



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:00 PM UTC

PDB ID : 5CIM / pdb_00005cim
Title : Structure of Mycobacterium thermoresistibile GlgE in complex with maltose (cocrystallisation with maltose-1-phosphate) at 3.32Å resolution
Authors : Mendes, V.; Blaszczyk, M.; Maranha, A.; Empadinhas, N.; Blundell, T.L.
Deposited on : 2015-07-13
Resolution : 3.32 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

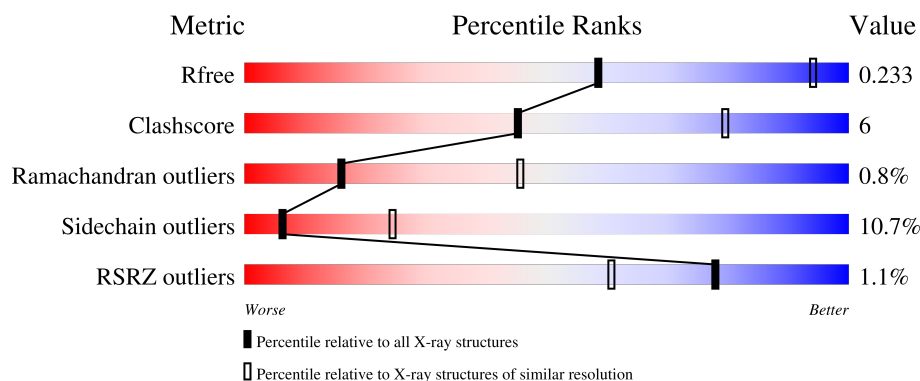
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.32 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	1153 (3.34-3.30)
Clashscore	190562	1193 (3.34-3.30)
Ramachandran outliers	187476	1172 (3.34-3.30)
Sidechain outliers	187428	1171 (3.34-3.30)
RSRZ outliers	180081	1153 (3.34-3.30)

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 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	698	
1	B	698	
2	C	2	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10246 atoms, of which 22 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	636	Total	C	N	O	S	0	0	0
			4924	3158	870	887	9			
1	B	675	Total	C	N	O	S	0	0	0
			5211	3347	901	954	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP G7CL00
A	0	SER	-	expression tag	UNP G7CL00
A	1	VAL	-	cloning artifact	UNP G7CL00
B	-1	GLY	-	expression tag	UNP G7CL00
B	0	SER	-	expression tag	UNP G7CL00
B	1	VAL	-	cloning artifact	UNP G7CL00

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	H	O	0	0	0
			45	12	22	11			

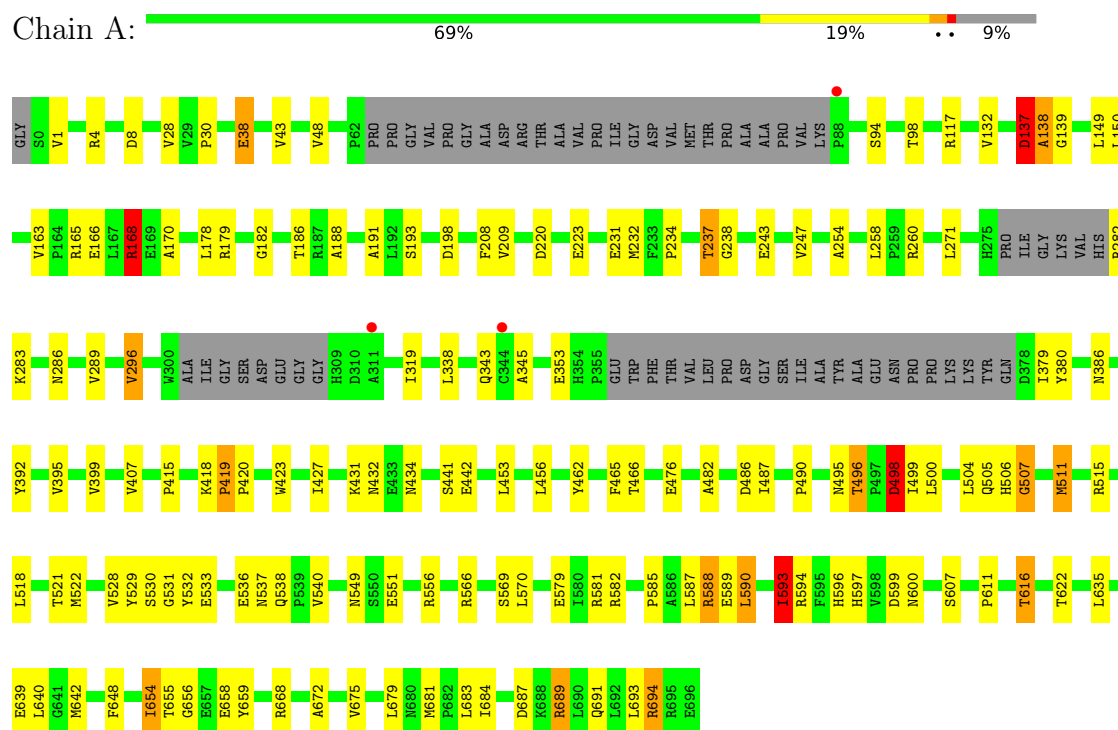
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total	O	0	0
			41	41		
3	B	25	Total	O	0	0
			25	25		

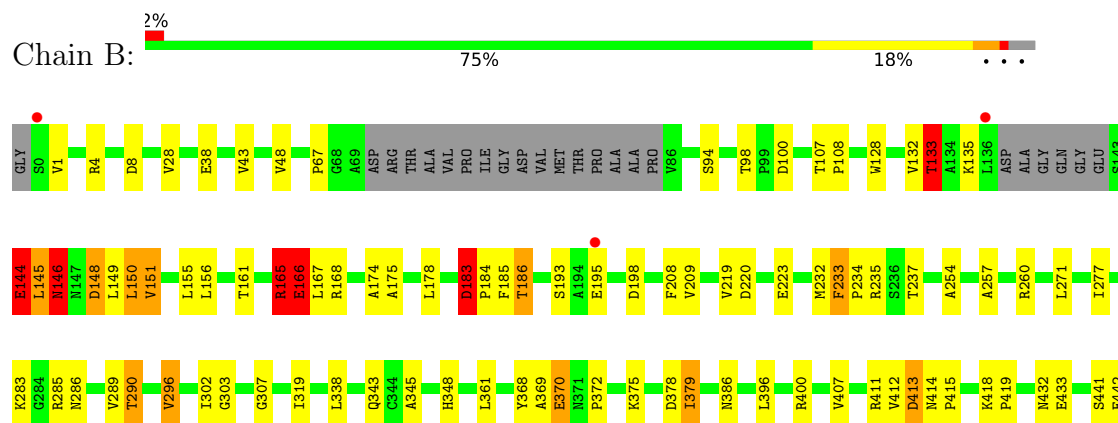
3 Residue-property plots

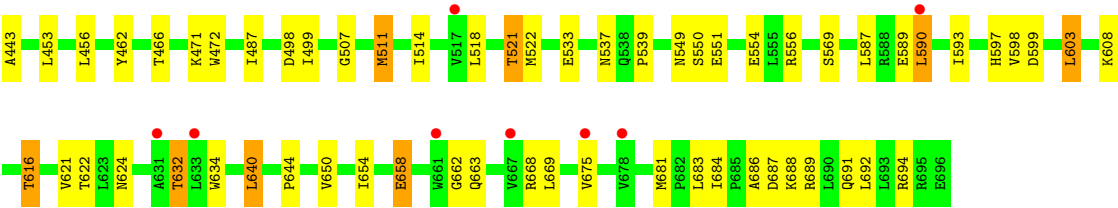
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Alpha-1,4-glucan:maltose-1-phosphate maltosyltransferase



- Molecule 1: Alpha-1,4-glucan:maltose-1-phosphate maltosyltransferase





● Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain C: 100%

GLC1
GLC2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.54Å 112.97Å 221.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.63 – 3.32 100.63 – 3.32	Depositor EDS
% Data completeness (in resolution range)	99.7 (100.63-3.32) 99.7 (100.63-3.32)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 3.33Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.10.0	Depositor
R, R_{free}	0.178 , 0.214 0.194 , 0.233	Depositor DCC
R_{free} test set	1497 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	94.5	Xtriage
Anisotropy	0.578	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 99.3	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	0.94	EDS
Total number of atoms	10246	wwPDB-VP
Average B, all atoms (Å ²)	110.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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

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	0.88	0/5071	1.41	43/6940 (0.6%)
1	B	0.87	4/5373 (0.1%)	1.42	49/7367 (0.7%)
All	All	0.88	4/10444 (0.0%)	1.42	92/14307 (0.6%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	144	GLU	CA-C	7.25	1.62	1.52
1	B	133	THR	CA-C	5.45	1.59	1.53
1	B	133	THR	CA-CB	5.25	1.58	1.53
1	B	183	ASP	CA-C	5.19	1.59	1.52

All (92) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	148	ASP	N-CA-C	-10.90	99.55	113.72
1	A	137	ASP	CA-C-N	9.03	137.95	121.70
1	A	137	ASP	C-N-CA	9.03	137.95	121.70
1	B	166	GLU	CA-C-N	8.85	137.63	121.70
1	B	166	GLU	C-N-CA	8.85	137.63	121.70
1	A	168	ARG	N-CA-C	8.67	120.35	111.07
1	B	133	THR	CA-C-N	8.43	136.86	121.70
1	B	133	THR	C-N-CA	8.43	136.86	121.70
1	B	186	THR	CA-C-N	7.83	136.50	121.54
1	B	186	THR	C-N-CA	7.83	136.50	121.54
1	A	182	GLY	N-CA-C	7.75	120.21	110.29
1	A	379	ILE	CA-C-N	7.07	132.86	121.83
1	A	379	ILE	C-N-CA	7.07	132.86	121.83
1	A	498	ASP	CA-CB-CG	6.88	119.48	112.60
1	B	165	ARG	CA-C-N	6.77	134.48	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	165	ARG	C-N-CA	6.77	134.48	121.54
1	A	521	THR	CA-C-N	6.36	128.70	120.44
1	A	521	THR	C-N-CA	6.36	128.70	120.44
1	A	537	ASN	CA-CB-CG	6.31	118.91	112.60
1	B	521	THR	CA-C-N	6.26	128.58	120.44
1	B	521	THR	C-N-CA	6.26	128.58	120.44
1	B	662	GLY	CA-C-N	6.21	129.75	120.31
1	B	662	GLY	C-N-CA	6.21	129.75	120.31
1	A	432	ASN	CA-C-N	6.21	128.59	120.28
1	A	432	ASN	C-N-CA	6.21	128.59	120.28
1	B	193	SER	CA-C-N	6.11	128.75	120.44
1	B	193	SER	C-N-CA	6.11	128.75	120.44
1	A	193	SER	CA-C-N	6.02	128.63	120.44
1	A	193	SER	C-N-CA	6.02	128.63	120.44
1	B	307	GLY	N-CA-C	5.99	119.31	110.63
1	B	146	ASN	N-CA-C	5.94	117.15	107.23
1	B	233	PHE	CA-CB-CG	5.92	119.72	113.80
1	B	1	VAL	N-CA-C	5.78	117.50	109.29
1	A	237	THR	CA-C-N	5.70	126.31	119.98
1	A	237	THR	C-N-CA	5.70	126.31	119.98
1	A	198	ASP	CA-CB-CG	5.69	118.29	112.60
1	A	28	VAL	N-CA-CB	5.68	117.98	111.39
1	A	28	VAL	CA-C-N	-5.67	118.75	122.60
1	A	28	VAL	C-N-CA	-5.67	118.75	122.60
1	A	570	LEU	N-CA-C	-5.66	106.38	113.28
1	B	257	ALA	CA-C-N	5.65	126.78	119.78
1	B	257	ALA	C-N-CA	5.65	126.78	119.78
1	B	296	VAL	N-CA-C	5.57	116.31	110.62
1	B	198	ASP	CA-CB-CG	5.56	118.16	112.60
1	B	599	ASP	N-CA-C	5.54	118.03	111.33
1	B	186	THR	CB-CA-C	5.54	120.67	112.03
1	B	432	ASN	CA-C-N	5.54	127.64	120.44
1	B	432	ASN	C-N-CA	5.54	127.64	120.44
1	B	185	PHE	N-CA-C	-5.49	106.42	113.01
1	A	392	TYR	CA-C-N	5.49	127.63	120.28
1	A	392	TYR	C-N-CA	5.49	127.63	120.28
1	A	419	PRO	N-CA-C	5.48	117.38	110.70
1	B	632	THR	CB-CA-C	5.47	119.06	110.19
1	A	486	ASP	CA-CB-CG	5.44	118.04	112.60
1	A	296	VAL	N-CA-C	5.42	116.15	110.62
1	A	597	HIS	N-CA-C	5.42	117.55	109.59
1	A	345	ALA	N-CA-C	-5.40	102.36	110.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	589	GLU	CB-CG-CD	5.39	121.77	112.60
1	A	1	VAL	N-CA-C	5.37	116.92	109.29
1	B	345	ALA	N-CA-C	-5.37	102.41	110.24
1	B	166	GLU	CB-CG-CD	5.34	121.68	112.60
1	A	379	ILE	N-CA-C	5.26	117.43	108.86
1	B	378	ASP	CA-CB-CG	5.24	117.84	112.60
1	A	258	LEU	N-CA-C	5.23	120.05	113.25
1	B	621	VAL	N-CA-C	5.23	115.65	108.12
1	B	472	TRP	CA-C-N	5.20	127.56	120.54
1	B	472	TRP	C-N-CA	5.20	127.56	120.54
1	A	507	GLY	N-CA-C	5.20	120.45	114.16
1	A	353	GLU	N-CA-C	5.16	116.71	111.14
1	A	170	ALA	CA-C-N	5.16	127.45	120.44
1	A	170	ALA	C-N-CA	5.16	127.45	120.44
1	B	598	VAL	CB-CA-C	5.14	118.24	110.63
1	A	672	ALA	N-CA-C	-5.14	104.64	112.04
1	B	419	PRO	N-CA-C	5.11	116.94	110.70
1	B	499	ILE	N-CA-C	5.10	114.83	106.72
1	B	166	GLU	N-CA-C	-5.09	99.96	110.80
1	B	195	GLU	CA-C-N	5.07	126.94	120.56
1	B	195	GLU	C-N-CA	5.07	126.94	120.56
1	B	514	ILE	CA-C-N	5.07	127.02	120.44
1	B	514	ILE	C-N-CA	5.07	127.02	120.44
1	A	689	ARG	N-CA-C	5.06	118.72	112.54
1	B	372	PRO	N-CA-C	5.06	116.87	110.70
1	B	290	THR	CB-CA-C	5.04	118.53	110.16
1	A	138	ALA	CA-C-N	5.02	130.73	121.70
1	A	138	ALA	C-N-CA	5.02	130.73	121.70
1	B	379	ILE	N-CA-C	5.02	117.23	108.95
1	A	593	ILE	N-CA-CB	5.01	116.48	110.31
1	A	656	GLY	CA-C-N	5.01	127.82	120.71
1	A	656	GLY	C-N-CA	5.01	127.82	120.71
1	B	550	SER	CA-C-N	5.00	128.60	120.60
1	B	550	SER	C-N-CA	5.00	128.60	120.60
1	A	38	GLU	CB-CG-CD	5.00	121.10	112.60

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	4924	0	4704	57	0
1	B	5211	0	4953	59	0
2	C	23	22	20	0	0
3	A	41	0	0	0	0
3	B	25	0	0	1	0
All	All	10224	22	9677	116	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 6.

All (116) 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:133:THR:HA	1:B:135:LYS:H	1.39	0.88
1:A:137:ASP:HB3	1:A:138:ALA:HB3	1.60	0.83
1:A:495:ASN:HD21	1:A:504:LEU:HD13	1.45	0.79
1:B:183:ASP:HB3	1:B:184:PRO:HD2	1.65	0.78
1:B:593:ILE:HG12	1:B:608:LYS:HE2	1.69	0.74
1:A:684:ILE:O	1:A:689:ARG:NH2	2.22	0.72
1:B:174:ALA:O	1:B:178:LEU:HD23	1.90	0.71
1:B:590:LEU:O	1:B:593:ILE:HG13	1.91	0.70
1:B:415:PRO:HG2	1:B:441:SER:HB2	1.76	0.67
1:B:183:ASP:CB	1:B:184:PRO:HD2	2.26	0.66
1:A:684:ILE:HG22	1:A:689:ARG:HB3	1.80	0.64
1:A:163:VAL:O	1:A:168:ARG:HD2	1.98	0.64
1:A:585:PRO:HA	1:A:588:ARG:HE	1.62	0.64
1:A:551:GLU:HG3	1:A:556:ARG:HB2	1.81	0.63
1:B:370:GLU:HG3	1:B:375:LYS:HA	1.82	0.61
1:B:144:GLU:OE2	1:B:145:LEU:HA	2.02	0.60
1:A:581:ARG:HA	1:A:587:LEU:HD12	1.83	0.60
1:B:148:ASP:C	1:B:150:LEU:H	2.08	0.60
1:B:640:LEU:HD22	1:B:681:MET:HE2	1.83	0.60
1:B:551:GLU:HG3	1:B:556:ARG:HB2	1.84	0.59
1:A:642:MET:HE1	1:A:648:PHE:HB3	1.84	0.59
1:A:654:ILE:HG22	1:A:655:THR:HG23	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:611:PRO:O	1:A:689:ARG:HD2	2.02	0.59
1:B:688:LYS:O	1:B:692:LEU:HG	2.04	0.58
1:A:178:LEU:HD21	1:A:191:ALA:HB2	1.86	0.57
1:B:128:TRP:CE2	1:B:132:VAL:HG21	2.39	0.57
1:B:686:ALA:HA	1:B:689:ARG:HG3	1.87	0.56
1:B:183:ASP:HB3	1:B:184:PRO:CD	2.35	0.55
1:B:303:GLY:HA3	1:B:348:HIS:CD2	2.42	0.55
1:A:282:ARG:HG2	1:A:283:LYS:H	1.72	0.55
1:B:400:ARG:HE	1:B:433:GLU:CD	2.15	0.55
1:B:148:ASP:C	1:B:150:LEU:N	2.67	0.53
1:B:165:ARG:HA	1:B:168:ARG:HE	1.73	0.53
1:A:234:PRO:HG3	1:A:271:LEU:HD22	1.90	0.53
1:A:415:PRO:HG2	1:A:441:SER:HB2	1.92	0.52
1:A:442:GLU:HA	1:A:462:TYR:CD1	2.45	0.51
1:B:234:PRO:HG3	1:B:271:LEU:HD22	1.92	0.50
1:A:596:HIS:HE2	1:A:607:SER:HB3	1.76	0.50
1:A:616:THR:HB	1:A:681:MET:HB2	1.93	0.50
1:B:597:HIS:HD1	1:B:634:TRP:CD1	2.30	0.50
1:B:411:ARG:NH1	1:B:413:ASP:OD1	2.45	0.49
1:A:531:GLY:O	1:A:536:GLU:HG2	2.11	0.49
1:B:412:VAL:HG12	1:B:415:PRO:HG3	1.94	0.49
1:B:166:GLU:O	1:B:166:GLU:HG2	2.12	0.49
1:A:338:LEU:HG	1:A:407:VAL:HG11	1.95	0.49
1:B:132:VAL:O	1:B:132:VAL:HG12	2.13	0.49
1:A:482:ALA:HB2	1:A:593:ILE:HG23	1.94	0.48
1:B:442:GLU:HA	1:B:462:TYR:CD1	2.48	0.48
1:B:684:ILE:HG22	1:B:689:ARG:HG2	1.95	0.48
1:B:537:ASN:O	1:B:539:PRO:HD3	2.13	0.48
1:A:528:VAL:HG13	1:A:532:TYR:HB3	1.96	0.47
1:B:616:THR:HB	1:B:681:MET:HB2	1.95	0.47
1:A:282:ARG:CG	1:A:283:LYS:H	2.26	0.47
1:A:635:LEU:HD22	1:A:640:LEU:HD11	1.96	0.47
1:B:338:LEU:HG	1:B:407:VAL:HG11	1.96	0.47
1:B:156:LEU:HB2	1:B:175:ALA:HB2	1.96	0.47
1:B:148:ASP:HA	1:B:151:VAL:HG13	1.97	0.47
1:A:237:THR:HG21	1:A:254:ALA:HA	1.97	0.47
1:B:233:PHE:HB2	1:B:551:GLU:OE2	2.15	0.46
1:B:28:VAL:HG22	1:B:107:THR:OG1	2.15	0.46
1:A:165:ARG:HA	1:A:168:ARG:HD3	1.97	0.46
1:A:622:THR:HG22	1:A:675:VAL:O	2.16	0.46
1:A:496:THR:C	1:A:498:ASP:H	2.24	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:622:THR:HG22	1:B:675:VAL:O	2.15	0.45
1:B:151:VAL:O	1:B:155:LEU:HG	2.17	0.45
1:B:237:THR:HG21	1:B:254:ALA:HA	1.99	0.45
1:B:471:LYS:HE3	1:B:603:LEU:O	2.17	0.45
1:B:650:VAL:O	1:B:658:GLU:HA	2.16	0.45
1:A:231:GLU:HB3	1:A:529:TYR:HA	1.99	0.45
1:A:500:LEU:HG	1:A:505:GLN:HE21	1.81	0.45
1:B:260:ARG:NH2	1:B:533:GLU:OE2	2.50	0.44
1:B:166:GLU:O	1:B:166:GLU:CG	2.65	0.44
1:B:644:PRO:O	1:B:663:GLN:HB2	2.17	0.44
1:A:507:GLY:HA3	1:A:511:MET:HG2	2.00	0.44
1:B:148:ASP:O	1:B:151:VAL:HG22	2.18	0.44
1:A:590:LEU:O	1:A:593:ILE:HG22	2.18	0.44
1:B:507:GLY:HA3	1:B:511:MET:HG2	1.99	0.44
1:A:178:LEU:HD21	1:A:191:ALA:CB	2.47	0.43
1:B:396:LEU:HG	1:B:400:ARG:HD2	2.00	0.43
1:B:302:ILE:O	1:B:348:HIS:HD2	2.01	0.43
1:B:286:ASN:HD21	1:B:549:ASN:HA	1.82	0.43
1:A:149:LEU:HD21	1:A:188:ALA:HB2	2.01	0.43
1:B:518:LEU:O	1:B:522:MET:HB2	2.18	0.43
1:A:238:GLY:HA3	1:A:247:VAL:O	2.19	0.43
1:A:593:ILE:O	1:A:594:ARG:NH2	2.52	0.43
1:A:260:ARG:NH1	1:A:533:GLU:OE2	2.52	0.42
1:A:518:LEU:O	1:A:522:MET:HB2	2.19	0.42
1:B:183:ASP:CB	1:B:184:PRO:CD	2.94	0.42
1:B:4:ARG:HG2	1:B:208:PHE:CG	2.54	0.42
1:B:369:ALA:HB3	1:B:379:ILE:HG13	2.00	0.42
1:A:495:ASN:HB3	1:A:499:ILE:O	2.19	0.42
1:A:286:ASN:HD21	1:A:549:ASN:HA	1.84	0.42
1:A:589:GLU:HG3	1:A:611:PRO:HD2	2.01	0.42
1:A:453:LEU:HA	1:A:456:LEU:HD12	2.02	0.42
1:A:419:PRO:HA	1:A:420:PRO:HD3	1.87	0.41
1:A:178:LEU:HD22	1:A:188:ALA:HA	2.01	0.41
1:A:490:PRO:CD	1:A:590:LEU:HD11	2.50	0.41
1:B:415:PRO:HD2	1:B:443:ALA:HB2	2.02	0.41
1:A:684:ILE:CG2	1:A:689:ARG:HB3	2.48	0.41
1:A:30:PRO:HG3	1:A:693:LEU:HG	2.02	0.41
1:A:465:PHE:O	1:A:515:ARG:NH2	2.53	0.41
1:A:137:ASP:HB3	1:A:138:ALA:CB	2.42	0.41
1:A:654:ILE:HG22	1:A:655:THR:N	2.35	0.41
1:B:108:PRO:HB2	1:B:219:VAL:HG11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:PRO:HD3	1:A:590:LEU:HD11	2.01	0.41
1:A:4:ARG:HG2	1:A:208:PHE:CG	2.56	0.41
1:A:495:ASN:HD21	1:A:504:LEU:CD1	2.25	0.41
1:B:453:LEU:HA	1:B:456:LEU:HD12	2.02	0.41
1:B:603:LEU:HD21	1:B:669:LEU:HD12	2.03	0.41
1:A:691:GLN:O	1:A:694:ARG:HB2	2.21	0.40
1:B:144:GLU:HA	1:B:146:ASN:HD22	1.85	0.40
1:B:624:ASN:HA	3:B:813:HOH:O	2.22	0.40
1:A:395:VAL:O	1:A:399:VAL:HG23	2.21	0.40
1:A:423:TRP:HA	1:A:423:TRP:CE3	2.57	0.40
1:A:596:HIS:NE2	1:A:607:SER:HB3	2.37	0.40
1:B:521:THR:HA	1:B:587:LEU:HD22	2.03	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	626/698 (90%)	589 (94%)	33 (5%)	4 (1%)	21	52
1	B	669/698 (96%)	616 (92%)	47 (7%)	6 (1%)	14	43
All	All	1295/1396 (93%)	1205 (93%)	80 (6%)	10 (1%)	16	45

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	166	GLU
1	B	183	ASP
1	A	694	ARG
1	B	67	PRO
1	B	144	GLU
1	B	165	ARG

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Mol	Chain	Res	Type
1	B	694	ARG
1	A	139	GLY
1	A	600	ASN
1	A	540	VAL

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	490/566 (87%)	435 (89%)	55 (11%)	6	22
1	B	523/566 (92%)	470 (90%)	53 (10%)	7	27
All	All	1013/1132 (90%)	905 (89%)	108 (11%)	6	24

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASP
1	A	38	GLU
1	A	43	VAL
1	A	48	VAL
1	A	94	SER
1	A	98	THR
1	A	117	ARG
1	A	132	VAL
1	A	137	ASP
1	A	150	LEU
1	A	166	GLU
1	A	168	ARG
1	A	179	ARG
1	A	186	THR
1	A	209	VAL
1	A	220	ASP
1	A	223	GLU
1	A	232	MET
1	A	243	GLU

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Mol	Chain	Res	Type
1	A	289	VAL
1	A	296	VAL
1	A	319	ILE
1	A	343	GLN
1	A	380	TYR
1	A	386	ASN
1	A	418	LYS
1	A	427	ILE
1	A	431	LYS
1	A	434	ASN
1	A	466	THR
1	A	476	GLU
1	A	487	ILE
1	A	496	THR
1	A	498	ASP
1	A	506	HIS
1	A	511	MET
1	A	530	SER
1	A	538	GLN
1	A	566	ARG
1	A	569	SER
1	A	579	GLU
1	A	582	ARG
1	A	588	ARG
1	A	590	LEU
1	A	593	ILE
1	A	599	ASP
1	A	616	THR
1	A	639	GLU
1	A	654	ILE
1	A	658	GLU
1	A	659	TYR
1	A	668	ARG
1	A	679	LEU
1	A	683	LEU
1	A	687	ASP
1	B	8	ASP
1	B	38	GLU
1	B	43	VAL
1	B	48	VAL
1	B	94	SER
1	B	98	THR

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Mol	Chain	Res	Type
1	B	100	ASP
1	B	133	THR
1	B	145	LEU
1	B	146	ASN
1	B	149	LEU
1	B	150	LEU
1	B	151	VAL
1	B	161	THR
1	B	167	LEU
1	B	186	THR
1	B	209	VAL
1	B	220	ASP
1	B	223	GLU
1	B	232	MET
1	B	235	ARG
1	B	277	ILE
1	B	283	LYS
1	B	285	ARG
1	B	289	VAL
1	B	290	THR
1	B	296	VAL
1	B	319	ILE
1	B	343	GLN
1	B	361	LEU
1	B	368	TYR
1	B	370	GLU
1	B	386	ASN
1	B	413	ASP
1	B	414	ASN
1	B	418	LYS
1	B	466	THR
1	B	487	ILE
1	B	498	ASP
1	B	511	MET
1	B	554	GLU
1	B	569	SER
1	B	590	LEU
1	B	603	LEU
1	B	616	THR
1	B	632	THR
1	B	640	LEU
1	B	654	ILE

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Mol	Chain	Res	Type
1	B	658	GLU
1	B	668	ARG
1	B	683	LEU
1	B	687	ASP
1	B	691	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	140	GLN
1	A	287	ASN
1	A	393	GLN
1	A	405	HIS
1	A	408	ASN
1	A	434	ASN
1	A	495	ASN
1	A	505	GLN
1	A	538	GLN
1	A	624	ASN
1	B	97	HIS
1	B	146	ASN
1	B	287	ASN
1	B	343	GLN
1	B	405	HIS
1	B	408	ASN
1	B	506	HIS
1	B	624	ASN
1	B	663	GLN
1	B	680	ASN

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

2 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	C	1	2	12,12,12	1.47	1 (8%)	17,17,17	0.90	0
2	GLC	C	2	2	11,11,12	1.51	2 (18%)	15,15,17	0.87	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	GLC	C	1	2	-	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	1/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	GLC	O5-C1	3.99	1.52	1.42
2	C	2	GLC	O5-C5	3.33	1.49	1.43
2	C	2	GLC	O5-C1	2.64	1.48	1.43

There are no bond angle outliers.

There are no chirality outliers.

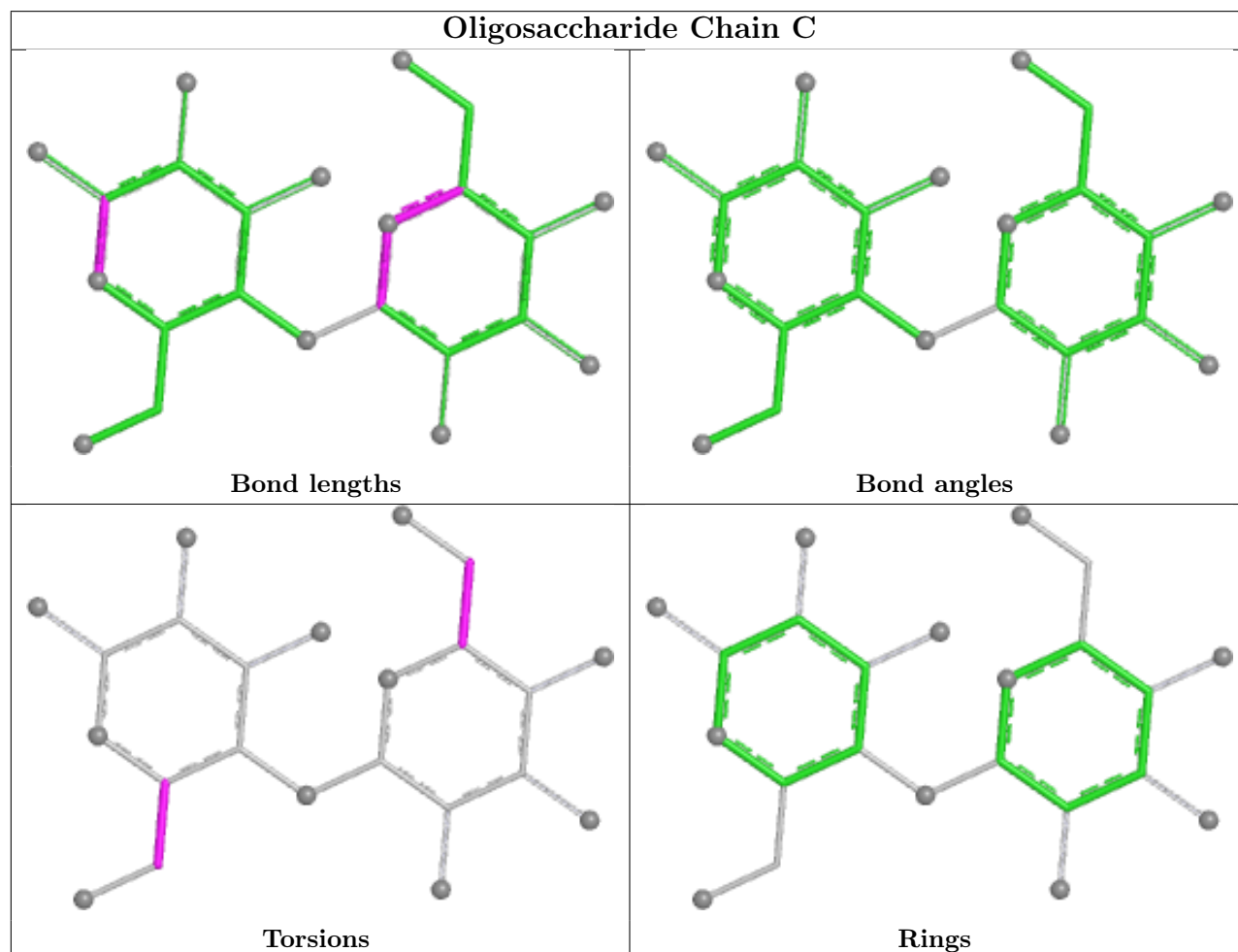
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	GLC	O5-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	C	2	GLC	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](#)

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 [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	636/698 (91%)	-0.36	3 (0%) 87 76	53, 91, 172, 196	0
1	B	675/698 (96%)	-0.08	11 (1%) 70 52	74, 118, 162, 190	0
All	All	1311/1396 (93%)	-0.22	14 (1%) 78 61	53, 107, 168, 196	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	88	PRO	2.9
1	B	667	VAL	2.8
1	B	633	LEU	2.7
1	B	517	VAL	2.7
1	B	675	VAL	2.6
1	A	344	CYS	2.4
1	B	195	GLU	2.4
1	B	661	TRP	2.4
1	B	136	LEU	2.3
1	B	678	VAL	2.2
1	B	631	ALA	2.1
1	A	311	ALA	2.1
1	B	0	SER	2.0
1	B	590	LEU	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.

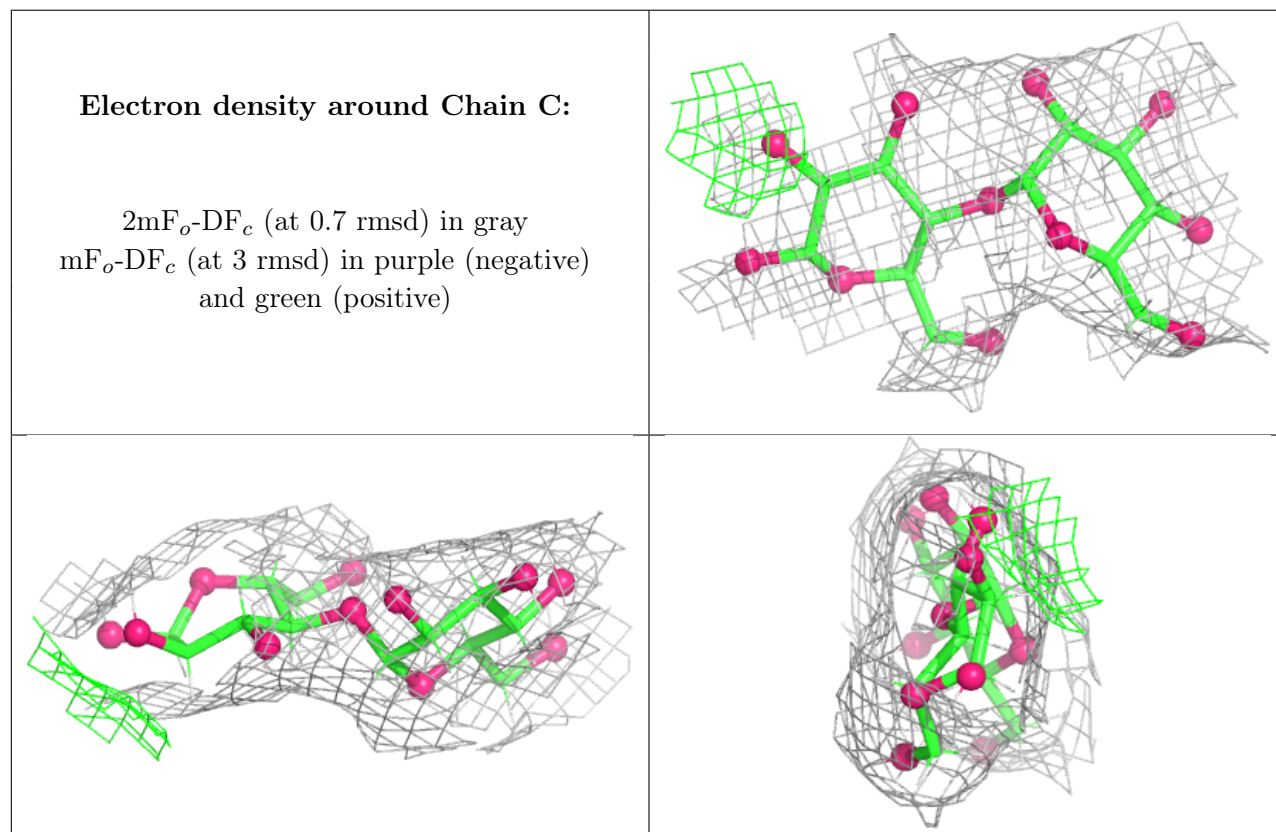
6.3 Carbohydrates [i](#)

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	GLC	C	1	12/12	0.89	0.08	116,124,132,133	0
2	GLC	C	2	11/12	0.94	0.05	112,114,116,119	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.