



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:20 PM UTC

PDB ID : 8ZRZ / pdb_00008zrz
Title : The 1.26 angstrom resolution structure of Bacillus cereus beta-amylase in complex with maltose
Authors : Mikami, B.; Hirata, A.
Deposited on : 2024-06-05
Resolution : 1.26 Å(reported)

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

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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

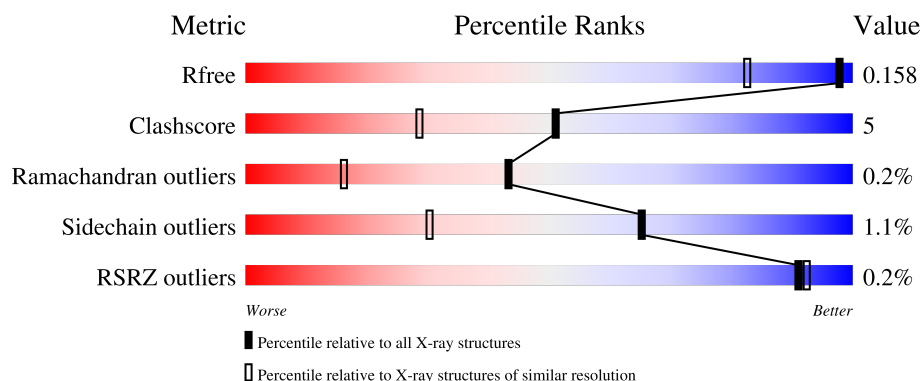
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.26 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	1693 (1.28-1.24)
Clashscore	190562	1730 (1.28-1.24)
Ramachandran outliers	187476	1695 (1.28-1.24)
Sidechain outliers	187428	1694 (1.28-1.24)
RSRZ outliers	180081	1693 (1.28-1.24)

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	516	89% 11%
2	C	2	50% 50%
2	D	2	100%
2	F	2	100%
3	E	2	50% 50%

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Mol	Chain	Length	Quality of chain
4	B	4	

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
4	GLC	B	3[B]	X	-	-	-
7	GOL	A	604	-	X	-	-
7	GOL	A	605	-	X	-	-
7	GOL	A	606[A]	-	X	-	-
7	GOL	A	606[B]	-	X	-	-
7	GOL	A	607	-	X	-	-
7	GOL	A	608	-	X	-	-
7	GOL	A	609	-	X	-	-
7	GOL	A	610	-	X	-	-
7	GOL	A	611	-	X	-	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 5480 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 Beta-amylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	516	Total	C	N	O	S	0	46	0
			4371	2819	706	824	22			

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	F	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	E	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 4 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

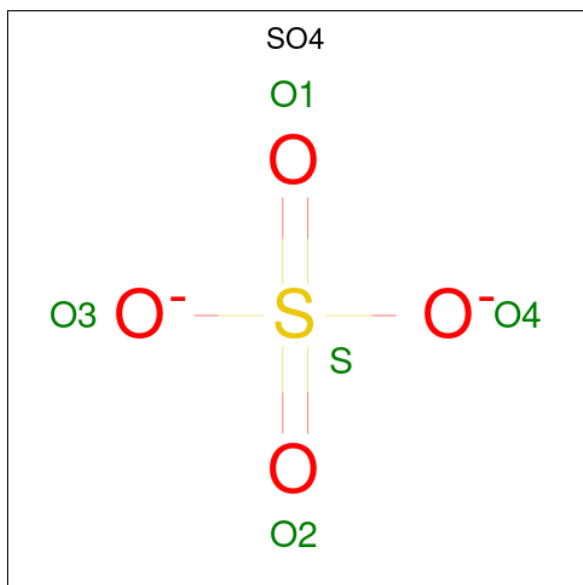


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	B	4	Total	C	O	0	2	0
			58	30	28			

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

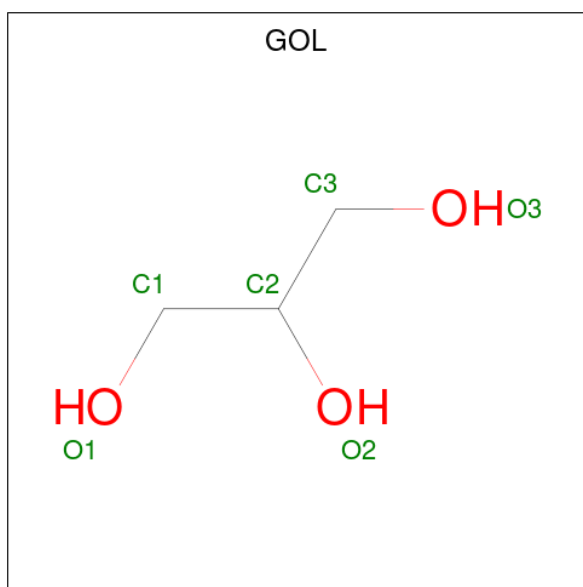
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



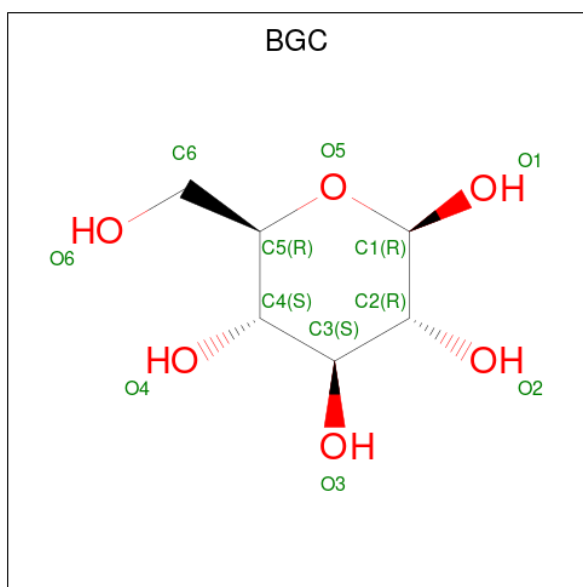
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	1
			12	6	6		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			12	6	6		

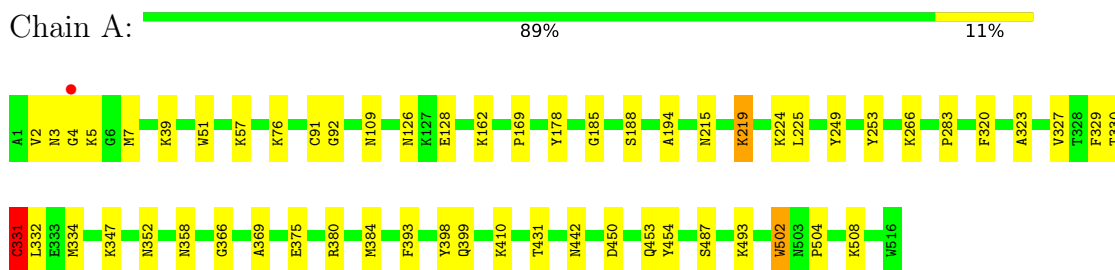
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	867	Total	O	0	20
			887	887		

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.

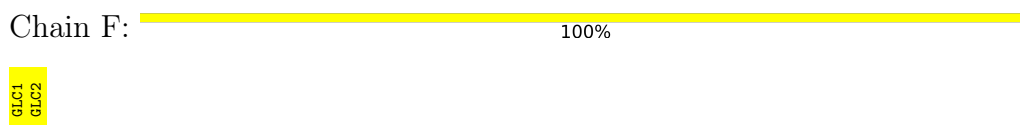
- Molecule 1: Beta-amylase



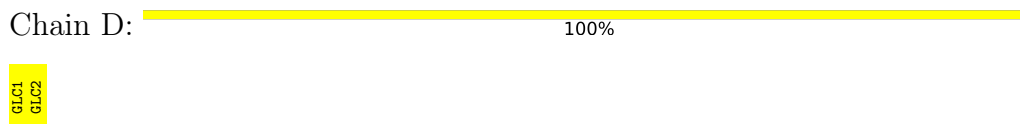
- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



- Molecule 3: alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose



- Molecule 4: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain B:  25% 25% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.97Å 89.84Å 66.03Å 90.00° 103.09° 90.00°	Depositor
Resolution (Å)	10.00 – 1.26 10.00 – 1.26	Depositor EDS
% Data completeness (in resolution range)	93.4 (10.00-1.26) 95.3 (10.00-1.26)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.63 (at 1.26Å)	Xtriage
Refinement program	SHELXL-97 2018/3	Depositor
R, R_{free}	0.124 , 0.165 0.121 , 0.158	Depositor DCC
R_{free} test set	8582 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	12.4	Xtriage
Anisotropy	0.611	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 80.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	5480	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.05% 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: GLC, CA, GOL, BGC, 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	0.91	3/4618 (0.1%)	1.26	17/6267 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	4	GLY	C-N	10.04	1.47	1.33
1	A	3	ASN	C-N	9.35	1.46	1.33
1	A	5	LYS	C-N	-6.64	1.26	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	358	ASN	OD1-CG-ND2	-10.50	112.10	122.60
1	A	331[A]	CYS	CA-CB-SG	8.26	133.40	114.40
1	A	331[B]	CYS	CA-CB-SG	8.26	133.40	114.40
1	A	109	ASN	CA-CB-CG	-7.42	105.18	112.60
1	A	393	PHE	CA-CB-CG	-6.52	107.28	113.80
1	A	332[A]	LEU	CA-C-N	-6.12	112.44	122.54
1	A	332[A]	LEU	C-N-CA	-6.12	112.44	122.54
1	A	332[B]	LEU	CA-C-N	-6.12	112.44	122.54
1	A	332[B]	LEU	C-N-CA	-6.12	112.44	122.54
1	A	215	ASN	OD1-CG-ND2	-6.05	116.55	122.60
1	A	92	GLY	CA-C-N	-5.97	114.14	122.86
1	A	92	GLY	C-N-CA	-5.97	114.14	122.86
1	A	380[A]	ARG	CA-C-O	5.75	126.65	120.55
1	A	380[B]	ARG	CA-C-O	5.75	126.65	120.55
1	A	431	THR	CA-C-O	5.47	127.76	121.47
1	A	178	TYR	CA-CB-CG	5.33	123.50	113.90
1	A	442	ASN	CA-CB-CG	5.30	117.90	112.60

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	4371	0	4263	41	0
2	C	23	0	21	1	0
2	D	23	0	21	0	0
2	F	23	0	21	0	0
3	E	23	0	21	0	0
4	B	58	0	47	3	0
5	A	1	0	0	0	0
6	A	5	0	0	0	0
7	A	54	0	36	7	0
8	A	12	0	12	2	0
9	A	887	0	0	14	0
All	All	5480	0	4442	47	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 (47) 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)
4:B:2[A]:GLC:H62	4:B:3[A]:GLC:O1	1.85	0.77
1:A:331[A]:CYS:HB2	1:A:369:ALA:HB2	1.67	0.76
1:A:493:LYS:HE2	9:A:1284:HOH:O	1.85	0.76
1:A:225:LEU:HA	9:A:1271:HOH:O	1.85	0.75
1:A:331[B]:CYS:HB2	9:A:946:HOH:O	1.93	0.69
1:A:487[B]:SER:OG	1:A:493:LYS:HD2	1.93	0.68
9:A:1104:HOH:O	4:B:3[A]:GLC:H3	1.92	0.68
7:A:605:GOL:H12	9:A:916:HOH:O	1.97	0.63
1:A:352[A]:ASN:ND2	7:A:608:GOL:O2	2.27	0.62
1:A:502:TRP:NE1	8:A:602:BGC:H6C1	2.14	0.62
1:A:266[C]:LYS:HD2	1:A:323:ALA:HB2	1.82	0.62
1:A:266[C]:LYS:HE2	1:A:323:ALA:N	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:GLU:HB2	1:A:410:LYS:HD3	1.82	0.60
1:A:508:LYS:HD3	9:A:1473[B]:HOH:O	2.00	0.59
7:A:606[A]:GOL:H12	9:A:728:HOH:O	2.03	0.57
1:A:502:TRP:HE1	8:A:602:BGC:H6C1	1.69	0.57
1:A:2[B]:VAL:CG1	1:A:454[B]:TYR:CE1	2.88	0.56
1:A:2[B]:VAL:CG1	1:A:454[B]:TYR:HE1	2.19	0.56
1:A:2[B]:VAL:HG12	1:A:454[B]:TYR:CE1	2.41	0.55
1:A:320:PHE:CD2	1:A:327[A]:VAL:HG22	2.40	0.55
1:A:347:LYS:HE3	1:A:384[B]:MET:SD	2.48	0.54
1:A:225:LEU:HD23	9:A:1271:HOH:O	2.09	0.52
1:A:162[A]:LYS:HG3	1:A:283:PRO:O	2.09	0.52
1:A:219[A]:LYS:HD3	9:A:809:HOH:O	2.11	0.49
1:A:224:LYS:HG3	7:A:610:GOL:C1	2.43	0.48
1:A:334[A]:MET:HE1	9:A:1250:HOH:O	2.13	0.48
1:A:126:ASN:OD1	1:A:128[B]:GLU:HG2	2.14	0.48
1:A:329[B]:PHE:O	1:A:366:GLY:HA2	2.13	0.48
7:A:609:GOL:H12	9:A:1081:HOH:O	2.12	0.48
1:A:51:TRP:CZ2	1:A:91:CYS:SG	3.07	0.48
1:A:57:LYS:HG3	9:A:775:HOH:O	2.14	0.47
1:A:185:GLY:O	1:A:188[B]:SER:OG	2.29	0.47
1:A:219[A]:LYS:HA	7:A:604:GOL:C2	2.46	0.46
1:A:194:ALA:HB3	1:A:253[A]:TYR:OH	2.16	0.46
1:A:76:LYS:HD2	9:A:1198:HOH:O	2.15	0.45
1:A:508:LYS:HB3	1:A:508:LYS:HE3	1.66	0.43
1:A:249:TYR:CE2	2:C:1:GLC:H4	2.53	0.43
1:A:450[B]:ASP:OD2	1:A:453:GLN:N	2.48	0.43
1:A:39:LYS:NZ	9:A:729:HOH:O	2.52	0.42
1:A:128[B]:GLU:CD	1:A:169:PRO:HB3	2.44	0.42
1:A:219[B]:LYS:HA	7:A:604:GOL:C2	2.51	0.41

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	561/516 (109%)	546 (97%)	14 (2%)	1 (0%)	43 16

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	398	TYR

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	487/440 (111%)	479 (98%)	8 (2%)	55 19

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7[A]	MET
1	A	7[B]	MET
1	A	219[A]	LYS
1	A	219[B]	LYS
1	A	331[A]	CYS
1	A	331[B]	CYS
1	A	502	TRP
1	A	504	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	ASN
1	A	109	ASN
1	A	215	ASN
1	A	239	GLN
1	A	243	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 ⓘ

14 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$
4	GLC	B	1	4	12,12,12	0.53	0	17,17,17	1.06	1 (5%)
4	GLC	B	2[A]	-	11,11,12	0.70	0	15,15,17	1.86	4 (26%)
4	GLC	B	2[B]	-	11,11,12	0.70	0	15,15,17	1.56	3 (20%)
4	GLC	B	3[A]	4	12,12,12	0.62	0	17,17,17	1.95	5 (29%)
4	GLC	B	3[B]	-	12,12,12	1.21	2 (16%)	17,17,17	2.98	7 (41%)
4	GLC	B	4	4	11,11,12	0.63	0	15,15,17	0.72	0
2	GLC	C	1	2	12,12,12	0.61	0	17,17,17	1.06	1 (5%)
2	GLC	C	2	2	11,11,12	0.79	1 (9%)	15,15,17	0.93	1 (6%)
2	GLC	D	1	2	12,12,12	0.68	0	17,17,17	0.98	1 (5%)
2	GLC	D	2	2	11,11,12	0.63	0	15,15,17	1.46	3 (20%)
3	BGC	E	1	3	12,12,12	0.58	0	17,17,17	1.16	1 (5%)
3	GLC	E	2	3	11,11,12	0.76	0	15,15,17	1.04	0
2	GLC	F	1	2	12,12,12	0.72	0	17,17,17	1.22	2 (11%)
2	GLC	F	2	2	11,11,12	0.63	0	15,15,17	1.29	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
4	GLC	B	1	4	-	0/2/22/22	0/1/1/1
4	GLC	B	2[A]	-	-	0/2/19/22	0/1/1/1
4	GLC	B	2[B]	-	-	0/2/19/22	0/1/1/1
4	GLC	B	3[A]	4	-	0/2/22/22	0/1/1/1
4	GLC	B	3[B]	-	1/1/5/5	0/2/22/22	0/1/1/1
4	GLC	B	4	4	-	0/2/19/22	0/1/1/1
2	GLC	C	1	2	-	0/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	2/2/19/22	0/1/1/1
3	BGC	E	1	3	-	0/2/22/22	0/1/1/1
3	GLC	E	2	3	-	0/2/19/22	0/1/1/1
2	GLC	F	1	2	-	0/2/22/22	0/1/1/1
2	GLC	F	2	2	-	0/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	3[B]	GLC	O5-C1	-2.71	1.36	1.42
2	C	2	GLC	O5-C1	-2.28	1.39	1.43
4	B	3[B]	GLC	O3-C3	2.15	1.48	1.43

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	3[B]	GLC	O1-C1-O5	6.93	130.98	110.41
4	B	3[B]	GLC	O5-C5-C4	-5.52	99.75	109.70
4	B	3[B]	GLC	C4-C3-C2	-5.08	101.91	110.83
4	B	2[A]	GLC	O6-C6-C5	3.92	124.67	111.33
4	B	3[A]	GLC	O4-C4-C3	3.53	118.69	110.38
4	B	2[A]	GLC	C1-C2-C3	3.50	114.73	109.64
4	B	2[B]	GLC	C1-C2-C3	3.50	114.73	109.64
4	B	3[B]	GLC	O1-C1-C2	3.37	118.76	108.98
4	B	3[A]	GLC	O5-C5-C4	-3.20	103.92	109.70
4	B	3[B]	GLC	O3-C3-C2	-3.16	102.93	110.38
4	B	3[B]	GLC	C1-O5-C5	3.15	119.74	113.65
4	B	3[A]	GLC	C4-C3-C2	3.04	116.17	110.83
3	E	1	BGC	O1-C1-O5	-2.98	101.55	110.41
4	B	2[A]	GLC	O3-C3-C2	-2.95	104.02	110.05
4	B	2[B]	GLC	O3-C3-C2	-2.95	104.02	110.05
4	B	3[A]	GLC	C6-C5-C4	-2.83	106.08	113.02
2	C	1	GLC	C4-C3-C2	-2.72	106.05	110.83
2	D	2	GLC	C1-O5-C5	2.72	115.83	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	GLC	C1-O5-C5	2.56	118.61	113.65
2	F	1	GLC	O4-C4-C3	-2.42	104.67	110.38
2	D	1	GLC	C3-C4-C5	2.31	114.42	110.23
4	B	2[A]	GLC	O5-C5-C6	-2.25	103.29	107.66
4	B	2[B]	GLC	O5-C5-C6	-2.25	103.29	107.66
2	F	2	GLC	O3-C3-C2	-2.19	105.59	110.05
2	C	2	GLC	O5-C5-C6	-2.16	103.45	107.66
4	B	3[A]	GLC	O4-C4-C5	-2.12	104.10	109.32
4	B	1	GLC	O3-C3-C4	-2.12	105.38	110.38
2	D	2	GLC	O5-C5-C4	-2.09	105.75	110.83
4	B	3[B]	GLC	C6-C5-C4	2.07	118.11	113.02
2	D	2	GLC	C1-C2-C3	-2.05	106.65	109.64

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	3[B]	GLC	C1

All (2) torsion outliers are listed below:

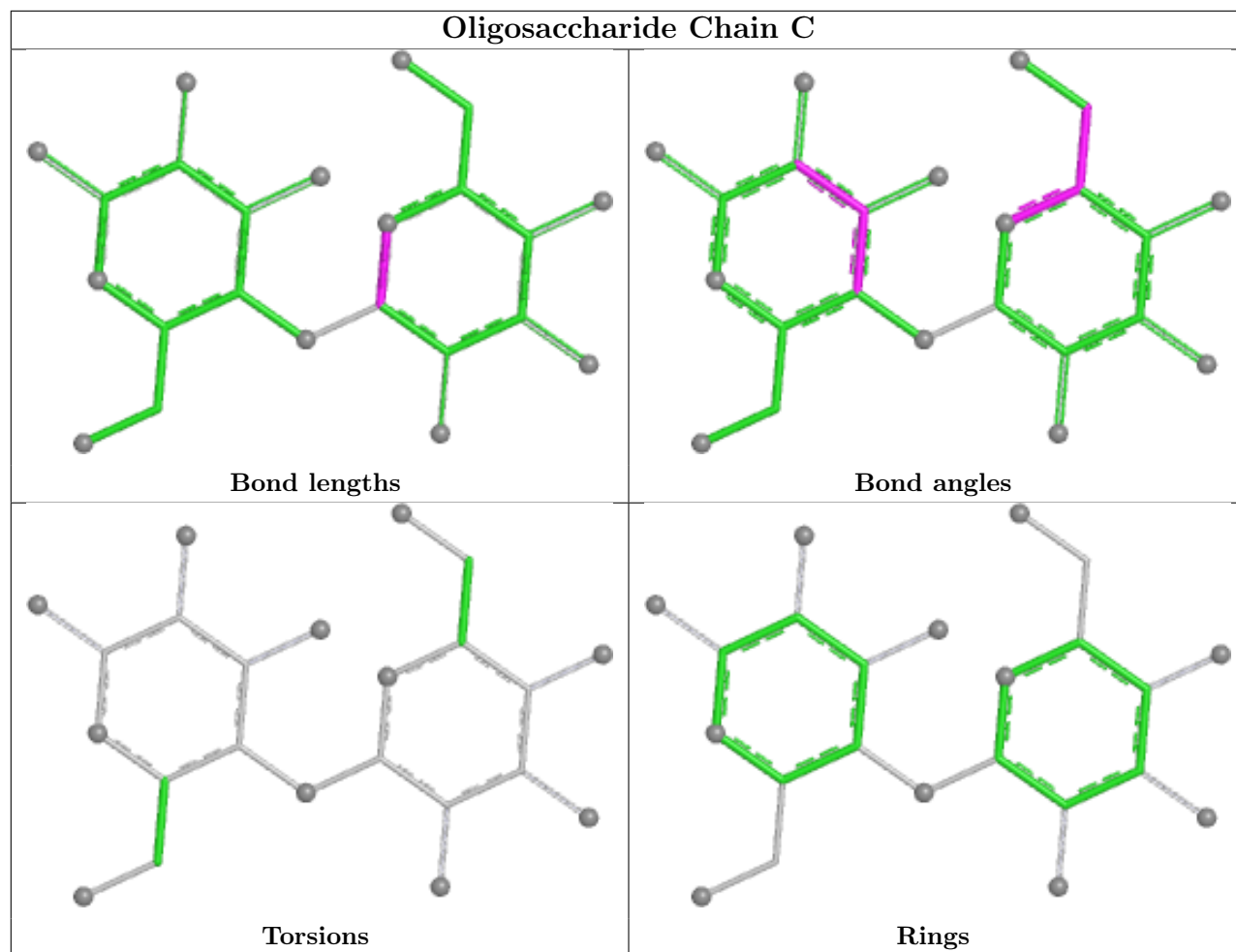
Mol	Chain	Res	Type	Atoms
2	D	2	GLC	O5-C5-C6-O6
2	D	2	GLC	C4-C5-C6-O6

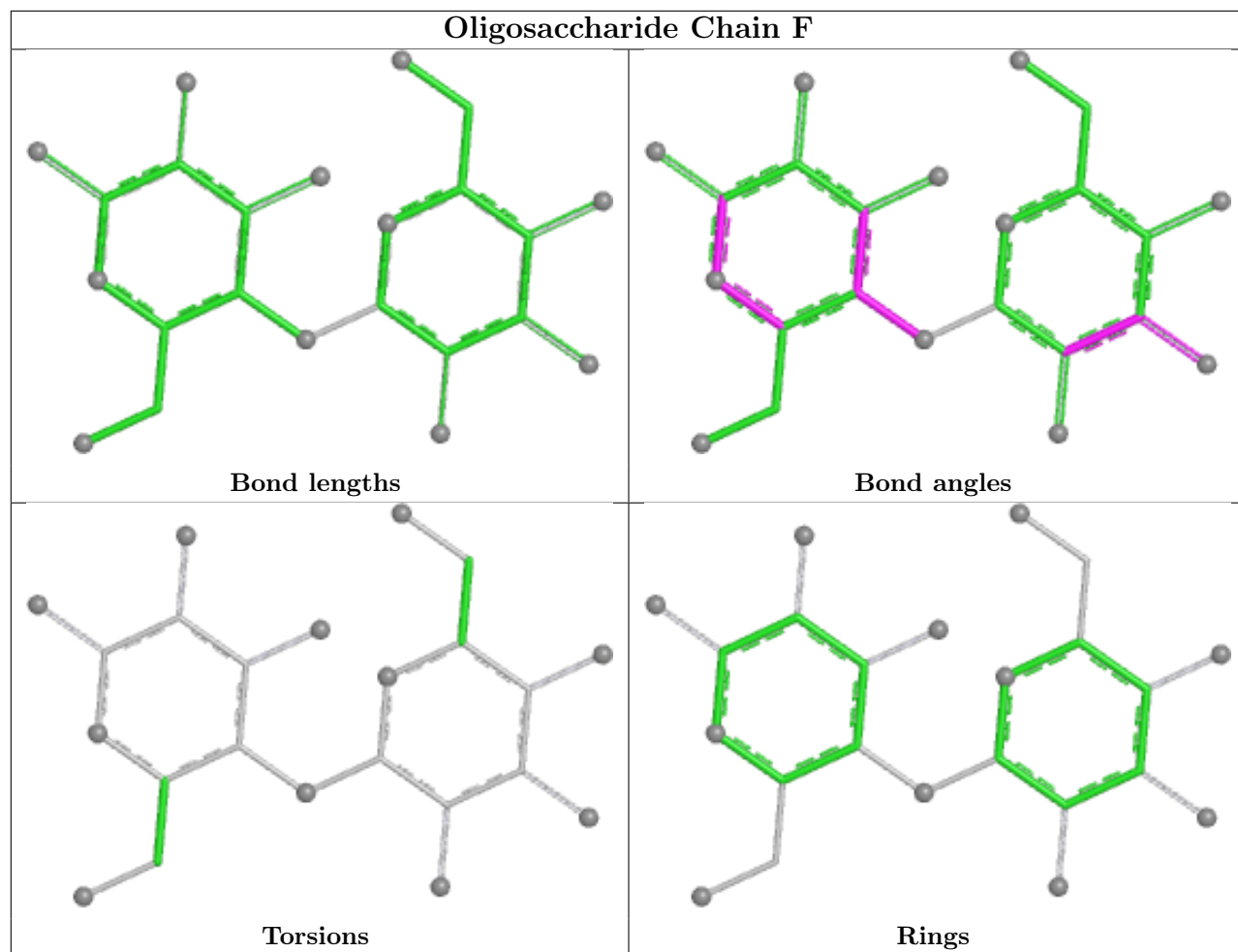
There are no ring outliers.

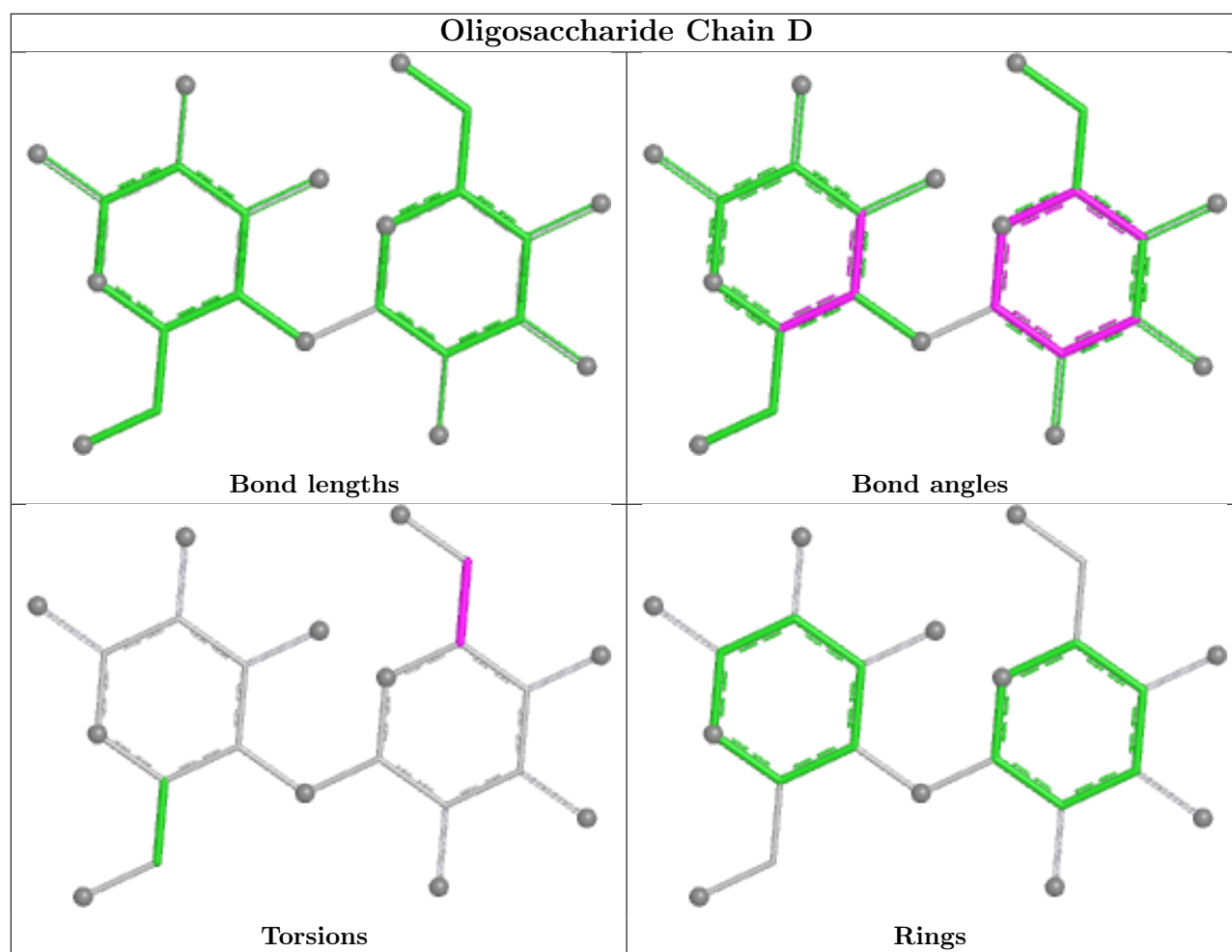
4 monomers are involved in 4 short contacts:

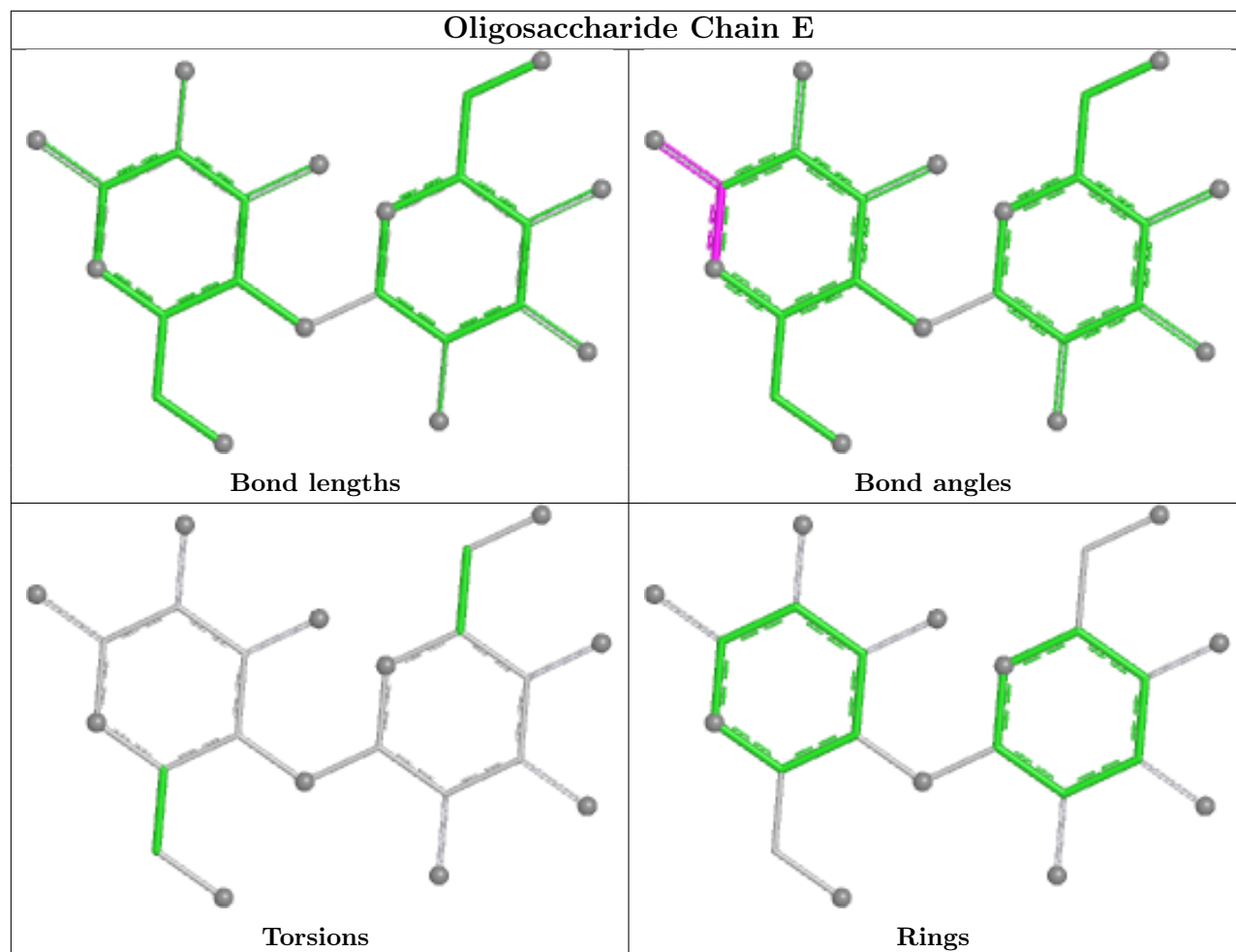
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	3[B]	GLC	1	0
4	B	3[A]	GLC	2	0
2	C	1	GLC	1	0
4	B	2[A]	GLC	1	0

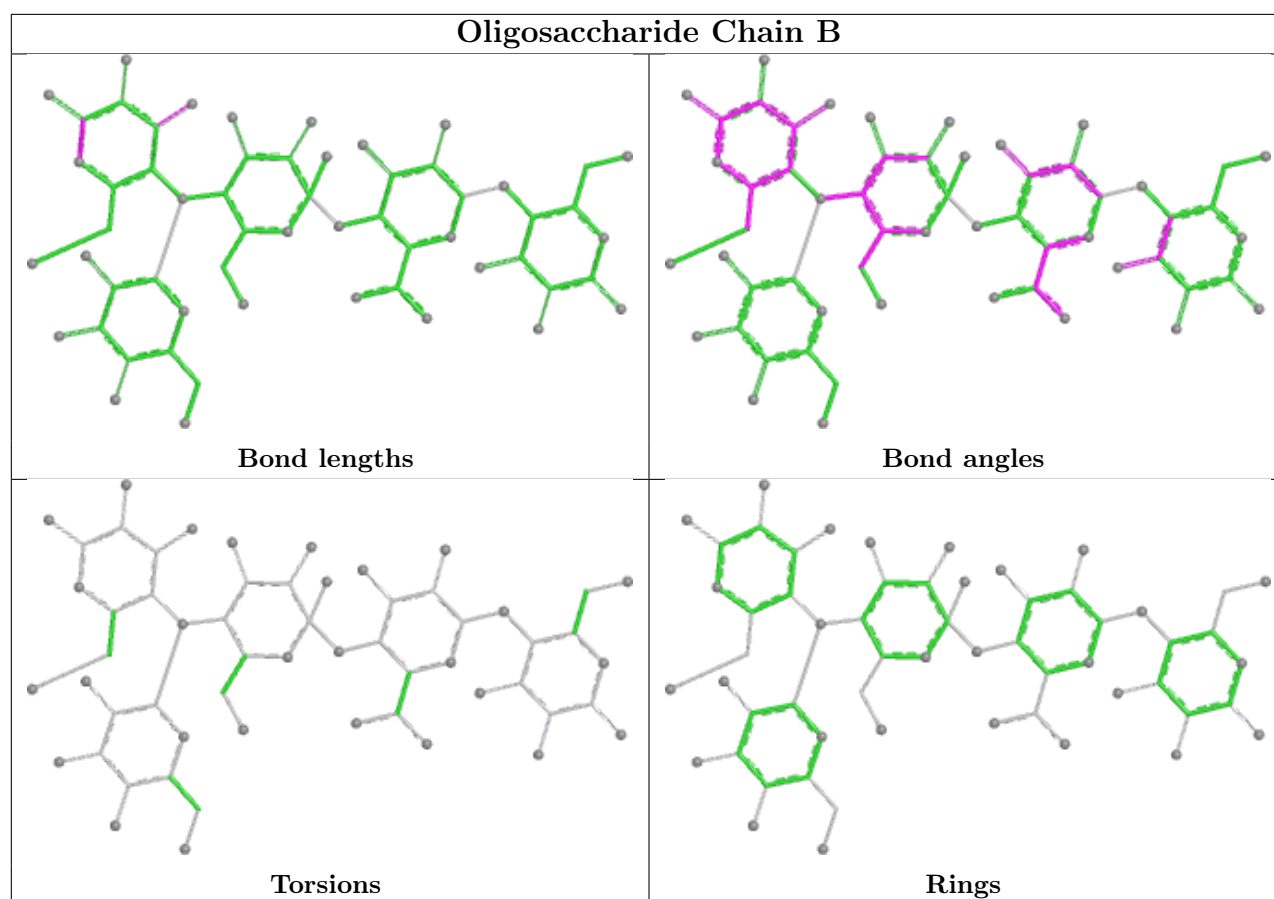
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 1 is monoatomic - leaving 11 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
7	GOL	A	606[B]	-	5,5,5	4.73	5 (100%)	5,5,5	4.42	3 (60%)
8	BGC	A	602	-	12,12,12	0.53	0	17,17,17	1.32	1 (5%)
7	GOL	A	607	-	5,5,5	4.71	5 (100%)	5,5,5	5.85	3 (60%)
7	GOL	A	610	-	5,5,5	4.71	5 (100%)	5,5,5	6.27	4 (80%)
7	GOL	A	604	-	5,5,5	4.52	5 (100%)	5,5,5	3.69	3 (60%)
7	GOL	A	609	-	5,5,5	4.65	5 (100%)	5,5,5	3.51	3 (60%)
6	SO4	A	603	-	4,4,4	0.81	0	6,6,6	0.88	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	GOL	A	608	-	5,5,5	4.50	5 (100%)	5,5,5	4.37	3 (60%)
7	GOL	A	606[A]	-	5,5,5	4.79	5 (100%)	5,5,5	4.76	3 (60%)
7	GOL	A	605	-	5,5,5	4.58	5 (100%)	5,5,5	3.92	3 (60%)
7	GOL	A	611	-	5,5,5	4.77	5 (100%)	5,5,5	6.29	3 (60%)

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	GOL	A	606[B]	-	-	1/4/4/4	-
8	BGC	A	602	-	-	2/2/22/22	0/1/1/1
7	GOL	A	607	-	-	1/4/4/4	-
7	GOL	A	610	-	-	1/4/4/4	-
7	GOL	A	604	-	-	4/4/4/4	-
7	GOL	A	609	-	-	3/4/4/4	-
7	GOL	A	608	-	-	2/4/4/4	-
7	GOL	A	606[A]	-	-	3/4/4/4	-
7	GOL	A	605	-	-	3/4/4/4	-
7	GOL	A	611	-	-	3/4/4/4	-

All (45) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	607	GOL	C3-C2	-7.92	1.21	1.51
7	A	611	GOL	C3-C2	-7.91	1.21	1.51
7	A	610	GOL	C3-C2	-7.79	1.22	1.51
7	A	606[A]	GOL	C3-C2	-7.75	1.22	1.51
7	A	606[B]	GOL	C3-C2	-7.45	1.23	1.51
7	A	605	GOL	C3-C2	-7.44	1.23	1.51
7	A	608	GOL	C3-C2	-7.34	1.23	1.51
7	A	609	GOL	C3-C2	-7.30	1.24	1.51
7	A	604	GOL	C3-C2	-7.25	1.24	1.51
7	A	606[A]	GOL	O1-C1	4.83	1.62	1.42
7	A	609	GOL	O1-C1	4.76	1.62	1.42
7	A	604	GOL	O1-C1	4.68	1.62	1.42
7	A	611	GOL	O1-C1	4.65	1.62	1.42
7	A	605	GOL	O1-C1	4.63	1.61	1.42
7	A	606[B]	GOL	O1-C1	4.60	1.61	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	610	GOL	C1-C2	-4.46	1.34	1.51
7	A	608	GOL	O1-C1	4.43	1.61	1.42
7	A	607	GOL	O1-C1	4.29	1.60	1.42
7	A	606[B]	GOL	O3-C3	4.05	1.59	1.42
7	A	606[A]	GOL	O3-C3	3.89	1.58	1.42
7	A	605	GOL	O3-C3	3.80	1.58	1.42
7	A	610	GOL	O3-C3	3.73	1.58	1.42
7	A	604	GOL	O3-C3	3.73	1.58	1.42
7	A	611	GOL	O3-C3	3.68	1.57	1.42
7	A	609	GOL	O3-C3	3.62	1.57	1.42
7	A	608	GOL	O3-C3	3.52	1.57	1.42
7	A	607	GOL	O3-C3	3.48	1.57	1.42
7	A	610	GOL	O1-C1	3.46	1.56	1.42
7	A	606[B]	GOL	C1-C2	-3.39	1.38	1.51
7	A	609	GOL	C1-C2	-3.29	1.39	1.51
7	A	608	GOL	C1-C2	-3.17	1.39	1.51
7	A	606[A]	GOL	C1-C2	-3.00	1.40	1.51
7	A	607	GOL	C1-C2	-2.99	1.40	1.51
7	A	607	GOL	O2-C2	-2.98	1.34	1.43
7	A	605	GOL	C1-C2	-2.94	1.40	1.51
7	A	609	GOL	O2-C2	-2.91	1.34	1.43
7	A	611	GOL	C1-C2	-2.90	1.40	1.51
7	A	604	GOL	O2-C2	-2.78	1.35	1.43
7	A	611	GOL	O2-C2	-2.76	1.35	1.43
7	A	606[B]	GOL	O2-C2	-2.71	1.35	1.43
7	A	606[A]	GOL	O2-C2	-2.67	1.35	1.43
7	A	604	GOL	C1-C2	-2.45	1.42	1.51
7	A	608	GOL	O2-C2	-2.35	1.36	1.43
7	A	605	GOL	O2-C2	-2.20	1.37	1.43
7	A	610	GOL	O2-C2	-2.11	1.37	1.43

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	611	GOL	O3-C3-C2	11.03	160.02	110.38
7	A	610	GOL	O1-C1-C2	10.97	159.77	110.38
7	A	607	GOL	O3-C3-C2	10.18	156.22	110.38
7	A	606[A]	GOL	O3-C3-C2	8.21	147.33	110.38
7	A	611	GOL	O2-C2-C3	8.01	142.34	109.18
7	A	608	GOL	O3-C3-C2	7.48	144.06	110.38
7	A	606[B]	GOL	O3-C3-C2	7.06	142.17	110.38
7	A	607	GOL	O2-C2-C3	6.93	137.89	109.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	604	GOL	O3-C3-C2	6.76	140.81	110.38
7	A	610	GOL	O3-C3-C2	6.45	139.43	110.38
7	A	605	GOL	O3-C3-C2	6.31	138.76	110.38
7	A	606[A]	GOL	O2-C2-C3	6.14	134.58	109.18
7	A	606[B]	GOL	O2-C2-C3	5.61	132.39	109.18
7	A	605	GOL	O2-C2-C3	5.43	131.66	109.18
7	A	610	GOL	O2-C2-C3	5.34	131.30	109.18
7	A	609	GOL	O2-C2-C3	5.07	130.15	109.18
7	A	609	GOL	O3-C3-C2	4.92	132.55	110.38
7	A	608	GOL	O1-C1-C2	4.50	130.64	110.38
7	A	608	GOL	O2-C2-C3	4.23	126.70	109.18
7	A	607	GOL	O1-C1-C2	4.14	129.01	110.38
7	A	606[B]	GOL	O1-C1-C2	3.90	127.95	110.38
7	A	604	GOL	O2-C2-C3	3.48	123.57	109.18
7	A	611	GOL	O1-C1-C2	3.36	125.49	110.38
7	A	609	GOL	O1-C1-C2	3.17	124.63	110.38
7	A	606[A]	GOL	O1-C1-C2	2.80	122.98	110.38
7	A	605	GOL	O1-C1-C2	2.51	121.69	110.38
7	A	604	GOL	O1-C1-C2	2.51	121.68	110.38
8	A	602	BGC	O5-C5-C6	-2.20	100.99	106.44
7	A	610	GOL	O2-C2-C1	2.14	118.04	109.18

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	611	GOL	C1-C2-C3-O3
7	A	604	GOL	O1-C1-C2-C3
7	A	604	GOL	C1-C2-C3-O3
7	A	605	GOL	O1-C1-C2-C3
7	A	606[A]	GOL	O1-C1-C2-C3
7	A	606[A]	GOL	C1-C2-C3-O3
7	A	608	GOL	C1-C2-C3-O3
7	A	609	GOL	O1-C1-C2-C3
7	A	604	GOL	O1-C1-C2-O2
7	A	604	GOL	O2-C2-C3-O3
7	A	607	GOL	C1-C2-C3-O3
7	A	610	GOL	O1-C1-C2-C3
7	A	606[B]	GOL	O2-C2-C3-O3
7	A	608	GOL	O2-C2-C3-O3
8	A	602	BGC	O5-C5-C6-O6
7	A	606[A]	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
7	A	605	GOL	O1-C1-C2-O2
7	A	605	GOL	O2-C2-C3-O3
7	A	609	GOL	O2-C2-C3-O3
7	A	609	GOL	O1-C1-C2-O2
7	A	611	GOL	O1-C1-C2-C3
8	A	602	BGC	C4-C5-C6-O6
7	A	611	GOL	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	602	BGC	2	0
7	A	610	GOL	1	0
7	A	604	GOL	2	0
7	A	609	GOL	1	0
7	A	608	GOL	1	0
7	A	606[A]	GOL	1	0
7	A	605	GOL	1	0

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	516/516 (100%)	-0.58	1 (0%) 91 92	7, 17, 30, 55	46 (8%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	4	GLY	2.3

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

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

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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.

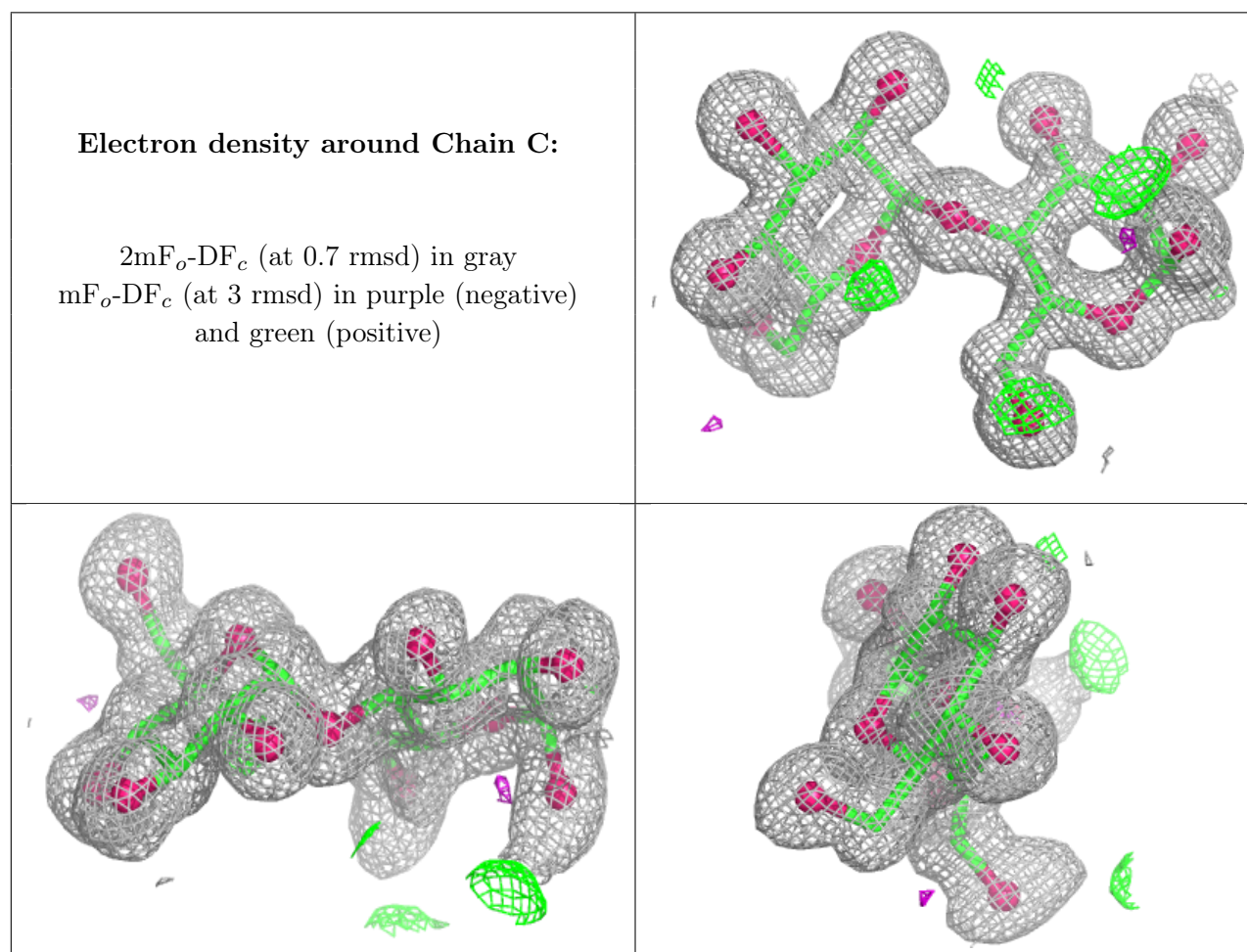
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	D	2	11/12	0.91	0.09	17,21,31,47	11
3	GLC	E	2	11/12	0.94	0.07	19,20,29,37	11
2	GLC	F	1	12/12	0.97	0.06	27,31,39,43	0
2	GLC	F	2	11/12	0.97	0.05	24,25,26,31	0
2	GLC	D	1	12/12	-	-	18,21,24,34	12
2	GLC	C	1	12/12	0.97	0.04	14,16,20,26	0
3	BGC	E	1	12/12	-	-	14,16,18,19	12
2	GLC	C	2	11/12	0.97	0.05	13,14,15,17	0
4	GLC	B	1	12/12	0.98	0.04	16,19,26,26	12
4	GLC	B	2[A]	11/12	0.99	0.03	13,15,16,23	11
4	GLC	B	2[B]	11/12	0.99	0.03	13,15,16,16	11

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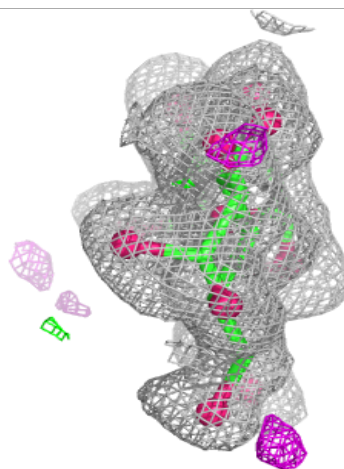
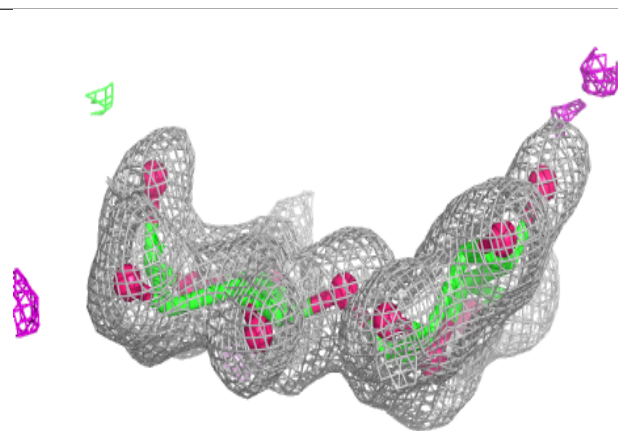
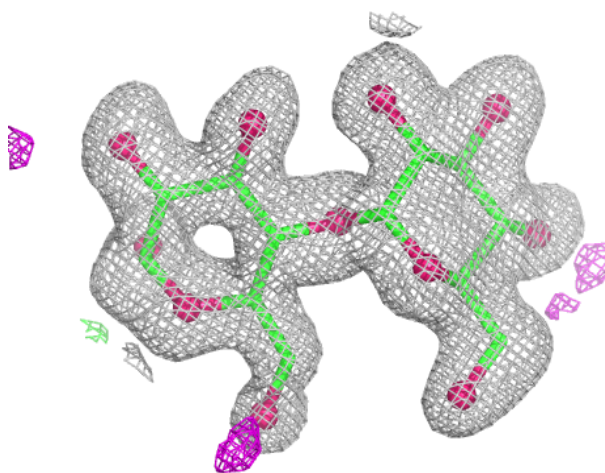
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GLC	B	3[A]	12/12	-	-	14,18,23,25	11
4	GLC	B	3[B]	12/12	-	-	14,21,29,39	11
4	GLC	B	4	11/12	-	-	12,13,15,15	0

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.



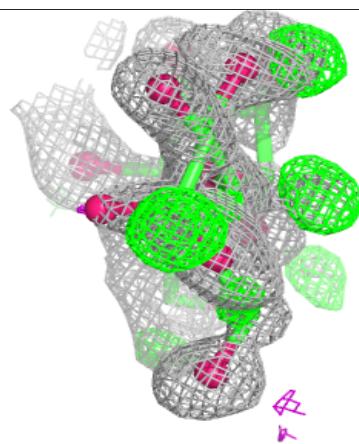
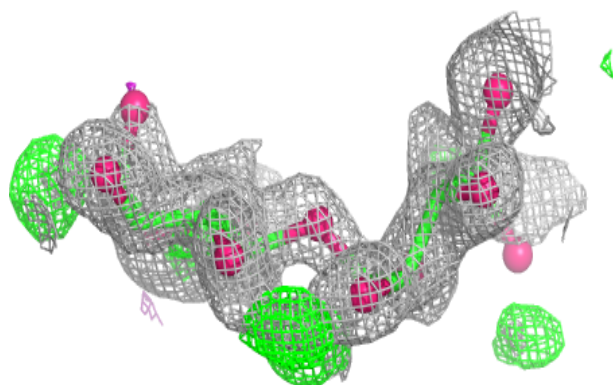
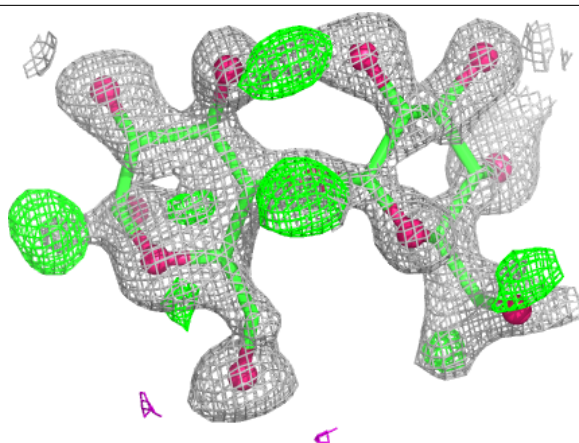
Electron density around Chain F:

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



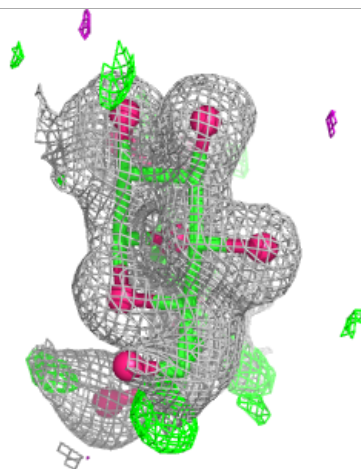
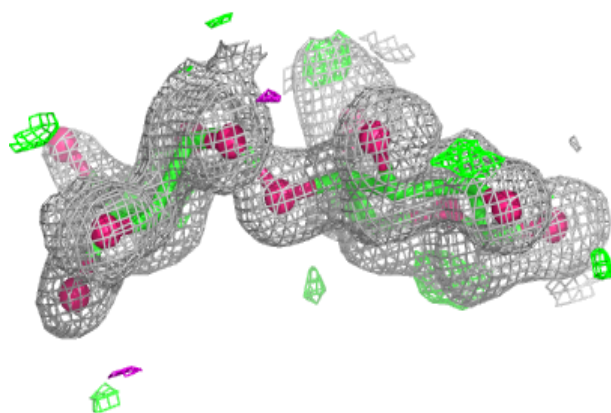
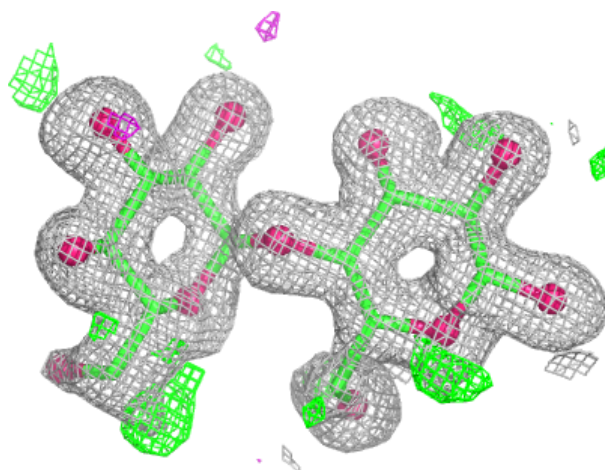
Electron density around Chain D:

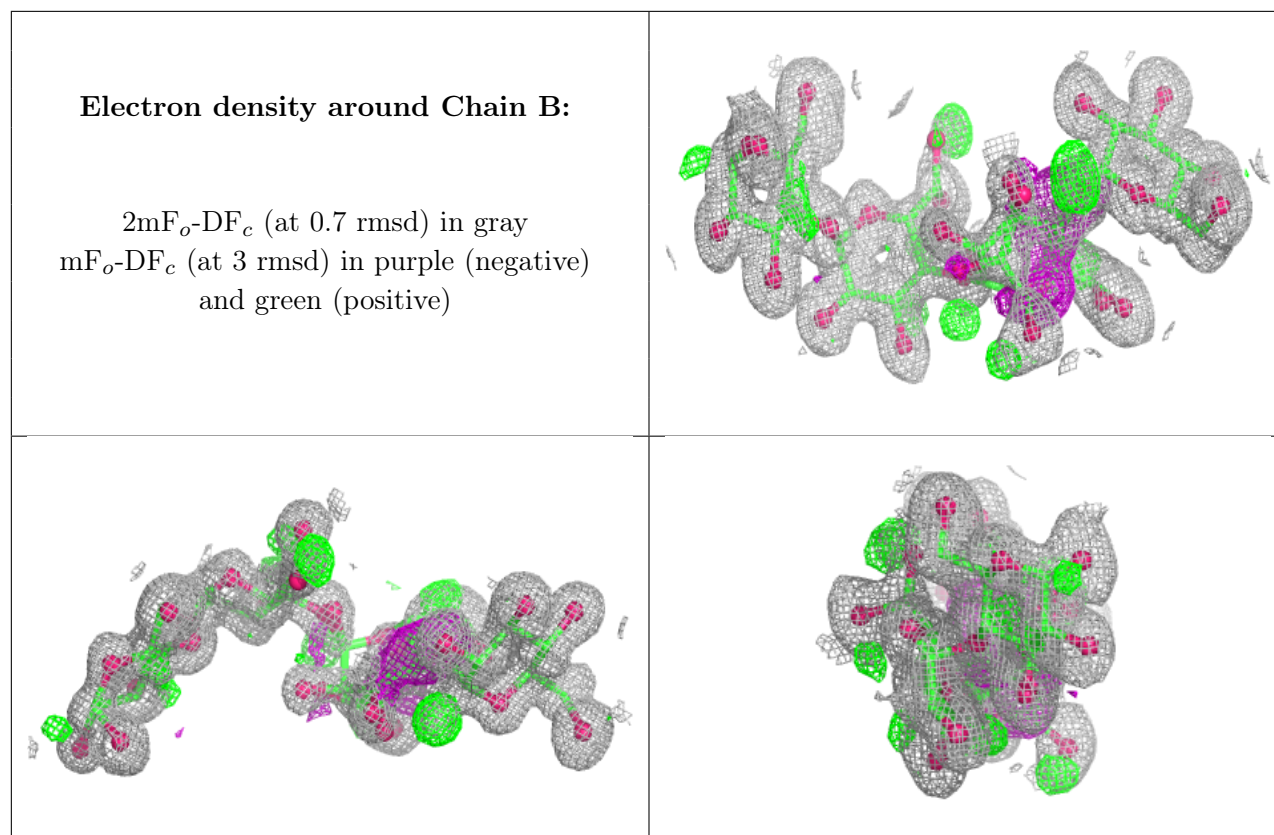
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



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
7	GOL	A	606[A]	6/6	0.75	0.14	33,39,46,57	6
7	GOL	A	606[B]	6/6	0.75	0.14	38,41,42,42	6
7	GOL	A	609	6/6	0.77	0.14	35,47,61,61	0
7	GOL	A	604	6/6	0.78	0.13	39,82,90,94	0
7	GOL	A	608	6/6	0.81	0.14	49,56,63,74	6
7	GOL	A	605	6/6	0.81	0.15	65,73,78,86	0
7	GOL	A	611	6/6	0.85	0.16	102,109,112,124	0
7	GOL	A	610	6/6	0.88	0.11	24,33,34,36	6
7	GOL	A	607	6/6	0.89	0.09	47,48,56,63	6
8	BGC	A	602	12/12	0.90	0.11	19,26,38,40	12
6	SO4	A	603	5/5	0.95	0.08	21,23,26,27	5
5	CA	A	601	1/1	1.00	0.03	18,18,18,18	1

6.5 Other polymers [i](#)

There are no such residues in this entry.