



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:45 AM UTC

PDB ID : 8IBT / pdb_00008ibt
Title : Crystal structure of GH42 beta-galactosidase BiBga42A from Bifidobacterium longum subspecies infantis E318S mutant in complex with lacto-N-tetraose
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Deposited on : 2023-02-10
Resolution : 2.20 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

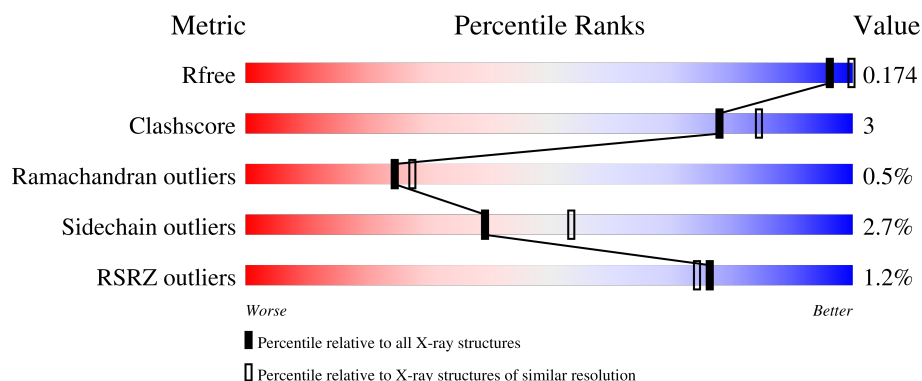
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	702	% 88% 10% ..
1	B	702	% 89% 9% ..
2	C	4	25% 75%
2	D	4	25% 75%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11711 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 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	694	Total	C	N	O	S	0	0	0
			5486	3489	937	1038	22			
1	B	693	Total	C	N	O	S	0	0	0
			5479	3483	936	1038	22			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	318	SER	GLU	engineered mutation	UNP B7GUD7
A	692	ALA	-	expression tag	UNP B7GUD7
A	693	ALA	-	expression tag	UNP B7GUD7
A	694	ALA	-	expression tag	UNP B7GUD7
A	695	LEU	-	expression tag	UNP B7GUD7
A	696	GLU	-	expression tag	UNP B7GUD7
A	697	HIS	-	expression tag	UNP B7GUD7
A	698	HIS	-	expression tag	UNP B7GUD7
A	699	HIS	-	expression tag	UNP B7GUD7
A	700	HIS	-	expression tag	UNP B7GUD7
A	701	HIS	-	expression tag	UNP B7GUD7
A	702	HIS	-	expression tag	UNP B7GUD7
B	318	SER	GLU	engineered mutation	UNP B7GUD7
B	692	ALA	-	expression tag	UNP B7GUD7
B	693	ALA	-	expression tag	UNP B7GUD7
B	694	ALA	-	expression tag	UNP B7GUD7
B	695	LEU	-	expression tag	UNP B7GUD7
B	696	GLU	-	expression tag	UNP B7GUD7
B	697	HIS	-	expression tag	UNP B7GUD7
B	698	HIS	-	expression tag	UNP B7GUD7
B	699	HIS	-	expression tag	UNP B7GUD7
B	700	HIS	-	expression tag	UNP B7GUD7
B	701	HIS	-	expression tag	UNP B7GUD7
B	702	HIS	-	expression tag	UNP B7GUD7

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	4	Total 48	C 26	N 1	O 21	0	0	0
2	D	4	Total 48	C 26	N 1	O 21	0	0	0

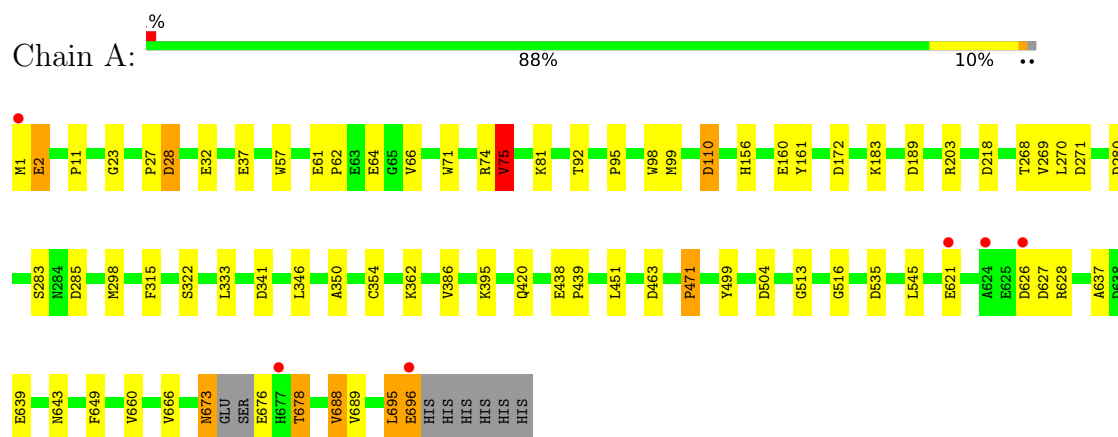
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	346	Total	O	0	0
			346	346		
3	B	304	Total	O	0	0
			304	304		

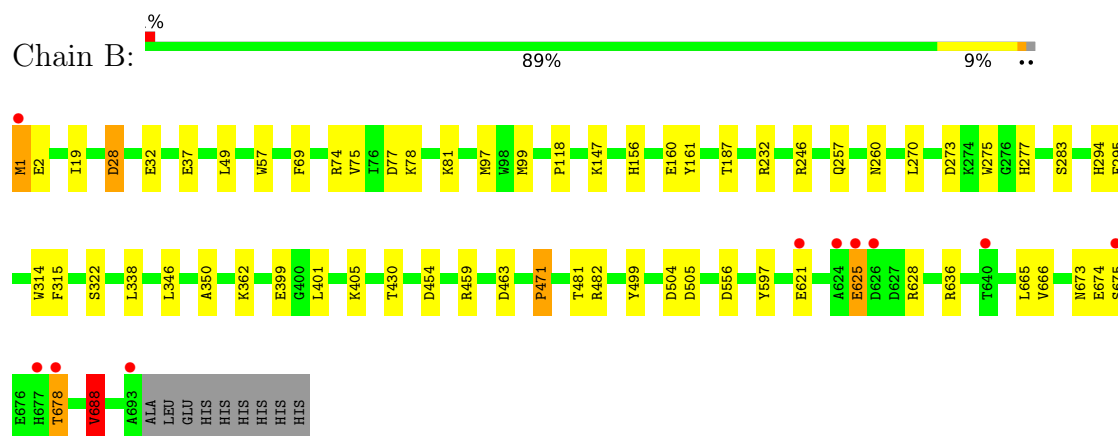
3 Residue-property plots [i](#)

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: Beta-galactosidase



- Molecule 1: Beta-galactosidase



- Molecule 2: beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose



- Molecule 2: beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain D:



8IBT
GAL2
MAG3
GAL4

4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	166.43Å 166.43Å 149.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.04 – 2.20 48.04 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.04-2.20) 99.9 (48.04-2.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.52 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0403	Depositor
R, R_{free}	0.146 , 0.175 0.151 , 0.174	Depositor DCC
R_{free} test set	6036 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	29.2	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 22.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11711	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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: BGC, NAG, GAL

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	0.84	1/5644 (0.0%)	1.21	17/7697 (0.2%)
1	B	0.80	0/5638	1.19	19/7690 (0.2%)
All	All	0.82	1/11282 (0.0%)	1.20	36/15387 (0.2%)

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	A	0	1
1	B	0	3
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	189	ASP	CG-OD1	5.14	1.35	1.25

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	37	GLU	CB-CG-CD	9.63	128.98	112.60
1	A	28	ASP	CA-CB-CG	8.82	121.42	112.60
1	B	28	ASP	CA-CB-CG	8.22	120.82	112.60
1	B	621	GLU	CB-CG-CD	7.89	126.01	112.60
1	A	621	GLU	CB-CG-CD	7.80	125.86	112.60
1	B	273	ASP	CA-CB-CG	-6.99	105.61	112.60
1	B	481	THR	CA-CB-OG1	-6.87	99.30	109.60
1	A	280	ASP	CA-CB-CG	6.82	119.42	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	69	PHE	CA-CB-CG	6.73	120.53	113.80
1	B	295	PHE	CA-CB-CG	-6.53	107.27	113.80
1	A	535	ASP	CB-CA-C	-6.39	96.63	109.99
1	A	696	GLU	CB-CG-CD	6.21	123.15	112.60
1	A	463	ASP	CA-CB-CG	6.14	118.75	112.60
1	B	260	ASN	CB-CA-C	5.98	119.60	109.80
1	A	95	PRO	N-CA-C	5.97	117.99	110.70
1	B	454	ASP	CA-CB-CG	5.97	118.57	112.60
1	B	156	HIS	CA-CB-CG	5.93	119.73	113.80
1	A	218	ASP	CB-CA-C	5.90	119.33	109.89
1	A	75	VAL	N-CA-CB	5.85	117.39	110.55
1	B	621	GLU	CB-CA-C	5.57	118.94	109.80
1	B	688	VAL	CB-CA-C	5.52	118.39	110.33
1	B	37	GLU	CB-CG-CD	5.43	121.83	112.60
1	A	627	ASP	CA-C-N	5.35	131.76	121.54
1	A	627	ASP	C-N-CA	5.35	131.76	121.54
1	B	187	THR	CA-CB-OG1	-5.34	101.59	109.60
1	B	505	ASP	CA-CB-CG	5.30	117.90	112.60
1	B	463	ASP	CA-CB-CG	5.30	117.90	112.60
1	A	75	VAL	CB-CA-C	-5.27	105.22	111.97
1	A	649	PHE	CB-CA-C	-5.27	101.71	110.14
1	A	172	ASP	CA-CB-CG	5.13	117.73	112.60
1	B	471	PRO	N-CA-CB	-5.10	97.89	103.25
1	B	78	LYS	CB-CA-C	5.09	119.51	110.85
1	B	294	HIS	CA-CB-CG	5.09	118.89	113.80
1	B	625	GLU	CB-CG-CD	5.02	121.13	112.60
1	A	678	THR	CA-CB-OG1	-5.02	102.08	109.60
1	A	438	GLU	CB-CG-CD	5.00	121.10	112.60

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	203	ARG	Sidechain
1	B	459	ARG	Sidechain
1	B	482	ARG	Sidechain
1	B	675	SER	Peptide

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	5486	0	5190	38	0
1	B	5479	0	5180	22	0
2	C	48	0	42	0	0
2	D	48	0	41	0	0
3	A	346	0	0	6	0
3	B	304	0	0	1	0
All	All	11711	0	10453	60	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 (60) 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:626:ASP:HB2	3:A:873:HOH:O	1.63	0.98
1:A:298:MET:HE2	3:A:881:HOH:O	1.86	0.75
1:B:32:GLU:CD	1:B:74:ARG:HH12	2.04	0.65
1:A:666:VAL:HB	1:A:688:VAL:HG13	1.83	0.61
1:A:1:MET:O	1:A:2:GLU:CB	2.50	0.60
1:A:673:ASN:ND2	1:A:676:GLU:HB3	2.17	0.59
1:A:673:ASN:HD22	1:A:673:ASN:C	2.10	0.59
1:A:32:GLU:OE1	1:A:74:ARG:NH1	2.36	0.58
1:B:666:VAL:HB	1:B:688:VAL:HG13	1.84	0.58
1:A:673:ASN:ND2	3:A:801:HOH:O	2.37	0.57
1:B:673:ASN:HB3	1:B:678:THR:HG22	1.86	0.57
1:A:626:ASP:CB	3:A:873:HOH:O	2.38	0.54
1:A:283:SER:HA	1:A:315:PHE:O	2.09	0.52
1:B:32:GLU:OE1	1:B:74:ARG:NH1	2.42	0.52
1:A:269:VAL:HG12	1:A:270:LEU:HG	1.92	0.52
1:A:23:GLY:HA2	1:A:354:CYS:HB3	1.92	0.51
1:B:1:MET:HG2	1:B:277:HIS:NE2	2.27	0.49
1:B:504:ASP:OD1	1:B:504:ASP:C	2.55	0.49
1:A:298:MET:HE1	1:A:341:ASP:CG	2.38	0.48
1:B:77:ASP:OD2	1:B:81:LYS:HE2	2.14	0.48
1:B:556:ASP:OD2	1:B:597:TYR:OH	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:O	1:A:2:GLU:HB3	2.15	0.47
1:B:346:LEU:HD12	1:B:350:ALA:O	2.15	0.47
1:A:32:GLU:CD	1:A:74:ARG:HH12	2.21	0.46
1:B:160:GLU:O	1:B:161:TYR:C	2.58	0.46
1:A:676:GLU:N	3:A:811:HOH:O	2.49	0.45
1:A:545:LEU:HD12	1:A:545:LEU:N	2.30	0.45
1:B:270:LEU:HD13	1:B:275:TRP:CE2	2.53	0.44
1:A:57:TRP:CZ2	1:A:99:MET:HE1	2.53	0.44
1:A:513:GLY:O	1:A:516:GLY:HA2	2.18	0.44
1:B:57:TRP:CZ2	1:B:99:MET:HE1	2.51	0.44
1:A:298:MET:HE1	1:A:341:ASP:HB3	2.00	0.44
1:A:61:GLU:HA	1:A:66:VAL:O	2.19	0.43
1:A:504:ASP:OD1	1:A:504:ASP:C	2.62	0.43
1:A:160:GLU:OE2	1:A:285:ASP:OD2	2.36	0.43
1:A:268:THR:HG21	1:A:420:GLN:HG2	2.01	0.43
1:B:49:LEU:HD23	1:B:49:LEU:C	2.44	0.43
1:A:71:TRP:O	1:A:75:VAL:HG23	2.18	0.42
1:A:439:PRO:HD3	1:A:471:PRO:HB2	2.02	0.42
1:A:64:GLU:HB2	1:A:98:TRP:CZ3	2.54	0.42
1:A:62:PRO:HD2	1:A:66:VAL:O	2.19	0.42
1:A:346:LEU:HD12	1:A:350:ALA:O	2.20	0.42
1:A:27:PRO:HG3	1:A:75:VAL:HG11	2.02	0.41
1:A:637:ALA:HA	1:A:643:ASN:O	2.19	0.41
1:B:338:LEU:HD23	1:B:338:LEU:C	2.45	0.41
1:A:160:GLU:O	1:A:161:TYR:C	2.61	0.41
1:B:399:GLU:HG3	1:B:665:LEU:HD13	2.02	0.41
1:A:110:ASP:OD1	1:A:110:ASP:C	2.63	0.41
1:B:97:MET:HE1	1:B:118:PRO:HG2	2.02	0.41
1:A:271:ASP:C	1:A:271:ASP:OD1	2.63	0.41
1:B:232:ARG:HD3	3:B:888:HOH:O	2.21	0.41
1:B:283:SER:HA	1:B:315:PHE:O	2.21	0.41
1:A:11:PRO:HD3	3:A:951:HOH:O	2.21	0.41
1:B:19:ILE:HD11	1:B:401:LEU:HD23	2.02	0.40
1:B:405:LYS:O	1:B:636:ARG:HA	2.21	0.40
1:A:666:VAL:HB	1:A:688:VAL:CG1	2.48	0.40
1:A:695:LEU:O	1:A:696:GLU:C	2.65	0.40
1:B:314:TRP:O	1:B:350:ALA:HA	2.22	0.40
1:B:246:ARG:HG3	1:B:257:GLN:CD	2.47	0.40
1:A:57:TRP:CE3	1:A:92:THR:HA	2.56	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	690/702 (98%)	669 (97%)	17 (2%)	4 (1%)	21	23
1	B	691/702 (98%)	664 (96%)	24 (4%)	3 (0%)	30	34
All	All	1381/1404 (98%)	1333 (96%)	41 (3%)	7 (0%)	24	27

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2	GLU
1	A	628	ARG
1	A	2	GLU
1	A	322	SER
1	A	471	PRO
1	B	322	SER
1	B	471	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	568/576 (99%)	549 (97%)	19 (3%)	33	45
1	B	568/576 (99%)	556 (98%)	12 (2%)	47	63
All	All	1136/1152 (99%)	1105 (97%)	31 (3%)	39	53

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	75	VAL
1	A	81	LYS
1	A	110	ASP
1	A	156	HIS
1	A	183	LYS
1	A	333	LEU
1	A	362	LYS
1	A	386	VAL
1	A	395	LYS
1	A	451	LEU
1	A	499	TYR
1	A	639	GLU
1	A	660	VAL
1	A	673	ASN
1	A	678	THR
1	A	688	VAL
1	A	689	VAL
1	A	695	LEU
1	B	1	MET
1	B	28	ASP
1	B	75	VAL
1	B	147	LYS
1	B	362	LYS
1	B	430	THR
1	B	499	TYR
1	B	625	GLU
1	B	628	ARG
1	B	674	GLU
1	B	678	THR
1	B	688	VAL

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	490	ASN
1	A	673	ASN
1	B	26	ASN
1	B	117	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 ⓘ

8 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	0.54	0	17,17,17	1.01	0
2	GAL	C	2	2	11,11,12	1.17	1 (9%)	15,15,17	1.41	4 (26%)
2	NAG	C	3	2	14,14,15	0.65	0	17,19,21	1.79	3 (17%)
2	GAL	C	4	2	11,11,12	0.86	1 (9%)	15,15,17	1.69	5 (33%)
2	BGC	D	1	2	12,12,12	0.50	0	17,17,17	0.77	0
2	GAL	D	2	2	11,11,12	1.08	1 (9%)	15,15,17	1.31	2 (13%)
2	NAG	D	3	2	14,14,15	0.69	1 (7%)	17,19,21	1.50	3 (17%)
2	GAL	D	4	2	11,11,12	1.13	1 (9%)	15,15,17	1.89	6 (40%)

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	GAL	C	2	2	-	2/2/19/22	0/1/1/1
2	NAG	C	3	2	-	1/6/23/26	0/1/1/1
2	GAL	C	4	2	-	0/2/19/22	0/1/1/1
2	BGC	D	1	2	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GAL	D	2	2	-	1/2/19/22	0/1/1/1
2	NAG	D	3	2	-	1/6/23/26	0/1/1/1
2	GAL	D	4	2	-	0/2/19/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	GAL	C2-C3	-3.37	1.47	1.52
2	D	2	GAL	O5-C5	2.67	1.48	1.43
2	D	4	GAL	C2-C3	2.39	1.56	1.52
2	C	4	GAL	C2-C3	2.11	1.55	1.52
2	D	3	NAG	C1-C2	2.01	1.55	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	NAG	C1-C2-N2	4.67	117.80	110.43
2	D	4	GAL	C1-C2-C3	3.85	115.25	109.64
2	D	3	NAG	C1-C2-N2	3.68	116.23	110.43
2	C	3	NAG	O3-C3-C2	-3.64	101.84	109.40
2	C	4	GAL	C1-C2-C3	3.15	114.23	109.64
2	D	3	NAG	C2-N2-C7	3.10	127.05	122.90
2	D	4	GAL	O4-C4-C3	-3.02	103.26	110.38
2	C	2	GAL	O3-C3-C2	-2.94	104.05	110.05
2	C	3	NAG	C2-N2-C7	2.71	126.53	122.90
2	D	4	GAL	O3-C3-C4	-2.51	104.45	110.38
2	C	4	GAL	O3-C3-C4	-2.46	104.57	110.38
2	C	2	GAL	O2-C2-C1	2.45	114.83	109.22
2	C	4	GAL	O5-C5-C4	-2.32	105.17	110.83
2	D	2	GAL	C1-C2-C3	-2.32	106.26	109.64
2	D	3	NAG	O3-C3-C2	-2.30	104.62	109.40
2	D	4	GAL	O5-C5-C4	-2.25	105.35	110.83
2	C	2	GAL	C2-C3-C4	-2.21	106.97	110.86
2	C	4	GAL	C2-C3-C4	2.11	114.57	110.86
2	D	4	GAL	O6-C6-C5	2.11	118.50	111.33
2	D	2	GAL	O2-C2-C1	2.09	114.01	109.22
2	C	2	GAL	C1-O5-C5	2.05	114.94	112.19
2	C	4	GAL	O4-C4-C3	-2.03	105.60	110.38
2	D	4	GAL	C6-C5-C4	2.01	117.95	113.02

There are no chirality outliers.

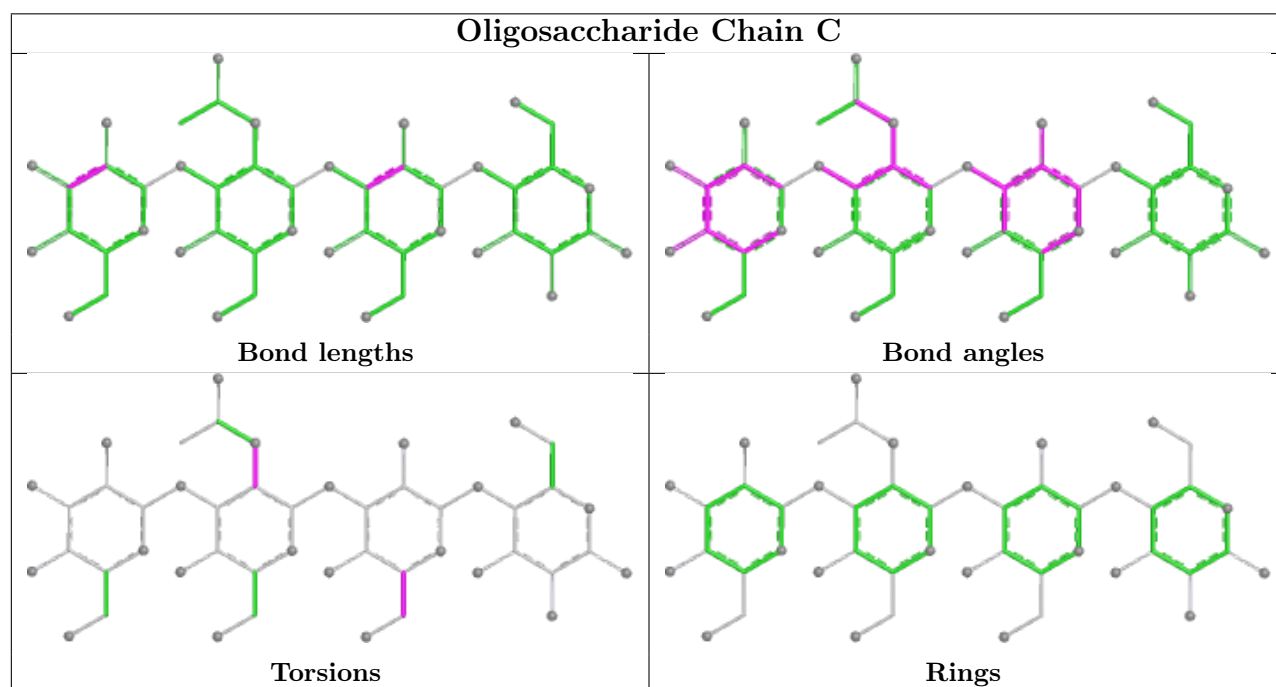
All (5) torsion outliers are listed below:

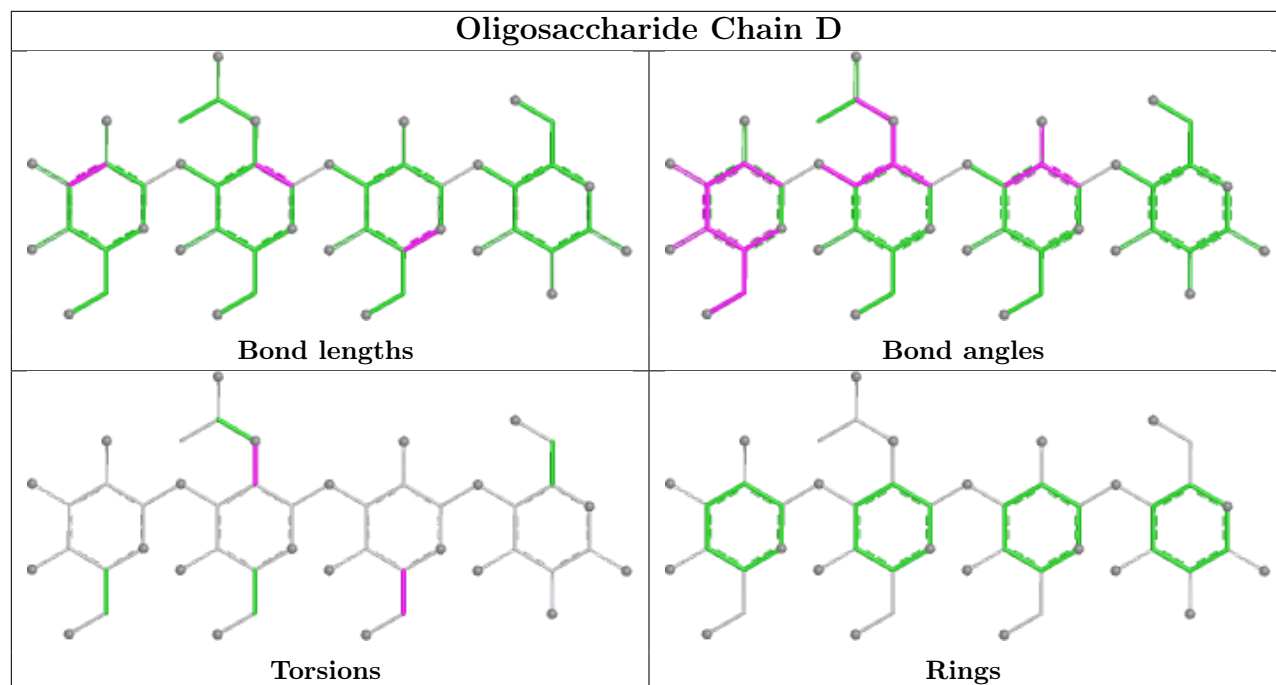
Mol	Chain	Res	Type	Atoms
2	C	3	NAG	C1-C2-N2-C7
2	D	3	NAG	C1-C2-N2-C7
2	C	2	GAL	O5-C5-C6-O6
2	C	2	GAL	C4-C5-C6-O6
2	D	2	GAL	O5-C5-C6-O6

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	694/702 (98%)	-0.48	6 (0%) 81 79	18, 26, 46, 78	0
1	B	693/702 (98%)	-0.40	10 (1%) 73 71	19, 29, 49, 83	0
All	All	1387/1404 (98%)	-0.44	16 (1%) 76 74	18, 28, 48, 83	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	6.2
1	B	1	MET	4.0
1	A	696	GLU	3.6
1	B	693	ALA	3.4
1	B	626	ASP	3.4
1	B	621	GLU	3.0
1	A	621	GLU	2.7
1	A	624	ALA	2.6
1	B	625	GLU	2.6
1	B	677	HIS	2.6
1	B	640	THR	2.5
1	A	677	HIS	2.5
1	A	626	ASP	2.3
1	B	675	SER	2.3
1	B	678	THR	2.2
1	B	624	ALA	2.0

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

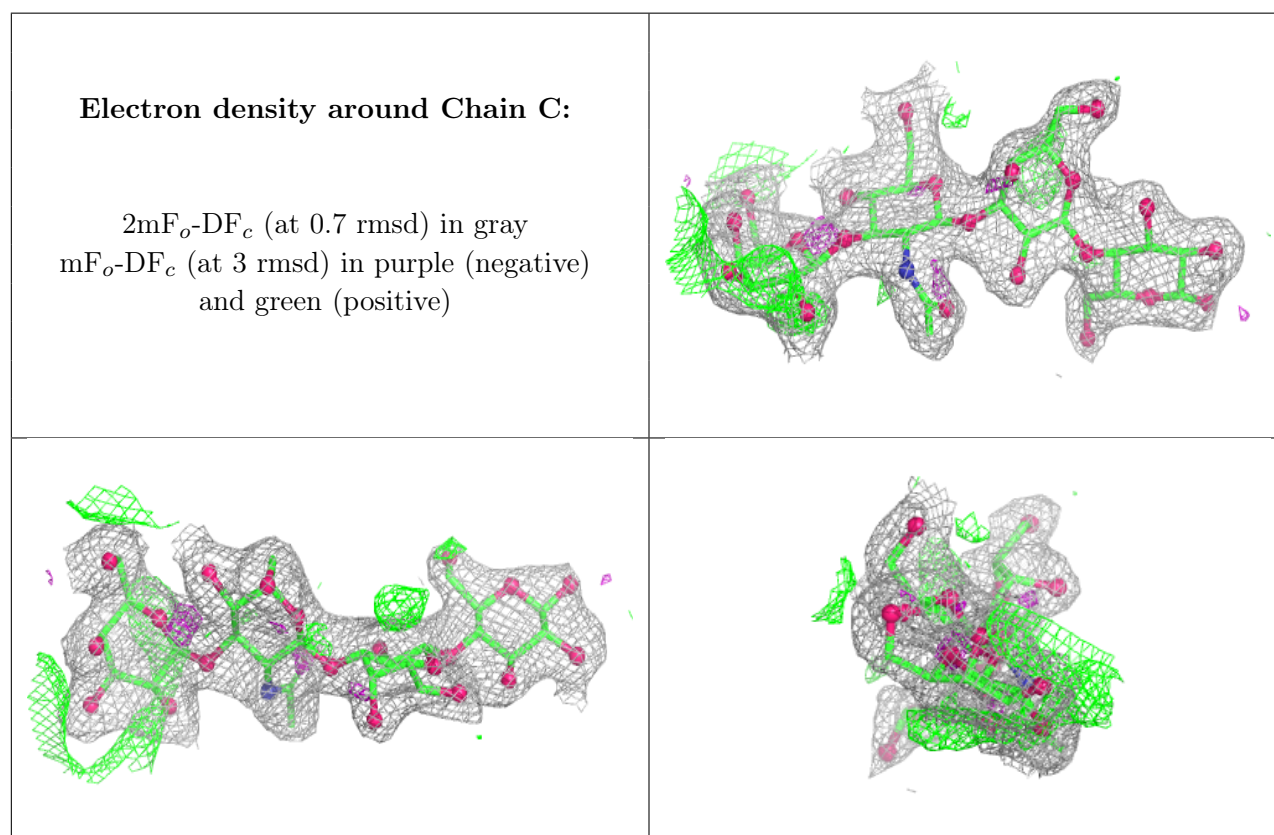
There are no non-standard protein/DNA/RNA residues in this entry.

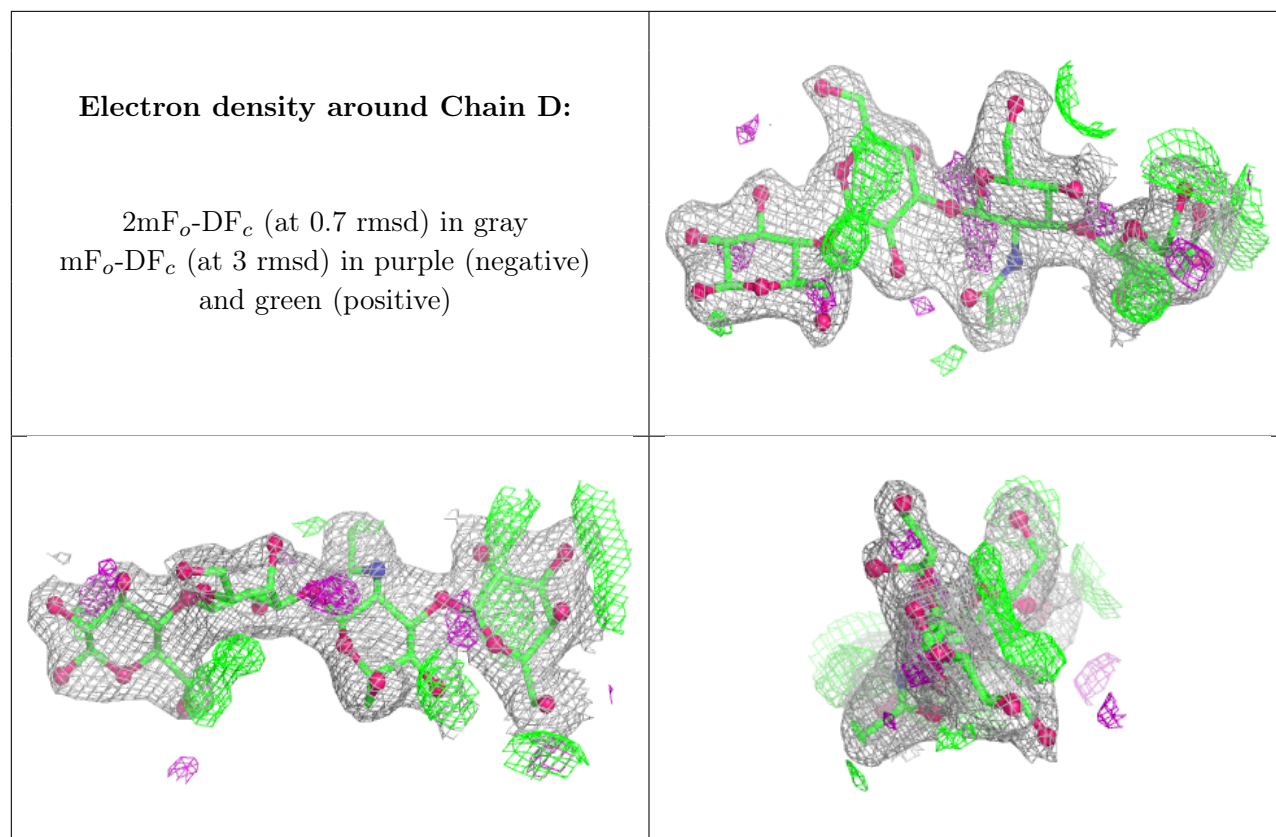
6.3 Carbohydrates [i](#)

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($\text{\AA}^2$)	Q<0.9
2	GAL	C	2	11/12	0.88	0.14	35,49,61,67	0
2	BGC	D	1	12/12	0.89	0.12	39,50,59,65	0
2	GAL	D	2	11/12	0.90	0.13	40,56,65,70	0
2	BGC	C	1	12/12	0.91	0.11	38,50,56,62	0
2	GAL	C	4	11/12	0.91	0.10	27,31,41,43	0
2	NAG	D	3	14/15	0.91	0.11	29,39,44,46	0
2	GAL	D	4	11/12	0.91	0.12	28,37,43,45	0
2	NAG	C	3	14/15	0.93	0.10	30,39,43,47	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.