



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 06:11 PM EDT

PDB ID : 5TA5 / pdb_00005ta5
Title : Crystal structure of BuGH86wt in complex with neoagarooctase
Authors : Pluvinage, B.; Boraston, A.B.
Deposited on : 2016-09-09
Resolution : 1.65 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

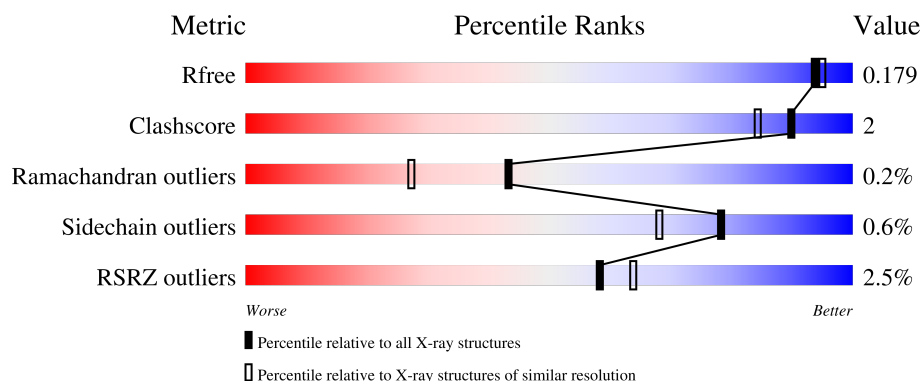
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	649	
1	B	649	
2	C	2	
3	D	2	

2 Entry composition [i](#)

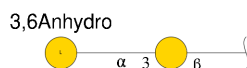
There are 7 unique types of molecules in this entry. The entry contains 11778 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	624	Total	C	N	O	S	11	24	0
			5262	3352	886	991	33			
1	B	613	Total	C	N	O	S	24	24	0
			5148	3278	869	968	33			

- Molecule 2 is an oligosaccharide called 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose.



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

- Molecule 3 is an oligosaccharide called beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-L-galactopyranose.



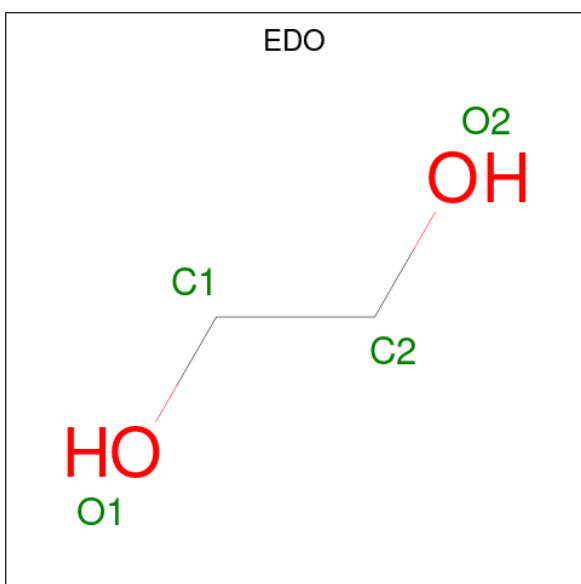
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	D	2	Total	C	O	0	0	0
			22	12	10			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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



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

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

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total 3	Ca 3	0	0
6	B	3	Total 3	Ca 3	0	0

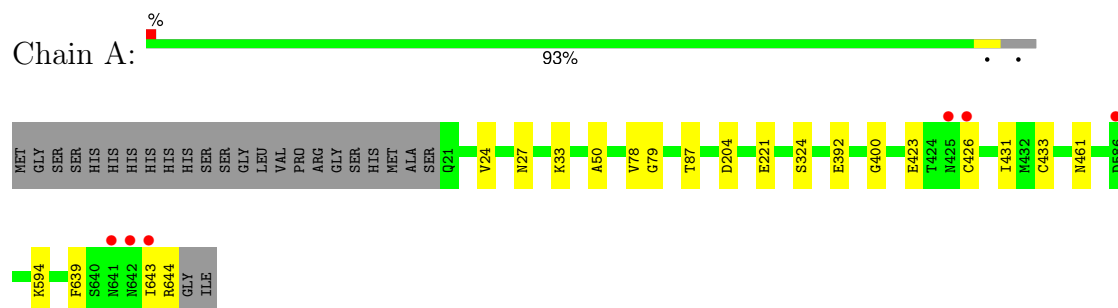
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	589	Total 589	O 589	0	0
7	B	536	Total 536	O 536	0	1

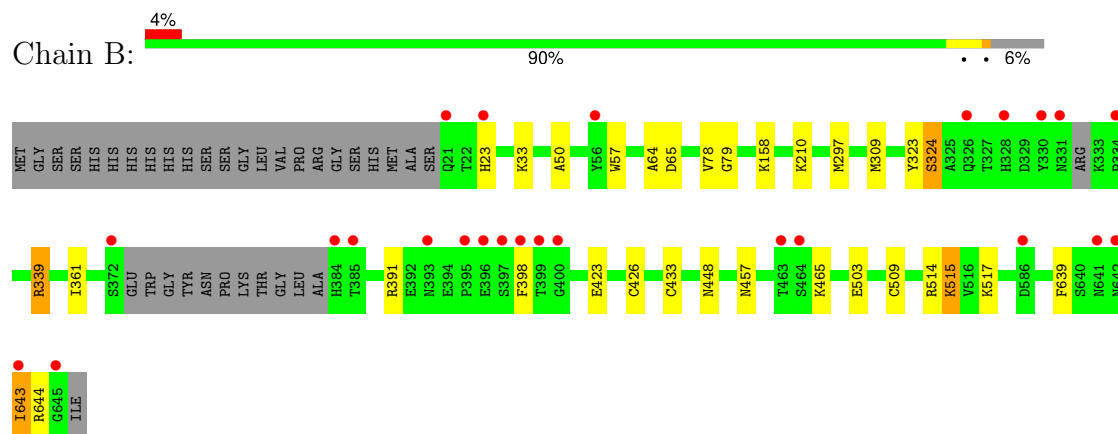
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Glycoside Hydrolase



- Molecule 1: Glycoside Hydrolase



- Molecule 2: 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose



- Molecule 3: beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-L-galactopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.74Å 73.60Å 83.39Å 85.84° 86.19° 71.82°	Depositor
Resolution (Å)	83.08 – 1.65 83.08 – 1.65	Depositor EDS
% Data completeness (in resolution range)	97.1 (83.08-1.65) 97.1 (83.08-1.65)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	50.00 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.144 , 0.169 0.157 , 0.179	Depositor DCC
R_{free} test set	8112 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	10.3	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 41.5	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	0.96	EDS
Total number of atoms	11778	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, CA, GAL, EDO, AAL

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	0.66	1/5390 (0.0%)	0.78	1/7280 (0.0%)
1	B	1.07	7/5268 (0.1%)	0.89	7/7110 (0.1%)
All	All	0.88	8/10658 (0.1%)	0.84	8/14390 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	514	ARG	CD-NE	-46.55	0.81	1.46
1	B	517	LYS	CE-NZ	-25.77	0.72	1.49
1	B	515	LYS	CD-CE	18.80	2.08	1.52
1	B	210	LYS	CD-CE	16.82	2.02	1.52
1	B	465	LYS	CD-CE	-15.10	1.07	1.52
1	A	594	LYS	CD-CE	-8.99	1.25	1.52
1	B	33	LYS	CE-NZ	8.98	1.76	1.49
1	B	158	LYS	CD-CE	7.77	1.75	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	514	ARG	CD-NE-CZ	22.54	155.96	124.40
1	B	514	ARG	CG-CD-NE	19.04	153.90	112.00
1	B	517	LYS	CD-CE-NZ	14.38	157.91	111.90
1	B	465	LYS	CG-CD-CE	11.56	137.89	111.30
1	B	210	LYS	CG-CD-CE	-10.66	86.79	111.30
1	A	594	LYS	CG-CD-CE	9.30	132.68	111.30
1	B	158	LYS	CG-CD-CE	-7.49	94.07	111.30
1	B	33	LYS	CD-CE-NZ	-7.41	88.19	111.90

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	5262	0	5096	15	0
1	B	5148	0	4988	22	0
2	C	22	0	19	0	0
3	D	22	0	19	0	0
4	A	5	0	0	0	0
4	B	20	0	0	1	0
5	A	80	0	120	2	0
5	B	88	0	132	3	0
6	A	3	0	0	0	0
6	B	3	0	0	0	0
7	A	589	0	0	1	0
7	B	536	0	0	4	0
All	All	11778	0	10374	37	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 2.

All (37) 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:339[A]:ARG:HH21	1:B:339[A]:ARG:HB3	1.37	0.89
1:B:65[B]:ASP:OD1	5:B:723:EDO:H21	1.74	0.86
1:B:64[A]:ALA:O	7:B:801:HOH:O	1.95	0.83
1:A:426[B]:CYS:HG	1:A:433[B]:CYS:HG	1.11	0.77
1:B:339[A]:ARG:HH21	1:B:339[A]:ARG:CB	2.04	0.71
1:A:426[B]:CYS:HG	1:A:433[B]:CYS:CB	2.09	0.65
1:B:423:GLU:OE2	1:B:644:ARG:NH1	2.36	0.58
1:B:426[B]:CYS:SG	1:B:433[B]:CYS:SG	3.02	0.57
1:B:323:TYR:O	1:B:324:SER:HB2	2.04	0.57
1:B:65[B]:ASP:OD1	5:B:723:EDO:C2	2.50	0.57
1:A:639[B]:PHE:CE1	1:A:643:ILE:HD11	2.41	0.55
1:B:391[B]:ARG:NH2	7:B:807:HOH:O	2.40	0.54
1:A:423:GLU:OE2	1:A:644:ARG:NH2	2.41	0.53
1:A:24[B]:VAL:HG12	1:A:426[B]:CYS:HB3	1.90	0.53
1:B:509:CYS:HB3	7:B:1272:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:426[A]:CYS:SG	1:A:431:ILE:HB	2.50	0.51
1:B:398:PHE:CB	7:B:905:HOH:O	2.58	0.51
1:B:23[B]:HIS:HD2	1:B:457[B]:ASN:O	1.94	0.50
1:A:33:LYS:NZ	4:B:702:SO4:O4	2.31	0.50
1:A:87:THR:HA	5:A:714:EDO:H21	1.94	0.49
1:A:639[B]:PHE:CD1	1:A:643:ILE:HG12	2.48	0.48
1:A:78:VAL:HB	1:A:79:GLY:HA3	1.94	0.48
5:A:709:EDO:H21	1:B:309:MET:HE3	1.96	0.48
1:A:423:GLU:OE2	1:A:644:ARG:NH1	2.47	0.47
1:B:639[B]:PHE:CD1	1:B:643:ILE:HG12	2.50	0.47
1:B:64[A]:ALA:O	1:B:65[A]:ASP:C	2.57	0.46
1:A:392:GLU:HG2	1:A:400:GLY:HA3	1.99	0.44
1:B:57:TRP:HB3	5:B:723:EDO:H12	1.99	0.44
1:B:339[A]:ARG:CB	1:B:339[A]:ARG:NH2	2.77	0.43
1:B:50:ALA:O	1:B:79:GLY:HA3	2.18	0.43
1:B:78:VAL:HB	1:B:79:GLY:HA3	2.01	0.43
1:A:50:ALA:O	1:A:79:GLY:HA3	2.19	0.43
1:B:448:ASN:O	1:B:503:GLU:HA	2.18	0.43
1:B:423:GLU:OE2	1:B:644:ARG:NH2	2.52	0.42
1:A:27:ASN:HD22	1:A:461:ASN:HB2	1.84	0.42
1:B:639[B]:PHE:CE1	1:B:643:ILE:HD11	2.54	0.42
1:A:221[B]:GLU:HG3	7:A:906:HOH:O	2.20	0.41

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	647/649 (100%)	625 (97%)	21 (3%)	1 (0%)	43	27
1	B	631/649 (97%)	609 (96%)	21 (3%)	1 (0%)	43	27
All	All	1278/1298 (98%)	1234 (97%)	42 (3%)	2 (0%)	43	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	324	SER
1	B	324	SER

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	576/575 (100%)	575 (100%)	1 (0%)	87	83
1	B	562/575 (98%)	555 (99%)	7 (1%)	63	45
All	All	1138/1150 (99%)	1130 (99%)	8 (1%)	78	64

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

Mol	Chain	Res	Type
1	A	204	ASP
1	B	297[A]	MET
1	B	297[B]	MET
1	B	339[A]	ARG
1	B	339[B]	ARG
1	B	361	ILE
1	B	515	LYS
1	B	643	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	91	ASN
1	A	252	GLN
1	A	441	ASN
1	A	447	ASN
1	A	461	ASN
1	A	511	ASN

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Mol	Chain	Res	Type
1	A	547	ASN
1	B	289	GLN
1	B	328	HIS
1	B	425	ASN
1	B	441	ASN
1	B	461	ASN
1	B	511	ASN
1	B	547	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](#)

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	GAL	C	1	2	12,12,12	0.57	0	17,17,17	0.88	1 (5%)
2	AAL	C	2	2	11,11,12	0.55	0	13,16,18	1.46	2 (15%)
3	AAL	D	1	3	12,12,12	0.76	0	16,18,18	1.13	2 (12%)
3	GAL	D	2	3	11,11,12	0.50	0	15,15,17	0.91	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	GAL	C	1	2	-	1/2/22/22	0/1/1/1
2	AAL	C	2	2	-	-	0/3/2/2
3	AAL	D	1	3	-	-	0/3/2/2
3	GAL	D	2	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	AAL	C1-O5-C5	4.07	117.65	112.19
2	C	1	GAL	O3-C3-C2	-2.32	104.90	110.38
3	D	1	AAL	C3-C4-C5	-2.22	97.22	101.99
3	D	2	GAL	C1-O5-C5	2.19	115.12	112.19
2	C	2	AAL	O5-C5-C6	-2.03	110.36	113.33
3	D	1	AAL	O5-C5-C4	-2.02	106.49	109.51

There are no chirality outliers.

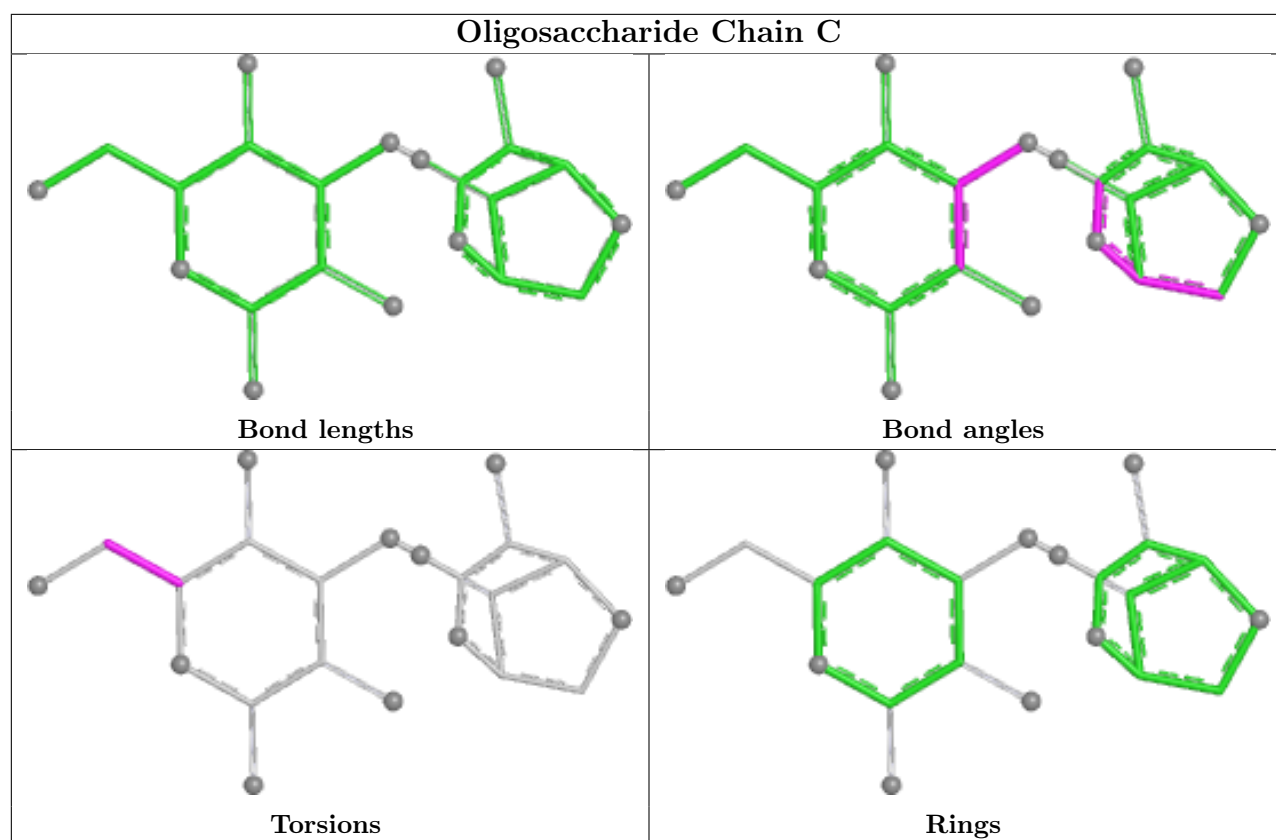
All (1) torsion outliers are listed below:

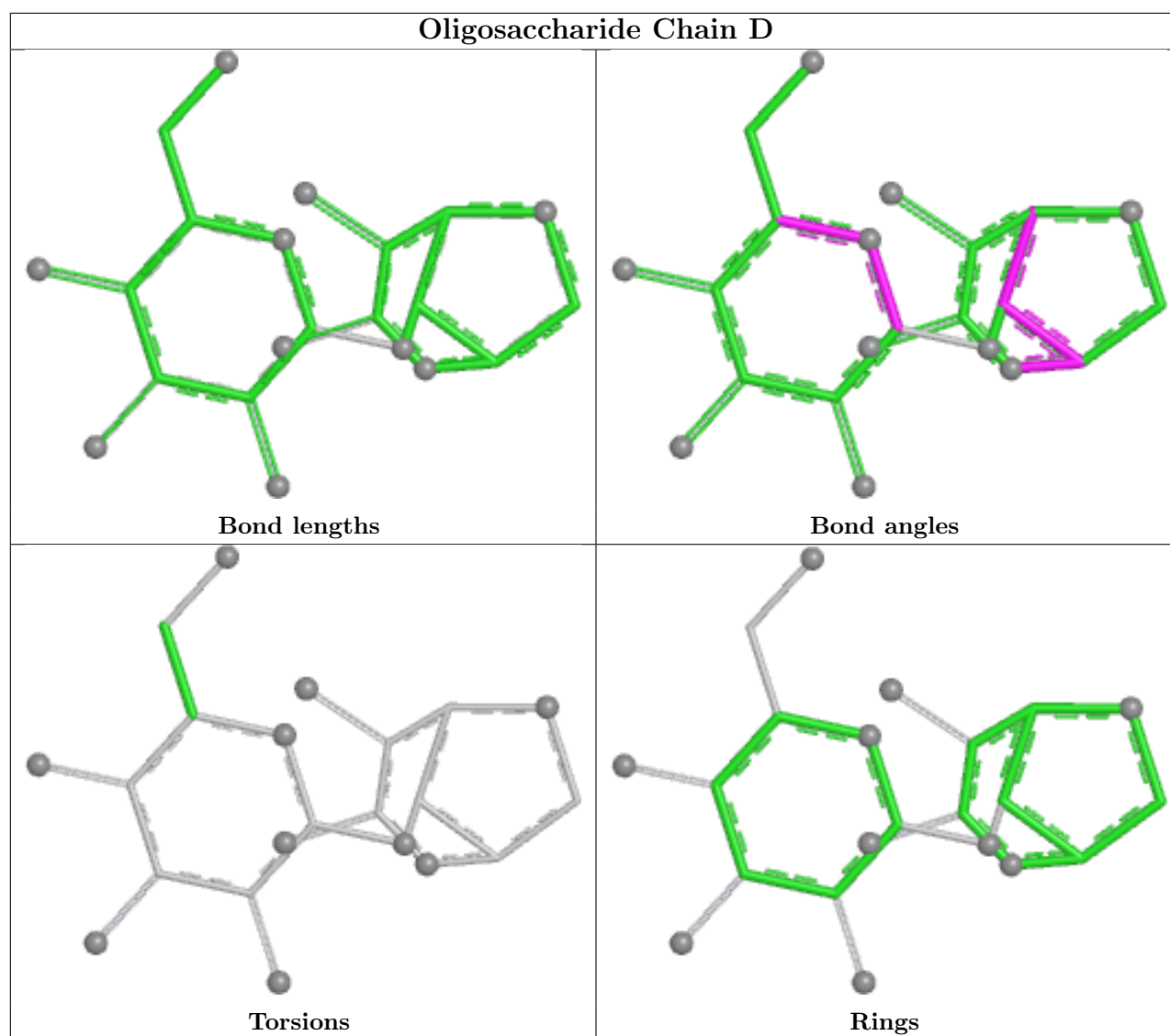
Mol	Chain	Res	Type	Atoms
2	C	1	GAL	O5-C5-C6-O6

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 53 ligands modelled in this entry, 6 are monoatomic - leaving 47 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	708	-	3,3,3	0.30	0	2,2,2	0.55	0
5	EDO	A	725	-	3,3,3	0.43	0	2,2,2	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	B	707	-	3,3,3	0.29	0	2,2,2	0.32	0
5	EDO	A	722	-	3,3,3	0.32	0	2,2,2	0.78	0
5	EDO	A	706	-	3,3,3	0.41	0	2,2,2	0.70	0
5	EDO	A	712	-	3,3,3	0.30	0	2,2,2	0.28	0
5	EDO	A	707	-	3,3,3	0.32	0	2,2,2	0.36	0
5	EDO	A	720	-	3,3,3	0.42	0	2,2,2	0.21	0
5	EDO	A	724	-	3,3,3	0.30	0	2,2,2	0.33	0
5	EDO	B	718	-	3,3,3	0.23	0	2,2,2	0.35	0
5	EDO	B	721	-	3,3,3	0.46	0	2,2,2	0.29	0
4	SO4	B	703	-	4,4,4	0.53	0	6,6,6	0.27	0
4	SO4	B	704	-	4,4,4	0.42	0	6,6,6	0.19	0
5	EDO	A	716	-	3,3,3	0.42	0	2,2,2	0.29	0
5	EDO	B	712	-	3,3,3	0.34	0	2,2,2	0.62	0
5	EDO	B	716	-	3,3,3	0.42	0	2,2,2	0.29	0
5	EDO	A	718	-	3,3,3	0.36	0	2,2,2	0.57	0
5	EDO	B	714	-	3,3,3	0.37	0	2,2,2	0.48	0
5	EDO	B	726	-	3,3,3	0.34	0	2,2,2	0.45	0
5	EDO	B	725	-	3,3,3	0.28	0	2,2,2	0.52	0
5	EDO	B	724	-	3,3,3	0.45	0	2,2,2	0.42	0
5	EDO	A	719	-	3,3,3	0.42	0	2,2,2	0.25	0
5	EDO	B	708	-	3,3,3	0.52	0	2,2,2	0.19	0
5	EDO	B	719	-	3,3,3	0.44	0	2,2,2	0.42	0
5	EDO	B	723	-	3,3,3	0.26	0	2,2,2	0.74	0
5	EDO	A	715	-	3,3,3	0.59	0	2,2,2	0.35	0
5	EDO	A	710	-	3,3,3	0.48	0	2,2,2	0.34	0
5	EDO	A	711	-	3,3,3	0.37	0	2,2,2	0.47	0
5	EDO	A	717	-	3,3,3	0.28	0	2,2,2	0.83	0
5	EDO	A	709	-	3,3,3	0.22	0	2,2,2	0.67	0
5	EDO	B	717	-	3,3,3	0.53	0	2,2,2	0.33	0
4	SO4	B	702	-	4,4,4	0.32	0	6,6,6	0.22	0
5	EDO	B	710	-	3,3,3	0.42	0	2,2,2	0.62	0
5	EDO	B	711	-	3,3,3	0.43	0	2,2,2	0.20	0
4	SO4	B	701	-	4,4,4	0.37	0	6,6,6	0.42	0
4	SO4	A	705	-	4,4,4	0.50	0	6,6,6	0.26	0
5	EDO	B	713	-	3,3,3	0.50	0	2,2,2	0.58	0
5	EDO	A	721	-	3,3,3	0.42	0	2,2,2	0.58	0
5	EDO	A	714	-	3,3,3	0.41	0	2,2,2	0.46	0
5	EDO	B	722	-	3,3,3	0.43	0	2,2,2	0.52	0
5	EDO	B	706	-	3,3,3	0.63	0	2,2,2	0.55	0
5	EDO	B	720	-	3,3,3	0.44	0	2,2,2	0.21	0
5	EDO	A	723	-	3,3,3	0.36	0	2,2,2	0.57	0
5	EDO	B	709	-	3,3,3	0.44	0	2,2,2	0.41	0
5	EDO	B	705	-	3,3,3	0.43	0	2,2,2	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	EDO	A	713	-	3,3,3	0.57	0	2,2,2	0.27	0
5	EDO	B	715	-	3,3,3	0.48	0	2,2,2	0.37	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	708	-	-	1/1/1/1	-
5	EDO	A	725	-	-	1/1/1/1	-
5	EDO	B	707	-	-	0/1/1/1	-
5	EDO	A	722	-	-	0/1/1/1	-
5	EDO	A	706	-	-	1/1/1/1	-
5	EDO	A	712	-	-	1/1/1/1	-
5	EDO	A	707	-	-	0/1/1/1	-
5	EDO	A	720	-	-	1/1/1/1	-
5	EDO	A	724	-	-	1/1/1/1	-
5	EDO	B	718	-	-	1/1/1/1	-
5	EDO	B	721	-	-	1/1/1/1	-
5	EDO	A	716	-	-	1/1/1/1	-
5	EDO	B	712	-	-	0/1/1/1	-
5	EDO	B	716	-	-	0/1/1/1	-
5	EDO	A	718	-	-	1/1/1/1	-
5	EDO	B	714	-	-	0/1/1/1	-
5	EDO	B	726	-	-	0/1/1/1	-
5	EDO	B	725	-	-	1/1/1/1	-
5	EDO	B	724	-	-	1/1/1/1	-
5	EDO	A	719	-	-	0/1/1/1	-
5	EDO	B	708	-	-	0/1/1/1	-
5	EDO	B	719	-	-	0/1/1/1	-
5	EDO	B	723	-	-	1/1/1/1	-
5	EDO	A	715	-	-	0/1/1/1	-
5	EDO	A	710	-	-	0/1/1/1	-
5	EDO	A	711	-	-	0/1/1/1	-
5	EDO	A	717	-	-	0/1/1/1	-
5	EDO	A	709	-	-	0/1/1/1	-
5	EDO	B	717	-	-	1/1/1/1	-
5	EDO	B	710	-	-	0/1/1/1	-
5	EDO	B	711	-	-	1/1/1/1	-
5	EDO	B	713	-	-	0/1/1/1	-
5	EDO	A	721	-	-	1/1/1/1	-
5	EDO	A	714	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	B	722	-	-	0/1/1/1	-
5	EDO	B	706	-	-	0/1/1/1	-
5	EDO	A	713	-	-	0/1/1/1	-
5	EDO	B	720	-	-	0/1/1/1	-
5	EDO	B	709	-	-	1/1/1/1	-
5	EDO	B	705	-	-	0/1/1/1	-
5	EDO	A	723	-	-	0/1/1/1	-
5	EDO	B	715	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	712	EDO	O1-C1-C2-O2
5	B	718	EDO	O1-C1-C2-O2
5	A	721	EDO	O1-C1-C2-O2
5	A	725	EDO	O1-C1-C2-O2
5	B	709	EDO	O1-C1-C2-O2
5	A	708	EDO	O1-C1-C2-O2
5	B	717	EDO	O1-C1-C2-O2
5	A	706	EDO	O1-C1-C2-O2
5	A	718	EDO	O1-C1-C2-O2
5	A	720	EDO	O1-C1-C2-O2
5	B	711	EDO	O1-C1-C2-O2
5	B	723	EDO	O1-C1-C2-O2
5	B	725	EDO	O1-C1-C2-O2
5	A	724	EDO	O1-C1-C2-O2
5	B	715	EDO	O1-C1-C2-O2
5	B	721	EDO	O1-C1-C2-O2
5	B	724	EDO	O1-C1-C2-O2
5	A	716	EDO	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	723	EDO	3	0
5	A	709	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	702	SO4	1	0
5	A	714	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	624/649 (96%)	-0.35	6 (0%) 79 84	3, 10, 20, 34	32 (5%)
1	B	613/649 (94%)	-0.18	25 (4%) 41 45	3, 10, 28, 53	36 (5%)
All	All	1237/1298 (95%)	-0.27	31 (2%) 58 64	3, 10, 23, 53	68 (5%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	398	PHE	5.0
1	B	384	HIS	3.8
1	A	426[A]	CYS	3.8
1	B	643	ILE	3.7
1	B	331	ASN	3.5
1	A	643	ILE	3.4
1	B	56	TYR	3.4
1	B	464	SER	3.3
1	B	385	THR	3.2
1	B	397	SER	3.1
1	B	586	ASP	3.0
1	B	395	PRO	3.0
1	B	396	GLU	2.9
1	B	641	ASN	2.8
1	B	330	TYR	2.8
1	B	328	HIS	2.8
1	B	21	GLN	2.6
1	B	372	SER	2.6
1	A	641	ASN	2.6
1	B	399	THR	2.5
1	B	393	ASN	2.5
1	B	463	THR	2.4
1	A	425	ASN	2.4
1	B	334	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	400	GLY	2.3
1	A	642	ASN	2.2
1	B	645	GLY	2.2
1	B	326	GLN	2.2
1	B	642	ASN	2.1
1	B	23[A]	HIS	2.1
1	A	586	ASP	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](#)

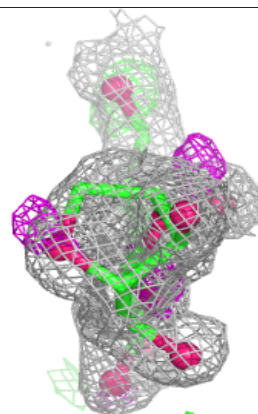
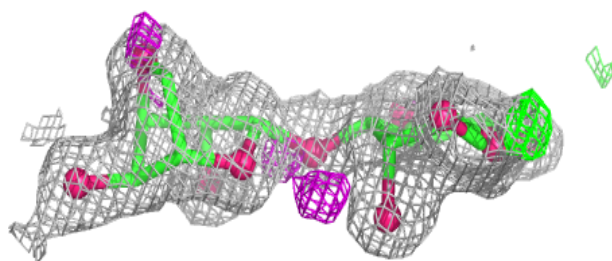
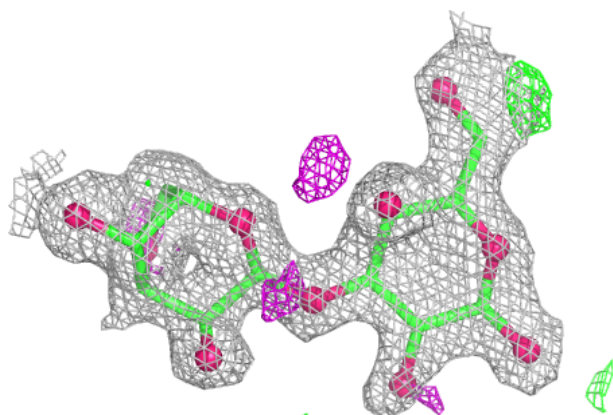
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	AAL	D	1	11/11	0.74	0.17	24,28,34,35	0
3	GAL	D	2	11/12	0.84	0.14	24,26,29,34	0
2	GAL	C	1	12/12	0.87	0.10	22,27,30,31	0
2	AAL	C	2	10/11	0.91	0.09	21,22,23,23	0

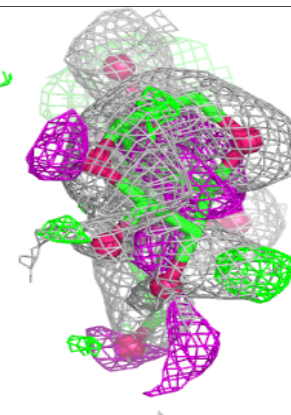
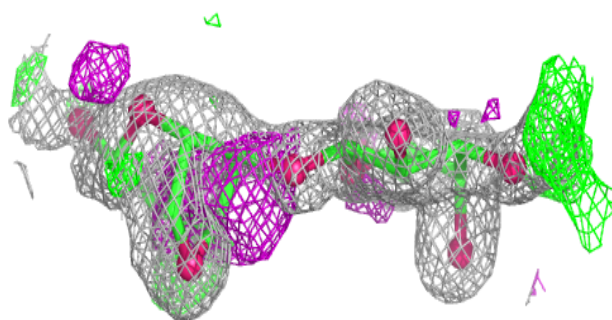
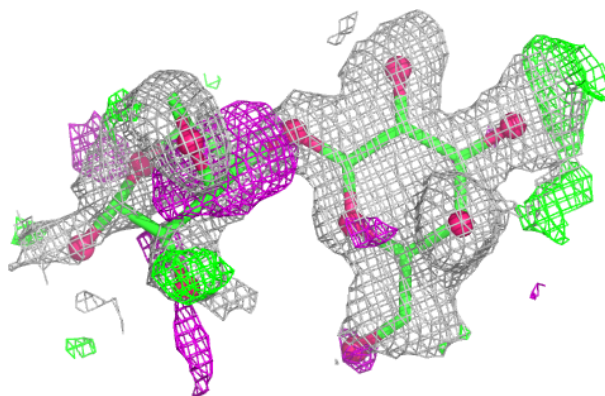
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

**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)



6.4 Ligands

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
5	EDO	B	723	4/4	0.67	0.26	38,41,43,46	0
5	EDO	A	722	4/4	0.69	0.28	45,46,48,50	0
5	EDO	A	723	4/4	0.75	0.22	38,39,40,41	0
5	EDO	A	718	4/4	0.77	0.23	37,38,38,42	0
5	EDO	B	717	4/4	0.78	0.19	30,32,32,32	0
5	EDO	B	709	4/4	0.78	0.21	29,30,32,35	0
5	EDO	A	721	4/4	0.80	0.31	41,45,46,47	0
5	EDO	B	724	4/4	0.80	0.21	41,43,45,46	0
5	EDO	B	726	4/4	0.80	0.21	35,39,39,41	0
5	EDO	B	712	4/4	0.81	0.21	44,45,46,48	0
5	EDO	B	714	4/4	0.81	0.17	34,36,37,40	0
5	EDO	A	725	4/4	0.81	0.17	40,41,41,42	0
5	EDO	B	710	4/4	0.83	0.15	29,30,31,32	0
5	EDO	B	716	4/4	0.83	0.15	33,33,34,35	0
5	EDO	B	721	4/4	0.84	0.18	30,31,32,33	0
5	EDO	B	715	4/4	0.84	0.15	30,30,30,34	0
5	EDO	A	720	4/4	0.84	0.14	27,31,32,33	0
5	EDO	B	706	4/4	0.84	0.14	18,19,19,20	0
5	EDO	A	714	4/4	0.85	0.16	22,25,27,30	0
5	EDO	B	722	4/4	0.86	0.17	22,22,24,24	0
5	EDO	A	715	4/4	0.86	0.18	27,27,28,29	0
5	EDO	B	719	4/4	0.86	0.13	22,23,23,26	0
5	EDO	B	725	4/4	0.86	0.17	32,35,36,37	0
5	EDO	A	710	4/4	0.86	0.14	27,27,29,29	0
5	EDO	A	708	4/4	0.87	0.15	20,24,26,32	0
5	EDO	A	724	4/4	0.87	0.16	23,27,30,32	0
5	EDO	B	713	4/4	0.88	0.12	21,21,22,24	0
5	EDO	A	706	4/4	0.88	0.15	18,20,21,22	0
5	EDO	A	719	4/4	0.88	0.13	25,28,29,30	0
5	EDO	A	712	4/4	0.88	0.16	26,28,30,33	0
5	EDO	B	718	4/4	0.89	0.16	25,25,28,30	0
5	EDO	A	717	4/4	0.89	0.11	18,21,21,23	0
4	SO4	B	703	5/5	0.89	0.11	11,12,13,13	5
5	EDO	A	709	4/4	0.89	0.15	26,28,30,34	0
5	EDO	A	711	4/4	0.90	0.11	22,22,23,24	0
4	SO4	A	705	5/5	0.90	0.17	31,31,31,31	0
5	EDO	B	720	4/4	0.90	0.12	19,22,25,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	EDO	B	711	4/4	0.91	0.12	15,19,22,25	0
5	EDO	A	716	4/4	0.92	0.11	26,27,28,28	0
5	EDO	A	713	4/4	0.93	0.11	19,19,19,20	0
4	SO4	B	704	5/5	0.93	0.09	38,38,44,44	0
5	EDO	B	708	4/4	0.93	0.11	19,19,19,20	0
5	EDO	B	707	4/4	0.94	0.08	17,19,19,21	0
5	EDO	A	707	4/4	0.95	0.07	17,18,19,20	0
5	EDO	B	705	4/4	0.97	0.06	12,12,12,12	0
4	SO4	B	702	5/5	0.97	0.08	27,30,32,32	0
4	SO4	B	701	5/5	0.98	0.08	18,20,21,22	0
6	CA	A	726	1/1	0.99	0.08	17,17,17,17	0
6	CA	B	728	1/1	0.99	0.10	18,18,18,18	0
6	CA	A	728	1/1	1.00	0.03	8,8,8,8	0
6	CA	B	727	1/1	1.00	0.05	10,10,10,10	0
6	CA	A	727	1/1	1.00	0.03	9,9,9,9	0
6	CA	B	729	1/1	1.00	0.03	10,10,10,10	0

6.5 Other polymers [i](#)

There are no such residues in this entry.