



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:14 AM UTC

PDB ID : 7C69 / pdb_00007c69
Title : Crystal structure of beta-glycosides-binding protein of ABC transporter in a closed state bound to sophorose
Authors : Kanaujia, S.P.; Chandravanshi, M.; Samanta, R.
Deposited on : 2020-05-21
Resolution : 2.35 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

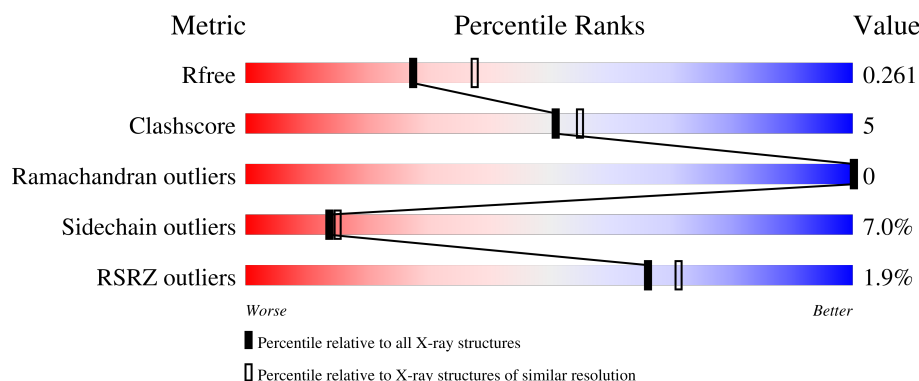
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	423	 83% 14% ..
1	B	423	 83% 12% ...
2	C	2	 50% 50%
2	D	2	 100%

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 6709 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 Sugar ABC transporter, periplasmic sugar-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	0	0
			3183	2046	550	577	10			
1	B	418	Total	C	N	O	S	0	0	0
			3204	2059	557	578	10			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q53W80
A	417	HIS	-	expression tag	UNP Q53W80
A	418	HIS	-	expression tag	UNP Q53W80
A	419	HIS	-	expression tag	UNP Q53W80
A	420	HIS	-	expression tag	UNP Q53W80
A	421	HIS	-	expression tag	UNP Q53W80
A	422	HIS	-	expression tag	UNP Q53W80
B	0	MET	-	initiating methionine	UNP Q53W80
B	417	HIS	-	expression tag	UNP Q53W80
B	418	HIS	-	expression tag	UNP Q53W80
B	419	HIS	-	expression tag	UNP Q53W80
B	420	HIS	-	expression tag	UNP Q53W80
B	421	HIS	-	expression tag	UNP Q53W80
B	422	HIS	-	expression tag	UNP Q53W80

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



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

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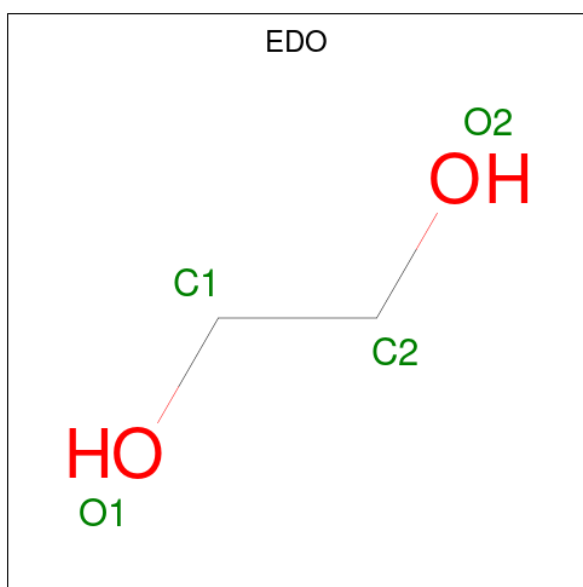
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Cl	0	0
			2	2		
3	B	2	Total	Cl	0	0
			2	2		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

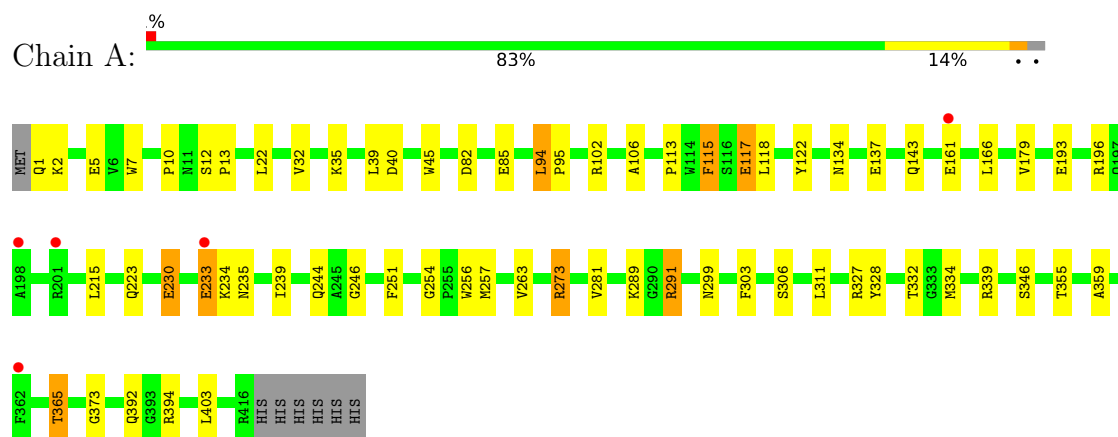
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	142	Total	O	0	0
			142	142		
5	B	122	Total	O	0	0
			122	122		

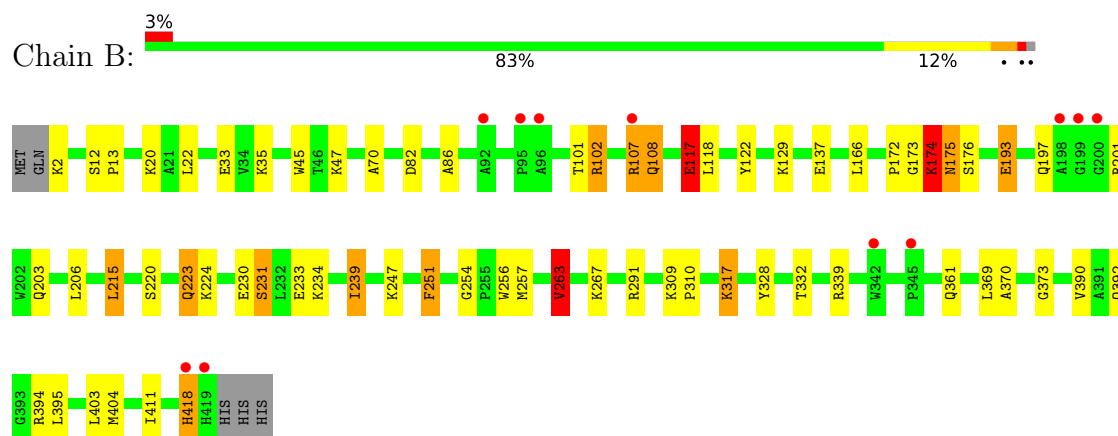
3 Residue-property plots

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: Sugar ABC transporter, periplasmic sugar-binding protein



- Molecule 1: Sugar ABC transporter, periplasmic sugar-binding protein



- Molecule 2: beta-D-glucopyranose-(1-2)-beta-D-glucopyranose



- Molecule 2: beta-D-glucopyranose-(1-2)-beta-D-glucopyranose



BGC1
BGC2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.79Å 108.93Å 109.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	55.06 – 2.35 55.06 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (55.06-2.35) 99.9 (55.06-2.35)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.65 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.200 , 0.262 0.204 , 0.261	Depositor DCC
R_{free} test set	1707 reflections (5.31%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.299	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.042 for -h,l,k	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6709	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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: BGC, CL, EDO

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.18	4/3271 (0.1%)	1.46	13/4453 (0.3%)
1	B	1.14	0/3295	1.47	10/4486 (0.2%)
All	All	1.16	4/6566 (0.1%)	1.46	23/8939 (0.3%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	10	PRO	C-O	-6.85	1.15	1.24
1	A	115	PHE	C-O	5.84	1.30	1.23
1	A	306	SER	C-O	5.52	1.30	1.23
1	A	303	PHE	C-O	5.21	1.30	1.23

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	291	ARG	CG-CD-NE	-6.73	97.20	112.00
1	B	137	GLU	CB-CG-CD	6.21	123.15	112.60
1	B	174	LYS	CA-C-N	5.77	133.48	123.91
1	B	174	LYS	C-N-CA	5.77	133.48	123.91
1	B	193	GLU	CB-CA-C	-5.75	99.90	111.17
1	B	263	VAL	CB-CA-C	5.66	116.43	110.37
1	A	106	ALA	CA-C-N	5.56	128.04	120.54
1	A	106	ALA	C-N-CA	5.56	128.04	120.54
1	A	95	PRO	CA-C-N	5.47	127.61	120.28
1	A	95	PRO	C-N-CA	5.47	127.61	120.28
1	A	233	GLU	CB-CG-CD	5.44	121.85	112.60
1	A	161	GLU	CB-CG-CD	5.44	121.84	112.60
1	A	196	ARG	CB-CG-CD	-5.42	98.83	111.30
1	B	369	LEU	CA-C-N	5.42	127.86	120.54
1	B	369	LEU	C-N-CA	5.42	127.86	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	193	GLU	CB-CA-C	-5.42	100.55	111.17
1	A	85	GLU	CB-CG-CD	5.32	121.64	112.60
1	B	82	ASP	CA-CB-CG	5.25	117.85	112.60
1	B	418	HIS	CA-CB-CG	5.24	119.04	113.80
1	B	117	GLU	CB-CA-C	5.23	118.73	109.89
1	A	179	VAL	N-CA-C	-5.20	104.70	110.62
1	A	82	ASP	CA-CB-CG	5.10	117.70	112.60
1	A	40	ASP	CB-CA-C	5.03	118.06	109.72

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	3183	0	3150	30	0
1	B	3204	0	3160	35	0
2	C	23	0	21	0	0
2	D	23	0	21	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	8	0	12	2	0
5	A	142	0	0	5	0
5	B	122	0	0	5	0
All	All	6709	0	6364	66	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 (66) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:THR:HG21	5:A:607:HOH:O	1.69	0.92
1:B:107:ARG:HG2	1:B:108:GLN:HG2	1.73	0.69
1:A:327:ARG:HG2	5:A:675:HOH:O	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:LEU:HD13	1:A:339:ARG:HD3	1.79	0.65
1:A:2:LYS:NZ	4:A:504:EDO:O2	2.32	0.62
1:A:5:GLU:OE1	1:A:35:LYS:HE2	2.00	0.61
1:A:32:VAL:HG21	1:A:311:LEU:HD11	1.82	0.61
1:A:246:GLY:HA2	5:A:629:HOH:O	1.99	0.60
1:A:254:GLY:HA3	1:A:256:TRP:CZ3	2.36	0.59
1:A:115:PHE:CE1	1:A:334:MET:HE3	2.38	0.58
1:B:86:ALA:CB	1:B:317:LYS:HE2	2.34	0.58
1:B:254:GLY:HA3	1:B:256:TRP:CZ3	2.39	0.57
1:A:230:GLU:HG2	1:A:234:LYS:HE3	1.89	0.54
1:B:70:ALA:HA	1:B:370:ALA:O	2.08	0.53
1:A:233:GLU:HG3	1:A:233:GLU:O	2.10	0.51
1:A:392:GLN:OE1	1:A:394:ARG:NH1	2.43	0.51
1:B:201:ARG:NH2	5:B:606:HOH:O	2.38	0.51
1:A:12:SER:HB2	1:A:13:PRO:HD2	1.91	0.51
1:B:166:LEU:HD12	1:B:166:LEU:N	2.27	0.50
1:B:220:SER:HA	1:B:223:GLN:HB3	1.93	0.50
1:B:392:GLN:OE1	1:B:394:ARG:NH1	2.45	0.49
1:A:134:ASN:HB3	1:A:137:GLU:HG2	1.93	0.49
1:A:166:LEU:HD12	1:A:166:LEU:N	2.28	0.49
1:A:234:LYS:HB2	1:A:239:ILE:HD11	1.95	0.48
1:B:175:ASN:O	1:B:175:ASN:ND2	2.42	0.48
1:B:231:SER:HA	1:B:234:LYS:HD2	1.95	0.48
1:B:403:LEU:HG	1:B:404:MET:HE2	1.94	0.48
1:B:102:ARG:NH2	5:B:607:HOH:O	2.41	0.48
1:B:122:TYR:CZ	1:B:257:MET:HB2	2.49	0.47
1:B:86:ALA:HB1	1:B:317:LYS:HE2	1.97	0.47
1:A:244:GLN:O	1:A:273:ARG:HG3	2.15	0.46
1:B:12:SER:HB2	1:B:13:PRO:HD2	1.96	0.46
1:B:239:ILE:HG22	1:B:251:PHE:CE1	2.51	0.46
1:A:7:TRP:CE3	1:A:39:LEU:HD11	2.51	0.46
1:B:101:THR:OG1	1:B:102:ARG:NH1	2.49	0.46
1:A:22:LEU:HD22	1:A:328:TYR:CD1	2.51	0.45
1:B:20:LYS:HE3	5:B:608:HOH:O	2.15	0.45
1:A:281:VAL:HG12	1:A:355:THR:HG21	1.98	0.45
1:B:173:GLY:O	1:B:174:LYS:HE3	2.17	0.45
1:A:118:LEU:HD11	1:A:359:ALA:HB3	1.98	0.45
1:A:117:GLU:CG	1:A:256:TRP:HZ3	2.31	0.44
4:A:503:EDO:H22	5:A:617:HOH:O	2.16	0.44
1:B:230:GLU:H	1:B:230:GLU:CD	2.24	0.44
1:B:309:LYS:HB2	1:B:310:PRO:HD3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:390:VAL:HG22	1:B:395:LEU:HD22	1.99	0.44
1:B:172:PRO:HB3	1:B:239:ILE:HD13	2.00	0.44
1:B:247:LYS:NZ	5:B:613:HOH:O	2.50	0.44
1:A:122:TYR:CZ	1:A:257:MET:HB2	2.52	0.43
1:A:45:TRP:CH2	1:A:373:GLY:HA3	2.53	0.43
1:B:224:LYS:HE2	5:B:715:HOH:O	2.17	0.43
1:A:235:ASN:O	1:A:239:ILE:HD13	2.19	0.43
1:B:22:LEU:HD22	1:B:328:TYR:CD1	2.54	0.43
1:A:113:PRO:HA	1:A:299:ASN:OD1	2.18	0.43
1:B:263:VAL:HG13	1:B:267:LYS:HB2	2.01	0.43
1:A:233:GLU:HB2	5:A:655:HOH:O	2.18	0.42
1:A:94:LEU:HD13	1:A:339:ARG:CD	2.49	0.42
1:B:206:LEU:HD12	1:B:411:ILE:HD12	2.01	0.42
1:B:117:GLU:CG	1:B:256:TRP:HZ3	2.33	0.42
1:B:215:LEU:HD23	1:B:215:LEU:HA	1.83	0.42
1:B:390:VAL:CG2	1:B:395:LEU:HD22	2.50	0.41
1:B:12:SER:HB3	1:B:332:THR:HG22	2.02	0.41
1:A:94:LEU:HD12	1:A:94:LEU:HA	1.94	0.41
1:B:45:TRP:CH2	1:B:373:GLY:HA3	2.55	0.41
1:A:12:SER:HB3	1:A:332:THR:HG22	2.03	0.41
1:B:175:ASN:ND2	1:B:175:ASN:C	2.79	0.41
1:B:172:PRO:CA	1:B:239:ILE:HD13	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	414/423 (98%)	409 (99%)	5 (1%)	0	100	100
1	B	416/423 (98%)	410 (99%)	6 (1%)	0	100	100
All	All	830/846 (98%)	819 (99%)	11 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	315/322 (98%)	299 (95%)	16 (5%)	21	26
1	B	317/322 (98%)	289 (91%)	28 (9%)	9	10
All	All	632/644 (98%)	588 (93%)	44 (7%)	14	15

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	GLN
1	A	94	LEU
1	A	102	ARG
1	A	117	GLU
1	A	143	GLN
1	A	215	LEU
1	A	223	GLN
1	A	230	GLU
1	A	251	PHE
1	A	263	VAL
1	A	273	ARG
1	A	289	LYS
1	A	291	ARG
1	A	346	SER
1	A	365	THR
1	A	403	LEU
1	B	2	LYS
1	B	33	GLU
1	B	35	LYS
1	B	47	LYS
1	B	102	ARG
1	B	107	ARG
1	B	108	GLN
1	B	117	GLU
1	B	118	LEU

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Mol	Chain	Res	Type
1	B	129	LYS
1	B	174	LYS
1	B	175	ASN
1	B	176	SER
1	B	193	GLU
1	B	197	GLN
1	B	203	GLN
1	B	215	LEU
1	B	223	GLN
1	B	231	SER
1	B	233	GLU
1	B	239	ILE
1	B	251	PHE
1	B	263	VAL
1	B	291	ARG
1	B	317	LYS
1	B	339	ARG
1	B	361	GLN
1	B	418	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	182	ASN
1	B	30	HIS
1	B	244	GLN
1	B	381	GLN
1	B	413	GLN

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

4 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	BGC	C	1	2	12,12,12	0.45	0	17,17,17	1.55	3 (17%)
2	BGC	C	2	2	11,11,12	0.61	0	15,15,17	0.97	0
2	BGC	D	1	2	12,12,12	0.51	0	17,17,17	1.52	3 (17%)
2	BGC	D	2	2	11,11,12	0.69	0	15,15,17	1.11	2 (13%)

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	BGC	C	1	2	-	1/2/22/22	0/1/1/1
2	BGC	C	2	2	-	0/2/19/22	0/1/1/1
2	BGC	D	1	2	-	0/2/22/22	0/1/1/1
2	BGC	D	2	2	-	1/2/19/22	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	D	1	BGC	O6-C6-C5	-3.26	100.22	111.33
2	D	1	BGC	C1-C2-C3	-3.21	103.81	110.36
2	C	1	BGC	C1-O5-C5	-3.05	107.75	113.65
2	D	1	BGC	C1-O5-C5	-2.63	108.57	113.65
2	C	1	BGC	O2-C2-C1	2.61	115.27	109.25
2	D	2	BGC	O5-C5-C6	2.45	112.43	107.66
2	D	2	BGC	C1-O5-C5	2.34	115.32	112.19
2	C	1	BGC	O3-C3-C2	-2.33	104.87	110.38

There are no chirality outliers.

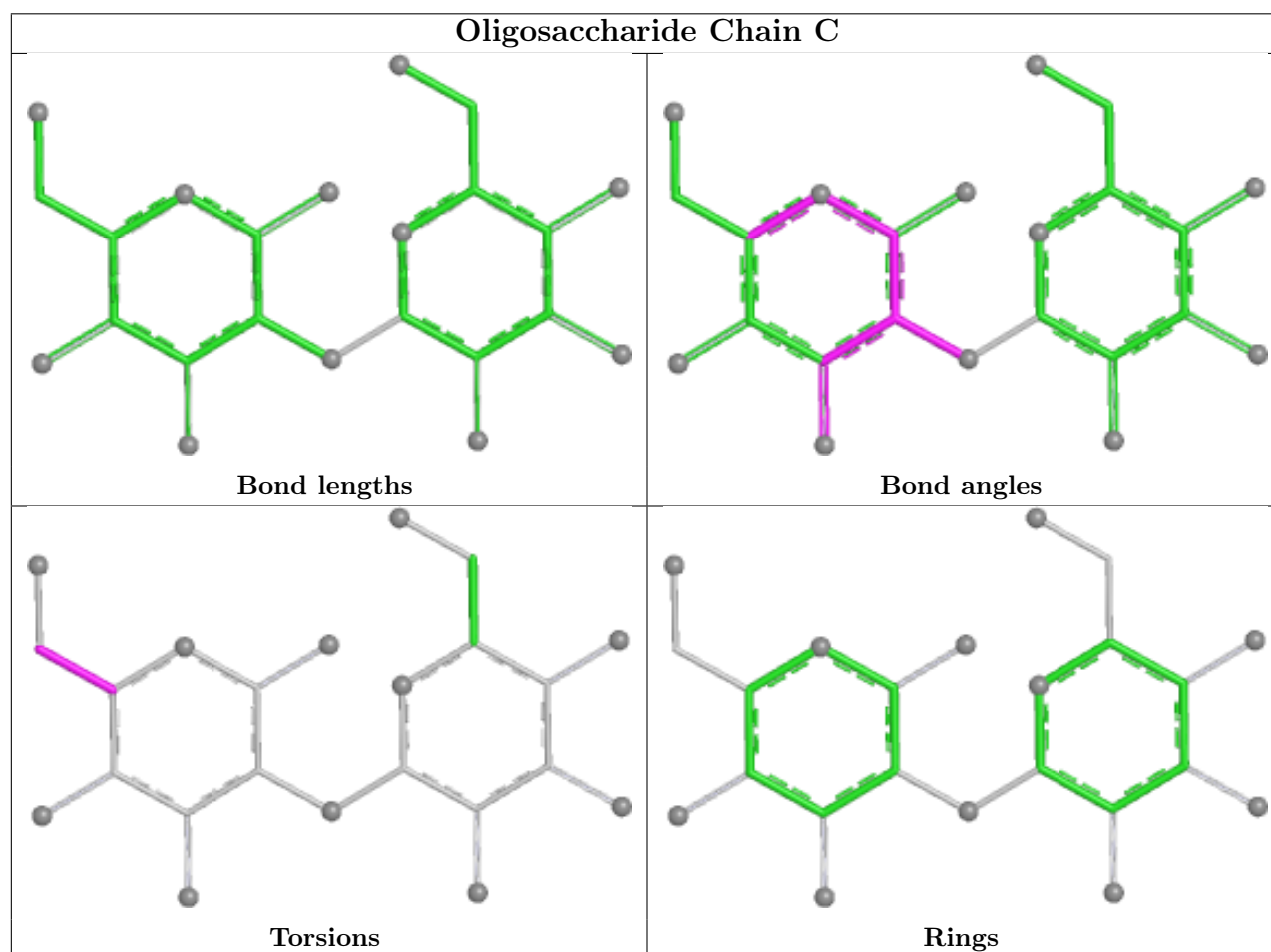
All (2) torsion outliers are listed below:

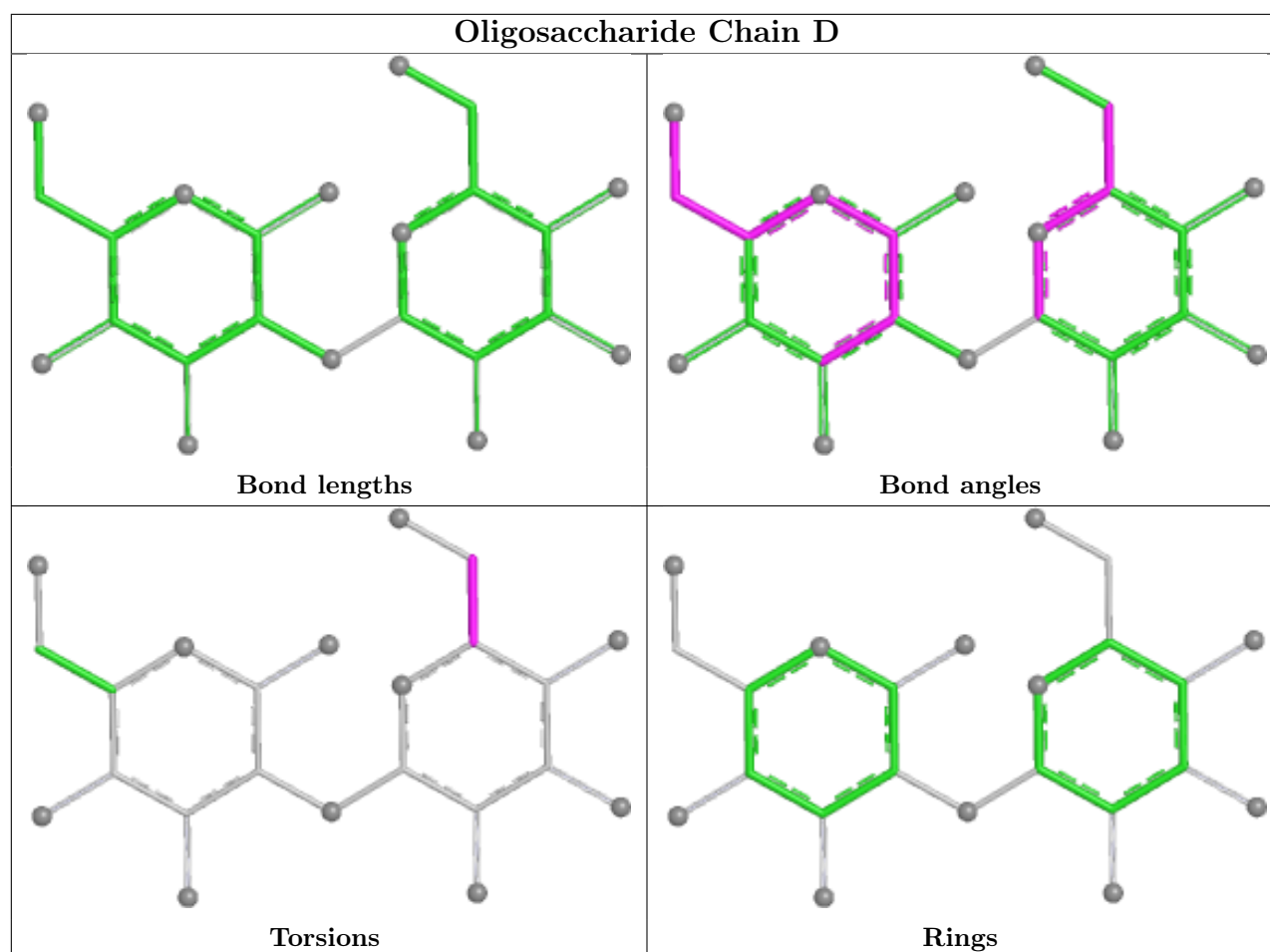
Mol	Chain	Res	Type	Atoms
2	D	2	BGC	C4-C5-C6-O6
2	C	1	BGC	O5-C5-C6-O6

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.





5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
4	EDO	A	504	-	3,3,3	0.08	0	2,2,2	0.16	0
4	EDO	A	503	-	3,3,3	0.09	0	2,2,2	0.13	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
4	EDO	A	504	-	-	1/1/1/1	-
4	EDO	A	503	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	504	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	504	EDO	1	0
4	A	503	EDO	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	416/423 (98%)	0.04	5 (1%) 76 81	10, 22, 37, 55	0
1	B	418/423 (98%)	0.25	11 (2%) 57 63	15, 26, 50, 79	0
All	All	834/846 (98%)	0.14	16 (1%) 66 71	10, 24, 44, 79	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	199	GLY	4.2
1	B	418	HIS	3.4
1	B	419	HIS	3.4
1	B	200	GLY	3.3
1	B	95	PRO	2.8
1	B	345	PRO	2.8
1	A	161	GLU	2.7
1	B	342	TRP	2.7
1	B	198	ALA	2.6
1	B	107	ARG	2.4
1	A	362	PHE	2.4
1	A	198	ALA	2.3
1	A	233	GLU	2.2
1	B	96	ALA	2.2
1	A	201	ARG	2.1
1	B	92	ALA	2.0

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

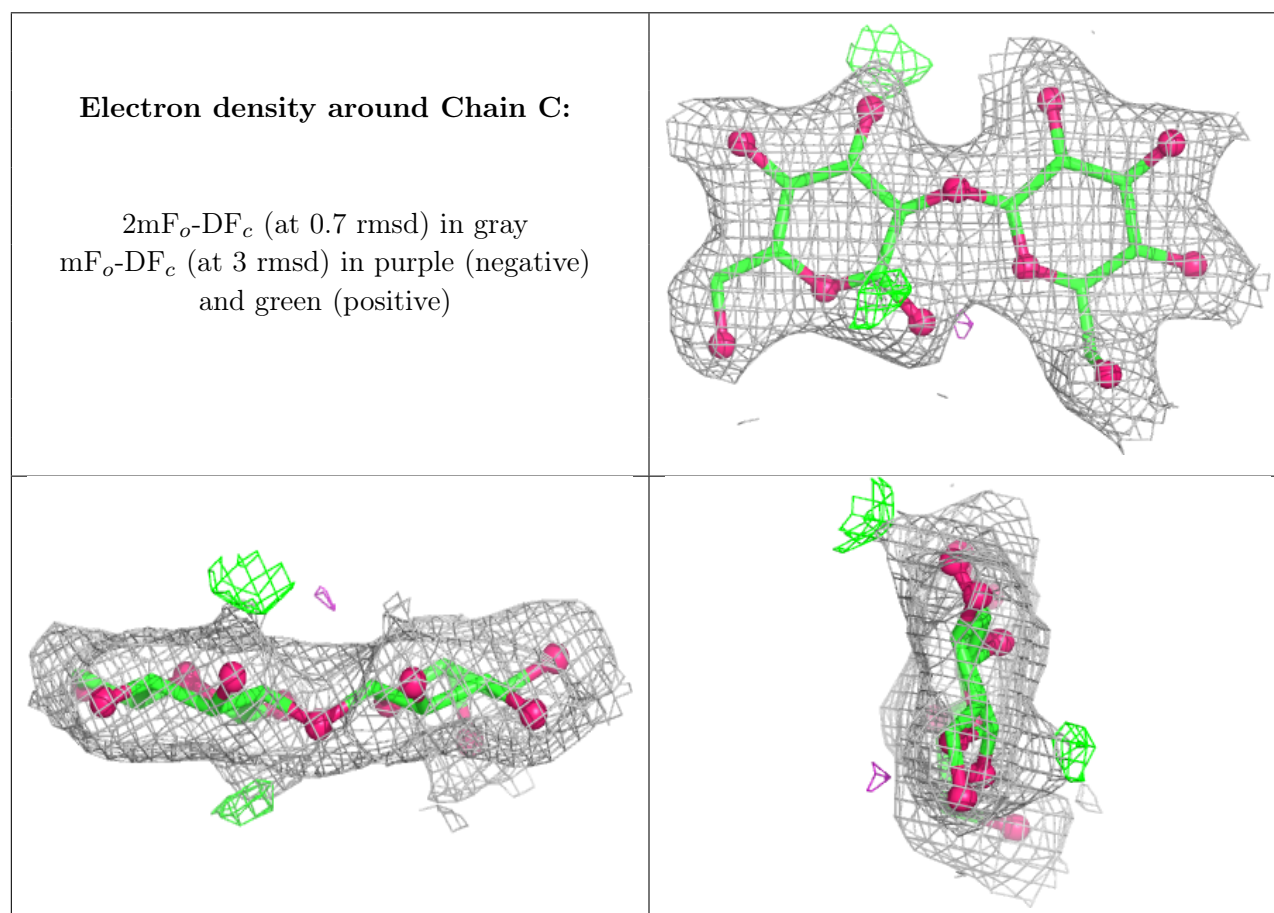
There are no non-standard protein/DNA/RNA residues in this entry.

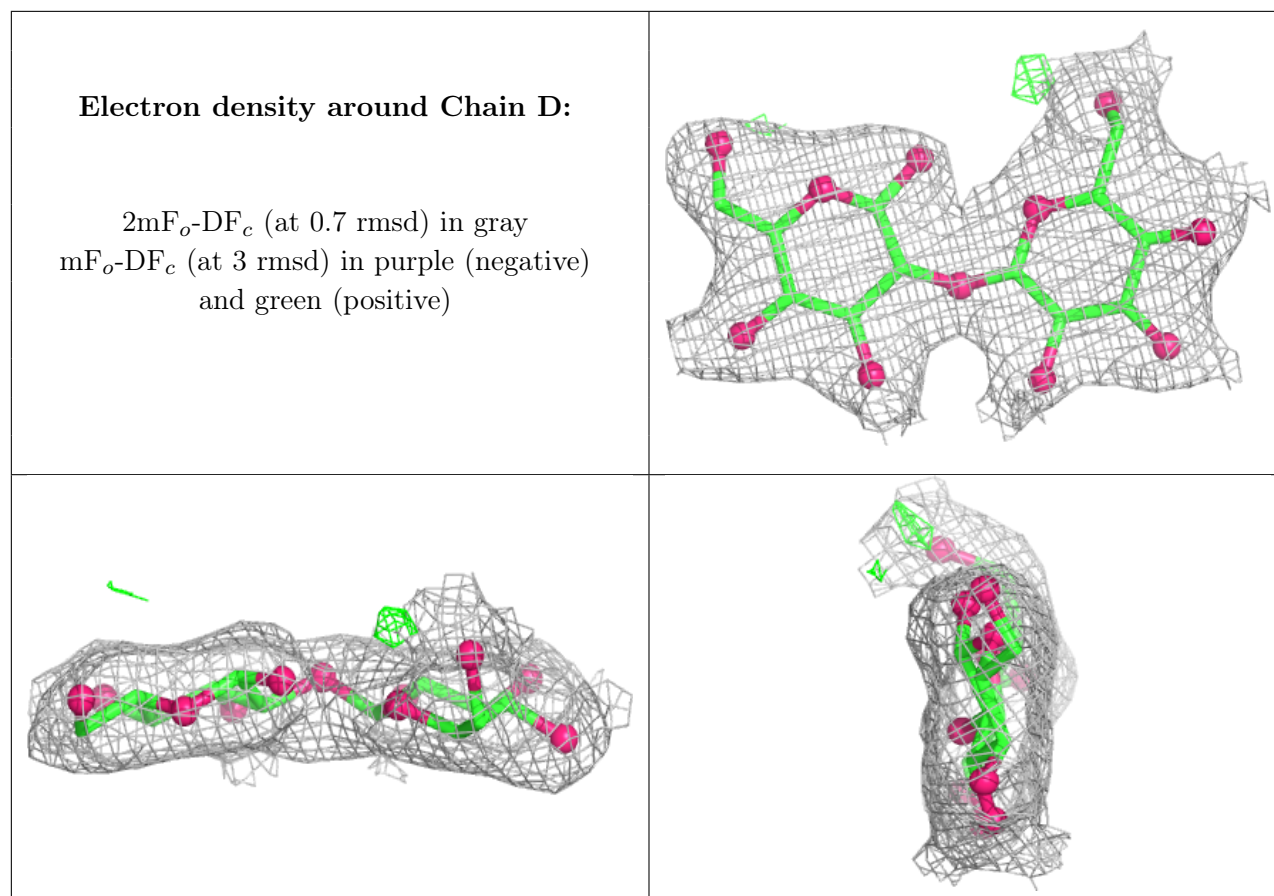
6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BGC	C	1	12/12	0.93	0.08	15,19,22,23	0
2	BGC	D	1	12/12	0.93	0.07	18,20,24,24	0
2	BGC	D	2	11/12	0.93	0.08	15,18,19,20	0
2	BGC	C	2	11/12	0.97	0.05	11,11,13,13	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, 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
4	EDO	A	504	4/4	0.76	0.20	38,39,40,40	0
4	EDO	A	503	4/4	0.85	0.19	37,40,40,42	0
3	CL	A	502	1/1	0.96	0.14	44,44,44,44	0
3	CL	B	502	1/1	0.98	0.04	36,36,36,36	0
3	CL	A	501	1/1	0.98	0.05	16,16,16,16	0
3	CL	B	501	1/1	0.98	0.04	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.