



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:12 AM UTC

PDB ID : 3HZ3 / pdb_00003hz3
Title : Lactobacillus reuteri N-terminally truncated glucansucrase GTF180(D1025N)
-sucrose complex
Authors : Vujicic-Zagar, A.; Pijning, T.; Kralj, S.; Eeuwema, W.; Dijkhuizen, L.; Dijkstra, B.W.
Deposited on : 2009-06-23
Resolution : 2.22 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

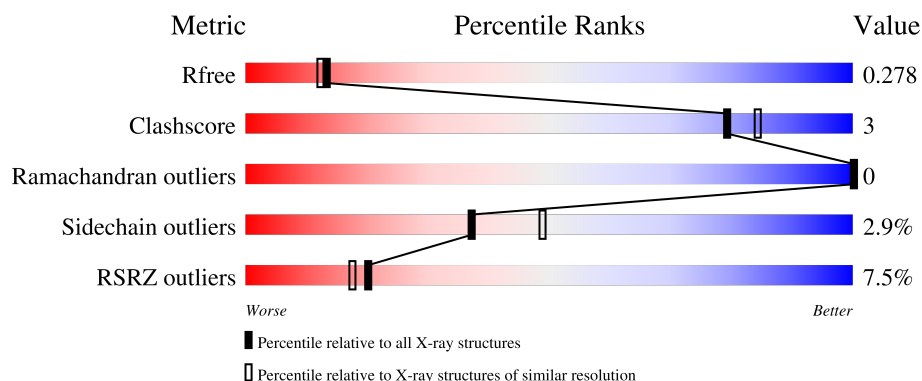
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	1039	7% 87% 9% .
2	B	2	100%
2	C	2	50% 50%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8251 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.

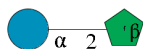
- Molecule 1 is a protein called Glucansucrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1003	7960	4979	1349	1611	21	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	740	MET	-	expression tag	UNP Q5SBN3
A	741	GLY	-	expression tag	UNP Q5SBN3
A	1025	ASN	ASP	engineered mutation	UNP Q5SBN3
A	1674	LEU	PHE	SEE REMARK 999	UNP Q5SBN3
A	1773	HIS	-	expression tag	UNP Q5SBN3
A	1774	HIS	-	expression tag	UNP Q5SBN3
A	1775	HIS	-	expression tag	UNP Q5SBN3
A	1776	HIS	-	expression tag	UNP Q5SBN3
A	1777	HIS	-	expression tag	UNP Q5SBN3
A	1778	HIS	-	expression tag	UNP Q5SBN3

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



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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	244	Total 244	O 244	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.78Å 66.23Å 82.72Å 107.13° 101.37° 85.40°	Depositor
Resolution (Å)	20.00 – 2.22 20.00 – 2.22	Depositor EDS
% Data completeness (in resolution range)	96.3 (20.00-2.22) 96.3 (20.00-2.22)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.21Å)	Xtriage
Refinement program	REFMAC 5.3.0011	Depositor
R, R_{free}	0.193 , 0.243 0.236 , 0.278	Depositor DCC
R_{free} test set	2723 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	31.4	Xtriage
Anisotropy	0.153	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8251	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.67% 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, CA, FRU

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	0.82	1/8136 (0.0%)	0.94	9/11066 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	953	THR	CA-CB	5.16	1.58	1.53

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1669	GLY	N-CA-C	6.11	122.07	114.37
1	A	1200	VAL	CA-C-N	-5.45	114.37	120.14
1	A	1200	VAL	C-N-CA	-5.45	114.37	120.14
1	A	1299	GLY	N-CA-C	5.15	118.30	111.70
1	A	1477	THR	CA-C-N	-5.07	114.73	119.85
1	A	1477	THR	C-N-CA	-5.07	114.73	119.85
1	A	1126	ILE	CA-C-N	5.03	124.79	119.76
1	A	1126	ILE	C-N-CA	5.03	124.79	119.76
1	A	943	GLY	N-CA-C	5.01	120.34	111.01

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	7960	0	7511	41	0
2	B	23	0	21	0	0
2	C	23	0	21	1	0
3	A	1	0	0	0	0
4	A	244	0	0	2	0
All	All	8251	0	7553	42	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 3.

All (42) 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:1390:THR:HG23	1:A:1392:THR:HG22	1.78	0.66
1:A:1544:ASN:HD21	1:A:1604:ARG:HH22	1.42	0.65
1:A:815:ASN:HD22	1:A:1518:ALA:H	1.53	0.54
1:A:854:ASN:HD22	1:A:857:THR:H	1.58	0.52
1:A:866:LYS:HB2	1:A:872:LEU:HB2	1.92	0.52
1:A:1061:ILE:HB	1:A:1129:TYR:CE2	2.45	0.51
1:A:815:ASN:ND2	1:A:1518:ALA:H	2.08	0.50
1:A:749:ILE:HA	1:A:756:PRO:HA	1.94	0.50
1:A:866:LYS:HE3	1:A:876:LEU:O	2.12	0.49
1:A:1409:PHE:CG	1:A:1445:ALA:HB2	2.47	0.49
1:A:1158:PHE:HB2	1:A:1162:ASP:HB2	1.94	0.49
1:A:1475:MET:HG3	1:A:1484:ASP:HB3	1.95	0.49
1:A:1258:LYS:HD3	1:A:1270:GLU:OE1	2.13	0.48
1:A:1203:VAL:HG21	1:A:1230:LEU:HD21	1.97	0.47
1:A:1358:ASN:HB2	1:A:1359:PRO:CD	2.44	0.47
1:A:944:PHE:O	1:A:1523:ARG:HD2	2.15	0.46
1:A:986:ILE:HD11	1:A:1510:ILE:HG13	1.97	0.46
1:A:854:ASN:ND2	1:A:857:THR:H	2.15	0.45
1:A:894:LEU:HD12	1:A:894:LEU:HA	1.80	0.45
1:A:891:TYR:O	1:A:895:VAL:HG23	2.17	0.45
1:A:1612:SER:OG	1:A:1614:ASP:OD2	2.34	0.45
1:A:1247:LYS:HG2	4:A:182:HOH:O	2.16	0.44
1:A:1611:ARG:HD2	4:A:225:HOH:O	2.18	0.44
1:A:829:ILE:HD13	1:A:843:ASP:HB3	2.00	0.44
1:A:1171:ILE:HD12	1:A:1210:GLN:HG2	2.00	0.44
1:A:1388:HIS:HE1	1:A:1393:ALA:O	2.01	0.43
1:A:1423:THR:HB	1:A:1481:TYR:HB3	2.00	0.43
2:C:1:GLC:O6	2:C:1:GLC:O4	2.36	0.43
1:A:848:LEU:HB2	1:A:916:MET:HE2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1604:ARG:HD3	1:A:1608:TYR:CE2	2.54	0.42
1:A:1720:TYR:HB3	1:A:1742:LEU:HD11	2.00	0.42
1:A:757:ARG:NH2	1:A:1751:GLU:OE2	2.53	0.42
1:A:889:ASN:O	1:A:893:GLU:HG3	2.20	0.42
1:A:1201:PRO:HD2	1:A:1403:ASN:O	2.20	0.41
1:A:1241:GLN:HA	1:A:1254:VAL:O	2.20	0.41
1:A:1019:PHE:CG	1:A:1503:ALA:HB2	2.55	0.41
1:A:1253:ASN:HB3	1:A:1278:VAL:HB	2.02	0.41
1:A:1034:LEU:HD23	1:A:1034:LEU:HA	1.86	0.41
1:A:1415:TRP:HA	1:A:1416:PRO:HD3	1.99	0.40
1:A:1650:HIS:HE1	1:A:1679:ASN:OD1	2.04	0.40
1:A:1270:GLU:O	1:A:1274:GLU:HG3	2.21	0.40
1:A:1042:PHE:HB3	1:A:1048:MET:HG3	2.04	0.40

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	999/1039 (96%)	955 (96%)	44 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	855/883 (97%)	830 (97%)	25 (3%)	37	49

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	785	GLN
1	A	788	LYS
1	A	809	LYS
1	A	829	ILE
1	A	854	ASN
1	A	880	SER
1	A	881	THR
1	A	893	GLU
1	A	921	THR
1	A	924	SER
1	A	1013	ASN
1	A	1338	LEU
1	A	1349	GLN
1	A	1362	THR
1	A	1539	VAL
1	A	1561	ASP
1	A	1564	GLN
1	A	1569	GLN
1	A	1611	ARG
1	A	1628	LEU
1	A	1634	VAL
1	A	1674	LEU
1	A	1690	SER
1	A	1708	ARG
1	A	1744	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	755	GLN
1	A	815	ASN
1	A	828	GLN
1	A	854	ASN
1	A	1013	ASN
1	A	1014	ASN
1	A	1174	GLN
1	A	1492	GLN

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Mol	Chain	Res	Type
1	A	1544	ASN
1	A	1650	HIS

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	GLC	B	1	2	11,11,12	1.81	3 (27%)	15,15,17	1.73	4 (26%)
2	FRU	B	2	2	11,12,12	1.18	1 (9%)	10,18,18	1.17	2 (20%)
2	GLC	C	1	2	11,11,12	1.39	2 (18%)	15,15,17	2.41	6 (40%)
2	FRU	C	2	2	11,12,12	2.49	4 (36%)	10,18,18	1.74	2 (20%)

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	GLC	B	1	2	-	2/2/19/22	0/1/1/1
2	FRU	B	2	2	-	0/5/24/24	0/1/1/1
2	GLC	C	1	2	-	1/2/19/22	0/1/1/1
2	FRU	C	2	2	-	3/5/24/24	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	FRU	O5-C2	4.86	1.51	1.43
2	C	2	FRU	O2-C2	4.80	1.49	1.40
2	B	1	GLC	O5-C1	3.78	1.50	1.43
2	B	1	GLC	O5-C5	2.94	1.49	1.43
2	B	2	FRU	O2-C2	2.85	1.45	1.40
2	C	1	GLC	O5-C1	2.57	1.48	1.43
2	C	2	FRU	O3-C3	2.46	1.47	1.42
2	C	2	FRU	O5-C5	2.45	1.49	1.43
2	B	1	GLC	C4-C3	2.14	1.57	1.52
2	C	1	GLC	C2-C3	2.07	1.55	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	GLC	C1-O5-C5	-4.99	105.50	112.19
2	C	1	GLC	C1-C2-C3	4.30	115.90	109.64
2	B	1	GLC	C1-O5-C5	3.95	117.48	112.19
2	C	2	FRU	O5-C5-C6	3.62	118.76	108.89
2	C	1	GLC	O5-C1-C2	3.27	118.60	110.79
2	C	1	GLC	O5-C5-C6	3.12	113.73	107.66
2	C	2	FRU	C6-C5-C4	-2.92	108.20	115.10
2	C	1	GLC	C3-C4-C5	-2.69	105.35	110.23
2	B	1	GLC	O5-C5-C6	-2.61	102.58	107.66
2	B	1	GLC	O5-C1-C2	-2.57	104.65	110.79
2	B	1	GLC	O2-C2-C1	2.38	114.66	109.22
2	C	1	GLC	C6-C5-C4	2.26	118.56	113.02
2	B	2	FRU	O4-C4-C3	-2.25	105.41	112.16
2	B	2	FRU	O5-C5-C6	-2.20	102.89	108.89

There are no chirality outliers.

All (6) torsion outliers are listed below:

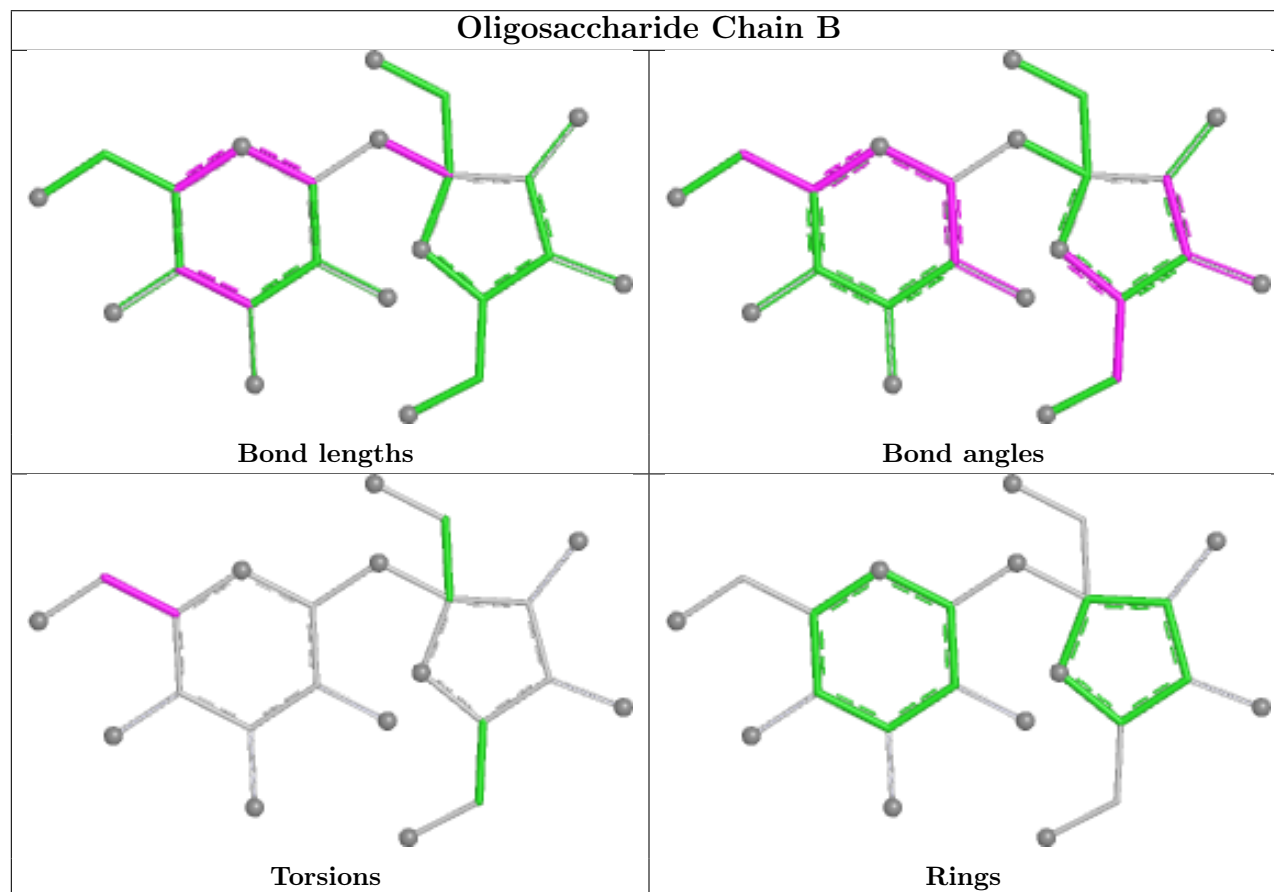
Mol	Chain	Res	Type	Atoms
2	C	2	FRU	O1-C1-C2-C3
2	C	2	FRU	O1-C1-C2-O2
2	C	2	FRU	O1-C1-C2-O5
2	C	1	GLC	O5-C5-C6-O6
2	B	1	GLC	C4-C5-C6-O6
2	B	1	GLC	O5-C5-C6-O6

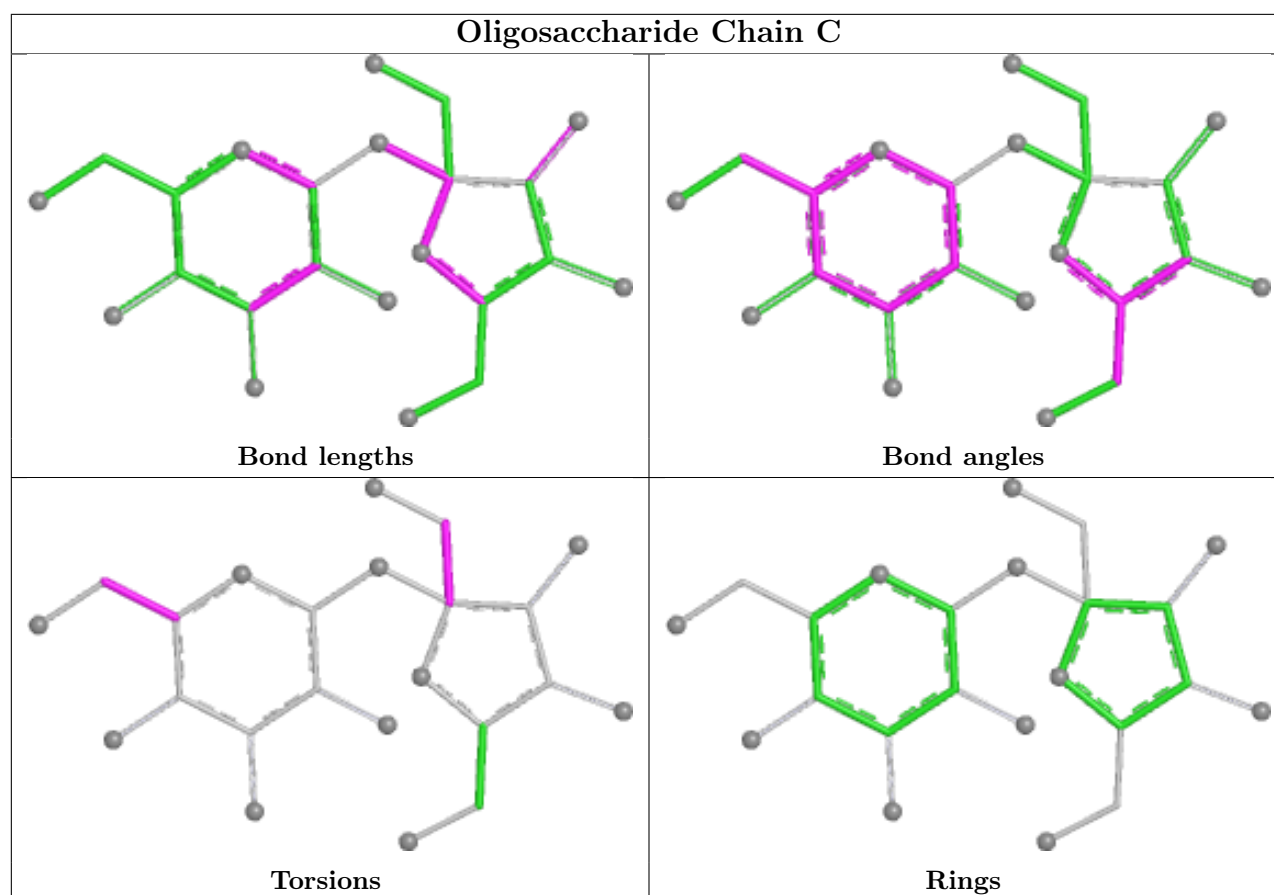
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	GLC	1	0

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 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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	1003/1039 (96%)	0.77	75 (7%) 20 17	35, 56, 84, 110	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	749	ILE	4.4
1	A	1695	SER	3.2
1	A	1387	ALA	3.1
1	A	1571	PHE	3.0
1	A	1014	ASN	3.0
1	A	1420	SER	2.9
1	A	1489	ASN	2.9
1	A	1568	PRO	2.8
1	A	1563	LEU	2.8
1	A	798	ARG	2.8
1	A	876	LEU	2.7
1	A	1428	ALA	2.7
1	A	1012	GLY	2.7
1	A	1733	VAL	2.7
1	A	1008	GLY	2.6
1	A	1385	ASP	2.6
1	A	1010	ILE	2.6
1	A	881	THR	2.6
1	A	1009	THR	2.5
1	A	782	LEU	2.5
1	A	891	TYR	2.4
1	A	1044	ALA	2.4
1	A	921	THR	2.4
1	A	1562	THR	2.4
1	A	1559	PHE	2.4
1	A	753	THR	2.4
1	A	1582	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	1487	LEU	2.4
1	A	1161	ASP	2.3
1	A	751	PRO	2.3
1	A	1045	ALA	2.3
1	A	1659	VAL	2.3
1	A	830	LEU	2.3
1	A	1640	TYR	2.3
1	A	874	ALA	2.3
1	A	886	ALA	2.3
1	A	879	PHE	2.3
1	A	1569	GLN	2.3
1	A	1047	ASN	2.3
1	A	1011	THR	2.3
1	A	1561	ASP	2.2
1	A	754	GLY	2.2
1	A	789	GLY	2.2
1	A	748	TYR	2.2
1	A	1691	SER	2.2
1	A	885	SER	2.2
1	A	888	LEU	2.2
1	A	1492	GLN	2.2
1	A	829	ILE	2.2
1	A	1574	VAL	2.2
1	A	1564	GLN	2.2
1	A	1696	GLY	2.2
1	A	836	TRP	2.1
1	A	1701	LEU	2.1
1	A	894	LEU	2.1
1	A	1017	ALA	2.1
1	A	1013	ASN	2.1
1	A	1664	ILE	2.1
1	A	752	THR	2.1
1	A	1738	MET	2.1
1	A	795	GLU	2.1
1	A	1570	LEU	2.1
1	A	1620	LEU	2.1
1	A	1002	TYR	2.1
1	A	884	ASP	2.0
1	A	1491	LEU	2.0
1	A	1560	LEU	2.0
1	A	1670	ASN	2.0
1	A	1634	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	778	GLY	2.0
1	A	783	LYS	2.0
1	A	855	THR	2.0
1	A	1389	THR	2.0
1	A	1579	GLN	2.0
1	A	791	ILE	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](#)

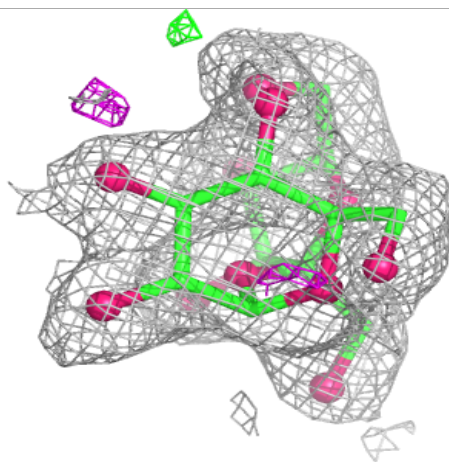
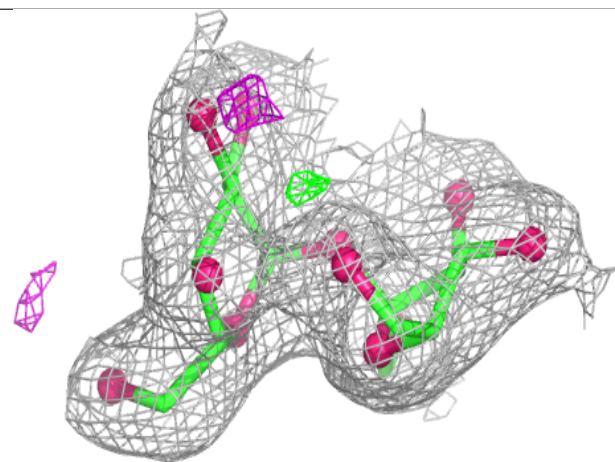
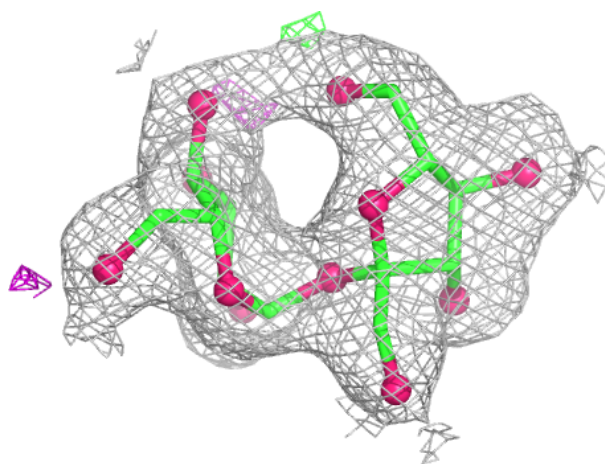
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

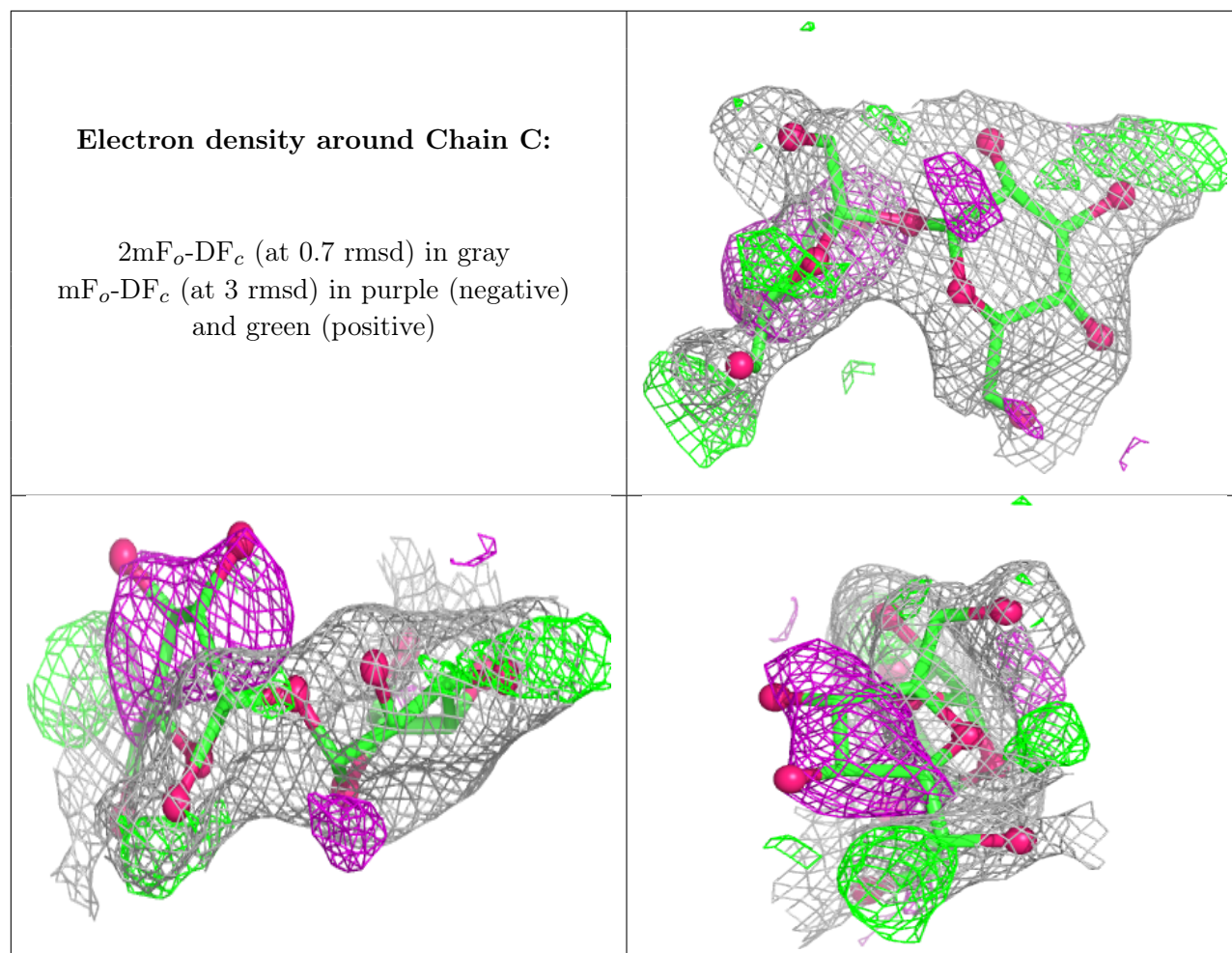
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FRU	C	2	12/12	0.43	0.24	59,67,71,72	0
2	GLC	C	1	11/12	0.81	0.15	49,54,64,67	0
2	GLC	B	1	11/12	0.93	0.08	37,39,43,44	0
2	FRU	B	2	12/12	0.94	0.07	36,40,43,43	0

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 B:

$2mF_o-DF_c$ (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($\text{\AA}^2$)	Q<0.9
3	CA	A	1	1/1	0.98	0.03	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.