



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 05:25 PM UTC

PDB ID : 6UB3 / pdb_00006ub3
Title : Crystal structure of a GH128 (subgroup IV) endo-beta-1,3-glucanase from Lentinula edodes (LeGH128_IV) with laminaribiose at the surface-binding site
Authors : Santos, C.R.; Lima, E.A.; Mandelli, F.; Murakami, M.T.
Deposited on : 2019-09-11
Resolution : 1.85 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

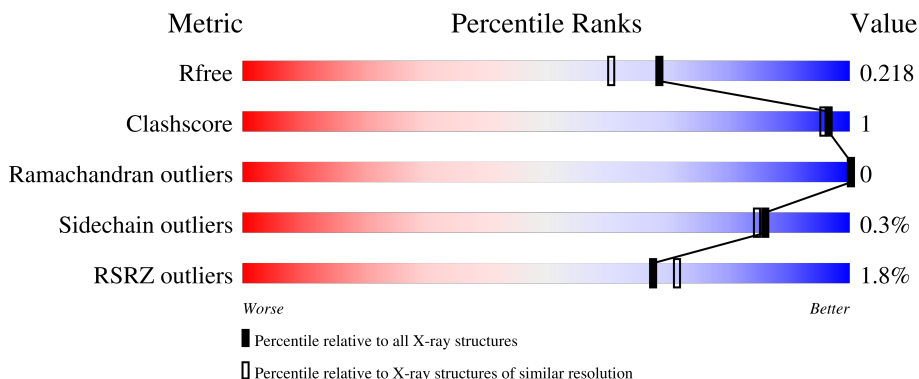
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

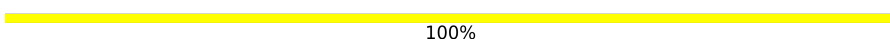
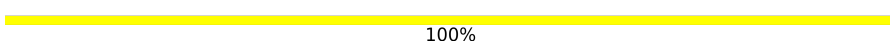
The reported resolution of this entry is 1.85 Å.

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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	270	
1	B	270	
2	C	2	
2	D	2	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4338 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	248	Total	C	N	O	S	0	1	0
			1874	1205	294	368	7			
1	B	248	Total	C	N	O	S	0	0	0
			1868	1202	293	366	7			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	initiating methionine	UNP G9M5R4
A	-1	GLY	-	expression tag	UNP G9M5R4
A	0	SER	-	expression tag	UNP G9M5R4
A	1	SER	-	expression tag	UNP G9M5R4
A	2	HIS	-	expression tag	UNP G9M5R4
A	3	HIS	-	expression tag	UNP G9M5R4
A	4	HIS	-	expression tag	UNP G9M5R4
A	5	HIS	-	expression tag	UNP G9M5R4
A	6	HIS	-	expression tag	UNP G9M5R4
A	7	HIS	-	expression tag	UNP G9M5R4
A	8	SER	-	expression tag	UNP G9M5R4
A	9	SER	-	expression tag	UNP G9M5R4
A	10	GLY	-	expression tag	UNP G9M5R4
A	11	LEU	-	expression tag	UNP G9M5R4
A	12	VAL	-	expression tag	UNP G9M5R4
A	13	PRO	-	expression tag	UNP G9M5R4
A	14	ARG	-	expression tag	UNP G9M5R4
A	15	GLY	-	expression tag	UNP G9M5R4
A	16	SER	-	expression tag	UNP G9M5R4
A	17	HIS	-	expression tag	UNP G9M5R4
A	18	MET	-	expression tag	UNP G9M5R4
A	19	ALA	-	expression tag	UNP G9M5R4
A	20	SER	-	expression tag	UNP G9M5R4
B	-2	MET	-	initiating methionine	UNP G9M5R4
B	-1	GLY	-	expression tag	UNP G9M5R4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	SER	-	expression tag	UNP G9M5R4
B	1	SER	-	expression tag	UNP G9M5R4
B	2	HIS	-	expression tag	UNP G9M5R4
B	3	HIS	-	expression tag	UNP G9M5R4
B	4	HIS	-	expression tag	UNP G9M5R4
B	5	HIS	-	expression tag	UNP G9M5R4
B	6	HIS	-	expression tag	UNP G9M5R4
B	7	HIS	-	expression tag	UNP G9M5R4
B	8	SER	-	expression tag	UNP G9M5R4
B	9	SER	-	expression tag	UNP G9M5R4
B	10	GLY	-	expression tag	UNP G9M5R4
B	11	LEU	-	expression tag	UNP G9M5R4
B	12	VAL	-	expression tag	UNP G9M5R4
B	13	PRO	-	expression tag	UNP G9M5R4
B	14	ARG	-	expression tag	UNP G9M5R4
B	15	GLY	-	expression tag	UNP G9M5R4
B	16	SER	-	expression tag	UNP G9M5R4
B	17	HIS	-	expression tag	UNP G9M5R4
B	18	MET	-	expression tag	UNP G9M5R4
B	19	ALA	-	expression tag	UNP G9M5R4
B	20	SER	-	expression tag	UNP G9M5R4

- Molecule 2 is an oligosaccharide called beta-D-glucopyranose-(1-3)-beta-D-glucopyranose.

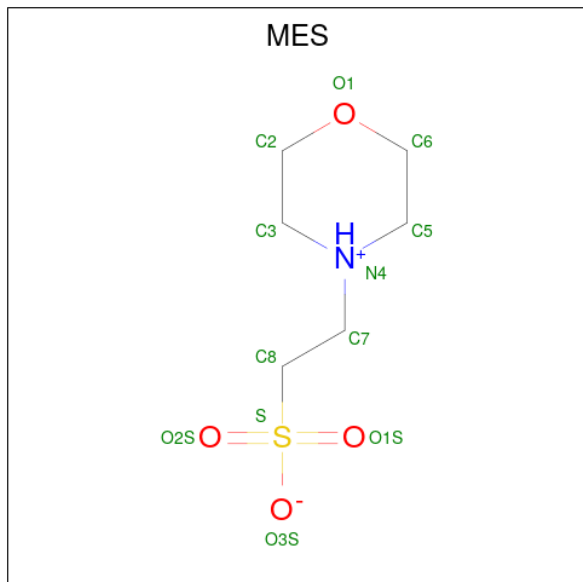


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

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

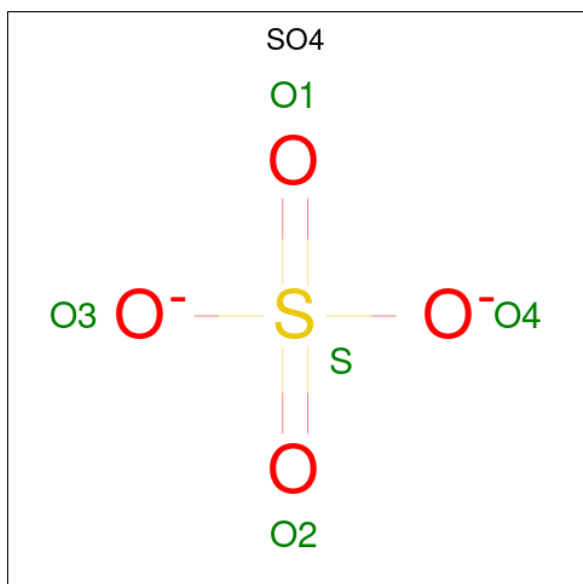
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

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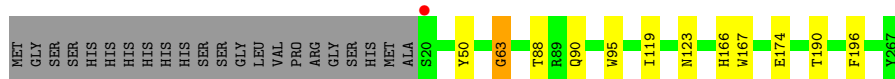
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	327	Total	O	0	0
			327	327		
6	B	199	Total	O	0	0
			199	199		

- Molecule 1: glycoside hydrolase



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- Sequence logo for position 257. The y-axis represents information content (bits) from 0 to 1. The x-axis shows amino acid positions from MET to G257. The logo indicates a high frequency of MET at the start, followed by a peak of HIS and SER around position 10. A large peak of MET is at position 257.

- BGC1
BGC2

- BGC1
BGC2

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	157.02Å 46.69Å 115.25Å 90.00° 119.70° 90.00°	Depositor
Resolution (Å)	44.18 – 1.85 44.18 – 1.85	Depositor EDS
% Data completeness (in resolution range)	96.4 (44.18-1.85) 96.4 (44.18-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 1.86Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.174 , 0.209 0.185 , 0.218	Depositor DCC
R_{free} test set	3010 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	24.1	Xtriage
Anisotropy	0.737	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4338	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.01% 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: MES, BGC, SO4, 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	1.19	6/1933 (0.3%)	1.29	2/2651 (0.1%)
1	B	1.29	7/1927 (0.4%)	1.37	4/2643 (0.2%)
All	All	1.24	13/3860 (0.3%)	1.33	6/5294 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	63	GLY	C-O	7.25	1.30	1.23
1	B	254	THR	C-O	6.98	1.32	1.23
1	B	257	GLY	C-O	6.04	1.32	1.24
1	A	166	HIS	CE1-NE2	5.86	1.38	1.32
1	A	174	GLU	CD-OE1	5.59	1.35	1.25

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	104	PRO	CA-C-N	5.38	128.03	120.28
1	B	104	PRO	C-N-CA	5.38	128.03	120.28
1	B	38	VAL	CA-C-O	-5.12	115.06	120.64
1	A	50	TYR	CA-C-O	-5.05	115.69	121.40
1	B	105	ASP	CA-CB-CG	5.04	117.64	112.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	257	GLY	Mainchain

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	1874	0	1750	3	0
1	B	1868	0	1746	4	0
2	C	23	0	21	0	0
2	D	23	0	21	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	12	0	13	0	0
5	A	10	0	0	1	0
6	A	327	0	0	0	0
6	B	199	0	0	0	0
All	All	4338	0	3551	7	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 1.

The worst 5 of 7 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:32:SER:N	1:B:33:PRO:CD	2.66	0.59
1:A:63:GLY:HA3	5:A:305:SO4:O4	2.05	0.56
1:B:167:TRP:O	1:B:196:PHE:HA	2.07	0.55
1:B:22:LYS:O	1:B:223:LEU:HA	2.13	0.48
1:A:167:TRP:O	1:A:196:PHE:HA	2.16	0.46

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	247/270 (92%)	242 (98%)	5 (2%)	0	100	100
1	B	246/270 (91%)	239 (97%)	7 (3%)	0	100	100
All	All	493/540 (91%)	481 (98%)	12 (2%)	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	195/212 (92%)	195 (100%)	0	100	100
1	B	194/212 (92%)	193 (100%)	1 (0%)	81	77
All	All	389/424 (92%)	388 (100%)	1 (0%)	86	85

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	20	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	92	GLN

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Mol	Chain	Res	Type
1	B	73	GLN
1	B	92	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	0.81	0	17,17,17	2.82	6 (35%)
2	BGC	C	2	2	11,11,12	0.98	0	15,15,17	1.73	3 (20%)
2	BGC	D	1	2	12,12,12	0.77	1 (8%)	17,17,17	1.59	6 (35%)
2	BGC	D	2	2	11,11,12	0.89	0	15,15,17	1.20	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	-	0/2/22/22	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	-	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	D	1	BGC	C4-C5	2.40	1.58	1.53

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	BGC	O5-C1-C2	-6.36	99.12	110.30
2	C	1	BGC	C1-O5-C5	-4.45	105.05	113.65
2	C	1	BGC	C6-C5-C4	4.21	123.35	113.02
2	C	1	BGC	O1-C1-C2	4.15	121.04	108.98
2	C	1	BGC	C1-C2-C3	-3.96	102.28	110.36

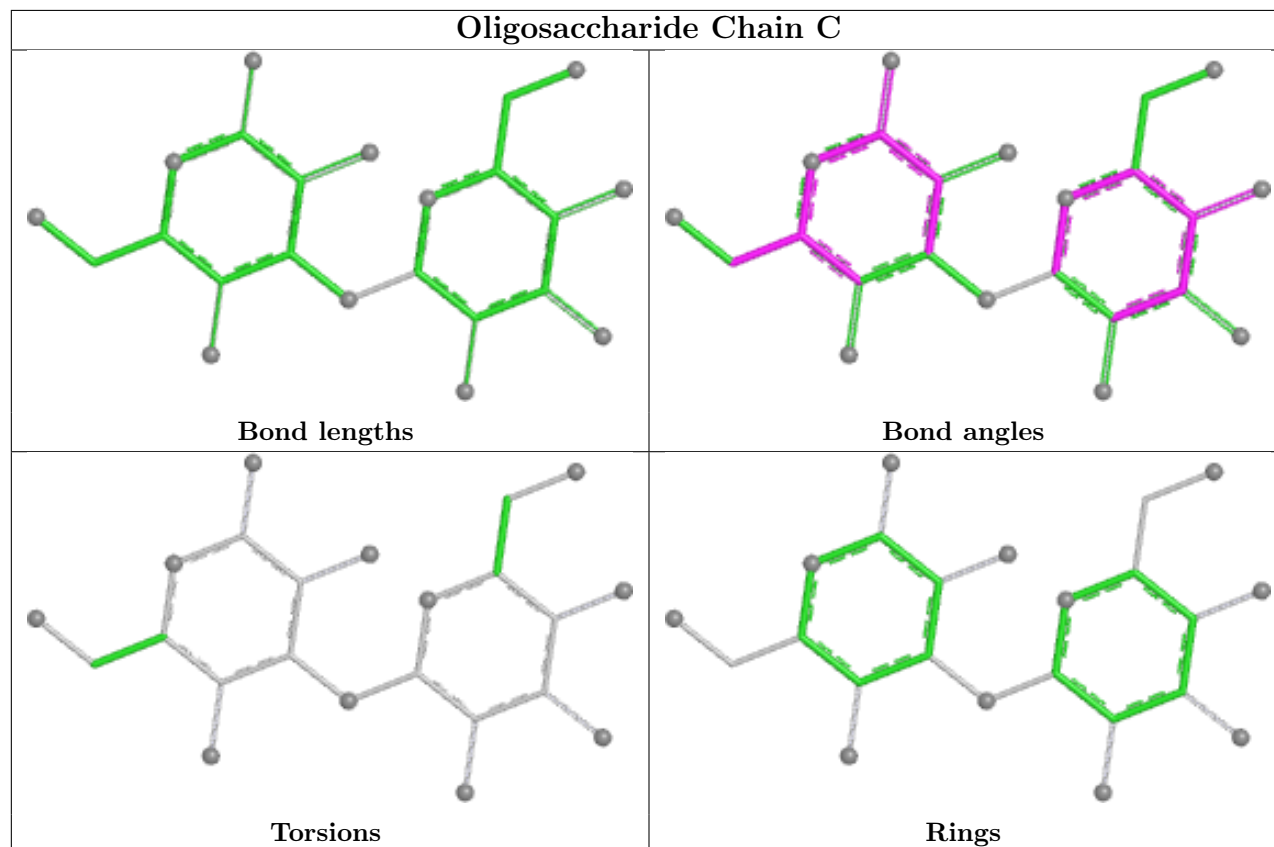
There are no chirality outliers.

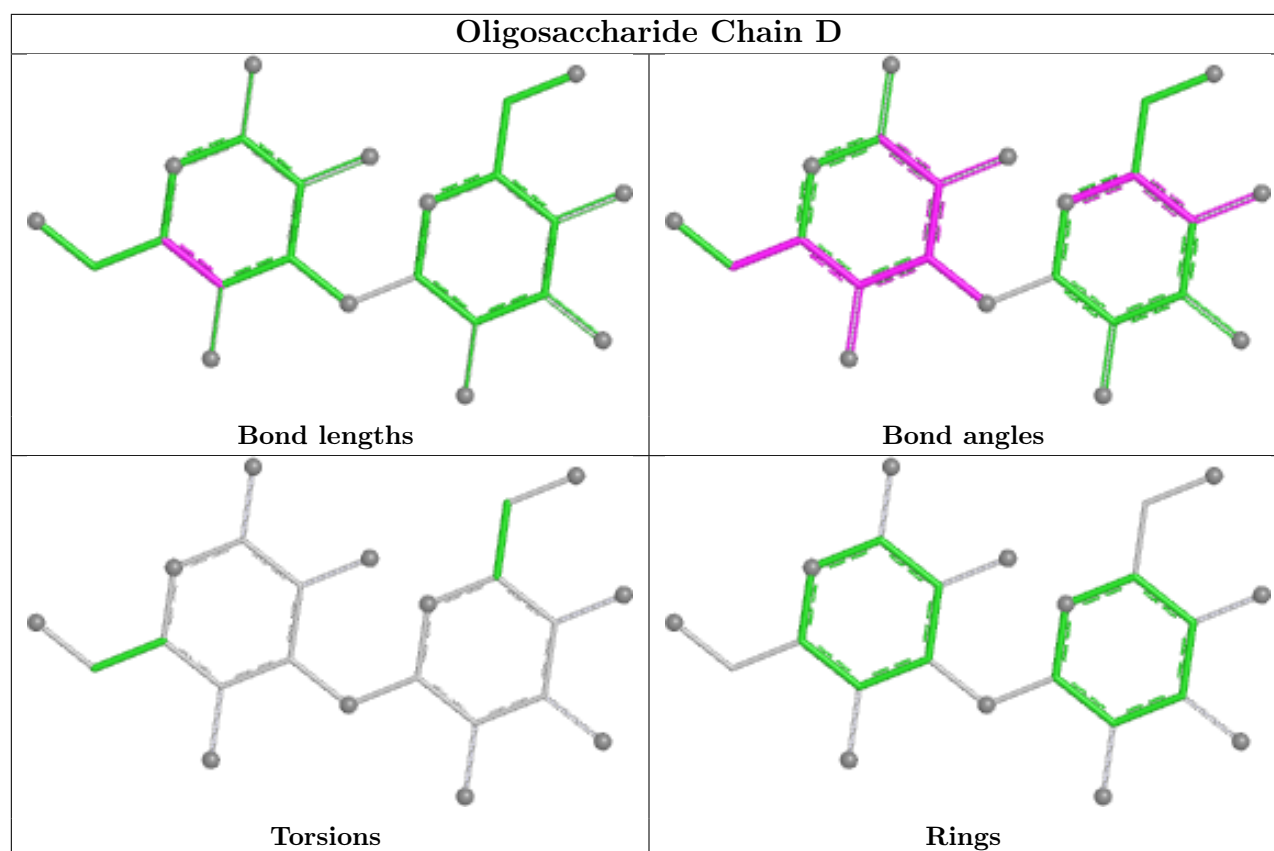
There are no torsion outliers.

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 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	SO4	A	305	-	4,4,4	0.30	0	6,6,6	0.31	0
5	SO4	A	306	-	4,4,4	0.66	0	6,6,6	0.33	0
4	MES	A	304	-	12,12,12	0.85	0	15,16,16	0.62	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	MES	A	304	-	-	0/6/14/14	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	305	SO4	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	248/270 (91%)	-0.28	1 (0%) 88 91	14, 28, 36, 57	1 (0%)
1	B	248/270 (91%)	0.33	8 (3%) 50 54	26, 39, 60, 71	0
All	All	496/540 (91%)	0.03	9 (1%) 67 71	14, 32, 57, 71	1 (0%)

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	20	SER	3.1
1	B	62	GLY	2.6
1	B	20	SER	2.5
1	B	257	GLY	2.4
1	B	58	PRO	2.3

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

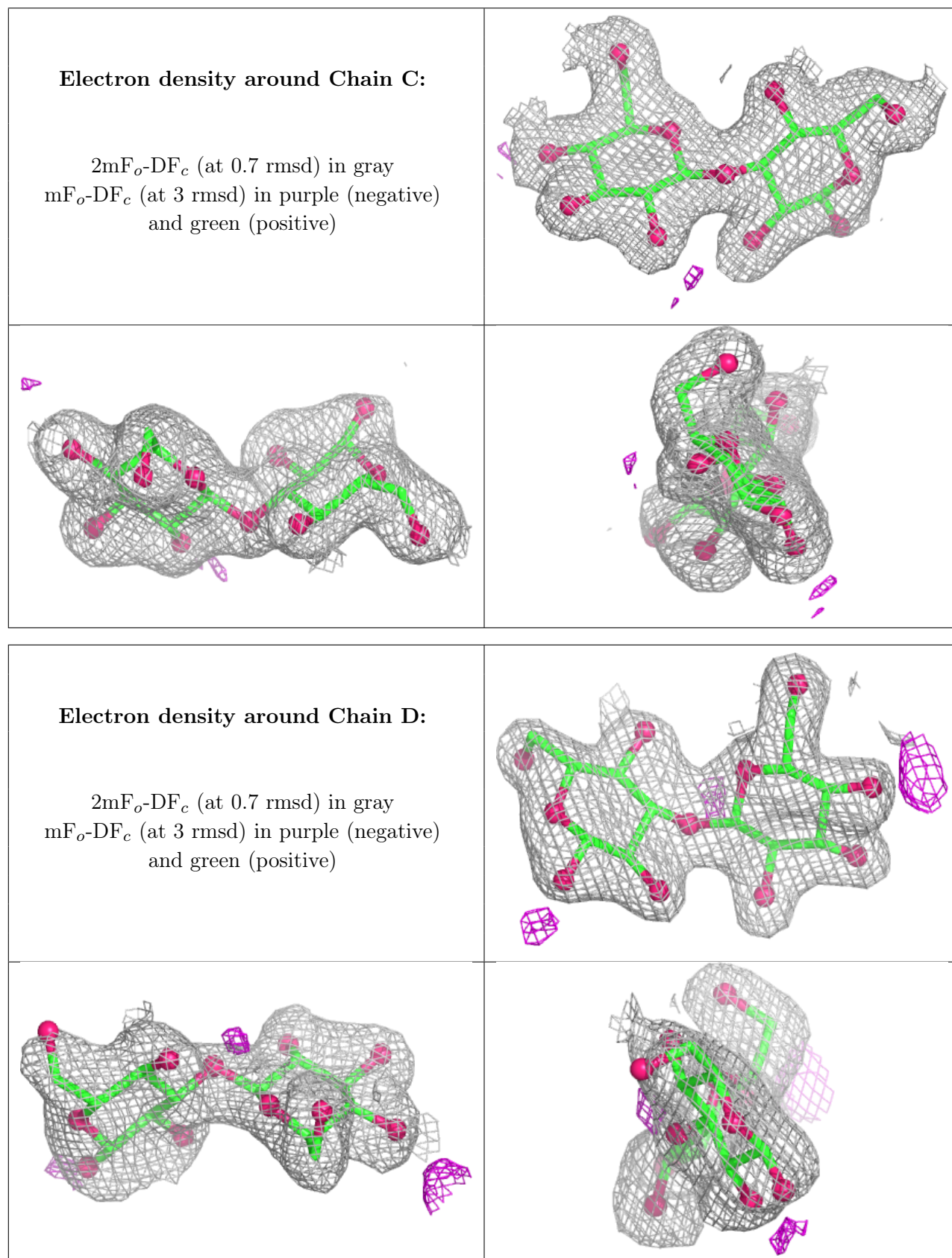
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BGC	D	1	12/12	0.89	0.10	51,58,60,64	0
2	BGC	D	2	11/12	0.92	0.08	43,47,49,50	0
2	BGC	C	1	12/12	0.93	0.08	38,43,48,58	0
2	BGC	C	2	11/12	0.98	0.04	31,35,37,39	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
5	SO4	A	306	5/5	0.79	0.14	42,58,66,72	0
5	SO4	A	305	5/5	0.84	0.14	68,78,84,90	0
4	MES	A	304	12/12	0.85	0.19	58,65,69,71	0
3	CL	B	303	1/1	0.99	0.12	42,42,42,42	0
3	CL	A	303	1/1	0.99	0.07	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.