



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 08:49 PM UTC

PDB ID : 7KR6 / pdb_00007kr6
Title : Glycoside hydrolase family 16 endo-glucanase from Bacteroides ovatus in complex with G4G3G-2F-DNP
Authors : Tamura, K.; Brumer, H.; van Petegem, F.
Deposited on : 2020-11-18
Resolution : 1.56 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

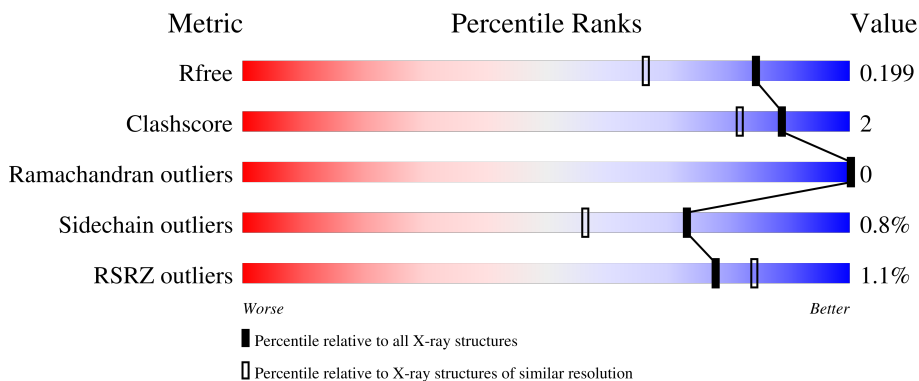
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


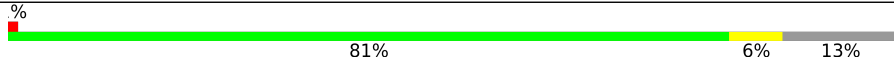
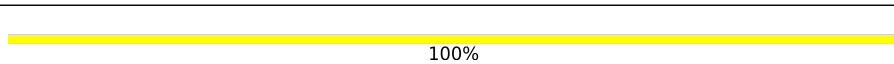
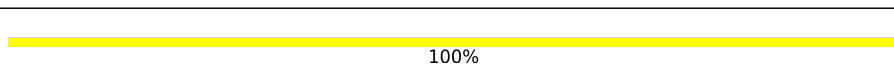
The reported resolution of this entry is 1.56 Å.

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	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	AAA	272	 82% 13%
1	BBB	272	 81% 6% 13%
2	AaA	3	 100%
2	BaB	3	 100%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7829 atoms, of which 3628 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 Glycoside hydrolase family 16 protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	AAA	236	3656	1197	1776	312	364	7	113	1	0
1	BBB	236	3673	1200	1792	314	360	7	113	1	0

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	0	MET	-	initiating methionine	UNP A0A1Y4PWX9
AAA	1	GLY	-	expression tag	UNP A0A1Y4PWX9
AAA	2	SER	-	expression tag	UNP A0A1Y4PWX9
AAA	3	SER	-	expression tag	UNP A0A1Y4PWX9
AAA	4	HIS	-	expression tag	UNP A0A1Y4PWX9
AAA	5	HIS	-	expression tag	UNP A0A1Y4PWX9
AAA	6	HIS	-	expression tag	UNP A0A1Y4PWX9
AAA	7	HIS	-	expression tag	UNP A0A1Y4PWX9
AAA	8	HIS	-	expression tag	UNP A0A1Y4PWX9
AAA	9	HIS	-	expression tag	UNP A0A1Y4PWX9
AAA	10	SER	-	expression tag	UNP A0A1Y4PWX9
AAA	11	SER	-	expression tag	UNP A0A1Y4PWX9
AAA	12	GLY	-	expression tag	UNP A0A1Y4PWX9
AAA	13	LEU	-	expression tag	UNP A0A1Y4PWX9
AAA	14	VAL	-	expression tag	UNP A0A1Y4PWX9
AAA	15	PRO	-	expression tag	UNP A0A1Y4PWX9
AAA	16	ARG	-	expression tag	UNP A0A1Y4PWX9
AAA	17	GLY	-	expression tag	UNP A0A1Y4PWX9
AAA	18	SER	-	expression tag	UNP A0A1Y4PWX9
AAA	19	HIS	-	expression tag	UNP A0A1Y4PWX9
AAA	20	MET	-	expression tag	UNP A0A1Y4PWX9
AAA	148	ALA	GLU	engineered mutation	UNP A0A1Y4PWX9
BBB	0	MET	-	initiating methionine	UNP A0A1Y4PWX9
BBB	1	GLY	-	expression tag	UNP A0A1Y4PWX9
BBB	2	SER	-	expression tag	UNP A0A1Y4PWX9

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Chain	Residue	Modelled	Actual	Comment	Reference
BBB	3	SER	-	expression tag	UNP A0A1Y4PWX9
BBB	4	HIS	-	expression tag	UNP A0A1Y4PWX9
BBB	5	HIS	-	expression tag	UNP A0A1Y4PWX9
BBB	6	HIS	-	expression tag	UNP A0A1Y4PWX9
BBB	7	HIS	-	expression tag	UNP A0A1Y4PWX9
BBB	8	HIS	-	expression tag	UNP A0A1Y4PWX9
BBB	9	HIS	-	expression tag	UNP A0A1Y4PWX9
BBB	10	SER	-	expression tag	UNP A0A1Y4PWX9
BBB	11	SER	-	expression tag	UNP A0A1Y4PWX9
BBB	12	GLY	-	expression tag	UNP A0A1Y4PWX9
BBB	13	LEU	-	expression tag	UNP A0A1Y4PWX9
BBB	14	VAL	-	expression tag	UNP A0A1Y4PWX9
BBB	15	PRO	-	expression tag	UNP A0A1Y4PWX9
BBB	16	ARG	-	expression tag	UNP A0A1Y4PWX9
BBB	17	GLY	-	expression tag	UNP A0A1Y4PWX9
BBB	18	SER	-	expression tag	UNP A0A1Y4PWX9
BBB	19	HIS	-	expression tag	UNP A0A1Y4PWX9
BBB	20	MET	-	expression tag	UNP A0A1Y4PWX9
BBB	148	ALA	GLU	engineered mutation	UNP A0A1Y4PWX9

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	AaA	3	Total	C	F	H	O	9	0	0
			63	18	1	30	14			
2	BaB	3	Total	C	F	H	O	8	0	0
			63	18	1	30	14			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	200	Total	O	0	0
			200	200		
3	BBB	174	Total	O	0	0
			174	174		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.45Å 84.58Å 67.11Å 90.00° 93.98° 90.00°	Depositor
Resolution (Å)	39.40 – 1.56 39.40 – 1.56	Depositor EDS
% Data completeness (in resolution range)	98.6 (39.40-1.56) 98.6 (39.40-1.56)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 1.56Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.172 , 0.199 0.172 , 0.199	Depositor DCC
R_{free} test set	3598 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	21.0	Xtrriage
Anisotropy	0.383	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 48.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7829	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.33% 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: G2F, 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	AAA	1.03	0/1928	1.15	0/2609
1	BBB	1.02	1/1929 (0.1%)	1.15	3/2608 (0.1%)
All	All	1.02	1/3857 (0.0%)	1.15	3/5217 (0.1%)

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	BBB	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	BBB	116	LYS	C-O	5.28	1.30	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BBB	190	LEU	N-CA-C	-5.08	106.77	113.17
1	BBB	153	SER	CA-C-N	5.04	125.58	119.98
1	BBB	153	SER	C-N-CA	5.04	125.58	119.98

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	BBB	64	GLN	Peptide

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	AAA	1880	1776	1757	7	0
1	BBB	1881	1792	1773	8	0
2	AaA	33	30	27	0	0
2	BaB	33	30	27	0	0
3	AAA	200	0	0	1	0
3	BBB	174	0	0	1	0
All	All	4201	3628	3584	14	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 2.

The worst 5 of 14 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:BBB:165:ASN:HB3	3:BBB:435:HOH:O	1.97	0.65
1:AAA:165:ASN:HB3	3:AAA:451:HOH:O	2.08	0.54
1:BBB:93:ILE:HD11	1:BBB:249:ILE:CD1	2.39	0.53
1:BBB:120:THR:HA	1:BBB:124:LEU:HB2	1.95	0.48
1:AAA:82:LEU:HD23	1:AAA:265:ILE:HD13	1.97	0.47

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	AAA	233/272 (86%)	230 (99%)	3 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	BBB	233/272 (86%)	229 (98%)	4 (2%)	0	100	100
All	All	466/544 (86%)	459 (98%)	7 (2%)	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	AAA	195/230 (85%)	194 (100%)	1 (0%)	81	68
1	BBB	195/230 (85%)	193 (99%)	2 (1%)	68	47
All	All	390/460 (85%)	387 (99%)	3 (1%)	73	56

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	143	GLU
1	BBB	143	GLU
1	BBB	151	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

6 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	G2F	AaA	1	2	11,11,12	1.10	0	12,15,17	1.08	1 (8%)
2	BGC	AaA	2	2	11,11,12	1.08	0	15,15,17	1.40	2 (13%)
2	BGC	AaA	3	2	11,11,12	0.65	0	15,15,17	1.12	1 (6%)
2	G2F	BaB	1	2	11,11,12	0.72	0	12,15,17	1.55	4 (33%)
2	BGC	BaB	2	2	11,11,12	1.55	3 (27%)	15,15,17	1.55	2 (13%)
2	BGC	BaB	3	2	11,11,12	0.67	0	15,15,17	1.09	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	G2F	AaA	1	2	-	0/2/19/22	0/1/1/1
2	BGC	AaA	2	2	-	0/2/19/22	0/1/1/1
2	BGC	AaA	3	2	-	0/2/19/22	0/1/1/1
2	G2F	BaB	1	2	-	0/2/19/22	0/1/1/1
2	BGC	BaB	2	2	-	0/2/19/22	0/1/1/1
2	BGC	BaB	3	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	BaB	2	BGC	C2-C3	3.51	1.57	1.52
2	BaB	2	BGC	O5-C5	2.37	1.48	1.43
2	BaB	2	BGC	O4-C4	2.02	1.48	1.43

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BaB	2	BGC	C1-C2-C3	-3.30	104.84	109.64
2	BaB	2	BGC	C6-C5-C4	2.84	120.00	113.02
2	BaB	1	G2F	C1-O5-C5	2.80	115.94	112.19
2	AaA	2	BGC	C1-C2-C3	-2.73	105.67	109.64
2	AaA	2	BGC	O3-C3-C2	-2.54	104.88	110.05

There are no chirality outliers.

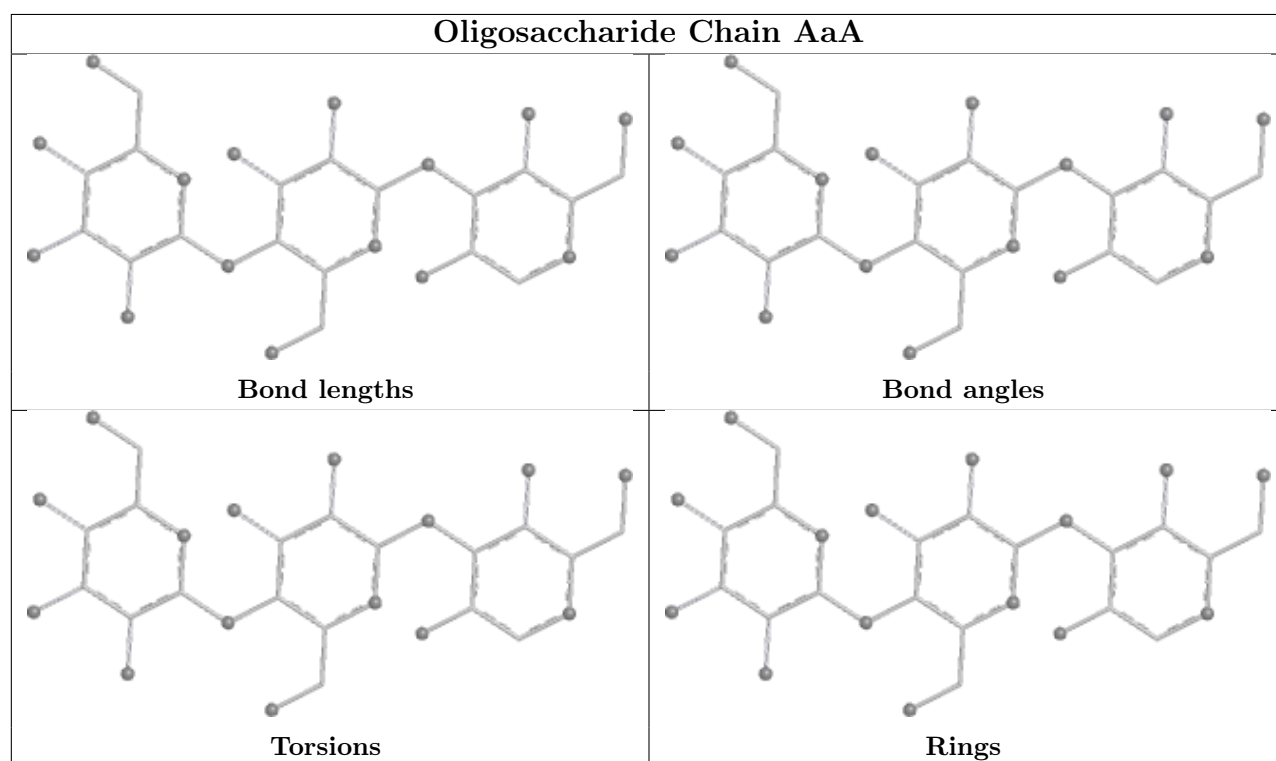
All (1) torsion outliers are listed below:

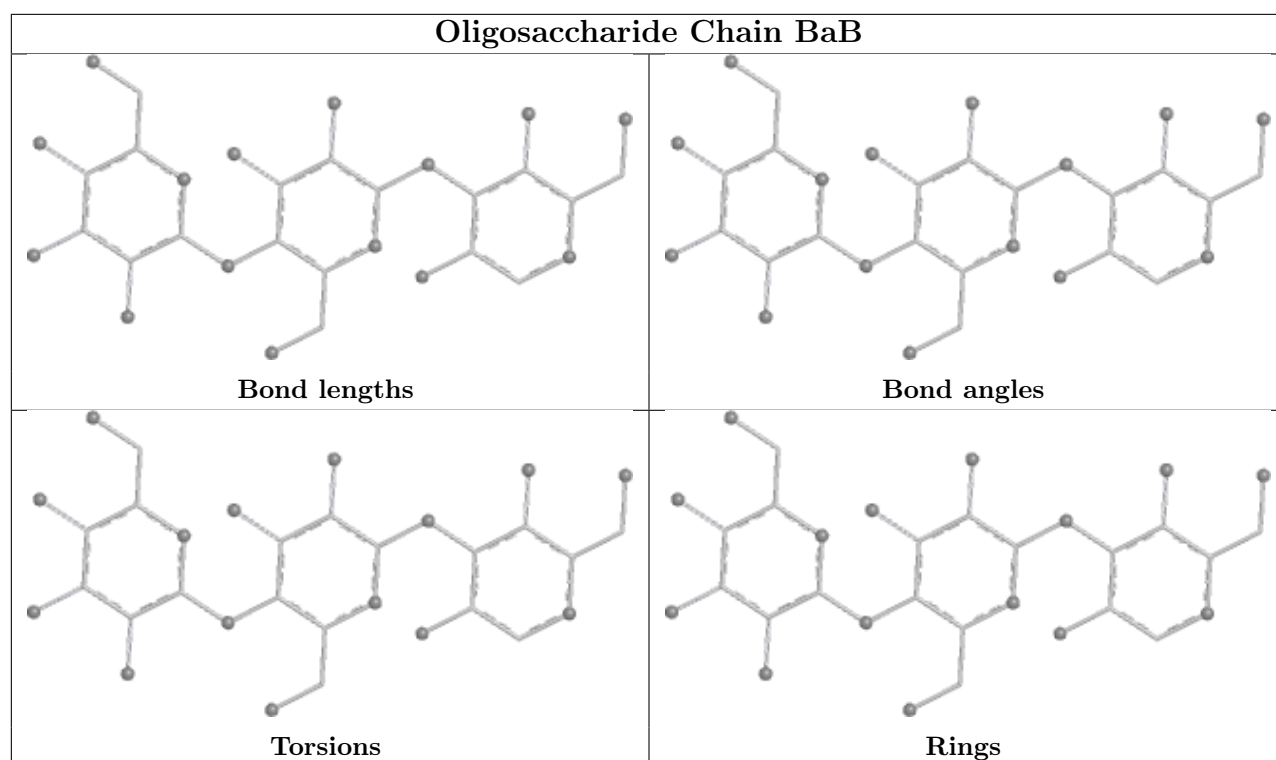
Mol	Chain	Res	Type	Atoms
2	BaB	3	BGC	C4-C5-C6-O6

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	AAA	236/272 (86%)	-0.27	1 (0%) 88 92	10, 22, 36, 46	1 (0%)
1	BBB	236/272 (86%)	0.08	4 (1%) 69 77	10, 26, 45, 72	1 (0%)
All	All	472/544 (86%)	-0.09	5 (1%) 78 84	10, 24, 41, 72	2 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	157	ALA	6.4
1	AAA	35	ILE	5.9
1	BBB	35	ILE	5.1
1	BBB	158	GLY	2.7
1	BBB	36	LEU	2.1

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.

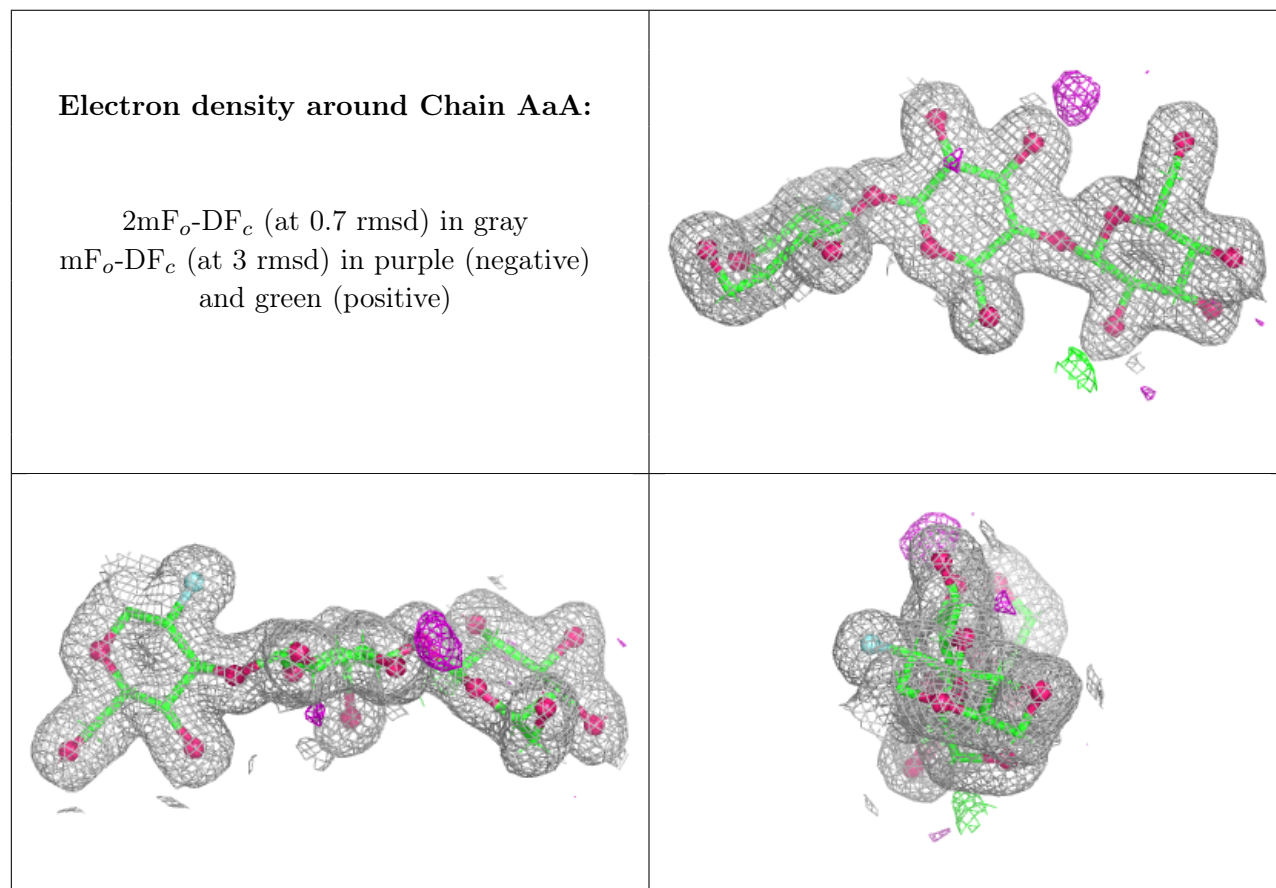
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	G2F	AaA	1	11/12	-	-	16,18,19,20	2
2	BGC	AaA	2	11/12	-	-	0,19,21,22	3
2	BGC	AaA	3	11/12	-	-	0,31,33,36	4
2	G2F	BaB	1	11/12	-	-	18,19,21,23	2

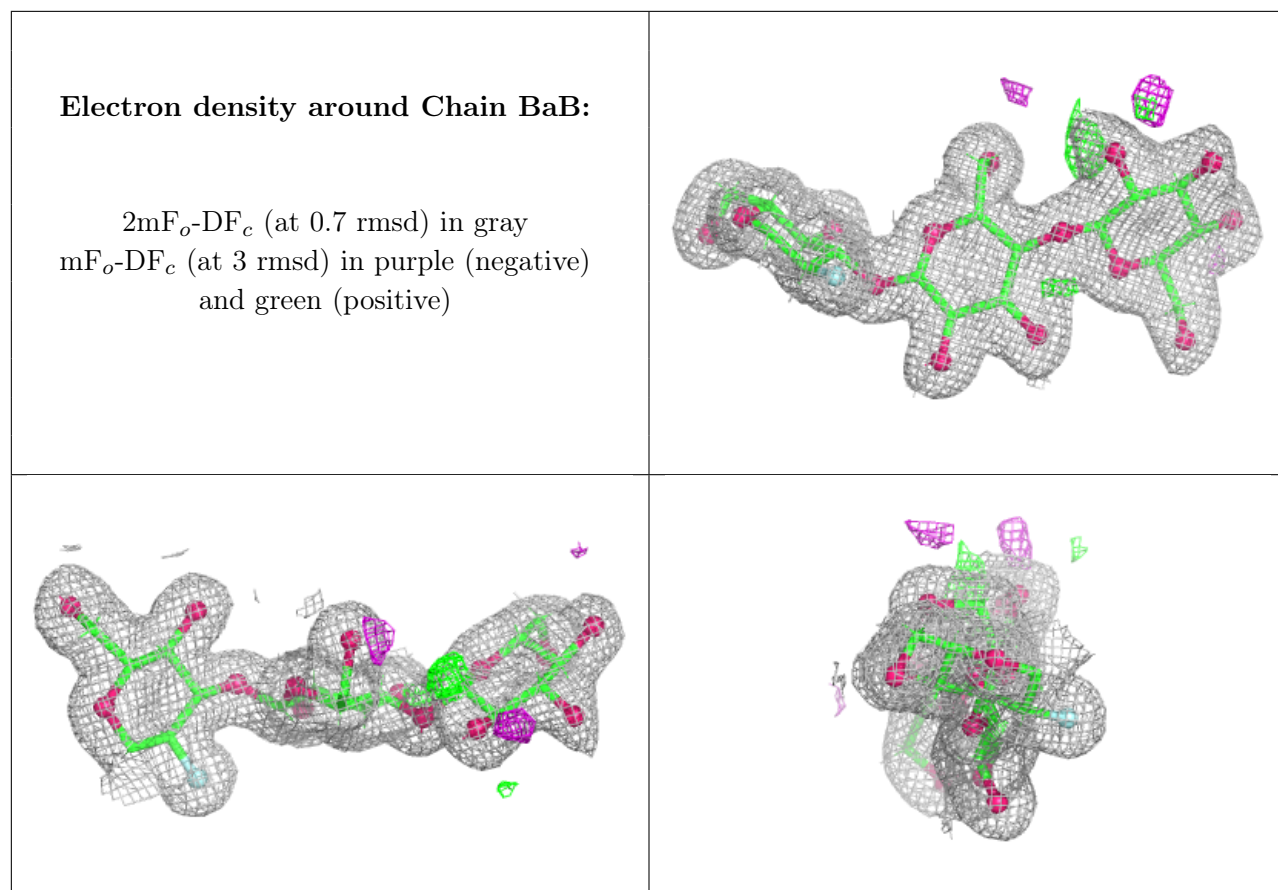
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BGC	BaB	2	11/12	-	-	0,21,25,25	3
2	BGC	BaB	3	11/12	-	-	0,35,39,41	3

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.