



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 2, 2025 – 01:17 pm BST

PDB ID : 9S8S / pdb_00009s8s
Title : Crystal structure of the BRI1 ectodomain from Arabidopsis thaliana in complex with typhasterol.
Authors : Caregnato, A.; Hothorn, M.
Deposited on : 2025-08-05
Resolution : 2.65 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 Phenix2.0rc1
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

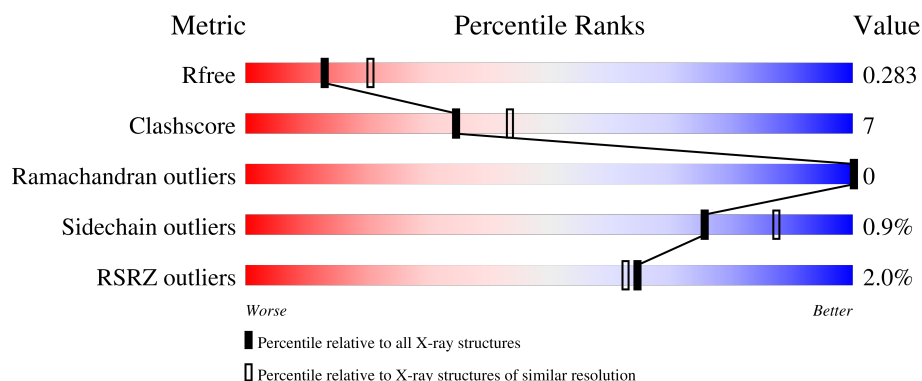
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

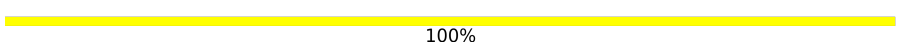

The reported resolution of this entry is 2.65 Å.

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	1003 (2.66-2.66)
Clashscore	180529	1063 (2.66-2.66)
Ramachandran outliers	177936	1052 (2.66-2.66)
Sidechain outliers	177891	1052 (2.66-2.66)
RSRZ outliers	164620	1003 (2.66-2.66)

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	797	
2	B	2	
2	C	2	
2	G	2	
2	H	2	

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Mol	Chain	Length	Quality of chain
3	D	2	 50%50%
4	E	4	 100%
5	F	3	 67%33%
5	I	3	 33%67%

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 11834 atoms, of which 5797 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 Protein BRASSINOSTEROID INSENSITIVE 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	742	Total	C	H	N	O	S	0	0	0
			11205	3549	5584	933	1107	32			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	789	ALA	-	expression tag	UNP O22476
A	790	ALA	-	expression tag	UNP O22476
A	791	ALA	-	expression tag	UNP O22476
A	792	GLU	-	expression tag	UNP O22476
A	793	ASN	-	expression tag	UNP O22476
A	794	LEU	-	expression tag	UNP O22476
A	795	TYR	-	expression tag	UNP O22476
A	796	PHE	-	expression tag	UNP O22476
A	797	GLN	-	expression tag	UNP O22476

- 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	B	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			
2	C	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	N	O		0	0	0
			28	16	2	10				

- Molecule 3 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.



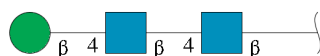
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	2	Total	C	H	N	O	0	0	0
			45	14	21	1	9			

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



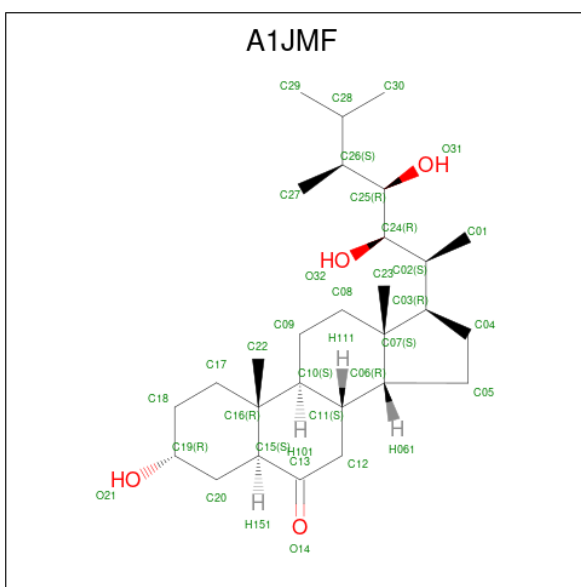
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	4	Total	C	N	O	0	0	0
			50	28	2	20			

- 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.



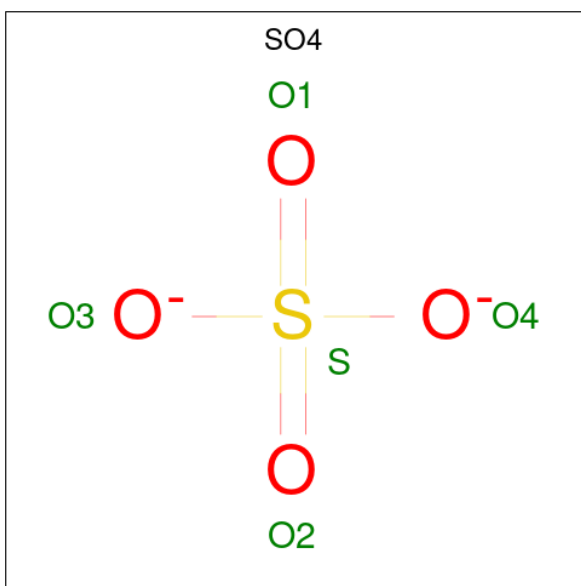
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	3	Total 73	C 22	H 34	N 2	O 15	0	0	0
5	I	3	Total 39	C 22	N 2	O 15		0	0	0

- Molecule 6 is Typhasterol (CCD ID: A1JMF) (formula: C₂₈H₄₈O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			80	28	48	4		

- Molecule 7 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



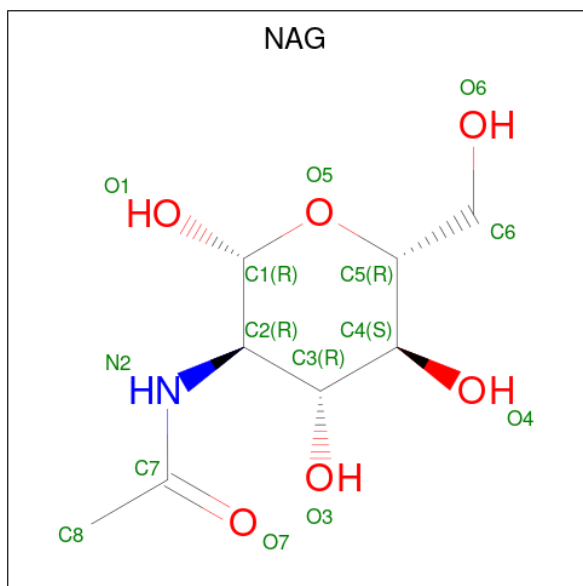
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 8 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



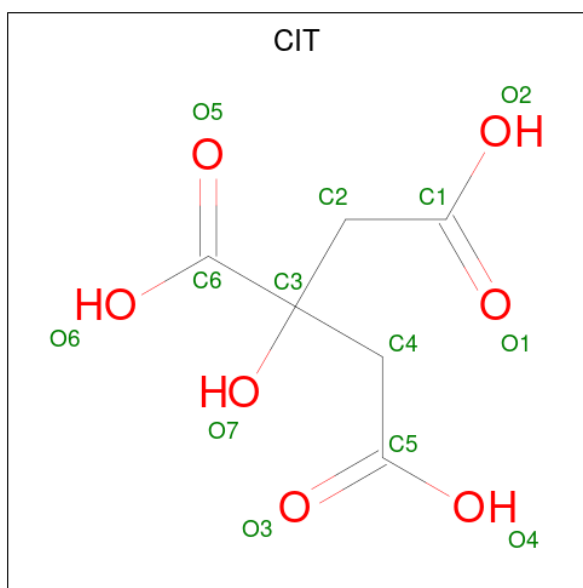
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	A	1	Total	C	H	N	O	0	0
			26	8	12	1	5		
9	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		

- Molecule 10 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	A	1	Total	C	H	O	0	0
			18	6	5	7		

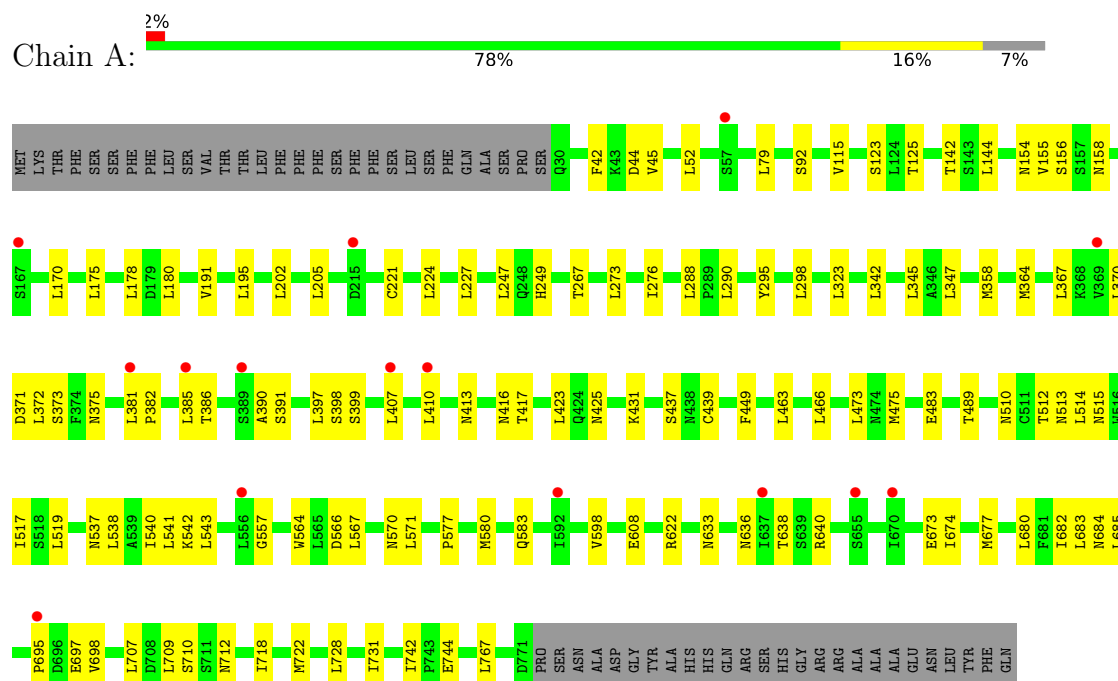
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	65	Total	O	0	0
			65	65		

3 Residue-property plots [i](#)

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

- Molecule 1: Protein BRASSINOSTEROID INSENSITIVE 1



- 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



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

Chain G:  100%

NAG1
NAG2

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

Chain H:  50% 50%

NAG1
NAG2

- Molecule 3: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  50% 50%


NAG1
FUC2

- Molecule 4: 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 E:  100%

NAG1
NAG2
BMA3
MAN4

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

Chain F:  67% 33%

NAG1
NAG2
BMA3

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

Chain I:  33% 67%

NAG1
NAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	174.17Å 66.65Å 119.70Å 90.00° 121.67° 90.00°	Depositor
Resolution (Å)	47.87 – 2.65 47.87 – 2.65	Depositor EDS
% Data completeness (in resolution range)	97.7 (47.87-2.65) 83.5 (47.87-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.44 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.252 , 0.284 0.251 , 0.283	Depositor DCC
R_{free} test set	1670 reflections (2.54%)	wwPDB-VP
Wilson B-factor (Å ²)	70.5	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11834	wwPDB-VP
Average B, all atoms (Å ²)	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.22% 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: A1JMF, FUC, SO4, EDO, CIT, MAN, NAG, BMA

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.11	0/5729	0.29	0/7773

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	5621	5584	5582	79	0
2	B	28	25	25	3	0
2	C	28	25	25	2	0
2	G	28	24	25	1	0
2	H	28	0	25	0	0
3	D	24	21	22	0	0
4	E	50	0	43	3	0
5	F	39	34	34	1	0
5	I	39	0	34	3	0
6	A	32	48	0	0	0
7	A	10	0	0	0	0
8	A	4	6	6	0	0
9	A	28	25	26	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	A	13	5	5	1	0
11	A	65	0	0	2	0
All	All	6037	5797	5852	87	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 (87) 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:413:ASN:O	1:A:416:ASN:ND2	2.13	0.81
1:A:742:ILE:HD11	1:A:767:LEU:HD12	1.64	0.79
1:A:680:LEU:HD21	1:A:683:LEU:HD13	1.69	0.75
2:B:2:NAG:O7	2:B:2:NAG:O3	2.08	0.71
1:A:577:PRO:HG2	1:A:580:MET:HE2	1.71	0.71
1:A:697:GLU:N	1:A:697:GLU:OE1	2.25	0.69
1:A:386:THR:HG22	1:A:410:LEU:HA	1.75	0.68
1:A:483:GLU:N	1:A:483:GLU:OE1	2.26	0.68
1:A:674:ILE:HD11	1:A:683:LEU:HD11	1.75	0.67
1:A:431:LYS:NZ	11:A:901:HOH:O	2.27	0.66
1:A:673:GLU:OE1	1:A:673:GLU:N	2.26	0.65
1:A:370:LEU:HD11	1:A:372:LEU:HD21	1.81	0.62
1:A:684:ASN:C	1:A:685:LEU:HD12	2.28	0.59
1:A:154:ASN:HD21	4:E:1:NAG:C1	2.18	0.56
1:A:728:LEU:HD13	1:A:731:ILE:HD12	1.86	0.56
1:A:685:LEU:HB2	1:A:709:LEU:HD23	1.89	0.55
1:A:674:ILE:HD12	1:A:677:MET:HE1	1.87	0.55
1:A:142:THR:HG23	1:A:170:LEU:O	2.07	0.54
1:A:744:GLU:N	11:A:903:HOH:O	2.40	0.53
2:B:2:NAG:HO3	2:B:2:NAG:C7	2.17	0.52
1:A:707:LEU:HD23	1:A:707:LEU:C	2.35	0.51
1:A:342:LEU:HD21	1:A:345:LEU:HD13	1.93	0.51
1:A:295:TYR:CE2	2:G:1:NAG:H82	2.46	0.51
1:A:636:ASN:O	1:A:640:ARG:NH1	2.44	0.51
1:A:52:LEU:HD21	1:A:79:LEU:HD22	1.92	0.50
1:A:249:HIS:CE1	1:A:273:LEU:HD22	2.47	0.49
1:A:323:LEU:HD12	1:A:347:LEU:CD2	2.42	0.49
1:A:513:ASN:O	1:A:537:ASN:ND2	2.45	0.49
1:A:540:ILE:HG23	1:A:564:TRP:HB3	1.93	0.49
4:E:2:NAG:H61	4:E:3:BMA:C1	2.43	0.49
1:A:515:ASN:HA	1:A:538:LEU:HA	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:541:LEU:CD1	1:A:543:LEU:HD21	2.43	0.48
5:I:2:NAG:O3	5:I:3:BMA:O5	2.28	0.48
1:A:397:LEU:HD12	1:A:423:LEU:CD2	2.43	0.48
1:A:375:ASN:O	1:A:622:ARG:NH2	2.47	0.48
1:A:510:ASN:O	1:A:512:THR:HG23	2.14	0.47
1:A:638:THR:O	1:A:638:THR:HG23	2.14	0.47
1:A:608:GLU:O	5:I:1:NAG:H83	2.15	0.47
1:A:710:SER:O	1:A:712:ASN:ND2	2.48	0.47
1:A:399:SER:O	1:A:622:ARG:NH2	2.47	0.47
1:A:514:LEU:HD13	1:A:517:ILE:HD11	1.98	0.46
1:A:449:PHE:HA	1:A:473:LEU:HD22	1.97	0.46
1:A:466:LEU:O	1:A:489:THR:OG1	2.32	0.45
1:A:267:THR:O	1:A:267:THR:HG22	2.16	0.45
1:A:718:ILE:H	1:A:718:ILE:HD12	1.81	0.45
1:A:191:VAL:O	1:A:195:LEU:HD12	2.17	0.45
1:A:570:ASN:O	1:A:571:LEU:HD23	2.16	0.45
2:B:1:NAG:O3	2:B:2:NAG:O5	2.25	0.45
1:A:510:ASN:HD21	2:C:1:NAG:H2	1.82	0.44
1:A:195:LEU:HD23	1:A:221:CYS:SG	2.58	0.44
4:E:3:BMA:H3	4:E:4:MAN:C5	2.46	0.43
1:A:276:ILE:HD11	1:A:298:LEU:HD22	2.00	0.43
5:F:1:NAG:O4	5:F:2:NAG:H83	2.18	0.43
1:A:636:ASN:HD22	9:A:807:NAG:C7	2.30	0.43
1:A:156:SER:O	1:A:158:ASN:ND2	2.51	0.43
1:A:695:PRO:O	1:A:698:VAL:HG22	2.18	0.43
1:A:42:PHE:O	1:A:45:VAL:HG22	2.19	0.43
1:A:381:LEU:O	1:A:407:LEU:HD11	2.18	0.43
1:A:115:VAL:HG11	1:A:144:LEU:HD13	2.01	0.42
1:A:247:LEU:HD11	1:A:249:HIS:O	2.20	0.42
1:A:364:MET:SD	1:A:367:LEU:HD22	2.58	0.42
1:A:633:ASN:C	5:I:2:NAG:H62	2.44	0.42
1:A:342:LEU:CD2	1:A:345:LEU:HD13	2.49	0.42
1:A:202:LEU:HD21	1:A:205:LEU:HB2	2.01	0.42
1:A:358:MET:HE1	1:A:385:LEU:HB2	2.01	0.42
1:A:382:PRO:O	1:A:407:LEU:HD11	2.20	0.42
1:A:390:ALA:O	1:A:417:THR:HG22	2.19	0.42
1:A:224:LEU:HD21	1:A:227:LEU:HB2	2.01	0.42
2:C:1:NAG:O7	2:C:1:NAG:C1	2.67	0.42
1:A:373:SER:HA	1:A:398:SER:O	2.20	0.41
1:A:175:LEU:HD13	1:A:178:LEU:HD13	2.03	0.41
1:A:370:LEU:CD1	1:A:372:LEU:HD21	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:707:LEU:HD21	1:A:709:LEU:HD12	2.02	0.41
1:A:437:SER:HA	1:A:463:LEU:HD21	2.01	0.41
1:A:557:GLY:O	1:A:583:GLN:HG2	2.19	0.41
1:A:381:LEU:O	1:A:407:LEU:HD21	2.21	0.41
1:A:425:ASN:HA	1:A:449:PHE:HB3	2.01	0.41
1:A:123:SER:O	1:A:125:THR:HG23	2.21	0.41
1:A:510:ASN:O	1:A:512:THR:N	2.54	0.41
1:A:718:ILE:HG23	1:A:722:MET:HE3	2.03	0.41
1:A:155:VAL:CG2	1:A:180:LEU:HD23	2.51	0.40
1:A:566:ASP:C	1:A:567:LEU:HD23	2.45	0.40
1:A:682:ILE:HG22	1:A:683:LEU:N	2.36	0.40
1:A:44:ASP:OD1	10:A:806:CIT:O1	2.38	0.40
1:A:542:LYS:O	1:A:543:LEU:HD23	2.21	0.40
1:A:288:LEU:HB2	1:A:290:LEU:HD21	2.02	0.40
1:A:371:ASP:C	1:A:372:LEU:HD23	2.47	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	740/797 (93%)	668 (90%)	72 (10%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	655/702 (93%)	649 (99%)	6 (1%)	75 87

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	SER
1	A	391	SER
1	A	439	CYS
1	A	475	MET
1	A	519	LEU
1	A	598	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	108	ASN
1	A	223	ASN
1	A	350	ASN
1	A	412	GLN
1	A	416	ASN
1	A	510	ASN
1	A	521	ASN
1	A	570	ASN
1	A	610	HIS
1	A	627	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 ⓘ

20 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	B	1	2,1	14,14,15	0.71	0	17,19,21	1.08	2 (11%)
2	NAG	B	2	2	14,14,15	0.72	0	17,19,21	0.86	0
2	NAG	C	1	2	14,14,15	0.75	0	17,19,21	1.47	3 (17%)
2	NAG	C	2	2	14,14,15	0.77	0	17,19,21	1.02	1 (5%)
3	NAG	D	1	1,3	14,14,15	0.67	0	17,19,21	1.11	1 (5%)
3	FUC	D	2	3	10,10,11	0.84	0	14,14,16	0.95	0
4	NAG	E	1	4	14,14,15	0.87	1 (7%)	17,19,21	1.40	1 (5%)
4	NAG	E	2	4	14,14,15	0.82	1 (7%)	17,19,21	0.95	1 (5%)
4	BMA	E	3	4	11,11,12	0.90	0	15,15,17	2.51	5 (33%)
4	MAN	E	4	4	11,11,12	0.65	0	15,15,17	1.45	1 (6%)
5	NAG	F	1	5,1	14,14,15	0.74	0	17,19,21	0.84	0
5	NAG	F	2	5	14,14,15	0.70	0	17,19,21	1.21	1 (5%)
5	BMA	F	3	5	11,11,12	0.79	0	15,15,17	3.13	8 (53%)
2	NAG	G	1	2,1	14,14,15	0.70	0	17,19,21	0.88	0
2	NAG	G	2	2	14,14,15	0.68	0	17,19,21	0.93	1 (5%)
2	NAG	H	1	2,1	14,14,15	0.94	1 (7%)	17,19,21	1.43	3 (17%)
2	NAG	H	2	2	14,14,15	0.77	0	17,19,21	0.82	0
5	NAG	I	1	5	14,14,15	0.76	0	17,19,21	1.15	1 (5%)
5	NAG	I	2	5	14,14,15	0.73	0	17,19,21	0.79	0
5	BMA	I	3	5	11,11,12	0.84	0	15,15,17	2.90	6 (40%)

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	B	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	3/6/23/26	0/1/1/1
2	NAG	C	1	2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	D	1	1,3	-	1/6/23/26	0/1/1/1
3	FUC	D	2	3	-	-	0/1/1/1
4	NAG	E	1	4	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	1/2/19/22	0/1/1/1
4	MAN	E	4	4	-	2/2/19/22	0/1/1/1
5	NAG	F	1	5,1	-	3/6/23/26	0/1/1/1
5	NAG	F	2	5	-	2/6/23/26	0/1/1/1
5	BMA	F	3	5	-	0/2/19/22	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	NAG	H	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	1/6/23/26	0/1/1/1
5	NAG	I	1	5	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	0/6/23/26	0/1/1/1
5	BMA	I	3	5	-	1/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1	NAG	O5-C1	-2.27	1.40	1.43
2	H	1	NAG	O5-C1	-2.16	1.40	1.43
4	E	2	NAG	O5-C1	-2.14	1.40	1.43

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	3	BMA	C1-O5-C5	8.62	123.87	112.19
5	I	3	BMA	C1-O5-C5	8.47	123.66	112.19
4	E	3	BMA	C1-O5-C5	6.76	121.35	112.19
4	E	4	MAN	C1-O5-C5	4.56	118.37	112.19
4	E	1	NAG	O5-C1-C2	-4.20	104.66	111.29
4	E	3	BMA	C3-C4-C5	4.17	117.69	110.24
5	F	3	BMA	C3-C4-C5	4.16	117.66	110.24
5	F	3	BMA	C2-C3-C4	4.02	117.86	110.89
5	I	3	BMA	C3-C4-C5	3.95	117.28	110.24
5	F	2	NAG	C2-N2-C7	3.55	127.95	122.90
5	F	3	BMA	C1-C2-C3	3.37	113.81	109.67
5	I	1	NAG	O5-C1-C2	-3.26	106.14	111.29
2	C	1	NAG	C2-N2-C7	3.15	127.39	122.90
2	B	1	NAG	C2-N2-C7	3.12	127.35	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	3	BMA	C2-C3-C4	2.99	116.06	110.89
4	E	3	BMA	C2-C3-C4	2.96	116.02	110.89
2	C	1	NAG	C4-C3-C2	-2.94	106.71	111.02
4	E	3	BMA	O4-C4-C3	-2.59	104.36	110.35
2	C	2	NAG	C2-N2-C7	2.59	126.59	122.90
5	I	3	BMA	O3-C3-C2	-2.54	105.12	109.99
5	F	3	BMA	O3-C3-C2	-2.54	105.13	109.99
5	I	3	BMA	O5-C5-C4	2.53	116.97	110.83
4	E	3	BMA	O5-C5-C4	2.47	116.83	110.83
5	F	3	BMA	O5-C5-C4	2.45	116.78	110.83
5	F	3	BMA	O4-C4-C3	-2.40	104.80	110.35
2	H	1	NAG	O5-C1-C2	-2.39	107.51	111.29
5	I	3	BMA	O4-C4-C3	-2.34	104.95	110.35
2	G	2	NAG	O5-C1-C2	-2.34	107.60	111.29
2	H	1	NAG	C1-O5-C5	-2.33	109.04	112.19
5	F	3	BMA	O5-C1-C2	2.31	114.33	110.77
4	E	2	NAG	O4-C4-C3	-2.30	105.03	110.35
2	C	1	NAG	C1-C2-N2	2.26	114.34	110.49
2	H	1	NAG	C6-C5-C4	2.23	118.22	113.00
2	B	1	NAG	O5-C1-C2	-2.12	107.94	111.29
3	D	1	NAG	C2-N2-C7	2.05	125.81	122.90

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C1-C2-N2-C7
4	E	4	MAN	O5-C5-C6-O6
2	B	1	NAG	C8-C7-N2-C2
2	B	1	NAG	O7-C7-N2-C2
4	E	1	NAG	C8-C7-N2-C2
4	E	1	NAG	O7-C7-N2-C2
5	F	2	NAG	C8-C7-N2-C2
5	F	2	NAG	O7-C7-N2-C2
2	B	2	NAG	C1-C2-N2-C7
4	E	4	MAN	C4-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
2	B	2	NAG	O5-C5-C6-O6
2	B	2	NAG	C3-C2-N2-C7
2	C	2	NAG	C3-C2-N2-C7
3	D	1	NAG	C3-C2-N2-C7
4	E	3	BMA	C4-C5-C6-O6

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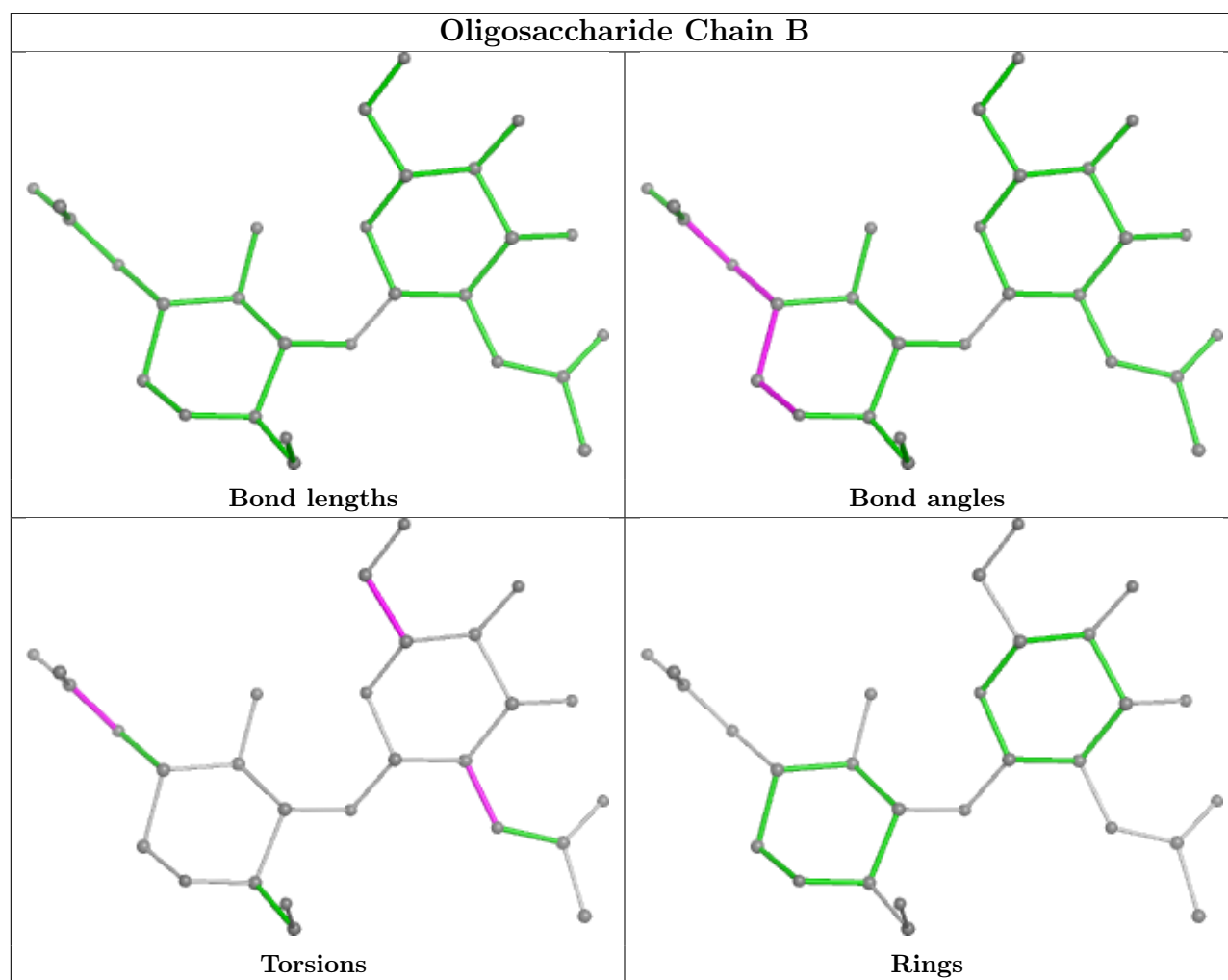
Mol	Chain	Res	Type	Atoms
5	F	1	NAG	C3-C2-N2-C7
5	F	1	NAG	O5-C5-C6-O6
5	I	3	BMA	C4-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
5	F	1	NAG	C1-C2-N2-C7

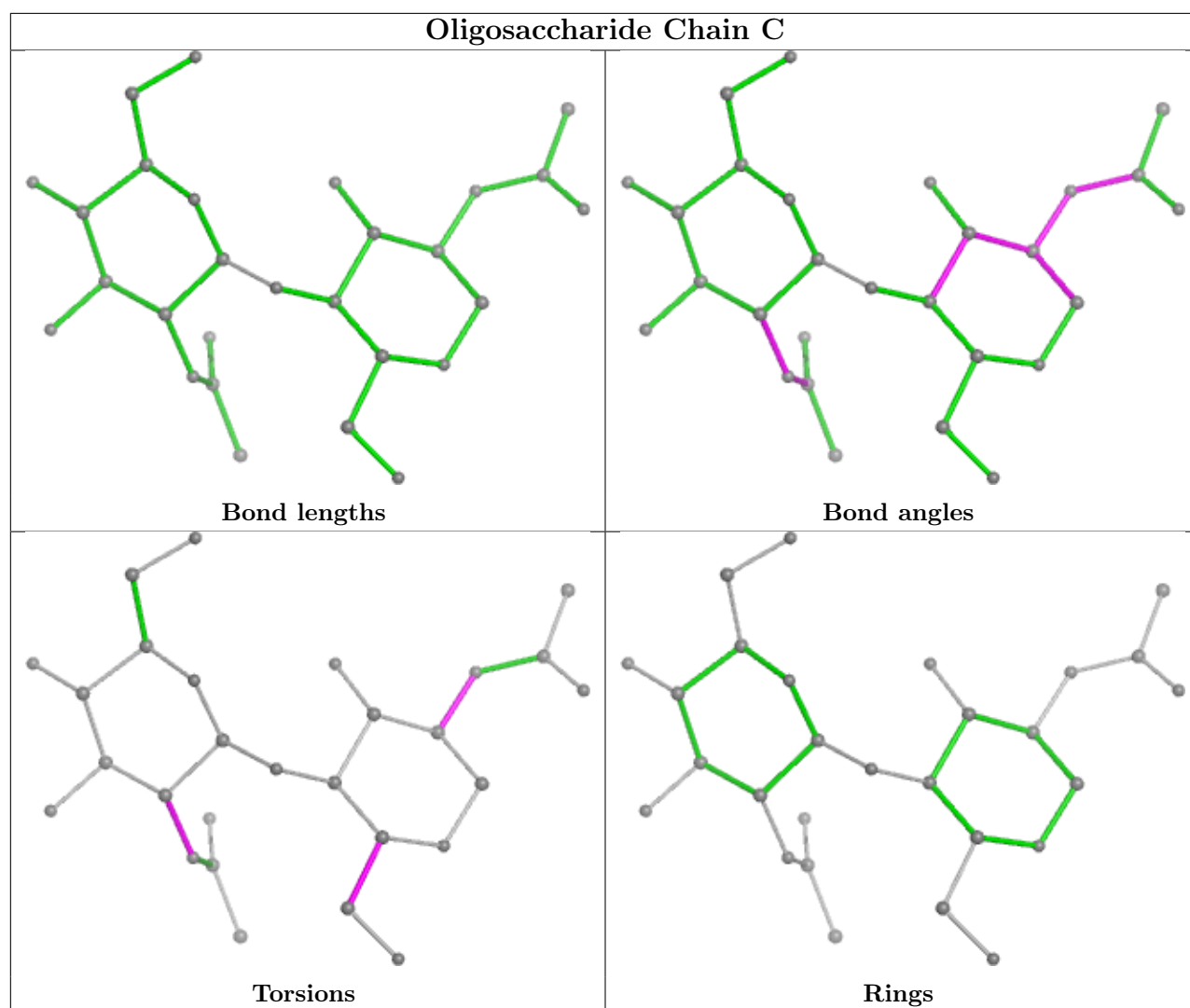
There are no ring outliers.

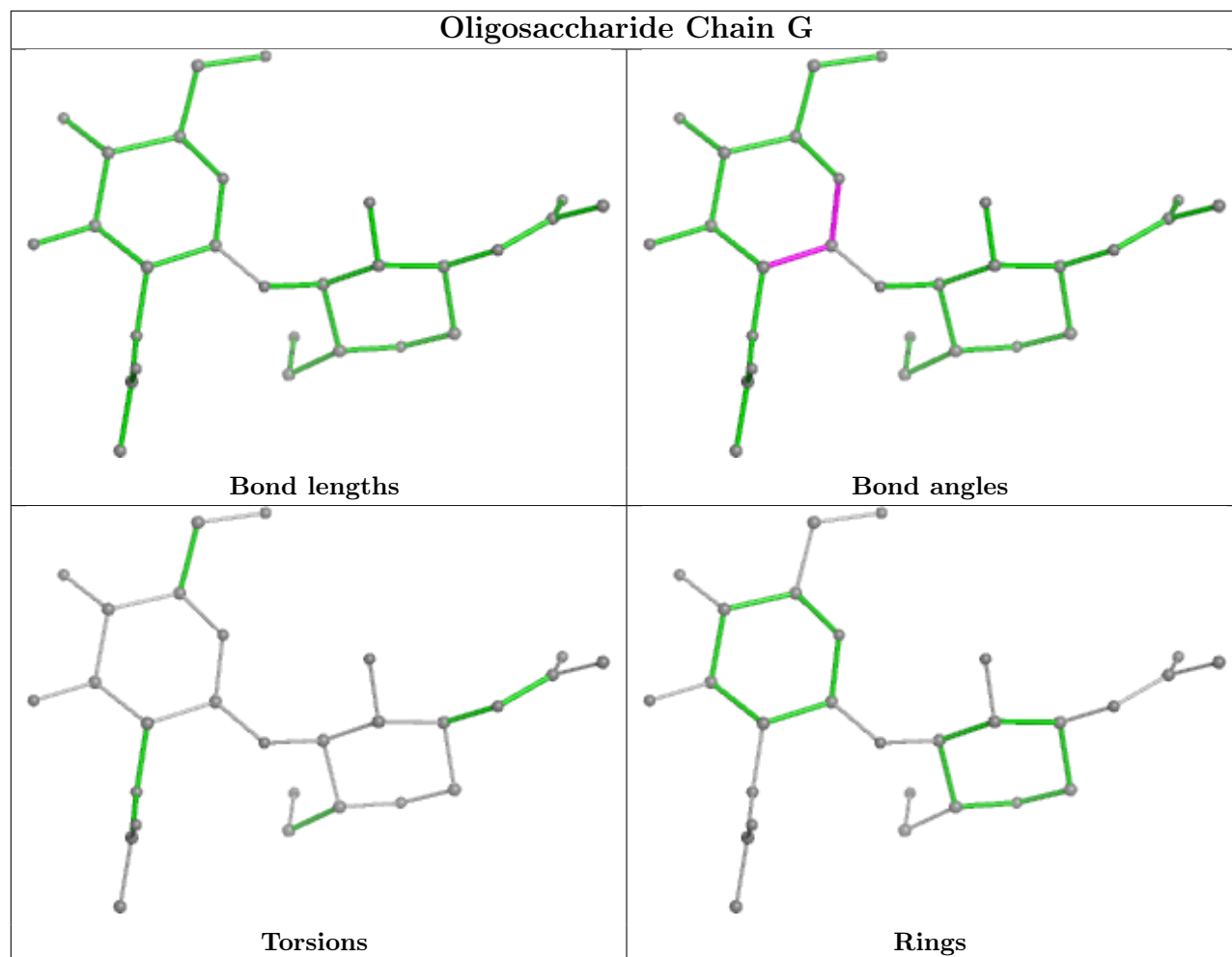
13 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	4	MAN	1	0
2	B	2	NAG	3	0
4	E	2	NAG	1	0
2	C	1	NAG	2	0
2	B	1	NAG	1	0
4	E	1	NAG	1	0
5	I	1	NAG	1	0
4	E	3	BMA	2	0
5	I	3	BMA	1	0
5	F	2	NAG	1	0
5	F	1	NAG	1	0
2	G	1	NAG	1	0
5	I	2	NAG	2	0

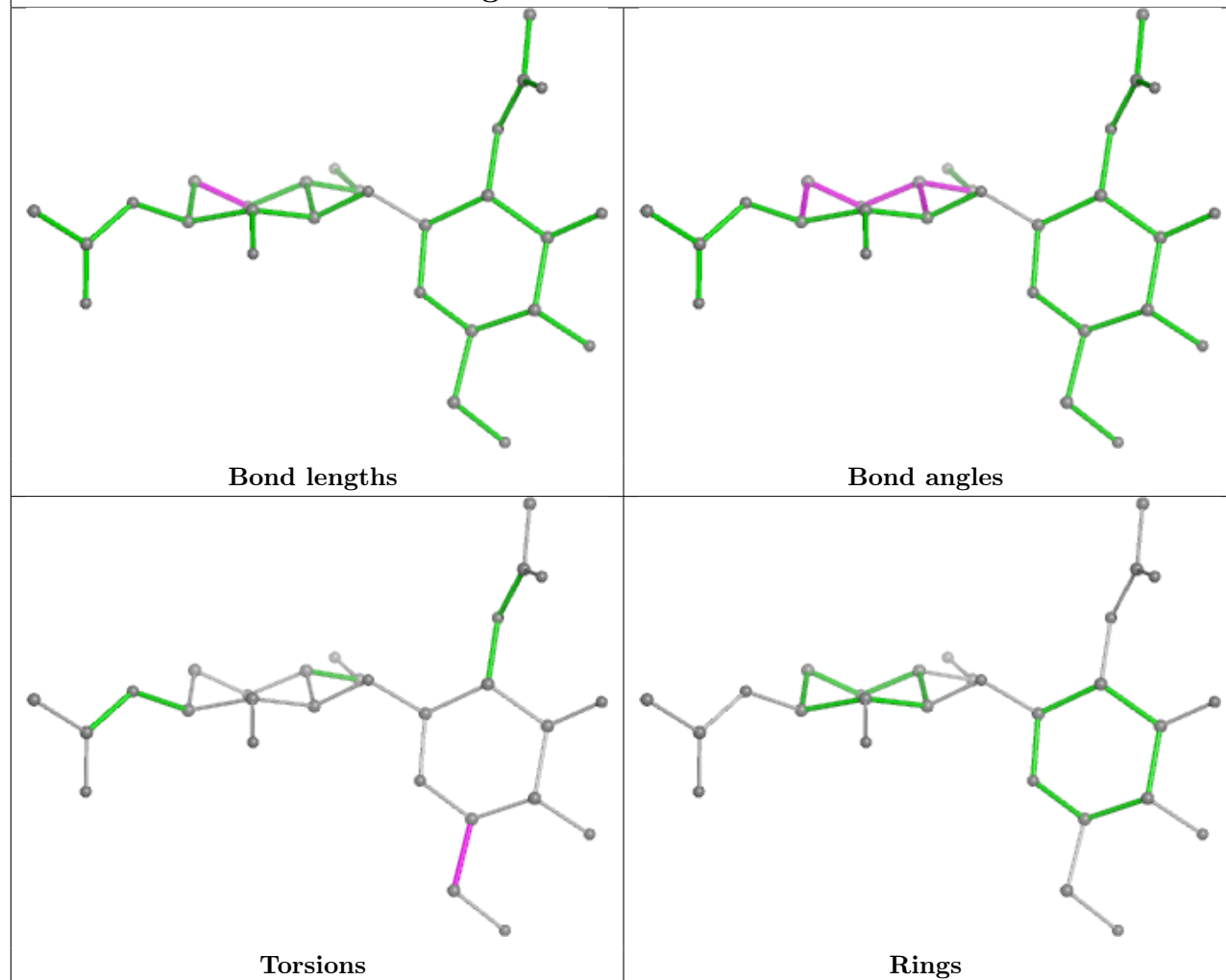
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

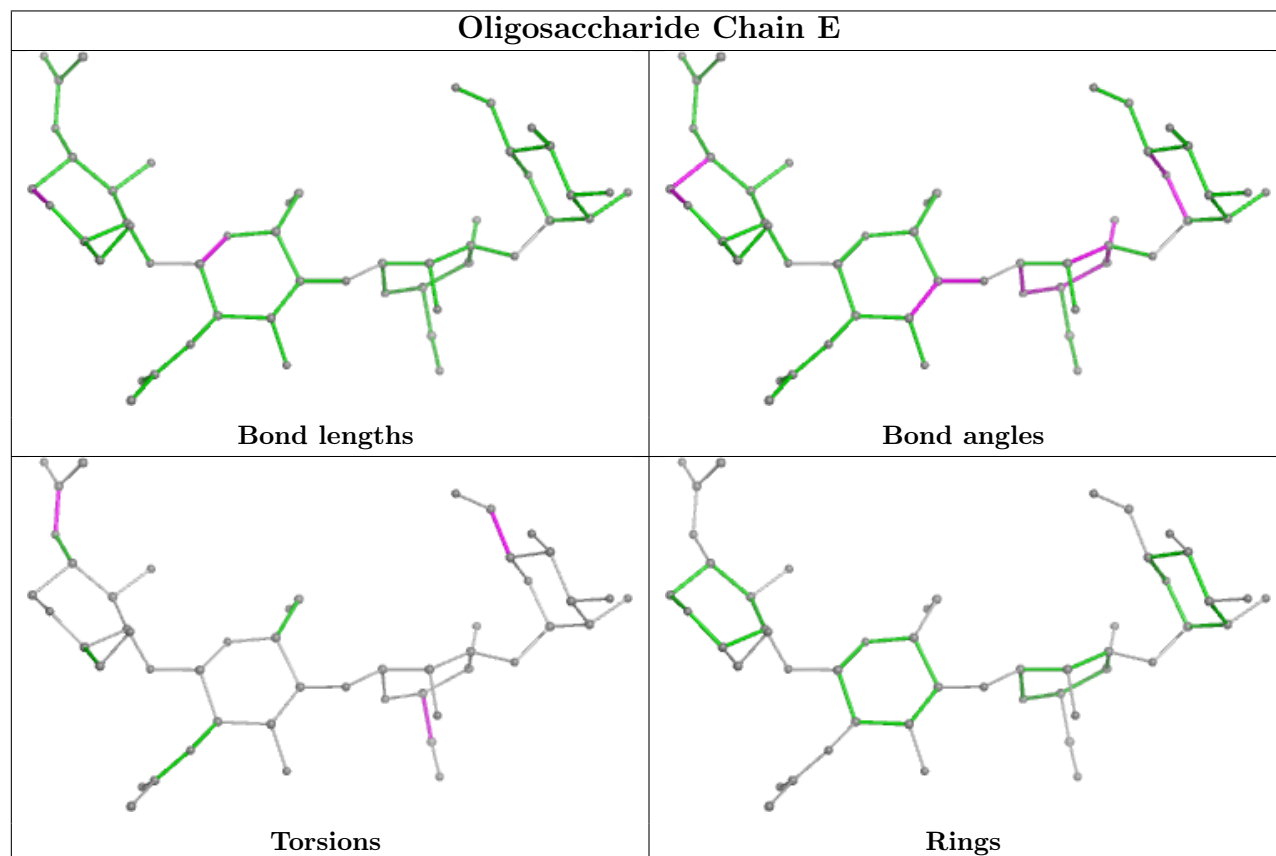
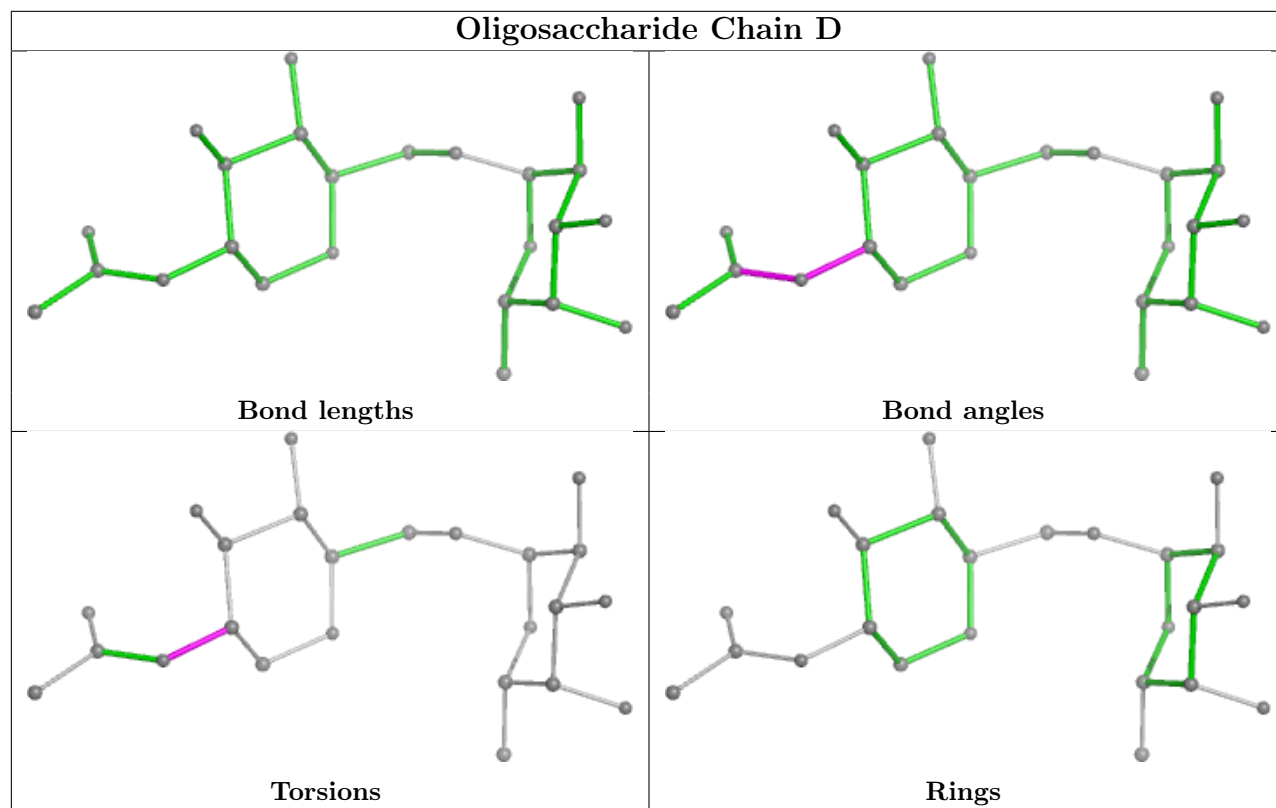


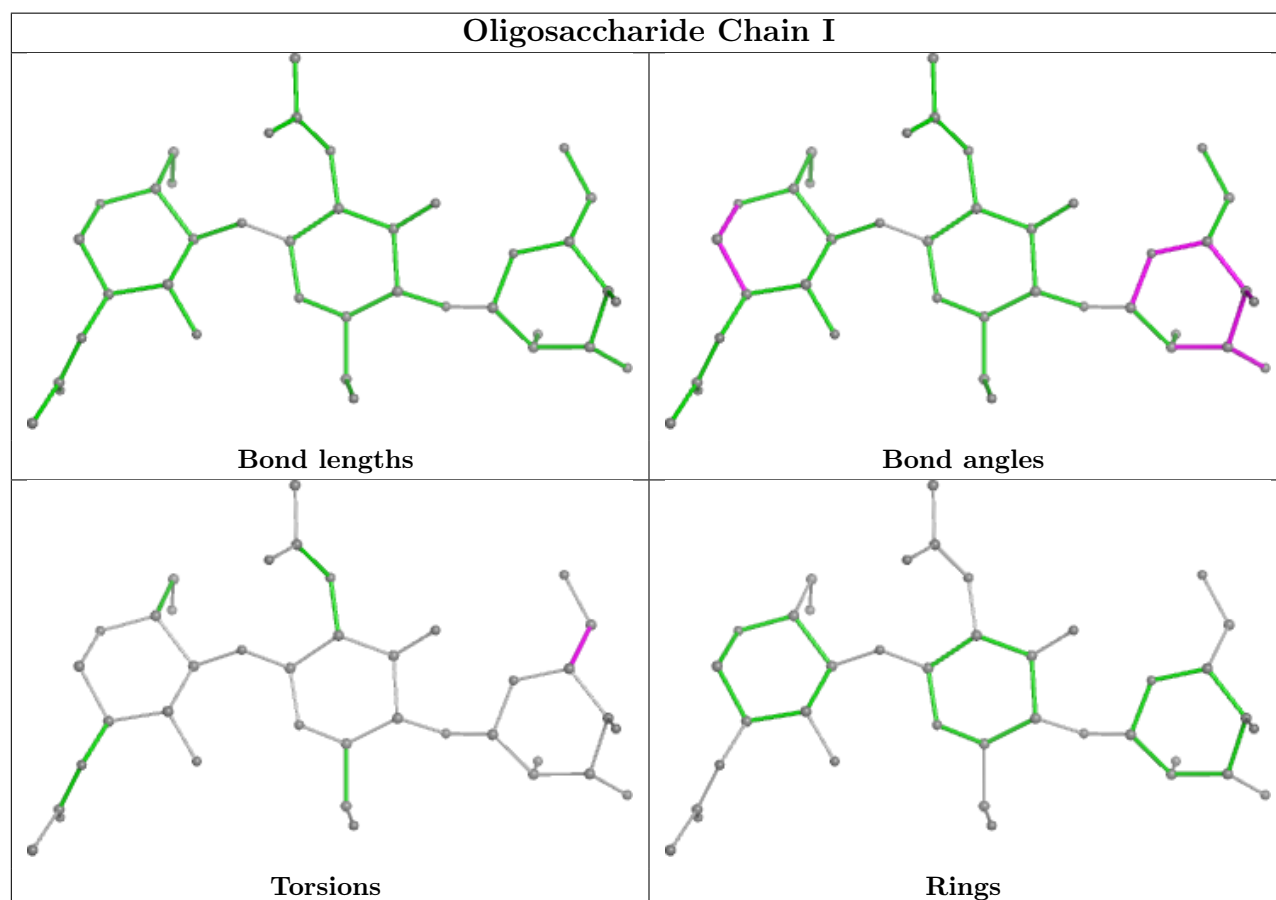
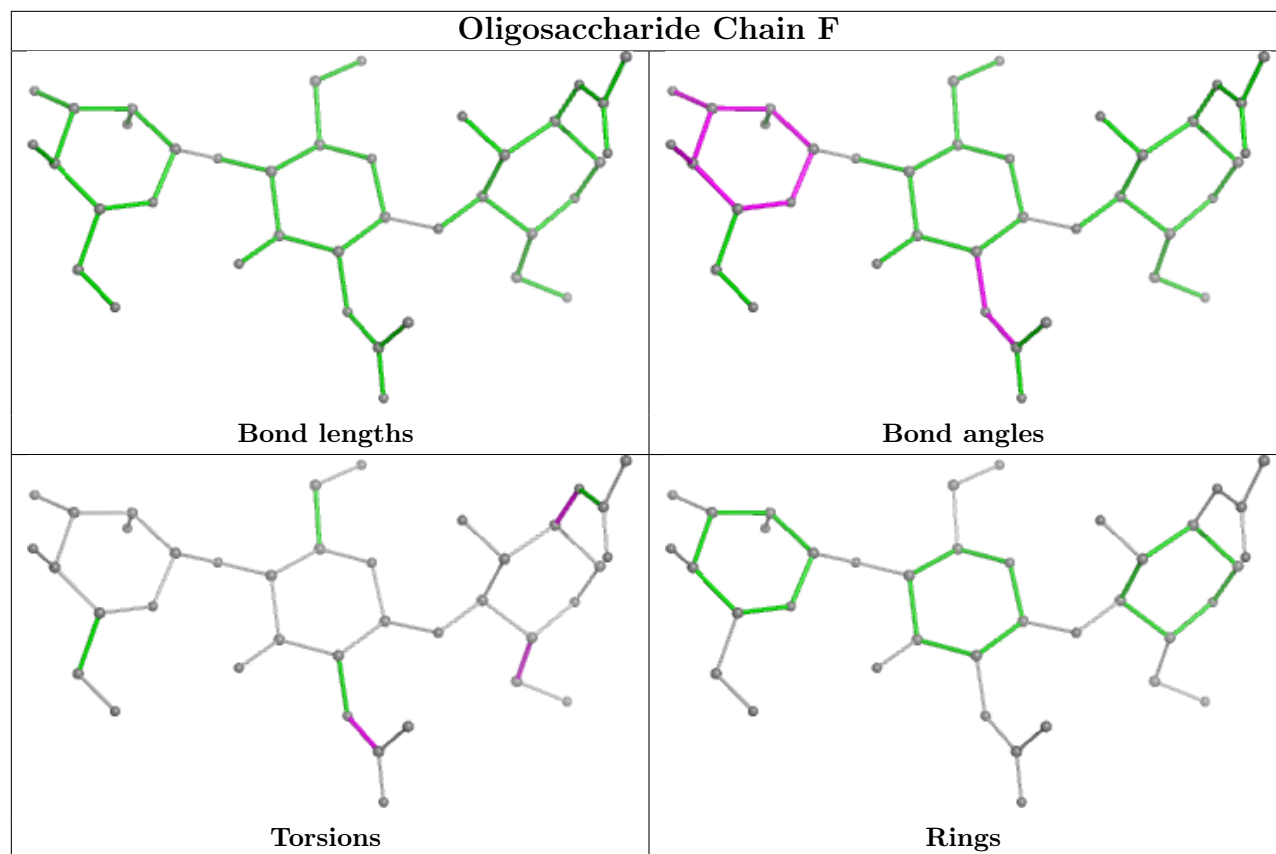




Oligosaccharide Chain H







5.6 Ligand geometry

7 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	$\# Z > 2$
9	NAG	A	807	1	14,14,15	0.68	0	17,19,21	0.87	0
9	NAG	A	805	-	14,14,15	0.72	0	17,19,21	0.98	0
7	SO4	A	802	-	4,4,4	0.59	0	6,6,6	0.09	0
8	EDO	A	804	-	3,3,3	0.28	0	2,2,2	0.14	0
10	CIT	A	806	-	12,12,12	1.27	1 (8%)	17,17,17	1.79	5 (29%)
7	SO4	A	803	-	4,4,4	0.60	0	6,6,6	0.08	0
6	A1JMF	A	801	-	35,35,35	2.00	9 (25%)	55,56,56	3.21	23 (41%)

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
9	NAG	A	807	1	-	1/6/23/26	0/1/1/1
9	NAG	A	805	-	-	2/6/23/26	0/1/1/1
8	EDO	A	804	-	-	0/1/1/1	-
10	CIT	A	806	-	-	7/16/16/16	-
6	A1JMF	A	801	-	-	2/20/81/81	0/4/4/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	801	A1JMF	C11-C10	5.14	1.63	1.53
6	A	801	A1JMF	C12-C11	4.40	1.61	1.53
6	A	801	A1JMF	C23-C07	-4.13	1.47	1.54
6	A	801	A1JMF	C08-C09	-3.32	1.46	1.53
6	A	801	A1JMF	C20-C19	3.07	1.57	1.51
6	A	801	A1JMF	C25-C24	3.00	1.59	1.53
10	A	806	CIT	C3-C6	2.69	1.56	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	801	A1JMF	C05-C06	-2.58	1.49	1.54
6	A	801	A1JMF	C08-C07	2.22	1.58	1.54
6	A	801	A1JMF	C15-C13	2.01	1.55	1.51

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	801	A1JMF	C08-C07-C03	-12.39	98.03	116.57
6	A	801	A1JMF	C05-C06-C11	-8.52	105.04	119.08
6	A	801	A1JMF	C23-C07-C08	8.23	123.58	110.59
6	A	801	A1JMF	C06-C07-C03	6.93	108.28	100.07
6	A	801	A1JMF	C23-C07-C06	-5.52	101.42	111.71
10	A	806	CIT	O5-C6-C3	-4.73	115.55	122.25
6	A	801	A1JMF	C08-C07-C06	4.47	114.20	107.27
6	A	801	A1JMF	C16-C15-C13	4.31	116.38	110.11
6	A	801	A1JMF	C01-C02-C03	-4.23	105.24	112.74
6	A	801	A1JMF	C17-C16-C15	3.66	112.08	107.46
6	A	801	A1JMF	C22-C16-C10	-3.56	106.27	111.18
6	A	801	A1JMF	C16-C10-C11	-3.48	108.77	112.42
6	A	801	A1JMF	C12-C11-C06	2.94	114.77	111.18
6	A	801	A1JMF	C07-C06-C11	2.84	118.59	114.38
6	A	801	A1JMF	C05-C04-C03	2.79	110.66	105.13
10	A	806	CIT	O6-C6-C3	2.76	117.84	113.05
6	A	801	A1JMF	C04-C03-C07	-2.74	100.54	103.84
6	A	801	A1JMF	C20-C15-C16	-2.60	109.23	112.80
10	A	806	CIT	O7-C3-C6	2.58	112.49	108.86
6	A	801	A1JMF	C20-C15-C13	2.56	116.53	113.23
6	A	801	A1JMF	C10-C11-C06	2.34	112.22	109.09
6	A	801	A1JMF	O14-C13-C15	2.30	125.57	122.44
6	A	801	A1JMF	O32-C24-C25	2.29	114.80	109.47
10	A	806	CIT	C2-C3-C6	-2.19	105.41	110.11
10	A	806	CIT	O1-C1-C2	-2.17	116.59	122.94
6	A	801	A1JMF	C03-C02-C24	2.11	115.26	111.51
6	A	801	A1JMF	O14-C13-C12	-2.07	118.61	122.08
6	A	801	A1JMF	C11-C12-C13	-2.06	107.40	111.56

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	A	806	CIT	C2-C3-C6-O6
9	A	807	NAG	O5-C5-C6-O6

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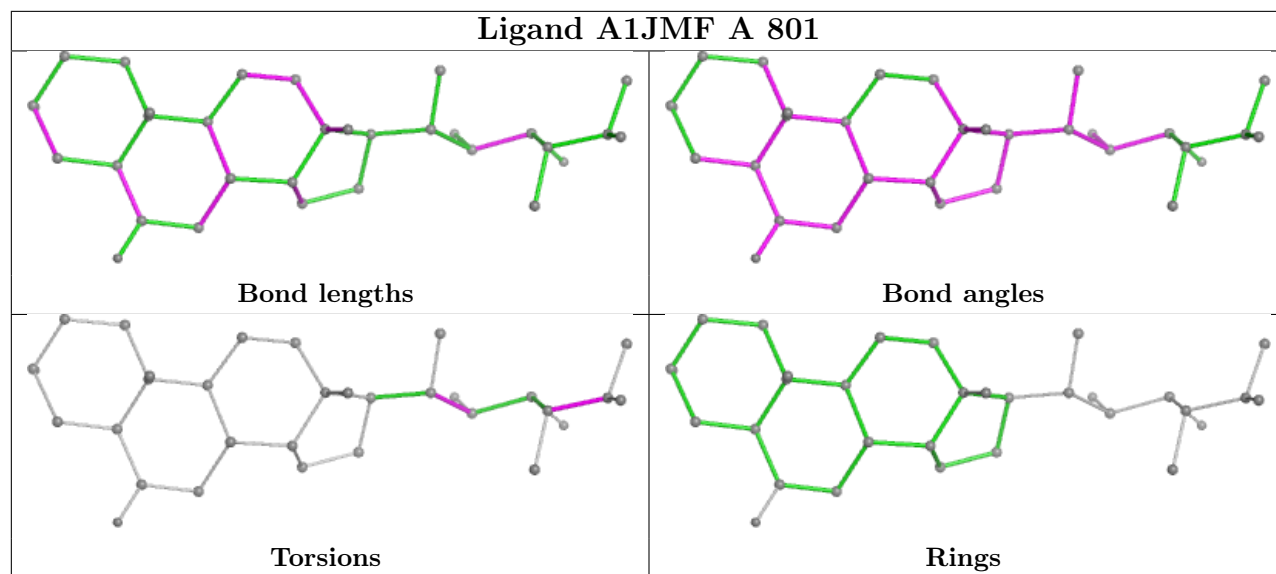
Mol	Chain	Res	Type	Atoms
10	A	806	CIT	C4-C3-C6-O5
10	A	806	CIT	C4-C3-C6-O6
10	A	806	CIT	C2-C3-C6-O5
10	A	806	CIT	C1-C2-C3-C4
9	A	805	NAG	C3-C2-N2-C7
9	A	805	NAG	C1-C2-N2-C7
10	A	806	CIT	O7-C3-C6-O5
10	A	806	CIT	O7-C3-C6-O6
6	A	801	A1JMF	C25-C26-C28-C29
6	A	801	A1JMF	C03-C02-C24-O32

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	807	NAG	1	0
10	A	806	CIT	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, 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	742/797 (93%)	0.18	15 (2%) 64 62	71, 111, 163, 203	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	637	ILE	3.0
1	A	592	ILE	2.9
1	A	655	SER	2.9
1	A	695	PRO	2.5
1	A	369	VAL	2.5
1	A	381	LEU	2.4
1	A	410	LEU	2.4
1	A	385	LEU	2.2
1	A	167	SER	2.2
1	A	670	ILE	2.2
1	A	407	LEU	2.1
1	A	556	LEU	2.1
1	A	215	ASP	2.1
1	A	57	SER	2.0
1	A	389	SER	2.0

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

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

6.3 Carbohydrates [i](#)

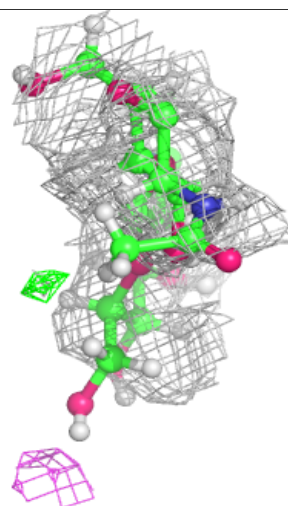
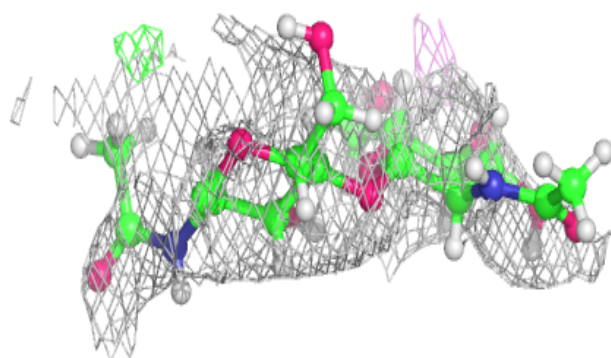
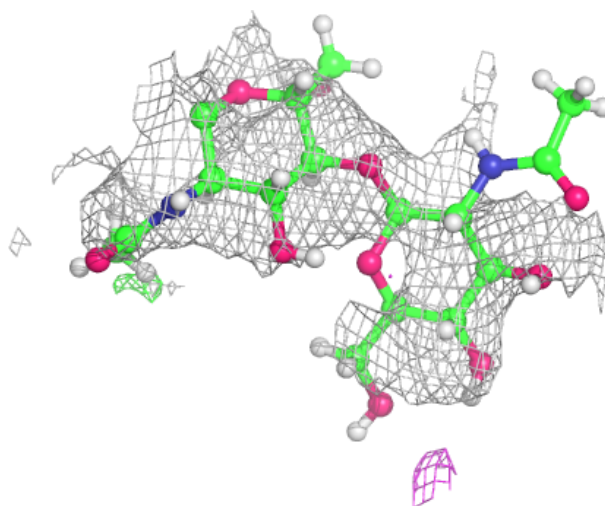
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	BMA	F	3	11/12	0.14	0.13	155,160,195,197	0
2	NAG	B	1	14/15	0.38	0.11	140,169,205,213	0
4	MAN	E	4	11/12	0.39	0.11	118,120,122,124	0
2	NAG	B	2	14/15	0.41	0.14	158,166,202,203	0
2	NAG	C	2	14/15	0.43	0.12	152,170,207,208	0
5	NAG	F	2	14/15	0.46	0.12	149,161,196,203	0
2	NAG	H	1	14/15	0.50	0.17	155,167,183,194	0
2	NAG	H	2	14/15	0.51	0.15	187,194,199,218	0
4	BMA	E	3	11/12	0.53	0.14	104,110,114,115	0
3	NAG	D	1	14/15	0.60	0.12	120,129,158,163	0
5	BMA	I	3	11/12	0.62	0.11	108,111,112,113	0
5	NAG	F	1	14/15	0.64	0.14	118,143,162,179	0
2	NAG	G	2	14/15	0.70	0.13	92,103,126,127	0
2	NAG	C	1	14/15	0.70	0.11	125,146,175,178	0
3	FUC	D	2	10/11	0.81	0.12	133,144,173,175	0
5	NAG	I	1	14/15	0.86	0.14	103,106,109,112	0
5	NAG	I	2	14/15	0.86	0.12	106,108,113,116	0
4	NAG	E	2	14/15	0.86	0.11	79,85,93,99	0
4	NAG	E	1	14/15	0.89	0.13	69,72,77,79	0
2	NAG	G	1	14/15	0.91	0.09	76,85,103,105	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.

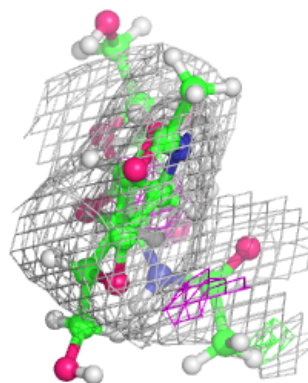
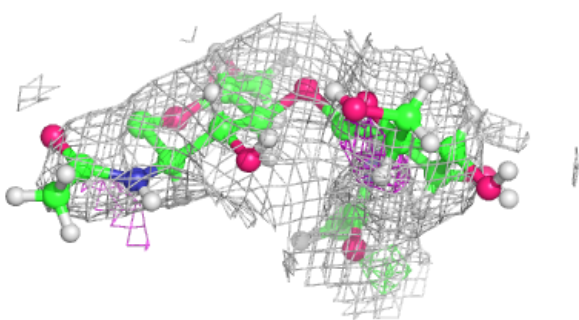
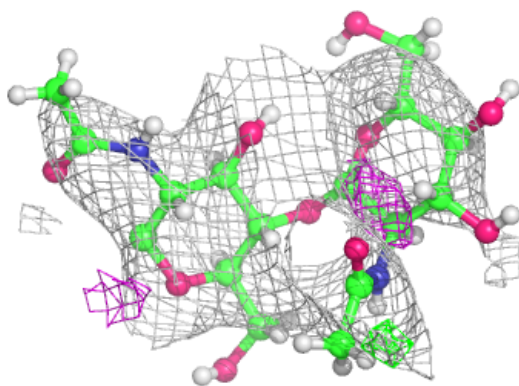
Electron density around Chain B:

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

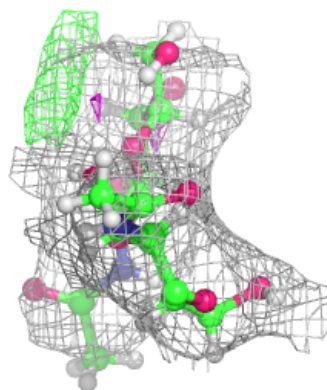
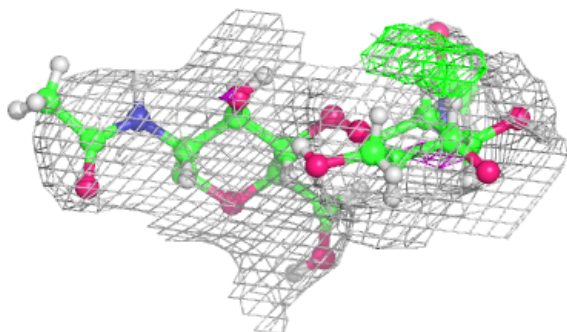
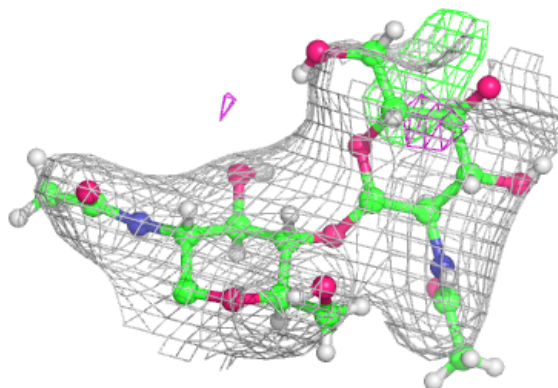


Electron density around Chain C:

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

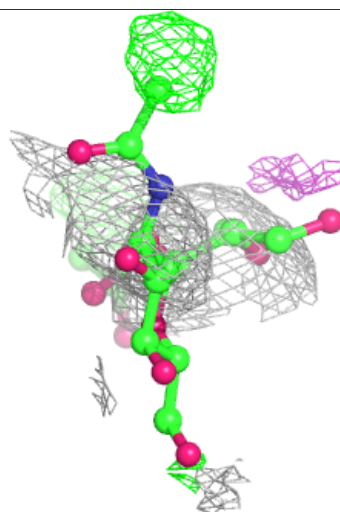
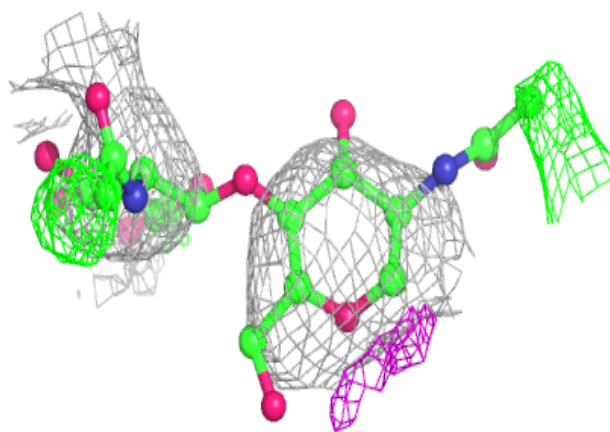
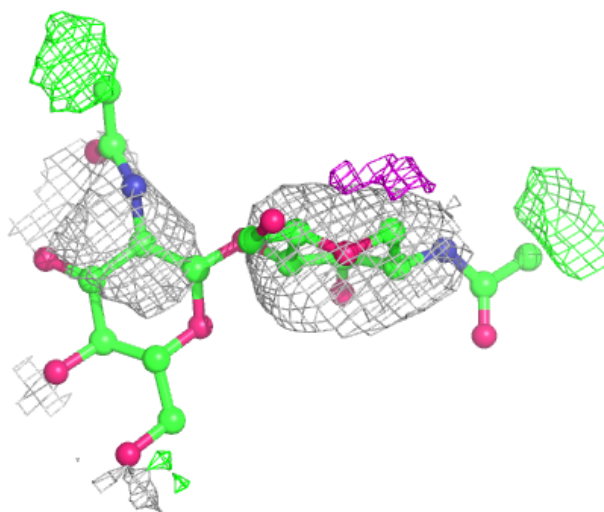
**Electron density around Chain G:**

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



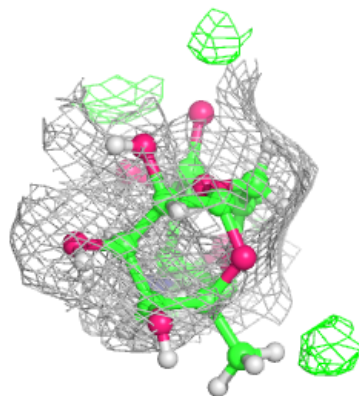
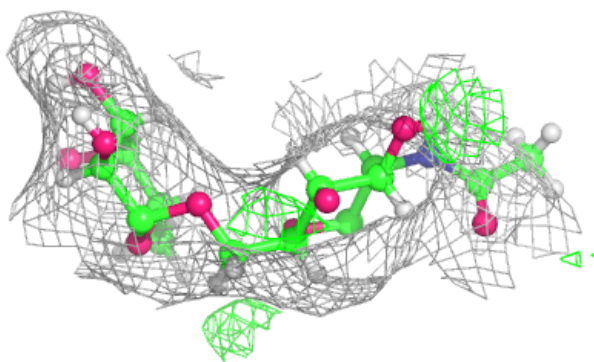
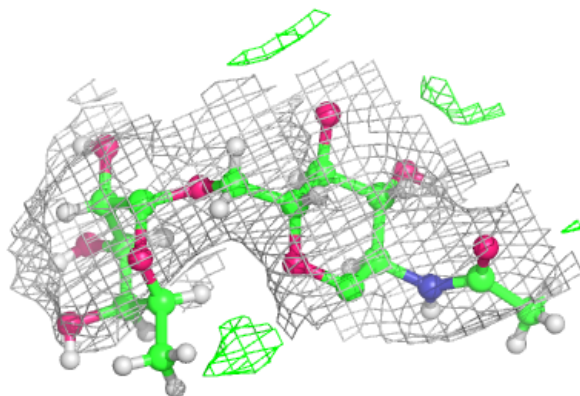
Electron density around Chain H:

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

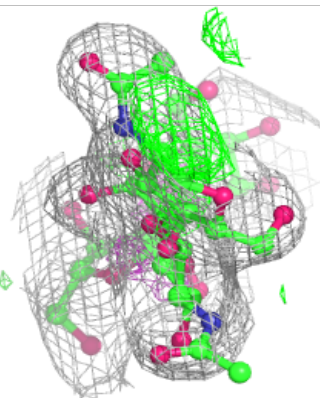
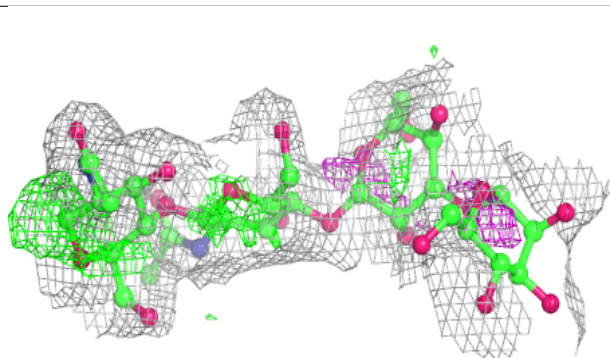
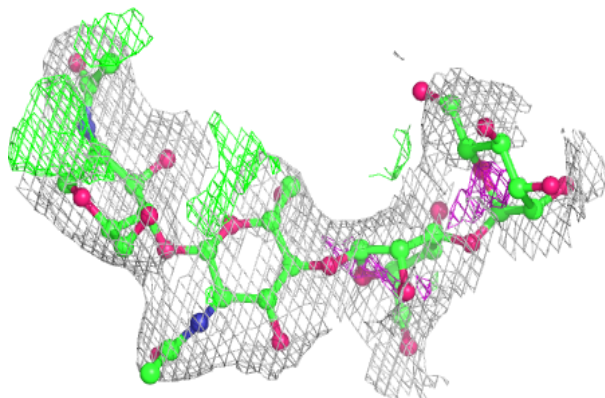


Electron density around Chain D:

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

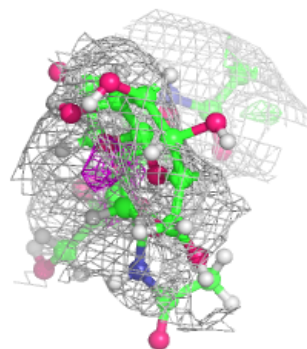
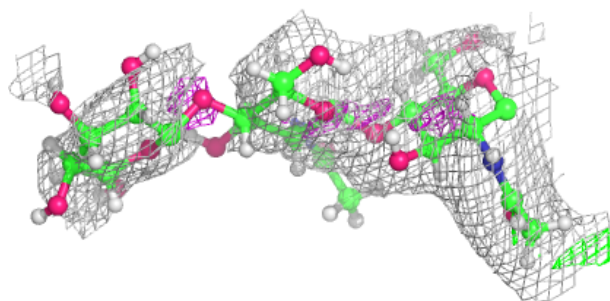
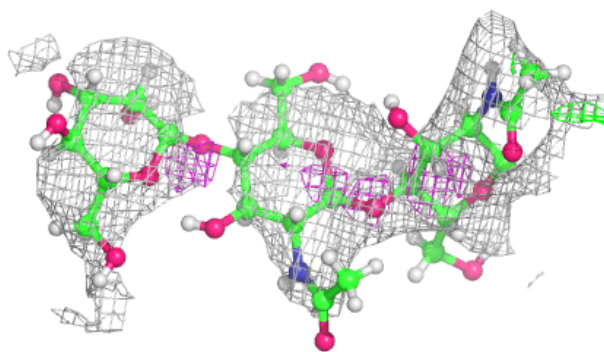
**Electron density around Chain E:**

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

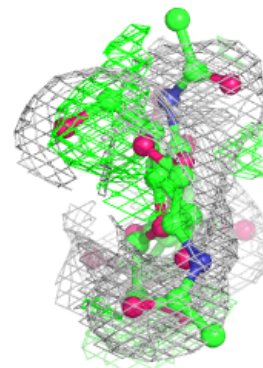
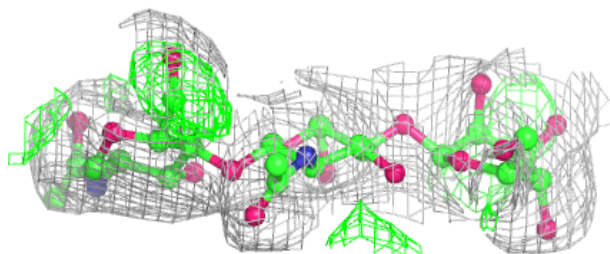
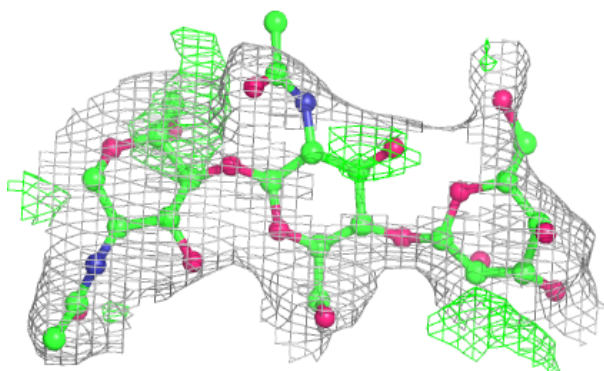


Electron density around Chain F:

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

**Electron density around Chain I:**

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



6.4 Ligands [i](#)

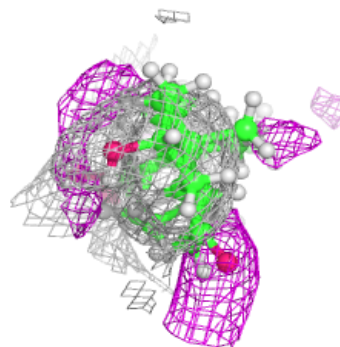
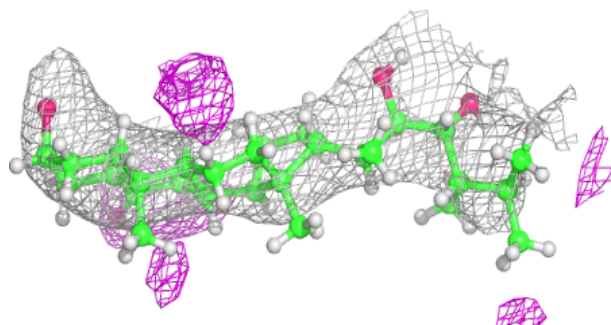
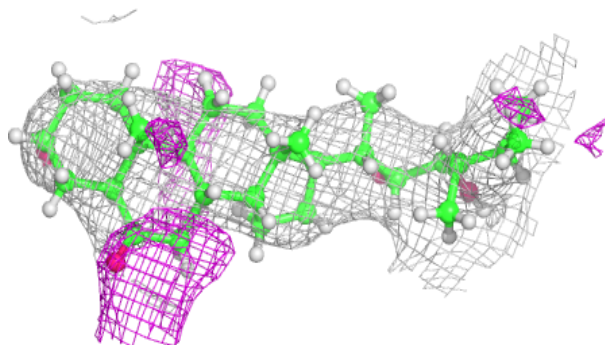
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
9	NAG	A	807	14/15	0.23	0.13	161,167,203,217	0
7	SO4	A	803	5/5	0.56	0.11	107,109,120,122	0
10	CIT	A	806	13/13	0.57	0.17	96,98,119,126	0
9	NAG	A	805	14/15	0.60	0.13	93,104,127,128	0
7	SO4	A	802	5/5	0.77	0.10	112,116,119,120	0
8	EDO	A	804	4/4	0.91	0.12	86,105,108,110	0
6	A1JMF	A	801	32/32	0.93	0.14	88,113,124,128	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 A1JMF 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.