



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:12 AM UTC

PDB ID : 2YGM / pdb_00002ygm
Title : THE X-RAY CRYSTAL STRUCTURE OF TANDEM CBM51 MODULES OF SP3GH98, THE FAMILY 98 GLYCOSIDE HYDROLASE FROM STREPTOCOCCUS PNEUMONIAE SP3-BS71, IN COMPLEX WITH THE BLOOD GROUP B ANTIGEN
Authors : Higgins, M.A.; Ficko-Blean, E.; Wright, C.; Meloncelli, P.J.; Lowary, T.L.; Boraston, A.B.
Deposited on : 2011-04-19
Resolution : 2.35 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)

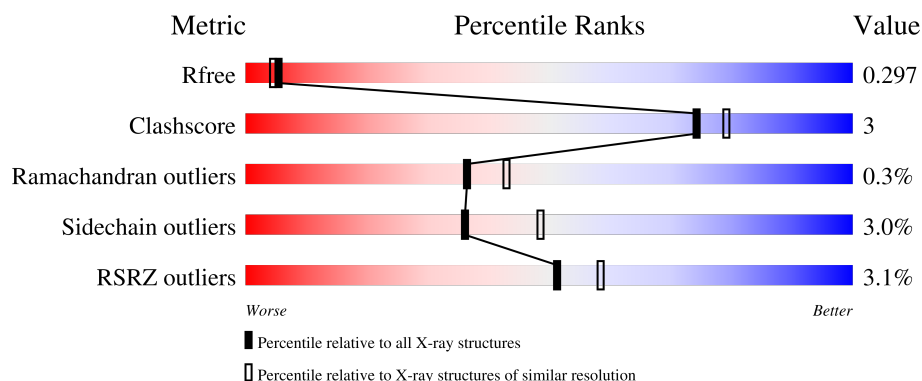
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	354	3% 89% 10% ..
1	B	354	3% 90% 9% ..
2	C	4	50% 50%
2	D	4	25% 75%

Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5697 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.

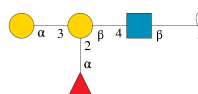
- Molecule 1 is a protein called BLOOD GROUP A-AND B-CLEAVING ENDO-BETA-GALACTOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	1	0
			2712	1707	447	555	3			
1	B	352	Total	C	N	O	S	0	0	0
			2712	1706	447	556	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	GLY	-	expression tag	UNP C1CB04
A	61	SER	-	expression tag	UNP C1CB04
A	62	HIS	-	expression tag	UNP C1CB04
A	63	MET	-	expression tag	UNP C1CB04
A	64	ALA	-	expression tag	UNP C1CB04
A	65	SER	-	expression tag	UNP C1CB04
B	60	GLY	-	expression tag	UNP C1CB04
B	61	SER	-	expression tag	UNP C1CB04
B	62	HIS	-	expression tag	UNP C1CB04
B	63	MET	-	expression tag	UNP C1CB04
B	64	ALA	-	expression tag	UNP C1CB04
B	65	SER	-	expression tag	UNP C1CB04

- Molecule 2 is an oligosaccharide called alpha-L-fucopyranose-(1-2)-[alpha-D-galactopyranose-(1-3)]beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	4	Total	C	N	O	0	0	0
			47	26	1	20			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	4	Total	C	N	O	0	0	0
			47	26	1	20			

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

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

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		
4	B	1	Total	Na	0	0
			1	1		

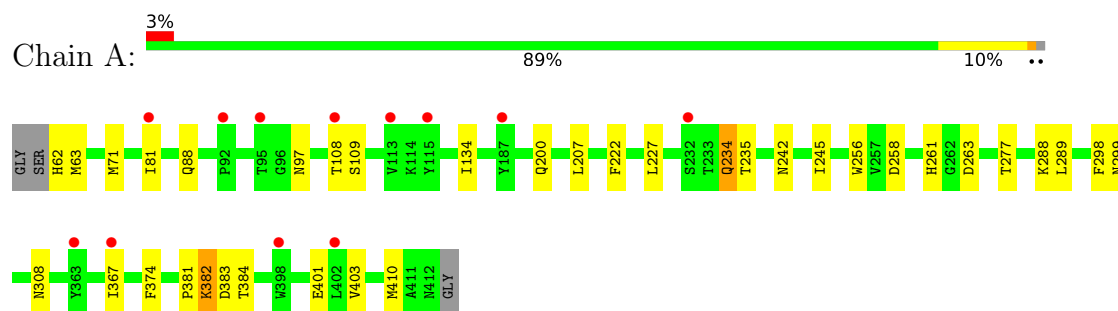
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	98	Total	O	0	0
			98	98		
5	B	77	Total	O	0	0
			77	77		

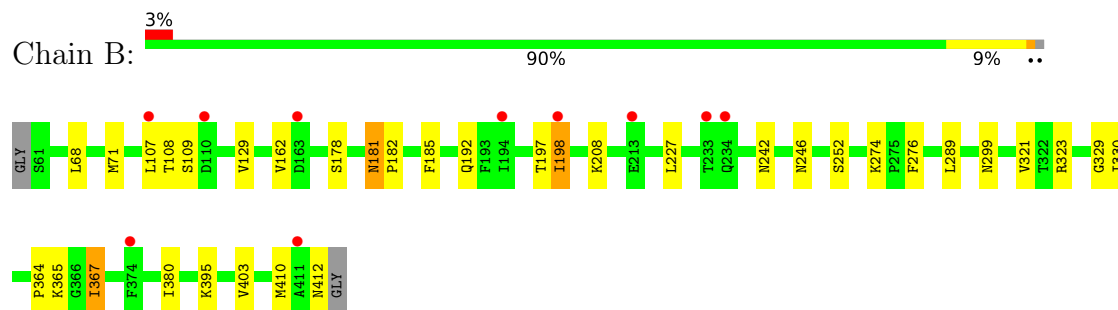
3 Residue-property plots [i](#)

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: BLOOD GROUP A-AND B-CLEAVING ENDO-BETA-GALACTOSIDASE



- Molecule 1: BLOOD GROUP A-AND B-CLEAVING ENDO-BETA-GALACTOSIDASE



- Molecule 2: alpha-L-fucopyranose-(1-2)-[alpha-D-galactopyranose-(1-3)]beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2: alpha-L-fucopyranose-(1-2)-[alpha-D-galactopyranose-(1-3)]beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-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	75.70Å 78.50Å 131.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.35 40.00 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-2.35) 99.9 (40.00-2.35)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.85 (at 2.36Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.245 , 0.297 0.245 , 0.297	Depositor DCC
R_{free} test set	1690 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	46.4	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 24.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.027 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5697	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GLA, NA, CA, FUC, 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	0.44	0/2768	0.71	0/3748
1	B	0.45	1/2765 (0.0%)	0.71	2/3744 (0.1%)
All	All	0.44	1/5533 (0.0%)	0.71	2/7492 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	198	ILE	CA-CB	5.02	1.58	1.54

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	181	ASN	CA-C-N	5.50	124.96	119.24
1	B	181	ASN	C-N-CA	5.50	124.96	119.24

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2712	0	2644	18	0
1	B	2712	0	2641	16	0
2	C	47	0	42	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	47	0	42	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	98	0	0	1	0
5	B	77	0	0	0	0
All	All	5697	0	5369	34	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 (34) 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:382:LYS:HD3	1:A:383:ASP:H	1.28	0.94
1:B:68:LEU:HA	1:B:71:MET:HE3	1.67	0.76
1:A:245:ILE:HG21	1:A:410:MET:HE1	1.68	0.75
1:B:108:THR:HG22	1:B:109:SER:H	1.52	0.73
1:B:181:ASN:HD21	1:B:192:GLN:HG3	1.55	0.71
1:A:71:MET:HE1	1:A:227:LEU:HD12	1.72	0.70
1:A:382:LYS:HD3	1:A:383:ASP:N	2.07	0.68
1:A:382:LYS:HA	1:A:382:LYS:HE2	1.81	0.62
1:A:108:THR:HG22	1:A:109:SER:H	1.67	0.59
1:A:382:LYS:CD	1:A:383:ASP:H	2.10	0.58
1:A:367:ILE:HG21	1:A:374:PHE:HE2	1.72	0.54
1:B:330:ILE:HG13	1:B:367:ILE:HD11	1.90	0.52
1:B:181:ASN:ND2	1:B:192:GLN:HG3	2.25	0.49
1:A:261:HIS:HD2	1:A:263:ASP:H	1.60	0.49
1:B:242:ASN:ND2	1:B:299:ASN:H	2.11	0.49
1:B:289:LEU:HD21	1:B:403:VAL:HG23	1.93	0.49
1:A:234:GLN:NE2	5:A:602:HOH:O	2.49	0.46
1:A:207:LEU:HD23	1:A:222:PHE:CG	2.52	0.45
1:A:242:ASN:ND2	1:A:299:ASN:H	2.15	0.45
1:A:242:ASN:HD21	1:A:298:PHE:HA	1.82	0.45
1:B:323:ARG:HB2	1:B:410:MET:HB2	1.99	0.45
1:A:277:THR:OG1	1:A:401:GLU:OE2	2.28	0.44
1:A:289:LEU:HD21	1:A:403:VAL:HG23	1.99	0.44
1:B:329:GLY:HA2	1:B:367:ILE:HD12	1.98	0.44
1:B:162:VAL:O	1:B:185:PHE:N	2.51	0.44
1:B:178:SER:O	1:B:182:PRO:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:MET:HE1	1:B:227:LEU:HD12	2.00	0.43
1:A:256:TRP:CD1	1:A:258:ASP:HA	2.53	0.43
1:B:129:VAL:HG22	1:B:208:LYS:HG2	1.99	0.43
1:B:108:THR:HG22	1:B:109:SER:N	2.27	0.42
1:A:88:GLN:HE21	1:A:97:ASN:CG	2.29	0.41
1:A:381:PRO:O	1:A:384:THR:OG1	2.25	0.40
1:B:364:PRO:HG2	1:B:365:LYS:HE2	2.02	0.40
1:B:321:VAL:HG12	1:B:380:ILE:HG13	2.02	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	350/354 (99%)	334 (95%)	15 (4%)	1 (0%)	36	43
1	B	350/354 (99%)	337 (96%)	12 (3%)	1 (0%)	36	43
All	All	700/708 (99%)	671 (96%)	27 (4%)	2 (0%)	36	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	276	PHE
1	A	308	ASN

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	303/303 (100%)	294 (97%)	9 (3%)	36	48
1	B	303/303 (100%)	294 (97%)	9 (3%)	36	48
All	All	606/606 (100%)	588 (97%)	18 (3%)	36	48

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	HIS
1	A	63	MET
1	A	81	ILE
1	A	134	ILE
1	A	200	GLN
1	A	234	GLN
1	A	235	THR
1	A	288	LYS
1	A	382	LYS
1	B	107	LEU
1	B	197	THR
1	B	198	ILE
1	B	246	ASN
1	B	252	SER
1	B	274	LYS
1	B	367	ILE
1	B	395	LYS
1	B	412	ASN

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	62	HIS
1	A	88	GLN
1	A	150	GLN
1	A	181	ASN
1	A	201	ASN
1	A	234	GLN
1	A	242	ASN
1	A	261	HIS
1	A	308	ASN
1	A	335	ASN

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Mol	Chain	Res	Type
1	A	336	HIS
1	A	396	HIS
1	A	412	ASN
1	B	153	ASN
1	B	181	ASN
1	B	214	HIS
1	B	242	ASN
1	B	261	HIS
1	B	335	ASN
1	B	412	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 ⓘ

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	NAG	C	1	2	15,15,15	0.45	0	21,21,21	1.15	3 (14%)
2	GAL	C	2	2	11,11,12	0.71	0	15,15,17	1.00	0
2	FUC	C	3	2	10,10,11	0.58	0	14,14,16	0.56	0
2	GLA	C	4	2	11,11,12	0.49	0	15,15,17	1.32	1 (6%)
2	NAG	D	1	2	15,15,15	0.44	0	21,21,21	0.97	1 (4%)
2	GAL	D	2	2	11,11,12	0.61	0	15,15,17	0.99	2 (13%)
2	FUC	D	3	2	10,10,11	0.66	0	14,14,16	0.77	0
2	GLA	D	4	2	11,11,12	0.55	0	15,15,17	0.70	1 (6%)

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	NAG	C	1	2	-	0/6/26/26	0/1/1/1
2	GAL	C	2	2	-	0/2/19/22	0/1/1/1
2	FUC	C	3	2	-	-	0/1/1/1
2	GLA	C	4	2	-	0/2/19/22	0/1/1/1
2	NAG	D	1	2	-	0/6/26/26	0/1/1/1
2	GAL	D	2	2	-	0/2/19/22	0/1/1/1
2	FUC	D	3	2	-	-	0/1/1/1
2	GLA	D	4	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4	GLA	C1-O5-C5	4.68	118.46	112.19
2	C	1	NAG	C4-C3-C2	2.82	114.51	110.40
2	C	1	NAG	C3-C4-C5	2.40	114.59	110.23
2	D	1	NAG	C4-C3-C2	2.32	113.78	110.40
2	D	4	GLA	C1-O5-C5	2.28	115.24	112.19
2	D	2	GAL	C1-C2-C3	2.19	112.84	109.64
2	D	2	GAL	O3-C3-C2	-2.15	105.66	110.05
2	C	1	NAG	O5-C5-C4	2.05	113.39	109.70

There are no chirality outliers.

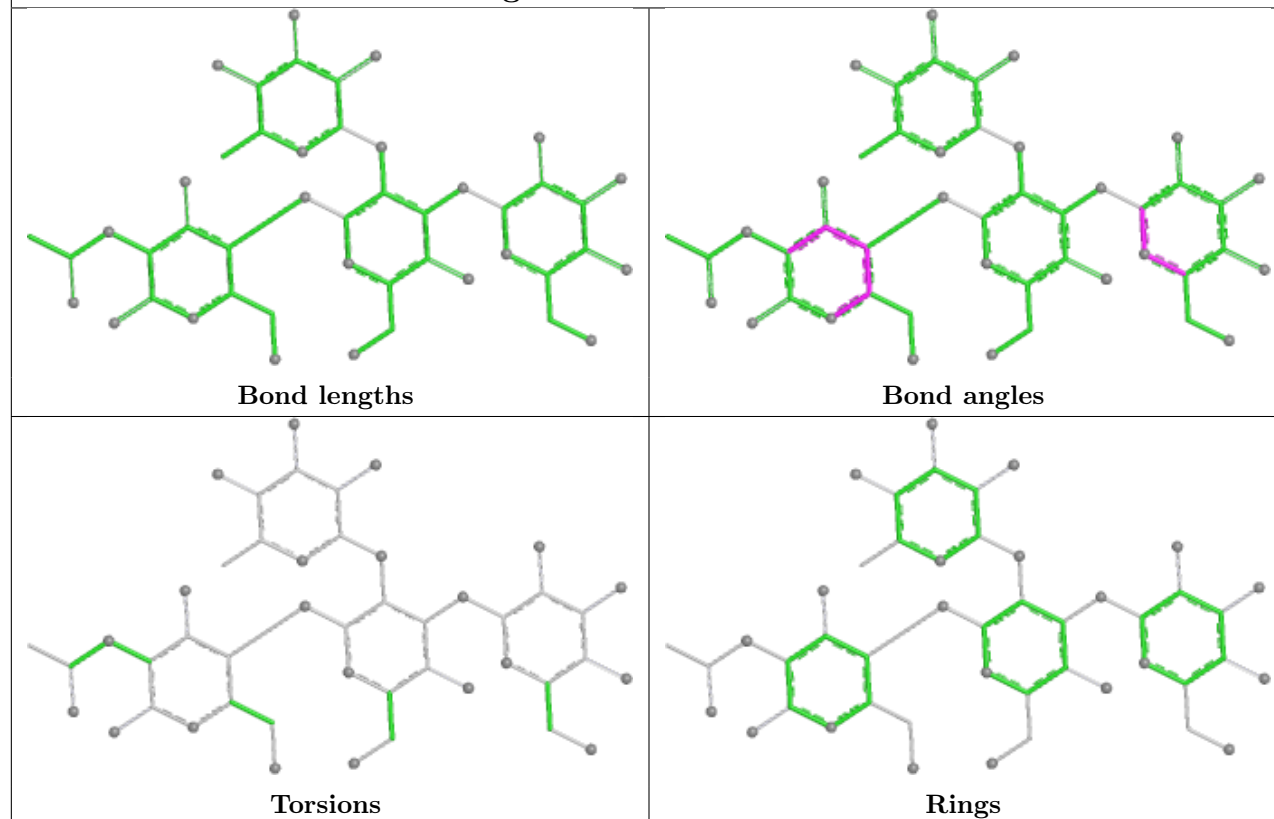
There are no torsion outliers.

There are no ring outliers.

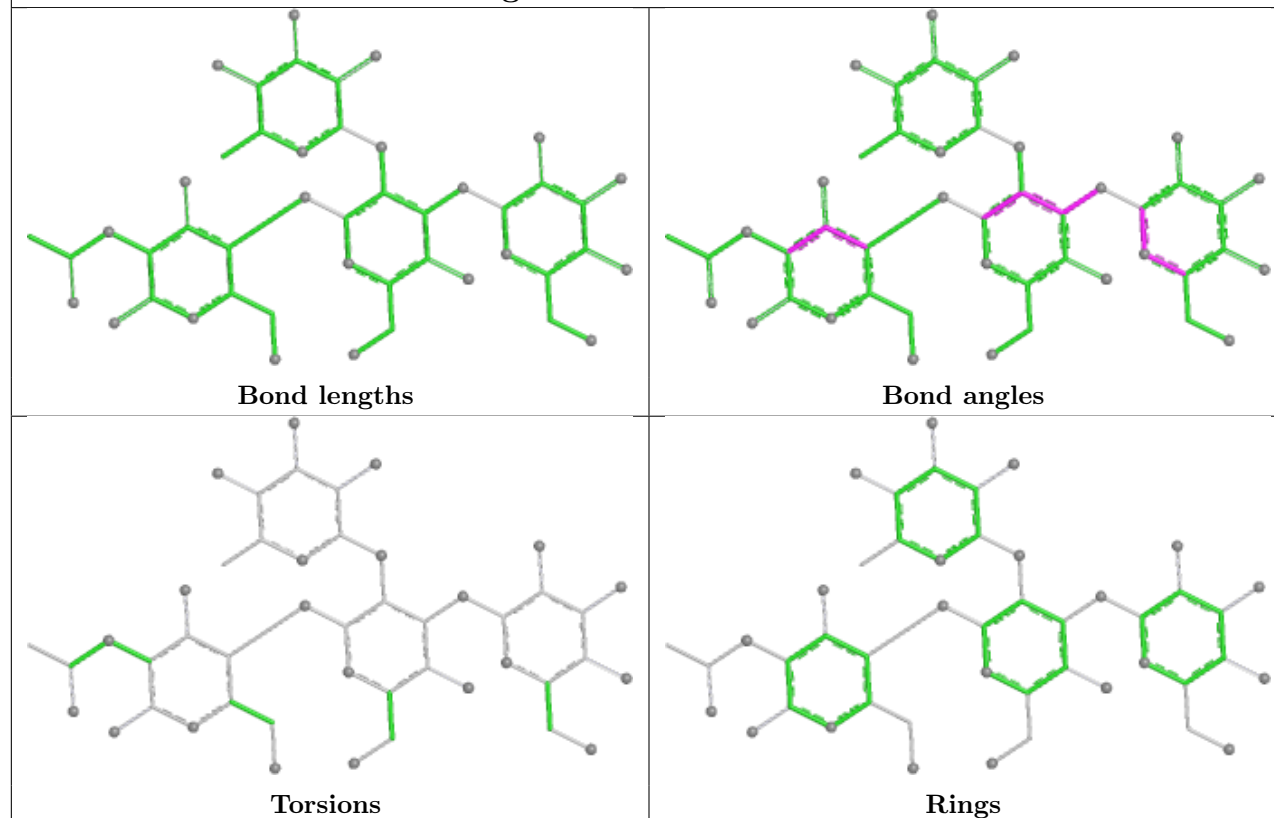
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain C



Oligosaccharide Chain D



5.6 Ligand geometry

Of 4 ligands modelled in this entry, 4 are 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

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	351/354 (99%)	0.42	12 (3%)	48 55	21, 45, 62, 67	1 (0%)
1	B	352/354 (99%)	0.47	10 (2%)	55 61	33, 47, 58, 61	0
All	All	703/708 (99%)	0.45	22 (3%)	51 58	21, 46, 59, 67	1 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	398	TRP	4.8
1	B	233	THR	3.5
1	B	411	ALA	3.1
1	A	115	TYR	2.6
1	B	234	GLN	2.6
1	A	113	VAL	2.5
1	A	367	ILE	2.5
1	B	107	LEU	2.5
1	B	163	ASP	2.4
1	B	110	ASP	2.3
1	A	187	TYR	2.3
1	B	213	GLU	2.3
1	A	402	LEU	2.3
1	B	198	ILE	2.2
1	A	108	THR	2.2
1	A	81	ILE	2.2
1	B	194	ILE	2.2
1	B	374	PHE	2.2
1	A	95	THR	2.1
1	A	363	TYR	2.1
1	A	232	SER	2.0
1	A	92	PRO	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 ⓘ

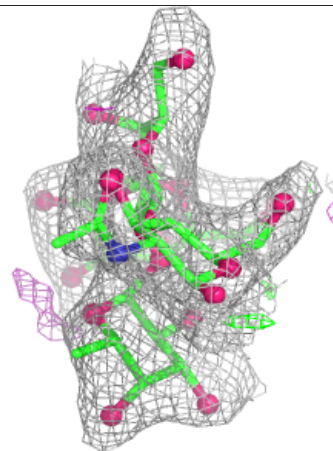
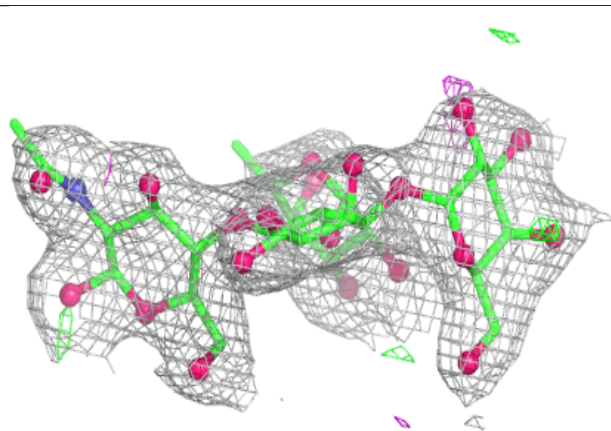
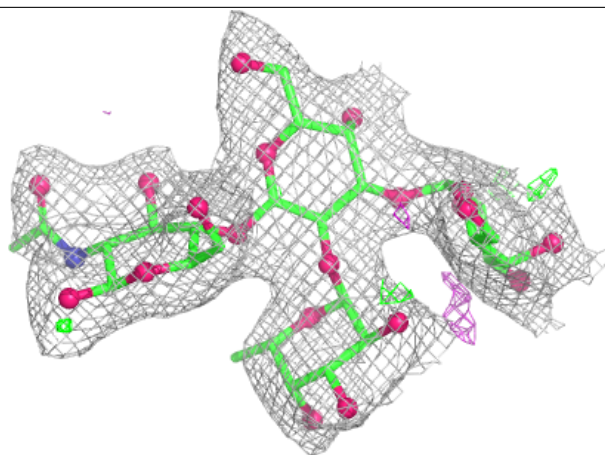
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

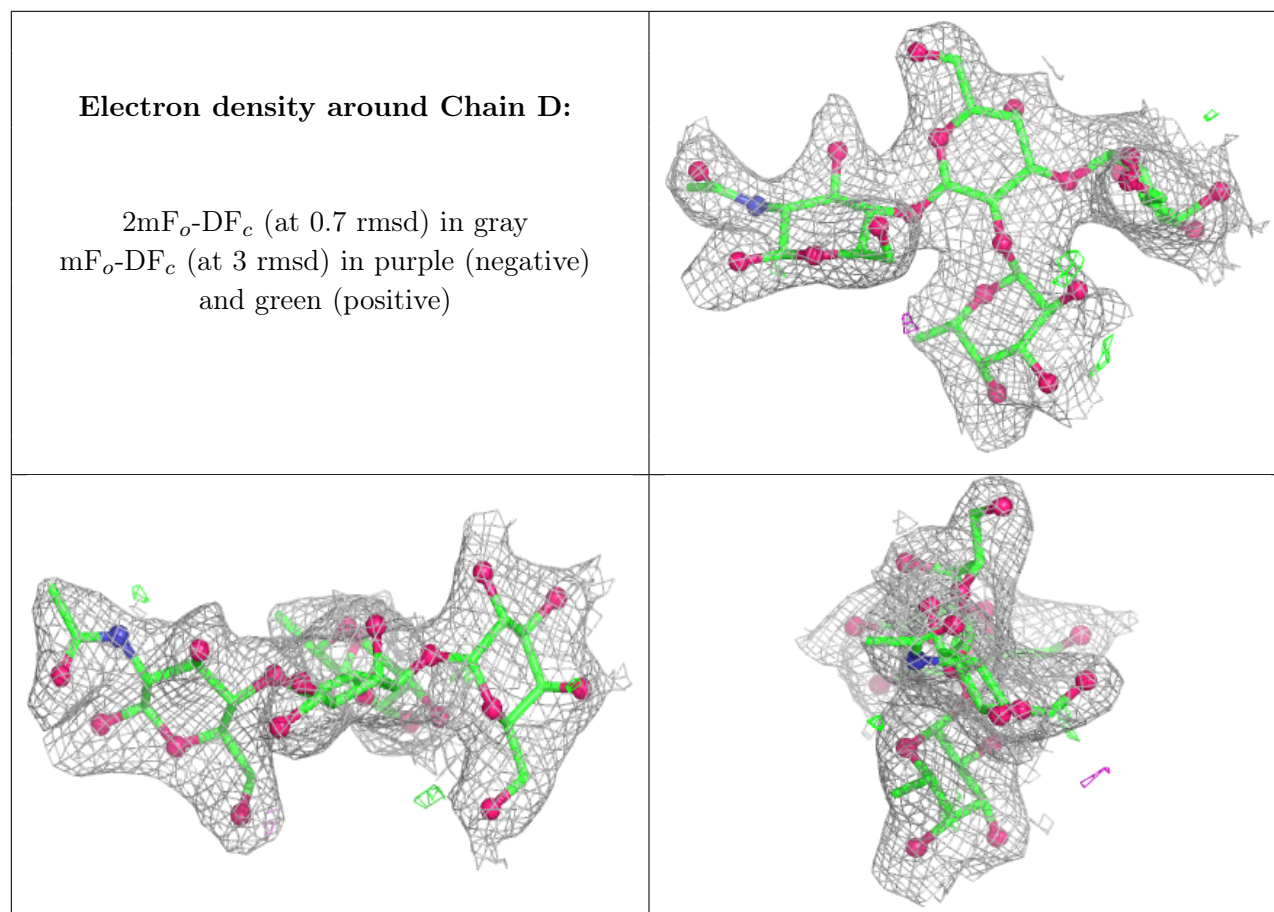
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	C	1	15/15	0.83	0.12	58,59,60,60	0
2	NAG	D	1	15/15	0.84	0.11	53,55,55,55	0
2	FUC	C	3	10/11	0.88	0.10	57,58,58,58	0
2	FUC	D	3	10/11	0.88	0.10	50,51,51,51	0
2	GLA	C	4	11/12	0.89	0.10	53,53,54,55	0
2	GAL	C	2	11/12	0.90	0.09	55,56,56,57	0
2	GAL	D	2	11/12	0.92	0.09	51,52,53,53	0
2	GLA	D	4	11/12	0.94	0.08	50,50,51,51	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 C:

$2mF_o - DF_c$ (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
4	NA	A	506	1/1	0.88	0.11	43,43,43,43	0
3	CA	B	506	1/1	0.98	0.08	53,53,53,53	0
4	NA	B	505	1/1	0.98	0.13	34,34,34,34	0
3	CA	A	505	1/1	0.99	0.02	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.