



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 18, 2025 – 05:23 PM EDT

PDB ID : 9BEU / pdb\_00009beu  
Title : Structure of GH110B in complex with a lambda-carrageenan oligosaccharide  
Authors : Hettle, J.A.; Vickers, C.; Boraston, A.B.  
Deposited on : 2024-04-16  
Resolution : 2.40 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0rc1  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.45.1

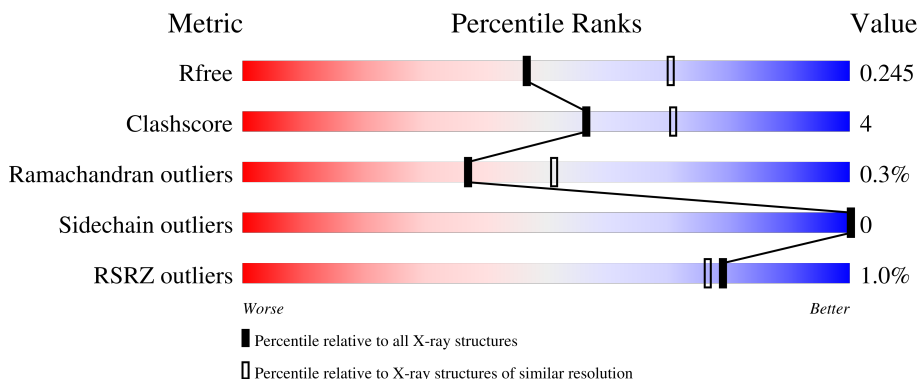
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

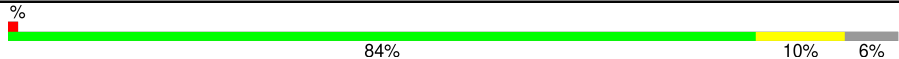

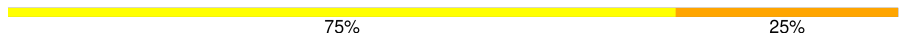
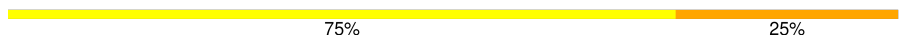
The reported resolution of this entry is 2.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	164625	4642 (2.40-2.40)
Clashscore	180529	5218 (2.40-2.40)
Ramachandran outliers	177936	5158 (2.40-2.40)
Sidechain outliers	177891	5159 (2.40-2.40)
RSRZ outliers	164620	4642 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	620	
1	B	620	
2	C	4	
2	D	4	

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9584 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 Glycoside hydrolase family 110.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	581	Total	C	N	O	S	0	1	0
			4577	2919	778	875	5			
1	B	583	Total	C	N	O	S	0	0	0
			4567	2914	775	873	5			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A8I3AZV9
A	2	GLY	-	expression tag	UNP A0A8I3AZV9
A	3	SER	-	expression tag	UNP A0A8I3AZV9
A	4	SER	-	expression tag	UNP A0A8I3AZV9
A	5	HIS	-	expression tag	UNP A0A8I3AZV9
A	6	HIS	-	expression tag	UNP A0A8I3AZV9
A	7	HIS	-	expression tag	UNP A0A8I3AZV9
A	8	HIS	-	expression tag	UNP A0A8I3AZV9
A	9	HIS	-	expression tag	UNP A0A8I3AZV9
A	10	HIS	-	expression tag	UNP A0A8I3AZV9
A	11	SER	-	expression tag	UNP A0A8I3AZV9
A	12	SER	-	expression tag	UNP A0A8I3AZV9
A	13	GLY	-	expression tag	UNP A0A8I3AZV9
A	14	LEU	-	expression tag	UNP A0A8I3AZV9
A	15	VAL	-	expression tag	UNP A0A8I3AZV9
A	16	PRO	-	expression tag	UNP A0A8I3AZV9
A	17	ARG	-	expression tag	UNP A0A8I3AZV9
A	18	GLY	-	expression tag	UNP A0A8I3AZV9
A	19	SER	-	expression tag	UNP A0A8I3AZV9
A	20	HIS	-	expression tag	UNP A0A8I3AZV9
B	1	MET	-	initiating methionine	UNP A0A8I3AZV9
B	2	GLY	-	expression tag	UNP A0A8I3AZV9
B	3	SER	-	expression tag	UNP A0A8I3AZV9
B	4	SER	-	expression tag	UNP A0A8I3AZV9
B	5	HIS	-	expression tag	UNP A0A8I3AZV9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	6	HIS	-	expression tag	UNP A0A8I3AZV9
B	7	HIS	-	expression tag	UNP A0A8I3AZV9
B	8	HIS	-	expression tag	UNP A0A8I3AZV9
B	9	HIS	-	expression tag	UNP A0A8I3AZV9
B	10	HIS	-	expression tag	UNP A0A8I3AZV9
B	11	SER	-	expression tag	UNP A0A8I3AZV9
B	12	SER	-	expression tag	UNP A0A8I3AZV9
B	13	GLY	-	expression tag	UNP A0A8I3AZV9
B	14	LEU	-	expression tag	UNP A0A8I3AZV9
B	15	VAL	-	expression tag	UNP A0A8I3AZV9
B	16	PRO	-	expression tag	UNP A0A8I3AZV9
B	17	ARG	-	expression tag	UNP A0A8I3AZV9
B	18	GLY	-	expression tag	UNP A0A8I3AZV9
B	19	SER	-	expression tag	UNP A0A8I3AZV9
B	20	HIS	-	expression tag	UNP A0A8I3AZV9

- Molecule 2 is an oligosaccharide called alpha-D-galactopyranose-(1-3)-2-O-sulfo-beta-D-galactopyranose-(1-4)-2,6-di-O-sulfo-alpha-D-galactopyranose-(1-3)-2-O-sulfo-beta-D-galactopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	4	Total	C	O	S	0	0	0
			61	24	33	4			
2	D	4	Total	C	O	S	0	0	0
			61	24	33	4			

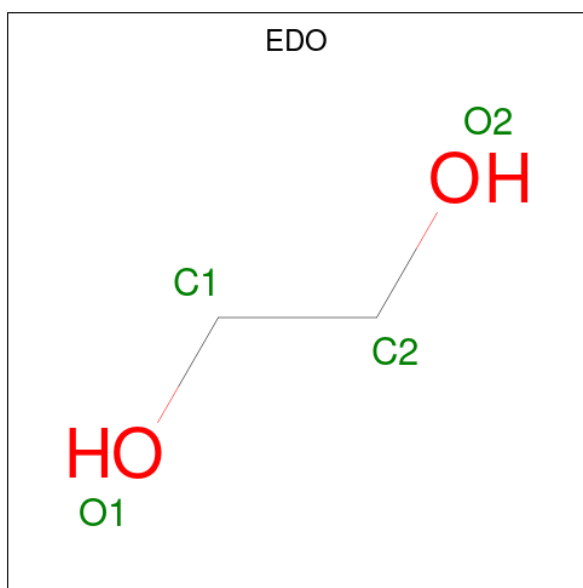
- Molecule 3 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	21	Total	I	0	0
			21	21		
3	B	22	Total	I	0	0
			22	22		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

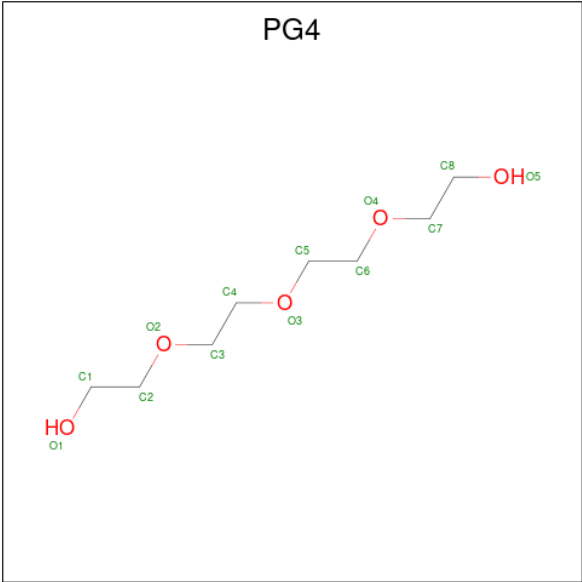
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	Cl	0	0
			12	12		
4	B	12	Total	Cl	0	0
			12	12		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula:  $C_2H_6O_2$ ).



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

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula:  $C_8H_{18}O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			13	8	5		

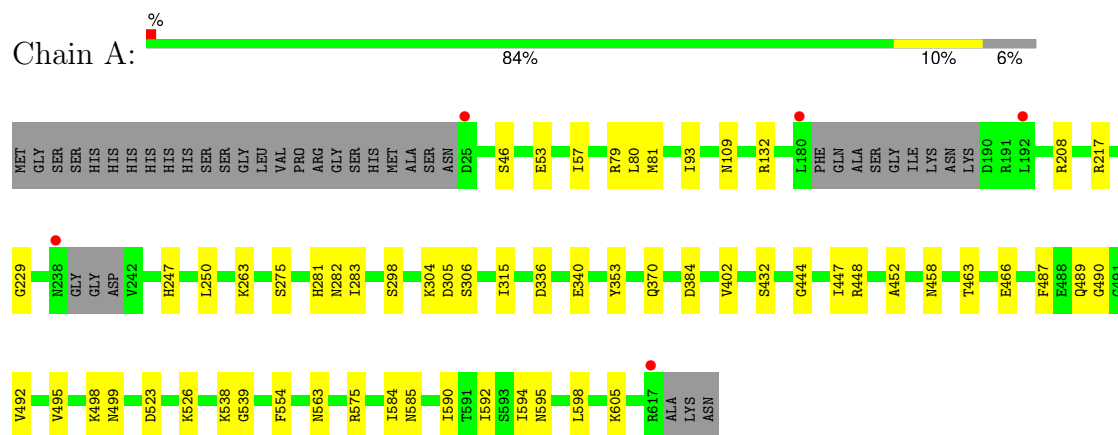
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	110	Total	O	0	0
			110	110		
7	B	100	Total	O	0	0
			100	100		

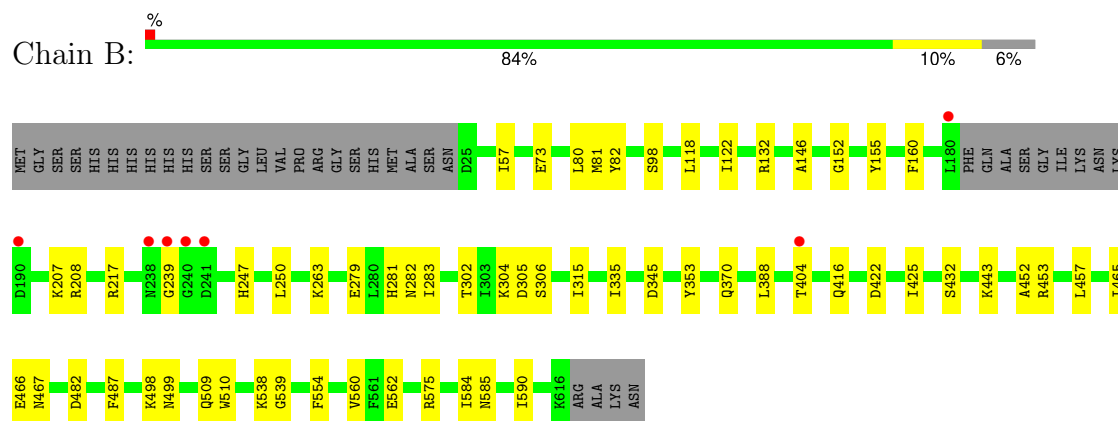
### 3 Residue-property plots

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.

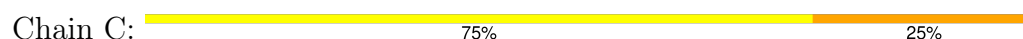
- Molecule 1: Glycoside hydrolase family 110



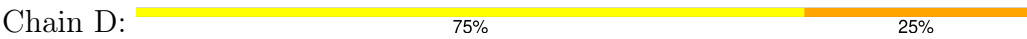
- Molecule 1: Glycoside hydrolase family 110



- Molecule 2: alpha-D-galactopyranose-(1-3)-2-O-sulfo-beta-D-galactopyranose-(1-4)-2,6-di-O-sulfo-alpha-D-galactopyranose-(1-3)-2-O-sulfo-beta-D-galactopyranose



- Molecule 2: alpha-D-galactopyranose-(1-3)-2-O-sulfo-beta-D-galactopyranose-(1-4)-2,6-di-O-sulfo-alpha-D-galactopyranose-(1-3)-2-O-sulfo-beta-D-galactopyranose



A1APB1
A1APD2
A1APB3
GLA4

## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	168.46Å 128.35Å 98.91Å 90.00° 122.15° 90.00°	Depositor
Resolution (Å)	29.96 – 2.40 29.96 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.96-2.40) 98.7 (29.96-2.40)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.23 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, $R_{free}$	0.197 , 0.244 0.198 , 0.245	Depositor DCC
$R_{free}$ test set	3503 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.5	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 31.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9584	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: IOD, A1APD, EDO, CL, PG4, GLA, A1APB

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	0.36	0/4671	0.56	0/6318
1	B	0.36	0/4659	0.56	0/6305
All	All	0.36	0/9330	0.56	0/12623

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	4577	0	4429	38	0
1	B	4567	0	4409	41	0
2	C	61	0	10	1	0
2	D	61	0	10	1	0
3	A	21	0	0	3	0
3	B	22	0	0	5	0
4	A	12	0	0	0	0
4	B	12	0	0	2	0
5	A	16	0	24	5	0
5	B	12	0	18	8	0
6	B	13	0	18	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	110	0	0	2	0
7	B	100	0	0	0	0
All	All	9584	0	8918	82	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 4.

All (82) 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:467:ASN:HD22	5:B:735:EDO:H12	1.40	0.85
1:A:523:ASP:HA	1:A:526:LYS:HD2	1.68	0.75
1:A:247:HIS:HB3	1:A:250:LEU:HG	1.69	0.74
1:B:247:HIS:HB3	1:B:250:LEU:HG	1.71	0.73
1:A:594:ILE:HG12	3:A:706:IOD:I	2.58	0.72
1:B:98:SER:OG	4:B:729:CL:CL	2.46	0.70
1:A:384:ASP:O	1:A:402:VAL:HG12	1.96	0.66
1:B:404:THR:HG23	3:B:702:IOD:I	2.67	0.65
1:B:217:ARG:HA	1:B:315:ILE:HD11	1.78	0.64
1:A:563:ASN:OD1	1:A:595:ASN:ND2	2.32	0.62
1:A:217:ARG:HA	1:A:315:ILE:HD11	1.82	0.61
1:B:584:ILE:HD13	1:B:590:ILE:HG21	1.82	0.61
1:A:217:ARG:NH2	7:A:802:HOH:O	2.32	0.60
5:B:736:EDO:O2	2:D:2:A1APD:O50	2.19	0.60
1:B:575:ARG:HB3	3:B:711:IOD:I	2.73	0.59
1:B:304:LYS:HD2	6:B:738:PG4:H71	1.84	0.58
1:A:283:ILE:O	1:A:306:SER:HB3	2.04	0.58
1:B:404:THR:HG22	1:B:416:GLN:HB3	1.86	0.57
1:B:132:ARG:HA	1:B:281:HIS:O	2.09	0.53
1:B:482:ASP:OD1	1:B:482:ASP:C	2.52	0.52
1:A:282:ASN:HA	1:A:305:ASP:O	2.10	0.52
1:B:283:ILE:O	1:B:306:SER:HB3	2.09	0.52
1:A:463:THR:HB	1:A:495:VAL:HG22	1.91	0.51
1:A:263:LYS:NZ	5:A:735:EDO:H12	2.26	0.51
1:A:80:LEU:HD12	1:A:81:MET:N	2.26	0.50
1:A:229:GLY:HA2	5:A:735:EDO:H21	1.92	0.50
1:A:575:ARG:HB3	3:A:718:IOD:I	2.81	0.50
1:B:560:VAL:HG12	1:B:562:GLU:HG2	1.93	0.50
1:B:263:LYS:NZ	5:B:737:EDO:H22	2.27	0.49
1:B:509:GLN:HB2	5:B:736:EDO:H22	1.94	0.49
1:A:584:ILE:HD13	1:A:590:ILE:HG21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:SER:OG	5:A:736:EDO:H22	2.13	0.48
1:B:304:LYS:HG3	1:B:335:ILE:HB	1.94	0.48
1:B:370:GLN:HG3	1:B:487:PHE:CZ	2.49	0.48
1:A:304:LYS:HD2	5:A:734:EDO:H12	1.95	0.47
1:B:388:LEU:HD23	1:B:425:ILE:HD13	1.96	0.47
1:B:282:ASN:HA	1:B:305:ASP:O	2.14	0.47
1:B:443:LYS:HB3	5:B:735:EDO:H22	1.96	0.47
1:A:336:ASP:HA	1:A:444:GLY:O	2.15	0.47
1:B:510:TRP:CD1	5:B:736:EDO:H12	2.49	0.47
1:A:109:ASN:OD1	5:A:737:EDO:H11	2.15	0.47
1:B:353:TYR:CD1	1:B:432:SER:HB2	2.51	0.46
1:A:208:ARG:HD3	2:C:2:A1APD:O55	2.15	0.46
3:B:710:IOD:I	4:B:733:CL:CL	3.40	0.46
1:B:467:ASN:ND2	5:B:735:EDO:H12	2.20	0.45
1:A:598:LEU:HD13	1:A:605:LYS:HB2	1.98	0.45
1:B:57:ILE:HD13	3:B:715:IOD:I	2.86	0.45
1:B:457:LEU:HD11	1:B:465:ILE:HG13	1.99	0.45
1:B:554:PHE:HA	1:B:585:ASN:O	2.16	0.45
1:A:499:ASN:HA	1:A:539:GLY:O	2.17	0.45
1:B:80:LEU:HD21	1:B:82:TYR:CE2	2.50	0.45
1:B:263:LYS:HZ2	5:B:737:EDO:H22	1.82	0.45
1:B:538:LYS:HG3	3:B:720:IOD:I	2.87	0.45
1:A:584:ILE:HD11	1:A:592:ILE:HD11	1.99	0.44
1:A:498:LYS:HA	1:A:538:LYS:O	2.17	0.44
1:B:443:LYS:HA	1:B:466:GLU:O	2.18	0.44
1:A:79:ARG:HG3	1:A:93:ILE:HD13	1.99	0.43
1:A:466:GLU:HA	1:A:498:LYS:O	2.17	0.43
1:A:353:TYR:CE2	1:A:489:GLN:HA	2.53	0.43
1:A:370:GLN:HG3	1:A:487:PHE:CZ	2.53	0.43
1:B:279:GLU:HG2	1:B:302:THR:HB	2.00	0.43
1:A:538:LYS:HG3	3:A:721:IOD:I	2.89	0.42
1:A:53:GLU:O	1:A:57:ILE:HG12	2.20	0.42
1:A:554:PHE:HA	1:A:585:ASN:O	2.19	0.42
1:B:146:ALA:HB2	1:B:160:PHE:CE1	2.55	0.42
1:A:132:ARG:HA	1:A:281:HIS:O	2.19	0.42
1:B:73:GLU:HA	1:B:118:LEU:HD21	2.02	0.42
1:B:422:ASP:O	1:B:425:ILE:HD12	2.19	0.42
1:A:458:ASN:HB3	1:A:490:GLY:HA2	2.02	0.41
1:A:492:VAL:HG13	1:A:495:VAL:HG23	2.01	0.41
1:B:499:ASN:HA	1:B:539:GLY:O	2.21	0.41
6:B:738:PG4:H12	6:B:738:PG4:H62	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:MET:HE2	1:B:122:ILE:HG21	2.02	0.41
1:A:340:GLU:HA	1:A:448:ARG:O	2.21	0.41
1:B:152:GLY:HA3	1:B:155:TYR:CE2	2.56	0.41
1:B:466:GLU:HA	1:B:498:LYS:O	2.21	0.41
1:B:345:ASP:HA	1:B:453:ARG:O	2.21	0.41
1:A:353:TYR:CD1	1:A:432:SER:HB2	2.55	0.40
1:B:207:LYS:HG3	1:B:208:ARG:HG2	2.02	0.40
1:A:447:ILE:HG23	7:A:823:HOH:O	2.20	0.40
1:B:304:LYS:HA	1:B:335:ILE:O	2.22	0.40
1:A:275:SER:O	1:A:298:SER:HB3	2.21	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	576/620 (93%)	548 (95%)	27 (5%)	1 (0%)	44	59
1	B	579/620 (93%)	549 (95%)	28 (5%)	2 (0%)	37	51
All	All	1155/1240 (93%)	1097 (95%)	55 (5%)	3 (0%)	37	51

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	452	ALA
1	B	452	ALA
1	B	239	GLY

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	485/541 (90%)	485 (100%)	0	100	100
1	B	481/541 (89%)	481 (100%)	0	100	100
All	All	966/1082 (89%)	966 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	ASN
1	A	373	GLN
1	A	563	ASN
1	A	595	ASN
1	B	370	GLN
1	B	467	ASN
1	B	563	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](#)

8 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	A1APB	C	1	2	16,16,16	1.70	4 (25%)	21,24,24	1.84	5 (23%)
2	A1APD	C	2	2	19,19,20	2.03	4 (21%)	23,29,31	2.01	5 (21%)
2	A1APB	C	3	2	15,15,16	1.82	5 (33%)	17,22,24	1.30	2 (11%)
2	GLA	C	4	2	11,11,12	1.77	2 (18%)	15,15,17	1.10	1 (6%)
2	A1APB	D	1	2	16,16,16	1.57	3 (18%)	21,24,24	1.80	5 (23%)
2	A1APD	D	2	2	19,19,20	1.97	4 (21%)	23,29,31	2.08	5 (21%)
2	A1APB	D	3	2	15,15,16	1.77	5 (33%)	17,22,24	1.07	2 (11%)
2	GLA	D	4	2	11,11,12	1.89	3 (27%)	15,15,17	1.02	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	A1APB	C	1	2	-	0/7/27/27	0/1/1/1
2	A1APD	C	2	2	-	4/11/28/31	0/1/1/1
2	A1APB	C	3	2	-	3/7/24/27	0/1/1/1
2	GLA	C	4	2	-	0/2/19/22	0/1/1/1
2	A1APB	D	1	2	-	2/7/27/27	0/1/1/1
2	A1APD	D	2	2	-	7/11/28/31	0/1/1/1
2	A1APB	D	3	2	-	3/7/24/27	0/1/1/1
2	GLA	D	4	2	-	0/2/19/22	0/1/1/1

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	A1APD	O2-S2	4.79	1.71	1.57
2	D	4	GLA	O5-C1	4.72	1.51	1.43
2	C	4	GLA	O5-C1	4.43	1.51	1.43
2	D	2	A1APD	O2-S2	4.39	1.70	1.57
2	C	2	A1APD	O6-S6	3.90	1.67	1.56
2	C	1	A1APB	O2-S1	3.77	1.68	1.57
2	D	2	A1APD	O4-C4	3.70	1.52	1.43
2	D	2	A1APD	O6-S6	3.68	1.66	1.56
2	C	2	A1APD	O4-C4	3.44	1.51	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	A1APB	O2-S1	3.40	1.67	1.57
2	C	3	A1APB	O2-S1	3.36	1.67	1.57
2	C	3	A1APB	O5-C1	3.02	1.48	1.43
2	C	1	A1APB	O5-C5	2.99	1.51	1.44
2	C	2	A1APD	O5-C5	2.98	1.49	1.43
2	C	1	A1APB	O3-C3	2.92	1.50	1.43
2	D	3	A1APB	O2-S1	2.89	1.66	1.57
2	D	2	A1APD	O5-C5	2.82	1.48	1.43
2	C	4	GLA	C2-C3	-2.81	1.48	1.52
2	D	1	A1APB	O5-C5	2.73	1.51	1.44
2	D	4	GLA	C2-C3	-2.70	1.48	1.52
2	D	3	A1APB	O5-C1	2.68	1.48	1.43
2	D	3	A1APB	O5-C5	2.54	1.48	1.43
2	D	3	A1APB	O3-C3	2.40	1.48	1.43
2	D	1	A1APB	O3-C3	2.36	1.48	1.43
2	D	3	A1APB	C1-C2	2.30	1.55	1.51
2	C	3	A1APB	C3-C2	-2.27	1.48	1.53
2	D	4	GLA	O5-C5	2.26	1.47	1.43
2	C	1	A1APB	O5-C1	2.24	1.48	1.42
2	C	3	A1APB	O3-C3	2.12	1.48	1.43
2	C	3	A1APB	C1-C2	2.06	1.54	1.51

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	A1APD	O2-C2-C3	7.11	116.89	106.95
2	C	2	A1APD	O2-C2-C3	6.12	115.49	106.95
2	D	1	A1APB	C4-C3-C2	4.40	119.67	109.68
2	C	1	A1APB	C1-O5-C5	-3.70	106.49	113.65
2	C	1	A1APB	C4-C3-C2	3.61	117.86	109.68
2	C	1	A1APB	C3-C4-C5	3.48	116.53	110.23
2	C	2	A1APD	O55-S2-O54	-3.46	98.89	112.24
2	C	1	A1APB	O8-S1-O7	-3.43	99.00	112.24
2	D	1	A1APB	O8-S1-O7	-3.31	99.46	112.24
2	D	1	A1APB	C1-O5-C5	-3.23	107.40	113.65
2	C	4	GLA	C1-O5-C5	3.21	116.48	112.19
2	D	2	A1APD	O55-S2-O54	-3.03	100.54	112.24
2	D	2	A1APD	C6-C5-C4	2.93	118.21	112.07
2	C	2	A1APD	O53-S2-O2	2.89	113.02	106.37
2	C	2	A1APD	C6-C5-C4	2.89	118.13	112.07
2	C	3	A1APB	O8-S1-O7	-2.71	101.77	112.24
2	C	3	A1APB	O9-S1-O2	2.71	112.61	106.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	A1APD	O51-S6-O50	-2.67	101.94	112.24
2	D	1	A1APB	C3-C4-C5	2.63	115.01	110.23
2	D	4	GLA	C1-O5-C5	2.53	115.58	112.19
2	D	2	A1APD	O6-C6-C5	2.46	111.95	107.57
2	C	2	A1APD	O51-S6-O50	-2.36	103.15	112.24
2	D	1	A1APB	O4-C4-C3	2.28	115.75	110.38
2	D	3	A1APB	O9-S1-O2	2.24	111.53	106.37
2	D	3	A1APB	O8-S1-O7	-2.15	103.95	112.24
2	C	1	A1APB	O2-C2-C1	-2.01	104.88	107.58

There are no chirality outliers.

All (19) torsion outliers are listed below:

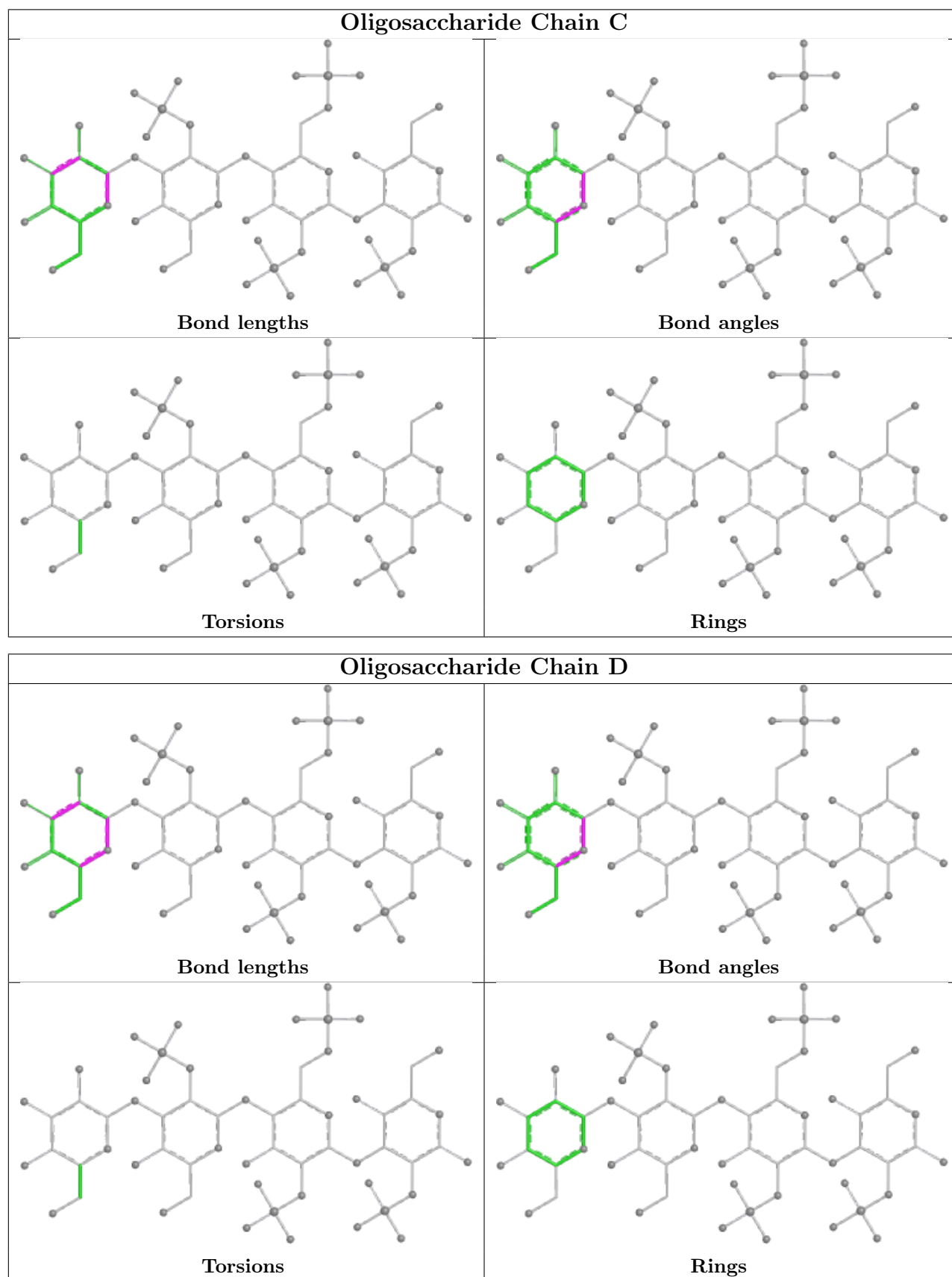
Mol	Chain	Res	Type	Atoms
2	C	2	A1APD	C6-O6-S6-O52
2	C	3	A1APB	C2-O2-S1-O8
2	D	2	A1APD	O5-C5-C6-O6
2	D	2	A1APD	C1-C2-O2-S2
2	D	2	A1APD	C3-C2-O2-S2
2	D	1	A1APB	O5-C5-C6-O6
2	D	1	A1APB	C4-C5-C6-O6
2	C	3	A1APB	C2-O2-S1-O7
2	D	2	A1APD	C6-O6-S6-O52
2	C	2	A1APD	C6-O6-S6-O51
2	D	2	A1APD	C6-O6-S6-O51
2	C	2	A1APD	C6-O6-S6-O50
2	C	3	A1APB	C2-O2-S1-O9
2	D	3	A1APB	C2-O2-S1-O8
2	D	3	A1APB	C2-O2-S1-O9
2	D	2	A1APD	C6-O6-S6-O50
2	C	2	A1APD	O5-C5-C6-O6
2	D	2	A1APD	C4-C5-C6-O6
2	D	3	A1APB	C2-O2-S1-O7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	A1APD	1	0
2	D	2	A1APD	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

Of 75 ligands modelled in this entry, 67 are monoatomic - leaving 8 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	$\# Z  > 2$
5	EDO	A	735	-	3,3,3	0.37	0	2,2,2	0.25	0
5	EDO	A	734	-	3,3,3	0.46	0	2,2,2	0.41	0
5	EDO	A	737	-	3,3,3	0.57	0	2,2,2	0.31	0
5	EDO	B	737	-	3,3,3	0.44	0	2,2,2	0.50	0
5	EDO	B	736	-	3,3,3	0.56	0	2,2,2	0.06	0
5	EDO	B	735	-	3,3,3	0.59	0	2,2,2	0.14	0
5	EDO	A	736	-	3,3,3	0.48	0	2,2,2	0.40	0
6	PG4	B	738	-	12,12,12	0.63	0	11,11,11	0.70	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
5	EDO	A	735	-	-	0/1/1/1	-
5	EDO	A	734	-	-	1/1/1/1	-
5	EDO	A	737	-	-	1/1/1/1	-
5	EDO	B	737	-	-	1/1/1/1	-
5	EDO	B	736	-	-	1/1/1/1	-
5	EDO	B	735	-	-	1/1/1/1	-
5	EDO	A	736	-	-	1/1/1/1	-
6	PG4	B	738	-	-	10/10/10/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	738	PG4	O3-C5-C6-O4
6	B	738	PG4	O2-C3-C4-O3
6	B	738	PG4	C6-C5-O3-C4
5	A	736	EDO	O1-C1-C2-O2
5	A	737	EDO	O1-C1-C2-O2
6	B	738	PG4	O4-C7-C8-O5
5	B	737	EDO	O1-C1-C2-O2
5	B	735	EDO	O1-C1-C2-O2
6	B	738	PG4	C1-C2-O2-C3
6	B	738	PG4	C8-C7-O4-C6
6	B	738	PG4	C4-C3-O2-C2
6	B	738	PG4	O1-C1-C2-O2
6	B	738	PG4	C3-C4-O3-C5
6	B	738	PG4	C5-C6-O4-C7
5	A	734	EDO	O1-C1-C2-O2
5	B	736	EDO	O1-C1-C2-O2

There are no ring outliers.

8 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	735	EDO	2	0
5	A	734	EDO	1	0
5	A	737	EDO	1	0
5	B	737	EDO	2	0
5	B	736	EDO	3	0
5	B	735	EDO	3	0
5	A	736	EDO	1	0
6	B	738	PG4	2	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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	581/620 (93%)	-0.26	5 (0%) 81 78	23, 40, 51, 68	2 (0%)
1	B	583/620 (94%)	-0.31	7 (1%) 76 73	30, 39, 51, 80	0
All	All	1164/1240 (93%)	-0.28	12 (1%) 79 76	23, 40, 51, 80	2 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	180	LEU	4.0
1	A	238	ASN	3.9
1	B	238	ASN	3.8
1	B	239	GLY	3.8
1	A	192	LEU	3.6
1	A	180	LEU	3.4
1	A	617	ARG	3.4
1	A	25	ASP	2.6
1	B	241	ASP	2.4
1	B	240	GLY	2.2
1	B	190	ASP	2.2
1	B	404	THR	2.0

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

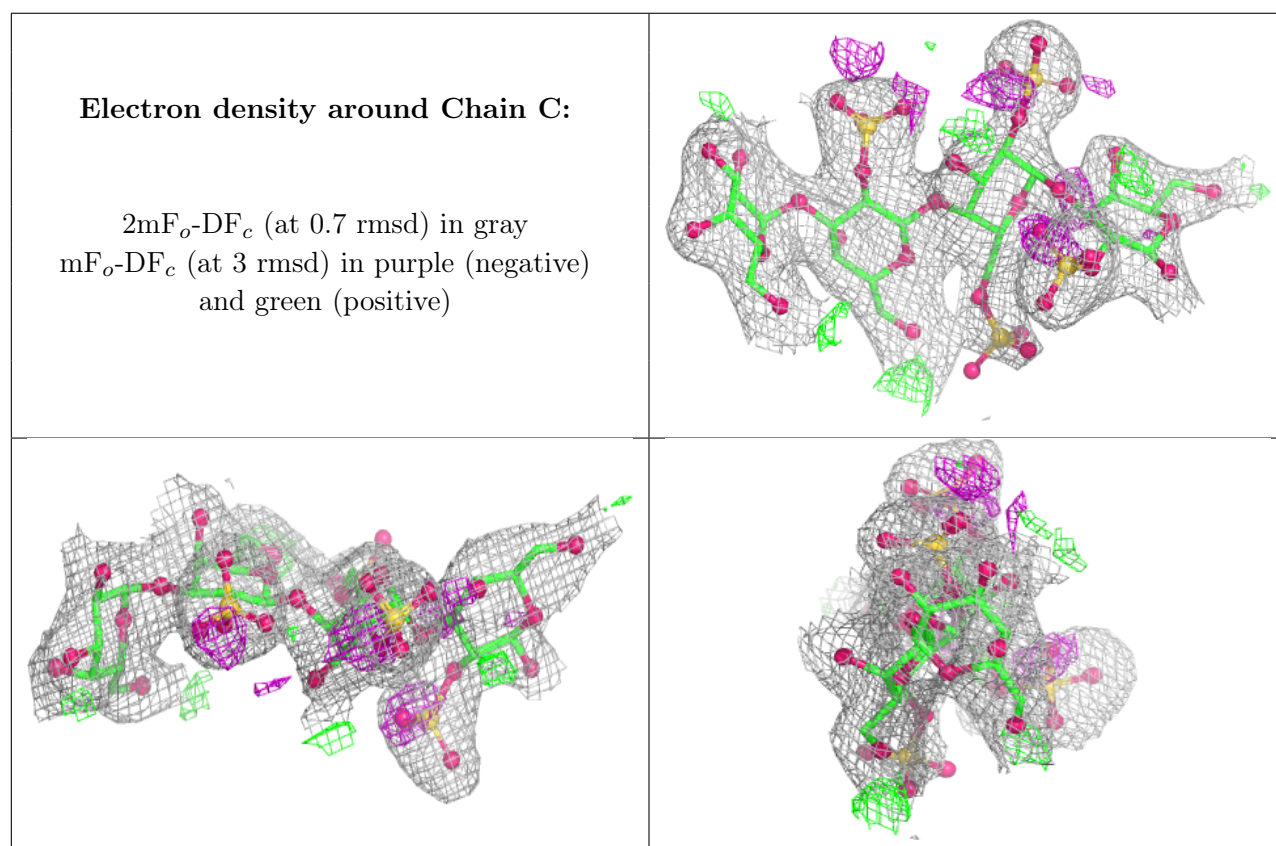
There are no non-standard protein/DNA/RNA residues in this entry.

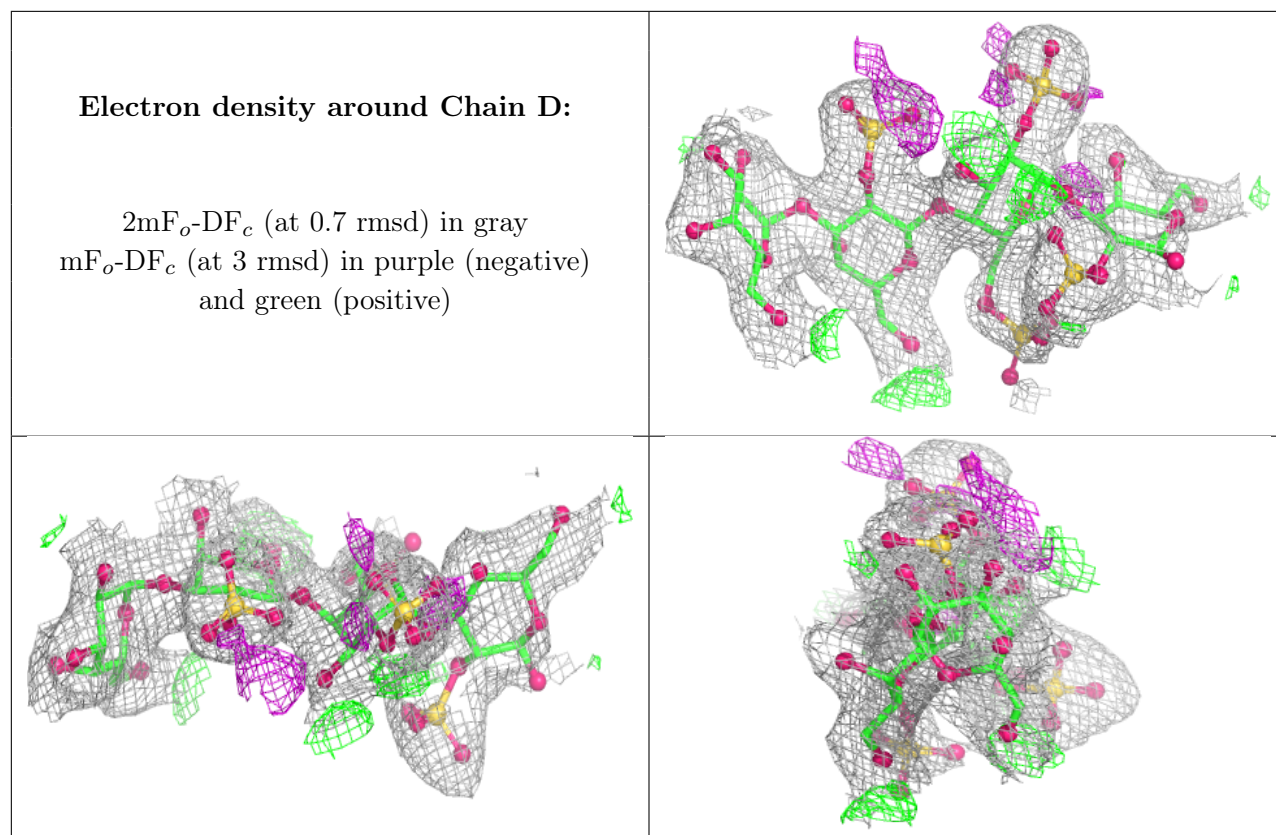
### 6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	A1APB	D	1	16/16	0.79	0.15	54,64,75,76	0
2	A1APB	C	1	16/16	0.82	0.16	55,65,73,76	0
2	A1APD	C	2	19/20	0.90	0.14	48,54,59,60	3
2	A1APD	D	2	19/20	0.90	0.14	44,54,58,61	4
2	A1APB	D	3	15/16	0.95	0.09	34,42,45,49	0
2	A1APB	C	3	15/16	0.96	0.09	38,41,45,47	0
2	GLA	D	4	11/12	0.96	0.08	35,38,41,42	0
2	GLA	C	4	11/12	0.97	0.09	35,38,40,41	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.





## 6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	IOD	A	721	1/1	0.80	0.13	77,77,77,77	1
6	PG4	B	738	13/13	0.80	0.19	35,43,53,56	0
5	EDO	A	736	4/4	0.81	0.14	46,49,51,55	0
5	EDO	B	735	4/4	0.83	0.17	36,41,42,47	0
5	EDO	B	736	4/4	0.87	0.19	46,49,54,54	0
4	CL	B	734	1/1	0.87	0.14	59,59,59,59	0
5	EDO	A	737	4/4	0.89	0.17	38,41,44,47	0
4	CL	A	733	1/1	0.89	0.12	62,62,62,62	0
3	IOD	B	722	1/1	0.89	0.12	75,75,75,75	1
4	CL	A	727	1/1	0.89	0.10	66,66,66,66	0
5	EDO	A	735	4/4	0.91	0.14	31,37,41,45	0
3	IOD	B	720	1/1	0.91	0.11	80,80,80,80	0
4	CL	B	730	1/1	0.92	0.12	62,62,62,62	0
3	IOD	A	704	1/1	0.93	0.09	80,80,80,80	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	IOD	B	710	1/1	0.93	0.09	75,75,75,75	1
5	EDO	B	737	4/4	0.93	0.13	31,38,40,44	0
5	EDO	A	734	4/4	0.93	0.15	27,34,35,36	0
4	CL	A	731	1/1	0.94	0.13	57,57,57,57	0
3	IOD	B	715	1/1	0.94	0.16	114,114,114,114	0
3	IOD	A	709	1/1	0.94	0.20	121,121,121,121	0
3	IOD	A	708	1/1	0.95	0.06	55,55,55,55	1
3	IOD	B	716	1/1	0.95	0.07	67,67,67,67	1
3	IOD	B	717	1/1	0.95	0.19	120,120,120,120	0
4	CL	A	730	1/1	0.96	0.12	51,51,51,51	0
4	CL	B	724	1/1	0.96	0.15	52,52,52,52	0
3	IOD	B	721	1/1	0.96	0.06	60,60,60,60	1
3	IOD	B	709	1/1	0.97	0.05	62,62,62,62	1
3	IOD	A	703	1/1	0.97	0.09	78,78,78,78	0
3	IOD	A	705	1/1	0.97	0.04	69,69,69,69	1
4	CL	B	726	1/1	0.97	0.08	41,41,41,41	0
3	IOD	A	715	1/1	0.97	0.05	58,58,58,58	1
4	CL	B	733	1/1	0.97	0.11	39,39,39,39	0
3	IOD	A	719	1/1	0.97	0.05	66,66,66,66	1
3	IOD	A	706	1/1	0.97	0.17	103,103,103,103	0
3	IOD	B	704	1/1	0.97	0.15	102,102,102,102	0
3	IOD	B	705	1/1	0.97	0.14	105,105,105,105	0
4	CL	A	723	1/1	0.97	0.09	55,55,55,55	0
4	CL	A	726	1/1	0.97	0.12	46,46,46,46	0
3	IOD	B	707	1/1	0.97	0.05	55,55,55,55	1
4	CL	A	728	1/1	0.97	0.07	48,48,48,48	0
3	IOD	B	708	1/1	0.97	0.06	51,51,51,51	1
3	IOD	B	719	1/1	0.98	0.04	59,59,59,59	1
3	IOD	A	716	1/1	0.98	0.05	59,59,59,59	1
4	CL	B	728	1/1	0.98	0.10	32,32,32,32	0
3	IOD	A	717	1/1	0.98	0.04	58,58,58,58	1
3	IOD	A	710	1/1	0.98	0.04	57,57,57,57	1
4	CL	A	722	1/1	0.98	0.06	29,29,29,29	0
3	IOD	A	720	1/1	0.98	0.05	69,69,69,69	1
3	IOD	B	711	1/1	0.98	0.04	61,61,61,61	1
3	IOD	B	714	1/1	0.98	0.06	58,58,58,58	1
3	IOD	A	711	1/1	0.98	0.15	88,88,88,88	0
4	CL	A	729	1/1	0.98	0.07	38,38,38,38	0
3	IOD	A	712	1/1	0.98	0.04	66,66,66,66	1
3	IOD	A	707	1/1	0.98	0.15	96,96,96,96	0
3	IOD	B	718	1/1	0.98	0.04	65,65,65,65	1
4	CL	A	724	1/1	0.99	0.09	30,30,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	CL	B	727	1/1	0.99	0.06	37,37,37,37	0
4	CL	A	725	1/1	0.99	0.12	38,38,38,38	0
4	CL	B	729	1/1	0.99	0.05	27,27,27,27	0
3	IOD	A	714	1/1	0.99	0.02	53,53,53,53	1
4	CL	B	731	1/1	0.99	0.14	40,40,40,40	0
4	CL	B	732	1/1	0.99	0.12	39,39,39,39	0
3	IOD	B	702	1/1	0.99	0.03	45,45,45,45	1
3	IOD	B	703	1/1	0.99	0.03	50,50,50,50	0
3	IOD	A	718	1/1	0.99	0.03	58,58,58,58	1
3	IOD	B	713	1/1	0.99	0.04	63,63,63,63	1
3	IOD	A	702	1/1	0.99	0.03	47,47,47,47	1
4	CL	A	732	1/1	0.99	0.06	39,39,39,39	0
3	IOD	B	706	1/1	0.99	0.03	61,61,61,61	1
4	CL	B	723	1/1	0.99	0.06	28,28,28,28	0
3	IOD	A	713	1/1	0.99	0.03	55,55,55,55	1
4	CL	B	725	1/1	0.99	0.09	44,44,44,44	0
3	IOD	A	701	1/1	1.00	0.02	42,42,42,42	0
3	IOD	B	701	1/1	1.00	0.01	40,40,40,40	0
3	IOD	B	712	1/1	1.00	0.08	69,69,69,69	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.