



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 03:19 AM UTC

PDB ID : 5LRB / pdb\_00005lrb  
Title : Plastidial phosphorylase from Barley in complex with acarbose  
Authors : Cuesta-Seijo, J.A.; Ruzanski, C.; Kruzewicz, K.; Palcic, M.M.  
Deposited on : 2016-08-18  
Resolution : 2.90 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](mailto: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>.

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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

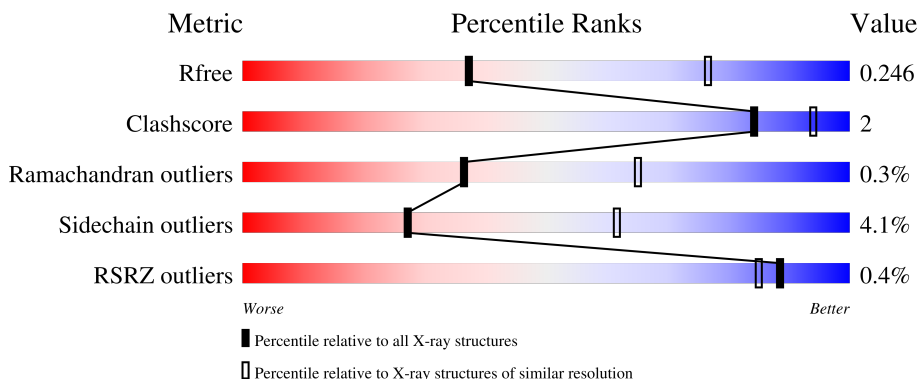
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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  $\geq 3$ , 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	938	 81% 8% 10%
1	B	938	 81% 9% 10%
2	C	3	 100%
2	D	3	 33% 33% 33%

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
3	GLC	B	1001	X	-	-	-

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 13689 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 Alpha-1,4 glucan phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	844	6744	4296	1148	1270	30	0	5	1
1	B	844	6747	4298	1146	1273	30	0	6	1

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	43	MET	-	initiating methionine	UNP F2E0G2
A	969	THR	-	expression tag	UNP F2E0G2
A	970	PHE	-	expression tag	UNP F2E0G2
A	971	PRO	-	expression tag	UNP F2E0G2
A	972	ASP	-	expression tag	UNP F2E0G2
A	973	ILE	-	expression tag	UNP F2E0G2
A	974	GLU	-	expression tag	UNP F2E0G2
A	975	ASN	-	expression tag	UNP F2E0G2
A	976	LEU	-	expression tag	UNP F2E0G2
A	977	TYR	-	expression tag	UNP F2E0G2
A	978	PHE	-	expression tag	UNP F2E0G2
A	979	GLN	-	expression tag	UNP F2E0G2
A	980	GLY	-	expression tag	UNP F2E0G2
B	43	MET	-	initiating methionine	UNP F2E0G2
B	969	THR	-	expression tag	UNP F2E0G2
B	970	PHE	-	expression tag	UNP F2E0G2
B	971	PRO	-	expression tag	UNP F2E0G2
B	972	ASP	-	expression tag	UNP F2E0G2
B	973	ILE	-	expression tag	UNP F2E0G2
B	974	GLU	-	expression tag	UNP F2E0G2
B	975	ASN	-	expression tag	UNP F2E0G2
B	976	LEU	-	expression tag	UNP F2E0G2
B	977	TYR	-	expression tag	UNP F2E0G2
B	978	PHE	-	expression tag	UNP F2E0G2
B	979	GLN	-	expression tag	UNP F2E0G2

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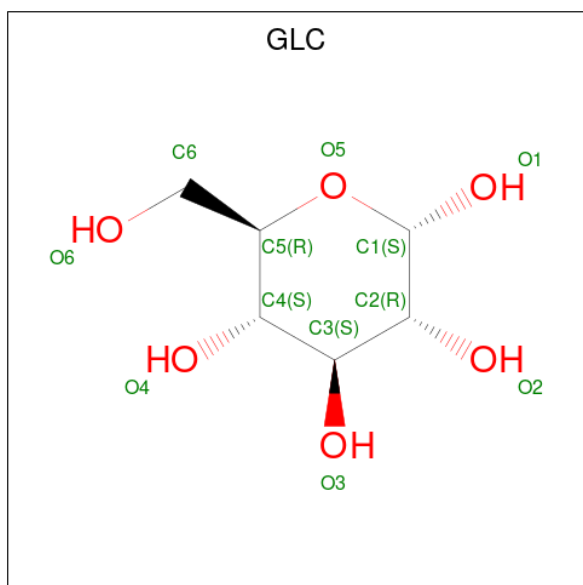
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Chain	Residue	Modelled	Actual	Comment	Reference
B	980	GLY	-	expression tag	UNP F2E0G2

- Molecule 2 is an oligosaccharide called 4,6-dideoxy-4-[[[(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

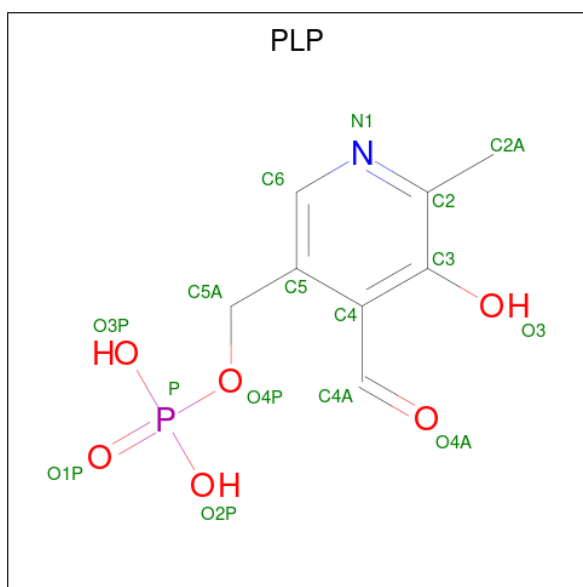
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace	
2	C	3	Total	C	N	O	0	0	0
			44	25	1	18			
2	D	3	Total	C	N	O	0	0	0
			44	25	1	18			

- Molecule 3 is alpha-D-glucopyranose (CCD ID: GLC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



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

- Molecule 4 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: C<sub>8</sub>H<sub>10</sub>NO<sub>6</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	15	8	1	5	1	0	0
4	B	1	15	8	1	5	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	28	28	28	0	0
5	B	28	28	28	0	0




- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}\}$ - $\alpha$ -D-glucopyranose-(1-4)- $\alpha$ -D-glucopyranose-(1-4)- $\alpha$ -D-glucopyranose

Chain C:  100%

GLC1  
GLC2  
AC13

- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}\}$ - $\alpha$ -D-glucopyranose-(1-4)- $\alpha$ -D-glucopyranose-(1-4)- $\alpha$ -D-glucopyranose

Chain D:  33% 33% 33%

GLC1  
GLC2  
AC13

## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	230.33Å 63.68Å 149.33Å 90.00° 115.07° 90.00°	Depositor
Resolution (Å)	48.92 – 2.90 48.92 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.2 (48.92-2.90) 98.2 (48.92-2.90)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.16 (at 2.91Å)	Xtrriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.213 , 0.246 0.218 , 0.246	Depositor DCC
$R_{free}$ test set	1297 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	76.2	Xtrriage
Anisotropy	0.046	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 41.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	13689	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	77.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.47 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.8395e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GLC, PLP, AC1

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	0.52	1/6914 (0.0%)	0.80	1/9365 (0.0%)
1	B	0.51	1/6920 (0.0%)	0.80	1/9374 (0.0%)
All	All	0.51	2/13834 (0.0%)	0.80	2/18739 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	968	PRO	C-N	-6.92	1.23	1.33
1	A	968	PRO	C-N	-6.79	1.23	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	463	GLY	N-CA-C	5.74	119.17	111.20
1	B	463	GLY	N-CA-C	5.42	119.40	110.77

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	6744	0	6642	29	0
1	B	6747	0	6641	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	44	0	30	0	0
2	D	44	0	30	1	0
3	A	12	0	12	0	0
3	B	12	0	11	0	0
4	A	15	0	6	1	0
4	B	15	0	6	0	0
5	A	28	0	0	0	0
5	B	28	0	0	1	0
All	All	13689	0	13378	63	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 63 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:422:THR:HG21	1:A:427:ALA:HB3	1.67	0.74
1:A:455:MET:HE1	1:A:485:LEU:HD22	1.72	0.70
1:A:140:ARG:O	1:A:619:ARG:NH1	2.28	0.66
1:B:140:ARG:O	1:B:619:ARG:NH1	2.29	0.66
1:A:455:MET:HE1	1:A:485:LEU:CD2	2.28	0.63

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	845/938 (90%)	808 (96%)	36 (4%)	1 (0%)	48 77
1	B	846/938 (90%)	807 (95%)	35 (4%)	4 (0%)	24 54
All	All	1691/1876 (90%)	1615 (96%)	71 (4%)	5 (0%)	36 65

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	258	ASP
1	A	259	GLY
1	B	259	GLY
1	B	257	THR
1	B	464	THR

### 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	729/810 (90%)	697 (96%)	32 (4%)	25 59
1	B	730/810 (90%)	700 (96%)	30 (4%)	27 61
All	All	1459/1620 (90%)	1397 (96%)	62 (4%)	27 60

5 of 62 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	746	THR
1	B	730	LYS
1	B	209	ARG
1	B	726	LYS
1	B	886	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	613	ASN
1	B	268	ASN
1	B	582	ASN
1	B	354	GLN
1	A	582	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
3	GLC	B	1001	-	12,12,12	0.59	0	17,17,17	2.37	5 (29%)
3	GLC	A	1001	-	12,12,12	0.56	0	17,17,17	0.97	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
3	GLC	B	1001	-	1/1/5/5	2/2/22/22	0/1/1/1
3	GLC	A	1001	-	-	2/2/22/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1001	GLC	O5-C1-C2	7.67	123.78	110.30
3	B	1001	GLC	O1-C1-C2	2.85	117.25	108.98
3	B	1001	GLC	O1-C1-O5	2.68	118.36	110.41
3	B	1001	GLC	C3-C4-C5	-2.51	105.68	110.23
3	B	1001	GLC	O5-C5-C6	2.21	111.92	106.44

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	1001	GLC	C1

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1001	GLC	O5-C5-C6-O6
3	B	1001	GLC	C4-C5-C6-O6
3	A	1001	GLC	O5-C5-C6-O6
3	A	1001	GLC	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

## 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	GLC	C	1	2	12,12,12	0.58	0	17,17,17	1.04	1 (5%)
2	GLC	C	2	2	11,11,12	0.45	0	15,15,17	1.26	2 (13%)
2	AC1	C	3	2	21,22,23	0.69	0	22,32,34	1.00	1 (4%)
2	GLC	D	1	2	12,12,12	0.56	0	17,17,17	0.54	0
2	GLC	D	2	2	11,11,12	0.53	0	15,15,17	1.14	2 (13%)
2	AC1	D	3	2	21,22,23	0.65	0	22,32,34	0.54	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	-	1/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	C	2	2	-	2/2/19/22	0/1/1/1
2	AC1	C	3	2	-	2/6/43/46	0/2/2/2
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	1/2/19/22	0/1/1/1
2	AC1	D	3	2	-	3/6/43/46	0/2/2/2

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	AC1	C1-C2-C3	3.64	114.95	109.64
2	D	2	GLC	C1-O5-C5	3.19	116.46	112.19
2	C	2	GLC	C1-C2-C3	2.89	113.85	109.64
2	C	2	GLC	C1-O5-C5	2.41	115.41	112.19
2	D	2	GLC	C1-C2-C3	2.21	112.86	109.64

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

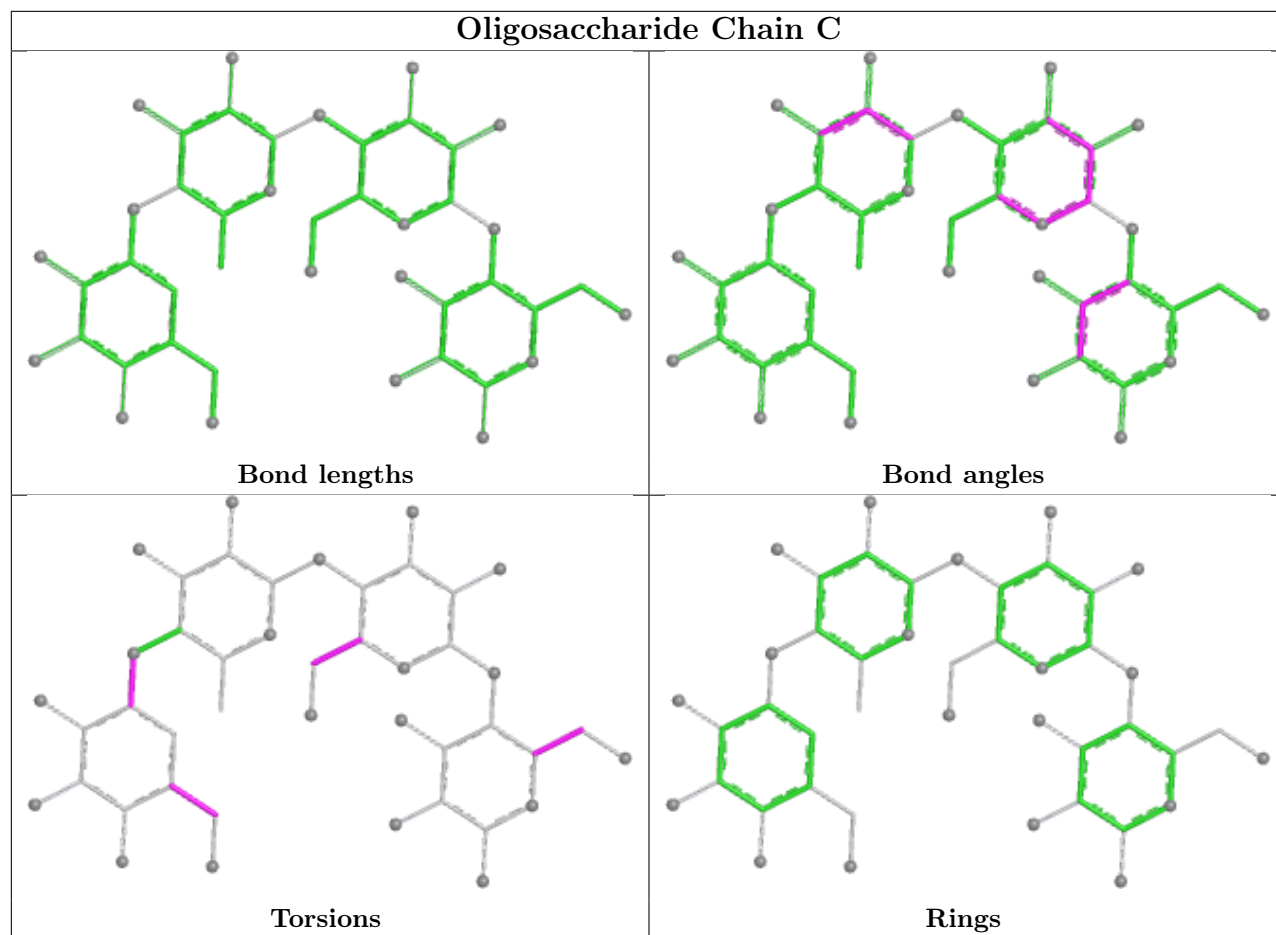
Mol	Chain	Res	Type	Atoms
2	C	3	AC1	C2B-C1B-N4A-C4
2	C	3	AC1	C7B-C5B-C6B-O6B
2	D	3	AC1	C3-C4-N4A-C1B
2	D	3	AC1	C5-C4-N4A-C1B
2	C	2	GLC	O5-C5-C6-O6

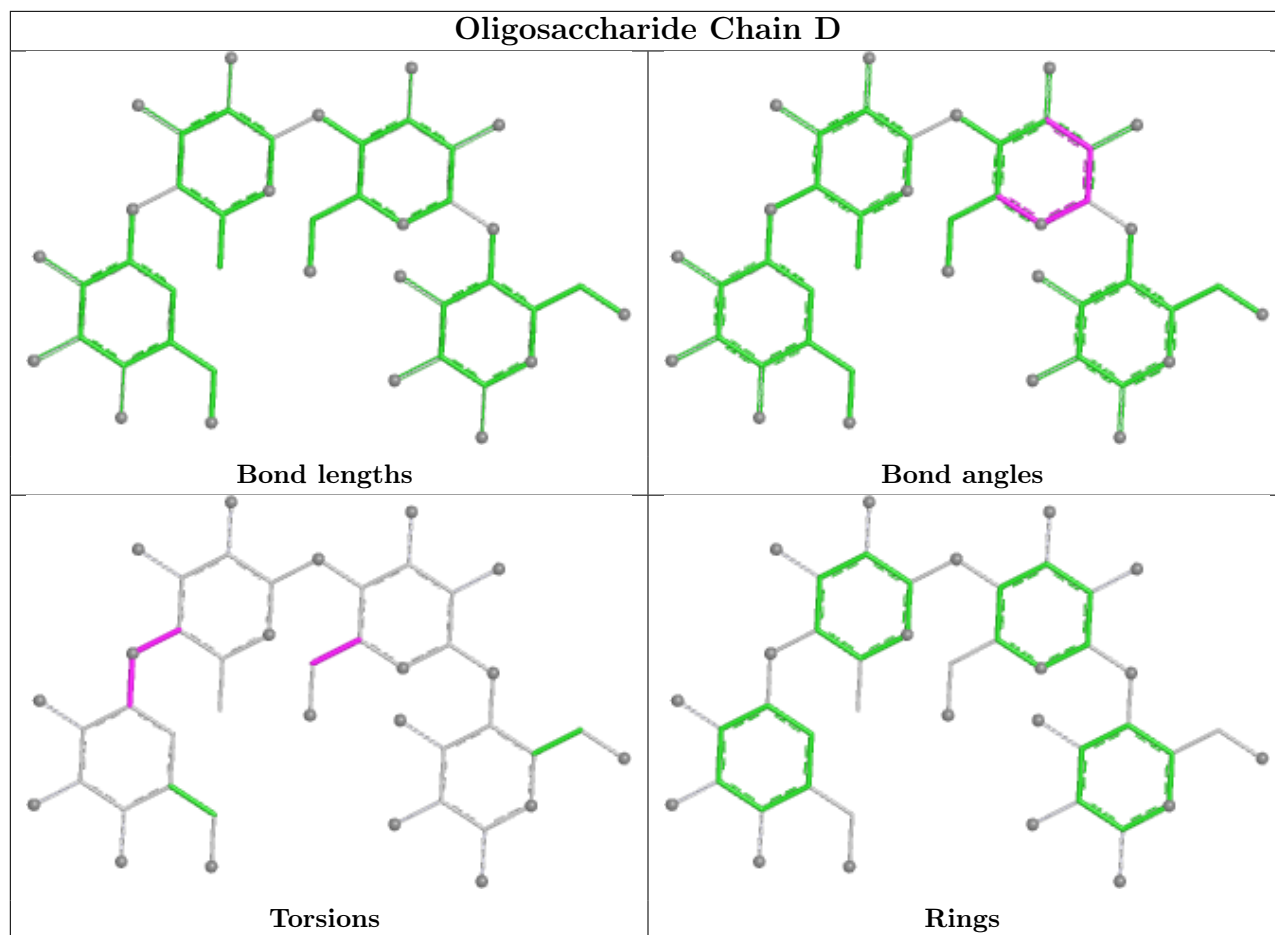
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	GLC	1	0
2	D	2	GLC	1	0

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](#)

4 ligands 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	$\# Z  > 2$
3	GLC	B	1001	-	12,12,12	0.59	0	17,17,17	2.37	5 (29%)
4	PLP	A	1002	1	15,15,16	2.70	3 (20%)	21,22,23	1.54	3 (14%)
4	PLP	B	1002	1	15,15,16	2.84	3 (20%)	21,22,23	1.30	3 (14%)
3	GLC	A	1001	-	12,12,12	0.56	0	17,17,17	0.97	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
3	GLC	B	1001	-	1/1/5/5	2/2/22/22	0/1/1/1
4	PLP	A	1002	1	-	4/6/6/8	0/1/1/1
4	PLP	B	1002	1	-	3/6/6/8	0/1/1/1
3	GLC	A	1001	-	-	2/2/22/22	0/1/1/1

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1002	PLP	C5-C4	7.40	1.48	1.40
4	B	1002	PLP	C3-C2	7.19	1.48	1.41
4	B	1002	PLP	C5-C4	6.89	1.48	1.40
4	A	1002	PLP	C3-C2	6.07	1.47	1.41
4	B	1002	PLP	C3-C4	3.89	1.47	1.40

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1001	GLC	O5-C1-C2	7.67	123.78	110.30
4	A	1002	PLP	C4A-C4-C5	3.98	125.04	120.94
3	B	1001	GLC	O1-C1-C2	2.85	117.25	108.98
3	B	1001	GLC	O1-C1-O5	2.68	118.36	110.41
3	B	1001	GLC	C3-C4-C5	-2.51	105.68	110.23

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	1001	GLC	C1

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1002	PLP	C5A-O4P-P-O2P
4	A	1002	PLP	C5A-O4P-P-O3P
4	B	1002	PLP	C5A-O4P-P-O2P
4	B	1002	PLP	C5A-O4P-P-O3P
3	B	1001	GLC	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1002	PLP	1	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	844/938 (89%)	-0.32	2 (0%) 91 89	37, 73, 106, 130	5 (0%)
1	B	844/938 (89%)	-0.32	4 (0%) 87 83	35, 75, 108, 134	6 (0%)
All	All	1688/1876 (89%)	-0.32	6 (0%) 88 85	35, 74, 107, 134	11 (0%)

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	424	LEU	2.9
1	B	425	PRO	2.8
1	B	789	THR	2.2
1	B	699[A]	ARG	2.2
1	A	553	GLU	2.2

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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.

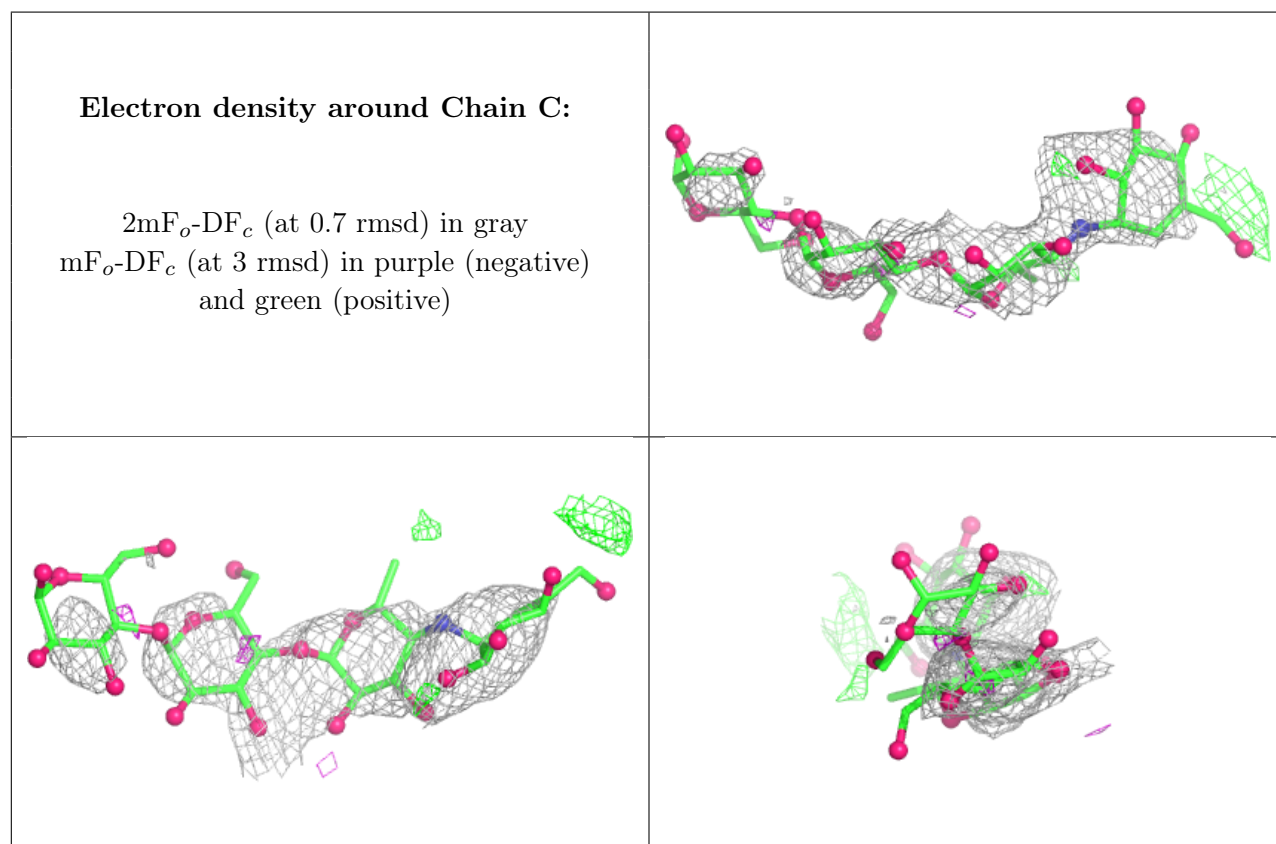
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	GLC	B	1001	12/12	0.77	0.22	106,121,127,135	0
3	GLC	A	1001	12/12	0.81	0.18	97,116,121,127	0

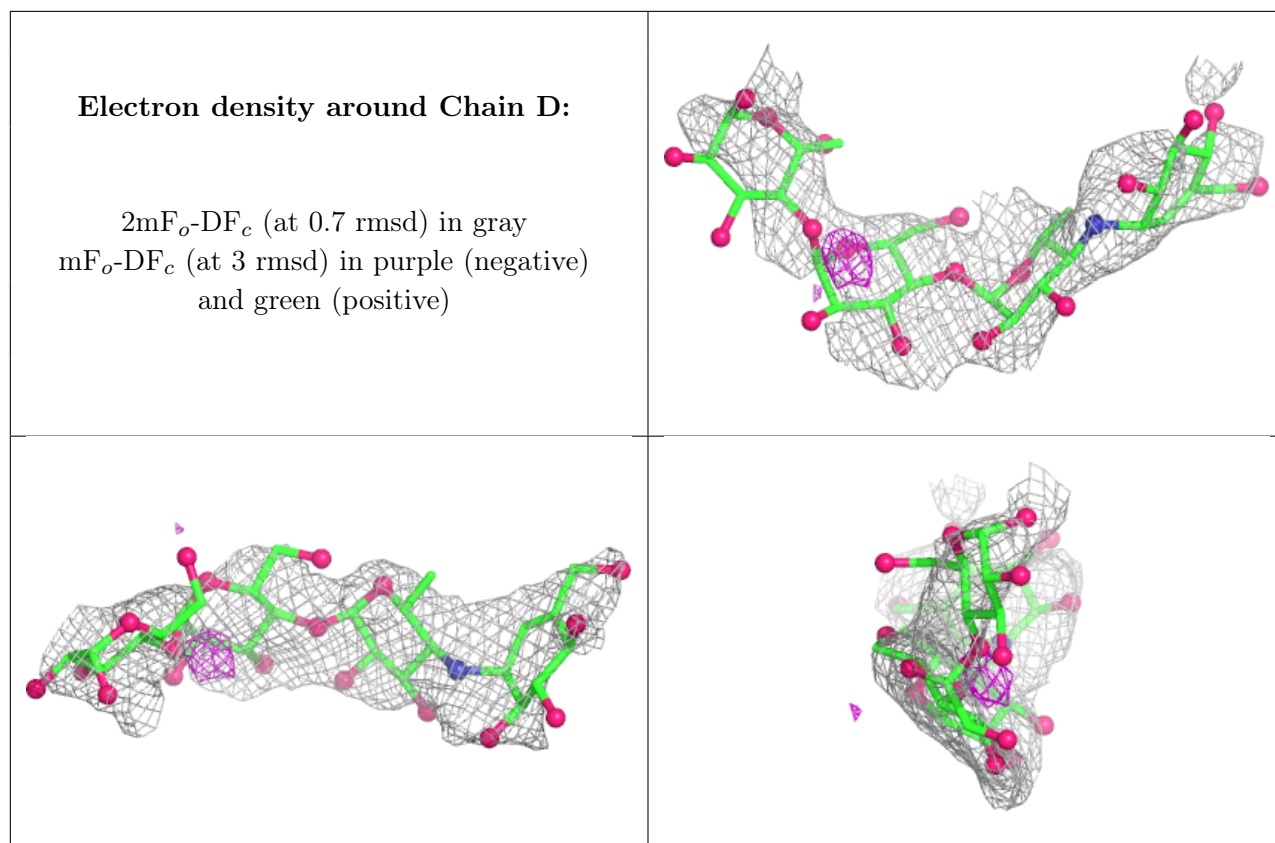
### 6.3 Carbohydrates [i](#)

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, 95<sup>th</sup> 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	GLC	D	1	12/12	0.44	0.11	150,157,165,166	0
2	GLC	C	1	12/12	0.46	0.11	159,167,172,173	0
2	AC1	C	3	21/22	0.66	0.16	127,151,168,171	0
2	GLC	C	2	11/12	0.66	0.14	148,165,168,169	0
2	GLC	D	2	11/12	0.71	0.16	146,152,157,158	0
2	AC1	D	3	21/22	0.74	0.14	121,135,145,148	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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	GLC	B	1001	12/12	0.77	0.22	106,121,127,135	0
3	GLC	A	1001	12/12	0.81	0.18	97,116,121,127	0
4	PLP	B	1002	15/16	0.91	0.10	59,62,69,75	0
4	PLP	A	1002	15/16	0.92	0.09	57,62,69,72	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.