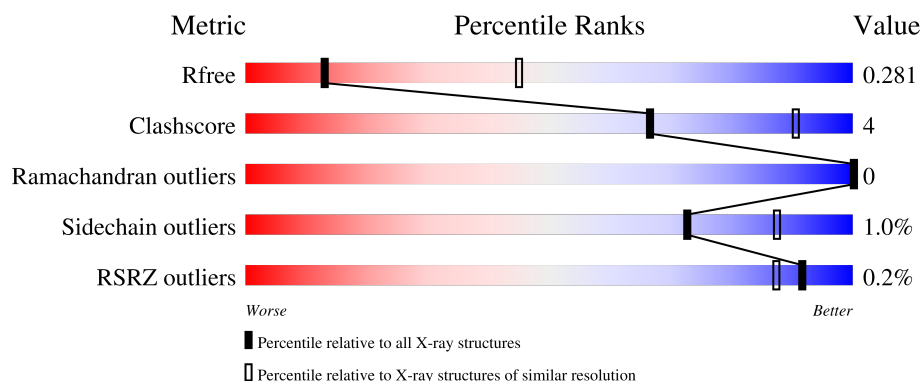
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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))
$R_{free}$	164625	1370 (3.20-3.20)
Clashscore	180529	1497 (3.20-3.20)
Ramachandran outliers	177936	1479 (3.20-3.20)
Sidechain outliers	177891	1478 (3.20-3.20)
RSRZ outliers	164620	1371 (3.20-3.20)

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  $\geq 3$ , 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	779	
1	B	779	
2	C	2	
2	E	2	
2	F	2	

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Mol	Chain	Length	Quality of chain
2	G	2	 50% 50%
2	H	2	 100%
2	I	2	 100%
2	J	2	 100%
2	M	2	 50% 50%
2	N	2	 50% 50%
2	O	2	 100%
2	P	2	 50% 50%
3	D	10	 100%
4	K	11	 18% 82%
5	L	3	 33% 67%

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 23142 atoms, of which 11502 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	1	0
			10753	3390	5355	908	1073	27			
1	B	727	Total	C	H	N	O	S	0	7	0
			10908	3436	5436	924	1083	29			

There are 18 discrepancies between the modelled and reference sequences:

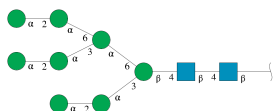
Chain	Residue	Modelled	Actual	Comment	Reference
A	771	ALA	-	expression tag	UNP Q9LJF3
A	772	ALA	-	expression tag	UNP Q9LJF3
A	773	ALA	-	expression tag	UNP Q9LJF3
A	774	GLU	-	expression tag	UNP Q9LJF3
A	775	ASN	-	expression tag	UNP Q9LJF3
A	776	LEU	-	expression tag	UNP Q9LJF3
A	777	TYR	-	expression tag	UNP Q9LJF3
A	778	PHE	-	expression tag	UNP Q9LJF3
A	779	GLN	-	expression tag	UNP Q9LJF3
B	771	ALA	-	expression tag	UNP Q9LJF3
B	772	ALA	-	expression tag	UNP Q9LJF3
B	773	ALA	-	expression tag	UNP Q9LJF3
B	774	GLU	-	expression tag	UNP Q9LJF3
B	775	ASN	-	expression tag	UNP Q9LJF3
B	776	LEU	-	expression tag	UNP Q9LJF3
B	777	TYR	-	expression tag	UNP Q9LJF3
B	778	PHE	-	expression tag	UNP Q9LJF3
B	779	GLN	-	expression tag	UNP Q9LJF3

- 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	E	2	Total	C	H	N	O	0	0	0
			52	16	24	2	10			
2	F	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	G	2	Total	C	H	N	O	0	0	0
			52	16	24	2	10			
2	H	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	I	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	M	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	N	2	Total	C	H	N	O	0	0	0
			52	16	24	2	10			
2	O	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			

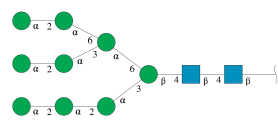
- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	10	Total	C	H	N	O	0	0	0
			212	64	96	2	50			

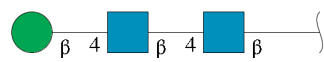
- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyra

nose-(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.



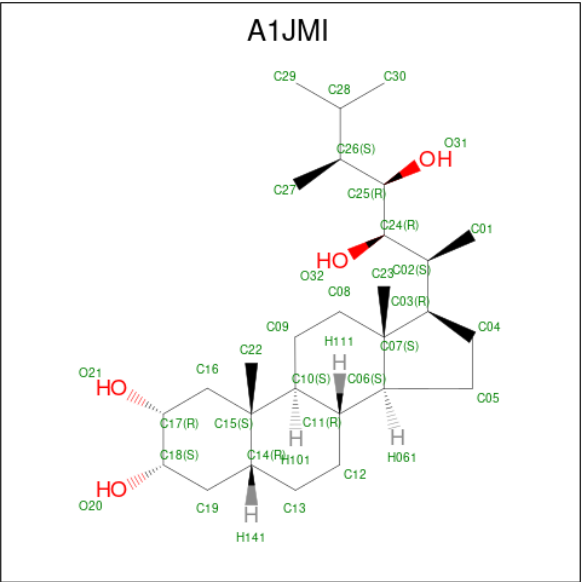
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	K	11	Total	C	H	N	O	0	0	0
			233	70	106	2	55			

- Molecule 5 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
5	L	3	Total	C	H	N	O	0	0	0
			73	22	34	2	15			

- Molecule 6 is 6-deoxocastasterone (CCD ID: A1JMI) (formula: C<sub>28</sub>H<sub>50</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			82	28	50	4		

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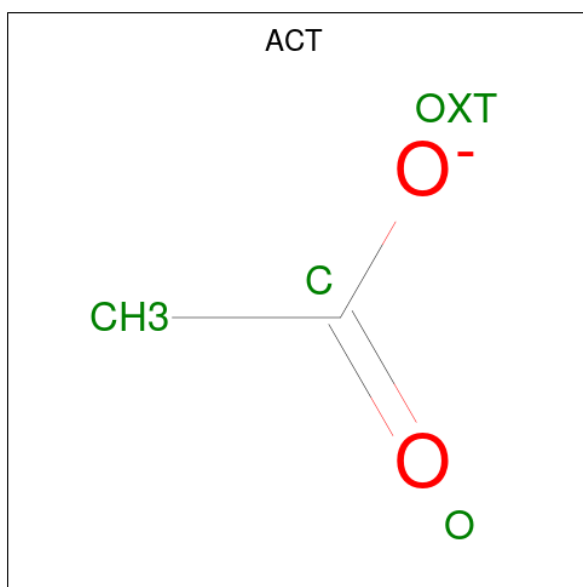
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	H	O	0	0
			82	28	50	4		

- 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 26	C 8	H 12	N 1	O 5	0	0
7	A	1	Total 26	C 8	H 12	N 1	O 5	0	0
7	A	1	Total 27	C 8	H 13	N 1	O 5	0	0
7	A	1	Total 27	C 8	H 13	N 1	O 5	0	0
7	B	1	Total 26	C 8	H 12	N 1	O 5	0	0
7	B	1	Total 27	C 8	H 13	N 1	O 5	0	0
7	B	1	Total 27	C 8	H 13	N 1	O 5	0	0
7	B	1	Total 26	C 8	H 12	N 1	O 5	0	0

- Molecule 8 is ACETATE ION (CCD ID: ACT) (formula:  $C_2H_3O_2$ ).



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



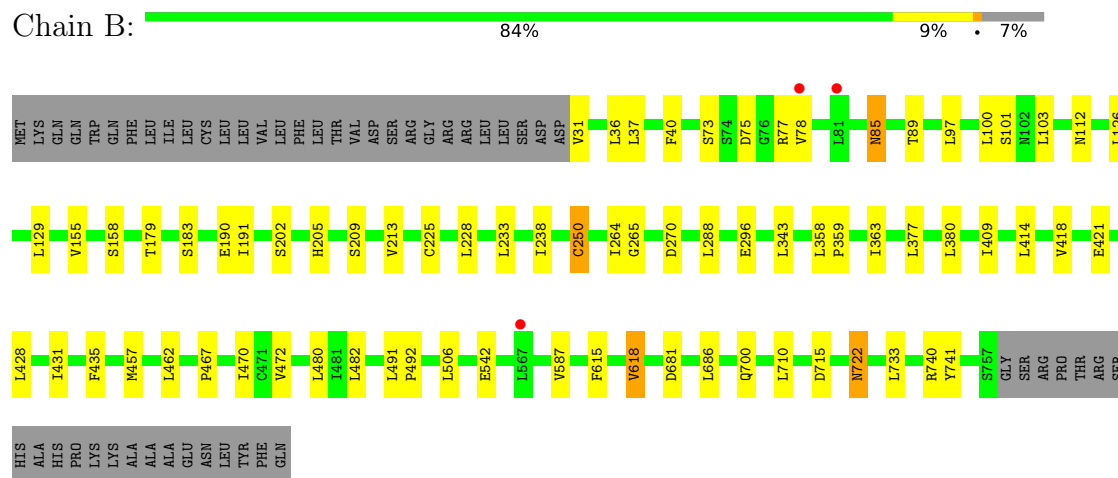
### 3 Residue-property plots

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



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



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

Chain E:  50% 50%

MAG1  
MAG2

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

Chain F:  50% 50%

MAG1  
MAG2

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

Chain G:  50% 50%

MAG1  
MAG2

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

Chain H:  100%

MAG1  
MAG2

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

Chain I:  100%

MAG1  
MAG2

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

Chain J:  100%

MAG1  
MAG2

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

Chain M:  50% 50%



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

Chain N: 50% 50%



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

Chain O: 100%



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

Chain P: 50% 50%



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

Chain D: 100%



- Molecule 4: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)-[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 K: 18% 82%



- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L: 33% 67%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.16Å 82.69Å 122.20Å 107.02° 91.35° 112.46°	Depositor
Resolution (Å)	48.11 – 3.20 48.11 – 3.20	Depositor EDS
% Data completeness (in resolution range)	89.2 (48.11-3.20) 76.1 (48.11-3.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.38 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, $R_{free}$	0.234 , 0.281 0.236 , 0.281	Depositor DCC
$R_{free}$ test set	1585 reflections (4.52%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	87.9	Xtriage
Anisotropy	0.508	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 93.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.031 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	23142	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	144.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.37% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, A1JMI, BMA, ACT, MAN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.09	0/5495	0.30	0/7466
1	B	0.11	0/5588	0.33	0/7592
All	All	0.10	0/11083	0.32	0/15058

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	5398	5355	5357	32	1
1	B	5472	5436	5433	48	1
2	C	28	25	25	0	0
2	E	28	24	25	0	0
2	F	28	25	25	1	0
2	G	28	24	25	0	0
2	H	28	25	25	2	0
2	I	28	25	25	0	0
2	J	28	25	25	0	0
2	M	28	25	25	0	0
2	N	28	24	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	O	28	25	25	1	0
2	P	28	25	25	0	0
3	D	116	96	97	1	0
4	K	127	106	106	0	0
5	L	39	34	34	0	0
6	A	32	50	0	1	0
6	B	32	50	0	0	0
7	A	56	50	52	0	0
7	B	56	50	52	1	0
8	B	4	3	3	0	0
All	All	11640	11502	11409	83	1

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 4.

All (83) 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	1.98	0.81
1:A:343:LEU:HD21	1:A:358:LEU:HD13	1.64	0.79
1:B:733:LEU:O	1:B:741:TYR:OH	2.05	0.75
2:O:1:NAG:O3	2:O:2:NAG:H83	1.89	0.71
1:B:715:ASP:OD1	1:B:740:ARG:NH2	2.24	0.70
1:A:130:ASP:OD1	1:A:132:SER:OG	2.11	0.67
1:B:421:GLU:N	1:B:421:GLU:OE1	2.27	0.67
1:B:77:ARG:NH1	1:B:101:SER:OG	2.28	0.66
1:B:40:PHE:CD2	1:B:97:LEU:HD21	2.32	0.65
1:A:152:LEU:HD21	1:A:155:VAL:HB	1.79	0.64
1:B:418:VAL:HG21	1:B:457:MET:HE1	1.80	0.64
1:A:201:ASN:ND2	1:A:226:GLU:OE1	2.33	0.61
1:A:34:THR:HG22	1:A:64:PRO:HB3	1.81	0.61
1:B:37:LEU:HD11	1:B:78:VAL:HG11	1.82	0.61
1:B:202:SER:HB2	7:B:803:NAG:H82	1.84	0.60
1:B:31:VAL:O	1:B:31:VAL:HG13	2.02	0.58
1:A:131:LEU:O	1:A:160:ASN:ND2	2.39	0.55
1:A:678:THR:HG22	1:A:679:ILE:N	2.21	0.55
1:B:428:LEU:HD13	1:B:431:ILE:HD11	1.88	0.55
1:B:233:LEU:HD13	1:B:238:ILE:HD13	1.88	0.54
1:B:233:LEU:HD13	1:B:238:ILE:CD1	2.37	0.54
1:A:343:LEU:HD21	1:A:358:LEU:CD1	2.36	0.54
1:B:103:LEU:HG	1:B:126:LEU:HD21	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:ILE:HG22	1:B:265:GLY:N	2.24	0.53
1:A:631:THR:HG22	6:A:801:A1JMI:C27	2.38	0.53
1:B:73:SER:OG	1:B:75:ASP:OD1	2.21	0.53
1:B:343:LEU:HD11	1:B:358:LEU:CD1	2.39	0.53
1:A:132:SER:O	1:A:158:SER:O	2.27	0.53
1:B:343:LEU:HD11	1:B:358:LEU:HD13	1.91	0.53
1:A:37:LEU:HD21	1:A:100:LEU:HD11	1.92	0.52
1:A:612:LEU:O	1:A:618:VAL:HG21	2.10	0.51
1:B:686:LEU:O	1:B:710:LEU:HD22	2.10	0.51
1:B:270:ASP:O	1:B:296:GLU:HB3	2.10	0.51
1:B:491[B]:LEU:HD21	1:B:506:LEU:HD13	1.92	0.51
1:A:686:LEU:O	1:A:710:LEU:HD22	2.12	0.50
1:B:179:THR:HG22	1:B:205[B]:HIS:CD2	2.49	0.48
1:A:409:ILE:HG21	1:A:414:LEU:HD11	1.94	0.48
1:B:491[B]:LEU:HD21	1:B:506:LEU:CD1	2.44	0.47
1:B:358:LEU:N	1:B:359:PRO:CD	2.78	0.47
1:B:75:ASP:OD1	1:B:75:ASP:N	2.48	0.47
1:B:191:ILE:HD11	1:B:213:VAL:CG2	2.45	0.46
1:B:183:SER:HA	1:B:209:SER:O	2.15	0.46
1:B:89:THR:HG23	1:B:112:ASN:ND2	2.31	0.46
1:B:409:ILE:HG21	1:B:414:LEU:HD11	1.97	0.46
1:B:480:LEU:HD11	1:B:482:LEU:CD1	2.45	0.46
1:A:358:LEU:N	1:A:359:PRO:CD	2.78	0.46
1:B:228:LEU:HB3	1:B:250[A]:CYS:SG	2.55	0.46
1:B:700:GLN:HG2	1:B:722:ASN:HB3	1.98	0.46
1:A:550:ILE:HD12	1:A:642:ILE:HD12	1.98	0.45
1:B:158:SER:HA	1:B:183:SER:O	2.17	0.45
1:A:333:LEU:HB2	1:A:358:LEU:HD23	1.97	0.45
1:B:85:ASN:N	1:B:85:ASN:OD1	2.50	0.45
1:A:260:ARG:HD2	3:D:2:NAG:H81	1.99	0.44
1:A:472:VAL:HG13	1:A:473:ASP:N	2.32	0.44
1:A:542:GLU:OE1	1:A:542:GLU:N	2.45	0.44
1:A:70:VAL:HG22	1:A:81:LEU:HD13	1.99	0.44
1:B:343:LEU:HD21	1:B:363:ILE:HD13	2.00	0.44
1:A:615:PHE:HD2	1:A:617:MET:HE3	1.82	0.44
1:B:457:MET:HE2	1:B:462:LEU:HD11	2.00	0.44
1:B:40:PHE:CZ	1:B:97:LEU:HD11	2.53	0.43
1:A:75:ASP:OD1	1:A:75:ASP:N	2.51	0.43
1:B:470:ILE:O	1:B:472:VAL:HG23	2.19	0.43
1:B:467:PRO:O	1:B:492:PRO:HG2	2.19	0.43
1:B:615:PHE:O	1:B:618:VAL:HG12	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:LEU:HD23	1:B:155:VAL:HG23	2.01	0.42
1:B:377:LEU:HD21	1:B:380:LEU:HB2	2.02	0.42
1:B:467:PRO:O	1:B:470:ILE:HD11	2.20	0.42
1:B:31:VAL:HG12	1:B:36:LEU:HD11	2.02	0.41
1:A:617:MET:C	1:A:617:MET:SD	3.03	0.41
1:A:150:LEU:HD22	1:A:173:SER:OG	2.19	0.41
1:B:233:LEU:HD22	1:B:238:ILE:HD11	2.03	0.41
2:F:1:NAG:O6	2:F:2:NAG:C7	2.68	0.41
1:A:678:THR:CG2	1:A:679:ILE:N	2.82	0.41
1:B:100:LEU:HD12	1:B:103:LEU:HD22	2.01	0.41
1:B:542:GLU:OE1	1:B:542:GLU:N	2.47	0.41
1:A:37:LEU:CD2	1:A:100:LEU:HD11	2.51	0.41
1:A:423:GLY:HA2	1:A:449:LEU:HD21	2.02	0.41
1:A:572:GLY:O	2:H:1:NAG:H83	2.21	0.41
2:H:1:NAG:O3	2:H:2:NAG:H83	2.20	0.41
1:B:587:VAL:HG11	1:B:618:VAL:HG23	2.02	0.40
1:A:727:ILE:O	1:A:728:PRO:C	2.64	0.40
1:A:618:VAL:O	1:A:624:THR:HG21	2.21	0.40
1:B:264:ILE:HG12	1:B:288:LEU:HD12	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:756:SER:OG	1:B:681:ASP:OD2[1_556]	2.19	0.01

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	718/779 (92%)	689 (96%)	29 (4%)	0	<b>100</b> <b>100</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	732/779 (94%)	702 (96%)	30 (4%)	0	100	100
All	All	1450/1558 (93%)	1391 (96%)	59 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	630/679 (93%)	624 (99%)	6 (1%)	73	87
1	B	639/679 (94%)	631 (99%)	8 (1%)	65	83
All	All	1269/1358 (93%)	1255 (99%)	14 (1%)	73	86

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	CYS
1	A	277	GLN
1	A	325	CYS
1	A	353	ILE
1	A	435	PHE
1	A	636	SER
1	B	85	ASN
1	B	190	GLU
1	B	225	CYS
1	B	250[A]	CYS
1	B	250[B]	CYS
1	B	435	PHE
1	B	618	VAL
1	B	722	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	95	ASN
1	B	134	ASN
1	B	589	ASN

### 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 ⓘ

46 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	0.70	0	17,19,21	1.10	1 (5%)
2	NAG	C	2	2	14,14,15	0.66	0	17,19,21	0.87	0
3	NAG	D	1	1,3	14,14,15	0.69	0	17,19,21	1.13	1 (5%)
3	MAN	D	10	3	11,11,12	0.85	1 (9%)	15,15,17	0.96	0
3	NAG	D	2	3	14,14,15	0.72	0	17,19,21	0.84	0
3	BMA	D	3	3	11,11,12	0.85	0	15,15,17	2.28	5 (33%)
3	MAN	D	4	3	11,11,12	0.80	1 (9%)	15,15,17	1.36	1 (6%)
3	MAN	D	5	3	11,11,12	0.86	1 (9%)	15,15,17	1.07	1 (6%)
3	MAN	D	6	3	11,11,12	0.89	1 (9%)	15,15,17	0.82	0
3	MAN	D	7	3	11,11,12	0.70	0	15,15,17	1.13	1 (6%)
3	MAN	D	8	3	11,11,12	0.80	1 (9%)	15,15,17	1.18	1 (6%)
3	MAN	D	9	3	11,11,12	0.75	0	15,15,17	1.07	1 (6%)
2	NAG	E	1	1,2	14,14,15	0.68	0	17,19,21	1.02	1 (5%)
2	NAG	E	2	2	14,14,15	0.72	0	17,19,21	0.84	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	F	1	1,2	14,14,15	0.79	0	17,19,21	0.84	0
2	NAG	F	2	2	14,14,15	0.70	0	17,19,21	1.10	1 (5%)
2	NAG	G	1	1,2	14,14,15	0.63	0	17,19,21	1.01	1 (5%)
2	NAG	G	2	2	14,14,15	0.70	0	17,19,21	0.83	0
2	NAG	H	1	1,2	14,14,15	0.62	0	17,19,21	1.53	2 (11%)
2	NAG	H	2	2	14,14,15	0.66	0	17,19,21	1.23	2 (11%)
2	NAG	I	1	1,2	14,14,15	0.72	0	17,19,21	0.89	0
2	NAG	I	2	2	14,14,15	0.70	0	17,19,21	0.95	0
2	NAG	J	1	1,2	14,14,15	0.69	0	17,19,21	1.00	1 (5%)
2	NAG	J	2	2	14,14,15	0.65	0	17,19,21	1.19	1 (5%)
4	NAG	K	1	1,4	14,14,15	0.69	0	17,19,21	0.94	1 (5%)
4	MAN	K	10	4	11,11,12	0.76	0	15,15,17	1.18	1 (6%)
4	MAN	K	11	4	11,11,12	0.80	0	15,15,17	0.94	1 (6%)
4	NAG	K	2	4	14,14,15	0.74	0	17,19,21	0.90	0
4	BMA	K	3	4	11,11,12	0.88	0	15,15,17	2.26	5 (33%)
4	MAN	K	4	4	11,11,12	0.77	0	15,15,17	1.03	1 (6%)
4	MAN	K	5	4	11,11,12	0.85	1 (9%)	15,15,17	0.86	0
4	MAN	K	6	4	11,11,12	0.77	0	15,15,17	0.98	0
4	MAN	K	7	4	11,11,12	0.95	1 (9%)	15,15,17	0.94	0
4	MAN	K	8	4	11,11,12	0.80	1 (9%)	15,15,17	0.98	1 (6%)
4	MAN	K	9	4	11,11,12	0.82	1 (9%)	15,15,17	0.92	0
5	NAG	L	1	1,5	14,14,15	0.73	0	17,19,21	0.93	1 (5%)
5	NAG	L	2	5	14,14,15	0.70	0	17,19,21	0.86	0
5	BMA	L	3	5	11,11,12	0.78	0	15,15,17	2.56	6 (40%)
2	NAG	M	1	1,2	14,14,15	0.65	0	17,19,21	1.18	1 (5%)
2	NAG	M	2	2	14,14,15	0.71	0	17,19,21	0.83	0
2	NAG	N	1	1,2	14,14,15	0.71	0	17,19,21	0.66	0
2	NAG	N	2	2	14,14,15	0.72	0	17,19,21	1.15	1 (5%)
2	NAG	O	1	1,2	14,14,15	0.68	0	17,19,21	1.19	2 (11%)
2	NAG	O	2	2	14,14,15	0.73	0	17,19,21	1.65	2 (11%)
2	NAG	P	1	1,2	14,14,15	0.67	0	17,19,21	0.90	0
2	NAG	P	2	2	14,14,15	0.69	0	17,19,21	0.92	1 (5%)

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

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	N	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	N	2	2	-	2/6/23/26	0/1/1/1
2	NAG	O	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	O	2	2	-	2/6/23/26	0/1/1/1
2	NAG	P	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	P	2	2	-	0/6/23/26	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	K	7	MAN	O5-C1	-2.66	1.39	1.43
3	D	6	MAN	O5-C1	-2.45	1.39	1.43
3	D	10	MAN	O5-C1	-2.31	1.40	1.43
3	D	5	MAN	O5-C1	-2.30	1.40	1.43
4	K	5	MAN	O5-C1	-2.27	1.40	1.43
4	K	9	MAN	O5-C1	-2.13	1.40	1.43
3	D	4	MAN	O5-C1	-2.06	1.40	1.43
3	D	8	MAN	O5-C1	-2.05	1.40	1.43
4	K	8	MAN	O5-C1	-2.03	1.40	1.43

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	3	BMA	C1-O5-C5	7.53	122.40	112.19
4	K	3	BMA	C1-O5-C5	5.79	120.04	112.19
3	D	3	BMA	C1-O5-C5	5.64	119.84	112.19
2	O	2	NAG	C2-N2-C7	4.97	129.97	122.90
2	H	1	NAG	C2-N2-C7	4.42	129.19	122.90
5	L	3	BMA	C3-C4-C5	3.81	117.04	110.24
2	J	2	NAG	C2-N2-C7	3.38	127.71	122.90
3	D	7	MAN	C1-O5-C5	3.25	116.60	112.19
3	D	3	BMA	C3-C4-C5	3.22	115.99	110.24
2	O	1	NAG	C2-N2-C7	3.21	127.47	122.90
3	D	4	MAN	C1-O5-C5	3.19	116.51	112.19
2	C	1	NAG	C2-N2-C7	3.14	127.37	122.90
4	K	10	MAN	C1-O5-C5	3.11	116.41	112.19
2	H	2	NAG	C2-N2-C7	3.10	127.32	122.90
3	D	3	BMA	O4-C4-C3	-2.94	103.55	110.35
4	K	3	BMA	C3-C4-C5	2.94	115.49	110.24
4	K	3	BMA	C2-C3-C4	2.93	115.97	110.89
3	D	3	BMA	C2-C3-C4	2.87	115.85	110.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	2	NAG	C2-N2-C7	2.78	126.86	122.90
3	D	9	MAN	C1-O5-C5	2.73	115.90	112.19
4	K	3	BMA	O4-C4-C3	-2.70	104.10	110.35
3	D	8	MAN	C1-O5-C5	2.70	115.85	112.19
5	L	3	BMA	C2-C3-C4	2.67	115.52	110.89
2	O	2	NAG	C8-C7-N2	2.67	120.61	116.10
3	D	1	NAG	C2-N2-C7	2.64	126.66	122.90
4	K	3	BMA	O6-C6-C5	-2.58	102.43	111.29
3	D	3	BMA	O6-C6-C5	-2.52	102.64	111.29
2	M	1	NAG	C2-N2-C7	2.49	126.45	122.90
4	K	4	MAN	C1-O5-C5	2.47	115.54	112.19
4	K	1	NAG	O5-C1-C2	-2.47	107.38	111.29
2	H	1	NAG	O4-C4-C3	-2.47	104.63	110.35
5	L	1	NAG	O5-C1-C2	-2.30	107.66	111.29
2	H	2	NAG	O5-C1-C2	-2.28	107.69	111.29
5	L	3	BMA	O4-C4-C3	-2.27	105.11	110.35
2	F	2	NAG	O5-C1-C2	-2.19	107.83	111.29
5	L	3	BMA	O5-C5-C4	2.18	116.14	110.83
2	E	1	NAG	O5-C1-C2	-2.18	107.85	111.29
2	P	2	NAG	C1-O5-C5	2.17	115.13	112.19
4	K	8	MAN	C1-O5-C5	2.12	115.06	112.19
3	D	5	MAN	O4-C4-C3	-2.12	105.46	110.35
5	L	3	BMA	O3-C3-C2	-2.10	105.98	109.99
2	O	1	NAG	O4-C4-C3	-2.07	105.56	110.35
2	J	1	NAG	O4-C4-C3	-2.04	105.64	110.35
4	K	11	MAN	C1-O5-C5	2.03	114.95	112.19
2	G	1	NAG	C2-N2-C7	2.02	125.78	122.90

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	K	4	MAN	O5-C5-C6-O6
4	K	10	MAN	O5-C5-C6-O6
4	K	4	MAN	C4-C5-C6-O6
4	K	1	NAG	C4-C5-C6-O6
3	D	7	MAN	O5-C5-C6-O6
4	K	10	MAN	C4-C5-C6-O6
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	H	1	NAG	C8-C7-N2-C2
2	H	1	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	H	2	NAG	C8-C7-N2-C2
2	H	2	NAG	O7-C7-N2-C2
2	J	2	NAG	C8-C7-N2-C2
2	J	2	NAG	O7-C7-N2-C2
2	O	1	NAG	C8-C7-N2-C2
2	O	1	NAG	O7-C7-N2-C2
2	O	2	NAG	C8-C7-N2-C2
2	O	2	NAG	O7-C7-N2-C2
4	K	1	NAG	O5-C5-C6-O6
4	K	8	MAN	O5-C5-C6-O6
3	D	7	MAN	C4-C5-C6-O6
4	K	9	MAN	C4-C5-C6-O6
4	K	6	MAN	O5-C5-C6-O6
4	K	8	MAN	C4-C5-C6-O6
4	K	9	MAN	O5-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
3	D	10	MAN	O5-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
2	M	2	NAG	O5-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
5	L	2	NAG	O5-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
2	M	1	NAG	O5-C5-C6-O6
2	N	2	NAG	O5-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
2	N	2	NAG	C3-C2-N2-C7
3	D	5	MAN	C4-C5-C6-O6
4	K	6	MAN	C4-C5-C6-O6
3	D	5	MAN	O5-C5-C6-O6
2	G	2	NAG	C3-C2-N2-C7
2	O	1	NAG	O5-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
4	K	5	MAN	C4-C5-C6-O6

There are no ring outliers.

7 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	1	NAG	2	0
2	O	2	NAG	1	0
2	O	1	NAG	1	0
2	H	2	NAG	1	0

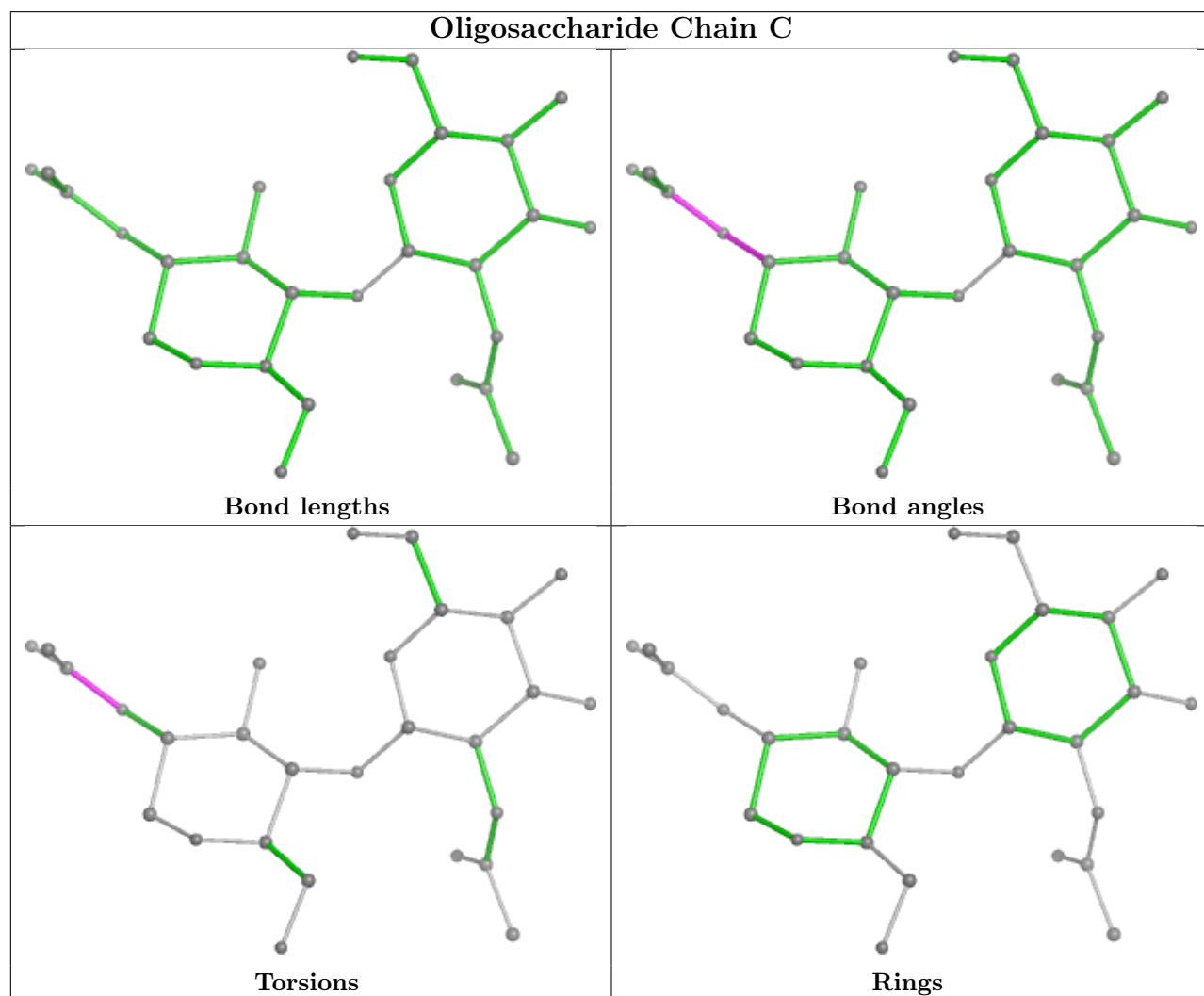
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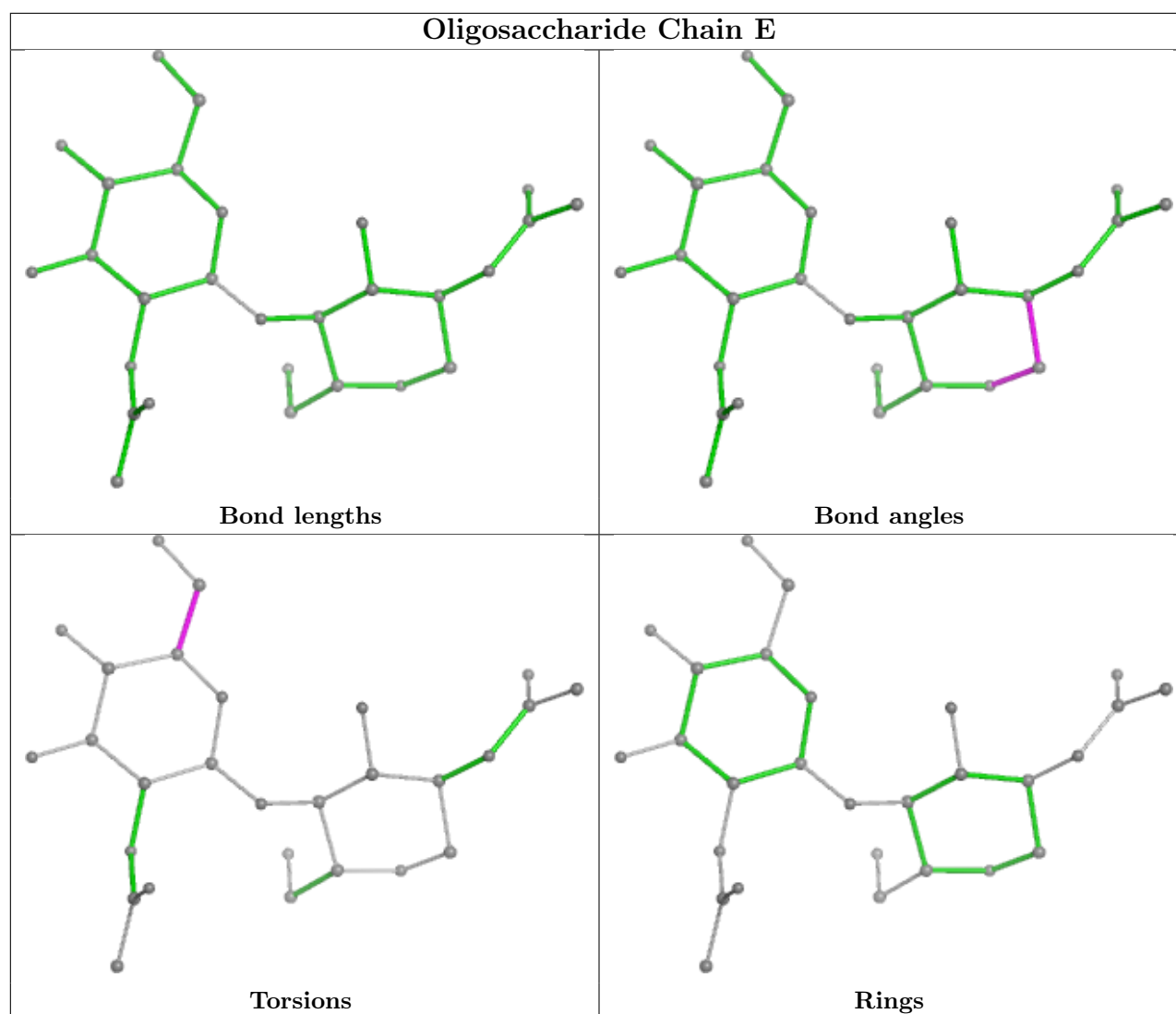


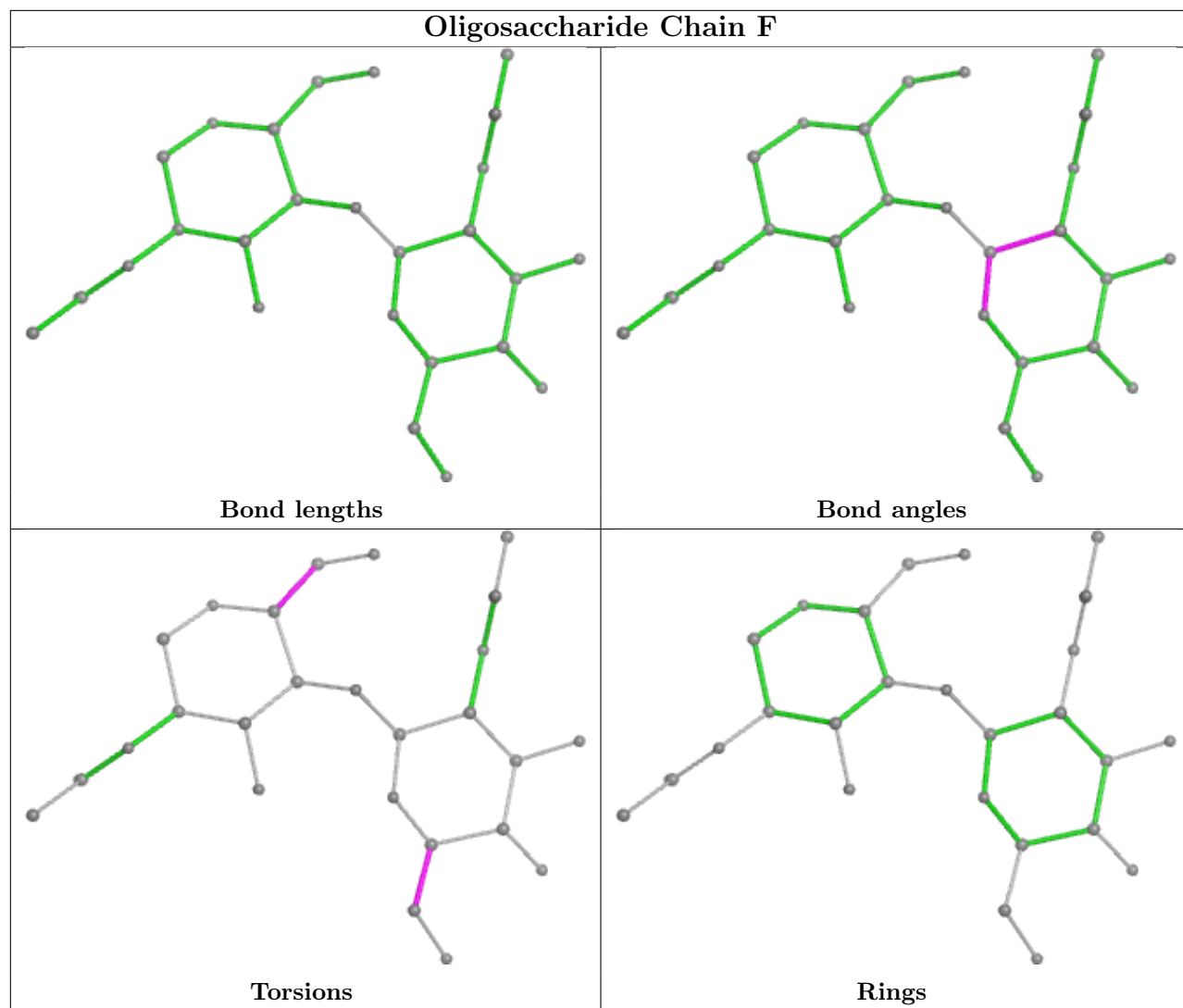
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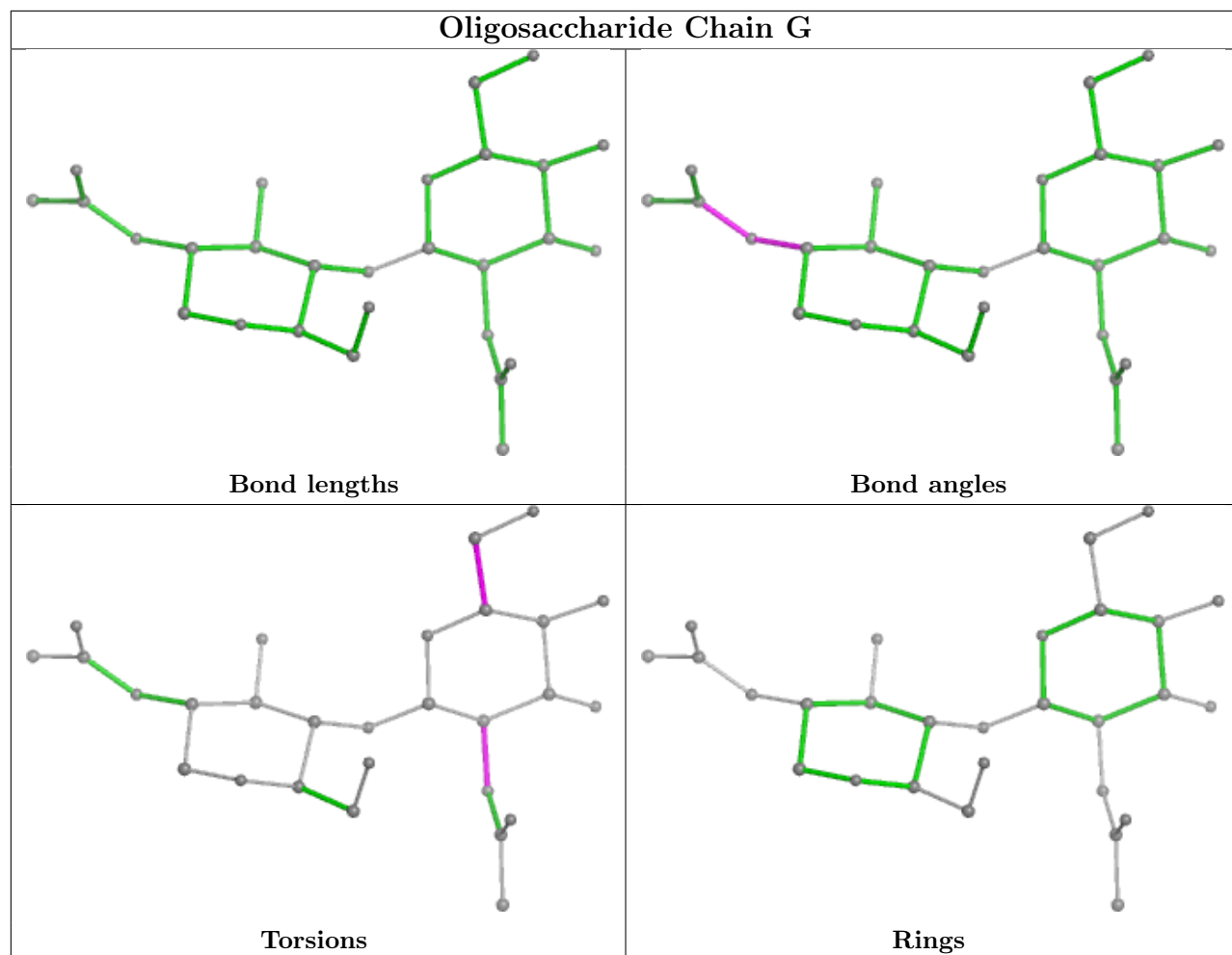
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	2	NAG	1	0
2	F	2	NAG	1	0
2	F	1	NAG	1	0

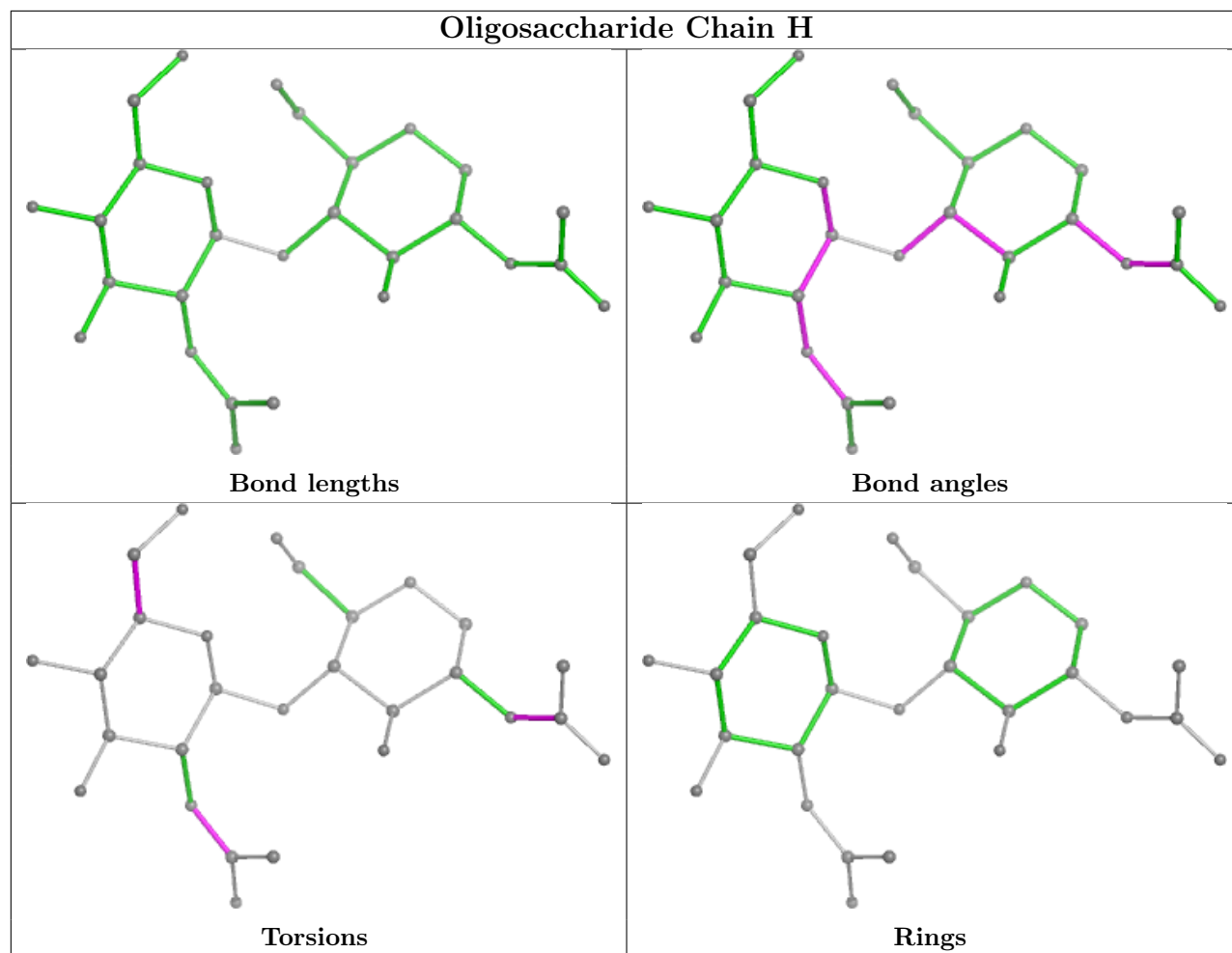
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

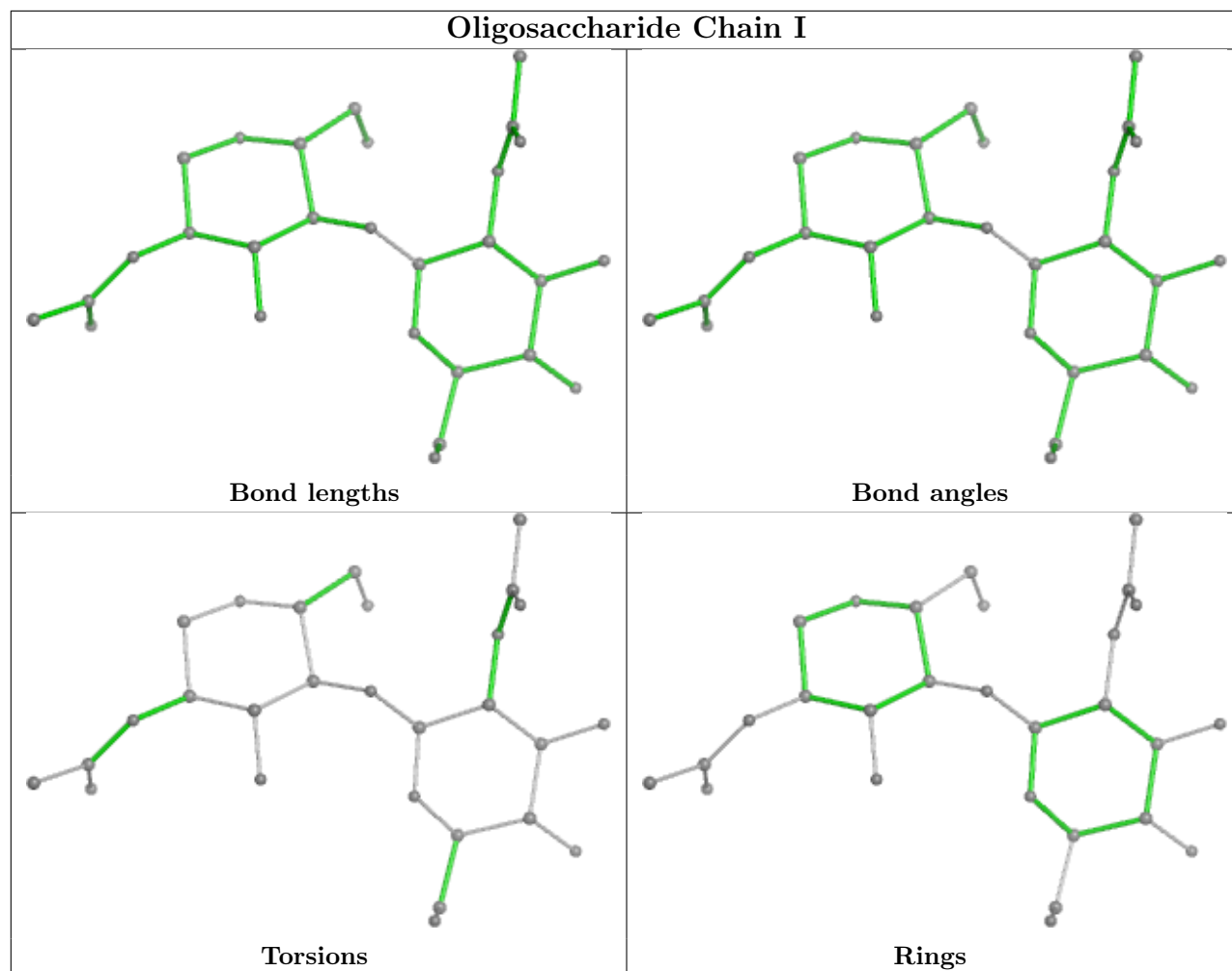


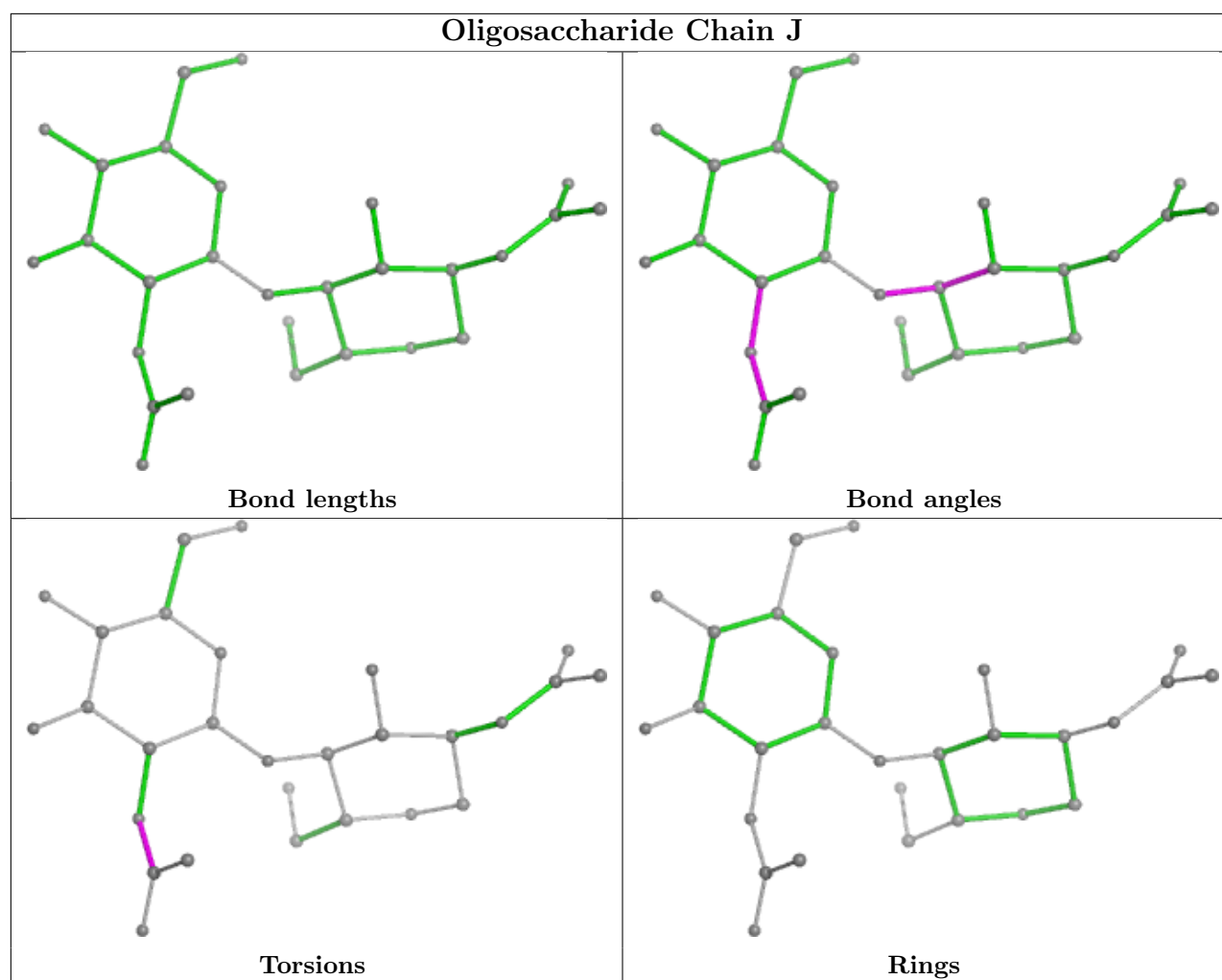


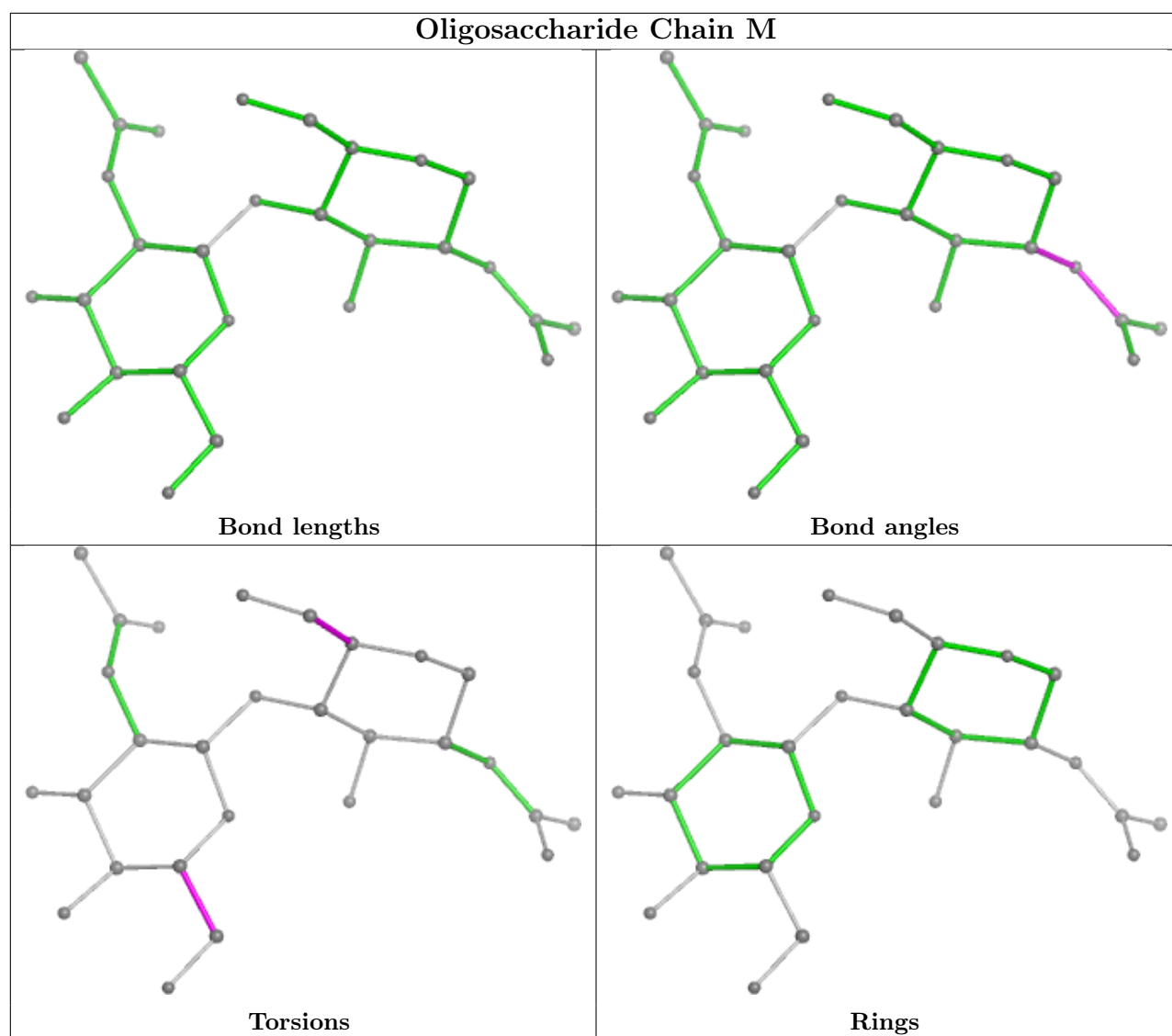




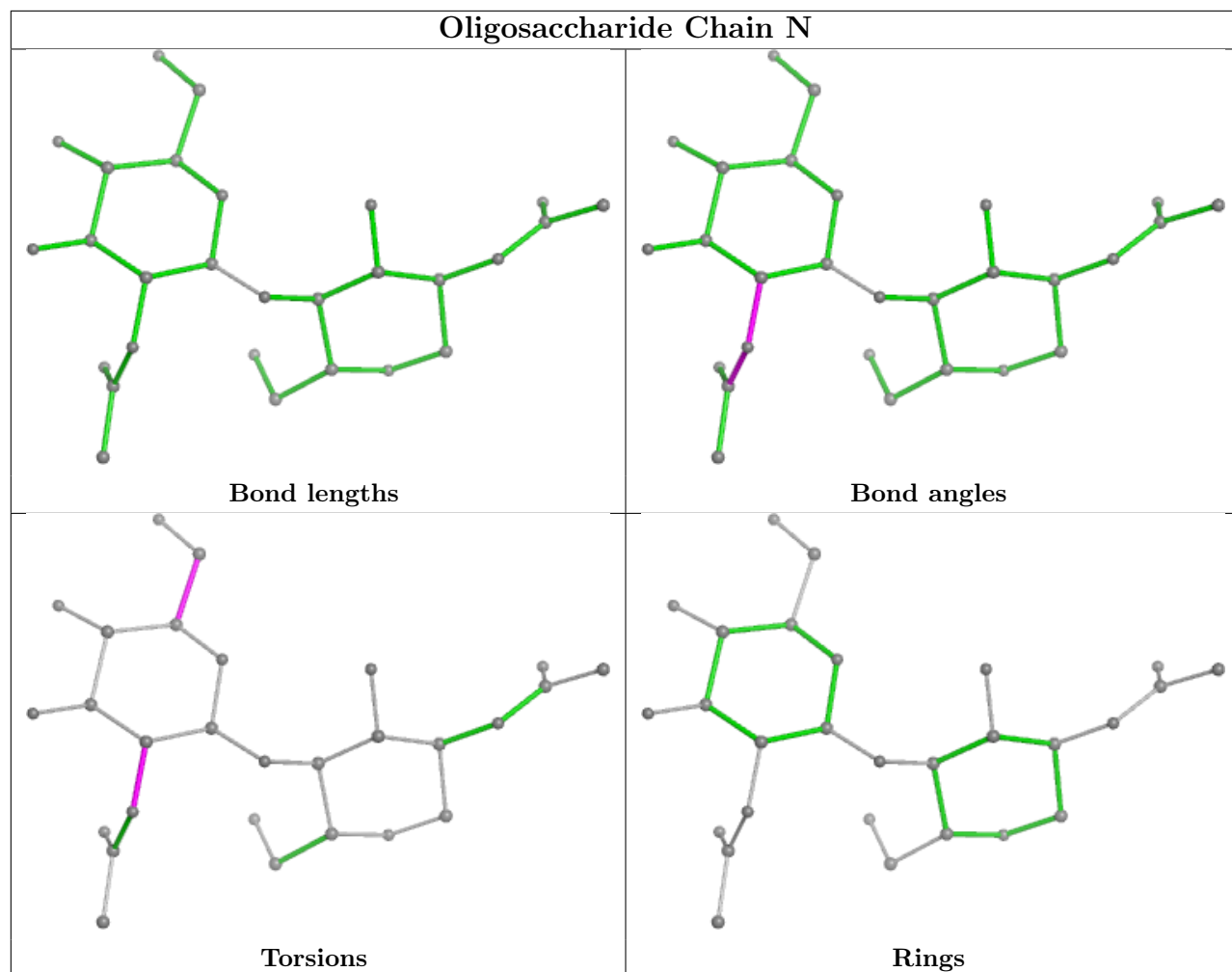




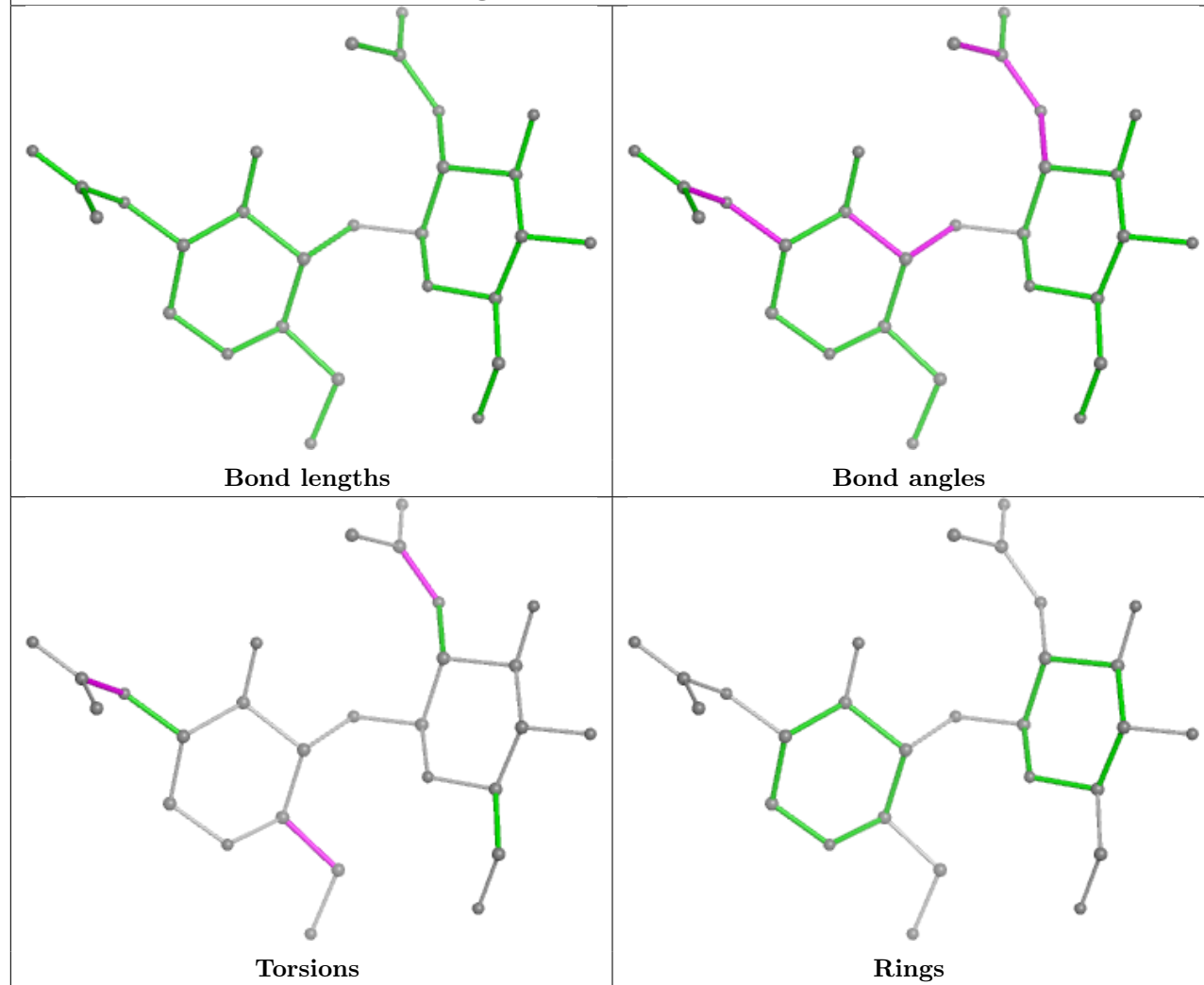


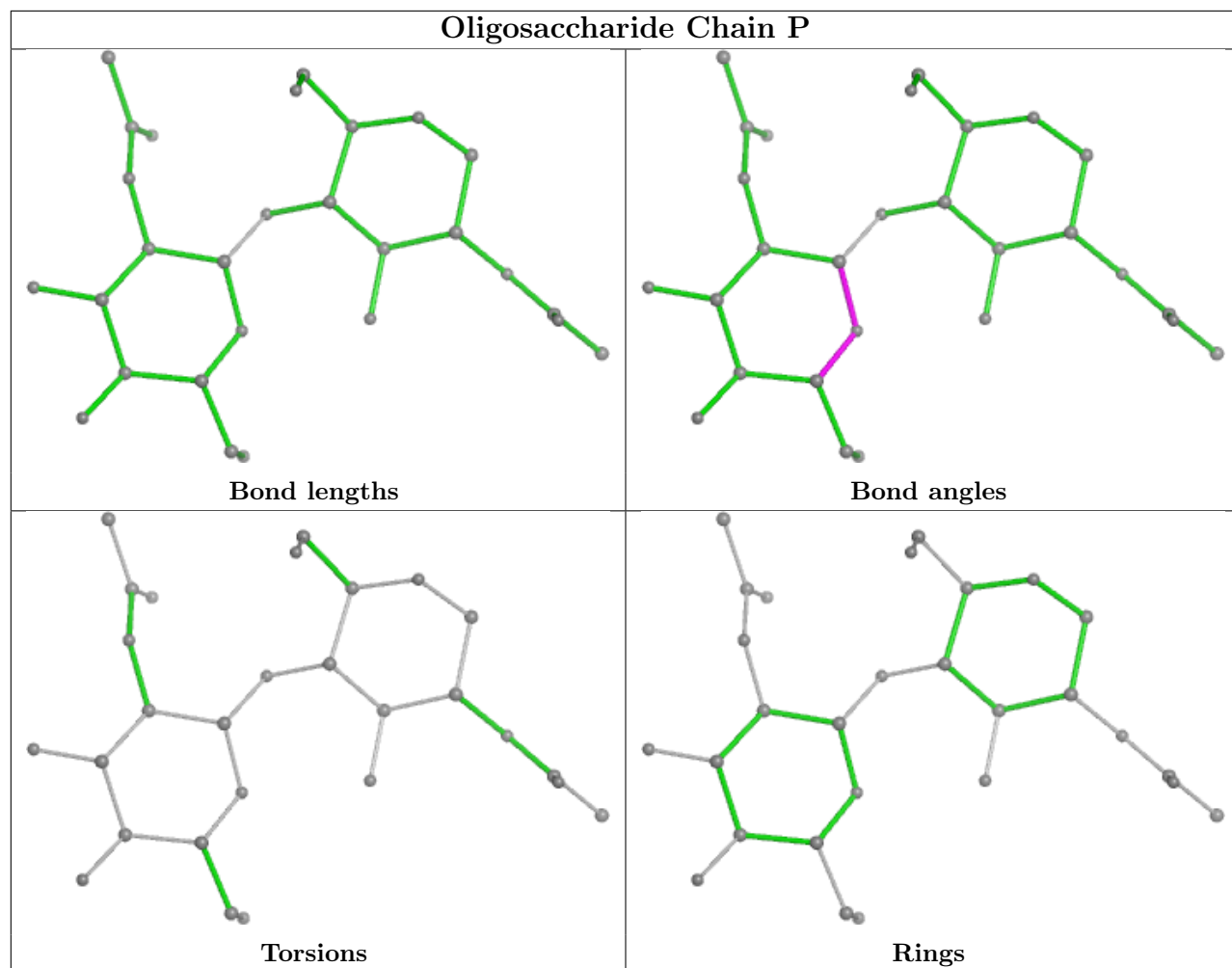


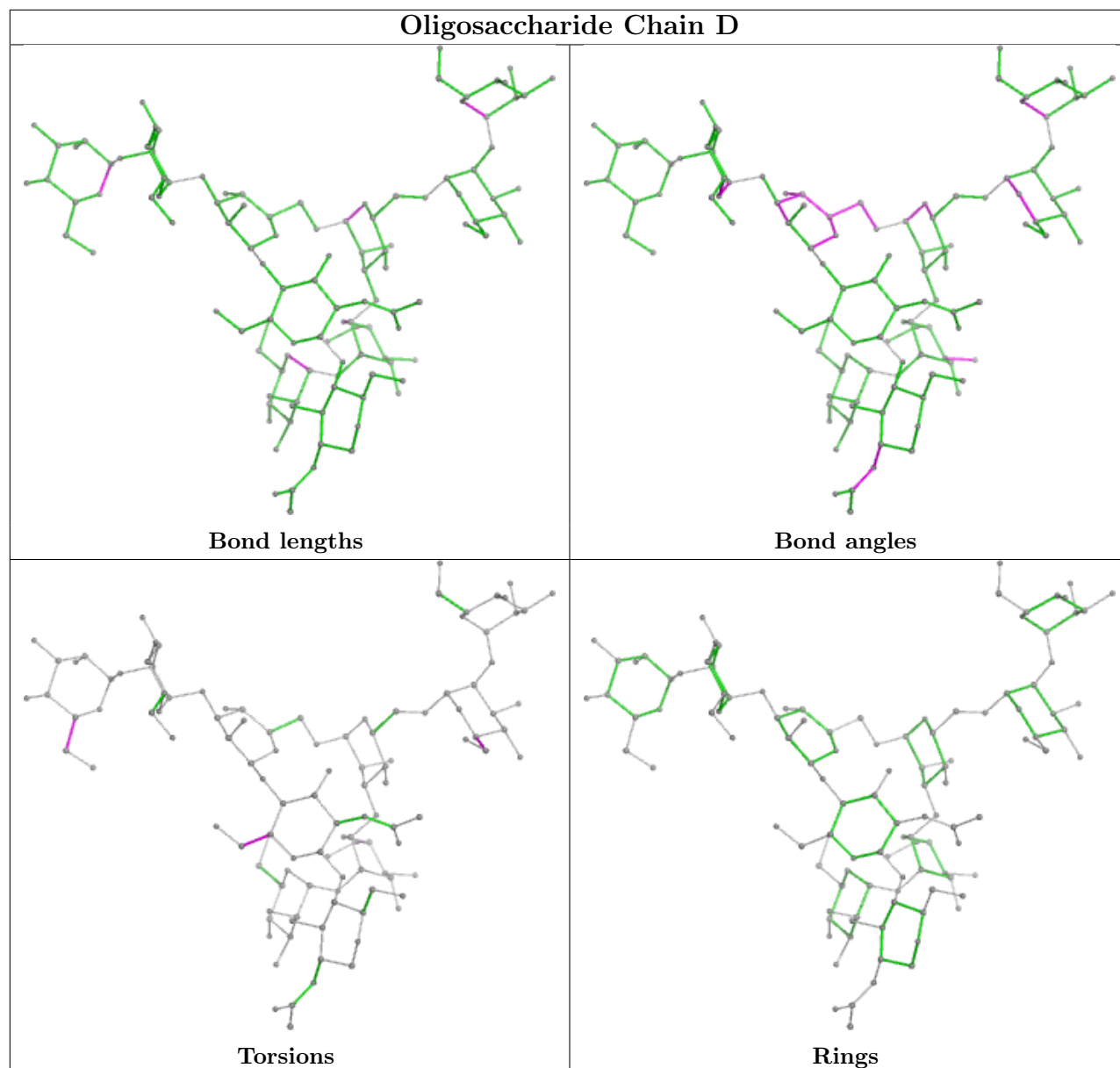




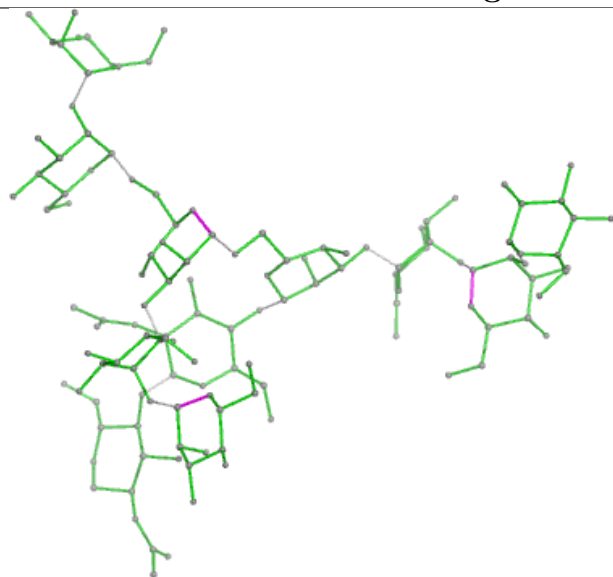
## Oligosaccharide Chain O



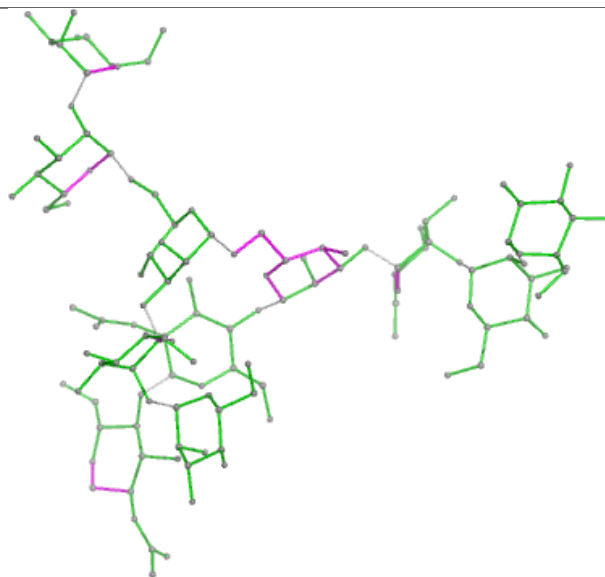




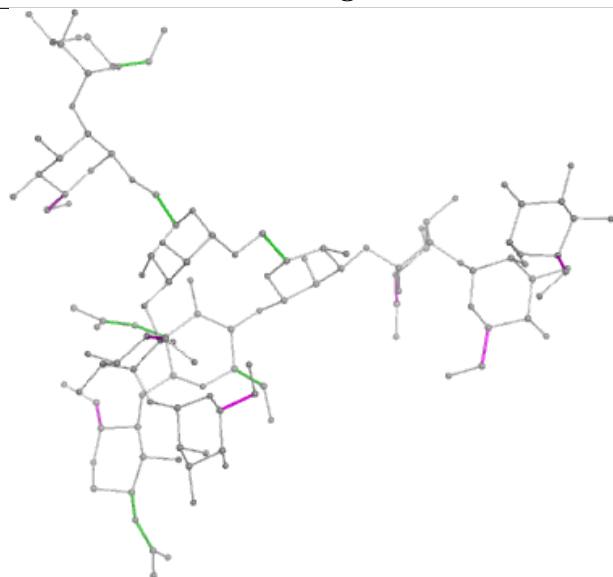
## Oligosaccharide Chain K



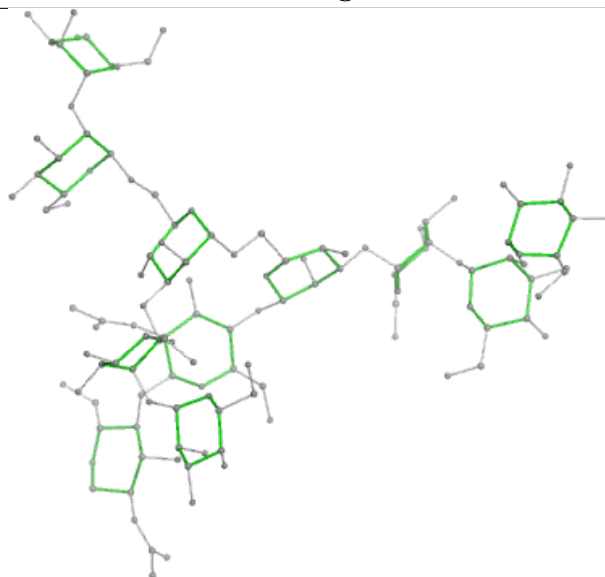
Bond lengths



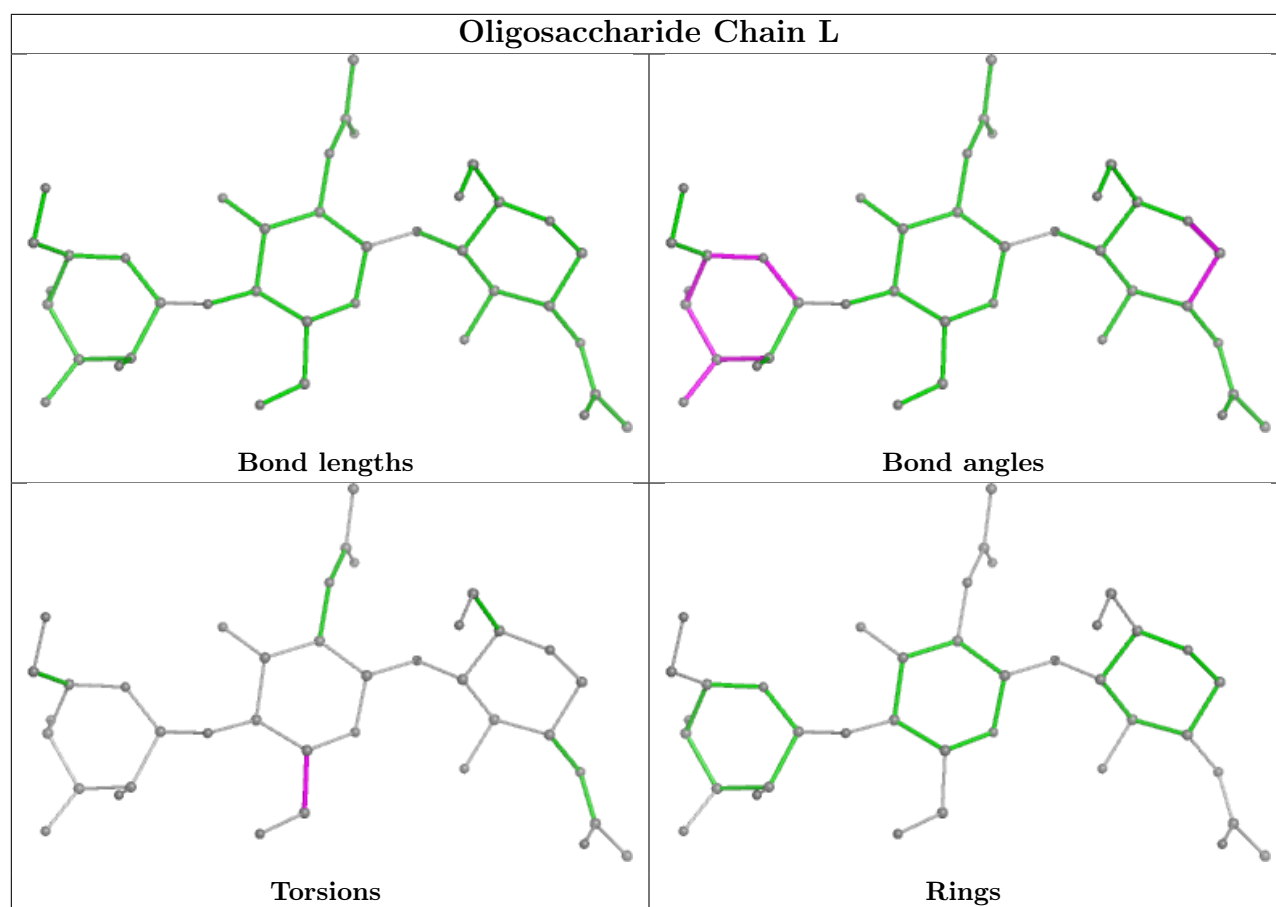
Bond angles



Torsions



Rings



## 5.6 Ligand geometry [i](#)

11 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	$\# Z  > 2$
7	NAG	A	805	1	14,14,15	0.71	0	17,19,21	1.05	1 (5%)
7	NAG	A	803	1	14,14,15	0.68	0	17,19,21	0.98	0
6	A1JMI	B	801	-	35,35,35	1.80	8 (22%)	55,56,56	2.22	21 (38%)
7	NAG	B	802	1	14,14,15	0.69	0	17,19,21	1.13	1 (5%)
6	A1JMI	A	801	-	35,35,35	1.69	8 (22%)	55,56,56	2.25	21 (38%)
7	NAG	B	803	1	14,14,15	0.72	0	17,19,21	0.81	0
7	NAG	B	805	1	14,14,15	0.78	0	17,19,21	1.00	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
7	NAG	A	802	1	14,14,15	0.70	0	17,19,21	0.89	0
7	NAG	B	804	1	14,14,15	0.68	0	17,19,21	1.41	3 (17%)
8	ACT	B	806	-	3,3,3	1.16	0	3,3,3	1.19	0
7	NAG	A	804	1	14,14,15	0.70	0	17,19,21	0.97	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	A	805	1	-	1/6/23/26	0/1/1/1
7	NAG	A	803	1	-	0/6/23/26	0/1/1/1
6	A1JMI	B	801	-	-	6/20/81/81	0/4/4/4
7	NAG	B	802	1	-	2/6/23/26	0/1/1/1
6	A1JMI	A	801	-	-	6/20/81/81	0/4/4/4
7	NAG	B	803	1	-	0/6/23/26	0/1/1/1
7	NAG	B	805	1	-	0/6/23/26	0/1/1/1
7	NAG	A	802	1	-	0/6/23/26	0/1/1/1
7	NAG	B	804	1	-	1/6/23/26	0/1/1/1
7	NAG	A	804	1	-	0/6/23/26	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	801	A1JMI	C15-C10	4.06	1.63	1.56
6	B	801	A1JMI	C25-C24	3.97	1.61	1.53
6	A	801	A1JMI	C15-C10	3.89	1.63	1.56
6	B	801	A1JMI	C11-C06	3.80	1.60	1.53
6	A	801	A1JMI	C07-C03	3.52	1.61	1.55
6	A	801	A1JMI	C25-C24	3.47	1.60	1.53
6	A	801	A1JMI	O20-C18	-3.15	1.36	1.43
6	B	801	A1JMI	O20-C18	-3.11	1.36	1.43
6	B	801	A1JMI	C07-C03	2.75	1.60	1.55
6	A	801	A1JMI	C02-C03	2.67	1.58	1.54
6	A	801	A1JMI	C11-C06	2.60	1.58	1.53
6	A	801	A1JMI	C19-C18	-2.60	1.48	1.52
6	B	801	A1JMI	C19-C18	-2.59	1.48	1.52
6	B	801	A1JMI	C02-C03	2.57	1.58	1.54
6	A	801	A1JMI	C16-C17	-2.10	1.50	1.53
6	B	801	A1JMI	C02-C24	2.00	1.59	1.54

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	801	A1JMI	C22-C15-C14	-7.50	97.63	110.36
6	A	801	A1JMI	C22-C15-C14	-7.42	97.78	110.36
6	B	801	A1JMI	C19-C14-C13	-4.64	103.23	111.74
6	A	801	A1JMI	C10-C11-C06	4.42	115.01	109.09
6	A	801	A1JMI	C19-C14-C13	-4.40	103.68	111.74
6	A	801	A1JMI	C14-C19-C18	-4.35	109.66	114.46
6	B	801	A1JMI	C14-C19-C18	-4.09	109.95	114.46
6	B	801	A1JMI	C10-C11-C06	4.03	114.48	109.09
6	A	801	A1JMI	C16-C17-C18	-3.89	106.90	111.36
6	B	801	A1JMI	C10-C15-C14	3.71	113.79	108.58
6	A	801	A1JMI	C10-C15-C14	3.70	113.78	108.58
6	B	801	A1JMI	C16-C17-C18	-3.66	107.17	111.36
6	A	801	A1JMI	C23-C07-C08	-3.58	104.93	110.59
7	B	802	NAG	C2-N2-C7	3.34	127.66	122.90
6	A	801	A1JMI	C16-C15-C10	3.22	114.90	110.08
6	B	801	A1JMI	C16-C15-C10	3.14	114.78	110.08
6	A	801	A1JMI	C22-C15-C16	-3.14	104.21	108.97
6	A	801	A1JMI	C08-C07-C03	3.10	121.22	116.57
6	B	801	A1JMI	C23-C07-C08	-3.09	105.70	110.59
6	A	801	A1JMI	C07-C03-C02	-2.93	115.51	118.89
7	B	804	NAG	C2-N2-C7	2.92	127.06	122.90
6	B	801	A1JMI	C05-C06-C11	2.90	123.85	119.08
6	B	801	A1JMI	C22-C15-C16	-2.83	104.67	108.97
6	B	801	A1JMI	C12-C11-C10	-2.79	107.03	110.49
7	A	805	NAG	C1-O5-C5	2.70	115.84	112.19
6	B	801	A1JMI	C05-C06-C07	-2.58	100.73	103.84
7	B	804	NAG	C1-O5-C5	2.57	115.67	112.19
6	A	801	A1JMI	C12-C11-C10	-2.54	107.34	110.49
6	A	801	A1JMI	C22-C15-C10	2.50	114.63	111.18
6	A	801	A1JMI	C16-C15-C14	2.44	109.89	107.14
6	B	801	A1JMI	C16-C15-C14	2.44	109.89	107.14
6	B	801	A1JMI	C06-C07-C03	2.43	102.96	100.07
7	B	804	NAG	O5-C1-C2	-2.40	107.50	111.29
6	B	801	A1JMI	C22-C15-C10	2.37	114.45	111.18
6	A	801	A1JMI	C01-C02-C24	2.34	116.12	111.39
6	B	801	A1JMI	C07-C03-C02	-2.29	116.25	118.89
6	A	801	A1JMI	C19-C18-C17	2.28	113.10	110.27
6	B	801	A1JMI	C09-C10-C11	-2.24	108.53	111.75
6	B	801	A1JMI	C30-C28-C26	2.20	116.81	112.47
6	A	801	A1JMI	O31-C25-C24	-2.18	104.39	109.47
6	A	801	A1JMI	C05-C06-C11	2.16	122.63	119.08
6	A	801	A1JMI	O20-C18-C17	-2.16	105.44	110.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	801	A1JMI	C19-C18-C17	2.16	112.94	110.27
6	B	801	A1JMI	C07-C06-C11	-2.12	111.25	114.38
6	A	801	A1JMI	C09-C10-C11	-2.11	108.71	111.75
6	A	801	A1JMI	C05-C06-C07	-2.09	101.32	103.84
7	A	804	NAG	O5-C1-C2	-2.07	108.02	111.29
6	B	801	A1JMI	C08-C07-C03	2.05	119.63	116.57
7	B	805	NAG	C1-O5-C5	2.04	114.96	112.19

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	801	A1JMI	C02-C24-C25-C26
6	A	801	A1JMI	C02-C24-C25-O31
6	A	801	A1JMI	O32-C24-C25-C26
7	B	802	NAG	C8-C7-N2-C2
7	B	802	NAG	O7-C7-N2-C2
6	A	801	A1JMI	O32-C24-C25-O31
6	B	801	A1JMI	O32-C24-C25-O31
7	B	804	NAG	O5-C5-C6-O6
7	A	805	NAG	O5-C5-C6-O6
6	A	801	A1JMI	O31-C25-C26-C27
6	B	801	A1JMI	O31-C25-C26-C27
6	B	801	A1JMI	C02-C24-C25-O31
6	B	801	A1JMI	C02-C24-C25-C26
6	B	801	A1JMI	O32-C24-C25-C26
6	A	801	A1JMI	O31-C25-C26-C28
6	B	801	A1JMI	O31-C25-C26-C28

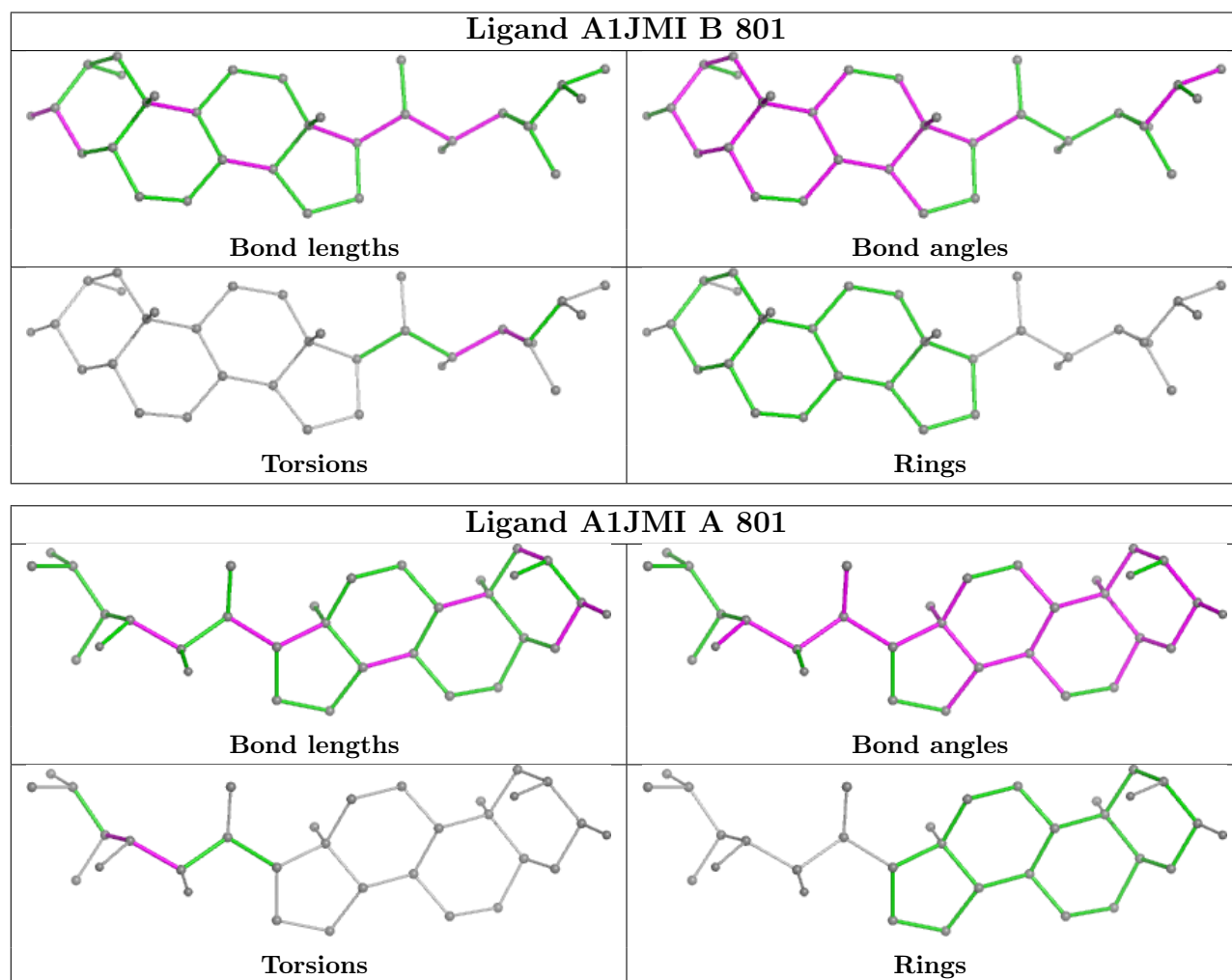
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	801	A1JMI	1	0
7	B	803	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for 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 [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	721/779 (92%)	-0.77	0 <span>100</span> <span>100</span>	83, 140, 217, 282	1 (0%)
1	B	727/779 (93%)	-0.65	3 (0%) <span>89</span> <span>81</span>	74, 136, 204, 269	6 (0%)
All	All	1448/1558 (92%)	-0.71	3 (0%) <span>92</span> <span>87</span>	74, 138, 213, 282	7 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	567	LEU	2.9
1	B	81	LEU	2.8
1	B	78	VAL	2.3

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	NAG	H	2	14/15	0.26	0.09	163,177,212,215	0
2	NAG	C	2	14/15	0.48	0.07	143,185,224,229	0
2	NAG	H	1	14/15	0.53	0.06	133,164,204,209	0
3	MAN	D	9	11/12	0.53	0.09	120,152,193,194	0
2	NAG	J	1	14/15	0.57	0.09	137,169,210,212	0
2	NAG	J	2	14/15	0.58	0.07	129,173,212,212	0

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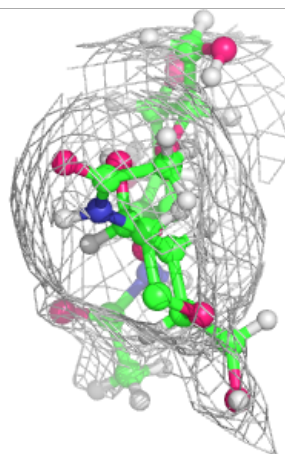
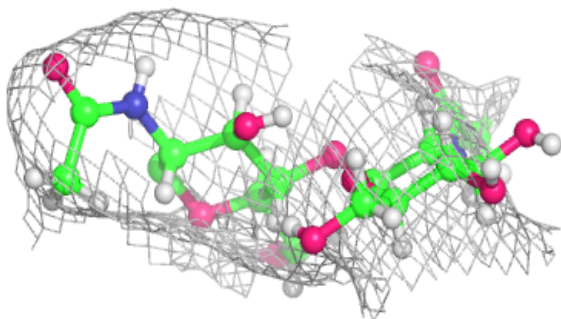
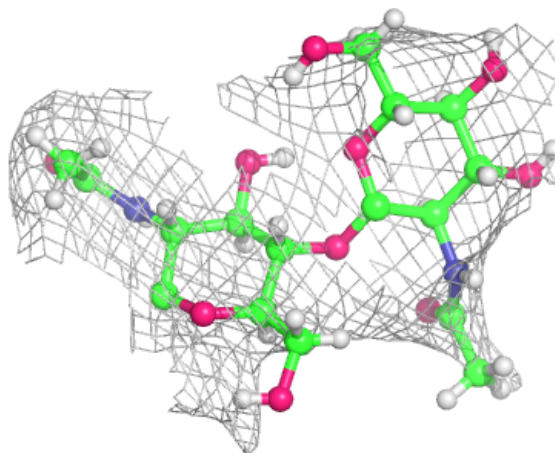
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAG	P	2	14/15	0.60	0.06	139,165,195,208	0
2	NAG	O	2	14/15	0.64	0.07	150,166,201,202	0
2	NAG	F	2	14/15	0.65	0.08	157,172,203,213	0
2	NAG	I	2	14/15	0.67	0.07	166,189,227,230	0
3	MAN	D	10	11/12	0.68	0.07	137,155,189,189	0
2	NAG	M	2	14/15	0.69	0.06	161,196,235,239	0
4	MAN	K	6	11/12	0.72	0.10	169,200,228,242	0
3	MAN	D	7	11/12	0.73	0.07	95,118,139,149	0
2	NAG	F	1	14/15	0.73	0.08	131,151,181,184	0
2	NAG	N	2	14/15	0.74	0.08	126,154,192,194	0
2	NAG	C	1	14/15	0.75	0.07	148,183,216,222	0
2	NAG	M	1	14/15	0.76	0.07	102,146,181,189	0
2	NAG	E	2	14/15	0.77	0.07	120,148,185,192	0
2	NAG	O	1	14/15	0.77	0.06	115,151,190,193	0
2	NAG	P	1	14/15	0.78	0.09	111,152,186,191	0
4	NAG	K	2	14/15	0.79	0.08	83,110,134,143	0
5	NAG	L	2	14/15	0.79	0.07	125,154,191,216	0
4	MAN	K	5	11/12	0.81	0.05	148,173,203,210	0
4	MAN	K	4	11/12	0.82	0.06	132,154,187,188	0
5	BMA	L	3	11/12	0.82	0.06	167,184,217,223	0
3	NAG	D	2	14/15	0.83	0.08	90,104,127,134	0
2	NAG	I	1	14/15	0.84	0.08	131,170,198,209	0
4	MAN	K	10	11/12	0.85	0.05	91,115,147,148	0
4	MAN	K	11	11/12	0.85	0.06	103,140,171,173	0
4	NAG	K	1	14/15	0.86	0.11	88,121,152,160	0
3	MAN	D	8	11/12	0.87	0.05	107,131,159,164	0
2	NAG	G	1	14/15	0.87	0.08	121,137,168,168	0
3	NAG	D	1	14/15	0.87	0.12	93,118,144,144	0
3	MAN	D	6	11/12	0.88	0.08	92,127,154,160	0
2	NAG	G	2	14/15	0.88	0.06	106,149,184,195	0
3	MAN	D	4	11/12	0.89	0.08	94,111,136,140	0
3	BMA	D	3	11/12	0.89	0.06	85,121,147,163	0
4	BMA	K	3	11/12	0.89	0.05	86,109,137,159	0
4	MAN	K	8	11/12	0.89	0.06	78,101,121,124	0
2	NAG	N	1	14/15	0.90	0.06	98,130,154,158	0
2	NAG	E	1	14/15	0.90	0.06	103,131,162,162	0
5	NAG	L	1	14/15	0.92	0.05	102,136,172,172	0
4	MAN	K	9	11/12	0.93	0.08	104,119,145,154	0
4	MAN	K	7	11/12	0.93	0.04	98,114,139,143	0
3	MAN	D	5	11/12	0.94	0.06	81,99,121,121	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.

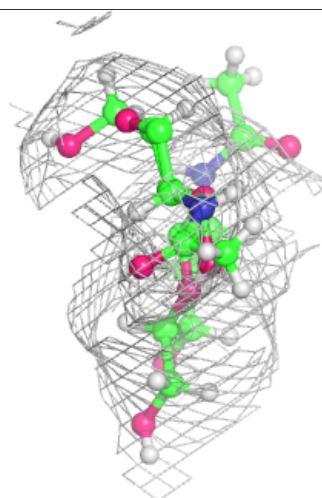
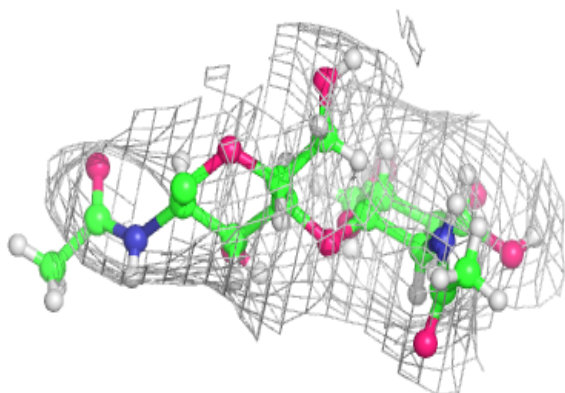
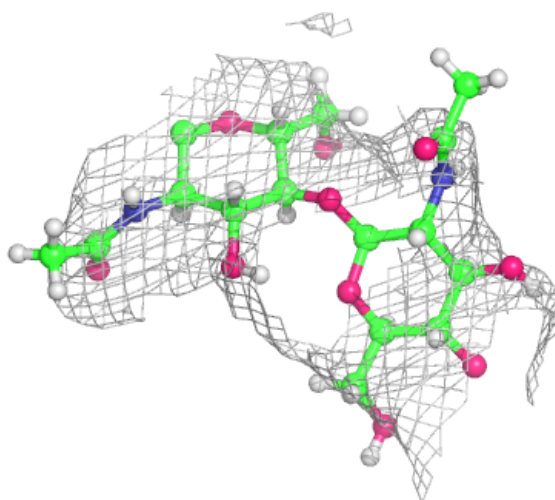
**Electron density around Chain C:**

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



**Electron density around Chain E:**

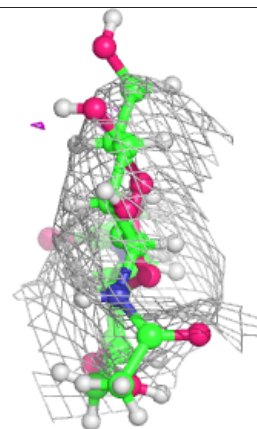
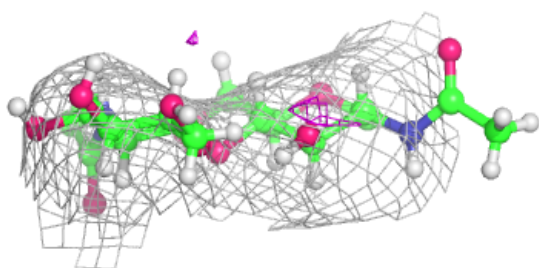
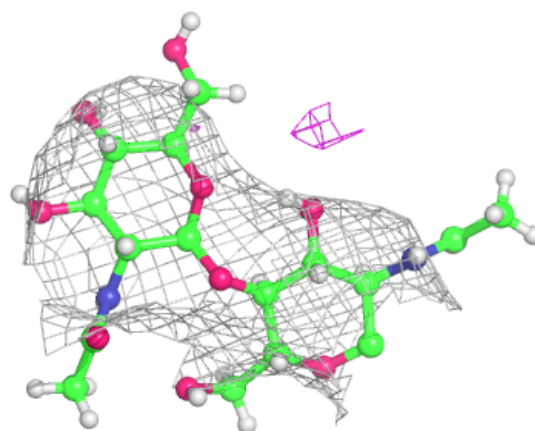
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



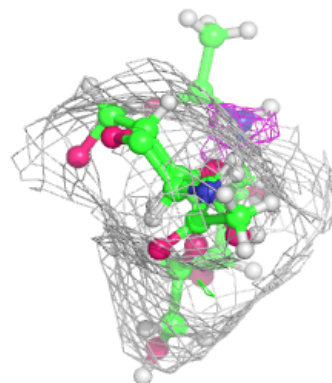
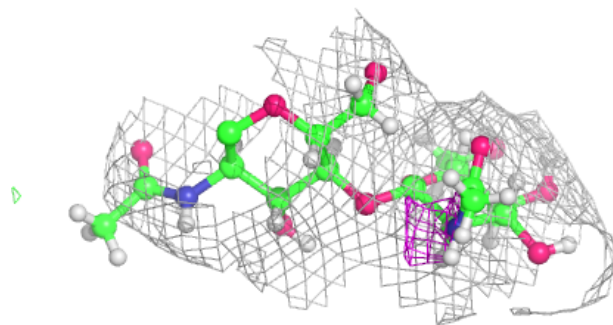
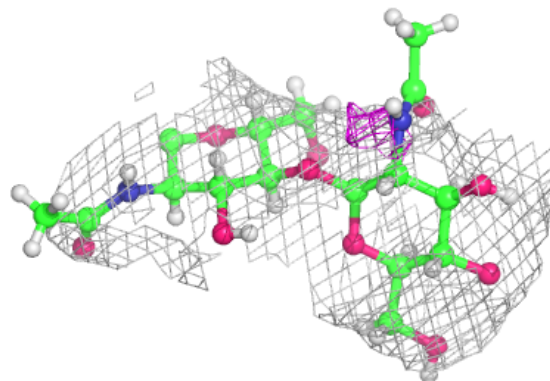


**Electron density around Chain F:**

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

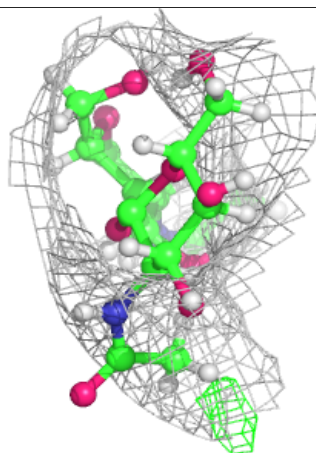
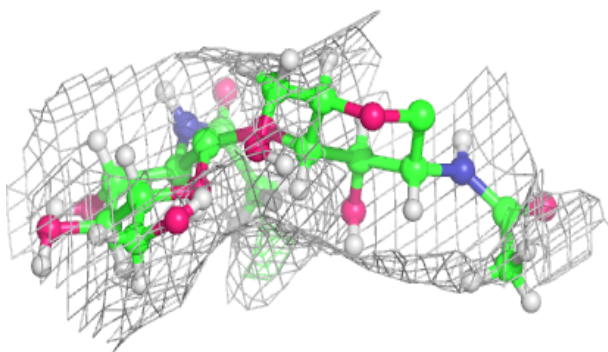
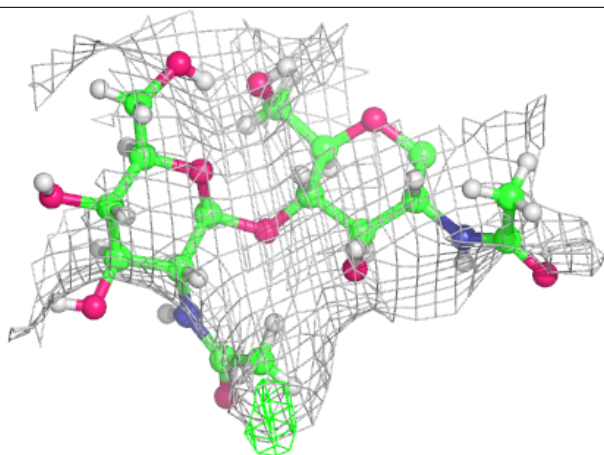
**Electron density around Chain G:**

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



**Electron density around Chain H:**

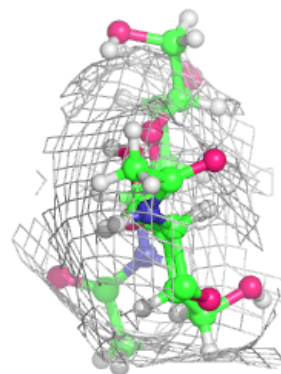
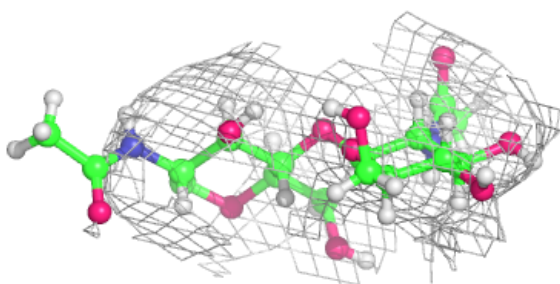
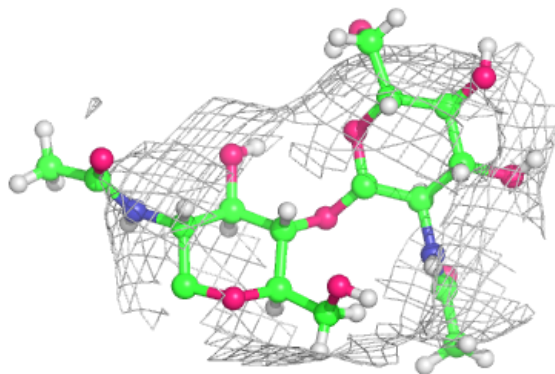
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



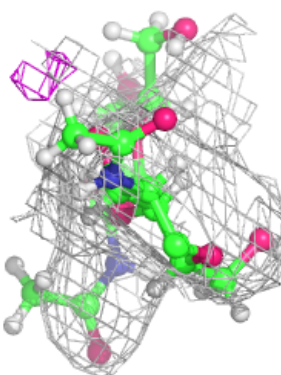
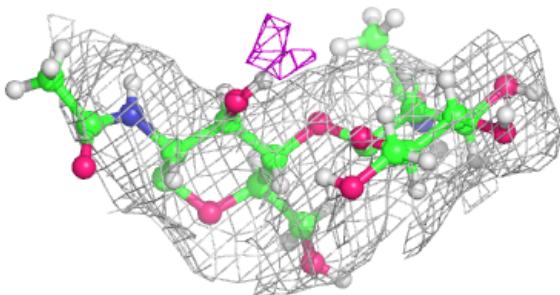
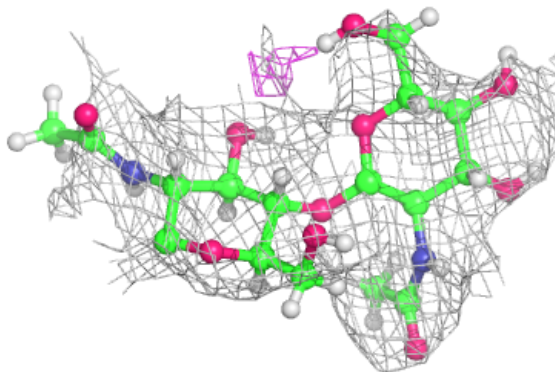


**Electron density around Chain I:**

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

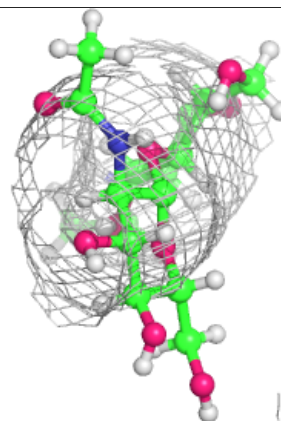
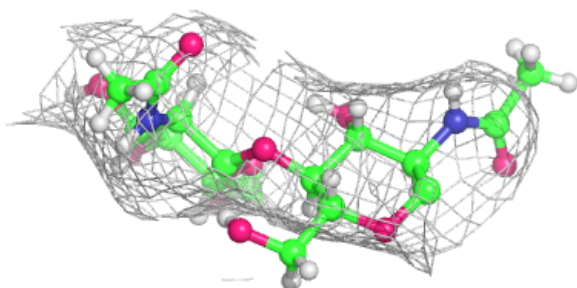
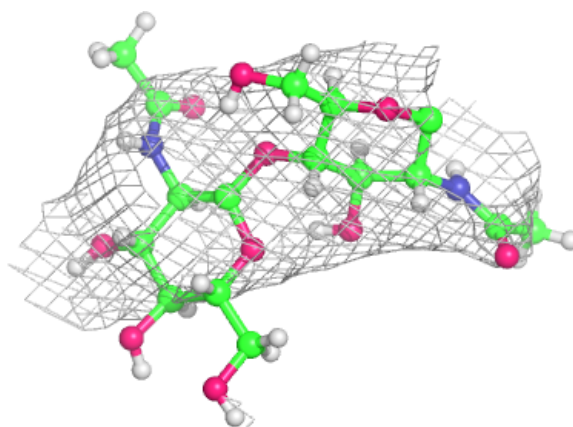
**Electron density around Chain J:**

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

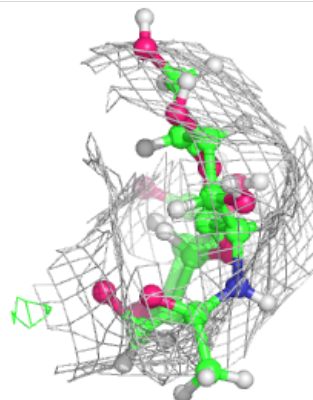
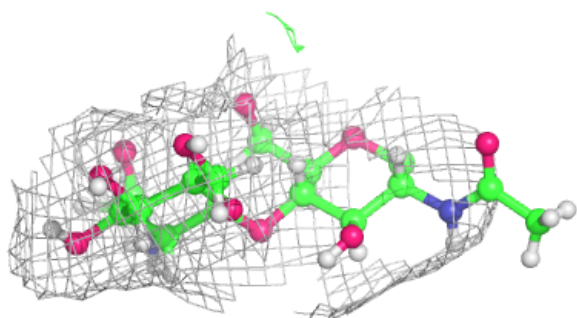
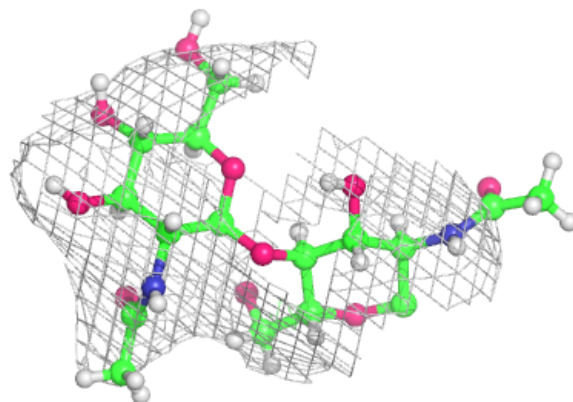


**Electron density around Chain M:**

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

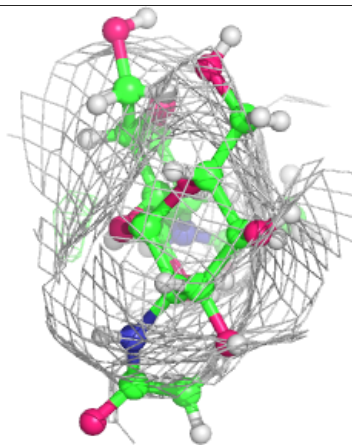
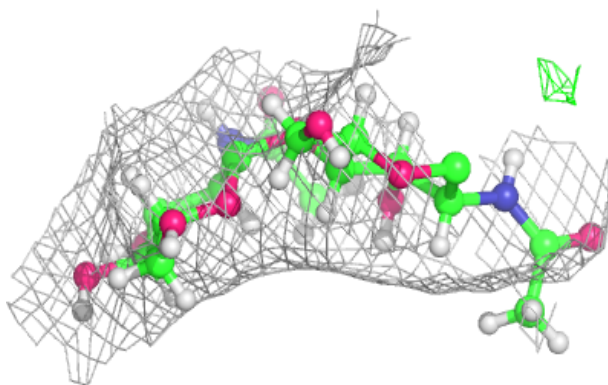
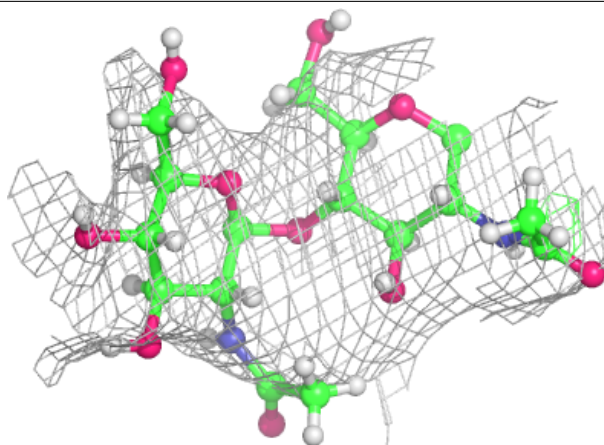
**Electron density around Chain N:**

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



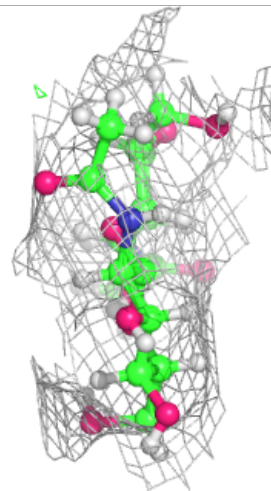
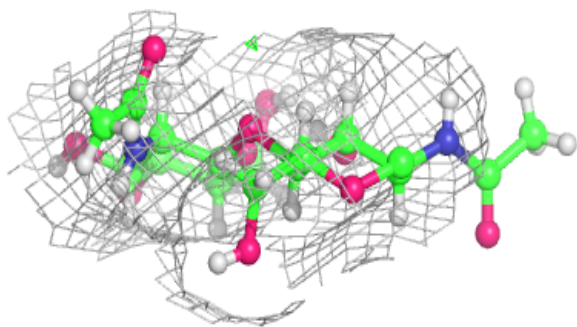
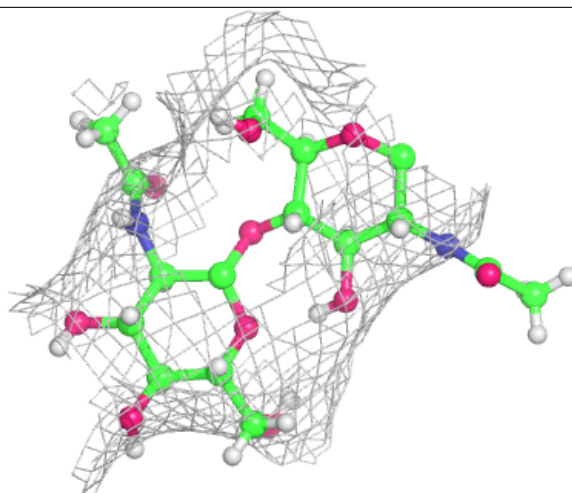
**Electron density around Chain O:**

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



**Electron density around Chain P:**

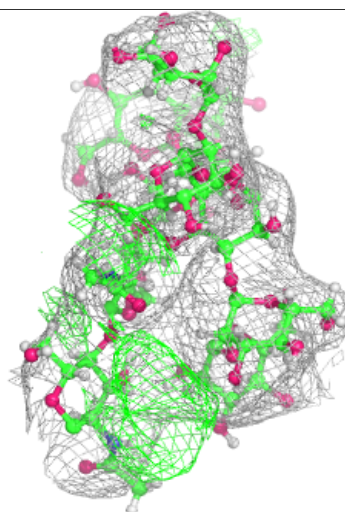
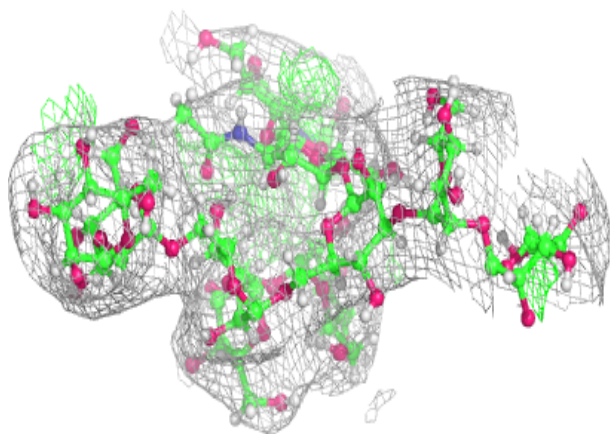
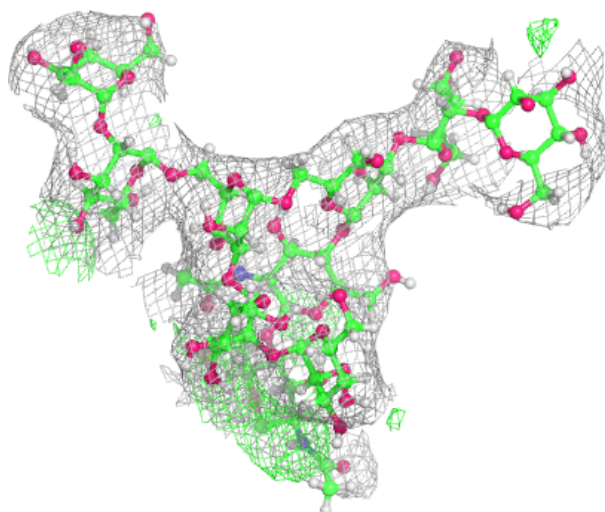
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





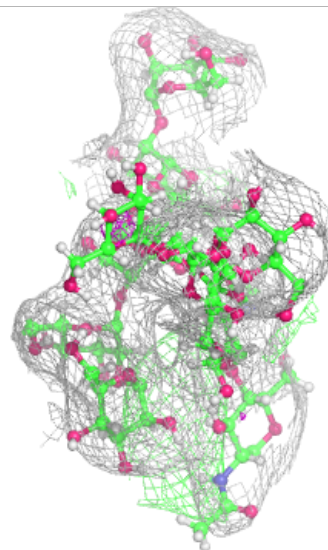
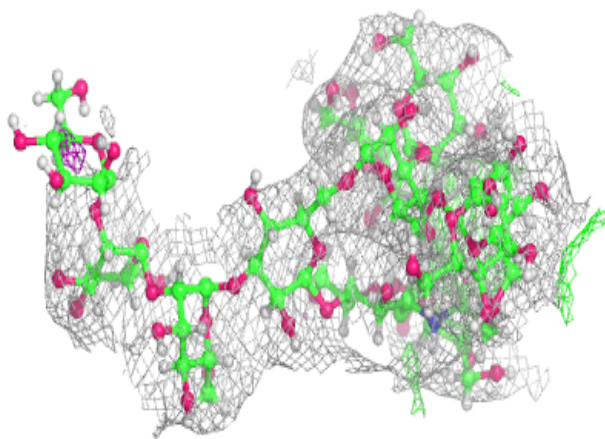
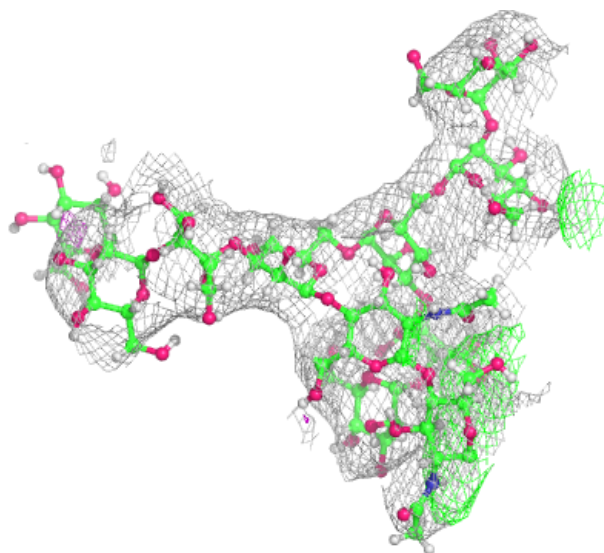
**Electron density around Chain D:**

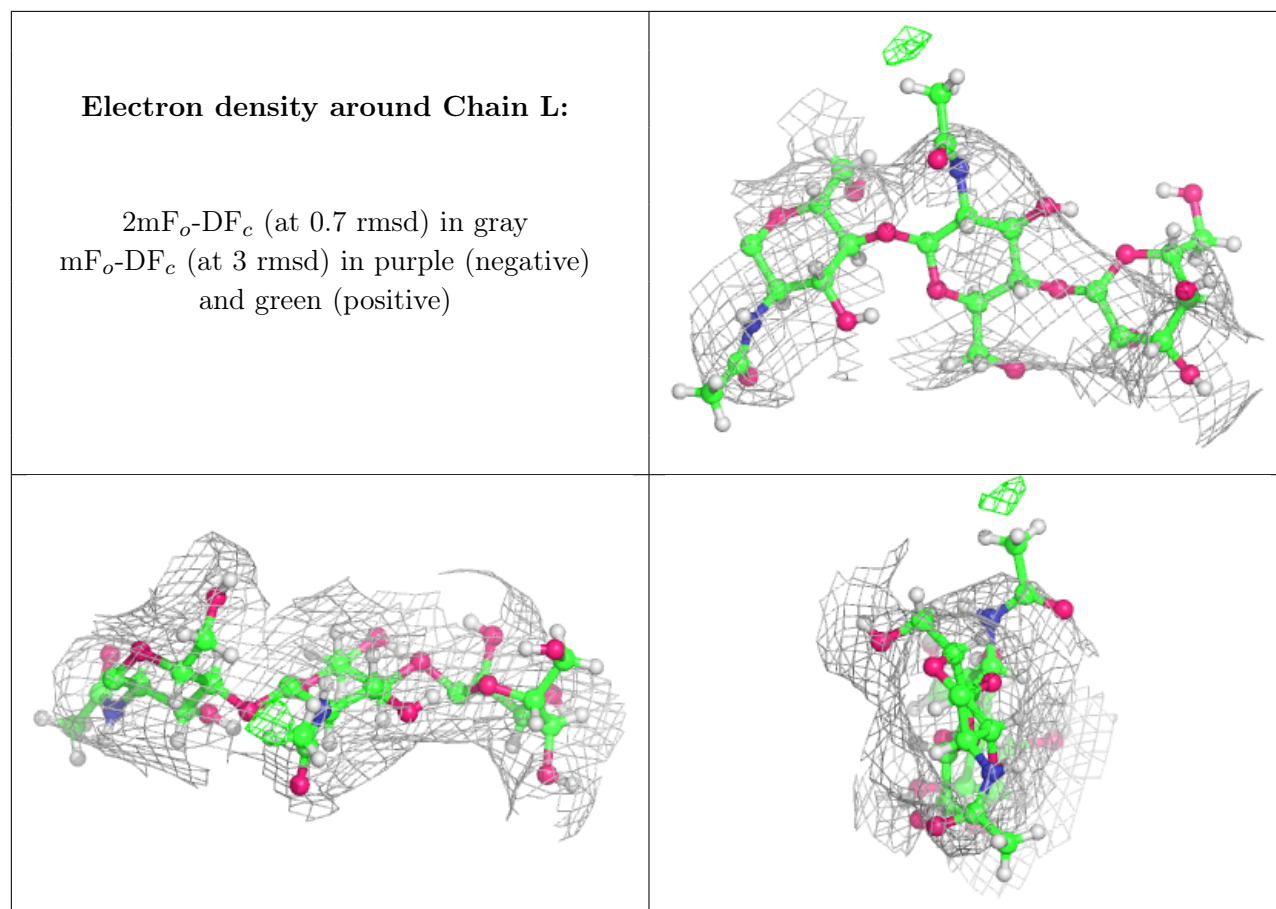
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around Chain K:**

$2mF_o-DF_c$  (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, 95<sup>th</sup> 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.

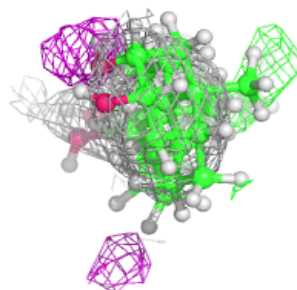
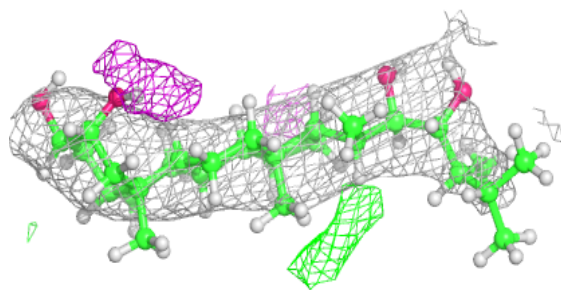
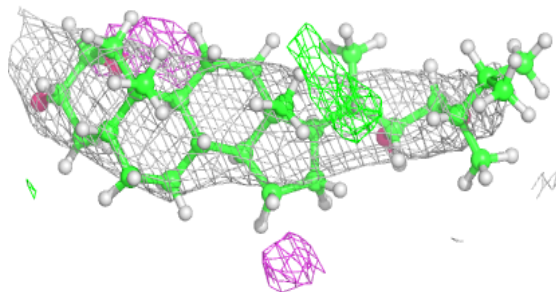
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
7	NAG	A	803	14/15	0.57	0.08	139,169,203,213	0
7	NAG	A	802	14/15	0.60	0.08	154,196,239,239	0
7	NAG	A	805	14/15	0.64	0.08	113,145,186,190	0
7	NAG	B	802	14/15	0.64	0.08	112,149,178,185	0
7	NAG	B	805	14/15	0.67	0.07	153,186,226,230	0
7	NAG	A	804	14/15	0.77	0.07	143,175,214,218	0
7	NAG	B	803	14/15	0.78	0.07	138,168,204,228	0
8	ACT	B	806	4/4	0.79	0.14	93,107,131,131	0
7	NAG	B	804	14/15	0.87	0.06	115,141,175,186	0
6	A1JMI	B	801	32/32	0.92	0.12	75,116,153,164	0
6	A1JMI	A	801	32/32	0.95	0.08	60,106,134,181	0

The following is a graphical depiction of the model fit to experimental electron density of all

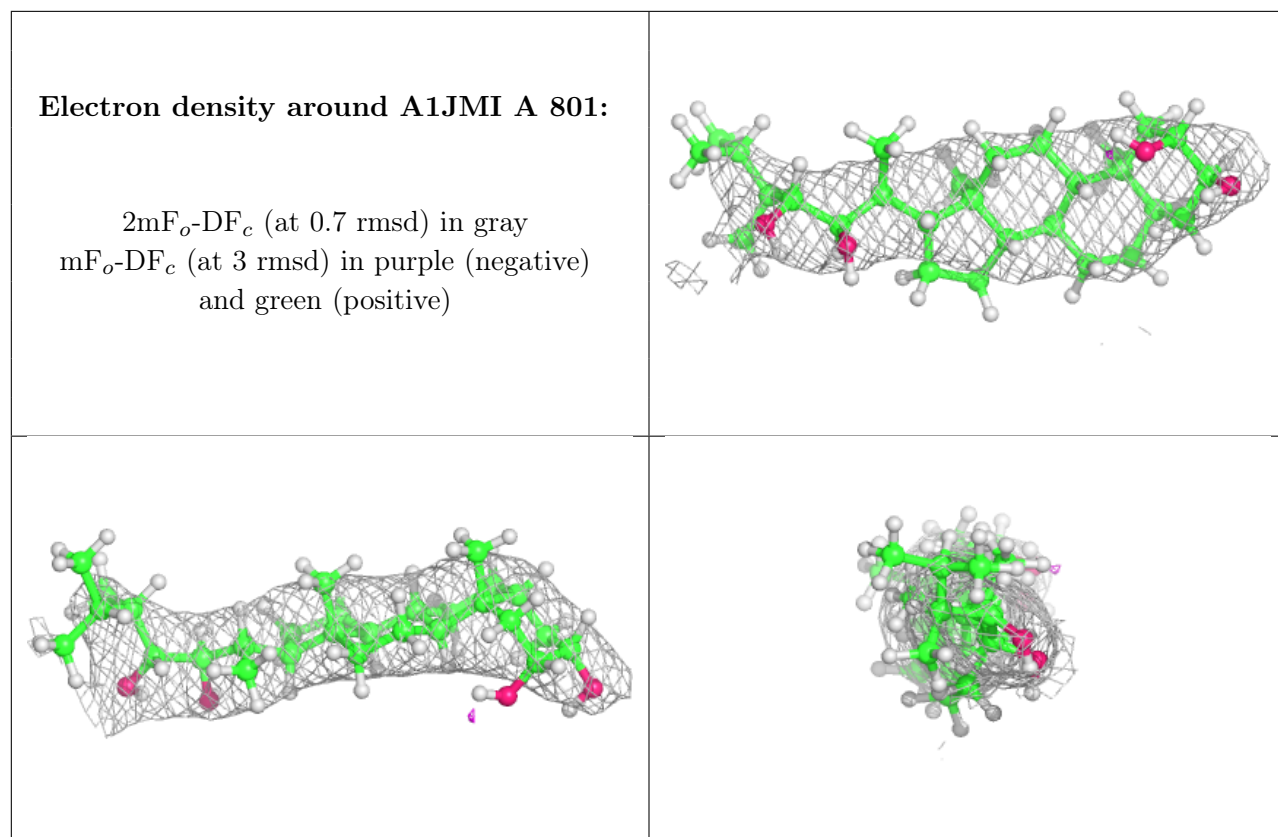
instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around A1JMI B 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.