



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:10 AM UTC

PDB ID : 2G7C / pdb_00002g7c
Title : Clostridium difficile Toxin A Fragment Bound to aGal(1,3)bGal(1,4)bGlcNAc
Authors : Greco, A.; Ho, J.G.S.; Lin, S.J.; Palcic, M.M.; Rupnik, M.; Ng, K.K.S.
Deposited on : 2006-02-28
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

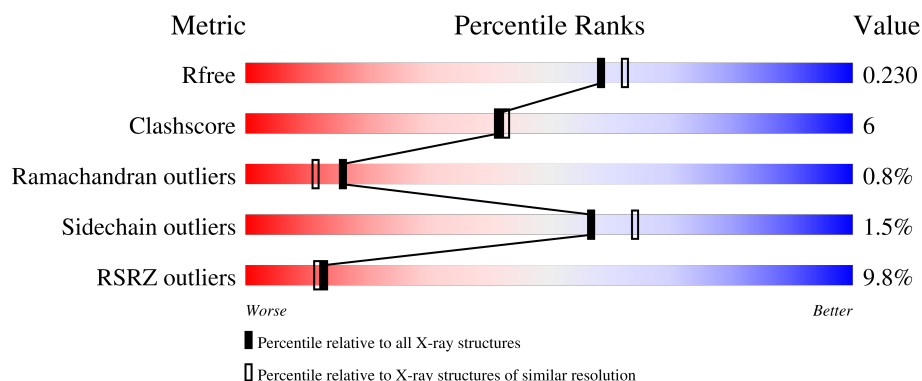
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	255	11% 79% 15% • •
1	B	255	8% 82% 15% •
2	C	3	33% 67%
2	D	3	33% 67%
2	E	3	33% 67%

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Mol	Chain	Length	Quality of chain
2	F	3	33% 67%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4588 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 Toxin A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1963	1267	326	365	5			
1	B	254	Total	C	N	O	S	0	0	0
			2028	1310	336	377	5			

There are 28 discrepancies between the modelled and reference sequences:

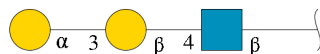
Chain	Residue	Modelled	Actual	Comment	Reference
A	14	MET	LEU	variant	UNP P16154
A	30	GLU	VAL	variant	UNP P16154
A	32	ALA	VAL	variant	UNP P16154
A	39	ASP	ASN	variant	UNP P16154
A	60	ASN	SER	variant	UNP P16154
A	61	ASP	GLY	variant	UNP P16154
A	133	ARG	ASN	variant	UNP P16154
A	143	ILE	MET	variant	UNP P16154
A	150	ILE	THR	variant	UNP P16154
A	171	PRO	SER	variant	UNP P16154
A	188	ASP	GLU	variant	UNP P16154
A	204	ASN	LYS	variant	UNP P16154
A	224	ASN	LYS	variant	UNP P16154
A	225	MET	VAL	variant	UNP P16154
B	14	MET	LEU	variant	UNP P16154
B	30	GLU	VAL	variant	UNP P16154
B	32	ALA	VAL	variant	UNP P16154
B	39	ASP	ASN	variant	UNP P16154
B	60	ASN	SER	variant	UNP P16154
B	61	ASP	GLY	variant	UNP P16154
B	133	ARG	ASN	variant	UNP P16154
B	143	ILE	MET	variant	UNP P16154
B	150	ILE	THR	variant	UNP P16154
B	171	PRO	SER	variant	UNP P16154
B	188	ASP	GLU	variant	UNP P16154

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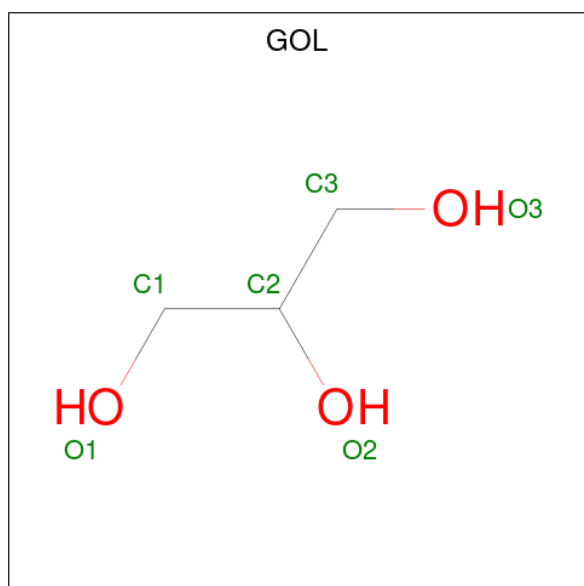
Chain	Residue	Modelled	Actual	Comment	Reference
B	204	ASN	LYS	variant	UNP P16154
B	224	ASN	LYS	variant	UNP P16154
B	225	MET	VAL	variant	UNP P16154

- Molecule 2 is an oligosaccharide called alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(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
			37	20	1	16			
2	D	3	Total	C	N	O	0	0	0
			37	20	1	16			
2	E	3	Total	C	N	O	0	0	0
			37	20	1	16			
2	F	3	Total	C	N	O	0	0	0
			37	20	1	16			

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 6	C 3	O 3	0	0
3	A	1	Total 6	C 3	O 3	0	0
3	A	1	Total 6	C 3	O 3	0	0
3	B	1	Total 6	C 3	O 3	0	0
3	B	1	Total 6	C 3	O 3	0	0

- Molecule 4 is water.

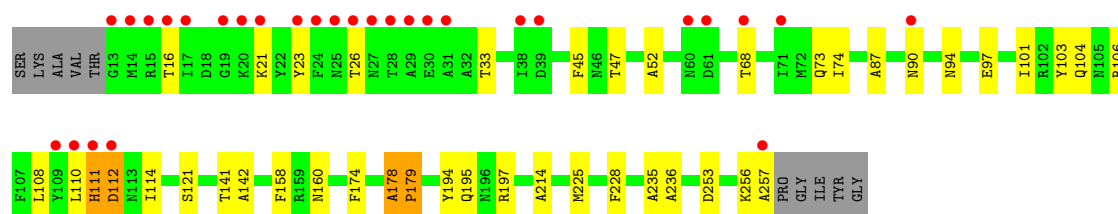
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	206	Total 206	O 206	0	0
4	B	207	Total 207	O 207	0	0

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.

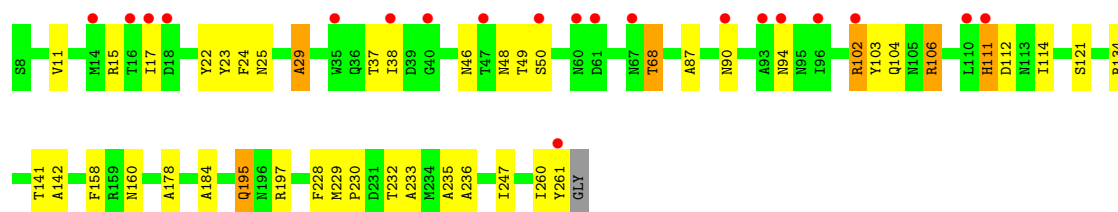
- Molecule 1: Toxin A

Chain A: 



- Molecule 1: Toxin A

Chain B: 



- Molecule 2: alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C: 



- Molecule 2: alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D: 



- Molecule 2: alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  33% 67%


MAG1
GAL2
GLA3

- Molecule 2: alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  33% 67%


MAG1
GAL2
GLA3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	38.68Å 132.82Å 65.92Å 90.00° 100.29° 90.00°	Depositor
Resolution (Å)	60.00 – 2.00 60.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (60.00-2.00) 99.9 (60.00-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.28 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.195 , 0.236 0.188 , 0.230	Depositor DCC
R_{free} test set	2210 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	19.9	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4588	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, GLA, GAL, NAG

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	0.40	0/2021	0.94	8/2740 (0.3%)
1	B	0.39	0/2088	0.91	9/2832 (0.3%)
All	All	0.40	0/4109	0.92	17/5572 (0.3%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	ALA	N-CA-C	9.30	122.83	110.36
1	B	178	ALA	N-CA-C	7.86	122.17	110.39
1	A	87	ALA	N-CA-C	7.63	120.40	110.40
1	B	104	GLN	N-CA-C	6.81	120.50	109.40
1	A	195	GLN	N-CA-C	6.73	120.37	109.40
1	B	195	GLN	N-CA-C	6.30	119.74	109.59
1	B	68	THR	N-CA-C	-6.01	105.21	112.54
1	A	104	GLN	N-CA-C	5.97	119.21	109.59
1	A	236	ALA	N-CA-C	-5.78	101.29	109.96
1	B	87	ALA	N-CA-C	5.76	119.56	110.58
1	A	158	PHE	N-CA-C	5.75	119.04	110.52
1	B	158	PHE	N-CA-C	5.23	118.59	110.17
1	B	184	ALA	CB-CA-C	-5.22	110.57	116.63
1	A	101	ILE	N-CA-C	5.16	116.29	108.96
1	B	29	ALA	N-CA-C	-5.10	106.65	112.92
1	B	236	ALA	N-CA-C	-5.09	102.74	110.28
1	A	179	PRO	N-CA-C	-5.03	103.39	111.38

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1963	0	1829	20	0
1	B	2028	0	1898	32	0
2	C	37	0	33	0	0
2	D	37	0	33	0	0
2	E	37	0	33	0	0
2	F	37	0	33	0	0
3	A	24	0	32	2	0
3	B	12	0	16	0	0
4	A	206	0	0	2	0
4	B	207	0	0	1	0
All	All	4588	0	3907	52	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 6.

All (52) 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:B:15:ARG:HH11	1:B:15:ARG:HG3	1.27	0.97
1:B:229:MET:HA	1:B:229:MET:HE3	1.66	0.77
1:B:195:GLN:O	1:B:197:ARG:HG3	1.93	0.67
1:B:15:ARG:HG3	1:B:15:ARG:NH1	2.03	0.63
1:B:11:VAL:HG12	1:B:29:ALA:HA	1.79	0.62
1:B:102:ARG:HH11	1:B:102:ARG:HB3	1.66	0.61
1:A:111:HIS:O	1:A:112:ASP:OD2	2.18	0.61
1:B:103:TYR:CZ	1:B:106:ARG:NH1	2.70	0.59
1:B:106:ARG:HG3	1:B:106:ARG:HH11	1.67	0.59
1:A:108:LEU:HG	1:A:110:LEU:CD2	2.34	0.58
1:A:16:THR:HB	1:A:21:LYS:HG2	1.85	0.58
1:B:260:ILE:O	1:B:261:TYR:HB2	2.04	0.57
1:B:46:ASN:OD1	1:B:48:ASN:HB2	2.05	0.56
1:B:17:ILE:O	1:B:17:ILE:HG13	2.05	0.55
1:B:228:PHE:CE1	1:B:235:ALA:HB2	2.42	0.55
1:B:37:THR:O	1:B:38:ILE:HD13	2.08	0.53
1:B:15:ARG:NH1	1:B:15:ARG:CG	2.71	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:THR:O	1:A:142:ALA:HB3	2.11	0.50
1:A:33:THR:HB	1:A:47:THR:HA	1.92	0.50
1:A:253:ASP:HB2	3:A:501:GOL:H32	1.92	0.50
1:B:114:ILE:HD12	1:B:160:ASN:HB3	1.93	0.50
1:A:194:TYR:CE2	1:A:197:ARG:HG3	2.48	0.49
1:A:108:LEU:HG	1:A:110:LEU:HD23	1.93	0.49
1:B:103:TYR:OH	1:B:106:ARG:NH1	2.46	0.49
1:A:73:GLN:C	1:A:74:ILE:HD12	2.37	0.48
1:A:228:PHE:CE1	1:A:235:ALA:HB2	2.49	0.48
1:B:90:ASN:OD1	1:B:94:ASN:HA	2.13	0.47
1:B:247:ILE:O	1:B:261:TYR:N	2.42	0.47
1:B:37:THR:C	1:B:38:ILE:HD13	2.39	0.47
1:B:111:HIS:O	1:B:112:ASP:HB2	2.15	0.46
1:B:106:ARG:NH1	1:B:106:ARG:HG3	2.31	0.46
1:B:229:MET:HE3	1:B:230:PRO:CD	2.45	0.46
1:B:17:ILE:HG12	1:B:22:TYR:CE1	2.51	0.45
1:A:90:ASN:HA	1:A:94:ASN:HA	1.98	0.45
1:B:49:THR:O	1:B:50:SER:HB2	2.16	0.45
1:A:23:TYR:HB2	1:A:45:PHE:CZ	2.52	0.44
1:B:232:THR:O	1:B:233:ALA:HB3	2.18	0.44
1:A:45:PHE:CE1	1:A:52:ALA:HB2	2.52	0.43
1:B:141:THR:O	1:B:142:ALA:HB3	2.18	0.43
1:A:257:ALA:HA	4:A:510:HOH:O	2.19	0.43
1:A:114:ILE:HD12	1:A:160:ASN:HB3	1.99	0.43
1:A:103:TYR:O	1:A:121:SER:HA	2.18	0.43
1:B:134:ARG:NH2	4:B:558:HOH:O	2.52	0.43
1:B:247:ILE:HB	1:B:261:TYR:HA	2.01	0.42
1:B:11:VAL:HG13	1:B:24:PHE:CD2	2.55	0.42
1:A:225:MET:HE1	3:A:503:GOL:O2	2.19	0.42
1:B:103:TYR:O	1:B:121:SER:HA	2.19	0.42
1:A:174:PHE:HB2	1:A:214:ALA:HB3	2.02	0.42
1:A:256:LYS:NZ	4:A:509:HOH:O	2.52	0.42
1:B:229:MET:HE3	1:B:229:MET:CA	2.44	0.41
1:A:178:ALA:HA	1:A:179:PRO:HD3	1.93	0.40
1:B:23:TYR:CE2	1:B:25:ASN:HB3	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	243/255 (95%)	235 (97%)	5 (2%)	3 (1%)	10	6
1	B	252/255 (99%)	243 (96%)	8 (3%)	1 (0%)	30	27
All	All	495/510 (97%)	478 (97%)	13 (3%)	4 (1%)	16	11

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	26	THR
1	A	111	HIS
1	B	111	HIS
1	A	112	ASP

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	194/201 (96%)	191 (98%)	3 (2%)	57	64
1	B	201/201 (100%)	198 (98%)	3 (2%)	57	64
All	All	395/402 (98%)	389 (98%)	6 (2%)	57	64

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

Mol	Chain	Res	Type
1	A	68	THR
1	A	97	GLU

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Mol	Chain	Res	Type
1	A	106	ARG
1	B	68	THR
1	B	102	ARG
1	B	106	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	36	GLN
1	A	140	ASN
1	A	190	GLN
1	A	196	ASN
1	A	210	ASN
1	B	48	ASN
1	B	104	GLN
1	B	140	ASN
1	B	164	GLN
1	B	172	ASN
1	B	181	ASN
1	B	196	ASN
1	B	210	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

12 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	15,15,15	0.42	0	21,21,21	0.59	0
2	GAL	C	2	2	11,11,12	0.60	0	15,15,17	0.99	1 (6%)
2	GLA	C	3	2	11,11,12	0.47	0	15,15,17	0.63	1 (6%)
2	NAG	D	1	2	15,15,15	0.46	0	21,21,21	0.69	0
2	GAL	D	2	2	11,11,12	0.57	0	15,15,17	1.01	1 (6%)
2	GLA	D	3	2	11,11,12	0.52	0	15,15,17	0.88	1 (6%)
2	NAG	E	1	2	15,15,15	0.43	0	21,21,21	0.60	0
2	GAL	E	2	2	11,11,12	0.61	0	15,15,17	1.13	1 (6%)
2	GLA	E	3	2	11,11,12	0.45	0	15,15,17	0.64	1 (6%)
2	NAG	F	1	2	15,15,15	0.38	0	21,21,21	0.70	0
2	GAL	F	2	2	11,11,12	0.44	0	15,15,17	0.84	1 (6%)
2	GLA	F	3	2	11,11,12	0.64	0	15,15,17	0.84	1 (6%)

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	2	-	0/6/26/26	0/1/1/1
2	GAL	C	2	2	-	0/2/19/22	0/1/1/1
2	GLA	C	3	2	-	0/2/19/22	0/1/1/1
2	NAG	D	1	2	-	0/6/26/26	0/1/1/1
2	GAL	D	2	2	-	0/2/19/22	0/1/1/1
2	GLA	D	3	2	-	0/2/19/22	0/1/1/1
2	NAG	E	1	2	-	0/6/26/26	0/1/1/1
2	GAL	E	2	2	-	1/2/19/22	0/1/1/1
2	GLA	E	3	2	-	0/2/19/22	0/1/1/1
2	NAG	F	1	2	-	0/6/26/26	0/1/1/1
2	GAL	F	2	2	-	0/2/19/22	0/1/1/1
2	GLA	F	3	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	2	GAL	C1-C2-C3	3.92	115.36	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	GAL	C1-C2-C3	3.38	114.57	109.64
2	C	2	GAL	C1-C2-C3	3.30	114.44	109.64
2	D	3	GLA	C1-O5-C5	2.74	115.86	112.19
2	F	2	GAL	C1-C2-C3	2.43	113.19	109.64
2	F	3	GLA	C1-O5-C5	2.18	115.11	112.19
2	C	3	GLA	C1-O5-C5	2.18	115.11	112.19
2	E	3	GLA	C1-O5-C5	2.11	115.02	112.19

There are no chirality outliers.

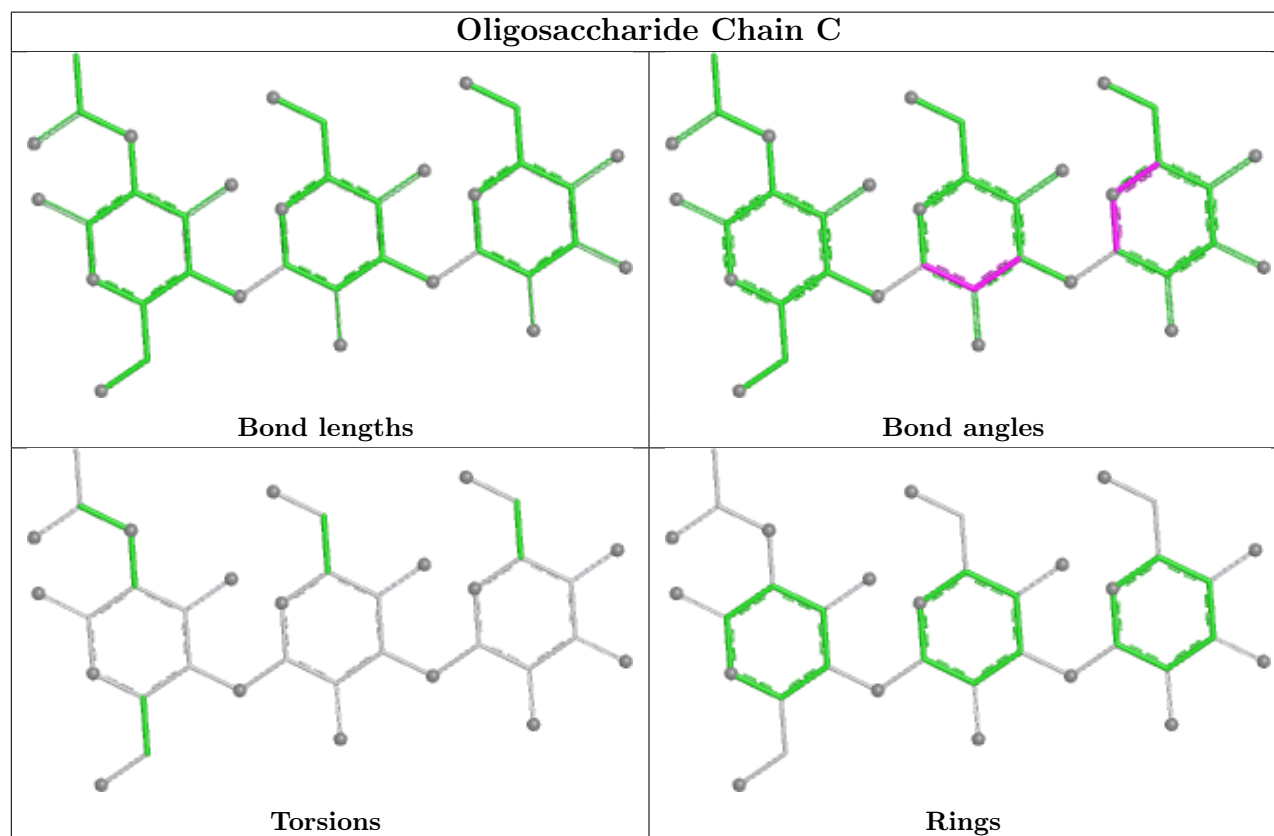
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	2	GAL	C4-C5-C6-O6

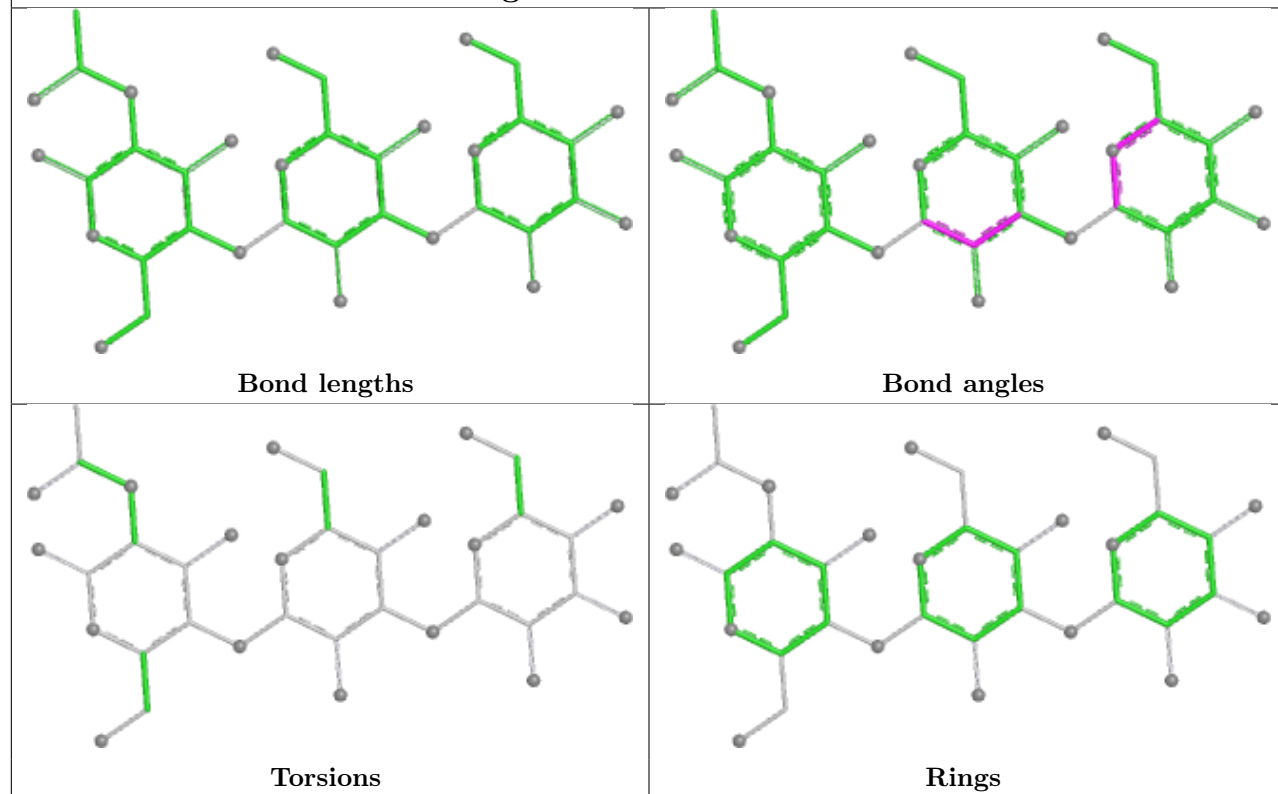
There are no ring outliers.

No monomer is involved in short contacts.

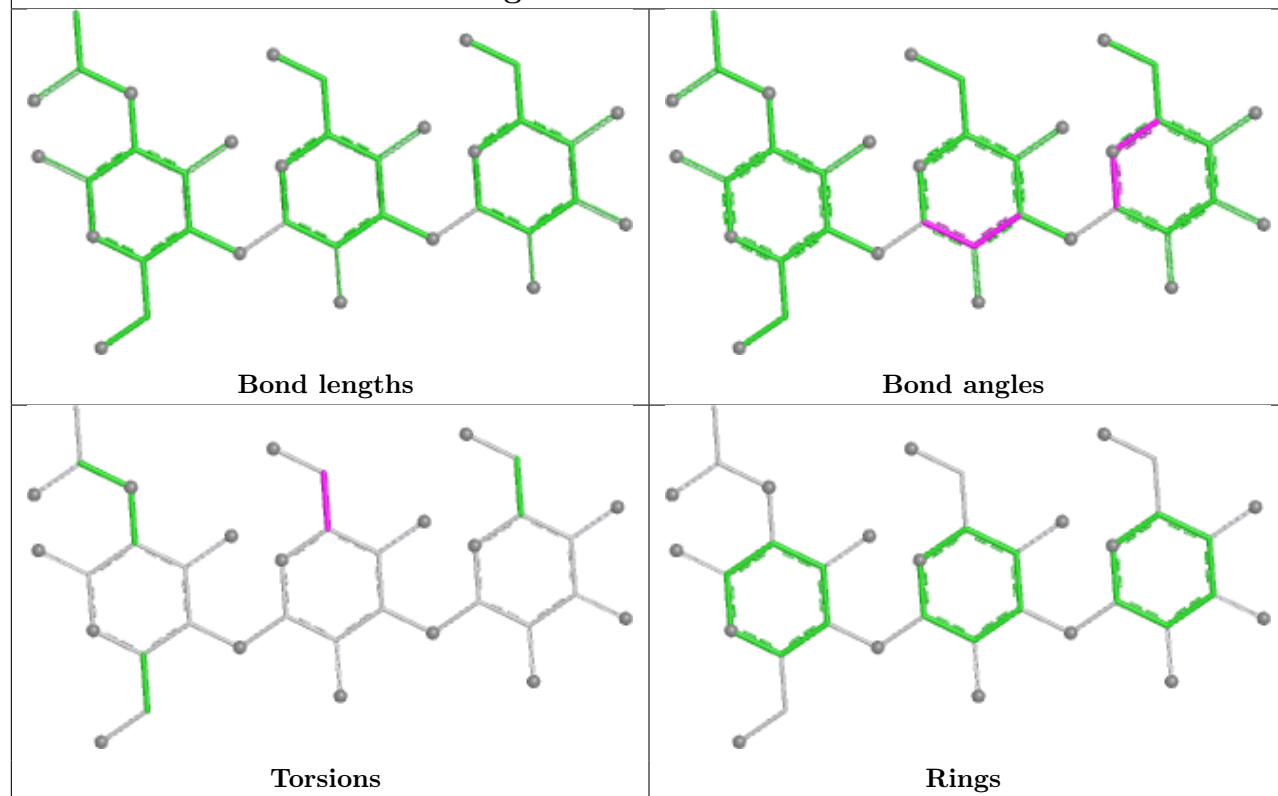
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

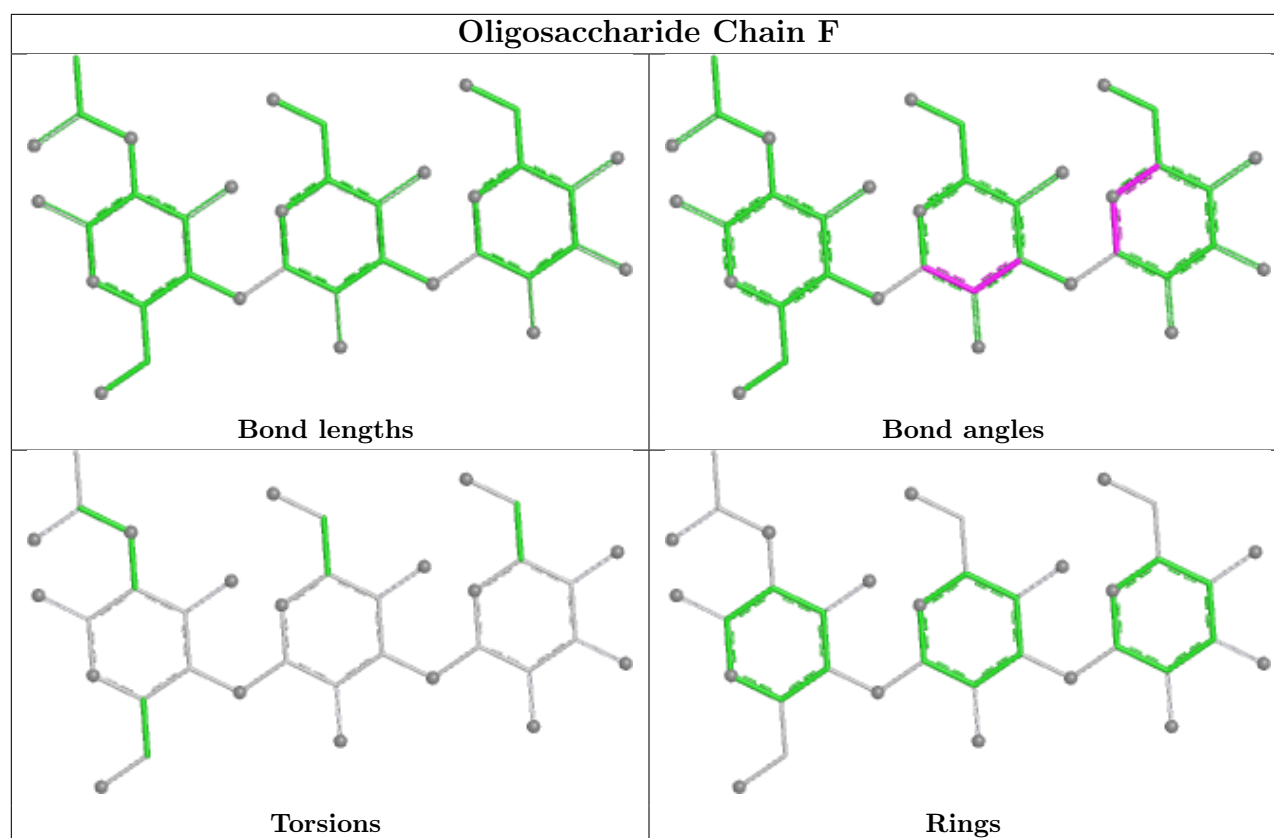


Oligosaccharide Chain D



Oligosaccharide Chain E





5.6 Ligand geometry [i](#)

6 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	$\# Z > 2$
3	GOL	A	503	-	5,5,5	0.15	0	5,5,5	0.45	0
3	GOL	A	501	-	5,5,5	0.35	0	5,5,5	0.25	0
3	GOL	B	502	-	5,5,5	0.32	0	5,5,5	0.30	0
3	GOL	A	504	-	5,5,5	0.37	0	5,5,5	0.27	0
3	GOL	B	506	-	5,5,5	0.36	0	5,5,5	0.28	0
3	GOL	A	505	-	5,5,5	0.31	0	5,5,5	0.30	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
3	GOL	A	503	-	-	2/4/4/4	-
3	GOL	A	501	-	-	2/4/4/4	-
3	GOL	B	502	-	-	0/4/4/4	-
3	GOL	A	504	-	-	4/4/4/4	-
3	GOL	B	506	-	-	2/4/4/4	-
3	GOL	A	505	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	GOL	C1-C2-C3-O3
3	A	501	GOL	O2-C2-C3-O3
3	A	503	GOL	O1-C1-C2-C3
3	A	505	GOL	C1-C2-C3-O3
3	A	505	GOL	O2-C2-C3-O3
3	A	504	GOL	O1-C1-C2-C3
3	A	504	GOL	C1-C2-C3-O3
3	A	505	GOL	O1-C1-C2-C3
3	B	506	GOL	C1-C2-C3-O3
3	A	503	GOL	O1-C1-C2-O2
3	A	504	GOL	O1-C1-C2-O2
3	B	506	GOL	O2-C2-C3-O3
3	A	504	GOL	O2-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	GOL	1	0
3	A	501	GOL	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	245/255 (96%)	0.32	29 (11%) 9 8	9, 20, 59, 74	0
1	B	254/255 (99%)	0.25	20 (7%) 18 17	11, 22, 43, 55	0
All	All	499/510 (97%)	0.28	49 (9%) 13 12	9, 21, 51, 74	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	17	ILE	6.8
1	A	29	ALA	5.9
1	A	26	THR	5.8
1	A	13	GLY	5.2
1	A	28	THR	5.2
1	A	38	ILE	5.1
1	A	27	ASN	5.0
1	B	261	TYR	4.9
1	A	111	HIS	4.7
1	A	25	ASN	4.6
1	A	14	MET	4.5
1	A	23	TYR	4.3
1	A	16	THR	4.1
1	A	15	ARG	4.0
1	A	39	ASP	3.8
1	A	19	GLY	3.8
1	A	110	LEU	3.4
1	B	17	ILE	3.4
1	B	111	HIS	3.3
1	B	90	ASN	3.2
1	A	112	ASP	3.2
1	A	24	PHE	3.2
1	B	110	LEU	3.1
1	B	94	ASN	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	67	ASN	2.9
1	B	18	ASP	2.7
1	B	38	ILE	2.7
1	A	31	ALA	2.6
1	A	21	LYS	2.6
1	B	96	ILE	2.5
1	A	30	GLU	2.5
1	A	71	ILE	2.5
1	A	61	ASP	2.5
1	A	60	ASN	2.5
1	B	40	GLY	2.4
1	B	50	SER	2.4
1	B	102	ARG	2.4
1	A	109	TYR	2.3
1	A	20	LYS	2.3
1	A	68	THR	2.3
1	B	16	THR	2.3
1	B	61	ASP	2.3
1	B	14	MET	2.3
1	B	93	ALA	2.2
1	B	47	THR	2.1
1	B	35	TRP	2.1
1	B	60	ASN	2.0
1	A	257	ALA	2.0
1	A	90	ASN	2.0

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

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

6.3 Carbohydrates ⓘ

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

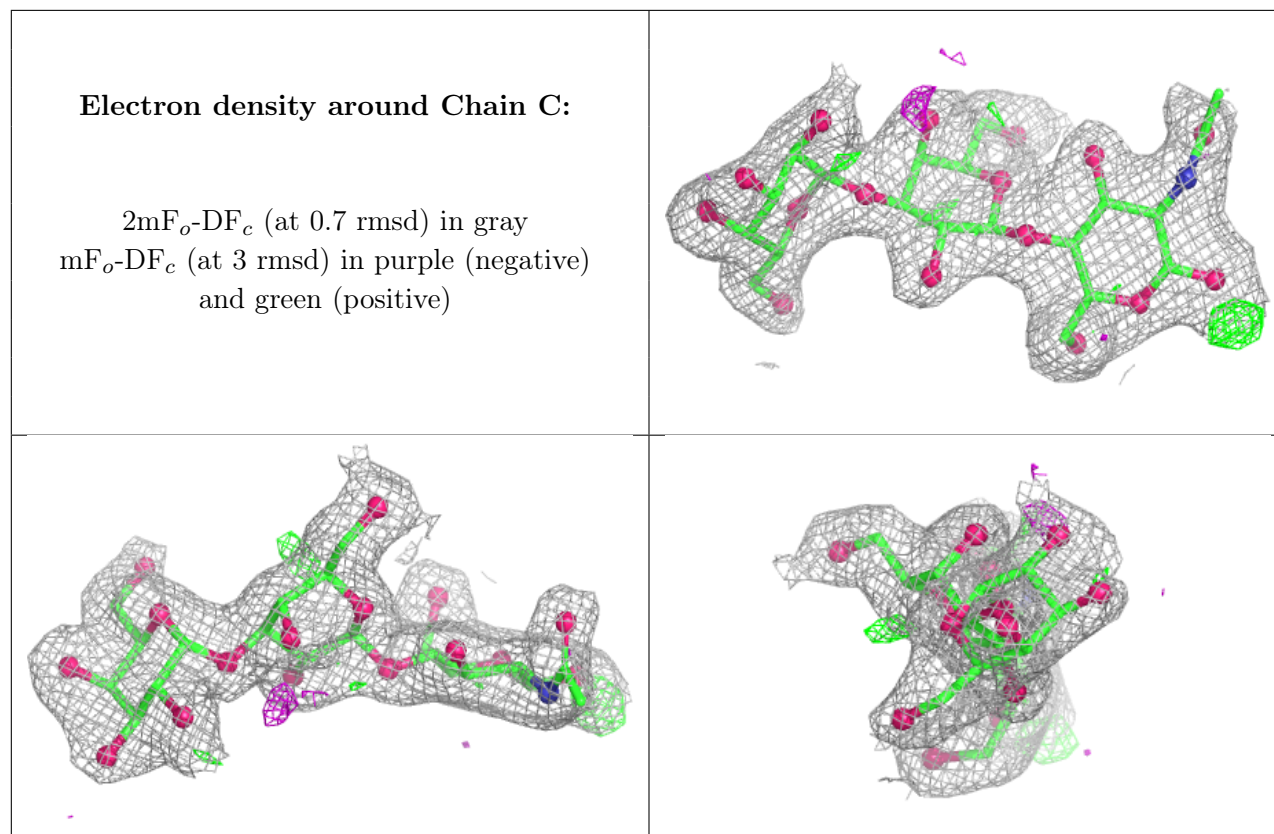
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GLA	E	3	11/12	0.82	0.14	34,35,39,40	0
2	GAL	E	2	11/12	0.83	0.13	32,34,35,36	0
2	GLA	C	3	11/12	0.87	0.11	33,35,39,39	0

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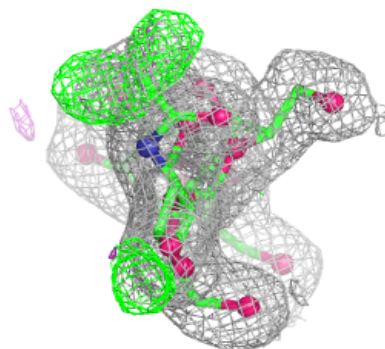
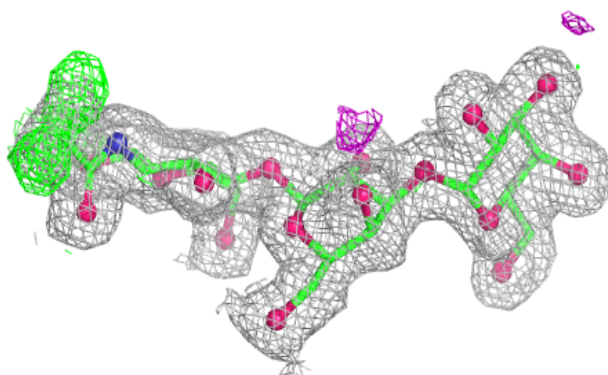
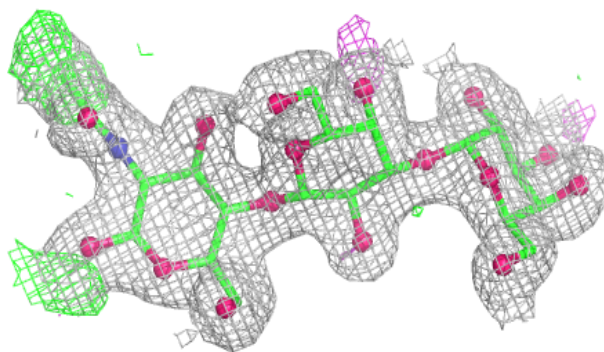
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GLA	F	3	11/12	0.87	0.11	21,24,30,31	0
2	NAG	E	1	15/15	0.90	0.12	30,32,40,43	0
2	GAL	C	2	11/12	0.90	0.09	25,27,31,32	0
2	NAG	F	1	15/15	0.91	0.08	18,20,24,25	0
2	NAG	C	1	15/15	0.91	0.11	26,30,40,40	0
2	NAG	D	1	15/15	0.94	0.06	9,12,14,15	0
2	GAL	F	2	11/12	0.96	0.06	15,18,21,23	0
2	GLA	D	3	11/12	0.96	0.06	12,14,16,17	0
2	GAL	D	2	11/12	0.97	0.04	6,10,12,14	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.

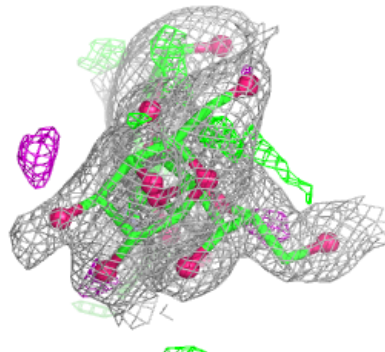
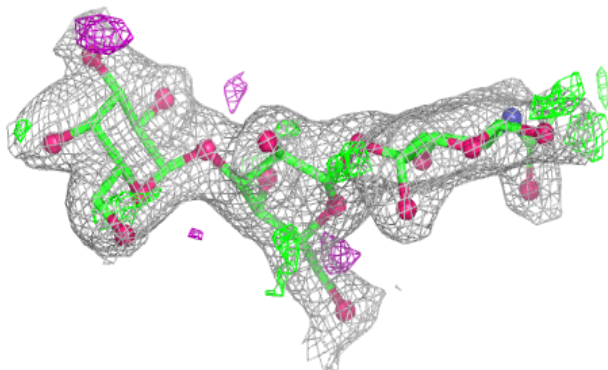
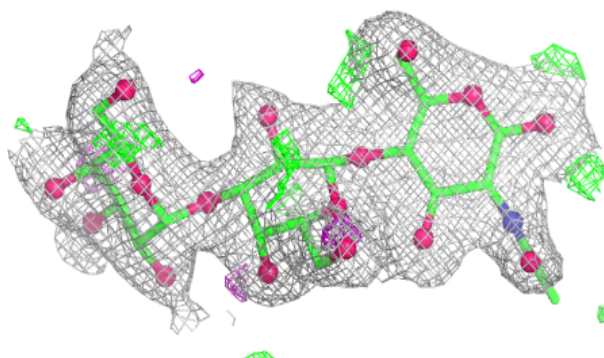


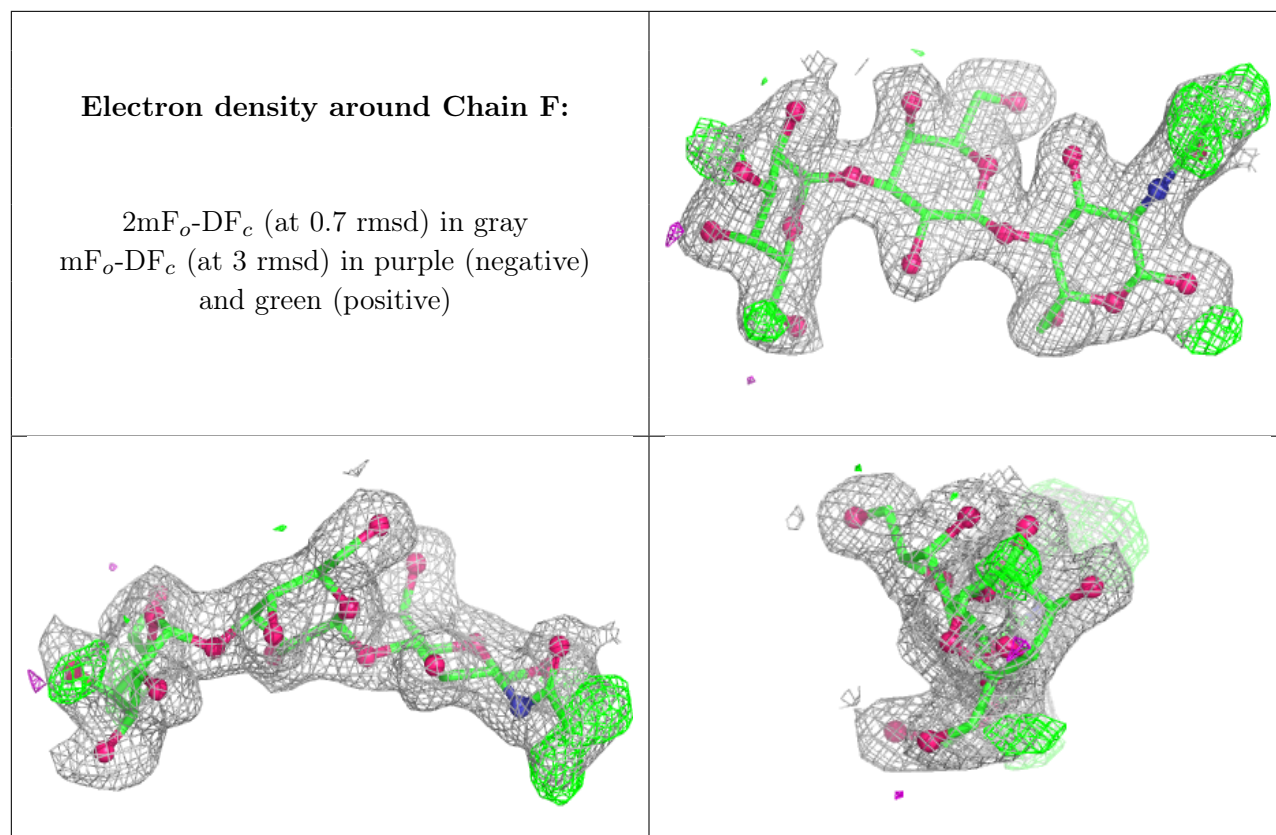
Electron density around Chain D:

$2mF_o-DF_c$ (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 [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
3	GOL	A	501	6/6	0.77	0.20	42,48,49,49	0
3	GOL	A	503	6/6	0.77	0.22	39,45,46,49	0
3	GOL	A	504	6/6	0.79	0.20	58,59,60,60	0
3	GOL	B	506	6/6	0.83	0.14	44,45,45,46	0
3	GOL	B	502	6/6	0.91	0.10	20,24,27,35	0
3	GOL	A	505	6/6	0.93	0.10	24,29,32,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.