



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 12:00 AM BST

PDB ID : 4CN6
Title : GlgE isoform 1 from *Streptomyces coelicolor* E423A mutant with maltose bound
Authors : Syson, K.; Stevenson, C.E.M.; Rashid, A.M.; Saalbach, G.; Tang, M.; Tuukainen, A.; Svergun, D.I.; Withers, S.G.; Lawson, D.M.; Bornemann, S.
Deposited on : 2014-01-21
Resolution : 2.29 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

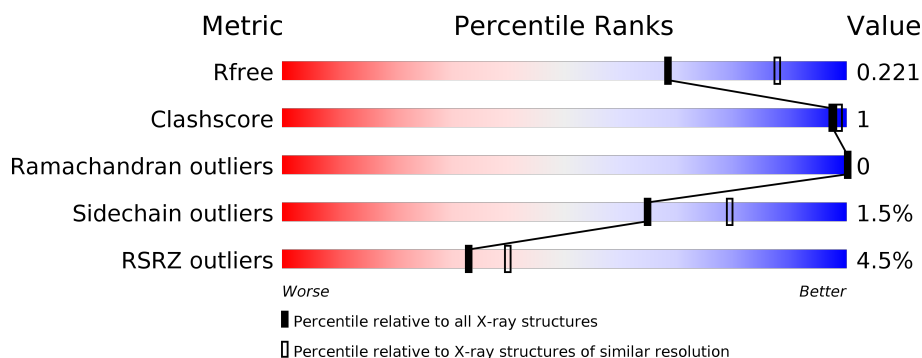
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	695	<div> <div>6%</div> <div> <div></div> <div>91%</div> <div>7%</div> </div> </div>
1	B	695	<div> <div>2%</div> <div> <div></div> <div>91%</div> <div>7%</div> </div> </div>
2	C	2	<div> <div></div> <div>100%</div> </div>
2	D	2	<div> <div></div> <div>100%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10719 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 ALPHA-1,4-GLUCAN\;MALTOSE-1-PHOSPHATE MALTOSYLTRANSFERASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	649	Total	C	N	O	S	0	6	0
			5134	3250	932	942	10			
1	B	649	Total	C	N	O	S	0	2	0
			5059	3205	906	938	10			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q9L1K2
A	-18	GLY	-	expression tag	UNP Q9L1K2
A	-17	SER	-	expression tag	UNP Q9L1K2
A	-16	SER	-	expression tag	UNP Q9L1K2
A	-15	HIS	-	expression tag	UNP Q9L1K2
A	-14	HIS	-	expression tag	UNP Q9L1K2
A	-13	HIS	-	expression tag	UNP Q9L1K2
A	-12	HIS	-	expression tag	UNP Q9L1K2
A	-11	HIS	-	expression tag	UNP Q9L1K2
A	-10	HIS	-	expression tag	UNP Q9L1K2
A	-9	SER	-	expression tag	UNP Q9L1K2
A	-8	SER	-	expression tag	UNP Q9L1K2
A	-7	GLY	-	expression tag	UNP Q9L1K2
A	-6	LEU	-	expression tag	UNP Q9L1K2
A	-5	VAL	-	expression tag	UNP Q9L1K2
A	-4	PRO	-	expression tag	UNP Q9L1K2
A	-3	ARG	-	expression tag	UNP Q9L1K2
A	-2	GLY	-	expression tag	UNP Q9L1K2
A	-1	SER	-	expression tag	UNP Q9L1K2
A	0	HIS	-	expression tag	UNP Q9L1K2
A	423	ALA	GLU	engineered mutation	UNP Q9L1K2
B	-19	MET	-	expression tag	UNP Q9L1K2
B	-18	GLY	-	expression tag	UNP Q9L1K2
B	-17	SER	-	expression tag	UNP Q9L1K2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	SER	-	expression tag	UNP Q9L1K2
B	-15	HIS	-	expression tag	UNP Q9L1K2
B	-14	HIS	-	expression tag	UNP Q9L1K2
B	-13	HIS	-	expression tag	UNP Q9L1K2
B	-12	HIS	-	expression tag	UNP Q9L1K2
B	-11	HIS	-	expression tag	UNP Q9L1K2
B	-10	HIS	-	expression tag	UNP Q9L1K2
B	-9	SER	-	expression tag	UNP Q9L1K2
B	-8	SER	-	expression tag	UNP Q9L1K2
B	-7	GLY	-	expression tag	UNP Q9L1K2
B	-6	LEU	-	expression tag	UNP Q9L1K2
B	-5	VAL	-	expression tag	UNP Q9L1K2
B	-4	PRO	-	expression tag	UNP Q9L1K2
B	-3	ARG	-	expression tag	UNP Q9L1K2
B	-2	GLY	-	expression tag	UNP Q9L1K2
B	-1	SER	-	expression tag	UNP Q9L1K2
B	0	HIS	-	expression tag	UNP Q9L1K2
B	423	ALA	GLU	engineered mutation	UNP Q9L1K2

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	290	Total	O	0	0
			290	290		
3	B	190	Total	O	0	0
			190	190		

B0C1
GLC2

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	114.18Å 114.18Å 313.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.78 – 2.29 37.78 – 2.29	Depositor EDS
% Data completeness (in resolution range)	99.9 (37.78-2.29) 99.9 (37.78-2.29)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.195 , 0.220 0.197 , 0.221	Depositor DCC
R_{free} test set	4697 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	45.3	Xtriage
Anisotropy	0.342	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 26.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10719	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.11% 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: GLC, BGC

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	0.61	0/5305	0.71	0/7257
1	B	0.53	0/5217	0.69	0/7148
All	All	0.57	0/10522	0.70	0/14405

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	5134	0	4961	8	0
1	B	5059	0	4824	7	0
2	C	23	0	21	0	0
2	D	23	0	21	0	0
3	A	290	0	0	4	0
3	B	190	0	0	1	0
All	All	10719	0	9827	15	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.

All (15) 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:340:HIS:ND1	3:A:2157:HOH:O	2.25	0.70
1:A:251:TYR:CZ	1:A:392:ARG:HD2	2.30	0.65
1:B:251:TYR:CZ	1:B:392:ARG:HD2	2.31	0.65
1:A:103:THR:HG21	3:A:2063:HOH:O	1.97	0.64
1:B:25[B]:ASP:OD2	1:B:27:ARG:NH2	2.32	0.63
1:B:25[B]:ASP:OD2	1:B:27:ARG:CZ	2.49	0.61
1:B:26:VAL:HG22	1:B:200:LEU:HD12	1.84	0.60
1:A:389:ARG:HB2	3:A:2081:HOH:O	2.11	0.50
1:B:180:GLN:N	1:B:180:GLN:OE1	2.46	0.49
1:A:212:GLY:O	1:A:507:THR:HA	2.14	0.47
1:A:595:ASN:HA	3:A:2280:HOH:O	2.16	0.46
1:B:212:GLY:O	1:B:507:THR:HA	2.16	0.46
1:A:628:VAL:HG22	1:A:629:PRO:HD2	1.98	0.46
1:B:103:THR:HG21	3:B:2066:HOH:O	2.17	0.43
1:A:141:ARG:NH2	1:A:192:ASP:OD1	2.54	0.41

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	653/695 (94%)	638 (98%)	15 (2%)	0	100	100
1	B	649/695 (93%)	636 (98%)	13 (2%)	0	100	100
All	All	1302/1390 (94%)	1274 (98%)	28 (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	531/574 (92%)	522 (98%)	9 (2%)	60	76
1	B	516/574 (90%)	509 (99%)	7 (1%)	67	81
All	All	1047/1148 (91%)	1031 (98%)	16 (2%)	65	79

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	180	GLN
1	A	196	SER
1	A	275	ASP
1	A	511	TYR
1	A	606	ARG
1	A	623	ASP
1	A	628	VAL
1	A	654	THR
1	B	26	VAL
1	B	31	GLN
1	B	167	ARG
1	B	180	GLN
1	B	196	SER
1	B	511	TYR
1	B	654	THR

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 ⓘ

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	1.48	1 (8%)	17,17,17	1.11	3 (17%)
2	GLC	C	2	2	11,11,12	2.11	2 (18%)	15,15,17	0.91	1 (6%)
2	BGC	D	1	2	12,12,12	1.13	1 (8%)	17,17,17	1.39	2 (11%)
2	GLC	D	2	2	11,11,12	1.93	3 (27%)	15,15,17	1.09	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	-	0/2/22/22	0/1/1/1
2	GLC	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	GLC	D	2	2	-	0/2/19/22	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	GLC	O2-C2	5.83	1.55	1.43
2	D	2	GLC	O2-C2	3.76	1.51	1.43
2	C	2	GLC	O5-C1	3.16	1.48	1.43
2	C	1	BGC	C4-C3	3.10	1.60	1.52
2	D	2	GLC	O5-C1	2.92	1.48	1.43
2	D	2	GLC	C4-C5	2.50	1.58	1.53
2	D	1	BGC	C4-C5	2.16	1.57	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	BGC	O5-C1-C2	-2.84	105.22	110.28
2	D	1	BGC	C1-C2-C3	-2.80	104.51	110.31
2	C	1	BGC	O5-C1-C2	-2.29	106.20	110.28
2	C	2	GLC	O2-C2-C1	2.08	113.40	109.15
2	C	1	BGC	C4-C3-C2	-2.05	107.25	110.82
2	C	1	BGC	C1-C2-C3	-2.01	106.13	110.31

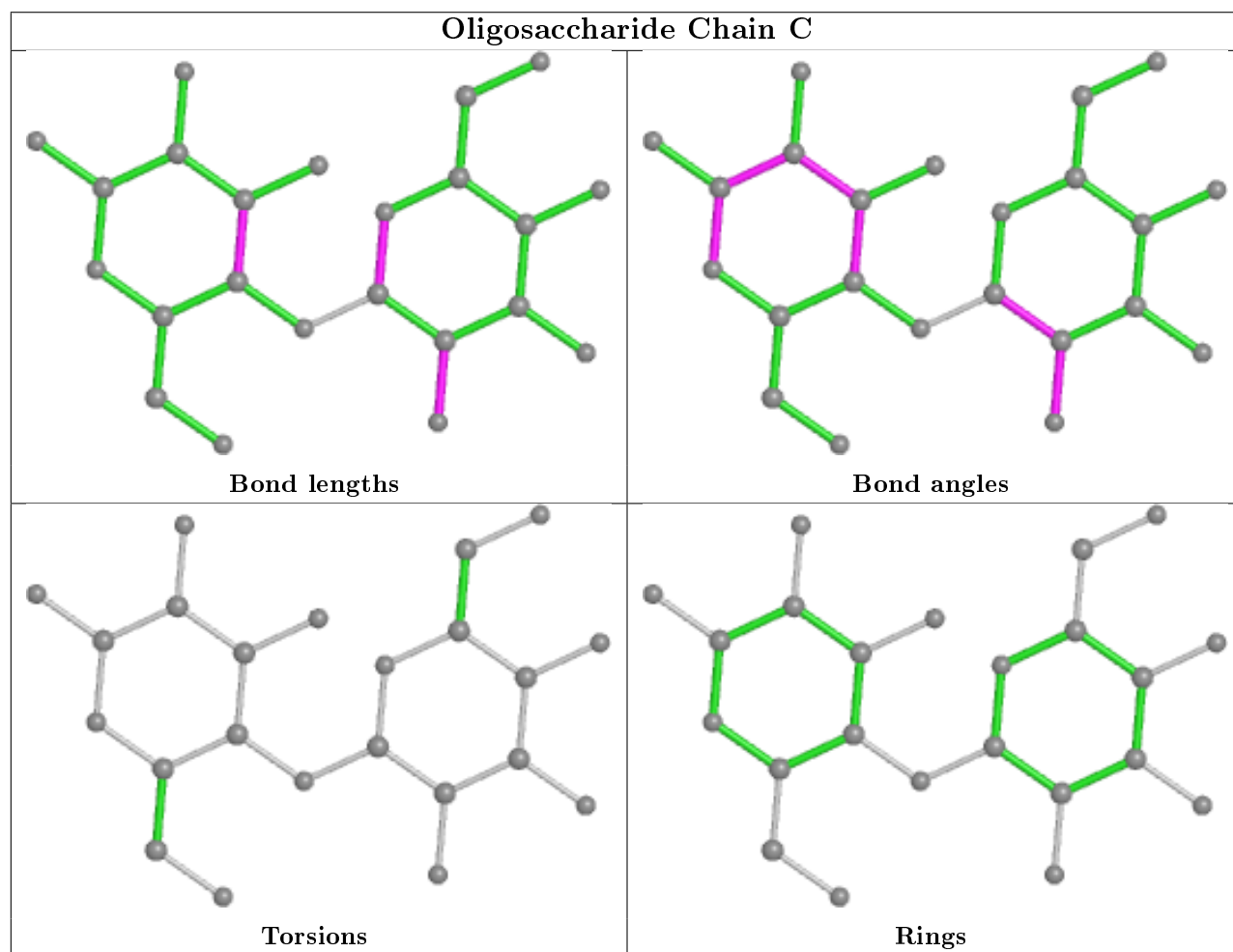
There are no chirality outliers.

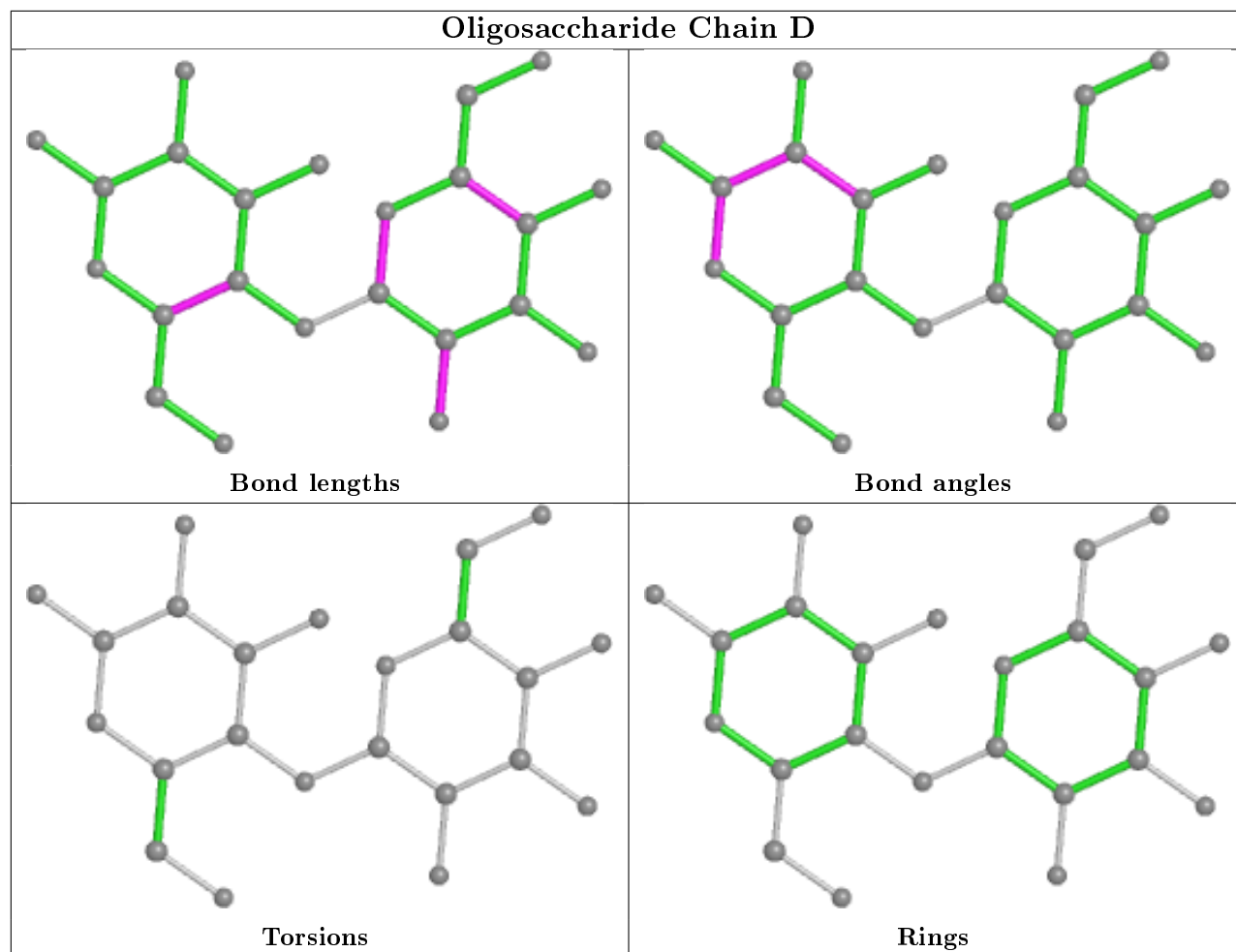
There are no torsion outliers.

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](#)

There are no ligands in this entry.

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	649/695 (93%)	0.05	44 (6%) 17 22	27, 43, 92, 107	0
1	B	649/695 (93%)	-0.13	15 (2%) 60 67	28, 53, 91, 126	0
All	All	1298/1390 (93%)	-0.04	59 (4%) 33 40	27, 47, 92, 126	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	176	ALA	6.8
1	A	181	VAL	6.4
1	A	184	VAL	5.5
1	A	147	PRO	5.2
1	A	178	THR	5.1
1	A	177	LEU	4.9
1	A	173	LEU	4.8
1	A	144	ALA	4.4
1	A	183	ALA	4.3
1	A	174	ALA	4.2
1	A	15	PRO	4.0
1	B	124	ALA	3.9
1	A	155	LEU	3.9
1	A	171	SER	3.9
1	A	169	ALA	3.9
1	B	625	HIS	3.8
1	B	15	PRO	3.7
1	A	125	GLY	3.7
1	A	159	VAL	3.7
1	A	119	ARG	3.7
1	B	624	TRP	3.6
1	A	122	ILE	3.4
1	A	185	LEU	3.4
1	A	180	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	152	ARG	3.3
1	A	175	ALA	3.3
1	B	622	LEU	3.3
1	A	142	ALA	3.1
1	A	123	PRO	3.1
1	A	157	ALA	3.1
1	B	226	HIS	3.1
1	B	126	LEU	3.1
1	B	123	PRO	3.0
1	A	421	LEU	3.0
1	A	226[A]	HIS	3.0
1	A	160	ASP	3.0
1	A	166	SER	2.9
1	A	153	ARG	2.9
1	A	145	ASP	2.8
1	A	168	PRO	2.8
1	A	130	LEU	2.8
1	A	129	ASP	2.7
1	A	148	GLY	2.7
1	B	225	PRO	2.7
1	A	124	ALA	2.7
1	B	641	TRP	2.6
1	B	663	ARG	2.6
1	B	119	ARG	2.6
1	A	118	ALA	2.6
1	B	227	THR	2.5
1	A	127	ASP	2.3
1	A	71	ARG	2.3
1	A	156	LEU	2.2
1	B	644	ALA	2.2
1	A	273	THR	2.2
1	A	225	PRO	2.1
1	A	434	LEU	2.1
1	A	179	PRO	2.1
1	B	122	ILE	2.0

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

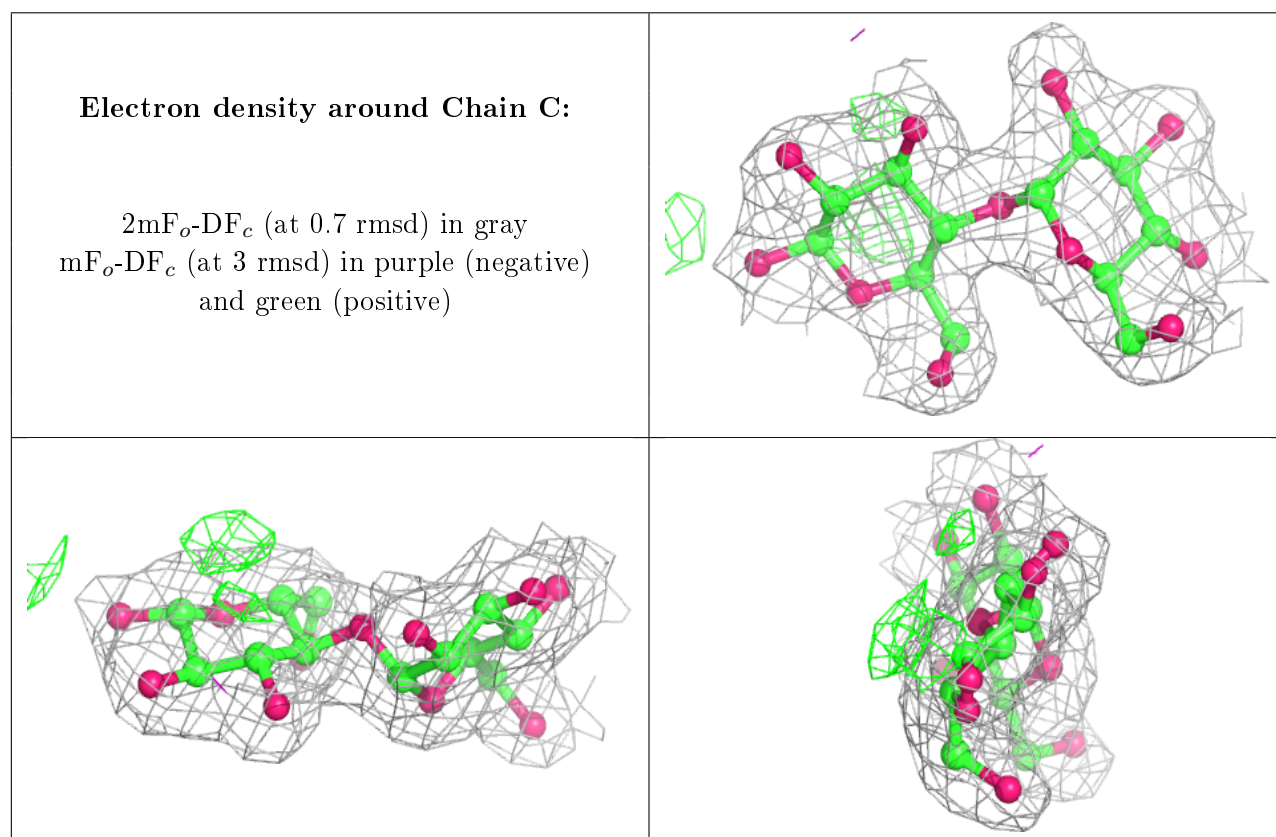
There are no non-standard protein/DNA/RNA residues in this entry.

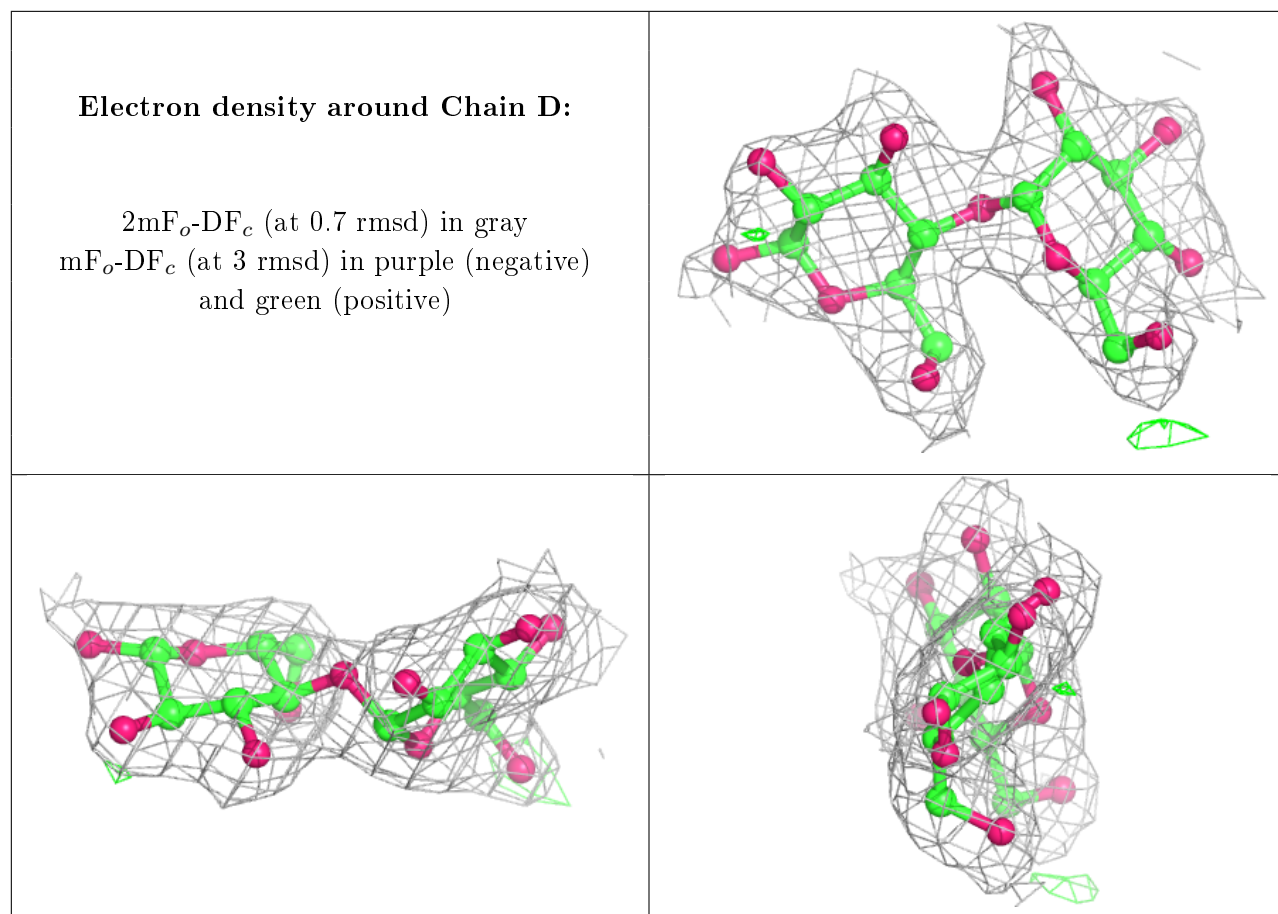
6.3 Carbohydrates ⓘ

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(\AA^2)	Q<0.9
2	BGC	D	1	12/12	0.91	0.15	44,54,61,64	0
2	BGC	C	1	12/12	0.94	0.15	35,41,48,49	0
2	GLC	D	2	11/12	0.95	0.09	40,44,47,48	0
2	GLC	C	2	11/12	0.97	0.13	34,37,39,40	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.