



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 05:19 PM UTC

PDB ID : 6IBT / pdb_00006ibt
Title : Crystal structure of human alpha-galactosidase A in complex with alpha-galactose configured cyclophellitol aziridine ME737
Authors : Rowland, R.J.; Wu, L.; Davies, G.J.
Deposited on : 2018-11-30
Resolution : 2.04 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

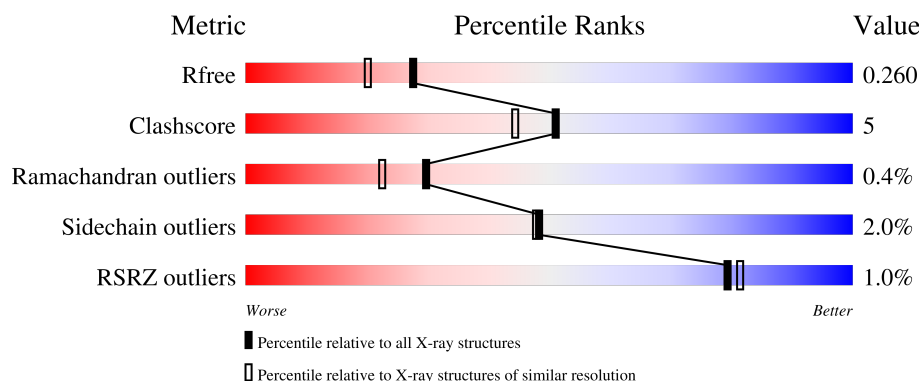
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

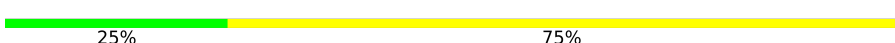
The reported resolution of this entry is 2.04 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	398	 87% 11% .
1	B	398	 86% 11% ..
2	C	3	 33% 67%
3	D	4	 50% 50%
3	E	4	 25% 75%

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Mol	Chain	Length	Quality of chain
4	F	2	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	PEG	A	506	-	-	X	-

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 6839 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Alpha-galactosidase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	390	Total	C	N	O	S	0	2	0
			3135	1996	537	576	26			
1	B	392	Total	C	N	O	S	0	2	0
			3147	2001	540	579	27			

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

- Molecule 3 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
3	D	4	Total	C	N	O	0	0	0
			50	28	2	20			
3	E	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 4 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
4	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



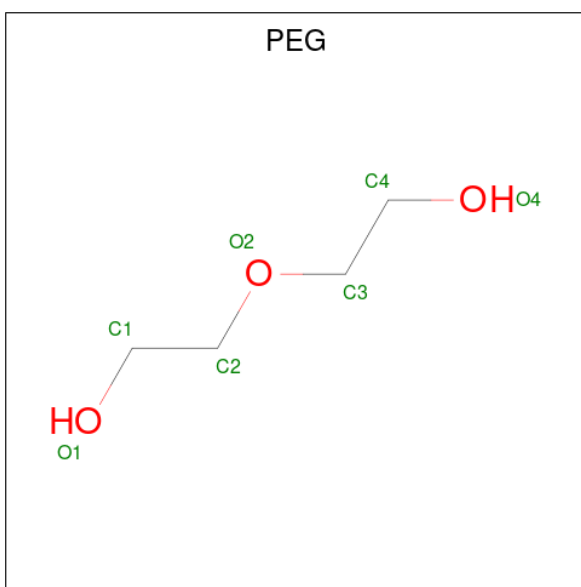
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



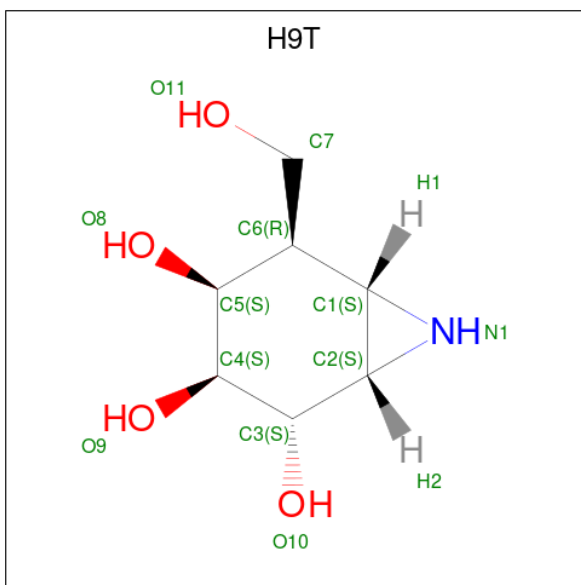
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



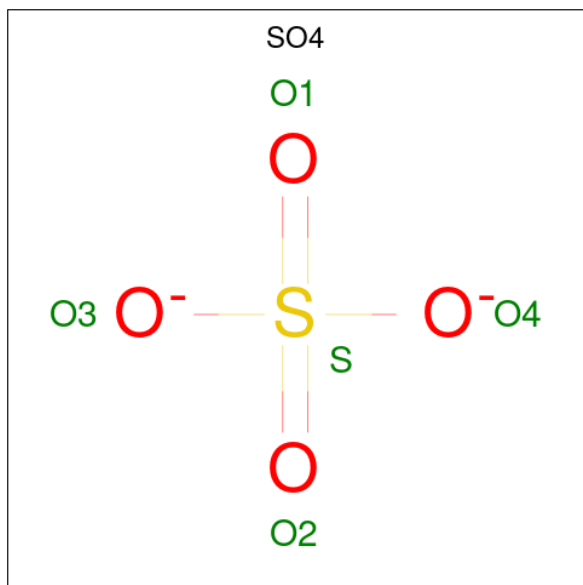
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			7	4	3		
7	A	1	Total	C	O	0	0
			7	4	3		
7	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is (1 {S},2 {S},3 {S},4 {S},5 {R},6 {S})-5-(hydroxymethyl)-7-azabicyclo[4.1.0]heptane-2,3,4-triol (CCD ID: H9T) (formula: $C_7H_{13}NO_4$) (labeled as "Ligand of Interest" by depositor).



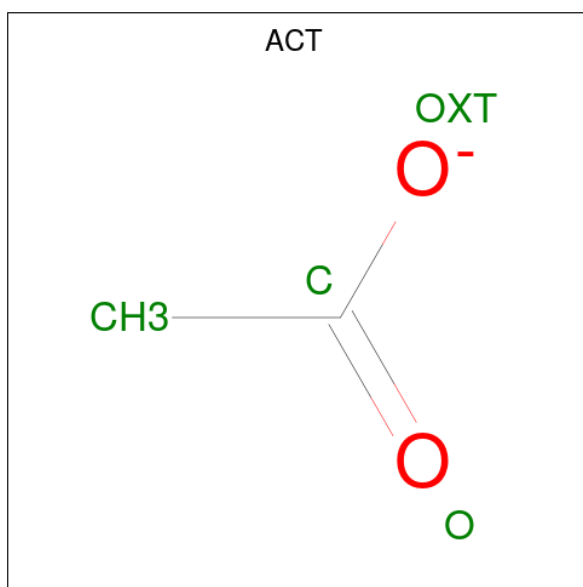
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			12	7	1	4		
8	B	1	Total	C	N	O	0	0
			12	7	1	4		

- Molecule 9 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 10 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		

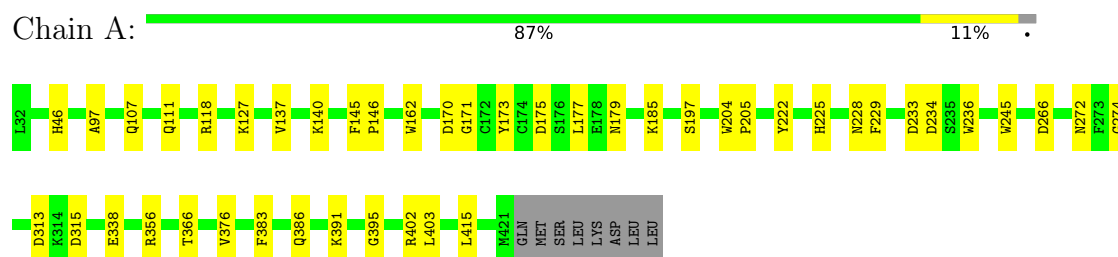
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	144	Total	O	0	0
			144	144		
11	B	96	Total	O	0	0
			96	96		

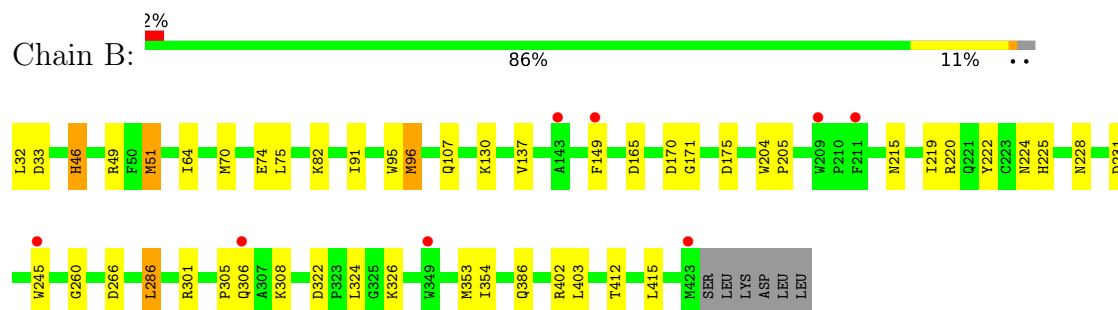
3 Residue-property plots

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

• Molecule 1: Alpha-galactosidase A



• Molecule 1: Alpha-galactosidase A



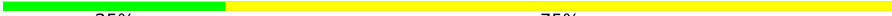
• Molecule 2: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 3: 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



• Molecule 3: 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:  25% 75%



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

Chain F:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	90.39Å 90.39Å 216.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	63.52 – 2.04 63.52 – 2.04	Depositor EDS
% Data completeness (in resolution range)	99.9 (63.52-2.04) 100.0 (63.52-2.04)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 2.03Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.195 , 0.260 0.200 , 0.260	Depositor DCC
R_{free} test set	3206 reflections (3.16%)	wwPDB-VP
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.566	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.020 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6839	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: H9T, ACT, EDO, PEG, BMA, MAN, NAG, SO4

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	1.01	2/3222 (0.1%)	1.31	3/4376 (0.1%)
1	B	1.04	1/3234 (0.0%)	1.32	0/4393
All	All	1.03	3/6456 (0.0%)	1.31	3/8769 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	170	ASP	CG-OD2	10.90	1.46	1.25
1	B	170	ASP	CG-OD2	10.34	1.45	1.25
1	A	272	ASN	N-CA	5.11	1.49	1.46

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	ASP	CB-CA-C	-5.82	100.00	110.36
1	A	313	ASP	CA-C-N	5.02	126.97	120.44
1	A	313	ASP	C-N-CA	5.02	126.97	120.44

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	3135	0	2982	31	0
1	B	3147	0	2989	33	0
2	C	39	0	34	0	0
3	D	50	0	43	0	0
3	E	50	0	43	0	0
4	F	28	0	25	1	0
5	A	14	0	13	0	0
5	B	14	0	13	2	0
6	A	8	0	12	0	0
6	B	20	0	30	3	0
7	A	21	0	30	10	0
8	A	12	0	0	0	0
8	B	12	0	0	1	0
9	A	15	0	0	1	0
9	B	10	0	0	2	0
10	A	12	0	9	1	0
10	B	12	0	9	1	0
11	A	144	0	0	1	0
11	B	96	0	0	0	0
All	All	6839	0	6232	67	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 5.

All (67) 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:236:TRP:HB3	7:A:506:PEG:H42	1.46	0.96
1:B:402:ARG:NH2	9:B:514:SO4:O4	2.17	0.77
1:B:231:ASP:O	10:B:517:ACT:H2	1.88	0.73
1:A:236:TRP:HB3	7:A:506:PEG:C4	2.20	0.71
1:B:49:ARG:HH11	6:B:501:EDO:C1	2.05	0.70
1:A:236:TRP:CB	7:A:506:PEG:H42	2.23	0.65
1:A:403:LEU:HD21	1:A:415:LEU:CD1	2.27	0.65
1:B:49:ARG:HH11	6:B:501:EDO:H11	1.63	0.62
1:B:286:LEU:HD21	1:B:354:ILE:HD11	1.83	0.61
1:A:185:LYS:HG2	1:A:222:TYR:CZ	2.36	0.61
1:A:118:ARG:H	7:A:507:PEG:H21	1.67	0.59
1:A:229:PHE:CE2	10:A:518:ACT:H1	2.40	0.57
1:A:402:ARG:NH2	9:A:515:SO4:O2	2.37	0.57
1:A:97:ALA:HB3	1:A:107:GLN:HG3	1.87	0.56
1:A:315:ASP:OD2	1:A:391:LYS:HE3	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:VAL:HG12	1:A:171:GLY:HA2	1.88	0.56
1:A:315:ASP:OD2	1:A:391:LYS:CE	2.54	0.56
1:B:403:LEU:HD21	1:B:415:LEU:CD1	2.37	0.55
1:B:228:ASN:HB3	1:B:245:TRP:CH2	2.42	0.54
1:A:145:PHE:HB3	1:A:146:PRO:HD2	1.89	0.53
1:A:204:TRP:HB3	1:A:205:PRO:HD3	1.90	0.53
1:B:137:VAL:HG12	1:B:171:GLY:HA2	1.90	0.52
1:B:205:PRO:HG3	1:B:219:ILE:HD13	1.91	0.52
7:A:506:PEG:H41	11:A:647:HOH:O	2.09	0.51
1:B:32:LEU:N	1:B:220:ARG:O	2.45	0.50
1:A:386:GLN:HA	1:A:415:LEU:HD23	1.92	0.50
1:B:91:ILE:HG21	1:B:95:TRP:HB3	1.95	0.49
1:A:236:TRP:N	7:A:506:PEG:H42	2.27	0.49
1:B:95:TRP:CD2	1:B:96:MET:HG3	2.48	0.49
1:A:118:ARG:HG3	1:A:162:TRP:HA	1.94	0.49
1:A:228:ASN:HB3	1:A:245:TRP:CH2	2.47	0.49
1:A:366:THR:HA	1:A:403:LEU:O	2.14	0.48
4:F:2:NAG:H83	4:F:2:NAG:H3	1.96	0.48
1:A:403:LEU:HD21	1:A:415:LEU:HD11	1.93	0.48
1:A:140:LYS:HB2	1:A:173:TYR:CD2	2.49	0.48
1:B:33:ASP:HB2	9:B:515:SO4:O1	2.13	0.47
1:B:215:ASN:C	1:B:215:ASN:OD1	2.57	0.47
1:B:386:GLN:HA	1:B:415:LEU:HD23	1.97	0.47
1:B:322:ASP:OD2	1:B:324:LEU:HB3	2.16	0.46
1:B:324:LEU:HG	1:B:326:LYS:HG3	1.97	0.45
1:A:177:LEU:HD12	1:A:177:LEU:H	1.81	0.45
1:B:149:PHE:CG	5:B:513:NAG:H82	2.51	0.45
1:B:70:MET:O	1:B:74:GLU:HG3	2.15	0.44
1:B:32:LEU:HD22	1:B:222:TYR:O	2.17	0.44
1:B:49:ARG:HH11	6:B:501:EDO:H12	1.79	0.44
1:B:204:TRP:HB3	1:B:205:PRO:HD3	2.00	0.44
1:A:175:ASP:OD1	1:A:179:ASN:ND2	2.49	0.43
1:B:306[B]:GLN:CD	1:B:306[B]:GLN:H	2.26	0.43
1:A:236:TRP:HB3	7:A:506:PEG:H11	1.99	0.43
1:B:130:LYS:HB3	1:B:165:ASP:HB2	2.01	0.43
1:A:383:PHE:CE2	1:A:395:GLY:HA2	2.54	0.42
1:B:175:ASP:OD2	5:B:513:NAG:C7	2.68	0.42
1:A:145:PHE:HB3	1:A:146:PRO:CD	2.49	0.42
1:A:234:ASP:O	1:A:274:GLY:HA3	2.20	0.42
1:B:51:MET:O	1:B:64:ILE:HD11	2.20	0.41
7:A:507:PEG:H31	7:A:507:PEG:H12	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:HIS:CD2	1:B:46:HIS:C	2.98	0.41
1:B:231:ASP:OD1	8:B:505:H9T:O10	2.39	0.41
1:B:286:LEU:CD2	1:B:354:ILE:HD11	2.49	0.41
1:A:338:GLU:OE2	1:A:356:ARG:NH1	2.54	0.41
1:B:75:LEU:CD1	1:B:301:ARG:HG2	2.51	0.41
1:B:224:ASN:HA	1:B:260:GLY:O	2.20	0.41
1:A:236:TRP:CA	7:A:506:PEG:H42	2.51	0.41
1:A:236:TRP:H	7:A:506:PEG:H42	1.86	0.41
1:A:315:ASP:OD2	1:A:391:LYS:HE2	2.21	0.40
1:B:353:MET:O	1:B:412:THR:HA	2.21	0.40
1:B:305:PRO:HD2	1:B:306[B]:GLN:OE1	2.21	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	390/398 (98%)	375 (96%)	13 (3%)	2 (0%)	24	16
1	B	392/398 (98%)	372 (95%)	19 (5%)	1 (0%)	36	30
All	All	782/796 (98%)	747 (96%)	32 (4%)	3 (0%)	30	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	266	ASP
1	B	266	ASP
1	A	197	SER

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	330/339 (97%)	325 (98%)	5 (2%)	57	58
1	B	332/339 (98%)	324 (98%)	8 (2%)	43	40
All	All	662/678 (98%)	649 (98%)	13 (2%)	48	48

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

Mol	Chain	Res	Type
1	A	46	HIS
1	A	111	GLN
1	A	127	LYS
1	A	225	HIS
1	A	376	VAL
1	B	46	HIS
1	B	51	MET
1	B	82	LYS
1	B	96	MET
1	B	107	GLN
1	B	225	HIS
1	B	286	LEU
1	B	308	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	336	ASN
1	A	357	GLN
1	B	111	GLN
1	B	179	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 ⓘ

13 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	2,1	14,14,15	0.54	0	17,19,21	1.21	1 (5%)
2	NAG	C	2	2	14,14,15	0.37	0	17,19,21	1.27	2 (11%)
2	BMA	C	3	2	11,11,12	0.44	0	15,15,17	0.74	0
3	NAG	D	1	3,1	14,14,15	0.62	0	17,19,21	0.77	0
3	NAG	D	2	3	14,14,15	0.38	0	17,19,21	1.49	4 (23%)
3	BMA	D	3	3	11,11,12	0.47	0	15,15,17	1.09	1 (6%)
3	MAN	D	4	3	11,11,12	0.47	0	15,15,17	0.87	0
3	NAG	E	1	3,1	14,14,15	0.30	0	17,19,21	0.93	1 (5%)
3	NAG	E	2	3	14,14,15	0.55	0	17,19,21	0.82	0
3	BMA	E	3	3	11,11,12	0.52	0	15,15,17	1.18	3 (20%)
3	MAN	E	4	3	11,11,12	0.43	0	15,15,17	1.30	1 (6%)
4	NAG	F	1	4,1	14,14,15	0.38	0	17,19,21	1.00	1 (5%)
4	NAG	F	2	4	14,14,15	0.42	0	17,19,21	1.99	4 (23%)

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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	1/2/19/22	0/1/1/1
3	NAG	D	1	3,1	-	3/6/23/26	0/1/1/1

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

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	2	NAG	C2-N2-C7	4.72	129.22	122.90
4	F	2	NAG	C8-C7-N2	4.10	122.91	116.12
2	C	1	NAG	C1-O5-C5	3.83	117.32	112.19
3	D	2	NAG	O4-C4-C5	3.42	117.76	109.32
4	F	2	NAG	C3-C4-C5	2.97	115.61	110.23
3	E	4	MAN	C3-C4-C5	2.96	115.60	110.23
2	C	2	NAG	C1-O5-C5	2.84	115.99	112.19
3	D	2	NAG	O4-C4-C3	-2.66	104.12	110.38
3	D	3	BMA	O5-C5-C6	2.53	112.59	107.66
4	F	1	NAG	C4-C3-C2	2.39	114.52	111.02
3	E	3	BMA	O4-C4-C3	-2.38	104.77	110.38
3	D	2	NAG	C6-C5-C4	2.30	118.67	113.02
3	E	3	BMA	C3-C4-C5	2.26	114.33	110.23
3	E	1	NAG	C1-C2-N2	2.24	113.96	110.43
3	D	2	NAG	C3-C4-C5	-2.19	106.25	110.23
2	C	2	NAG	O4-C4-C3	-2.17	105.27	110.38
4	F	2	NAG	O7-C7-N2	-2.17	118.16	121.98
3	E	3	BMA	C1-O5-C5	-2.06	109.42	112.19

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	4	MAN	O5-C5-C6-O6
3	D	4	MAN	C4-C5-C6-O6
4	F	2	NAG	C8-C7-N2-C2

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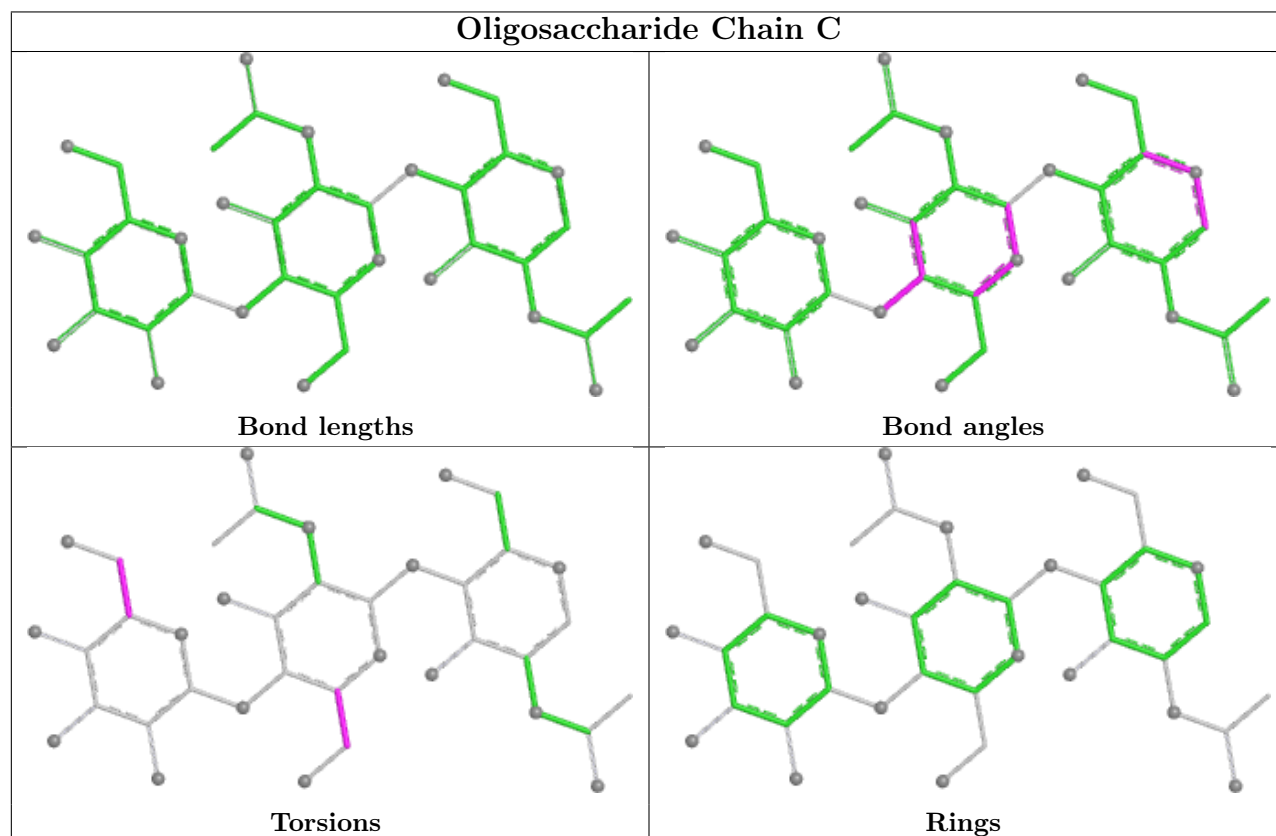
Mol	Chain	Res	Type	Atoms
4	F	2	NAG	O7-C7-N2-C2
3	E	3	BMA	C4-C5-C6-O6
3	E	3	BMA	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	C	3	BMA	C4-C5-C6-O6
4	F	2	NAG	C3-C2-N2-C7
3	D	1	NAG	C1-C2-N2-C7
3	E	2	NAG	C1-C2-N2-C7

There are no ring outliers.

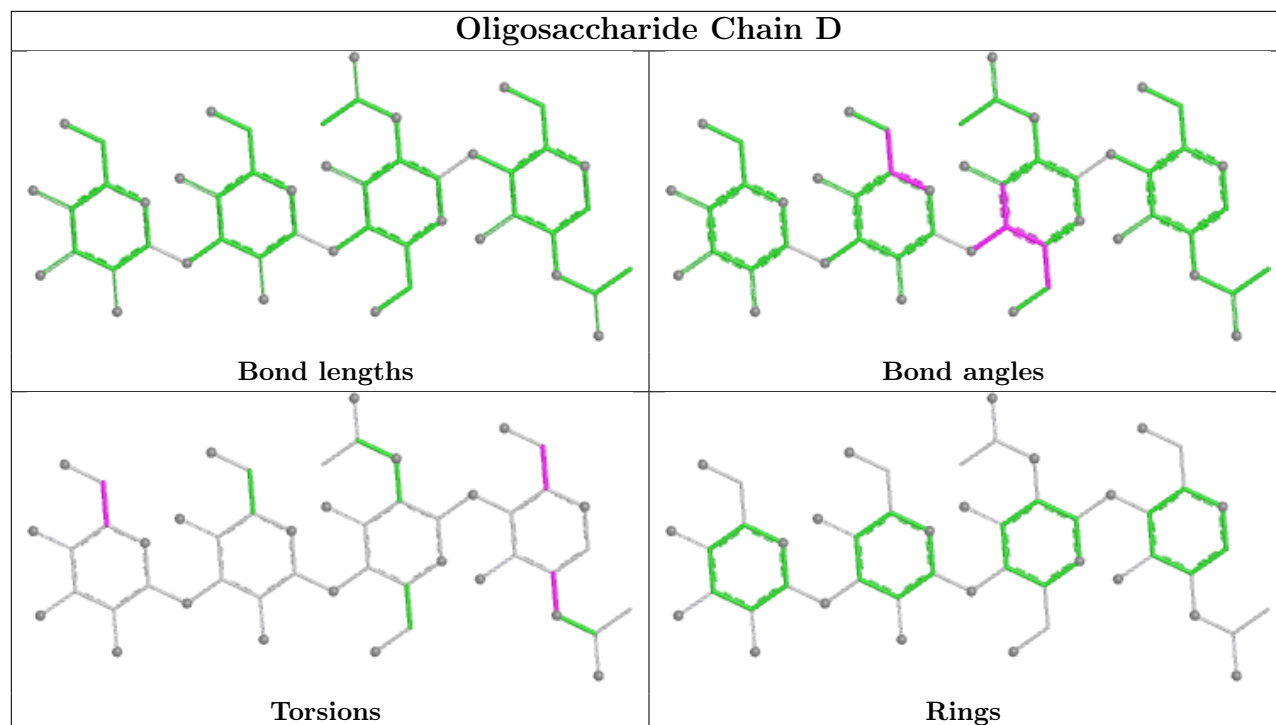
1 monomer is involved in 1 short contact:

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

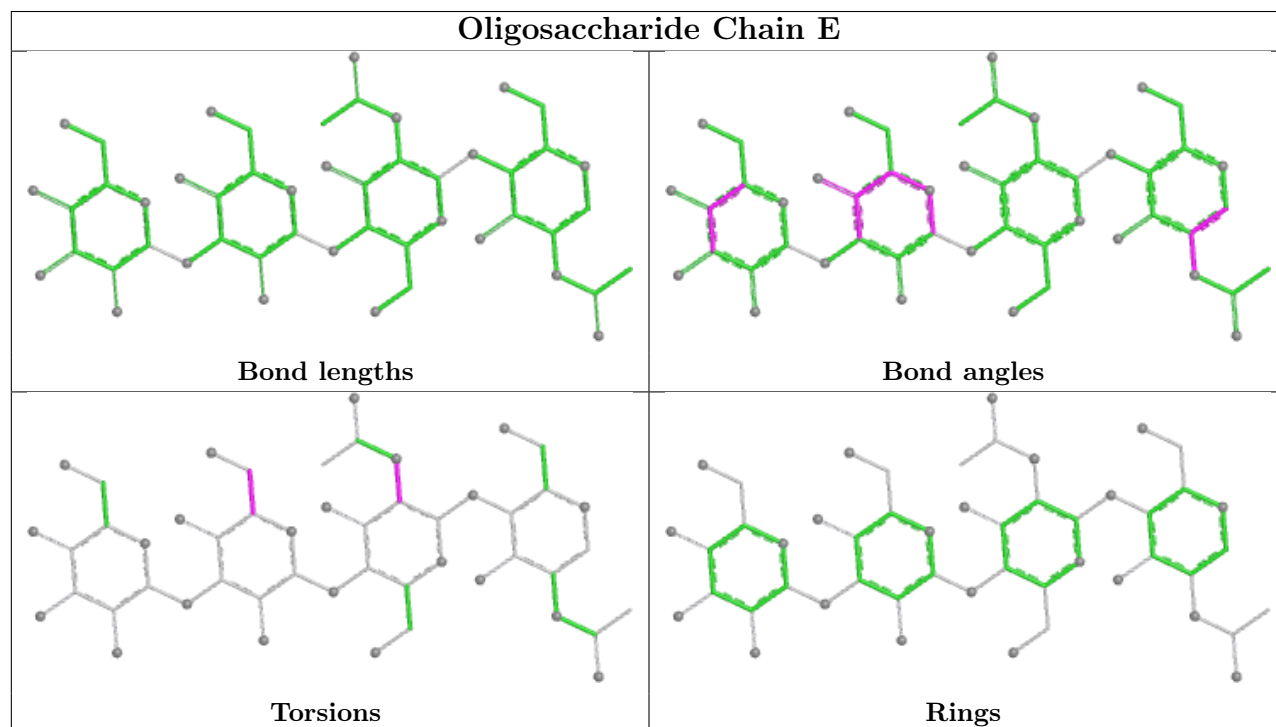
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

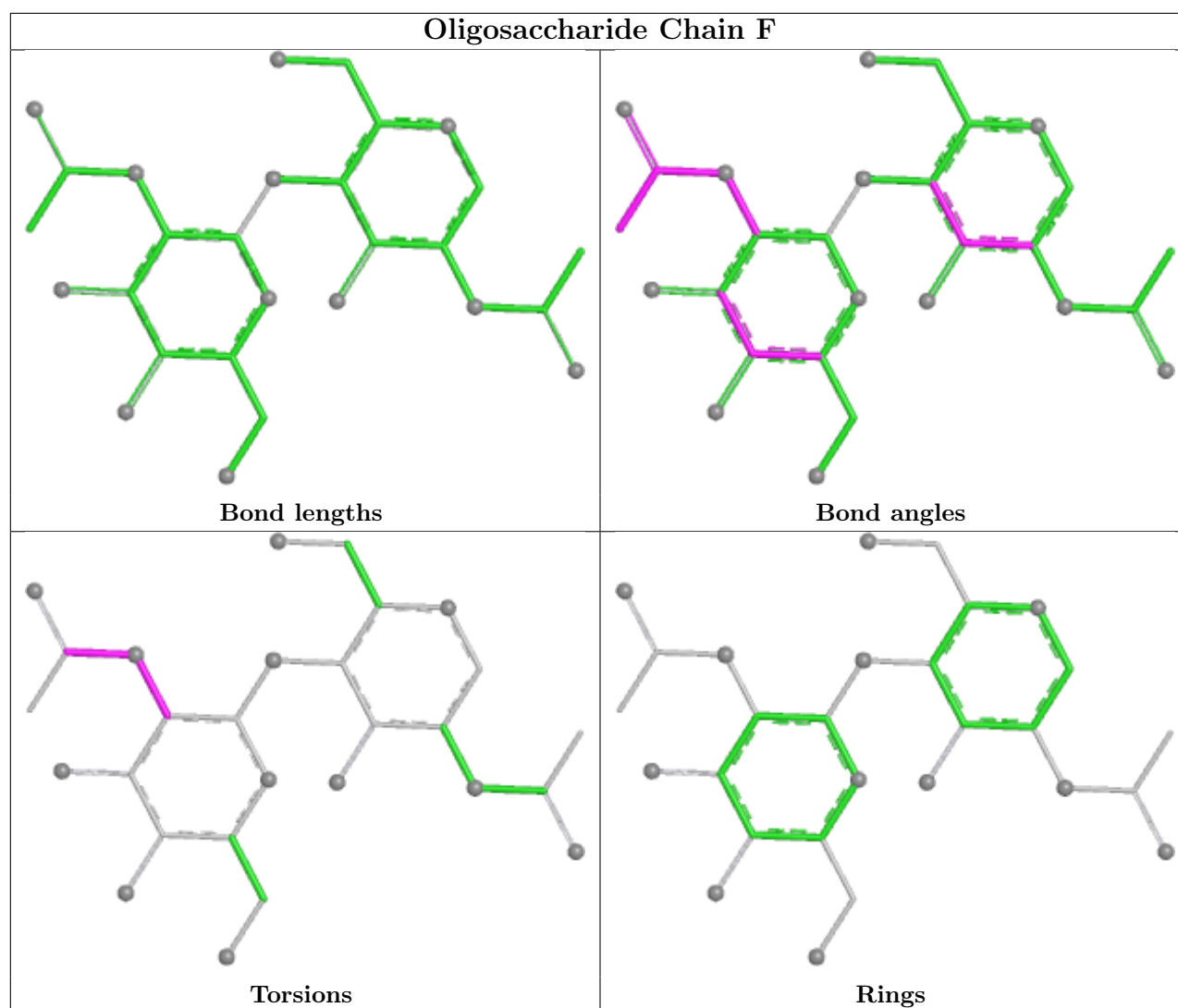


Oligosaccharide Chain D



Oligosaccharide Chain E





5.6 Ligand geometry [i](#)

25 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	SO4	B	514	-	4,4,4	0.39	0	6,6,6	0.15	0
9	SO4	A	515	-	4,4,4	0.37	0	6,6,6	0.11	0
6	EDO	B	502	-	3,3,3	0.12	0	2,2,2	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	SO4	B	515	-	4,4,4	0.31	0	6,6,6	0.12	0
6	EDO	B	503	-	3,3,3	0.27	0	2,2,2	0.27	0
10	ACT	A	520	-	3,3,3	1.14	0	3,3,3	0.71	0
10	ACT	B	517	-	3,3,3	0.93	0	3,3,3	0.88	0
6	EDO	B	501	-	3,3,3	0.10	0	2,2,2	0.61	0
5	NAG	A	501	1	14,14,15	0.65	0	17,19,21	1.53	2 (11%)
9	SO4	A	516	-	4,4,4	0.28	0	6,6,6	0.16	0
10	ACT	A	519	-	3,3,3	1.12	0	3,3,3	0.80	0
7	PEG	A	506	-	6,6,6	0.50	0	5,5,5	0.80	0
5	NAG	B	513	1	14,14,15	0.52	0	17,19,21	1.43	3 (17%)
10	ACT	B	516	-	3,3,3	1.06	0	3,3,3	0.84	0
10	ACT	A	518	-	3,3,3	0.74	0	3,3,3	0.99	0
7	PEG	A	507	-	6,6,6	0.42	0	5,5,5	0.41	0
6	EDO	A	508	-	3,3,3	0.10	0	2,2,2	0.15	0
8	H9T	B	505	1	13,13,13	38.07	3 (23%)	15,20,20	41.84	5 (33%)
10	ACT	B	518	-	3,3,3	0.69	0	3,3,3	1.10	0
7	PEG	A	514	-	6,6,6	0.16	0	5,5,5	0.17	0
6	EDO	A	505	-	3,3,3	0.11	0	2,2,2	0.18	0
6	EDO	B	504	-	3,3,3	0.16	0	2,2,2	0.27	0
6	EDO	B	506	-	3,3,3	0.10	0	2,2,2	0.24	0
8	H9T	A	509	1	13,13,13	36.73	2 (15%)	15,20,20	40.71	5 (33%)
9	SO4	A	517	-	4,4,4	0.31	0	6,6,6	0.08	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	PEG	A	507	-	-	2/4/4/4	-
6	EDO	A	508	-	-	1/1/1/1	-
7	PEG	A	506	-	-	3/4/4/4	-
8	H9T	B	505	1	-	0/2/27/27	0/2/2/2
5	NAG	B	513	1	-	0/6/23/26	0/1/1/1
6	EDO	B	502	-	-	1/1/1/1	-
6	EDO	B	503	-	-	1/1/1/1	-
6	EDO	B	501	-	-	0/1/1/1	-
6	EDO	A	505	-	-	1/1/1/1	-
6	EDO	B	504	-	-	1/1/1/1	-
6	EDO	B	506	-	-	1/1/1/1	-
7	PEG	A	514	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	501	1	-	0/6/23/26	0/1/1/1
8	H9T	A	509	1	-	0/2/27/27	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	505	H9T	C2-N1	137.16	2.41	1.48
8	A	509	H9T	C2-N1	132.36	2.38	1.48
8	B	505	H9T	C2-C1	4.21	1.52	1.48
8	A	509	H9T	C2-C1	4.07	1.52	1.48
8	B	505	H9T	C1-N1	2.63	1.49	1.48

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	505	H9T	C2-C1-N1	133.94	106.00	59.81
8	A	509	H9T	C2-C1-N1	130.27	104.74	59.81
8	B	505	H9T	C1-C2-N1	-67.35	36.58	59.81
8	A	509	H9T	C1-C2-N1	-66.11	37.01	59.81
8	B	505	H9T	C2-N1-C1	-61.32	37.42	60.38
8	A	509	H9T	C2-N1-C1	-59.08	38.25	60.38
5	A	501	NAG	C1-O5-C5	4.49	118.21	112.19
5	B	513	NAG	C1-O5-C5	3.73	117.18	112.19
5	B	513	NAG	C2-N2-C7	2.87	126.74	122.90
8	A	509	H9T	C6-C1-N1	-2.80	111.77	118.31
8	A	509	H9T	C4-C3-C2	2.65	114.25	110.40
8	B	505	H9T	O8-C5-C6	2.41	114.85	109.94
8	B	505	H9T	C4-C3-C2	2.38	113.87	110.40
5	B	513	NAG	C4-C3-C2	2.35	114.45	111.02
5	A	501	NAG	C4-C3-C2	2.09	114.08	111.02

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	507	PEG	O1-C1-C2-O2
6	B	504	EDO	O1-C1-C2-O2
7	A	506	PEG	O1-C1-C2-O2
7	A	514	PEG	O2-C3-C4-O4
7	A	507	PEG	C1-C2-O2-C3
6	B	503	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
6	A	508	EDO	O1-C1-C2-O2
7	A	506	PEG	C1-C2-O2-C3
6	A	505	EDO	O1-C1-C2-O2
6	B	506	EDO	O1-C1-C2-O2
7	A	506	PEG	C4-C3-O2-C2
7	A	514	PEG	C4-C3-O2-C2
6	B	502	EDO	O1-C1-C2-O2

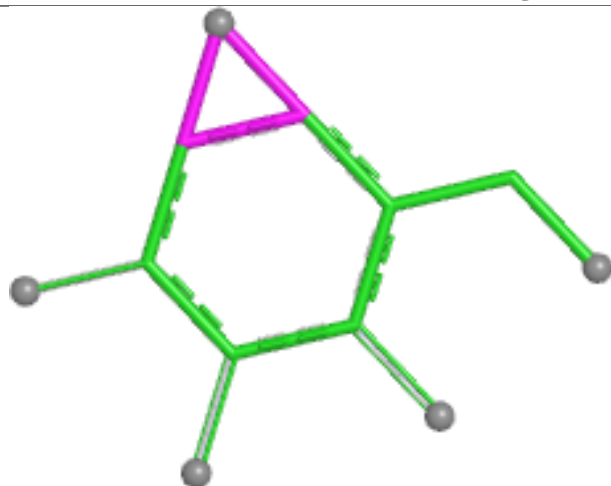
There are no ring outliers.

10 monomers are involved in 21 short contacts:

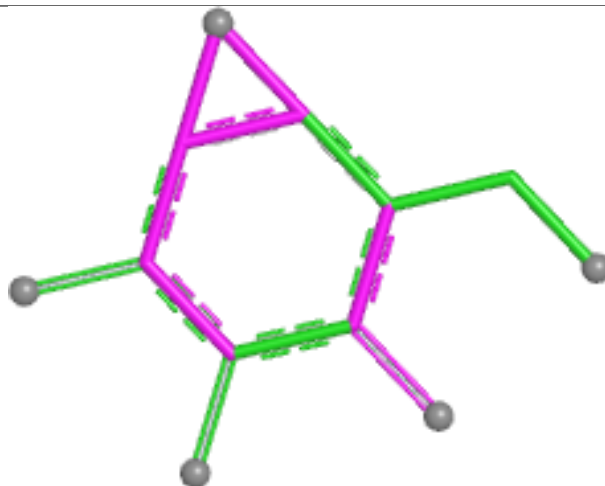
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B	514	SO4	1	0
9	A	515	SO4	1	0
9	B	515	SO4	1	0
10	B	517	ACT	1	0
6	B	501	EDO	3	0
7	A	506	PEG	8	0
5	B	513	NAG	2	0
10	A	518	ACT	1	0
7	A	507	PEG	2	0
8	B	505	H9T	1	0

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.

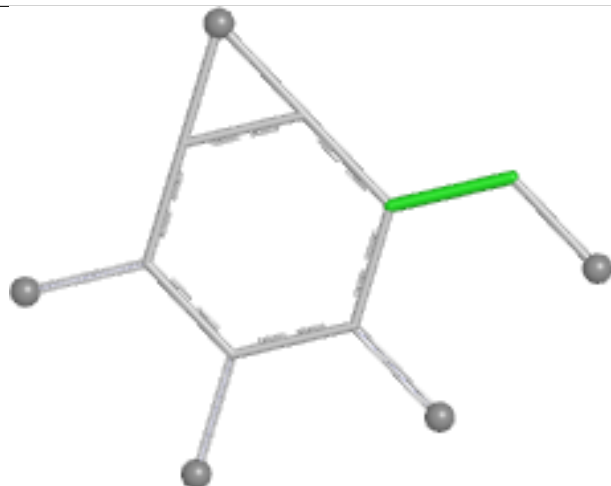
Ligand H9T B 505



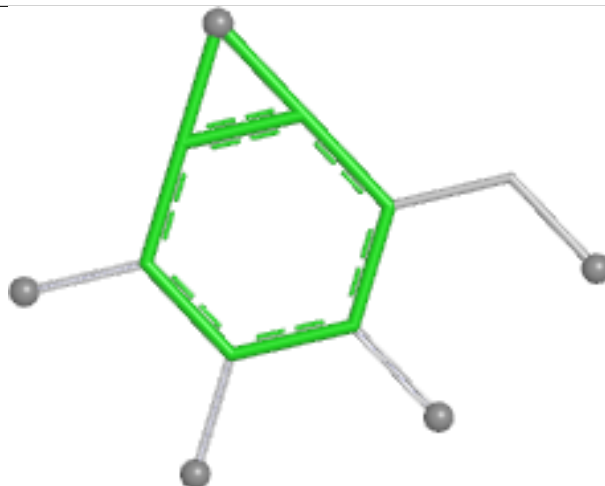
Bond lengths



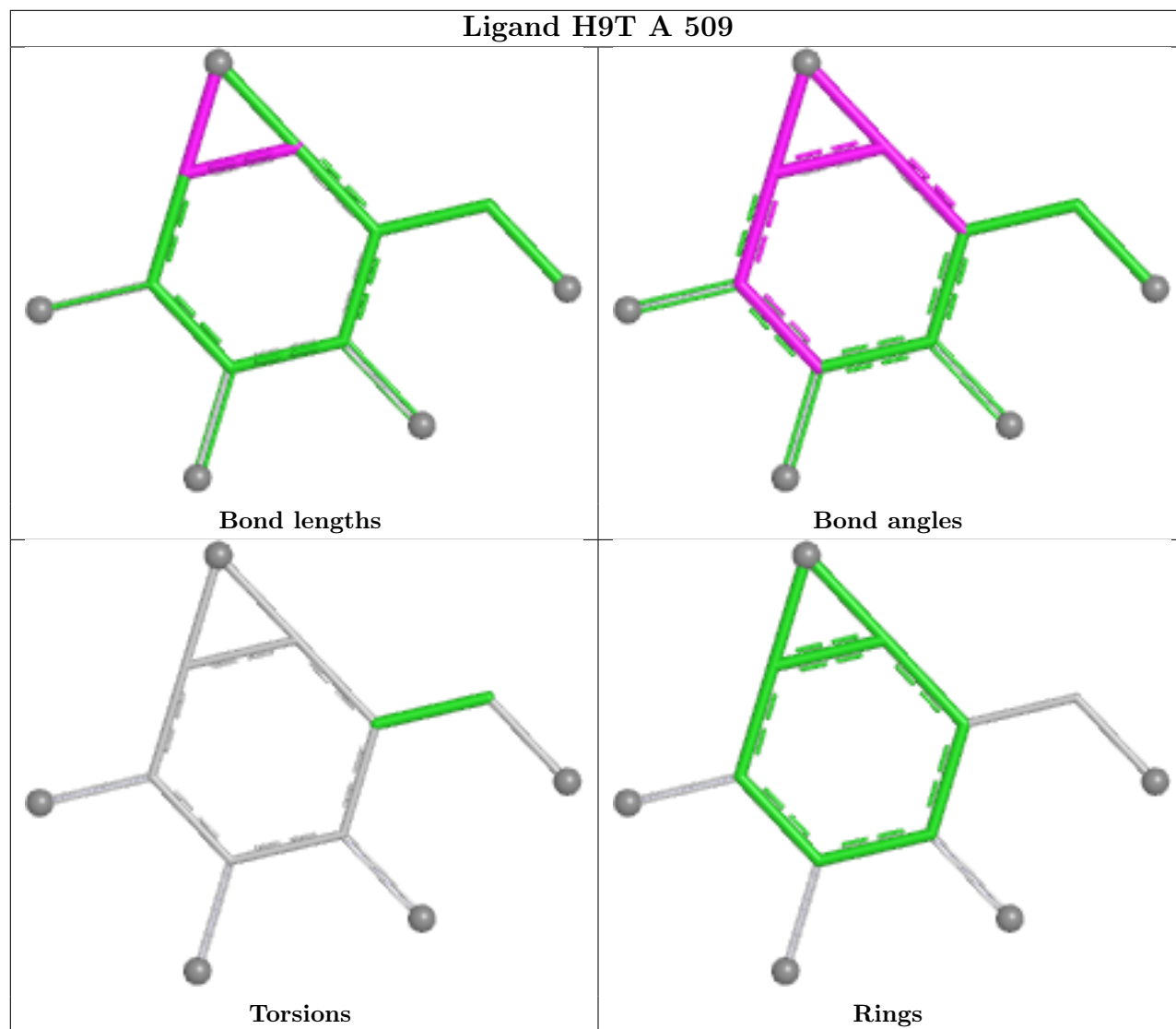
Bond angles



Torsions



Rings



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	390/398 (97%)	-0.11	0	100 100	23, 52, 69, 99	8 (2%)
1	B	392/398 (98%)	0.14	8 (2%)	65 67	31, 58, 83, 110	6 (1%)
All	All	782/796 (98%)	0.02	8 (1%)	79 81	23, 55, 80, 110	14 (1%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	149	PHE	3.8
1	B	211	PHE	2.9
1	B	306[A]	GLN	2.7
1	B	423	MET	2.6
1	B	245	TRP	2.4
1	B	209	TRP	2.2
1	B	349	TRP	2.1
1	B	143	ALA	2.1

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

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

6.3 Carbohydrates [i](#)

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

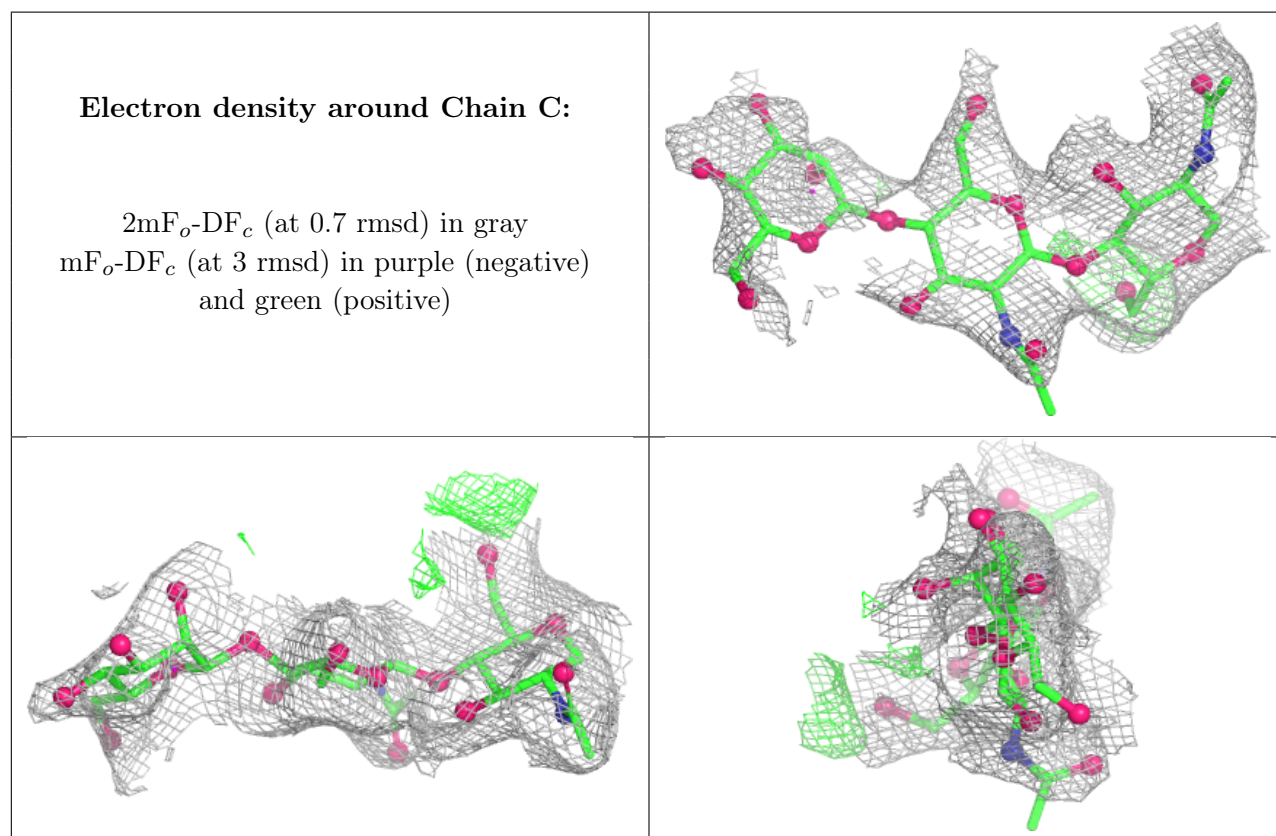
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BMA	C	3	11/12	0.62	0.09	126,132,138,139	0

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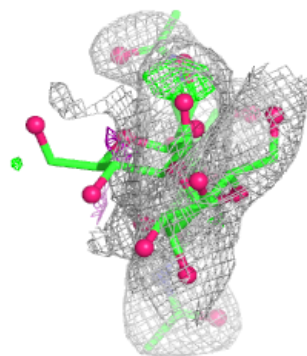
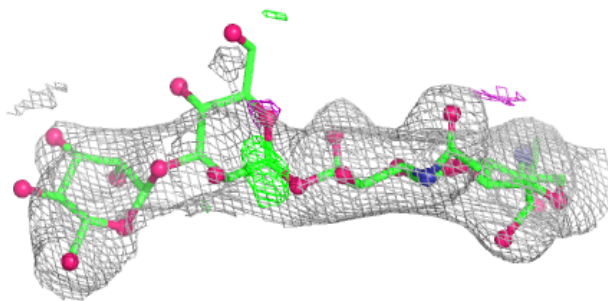
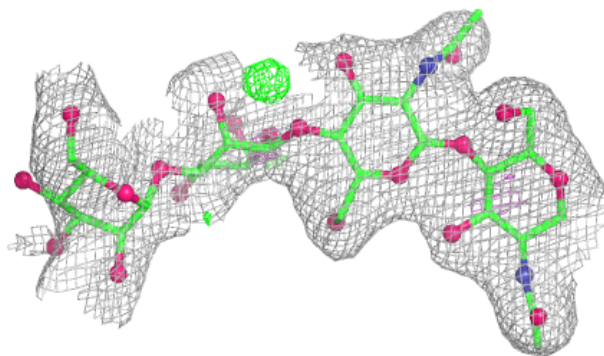
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	BMA	D	3	11/12	0.62	0.12	104,125,140,142	0
4	NAG	F	2	14/15	0.64	0.10	105,122,132,133	0
4	NAG	F	1	14/15	0.70	0.11	96,108,118,129	0
3	MAN	D	4	11/12	0.70	0.09	93,120,128,130	0
2	NAG	C	2	14/15	0.74	0.10	93,103,115,124	0
3	BMA	E	3	11/12	0.79	0.08	80,99,109,111	0
3	MAN	E	4	11/12	0.80	0.08	88,98,103,106	0
2	NAG	C	1	14/15	0.85	0.09	72,84,92,102	0
3	NAG	E	1	14/15	0.88	0.09	65,73,77,83	0
3	NAG	E	2	14/15	0.88	0.09	79,85,89,90	0
3	NAG	D	2	14/15	0.92	0.08	66,74,84,94	0
3	NAG	D	1	14/15	0.95	0.07	54,59,64,65	0

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

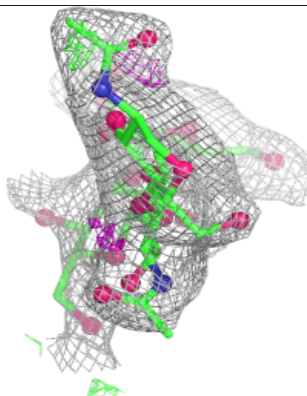
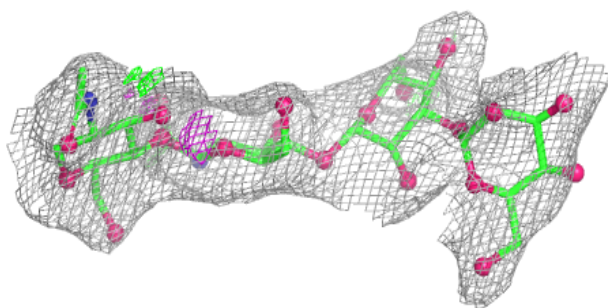
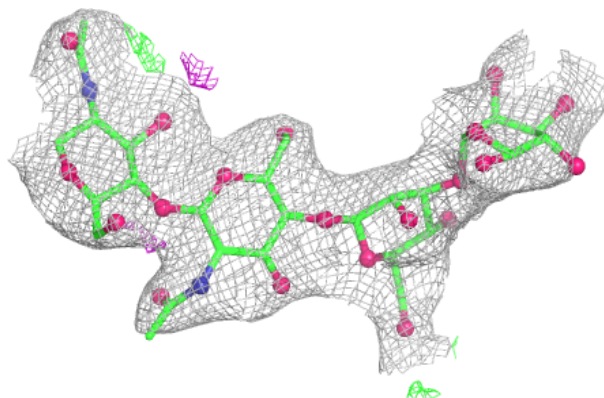


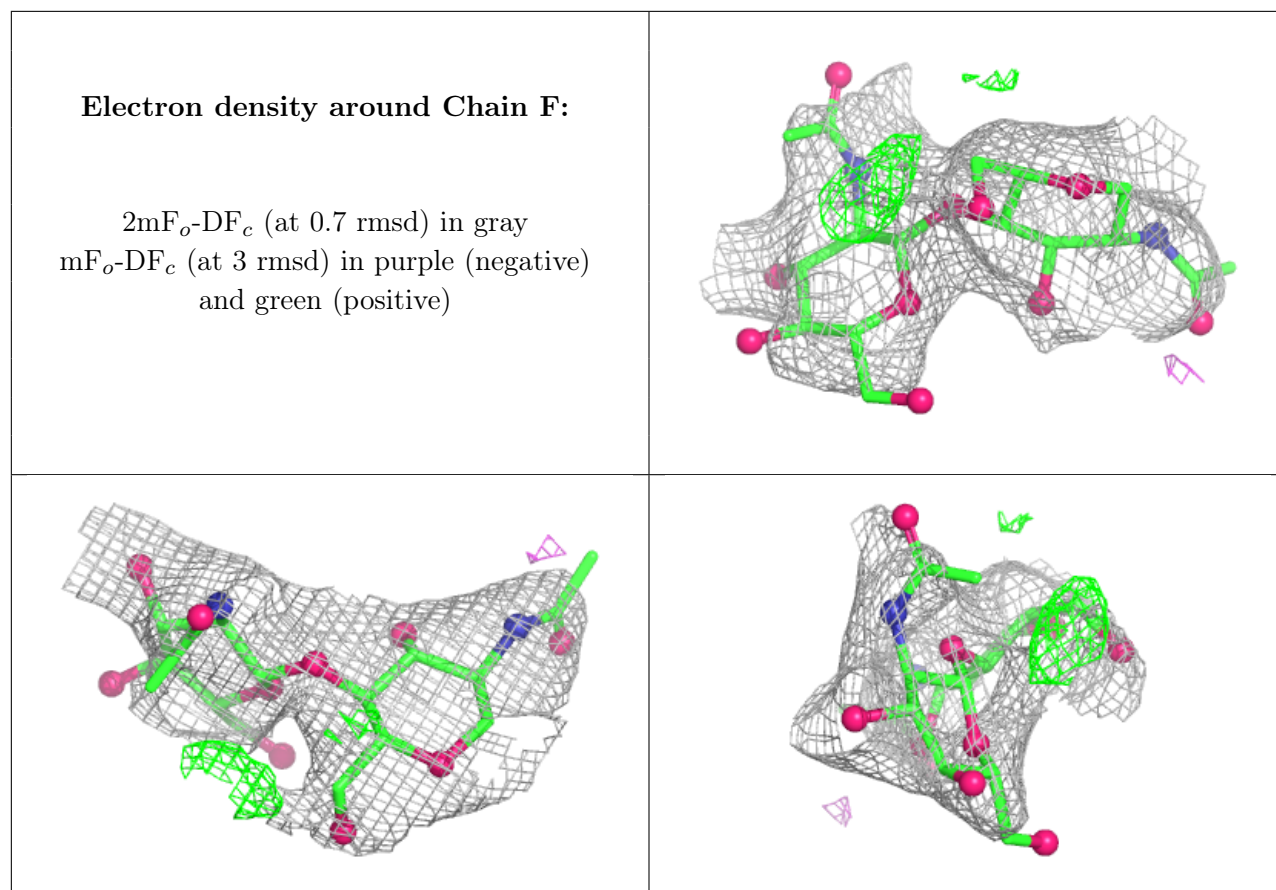
Electron density around Chain D:

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

**Electron density around Chain E:**

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





6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	B	513	14/15	0.68	0.12	93,109,119,120	0
9	SO4	A	517	5/5	0.68	0.10	112,123,126,128	0
7	PEG	A	507	7/7	0.69	0.20	67,87,93,95	0
10	ACT	A	519	4/4	0.74	0.17	67,73,80,82	0
7	PEG	A	506	7/7	0.76	0.21	22,24,30,32	7
5	NAG	A	501	14/15	0.79	0.09	68,75,88,98	0
6	EDO	B	506	4/4	0.82	0.14	81,84,84,86	0
6	EDO	B	503	4/4	0.84	0.18	69,74,77,77	0
10	ACT	B	518	4/4	0.85	0.15	74,75,76,78	0
9	SO4	B	515	5/5	0.86	0.07	102,106,112,117	0
7	PEG	A	514	7/7	0.86	0.14	68,73,75,80	0
10	ACT	A	520	4/4	0.86	0.14	64,67,73,78	3
6	EDO	A	508	4/4	0.86	0.19	105,107,110,112	0

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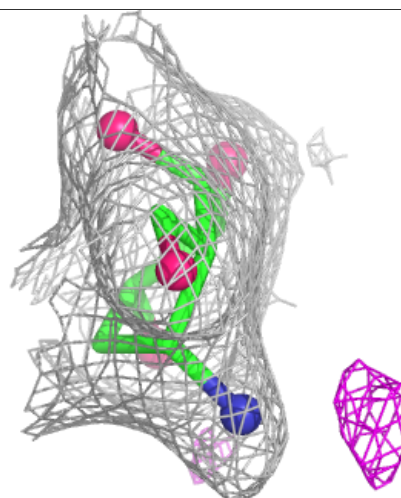
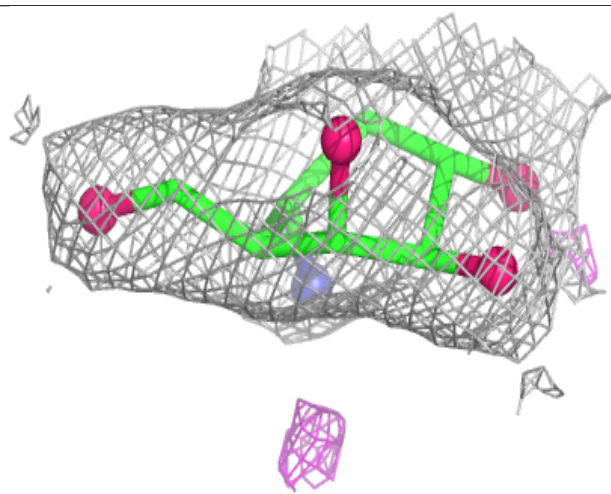
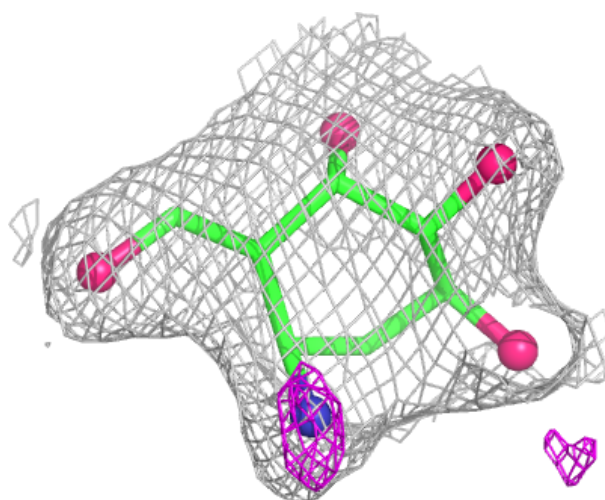
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
10	ACT	B	517	4/4	0.87	0.12	62,73,74,80	0
9	SO4	A	516	5/5	0.88	0.08	77,82,87,102	0
10	ACT	B	516	4/4	0.89	0.11	71,75,79,81	0
10	ACT	A	518	4/4	0.92	0.14	77,77,79,83	0
6	EDO	B	504	4/4	0.92	0.11	74,76,77,78	0
6	EDO	A	505	4/4	0.93	0.11	65,66,68,72	0
6	EDO	B	501	4/4	0.93	0.13	66,71,71,72	0
8	H9T	B	505	12/12	0.93	0.07	50,53,55,58	0
6	EDO	B	502	4/4	0.94	0.10	67,72,73,75	0
8	H9T	A	509	12/12	0.95	0.05	45,48,52,53	0
9	SO4	B	514	5/5	0.96	0.05	67,70,80,89	0
9	SO4	A	515	5/5	0.96	0.06	64,65,72,93	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.

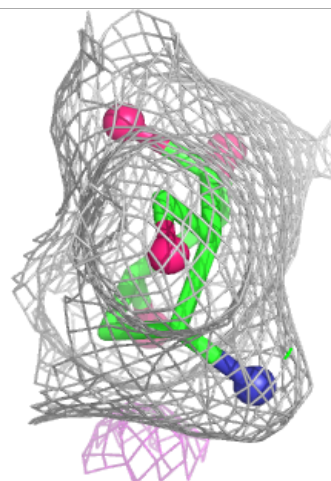
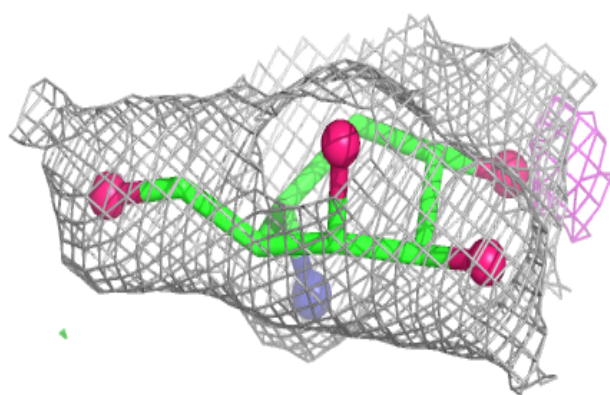
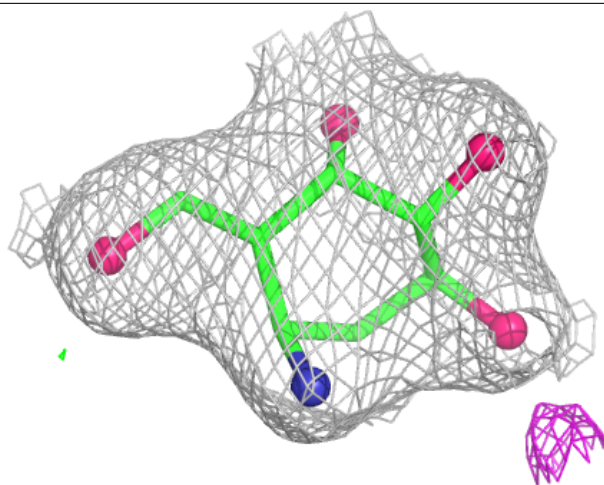
Electron density around H9T B 505:

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



Electron density around H9T A 509:

$2mF_o-DF_c$ (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.