



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 03:02 PM BST

PDB ID : 5E1Q
Title : Mutant (D415G) GH97 alpha-galactosidase in complex with Gal-Lac
Authors : Matsunaga, K.; Yamashita, K.; Tagami, T.; Yao, M.; Okuyama, M.; Kimura, A.
Deposited on : 2015-09-30
Resolution : 1.94 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

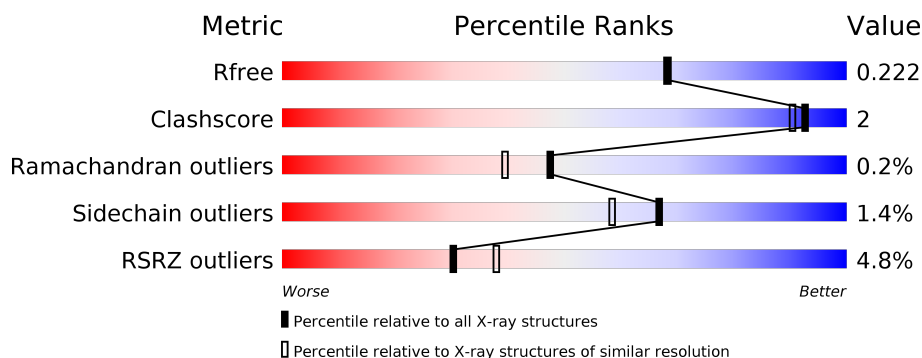
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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}	130704	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

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 on 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	655	<div> <div>7%</div> <div> <div></div> <div>93%</div> <div>5%</div> </div> </div>
1	B	655	<div> <div>3%</div> <div> <div></div> <div>93%</div> <div>5%</div> </div> </div>
2	C	3	<div> <div>33%</div> <div>67%</div> </div>
2	D	3	<div> <div>67%</div> <div>33%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11104 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 Retaining alpha-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	642	Total	C	N	O	S	0	1	0
			5120	3253	869	966	32			
1	B	644	Total	C	N	O	S	0	1	0
			5133	3262	871	968	32			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	MET	-	expression tag	UNP Q8A6L0
A	9	ASN	-	expression tag	UNP Q8A6L0
A	10	HIS	-	expression tag	UNP Q8A6L0
A	11	LYS	-	expression tag	UNP Q8A6L0
A	12	VAL	-	expression tag	UNP Q8A6L0
A	13	HIS	-	expression tag	UNP Q8A6L0
A	14	HIS	-	expression tag	UNP Q8A6L0
A	15	HIS	-	expression tag	UNP Q8A6L0
A	16	HIS	-	expression tag	UNP Q8A6L0
A	17	HIS	-	expression tag	UNP Q8A6L0
A	18	HIS	-	expression tag	UNP Q8A6L0
A	19	ILE	-	expression tag	UNP Q8A6L0
A	20	GLU	-	expression tag	UNP Q8A6L0
A	21	GLY	-	expression tag	UNP Q8A6L0
A	22	ARG	-	expression tag	UNP Q8A6L0
A	23	HIS	-	expression tag	UNP Q8A6L0
A	24	MET	-	expression tag	UNP Q8A6L0
A	25	GLU	-	expression tag	UNP Q8A6L0
A	26	LEU	-	expression tag	UNP Q8A6L0
A	415	GLY	ASP	engineered mutation	UNP Q8A6L0
B	8	MET	-	expression tag	UNP Q8A6L0
B	9	ASN	-	expression tag	UNP Q8A6L0
B	10	HIS	-	expression tag	UNP Q8A6L0
B	11	LYS	-	expression tag	UNP Q8A6L0
B	12	VAL	-	expression tag	UNP Q8A6L0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	13	HIS	-	expression tag	UNP Q8A6L0
B	14	HIS	-	expression tag	UNP Q8A6L0
B	15	HIS	-	expression tag	UNP Q8A6L0
B	16	HIS	-	expression tag	UNP Q8A6L0
B	17	HIS	-	expression tag	UNP Q8A6L0
B	18	HIS	-	expression tag	UNP Q8A6L0
B	19	ILE	-	expression tag	UNP Q8A6L0
B	20	GLU	-	expression tag	UNP Q8A6L0
B	21	GLY	-	expression tag	UNP Q8A6L0
B	22	ARG	-	expression tag	UNP Q8A6L0
B	23	HIS	-	expression tag	UNP Q8A6L0
B	24	MET	-	expression tag	UNP Q8A6L0
B	25	GLU	-	expression tag	UNP Q8A6L0
B	26	LEU	-	expression tag	UNP Q8A6L0
B	415	GLY	ASP	engineered mutation	UNP Q8A6L0

- Molecule 2 is an oligosaccharide called alpha-D-galactopyranose-(1-1)-[beta-D-galactopyranose-(1-4)]beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	3	Total	C	O	0	0	0
			34	18	16			
2	D	3	Total	C	O	0	0	0
			34	18	16			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

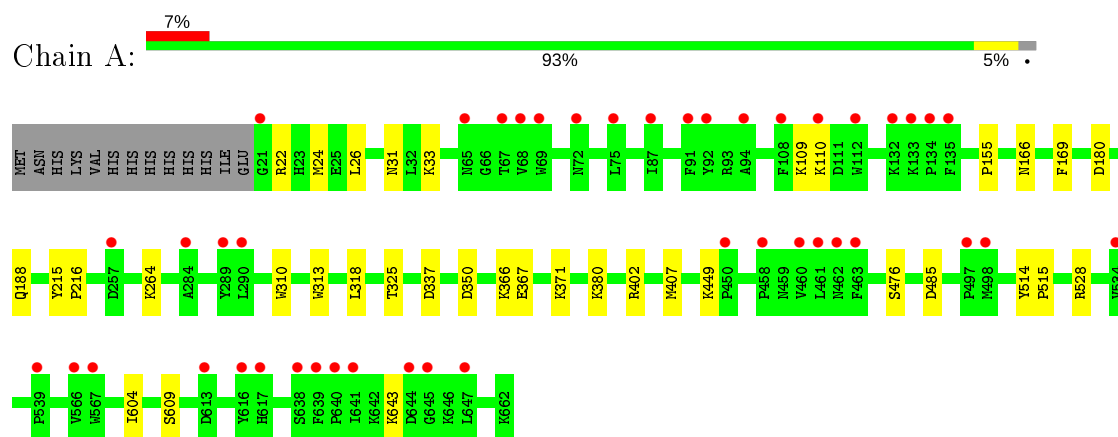
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	401	Total	O	0	0
			401	401		
5	B	362	Total	O	0	0
			362	362		

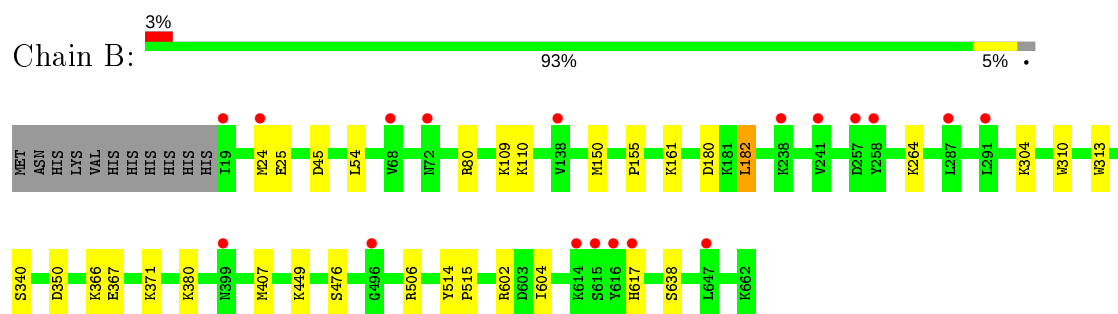
3 Residue-property plots [i](#)

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: Retaining alpha-galactosidase



- Molecule 1: Retaining alpha-galactosidase



- Molecule 2: alpha-D-galactopyranose-(1-1)-[beta-D-galactopyranose-(1-4)]beta-D-glucopyranose



- Molecule 2: alpha-D-galactopyranose-(1-1)-[beta-D-galactopyranose-(1-4)]beta-D-glucopyranose





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.62Å 100.62Å 237.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.44 – 1.94 48.10 – 1.94	Depositor EDS
% Data completeness (in resolution range)	99.2 (42.44-1.94) 99.2 (48.10-1.94)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.99 (at 1.94Å)	Xtriage
Refinement program	PHENIX 1.9_1690	Depositor
R, R_{free}	0.185 , 0.223 0.190 , 0.222	Depositor DCC
R_{free} test set	2140 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.838	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11104	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% 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: GOL, CA, BGC, GAL, GLA

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	0.40	0/5250	0.54	0/7115
1	B	0.40	0/5263	0.53	0/7133
All	All	0.40	0/10513	0.54	0/14248

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5120	0	4968	17	0
1	B	5133	0	4981	14	0
2	C	34	0	28	0	0
2	D	34	0	28	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	6	0	8	0	0
4	B	12	0	16	0	0
5	A	401	0	0	4	0
5	B	362	0	0	0	0
All	All	11104	0	10029	31	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 (31) 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:366:LYS:NZ	1:A:367:GLU:OE2	2.33	0.58
1:A:402:ARG:HD2	5:A:1080:HOH:O	2.04	0.57
1:A:514[A]:TYR:CD2	1:A:515:PRO:HD2	2.41	0.55
1:A:264:LYS:NZ	5:A:803:HOH:O	2.33	0.54
1:A:514[A]:TYR:CG	1:A:515:PRO:HD2	2.45	0.52
1:B:366:LYS:NZ	1:B:367:GLU:OE2	2.24	0.51
1:A:310:TRP:CZ3	1:A:313:TRP:HB2	2.46	0.51
1:A:26:LEU:HG	1:A:33:LYS:HB3	1.94	0.48
1:B:150:MET:O	1:B:182:LEU:HD13	2.13	0.48
1:B:25:GLU:OE1	1:B:80:ARG:NH1	2.47	0.48
1:B:45:ASP:HB2	1:B:54:LEU:O	2.14	0.47
1:A:188:GLN:NE2	5:A:812:HOH:O	2.42	0.47
1:A:109:LYS:O	1:A:110:LYS:HB2	2.15	0.47
1:B:109:LYS:O	1:B:110:LYS:HB2	2.15	0.47
1:A:31:ASN:HB2	5:A:969:HOH:O	2.16	0.46
1:A:485:ASP:HB2	1:A:528:ARG:HH11	1.80	0.46
1:A:609:SER:HA	1:A:643:LYS:HG3	1.97	0.46
1:B:340:SER:HB2	1:B:380:LYS:HB3	1.98	0.44
1:B:514[B]:TYR:CG	1:B:515:PRO:HD2	2.52	0.44
1:B:617:HIS:CE1	1:B:638:SER:HB3	2.53	0.44
1:A:318:LEU:HB2	1:A:325:THR:HG21	2.00	0.44
1:B:506:ARG:NH2	1:B:514[A]:TYR:OH	2.51	0.43
1:B:161:LYS:HD3	1:B:161:LYS:HA	1.87	0.42
1:B:264:LYS:HE3	1:B:264:LYS:HB2	1.66	0.42
1:B:310:TRP:CZ3	1:B:313:TRP:HB2	2.54	0.42
1:A:166:ASN:HA	1:A:169:PHE:CD2	2.55	0.41
1:A:371:LYS:HE2	1:A:407:MET:HG3	2.01	0.41
1:A:215:TYR:CG	1:A:216:PRO:HD2	2.55	0.41
1:B:304:LYS:HD3	1:B:304:LYS:HA	1.86	0.41
1:B:371:LYS:HE2	1:B:407:MET:HG3	2.03	0.41
1:A:337:ASP:OD1	1:A:380:LYS:NZ	2.55	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	641/655 (98%)	620 (97%)	20 (3%)	1 (0%)	47	39
1	B	643/655 (98%)	624 (97%)	18 (3%)	1 (0%)	47	39
All	All	1284/1310 (98%)	1244 (97%)	38 (3%)	2 (0%)	47	39

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	155	PRO
1	A	155	PRO

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	552/564 (98%)	545 (99%)	7 (1%)	69	62
1	B	553/564 (98%)	545 (99%)	8 (1%)	67	58
All	All	1105/1128 (98%)	1090 (99%)	15 (1%)	67	58

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ARG
1	A	24	MET
1	A	180	ASP
1	A	350	ASP

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Mol	Chain	Res	Type
1	A	449	LYS
1	A	476	SER
1	A	604	ILE
1	B	24	MET
1	B	180	ASP
1	B	182	LEU
1	B	350	ASP
1	B	449	LYS
1	B	476	SER
1	B	602	ARG
1	B	604	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

6 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	3,2	12,12,12	0.53	0	17,17,17	0.76	0
2	GLA	C	2	3,2	11,11,12	0.28	0	15,15,17	1.20	1 (6%)
2	GAL	C	3	2	11,11,12	0.29	0	15,15,17	1.09	1 (6%)
2	BGC	D	1	3,2	12,12,12	0.52	0	17,17,17	0.90	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GLA	D	2	3,2	11,11,12	0.28	0	15,15,17	1.46	1 (6%)
2	GAL	D	3	2	11,11,12	0.32	0	15,15,17	0.55	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	3,2	-	0/2/22/22	0/1/1/1
2	GLA	C	2	3,2	-	0/2/19/22	0/1/1/1
2	GAL	C	3	2	-	0/2/19/22	0/1/1/1
2	BGC	D	1	3,2	-	0/2/22/22	0/1/1/1
2	GLA	D	2	3,2	-	0/2/19/22	0/1/1/1
2	GAL	D	3	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	GLA	C1-O5-C5	4.49	118.27	112.19
2	C	2	GLA	C1-O5-C5	3.34	116.72	112.19
2	C	3	GAL	C6-C5-C4	-2.28	107.65	113.00

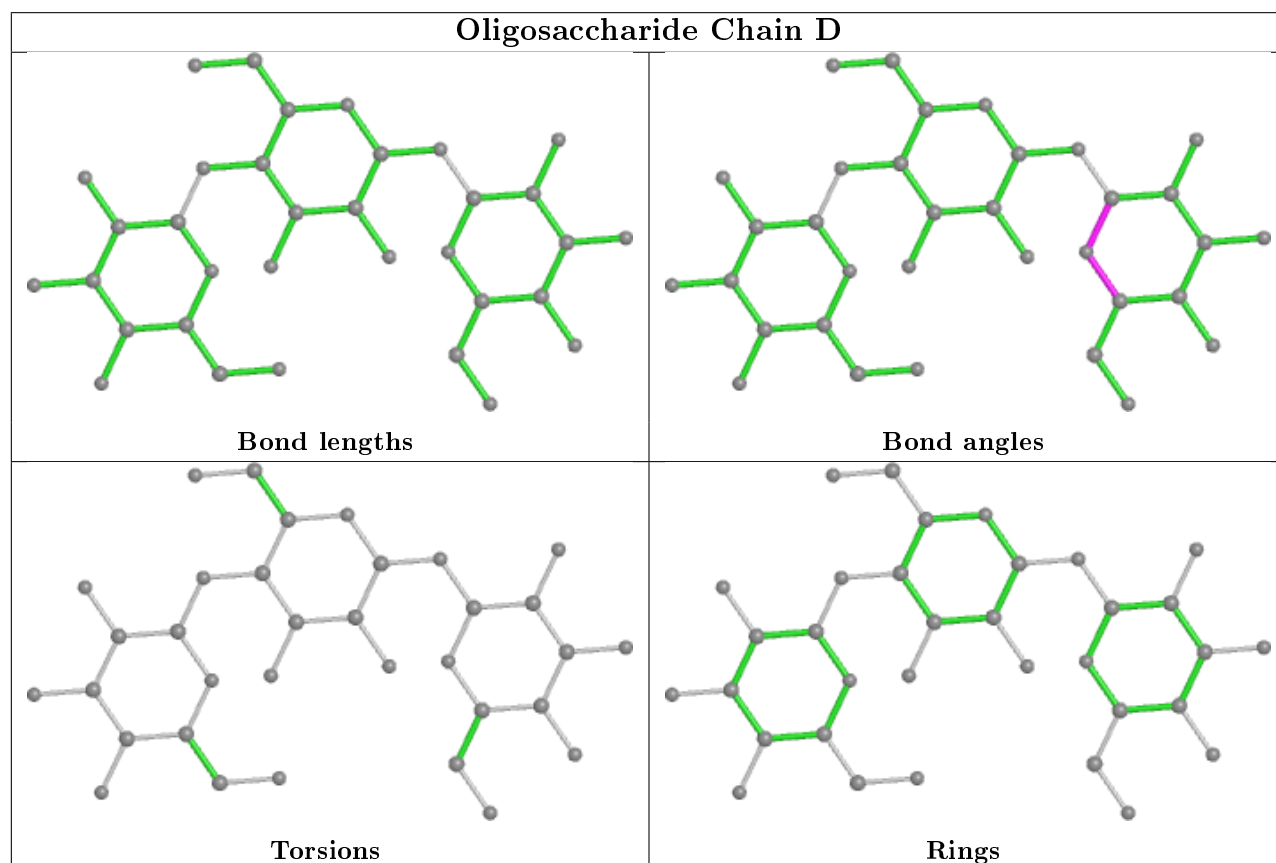
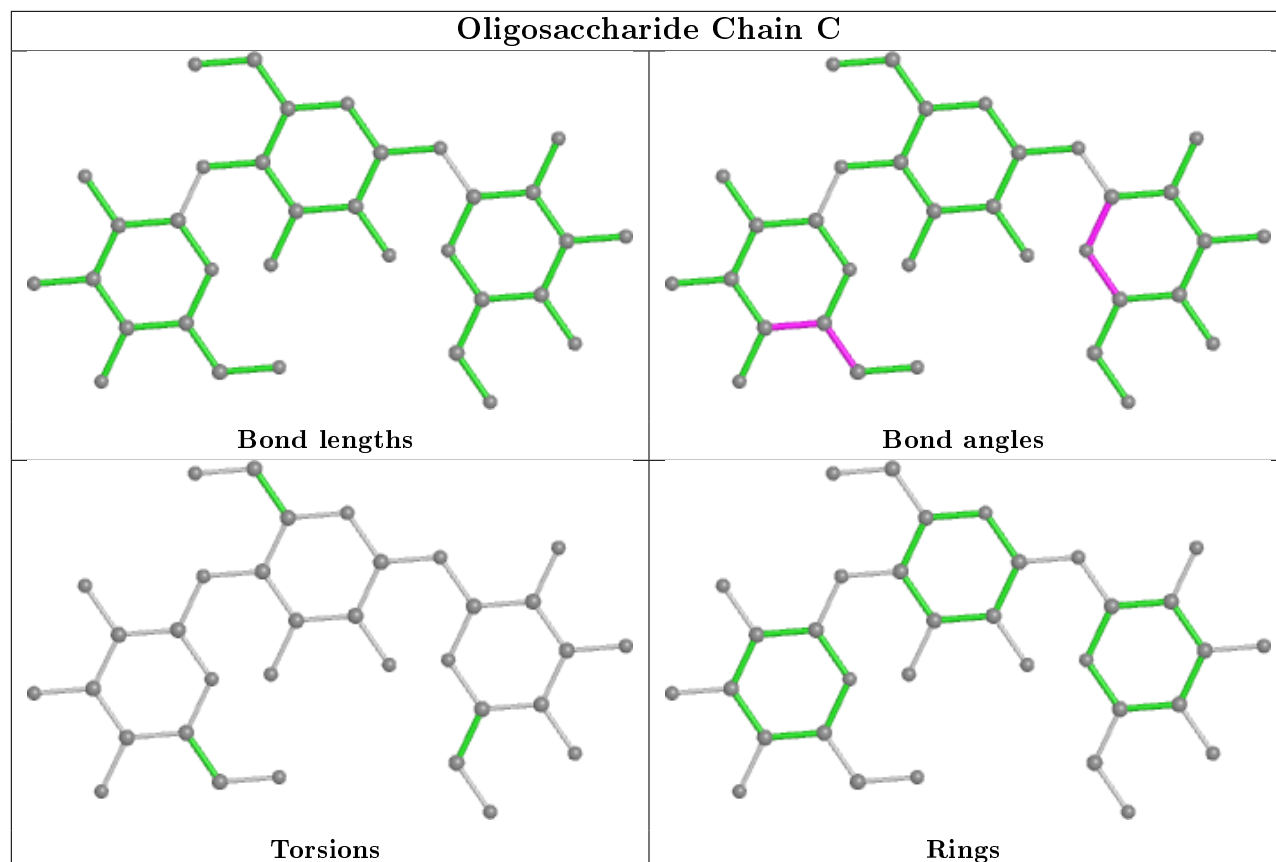
There are no chirality outliers.

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

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	$\# Z > 2$
4	GOL	B	704	-	5,5,5	0.42	0	5,5,5	0.37	0
4	GOL	A	704	-	5,5,5	0.47	0	5,5,5	0.71	0
4	GOL	B	705	-	5,5,5	0.34	0	5,5,5	0.50	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	GOL	B	704	-	-	2/4/4/4	-
4	GOL	A	704	-	-	2/4/4/4	-
4	GOL	B	705	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	704	GOL	O1-C1-C2-C3
4	B	705	GOL	O1-C1-C2-C3
4	B	704	GOL	O1-C1-C2-C3
4	B	705	GOL	O1-C1-C2-O2
4	A	704	GOL	O1-C1-C2-O2
4	B	704	GOL	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	642/655 (98%)	0.32	44 (6%)	16 23	26, 41, 62, 86	0
1	B	644/655 (98%)	0.24	18 (2%)	53 60	25, 41, 63, 93	0
All	All	1286/1310 (98%)	0.28	62 (4%)	30 38	25, 41, 63, 93	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	644	ASP	5.0
1	A	21	GLY	4.5
1	A	68	VAL	4.1
1	B	19	ILE	3.9
1	A	566	VAL	3.5
1	B	241	VAL	3.3
1	A	641	ILE	3.2
1	B	614	LYS	3.2
1	A	92	TYR	3.2
1	A	461	LEU	3.1
1	B	257	ASP	3.0
1	A	497	PRO	3.0
1	A	617	HIS	3.0
1	A	639	PHE	2.9
1	A	289	TYR	2.8
1	B	615	SER	2.8
1	A	110	LYS	2.8
1	A	567	TRP	2.8
1	A	133	LYS	2.7
1	A	638	SER	2.7
1	B	138	VAL	2.7
1	A	257	ASP	2.7
1	A	645	GLY	2.7
1	B	617	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	72	ASN	2.5
1	A	72	ASN	2.5
1	A	450	PRO	2.5
1	B	616	TYR	2.5
1	B	68	VAL	2.5
1	A	616	TYR	2.4
1	A	534	VAL	2.4
1	A	75	LEU	2.4
1	B	24	MET	2.4
1	A	87	ILE	2.4
1	A	112	TRP	2.4
1	A	91	PHE	2.4
1	A	539	PRO	2.3
1	A	462	ASN	2.3
1	B	258	TYR	2.3
1	A	94	ALA	2.3
1	A	132	LYS	2.3
1	B	496	GLY	2.3
1	A	65	ASN	2.3
1	A	460	VAL	2.2
1	B	287	LEU	2.2
1	A	67	THR	2.2
1	A	134	PRO	2.1
1	A	647	LEU	2.1
1	A	458	PRO	2.1
1	A	108	PHE	2.1
1	A	463	PHE	2.1
1	B	238	LYS	2.1
1	B	647	LEU	2.1
1	A	69	TRP	2.1
1	A	290	LEU	2.1
1	A	498	MET	2.1
1	A	640	PRO	2.1
1	B	399	ASN	2.0
1	A	135	PHE	2.0
1	A	284	ALA	2.0
1	A	613	ASP	2.0
1	B	291	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

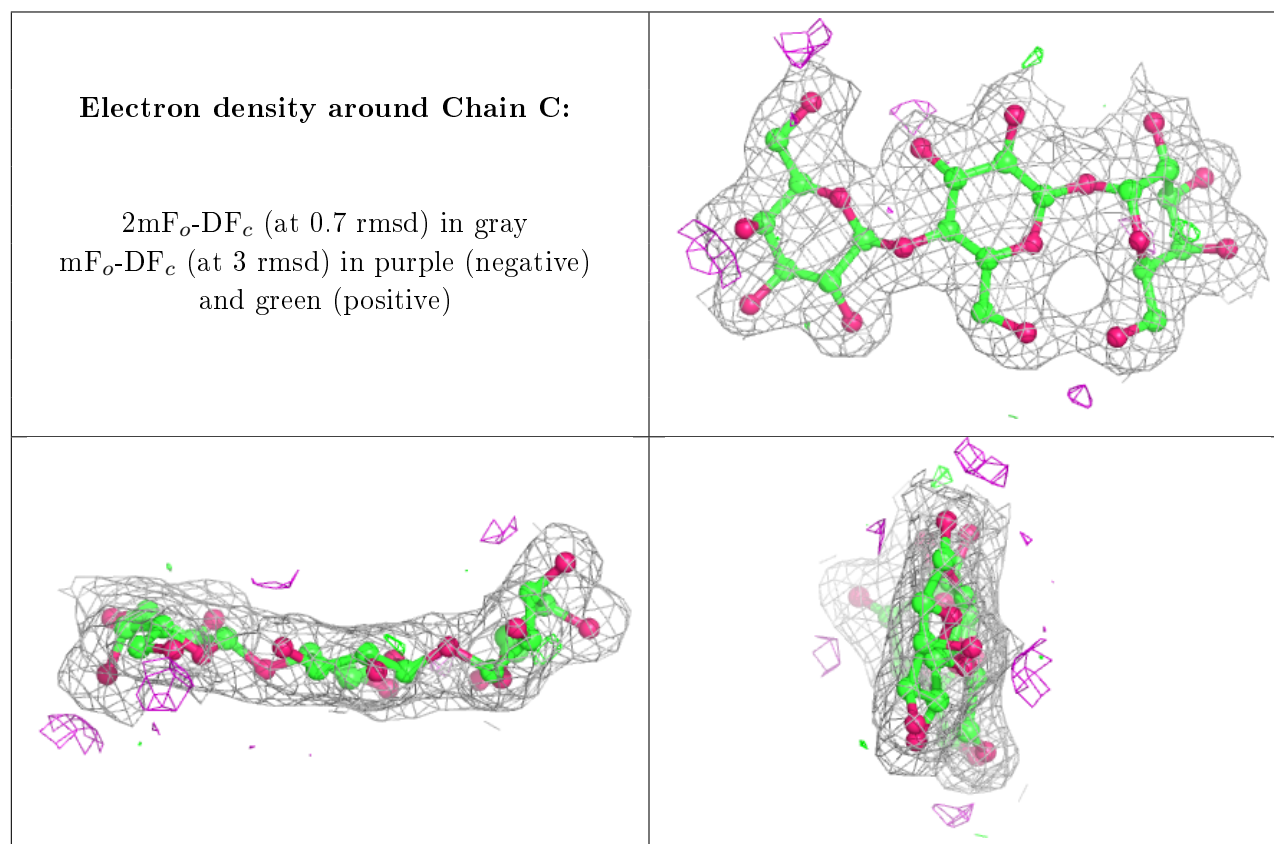
There are no non-standard protein/DNA/RNA residues in this entry.

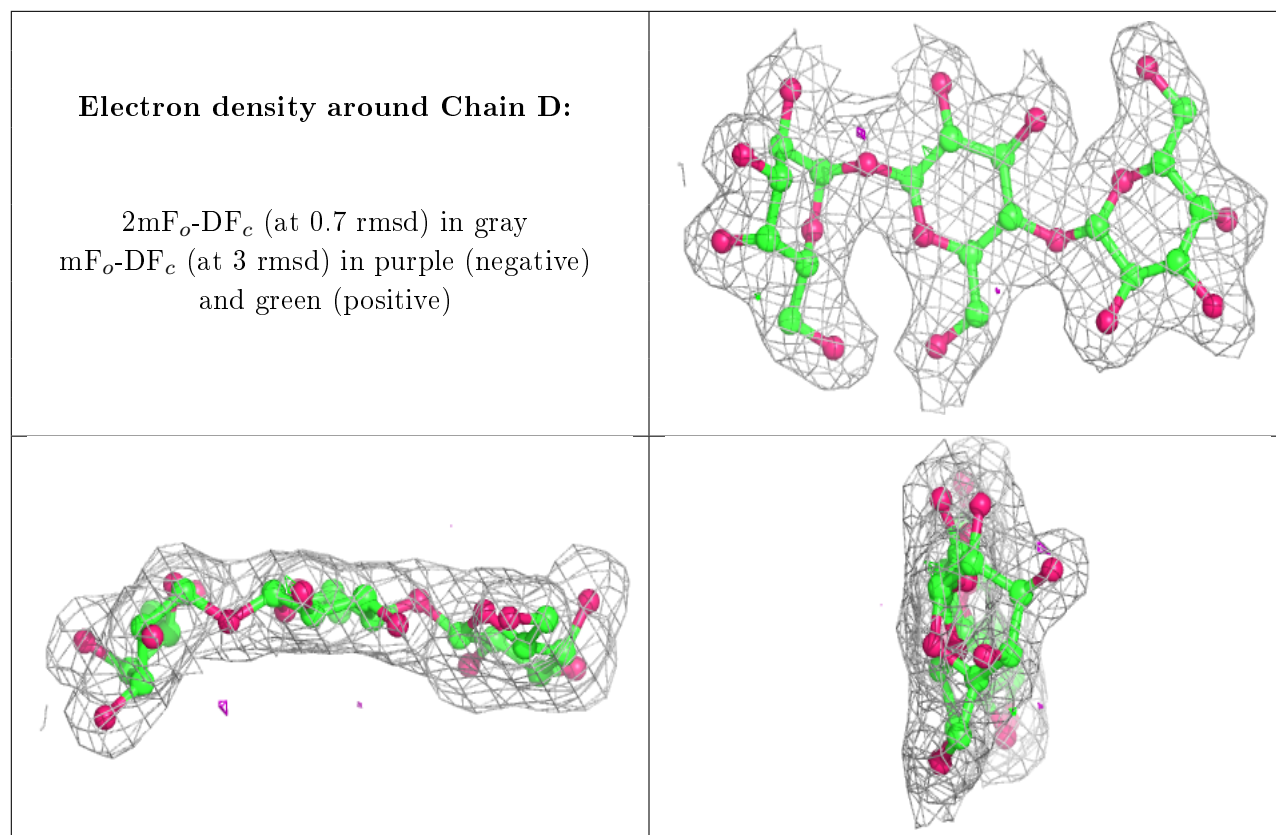
6.3 Carbohydrates ⓘ

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	GAL	D	3	11/12	0.96	0.09	33,34,36,37	0
2	BGC	D	1	12/12	0.96	0.12	28,30,32,33	0
2	GAL	C	3	11/12	0.97	0.08	32,32,34,34	0
2	GLA	D	2	11/12	0.97	0.12	25,28,29,29	0
2	GLA	C	2	11/12	0.97	0.12	27,30,32,33	0
2	BGC	C	1	12/12	0.98	0.09	29,31,32,32	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(\AA^2)	Q<0.9
4	GOL	B	705	6/6	0.85	0.23	58,61,62,62	0
4	GOL	A	704	6/6	0.92	0.11	33,37,38,38	0
4	GOL	B	704	6/6	0.92	0.13	40,40,40,41	0
3	CA	A	701	1/1	0.99	0.14	28,28,28,28	0
3	CA	B	701	1/1	1.00	0.08	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.