



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:20 PM UTC

PDB ID : 2YJQ / pdb_00002yjq
Title : Structure of a Paenibacillus Polymyxa Xyloglucanase from Glycoside Hydro-
lase Family 44
Authors : Ariza, A.; Eklof, J.M.; Spadiut, O.; Offen, W.A.; Roberts, S.M.; Besenmatter,
W.; Friis, E.P.; Skjot, M.; Wilson, K.S.; Brumer, H.; Davies, G.
Deposited on : 2011-05-23
Resolution : 2.25 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

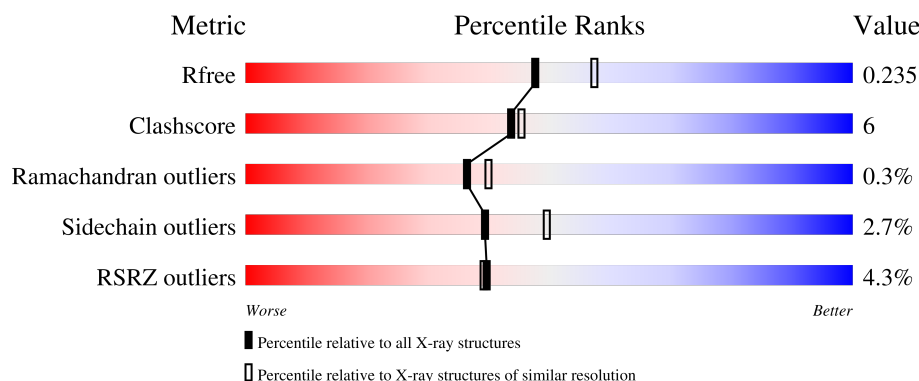
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	524	2% 83% 13% ..
1	B	524	6% 79% 17% ..
2	C	2	100%
2	D	2	50% 50%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 8563 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 CEL44C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	510	Total	C	N	O	S	0	5	0
			4022	2548	676	783	15			
1	B	506	Total	C	N	O	S	0	2	0
			3967	2512	666	775	14			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	PHE	SER	conflict	UNP Q1A2D0
A	19	ASP	GLU	conflict	UNP Q1A2D0
A	68	HIS	GLN	engineered mutation	UNP Q1A2D0
A	92	VAL	THR	engineered mutation	UNP Q1A2D0
A	118	ALA	LYS	engineered mutation	UNP Q1A2D0
A	129	ALA	LYS	engineered mutation	UNP Q1A2D0
A	144	ALA	THR	conflict	UNP Q1A2D0
A	156	TYR	ARG	engineered mutation	UNP Q1A2D0
A	200	PRO	GLY	engineered mutation	UNP Q1A2D0
A	228	VAL	ILE	conflict	UNP Q1A2D0
A	272	VAL	ALA	conflict	UNP Q1A2D0
A	331	PHE	ASN	engineered mutation	UNP Q1A2D0
A	358	SER	GLU	engineered mutation	UNP Q1A2D0
A	403	ALA	PRO	conflict	UNP Q1A2D0
A	473	ILE	SER	conflict	UNP Q1A2D0
B	17	PHE	SER	conflict	UNP Q1A2D0
B	19	ASP	GLU	conflict	UNP Q1A2D0
B	68	HIS	GLN	engineered mutation	UNP Q1A2D0
B	92	VAL	THR	engineered mutation	UNP Q1A2D0
B	118	ALA	LYS	engineered mutation	UNP Q1A2D0
B	129	ALA	LYS	engineered mutation	UNP Q1A2D0
B	144	ALA	THR	conflict	UNP Q1A2D0
B	156	TYR	ARG	engineered mutation	UNP Q1A2D0
B	200	PRO	GLY	engineered mutation	UNP Q1A2D0
B	228	VAL	ILE	conflict	UNP Q1A2D0

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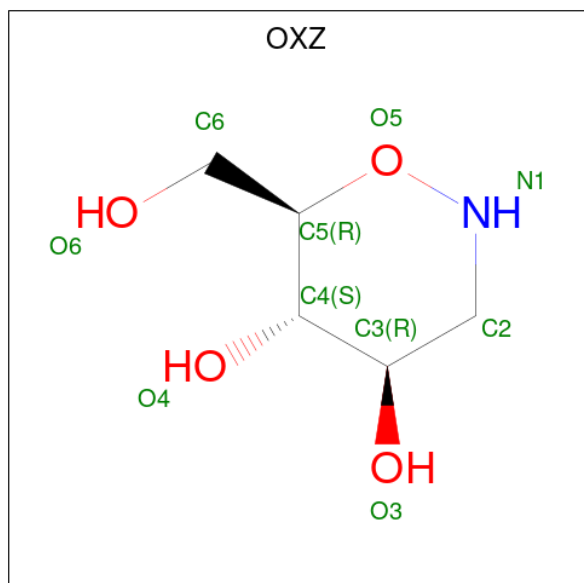
Chain	Residue	Modelled	Actual	Comment	Reference
B	272	VAL	ALA	conflict	UNP Q1A2D0
B	331	PHE	ASN	engineered mutation	UNP Q1A2D0
B	358	SER	GLU	engineered mutation	UNP Q1A2D0
B	403	ALA	PRO	conflict	UNP Q1A2D0
B	473	ILE	SER	conflict	UNP Q1A2D0

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			22	12	10			
2	D	2	Total	C	O	0	0	0
			22	12	10			

- Molecule 3 is TETRAHYDROOXAZINE (CCD ID: OXZ) (formula: C₅H₁₁NO₄).

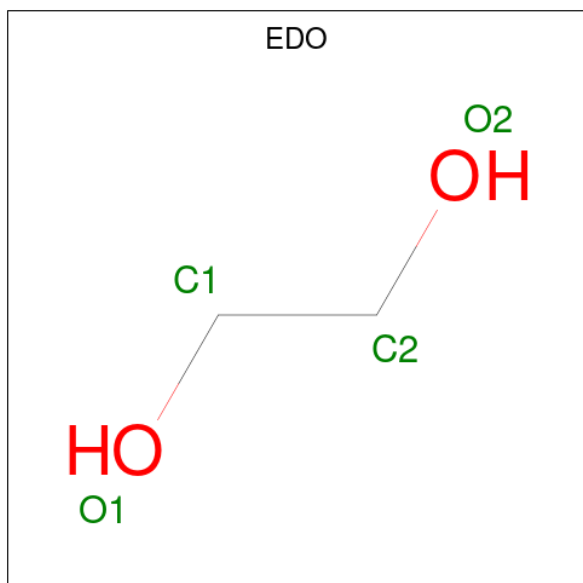


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	5	1	4		
3	B	1	Total	C	N	O	0	0
			10	5	1	4		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		
4	B	1	Total	Ca	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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	1
			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		
5	B	1	Total	C	O	0	0
			4	2	2		

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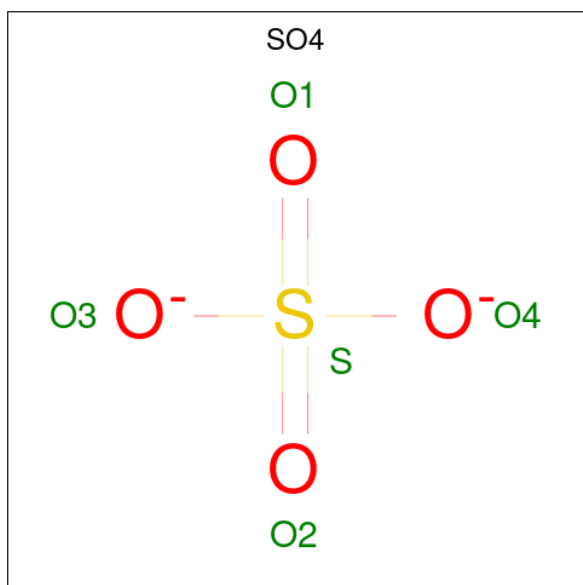
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	1
			4	2	2		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

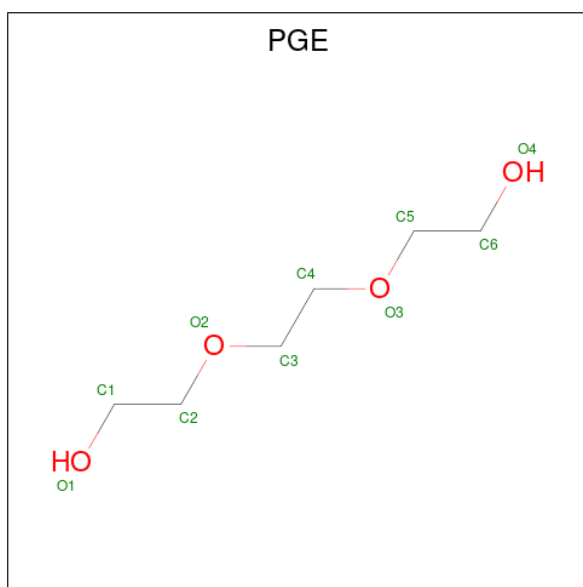
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cl	0	0
			1	1		
6	B	1	Total	Cl	0	0
			1	1		

- Molecule 7 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 8 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	1
			10	6	4		

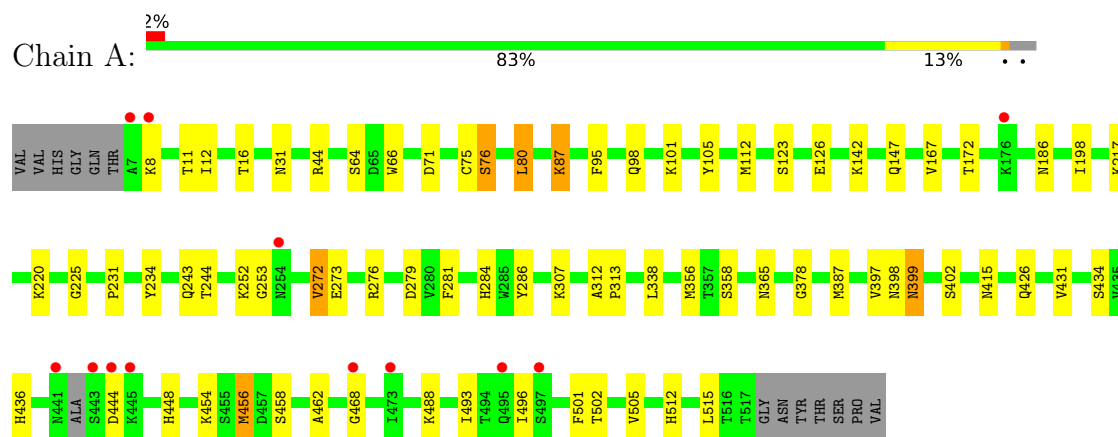
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	258	Total	O	0	4
			262	262		
9	B	181	Total	O	0	3
			184	184		

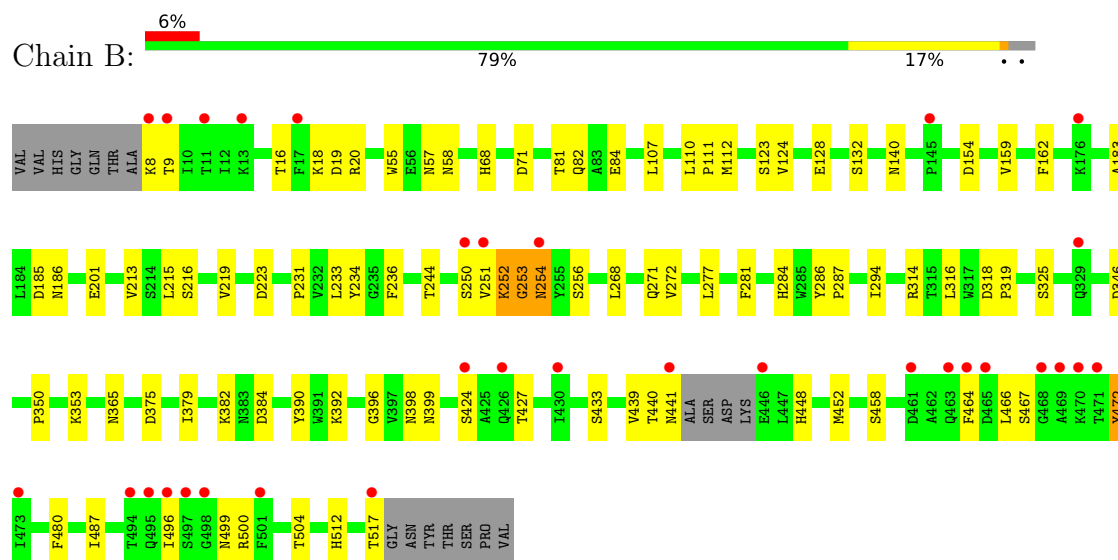
3 Residue-property plots

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: CEL44C



• Molecule 1: CEL44C



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



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

Chain D:  50% 50%

BG1
BG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	84.20Å 84.20Å 320.05Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	106.68 – 2.25 106.68 – 2.25	Depositor EDS
% Data completeness (in resolution range)	98.2 (106.68-2.25) 93.2 (106.68-2.25)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.25Å)	Xtriage
Refinement program	REFMAC 5.6.0086	Depositor
R, R_{free}	0.197 , 0.259 (Not available) , 0.235	Depositor DCC
R_{free} test set	3036 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.004	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.047 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8563	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.54 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.3737e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, EDO, BGC, OXZ, SO4, CA, CL

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	1.22	4/4135 (0.1%)	1.14	10/5618 (0.2%)
1	B	1.16	3/4072 (0.1%)	1.13	7/5534 (0.1%)
All	All	1.19	7/8207 (0.1%)	1.13	17/11152 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	338	LEU	N-CA	6.86	1.52	1.46
1	B	223	ASP	CA-C	-5.82	1.47	1.53
1	A	167	VAL	CA-C	5.80	1.60	1.52
1	B	213	VAL	CA-CB	5.79	1.61	1.54
1	A	31	ASN	C-O	-5.35	1.19	1.23
1	A	505	VAL	CA-CB	5.34	1.59	1.54
1	B	251	VAL	N-CA	5.19	1.51	1.46

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	71	ASP	N-CA-C	6.57	117.14	107.88
1	B	325	SER	N-CA-C	6.26	118.37	108.79
1	A	431	VAL	N-CA-C	-6.19	105.69	111.45
1	A	8	LYS	N-CA-C	5.70	118.28	110.24
1	B	472	TYR	N-CA-C	5.68	117.47	108.67
1	B	294	ILE	N-CA-C	5.62	116.21	108.12
1	A	225	GLY	N-CA-C	5.55	122.59	115.32
1	B	252	LYS	N-CA-C	5.44	119.66	113.19
1	A	142	LYS	N-CA-C	5.31	117.76	111.33
1	A	64	SER	N-CA-C	5.29	118.83	112.38
1	B	140	ASN	N-CA-C	5.26	117.44	111.02
1	A	101	LYS	N-CA-C	-5.25	105.64	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	399	ASN	N-CA-C	5.20	117.35	111.11
1	B	236	PHE	N-CA-C	5.09	117.22	111.11
1	A	272	VAL	N-CA-C	5.07	115.79	110.62
1	A	252	LYS	N-CA-C	5.03	117.50	111.71
1	A	198	ILE	N-CA-C	-5.02	106.90	111.67

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	4022	0	3864	45	0
1	B	3967	0	3806	55	0
2	C	22	0	19	0	0
2	D	22	0	19	0	0
3	A	10	0	10	1	0
3	B	10	0	10	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	20	0	30	5	0
5	B	20	0	30	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	5	0	0	0	0
7	B	5	0	0	0	0
8	A	10	0	14	3	0
9	A	262	0	0	3	0
9	B	184	0	0	5	0
All	All	8563	0	7802	95	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 (95) 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:396:GLY:H	5:B:1529[B]:EDO:H11	1.34	0.89
1:A:448:HIS:HD2	1:A:512:HIS:NE2	1.80	0.79
1:A:243:GLN:HE22	8:A:1529[A]:PGE:H22	1.48	0.79
1:A:415:ASN:HD21	1:B:82:GLN:HE22	1.29	0.78
1:A:243:GLN:HE22	8:A:1529[A]:PGE:C2	1.99	0.75
1:A:126[B]:GLU:OE2	9:A:2090:HOH:O	2.04	0.75
1:B:440:THR:HG22	1:B:441:ASN:H	1.52	0.73
1:B:254:ASN:ND2	1:B:254:ASN:H	1.93	0.66
1:A:488:LYS:NZ	1:B:123:SER:HB3	2.12	0.65
1:A:402:SER:HB2	5:A:1525:EDO:H22	1.79	0.64
1:B:448:HIS:HD2	1:B:512:HIS:NE2	1.96	0.64
1:B:472:TYR:HA	1:B:517:THR:HG22	1.80	0.64
1:A:147:GLN:HA	5:A:1526:EDO:H11	1.82	0.62
1:A:273[B]:GLU:HG2	9:A:2165:HOH:O	2.00	0.61
1:B:231:PRO:HG2	1:B:233:LEU:HG	1.82	0.61
1:A:365:ASN:ND2	1:A:399:ASN:H	1.99	0.60
1:A:448:HIS:CD2	1:A:512:HIS:NE2	2.68	0.60
1:B:112:MET:HB2	1:B:185:ASP:HB2	1.84	0.60
1:B:353:LYS:HE2	9:B:2151:HOH:O	2.01	0.59
1:B:365:ASN:ND2	1:B:399:ASN:H	1.99	0.59
1:B:254:ASN:H	1:B:254:ASN:HD22	1.47	0.59
1:A:307:LYS:HE3	1:A:456:MET:HG3	1.84	0.58
1:A:415:ASN:HD21	1:B:82:GLN:NE2	1.99	0.58
1:B:439:VAL:HG23	1:B:440:THR:O	2.03	0.58
1:B:286:TYR:O	1:B:287:PRO:C	2.49	0.56
1:A:66[C]:TRP:HA	1:A:66[C]:TRP:CE3	2.39	0.56
1:A:16:THR:OG1	1:A:468:GLY:HA3	2.06	0.55
1:A:397:VAL:HB	5:A:1525:EDO:H12	1.88	0.55
1:A:234:TYR:OH	8:A:1529[A]:PGE:H42	2.07	0.55
1:B:346:ASP:HB2	9:B:2145:HOH:O	2.09	0.52
1:B:382:LYS:HE2	9:B:2159:HOH:O	2.10	0.51
1:A:76:SER:HB2	5:A:1522:EDO:H21	1.92	0.51
1:B:124:VAL:HG13	1:B:128:GLU:HB2	1.92	0.51
1:B:375:ASP:O	1:B:379:ILE:HG13	2.11	0.51
1:B:231:PRO:HG2	1:B:233:LEU:CG	2.41	0.50
1:A:365:ASN:HD22	1:A:399:ASN:HB2	1.77	0.49
1:B:55:TRP:CE2	1:B:162:PHE:HB2	2.48	0.49
1:B:392:LYS:HG2	1:B:398:ASN:ND2	2.28	0.49
1:B:440:THR:HG22	1:B:441:ASN:N	2.24	0.48
1:B:81:THR:OG1	1:B:84:GLU:HG3	2.13	0.48
1:A:286:TYR:OH	3:A:1518:OXZ:N1	2.47	0.48
1:A:415:ASN:ND2	1:B:82:GLN:HE22	2.06	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:448:HIS:CD2	1:B:512:HIS:NE2	2.80	0.48
1:A:434:SER:OG	1:A:454:LYS:HE2	2.14	0.47
1:A:87:LYS:HE3	1:A:87:LYS:HB2	1.72	0.47
1:A:66[C]:TRP:HA	1:A:66[C]:TRP:HE3	1.80	0.47
1:A:365:ASN:HD21	1:A:399:ASN:H	1.61	0.47
1:B:186:ASN:HD21	1:B:284:HIS:HE1	1.63	0.47
1:B:159:VAL:HG11	1:B:215:LEU:HD11	1.97	0.46
1:B:16:THR:O	1:B:20:ARG:NH1	2.44	0.46
1:A:95:PHE:HA	1:A:98:GLN:OE1	2.14	0.46
1:A:496:ILE:HD11	1:A:515:LEU:HD22	1.97	0.46
1:A:172:THR:HB	9:A:2121:HOH:O	2.16	0.46
1:B:346:ASP:HA	1:B:350:PRO:HA	1.98	0.46
1:B:353:LYS:CE	9:B:2151:HOH:O	2.61	0.46
1:B:252:LYS:O	1:B:253:GLY:C	2.59	0.45
1:A:462:ALA:O	1:A:502:THR:HA	2.17	0.45
1:B:18:LYS:HG2	1:B:424:SER:HB2	1.98	0.45
1:A:217:LYS:HE2	1:A:273[B]:GLU:HG3	1.98	0.45
1:A:312:ALA:N	1:A:313:PRO:CD	2.80	0.45
1:B:107:LEU:HD11	1:B:183:ALA:HB2	1.99	0.44
1:B:215:LEU:O	1:B:219:VAL:HG23	2.18	0.43
1:B:472:TYR:O	1:B:496:ILE:HB	2.18	0.43
1:B:480:PHE:HD2	1:B:487:ILE:CD1	2.31	0.43
1:A:71:ASP:C	1:A:71:ASP:OD1	2.61	0.43
1:B:464:PHE:O	1:B:500:ARG:HA	2.19	0.43
1:B:110:LEU:HB3	1:B:111:PRO:HD2	2.01	0.43
1:B:128:GLU:O	1:B:132:SER:OG	2.34	0.43
1:B:466:LEU:HB2	1:B:499:ASN:HB3	2.01	0.42
1:B:154:ASP:C	1:B:154:ASP:OD1	2.62	0.42
1:A:231:PRO:HD3	1:A:281:PHE:CE1	2.54	0.42
1:B:244:THR:HG22	1:B:244:THR:O	2.19	0.42
1:B:68:HIS:HD2	9:B:2102:HOH:O	2.01	0.42
1:A:75:CYS:O	1:A:80:LEU:HB2	2.19	0.42
1:A:356:MET:HE3	1:A:356:MET:HB2	1.91	0.42
1:A:488:LYS:HZ1	1:B:123:SER:HB3	1.80	0.42
1:B:268:LEU:O	1:B:271:GLN:HB3	2.20	0.42
1:A:12:ILE:HA	1:A:426:GLN:O	2.20	0.42
1:A:284:HIS:CG	1:A:358:SER:HB3	2.55	0.41
1:A:378:GLY:HA3	1:A:436:HIS:CG	2.55	0.41
1:A:493:ILE:HD13	1:A:501:PHE:CE2	2.55	0.41
1:B:234:TYR:CD1	1:B:234:TYR:C	2.98	0.41
1:A:220:LYS:HE3	1:A:276:ARG:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:ARG:C	1:B:316:LEU:N	2.77	0.41
1:B:318:ASP:HA	1:B:319:PRO:HD2	1.85	0.41
1:B:452:MET:HE2	1:B:452:MET:HB3	1.72	0.41
1:B:216:SER:OG	1:B:277:LEU:HB3	2.19	0.41
1:A:105:TYR:CE1	1:A:387:MET:HE1	2.56	0.41
1:A:220:LYS:NZ	1:A:279:ASP:OD2	2.41	0.40
1:B:390:TYR:OH	1:B:398:ASN:ND2	2.54	0.40
1:B:231:PRO:HD3	1:B:281:PHE:CE1	2.56	0.40
1:B:268:LEU:O	1:B:272:VAL:HG23	2.22	0.40
1:A:398:ASN:H	5:A:1525:EDO:H21	1.87	0.40
1:B:57:ASN:O	1:B:58:ASN:HB2	2.21	0.40
1:A:44:ARG:HH22	1:A:186:ASN:HD22	1.68	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	511/524 (98%)	489 (96%)	21 (4%)	1 (0%)	43	50
1	B	504/524 (96%)	476 (94%)	26 (5%)	2 (0%)	30	31
All	All	1015/1048 (97%)	965 (95%)	47 (5%)	3 (0%)	36	40

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	253	GLY
1	B	250	SER
1	A	253	GLY

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	431/437 (99%)	420 (97%)	11 (3%)	40	50
1	B	425/437 (97%)	412 (97%)	13 (3%)	35	44
All	All	856/874 (98%)	832 (97%)	24 (3%)	39	48

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	THR
1	A	76	SER
1	A	80	LEU
1	A	87	LYS
1	A	112	MET
1	A	123	SER
1	A	244	THR
1	A	272	VAL
1	A	444	ASP
1	A	456	MET
1	A	458	SER
1	B	8	LYS
1	B	9	THR
1	B	19	ASP
1	B	201[A]	GLU
1	B	201[B]	GLU
1	B	254	ASN
1	B	256	SER
1	B	384	ASP
1	B	427	THR
1	B	433	SER
1	B	458	SER
1	B	467	SER
1	B	504	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	A	121	ASN
1	A	143	ASN
1	A	153	ASN
1	A	186	ASN
1	A	193	HIS
1	A	243	GLN
1	A	298	ASN
1	A	343	GLN
1	A	365	ASN
1	A	389	ASN
1	A	415	ASN
1	A	448	HIS
1	A	486	GLN
1	B	68	HIS
1	B	121	ASN
1	B	155	ASN
1	B	186	ASN
1	B	243	GLN
1	B	254	ASN
1	B	265	GLN
1	B	298	ASN
1	B	365	ASN
1	B	389	ASN
1	B	398	ASN
1	B	399	ASN
1	B	432	ASN
1	B	441	ASN
1	B	448	HIS
1	B	486	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	2,3	11,11,12	0.73	1 (9%)	15,15,17	1.53	4 (26%)
2	BGC	C	2	2	11,11,12	0.64	0	15,15,17	1.82	3 (20%)
2	BGC	D	1	2,3	11,11,12	0.83	0	15,15,17	1.62	3 (20%)
2	BGC	D	2	2	11,11,12	0.75	0	15,15,17	0.86	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
2	BGC	C	1	2,3	-	0/2/19/22	0/1/1/1
2	BGC	C	2	2	-	2/2/19/22	0/1/1/1
2	BGC	D	1	2,3	-	0/2/19/22	0/1/1/1
2	BGC	D	2	2	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	BGC	O4-C4	-2.05	1.37	1.43

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	BGC	O5-C5-C6	4.93	117.27	107.66
2	D	1	BGC	C1-O5-C5	3.64	117.07	112.19
2	D	1	BGC	C1-C2-C3	-3.26	104.90	109.64
2	C	1	BGC	O3-C3-C4	-2.64	104.14	110.38
2	C	1	BGC	O4-C4-C5	-2.61	102.90	109.32
2	C	1	BGC	C3-C4-C5	2.36	114.50	110.23
2	C	2	BGC	O5-C1-C2	-2.19	105.56	110.79
2	C	2	BGC	C6-C5-C4	-2.17	107.69	113.02
2	C	1	BGC	O6-C6-C5	-2.10	104.18	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	BGC	C2-C3-C4	2.07	114.51	110.86

There are no chirality outliers.

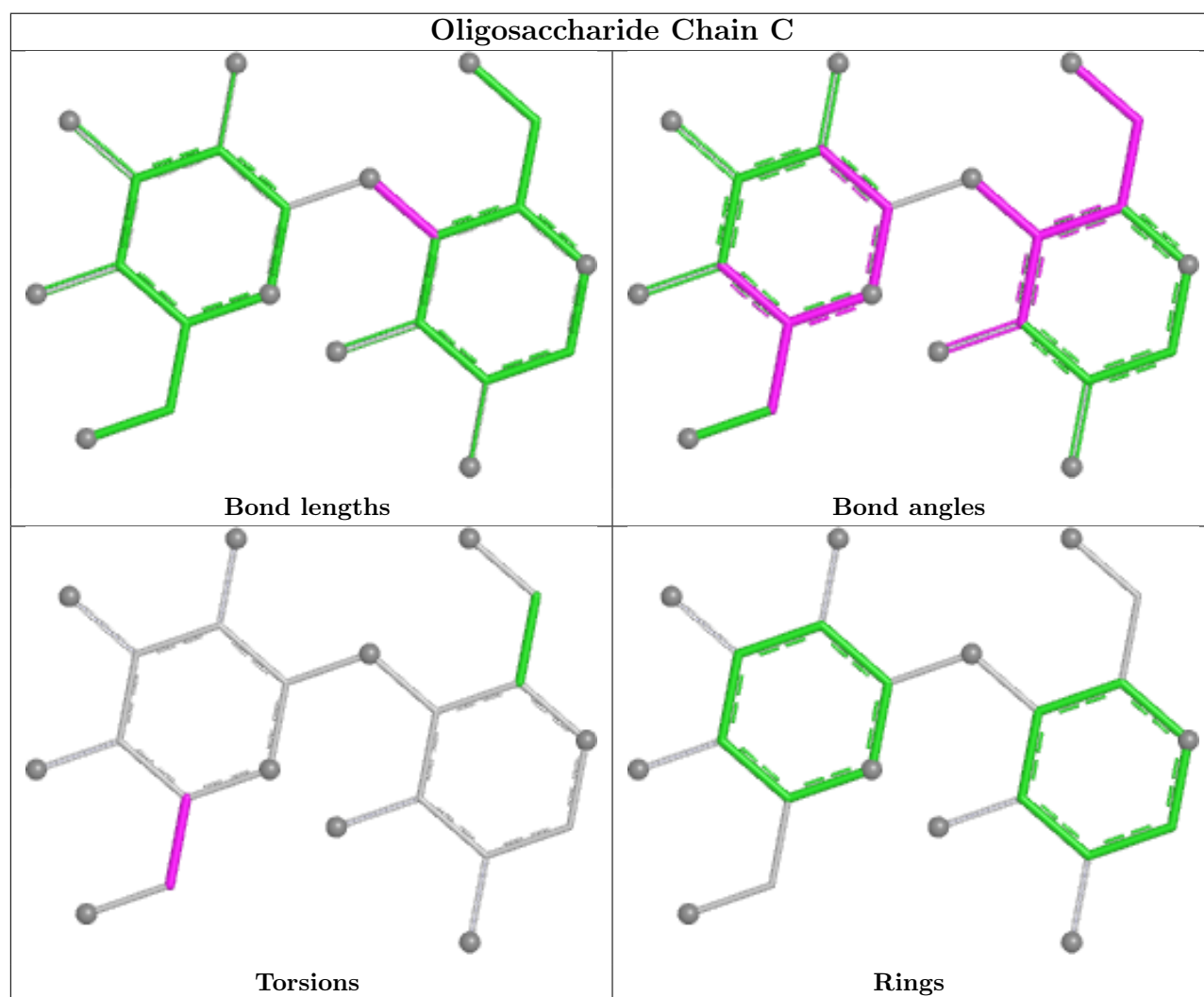
All (2) torsion outliers are listed below:

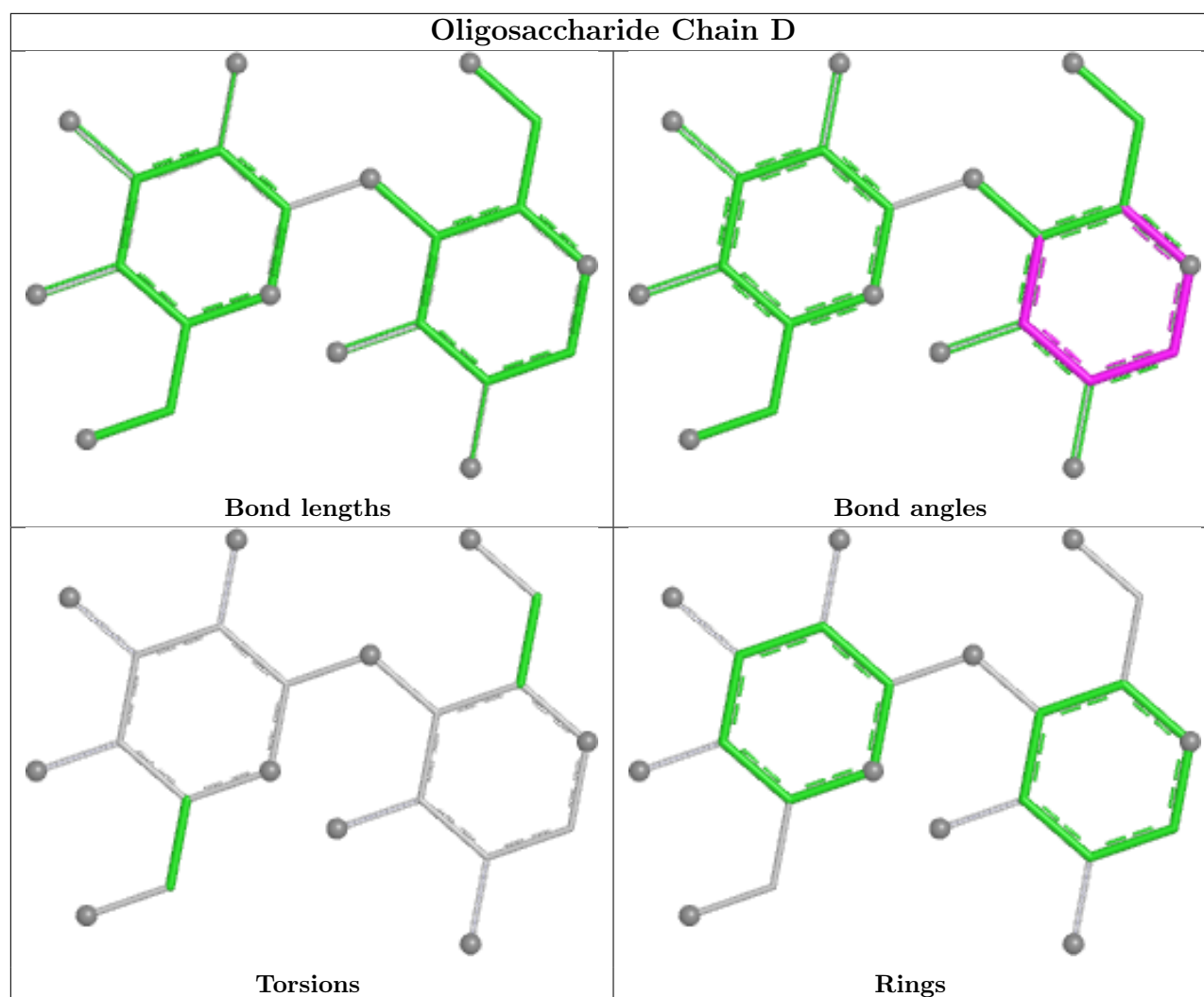
Mol	Chain	Res	Type	Atoms
2	C	2	BGC	O5-C5-C6-O6
2	C	2	BGC	C4-C5-C6-O6

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.





5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 4 are monoatomic - leaving 15 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	SO4	A	1528	-	4,4,4	0.17	0	6,6,6	0.61	0
5	EDO	A	1523	-	3,3,3	0.25	0	2,2,2	1.22	0
8	PGE	A	1529[A]	-	9,9,9	0.55	0	8,8,8	0.47	0
5	EDO	A	1522	-	3,3,3	0.43	0	2,2,2	0.71	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	OXZ	B	1518	2	8,10,10	0.67	0	8,13,13	1.60	2 (25%)
5	EDO	A	1525	-	3,3,3	0.51	0	2,2,2	0.11	0
5	EDO	B	1526	-	3,3,3	0.30	0	2,2,2	0.56	0
5	EDO	A	1530[B]	-	3,3,3	0.45	0	2,2,2	0.40	0
5	EDO	B	1524	-	3,3,3	0.48	0	2,2,2	0.87	0
5	EDO	B	1527	-	3,3,3	0.37	0	2,2,2	0.52	0
5	EDO	B	1529[B]	-	3,3,3	0.61	0	2,2,2	0.08	0
7	SO4	B	1528[A]	-	4,4,4	0.20	0	6,6,6	0.32	0
3	OXZ	A	1518	2	8,10,10	1.66	1 (12%)	8,13,13	1.39	2 (25%)
5	EDO	B	1522	-	3,3,3	0.50	0	2,2,2	0.48	0
5	EDO	A	1526	-	3,3,3	0.34	0	2,2,2	0.43	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	1523	-	-	1/1/1/1	-
8	PGE	A	1529[A]	-	-	4/7/7/7	-
5	EDO	A	1522	-	-	0/1/1/1	-
3	OXZ	B	1518	2	-	0/2/16/16	0/1/1/1
5	EDO	A	1525	-	-	1/1/1/1	-
5	EDO	B	1526	-	-	1/1/1/1	-
5	EDO	A	1530[B]	-	-	0/1/1/1	-
5	EDO	B	1524	-	-	0/1/1/1	-
5	EDO	B	1527	-	-	0/1/1/1	-
5	EDO	B	1529[B]	-	-	0/1/1/1	-
3	OXZ	A	1518	2	-	0/2/16/16	0/1/1/1
5	EDO	B	1522	-	-	0/1/1/1	-
5	EDO	A	1526	-	-	1/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1518	OXZ	C2-C3	4.01	1.56	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1518	OXZ	C3-C4-C5	2.68	112.68	109.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1518	OXZ	O3-C3-C2	-2.57	104.67	109.65
3	B	1518	OXZ	O4-C4-C5	-2.26	103.76	109.32
3	A	1518	OXZ	C2-C3-C4	2.03	112.73	110.25

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	1529[A]	PGE	C4-C3-O2-C2
5	A	1523	EDO	O1-C1-C2-O2
5	A	1525	EDO	O1-C1-C2-O2
5	A	1526	EDO	O1-C1-C2-O2
8	A	1529[A]	PGE	O2-C3-C4-O3
8	A	1529[A]	PGE	O3-C5-C6-O4
5	B	1526	EDO	O1-C1-C2-O2
8	A	1529[A]	PGE	C6-C5-O3-C4

There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	1529[A]	PGE	3	0
5	A	1522	EDO	1	0
5	A	1525	EDO	3	0
5	B	1529[B]	EDO	1	0
3	A	1518	OXZ	1	0
5	A	1526	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 ⓘ

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	510/524 (97%)	-0.28	12 (2%)	59 61	14, 24, 39, 52	34 (6%)
1	B	506/524 (96%)	0.33	32 (6%)	26 24	15, 30, 47, 57	69 (13%)
All	All	1016/1048 (96%)	0.03	44 (4%)	40 39	14, 27, 44, 57	103 (10%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	17	PHE	7.7
1	B	441	ASN	7.7
1	B	498	GLY	7.0
1	B	254	ASN	6.7
1	B	473	ILE	6.4
1	B	495	GLN	5.8
1	B	494	THR	5.1
1	B	8	LYS	4.8
1	B	471	THR	4.8
1	A	7	ALA	4.7
1	A	473	ILE	4.6
1	B	501	PHE	4.4
1	B	463	GLN	4.2
1	B	250	SER	4.1
1	B	497	SER	4.1
1	B	251	VAL	4.0
1	B	496	ILE	4.0
1	B	176	LYS	3.9
1	B	329	GLN	3.8
1	B	13	LYS	3.5
1	A	443	SER	3.4
1	A	444	ASP	3.1
1	B	430	ILE	3.1
1	A	441	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	8	LYS	2.9
1	B	468	GLY	2.8
1	B	517	THR	2.8
1	A	445	LYS	2.7
1	A	254	ASN	2.7
1	B	446	GLU	2.7
1	B	461	ASP	2.6
1	B	424	SER	2.6
1	B	145	PRO	2.6
1	B	464	PHE	2.4
1	A	176	LYS	2.4
1	B	426	GLN	2.4
1	A	497	SER	2.3
1	B	469	ALA	2.3
1	B	465	ASP	2.2
1	B	9	THR	2.2
1	B	470	LYS	2.2
1	B	11	THR	2.2
1	A	495	GLN	2.1
1	A	468	GLY	2.1

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

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

6.3 Carbohydrates [i](#)

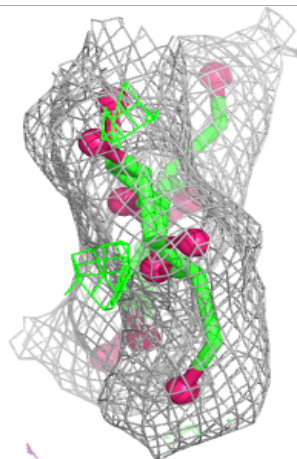
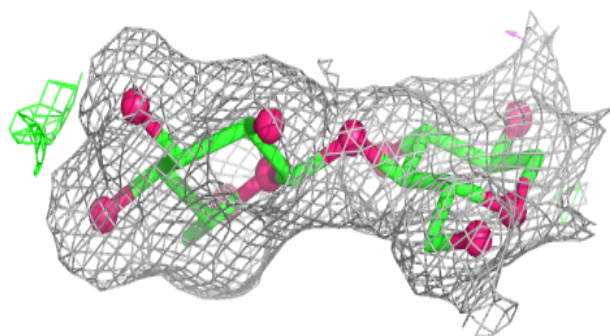
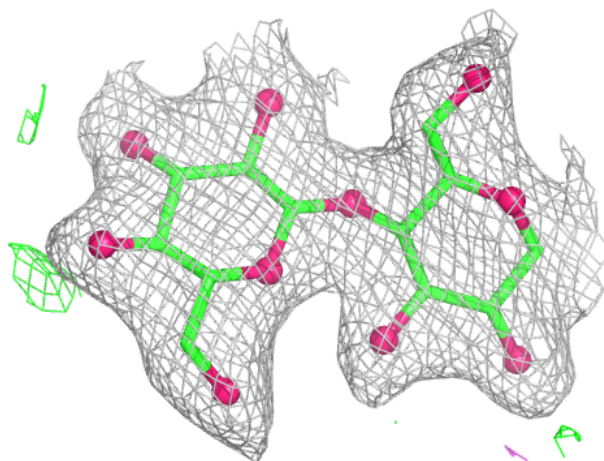
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

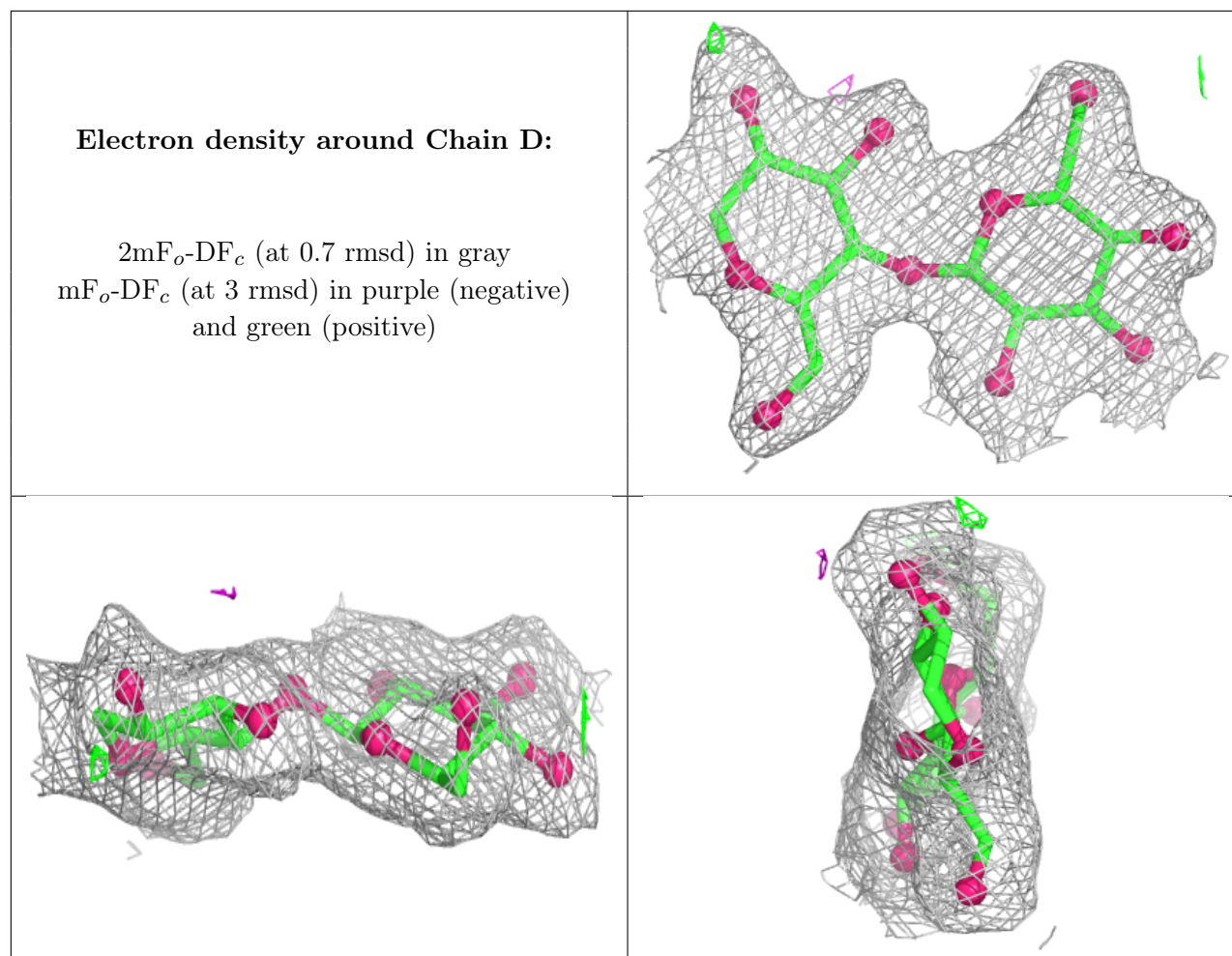
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BGC	C	1	11/12	0.97	0.06	19,20,21,23	0
2	BGC	C	2	11/12	0.97	0.05	21,22,24,27	0
2	BGC	D	1	11/12	0.97	0.06	23,25,27,27	0
2	BGC	D	2	11/12	0.97	0.06	23,26,29,35	0

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 C:

$2mF_o-DF_c$ (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
8	PGE	A	1529[A]	10/10	0.80	0.16	30,32,33,33	10
5	EDO	B	1524	4/4	0.88	0.14	45,47,47,50	0
5	EDO	A	1522	4/4	0.90	0.09	41,44,45,45	0
5	EDO	B	1527	4/4	0.91	0.17	21,22,22,22	4
5	EDO	A	1525	4/4	0.93	0.18	38,38,38,41	0
5	EDO	B	1526	4/4	0.93	0.21	18,19,19,20	4
5	EDO	B	1529[B]	4/4	0.94	0.07	25,27,27,27	4
5	EDO	A	1530[B]	4/4	0.94	0.15	51,52,52,52	4
7	SO4	A	1528	5/5	0.95	0.07	22,22,23,23	5
5	EDO	A	1523	4/4	0.96	0.09	47,47,48,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	EDO	B	1522	4/4	0.96	0.07	46,46,47,49	0
3	OXZ	B	1518	10/10	0.96	0.06	20,22,22,25	0
5	EDO	A	1526	4/4	0.96	0.10	38,40,41,41	1
7	SO4	B	1528[A]	5/5	0.97	0.08	34,34,36,36	5
3	OXZ	A	1518	10/10	0.97	0.06	15,16,17,18	0
6	CL	B	1525	1/1	0.98	0.04	28,28,28,28	0
4	CA	B	1521	1/1	0.98	0.05	25,25,25,25	0
4	CA	A	1521	1/1	0.99	0.04	21,21,21,21	0
6	CL	A	1527	1/1	0.99	0.04	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.