



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 09:33 AM UTC

PDB ID : 5GZK / pdb_00005gzk
Title : Endo-beta-1,2-glucanase from Chitinophaga pinensis - sophorotriose and glucose complex
Authors : Abe, K.; Nakajima, M.; Arakawa, T.; Fushinobu, S.; Taguchi, H.
Deposited on : 2016-09-28
Resolution : 1.70 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

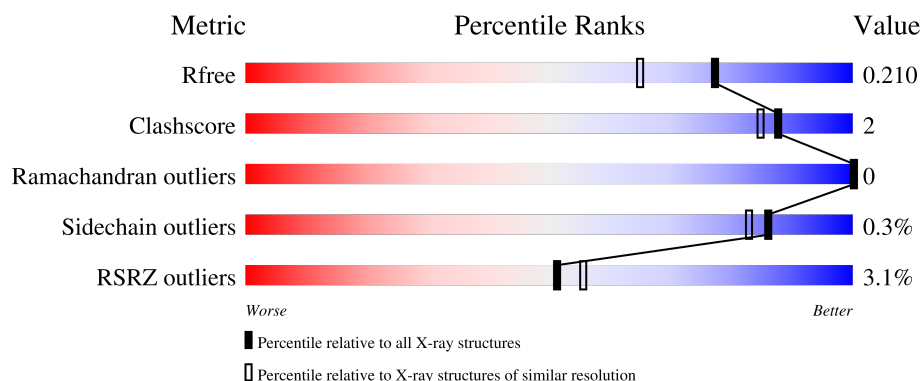
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

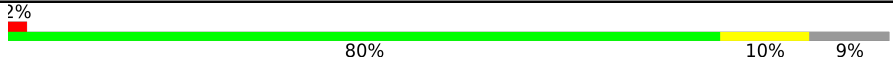
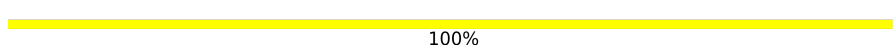
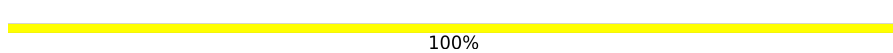
The reported resolution of this entry is 1.70 Å.

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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	461	
1	B	461	
2	C	3	
2	D	3	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
8	CL	B	503	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 7577 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	418	Total	C	N	O	S	0	0	0
			3385	2188	564	616	17			
1	B	418	Total	C	N	O	S	0	0	0
			3385	2188	564	616	17			

There are 40 discrepancies between the modelled and reference sequences:

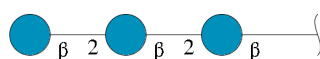
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP C7PIC2
A	-18	GLY	-	expression tag	UNP C7PIC2
A	-17	SER	-	expression tag	UNP C7PIC2
A	-16	SER	-	expression tag	UNP C7PIC2
A	-15	HIS	-	expression tag	UNP C7PIC2
A	-14	HIS	-	expression tag	UNP C7PIC2
A	-13	HIS	-	expression tag	UNP C7PIC2
A	-12	HIS	-	expression tag	UNP C7PIC2
A	-11	HIS	-	expression tag	UNP C7PIC2
A	-10	HIS	-	expression tag	UNP C7PIC2
A	-9	SER	-	expression tag	UNP C7PIC2
A	-8	SER	-	expression tag	UNP C7PIC2
A	-7	GLY	-	expression tag	UNP C7PIC2
A	-6	LEU	-	expression tag	UNP C7PIC2
A	-5	VAL	-	expression tag	UNP C7PIC2
A	-4	PRO	-	expression tag	UNP C7PIC2
A	-3	ARG	-	expression tag	UNP C7PIC2
A	-2	GLY	-	expression tag	UNP C7PIC2
A	-1	SER	-	expression tag	UNP C7PIC2
A	0	HIS	-	expression tag	UNP C7PIC2
B	-19	MET	-	initiating methionine	UNP C7PIC2
B	-18	GLY	-	expression tag	UNP C7PIC2
B	-17	SER	-	expression tag	UNP C7PIC2
B	-16	SER	-	expression tag	UNP C7PIC2
B	-15	HIS	-	expression tag	UNP C7PIC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP C7PIC2
B	-13	HIS	-	expression tag	UNP C7PIC2
B	-12	HIS	-	expression tag	UNP C7PIC2
B	-11	HIS	-	expression tag	UNP C7PIC2
B	-10	HIS	-	expression tag	UNP C7PIC2
B	-9	SER	-	expression tag	UNP C7PIC2
B	-8	SER	-	expression tag	UNP C7PIC2
B	-7	GLY	-	expression tag	UNP C7PIC2
B	-6	LEU	-	expression tag	UNP C7PIC2
B	-5	VAL	-	expression tag	UNP C7PIC2
B	-4	PRO	-	expression tag	UNP C7PIC2
B	-3	ARG	-	expression tag	UNP C7PIC2
B	-2	GLY	-	expression tag	UNP C7PIC2
B	-1	SER	-	expression tag	UNP C7PIC2
B	0	HIS	-	expression tag	UNP C7PIC2

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	3	Total	C	O	0	0	0
			34	18	16			
2	D	3	Total	C	O	0	0	0
			34	18	16			

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).

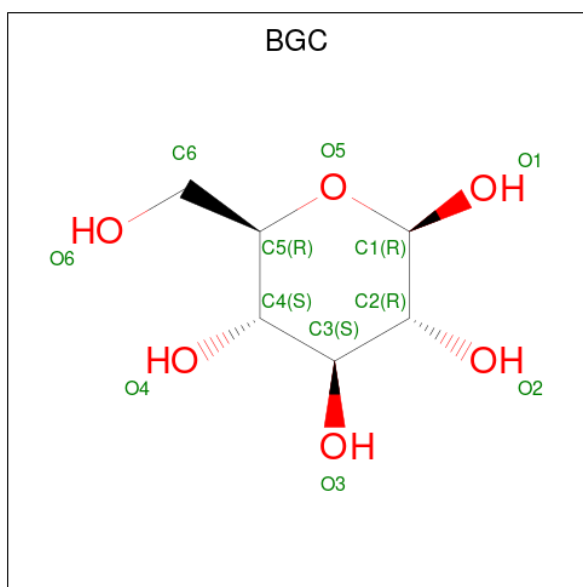


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

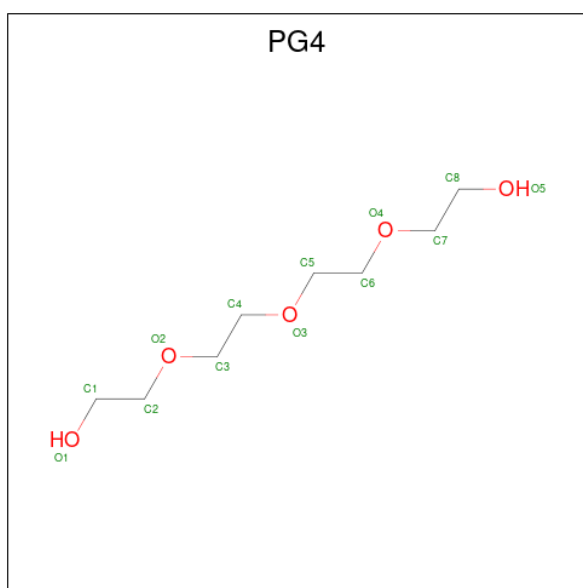
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	K	0	0
			1	1		
4	B	1	Total	K	0	0
			1	1		

- Molecule 5 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			12	6	6		
5	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).

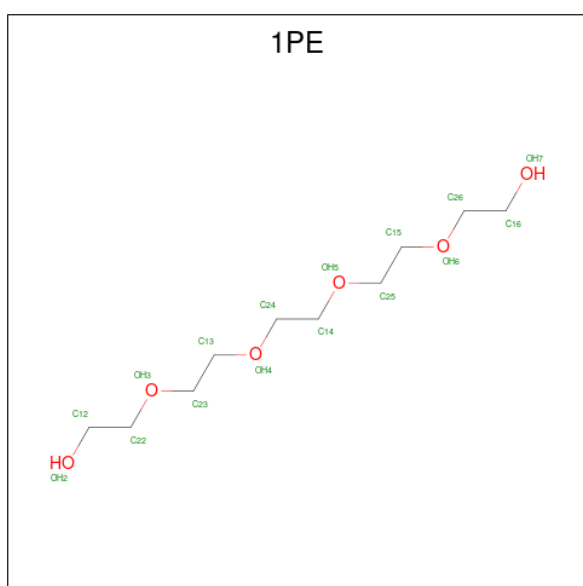


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Cl	0	0
			1	1		

- Molecule 9 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).

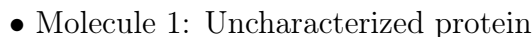


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	334	Total	O	0	0
			334	334		
10	B	332	Total	O	0	0
			332	332		

- Molecule 1: Uncharacterized protein



Chain D:

100%

B6C1
B6C2
B6C3

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	126.11Å 117.40Å 75.55Å 90.00° 92.56° 90.00°	Depositor
Resolution (Å)	49.46 – 1.70 49.46 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.46-1.70) 99.9 (49.46-1.70)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.48 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.166 , 0.199 0.181 , 0.210	Depositor DCC
R_{free} test set	5990 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	14.3	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.083 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7577	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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: PG4, K, PO4, 1PE, BGC, PEG, CL

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	1.59	30/3499 (0.9%)	1.29	11/4753 (0.2%)
1	B	1.58	32/3499 (0.9%)	1.32	8/4753 (0.2%)
All	All	1.59	62/6998 (0.9%)	1.31	19/9506 (0.2%)

The worst 5 of 62 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	419	LEU	N-CA	-8.44	1.36	1.46
1	A	26	ALA	CA-C	-8.39	1.41	1.52
1	A	27	ASP	N-CA	-7.15	1.37	1.46
1	A	149	GLY	N-CA	-7.15	1.37	1.45
1	A	307	ASN	C-O	6.79	1.27	1.23

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	87	GLU	N-CA-C	6.76	118.65	111.28
1	A	97	LEU	N-CA-C	6.24	117.74	111.07
1	B	375	ILE	N-CA-C	6.20	119.11	113.10
1	B	96	GLY	N-CA-C	6.19	120.13	112.64
1	A	378	LYS	N-CA-C	6.08	118.70	111.71

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	3385	0	3193	14	0
1	B	3385	0	3193	12	0
2	C	34	0	30	0	0
2	D	34	0	29	0	0
3	A	5	0	0	1	0
3	B	5	0	0	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	12	0	12	0	0
5	B	12	0	12	1	0
6	A	13	0	18	4	0
7	A	7	0	10	0	0
8	B	1	0	0	3	0
9	B	16	0	22	3	0
10	A	334	0	0	2	1
10	B	332	0	0	2	2
All	All	7577	0	6519	27	3

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 27 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)
5:B:504:BGC:H3	10:B:834:HOH:O	1.66	0.93
1:A:97:LEU:HG	1:A:101:MET:HE2	1.73	0.70
1:A:418:LYS:NZ	6:A:504:PG4:H22	2.07	0.70
1:A:291:GLU:OE2	3:A:501:PO4:O4	2.17	0.62
1:B:291:GLU:OE2	3:B:501:PO4:O4	2.20	0.60

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:B:903:HOH:O	10:B:903:HOH:O[2_655]	1.03	1.17
10:A:891:HOH:O	10:A:891:HOH:O[2_555]	1.14	1.06
10:B:875:HOH:O	10:B:875:HOH:O[2_655]	1.14	1.06

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	416/461 (90%)	399 (96%)	17 (4%)	0	100	100
1	B	416/461 (90%)	402 (97%)	14 (3%)	0	100	100
All	All	832/922 (90%)	801 (96%)	31 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	348/385 (90%)	347 (100%)	1 (0%)	86	83
1	B	348/385 (90%)	347 (100%)	1 (0%)	86	83
All	All	696/770 (90%)	694 (100%)	2 (0%)	86	83

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	THR
1	B	162	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	251	HIS

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Mol	Chain	Res	Type
1	B	395	GLN
1	B	436	GLN
1	A	395	GLN
1	A	339	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 ⓘ

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	BGC	C	1	2	12,12,12	0.78	0	17,17,17	1.64	5 (29%)
2	BGC	C	2	2	11,11,12	1.77	3 (27%)	15,15,17	2.02	4 (26%)
2	BGC	C	3	2	11,11,12	1.16	0	15,15,17	1.12	1 (6%)
2	BGC	D	1	2	12,12,12	0.92	0	17,17,17	1.10	1 (5%)
2	BGC	D	2	2	11,11,12	1.78	3 (27%)	15,15,17	2.26	5 (33%)
2	BGC	D	3	2	11,11,12	1.88	3 (27%)	15,15,17	1.33	2 (13%)

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	-	2/2/22/22	0/1/1/1
2	BGC	C	2	2	-	0/2/19/22	0/1/1/1
2	BGC	C	3	2	-	0/2/19/22	0/1/1/1
2	BGC	D	1	2	-	0/2/22/22	0/1/1/1
2	BGC	D	2	2	-	0/2/19/22	0/1/1/1
2	BGC	D	3	2	-	0/2/19/22	0/1/1/1

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	BGC	O4-C4	-3.47	1.34	1.43
2	C	2	BGC	C4-C5	3.39	1.60	1.53
2	D	2	BGC	O3-C3	2.97	1.50	1.43
2	D	3	BGC	C2-C3	2.94	1.57	1.52
2	D	3	BGC	C4-C5	2.62	1.58	1.53

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	BGC	C2-C3-C4	-4.56	102.83	110.86
2	C	2	BGC	C1-O5-C5	4.52	118.24	112.19
2	C	2	BGC	O3-C3-C4	3.94	119.66	110.38
2	C	1	BGC	C1-O5-C5	3.55	120.53	113.65
2	D	2	BGC	O3-C3-C4	3.48	118.59	110.38

There are no chirality outliers.

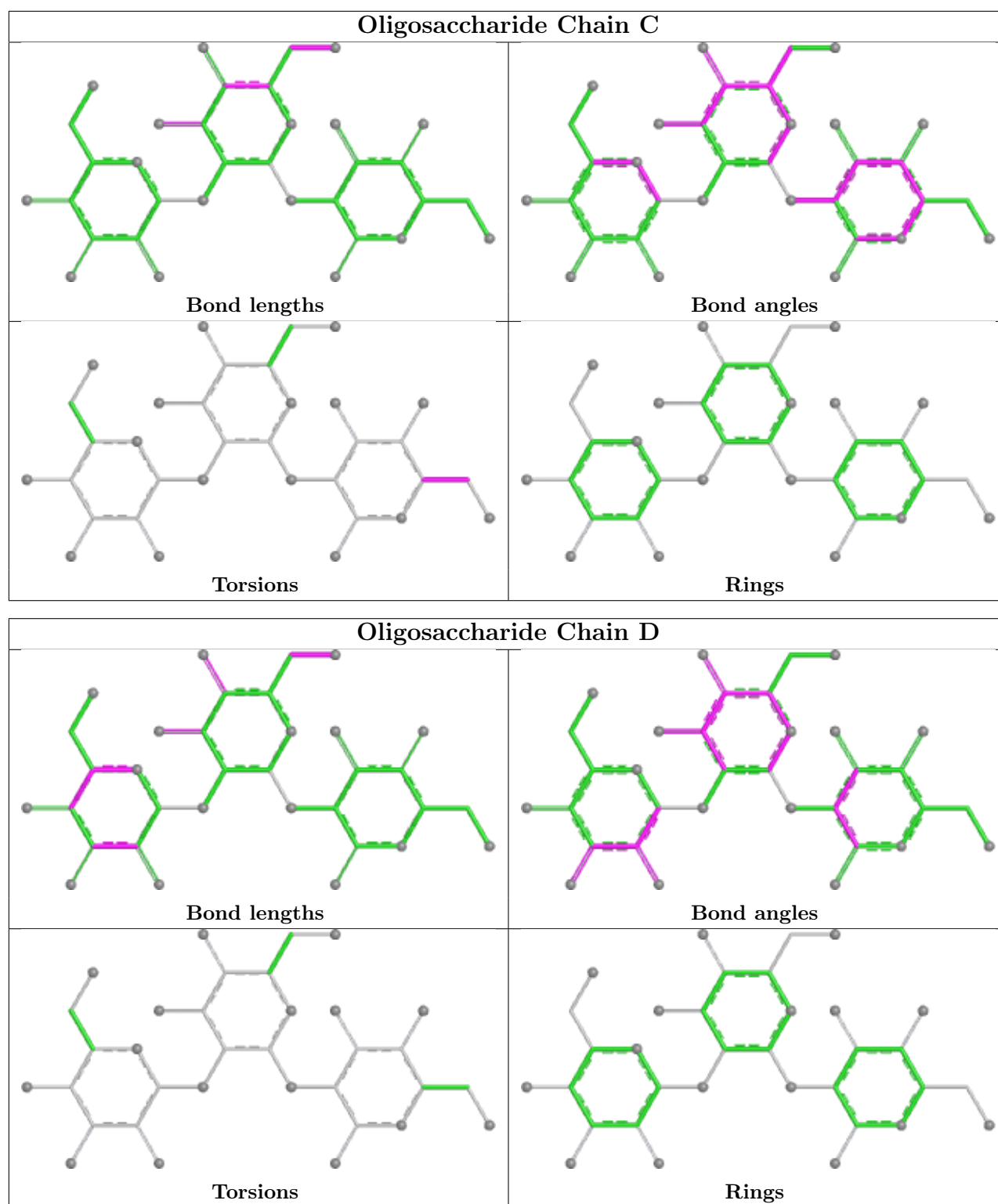
All (2) torsion outliers are listed below:

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

Of 10 ligands modelled in this entry, 3 are monoatomic - leaving 7 for Mogul analysis.

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	$\# Z > 2$
5	BGC	B	504	-	12,12,12	1.03	0	17,17,17	0.99	1 (5%)
3	PO4	A	501	-	4,4,4	1.45	1 (25%)	6,6,6	1.25	1 (16%)
7	PEG	A	505	-	6,6,6	0.82	0	5,5,5	1.21	1 (20%)
9	1PE	B	505	-	15,15,15	0.71	0	14,14,14	1.07	1 (7%)
3	PO4	B	501	-	4,4,4	1.54	1 (25%)	6,6,6	0.99	0
5	BGC	A	503	-	12,12,12	1.06	0	17,17,17	1.59	2 (11%)
6	PG4	A	504	-	12,12,12	0.88	0	11,11,11	1.44	1 (9%)

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
5	BGC	B	504	-	-	0/2/22/22	0/1/1/1
7	PEG	A	505	-	-	1/4/4/4	-
9	1PE	B	505	-	-	4/13/13/13	-
5	BGC	A	503	-	-	0/2/22/22	0/1/1/1
6	PG4	A	504	-	-	3/10/10/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	PO4	P-O3	2.67	1.62	1.54
3	A	501	PO4	P-O2	2.12	1.60	1.54

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	504	PG4	O2-C2-C1	3.79	126.83	110.11
5	A	503	BGC	C3-C4-C5	3.40	116.39	110.23
5	B	504	BGC	C4-C3-C2	2.96	116.03	110.83
5	A	503	BGC	C1-O5-C5	-2.47	108.88	113.65
7	A	505	PEG	O2-C3-C4	2.35	120.45	110.11

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	504	PG4	O1-C1-C2-O2
9	B	505	1PE	OH6-C15-C25-OH5
7	A	505	PEG	O2-C3-C4-O4
6	A	504	PG4	O3-C5-C6-O4
6	A	504	PG4	O4-C7-C8-O5

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	504	BGC	1	0
3	A	501	PO4	1	0
9	B	505	1PE	3	0
3	B	501	PO4	1	0
6	A	504	PG4	4	0

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	418/461 (90%)	-0.01	10 (2%) 59 64	9, 16, 28, 50	0
1	B	418/461 (90%)	0.07	16 (3%) 44 48	9, 17, 31, 57	0
All	All	836/922 (90%)	0.03	26 (3%) 51 55	9, 17, 30, 57	0

The worst 5 of 26 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	23	ALA	4.5
1	B	23	ALA	4.5
1	B	160	GLY	3.7
1	B	161	ASN	3.2
1	B	158	ALA	3.1

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	BGC	D	3	11/12	0.85	0.13	22,26,31,32	0
2	BGC	C	3	11/12	0.88	0.12	23,28,31,34	0
2	BGC	C	1	12/12	0.88	0.12	27,38,41,45	0
2	BGC	D	1	12/12	0.90	0.10	23,30,34,34	0
2	BGC	C	2	11/12	0.92	0.11	17,25,26,28	0

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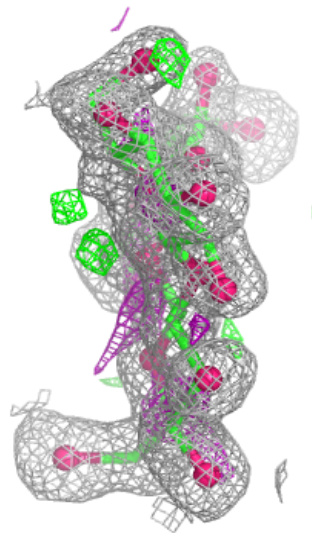
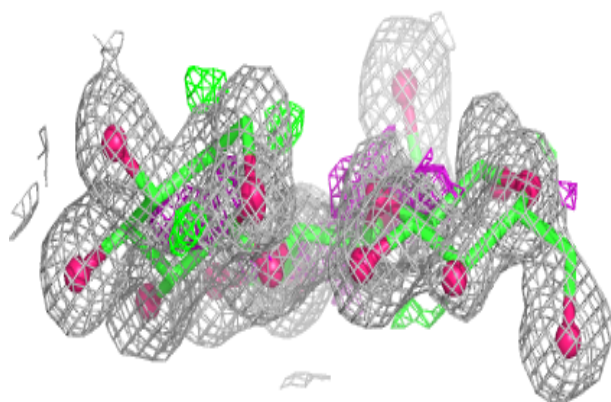
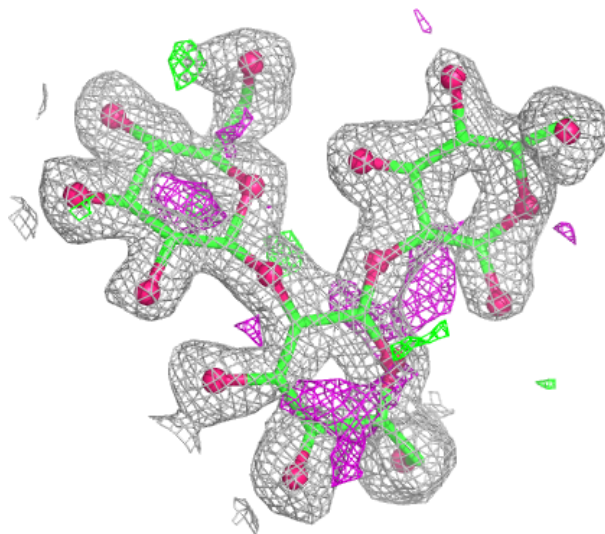
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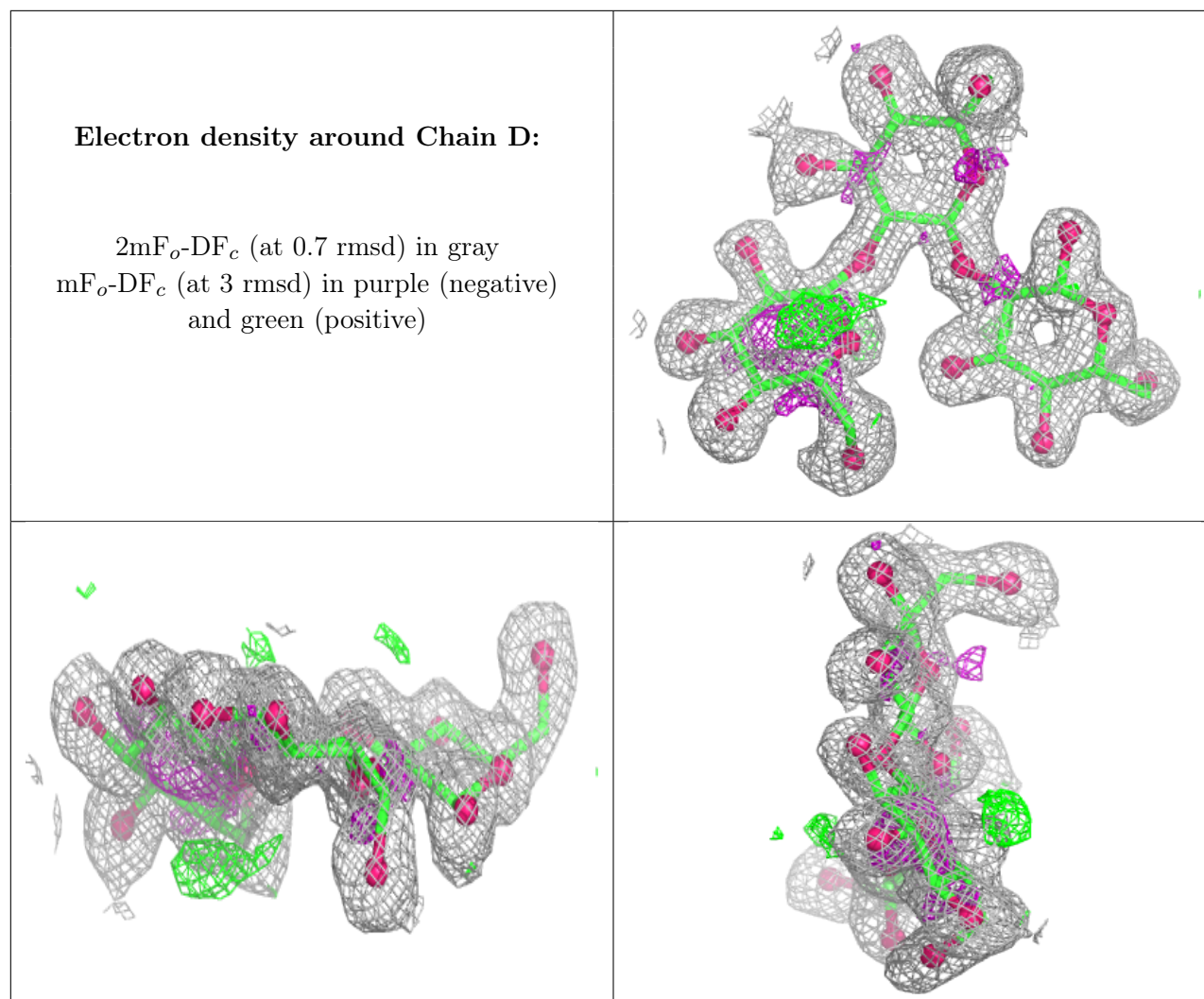
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BGC	D	2	11/12	0.94	0.09	15,25,28,29	0

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 ⓘ

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
5	BGC	A	503	12/12	0.85	0.14	28,40,46,48	0
5	BGC	B	504	12/12	0.85	0.14	32,46,51,51	0
6	PG4	A	504	13/13	0.85	0.15	30,32,39,40	0
9	1PE	B	505	16/16	0.85	0.14	31,35,42,47	0
7	PEG	A	505	7/7	0.86	0.15	32,33,36,41	0
8	CL	B	503	1/1	0.93	0.20	43,43,43,43	0
3	PO4	A	501	5/5	0.95	0.14	23,29,33,33	0
3	PO4	B	501	5/5	0.96	0.12	22,26,30,31	0
4	K	B	502	1/1	0.96	0.23	42,42,42,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	K	A	502	1/1	0.97	0.21	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.