

# Full wwPDB X-ray Structure Validation Report (i)

#### Nov 22, 2023 – 05:33 PM JST

PDB ID	:	7VWB
Title	:	Phloem lectin (PP2) structure -complex with N-Acetyllactosamine (LacNAc)
Authors	:	Sivaji, N.; Bobbili, K.B.; Suguna, K.; Surolia, A.
Deposited on		
Resolution	:	1.90 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 (i) 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 (1)) were used in the production of this report:

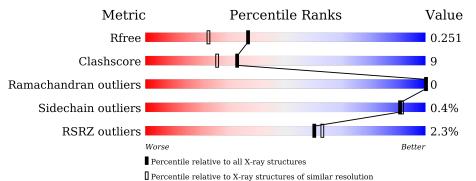
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 1.90 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R <sub>free</sub>	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 >=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 <=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	А	154	3% 86%	12% •
1	В	154	% <b>8</b> 6%	12% •
2	С	2	50% 50%	
2	D	2	100%	



#### 7VWB

# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2718 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 phloem lectin.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Δ	150	Total	С	Ν	0	S	0	0	0
1		150	1213	782	208	219	4	0		
1	P	150	Total	С	Ν	Ο	S	0	0	0
	D	100	1209	780	207	218	4	0	0	U

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	149	ILE	THR	conflict	UNP Q8LK69
В	149	ILE	THR	conflict	UNP Q8LK69

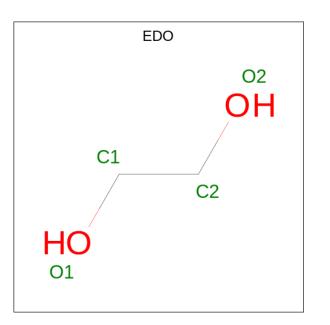
• Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-a lpha-D-glucopyranose.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
2	С	2	Total C N O 0	0	0	
		2	26 14 1 11	0	Ū	0
2	Л	n	Total C N O	0	0	0
2	D	2	26  14  1  11	0	0 0	0

• Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	А	1	Total 4	$\begin{array}{c} \mathrm{C} \\ \mathrm{2} \end{array}$	O 2	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	115	Total O 115 115	0	0
4	В	125	Total O 125 125	0	0



# 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.

Chain A:	86%	12%	·		
MET ALA GLY GLN S5 F11 F11	D59 N77 N82 N82 N82 N82 P92 P93 P93 P93 P93 P93 P93 P93 P13 P13 P13 P13 P13 P13 P13 P13 P13 P1				
• Molecule	1: phloem lectin				
Chain B:	86%	12%	·		
MET ALA GLY GLN S5 F11 F11	C34 836 836 836 836 845 845 817 817 817 812 812 812 812 812 812 812 812 812 812				
• Molecule	2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-a	alpha-I	)-gluc	opyra	nose
Chain C:	50% 50%		_		
GAL2					

• Molecule 1: phloem lectin

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

Chain D:

100%



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	37.69Å 60.39Å 82.98Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $99.13^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	24.90 - 1.90	Depositor
Resolution (A)	24.88 - 1.90	EDS
% Data completeness	98.0 (24.90-1.90)	Depositor
(in resolution range)	98.1 (24.88-1.90)	EDS
R <sub>merge</sub>	0.13	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	$4.65 (at 1.90 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
D D.	0.218 , $0.248$	Depositor
$R, R_{free}$	0.225 , $0.251$	DCC
$R_{free}$ test set	1468 reflections $(5.14\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	13.7	Xtriage
Anisotropy	0.771	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.38, 45.8	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.43, < L^2>=0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	2718	wwPDB-VP
Average B, all atoms $(Å^2)$	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 13.96% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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: GAL, NDG, EDO

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	Bo	nd lengths	Bond angles		
	Unam	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.74	1/1250~(0.1%)	0.84	0/1704	
1	В	0.73	1/1246~(0.1%)	0.88	0/1699	
All	All	0.74	2/2496~(0.1%)	0.86	0/3403	

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
1	А	98	GLU	CD-OE1	-5.91	1.19	1.25
1	В	98	GLU	CD-OE1	-5.48	1.19	1.25

There are no bond angle outliers.

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	А	1213	0	1158	21	0
1	В	1209	0	1152	22	0
2	С	26	0	21	1	0
2	D	26	0	21	0	0
3	А	4	0	6	0	0
4	А	115	0	0	4	0
4	В	125	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	2718	0	2358	42	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 9.

All (42) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance $(\text{\AA})$	overlap (Å)
1:A:11:PHE:CE1	1:A:149:ILE:HD13	2.05	0.91
1:B:11:PHE:CE1	1:B:149:ILE:HD13	2.13	0.84
4:A:407:HOH:O	1:B:129:THR:HG22	1.87	0.75
1:A:11:PHE:CD1	1:A:149:ILE:HD13	2.22	0.74
1:B:11:PHE:CD1	1:B:149:ILE:HD13	2.28	0.69
1:A:129:THR:HG22	4:A:411:HOH:O	1.94	0.68
1:B:149:ILE:HG13	1:B:150:LEU:N	2.15	0.62
1:A:149:ILE:HG13	1:A:150:LEU:N	2.17	0.60
1:A:11:PHE:CZ	1:A:149:ILE:HD13	2.38	0.58
1:B:45:GLN:HG2	4:B:311:HOH:O	2.03	0.57
1:A:149:ILE:HG13	1:A:150:LEU:O	2.04	0.57
1:A:11:PHE:CD1	1:A:149:ILE:CD1	2.86	0.57
1:B:11:PHE:CZ	1:B:149:ILE:HD13	2.38	0.57
1:B:149:ILE:HG13	1:B:150:LEU:O	2.05	0.57
1:B:10:ALA:N	1:B:149:ILE:HD11	2.20	0.56
1:B:11:PHE:CD1	1:B:149:ILE:CD1	2.91	0.53
1:A:10:ALA:N	1:A:149:ILE:HD11	2.25	0.52
1:A:86:HIS:O	1:A:134:PHE:HA	2.10	0.52
1:A:151:ALA:HB1	1:A:153:GLN:OE1	2.12	0.50
1:B:125:LEU:HD23	4:B:288:HOH:O	2.12	0.50
1:B:86:HIS:O	1:B:134:PHE:HA	2.11	0.49
1:A:113:LYS:HE2	1:B:36:TYR:CZ	2.47	0.48
1:B:151:ALA:HB1	1:B:153:GLN:OE1	2.13	0.47
1:B:10:ALA:C	1:B:149:ILE:HD12	2.35	0.47
1:B:92:PRO:HG2	1:B:125:LEU:HD11	1.97	0.47
1:A:10:ALA:C	1:A:149:ILE:HD12	2.35	0.46
1:A:11:PHE:HA	1:A:149:ILE:HD12	1.97	0.46
1:A:92:PRO:HG2	1:A:125:LEU:HD11	1.98	0.45
1:B:10:ALA:C	1:B:149:ILE:CD1	2.85	0.45
1:A:59:ASP:OD2	1:B:59:ASP:OD2	2.34	0.45
1:A:11:PHE:CA	1:A:149:ILE:HD12	2.47	0.45
1:B:11:PHE:CA	1:B:149:ILE:HD12	2.46	0.45
1:A:10:ALA:C	1:A:149:ILE:CD1	2.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:515:HOH:O	2:C:2:GAL:H4	2.16	0.44
1:B:11:PHE:HA	1:B:149:ILE:HD12	1.99	0.44
1:A:103:LEU:HA	1:A:106:ARG:HD2	2.00	0.43
1:A:134:PHE:HZ	1:A:147:ILE:HD13	1.83	0.43
1:A:82:ASN:HB2	4:A:490:HOH:O	2.18	0.43
1:B:11:PHE:N	1:B:149:ILE:HD12	2.32	0.43
1:B:134:PHE:HZ	1:B:147:ILE:HD13	1.85	0.41
1:A:11:PHE:N	1:A:149:ILE:HD12	2.36	0.40
1:B:80:GLY:HA2	4:B:287:HOH:O	2.20	0.40

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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	А	148/154~(96%)	144 (97%)	4(3%)	0	100 100
1	В	148/154~(96%)	144 (97%)	4 (3%)	0	100 100
All	All	296/308~(96%)	288~(97%)	8(3%)	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	А	127/131~(97%)	127~(100%)	0	100 100		
1	В	126/131~(96%)	125~(99%)	1 (1%)	81 82		
All	All	253/262~(97%)	252 (100%)	1 (0%)	91 91		

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	34	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	45	GLN
1	А	97	GLN
1	А	100	GLN
1	В	45	GLN
1	В	97	GLN

#### 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)

4 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	Turne	Chain	Res	Link	Bo	Bond lengths			Bond angles		
10101	Type	Unam	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
2	NDG	С	1	2	15,15,15	1.45	3 (20%)	21,21,21	0.90	0	
2	GAL	С	2	2	11,11,12	1.55	2 (18%)	$15,\!15,\!17$	1.07	1 (6%)	
2	NDG	D	1	2	15,15,15	1.75	3 (20%)	21,21,21	1.15	2 (9%)	
2	GAL	D	2	2	11,11,12	0.80	0	$15,\!15,\!17$	1.13	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	NDG	С	1	2	-	0/6/26/26	0/1/1/1
2	GAL	С	2	2	-	1/2/19/22	0/1/1/1
2	NDG	D	1	2	-	0/6/26/26	0/1/1/1
2	GAL	D	2	2	-	1/2/19/22	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	NDG	C1-C2	4.73	1.58	1.52
2	С	2	GAL	O4-C4	3.59	1.51	1.43
2	С	1	NDG	C1-C2	2.91	1.56	1.52
2	С	1	NDG	O1-C1	2.47	1.47	1.39
2	С	1	NDG	C3-C2	-2.27	1.48	1.53
2	D	1	NDG	O5-C1	2.22	1.48	1.42
2	С	2	GAL	O5-C1	-2.21	1.40	1.43
2	D	1	NDG	C2-N2	2.13	1.49	1.45

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	С	2	GAL	C3-C4-C5	-2.41	105.95	110.24
2	D	1	NDG	C2-N2-C7	-2.18	117.86	123.18
2	D	1	NDG	O3-C3-C2	-2.16	105.29	109.66
2	D	2	GAL	C1-O5-C5	-2.08	109.38	112.19
2	D	2	GAL	C3-C4-C5	-2.01	106.66	110.24

There are no chirality outliers.

All (2) torsion outliers are listed below:



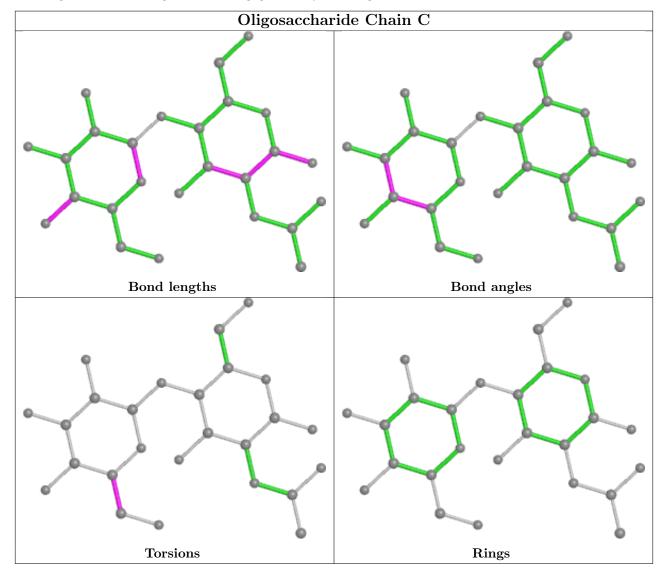
Mol	Chain	Res	Type	Atoms
2	С	2	GAL	O5-C5-C6-O6
2	D	2	GAL	O5-C5-C6-O6

There are no ring outliers.

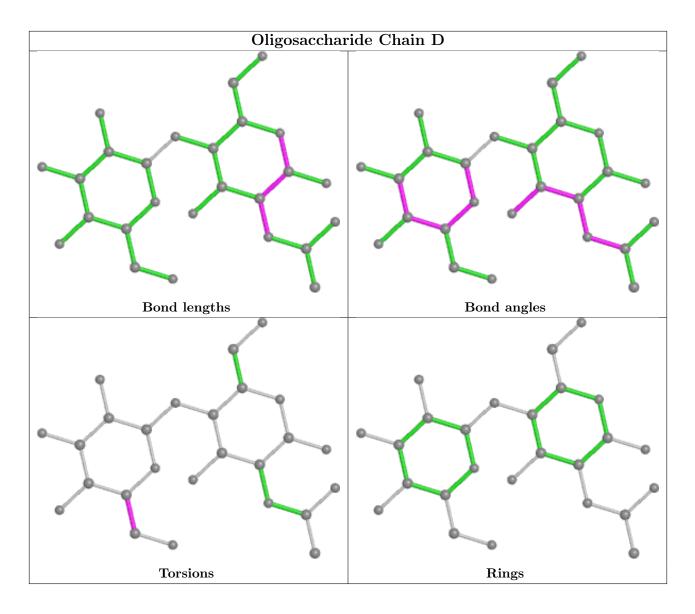
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	С	2	GAL	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)

1 ligand is 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		
	woi Type Chain I	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
3	EDO	А	301	-	3,3,3	0.21	0	$2,\!2,\!2$	0.28	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	EDO	А	301	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	А	301	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

#### 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,  $95^{th}$  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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	Q<0.9
1	А	150/154~(97%)	0.28	5 (3%) 46 49	7, 13, 30, 45	0
1	В	150/154~(97%)	0.18	2 (1%) 77 79	7, 13, 24, 35	0
All	All	300/308~(97%)	0.23	7 (2%) 60 63	7, 13, 28, 45	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	126	THR	3.4
1	А	77	ASN	3.3
1	А	124	SER	3.3
1	В	123	GLY	3.2
1	В	77	ASN	3.1
1	А	93	ASP	2.8
1	А	123	GLY	2.2

## 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,  $95^{th}$  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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
2	GAL	D	2	11/12	0.83	0.21	23,26,29,31	0
2	GAL	С	2	11/12	0.84	0.21	19,21,24,25	0

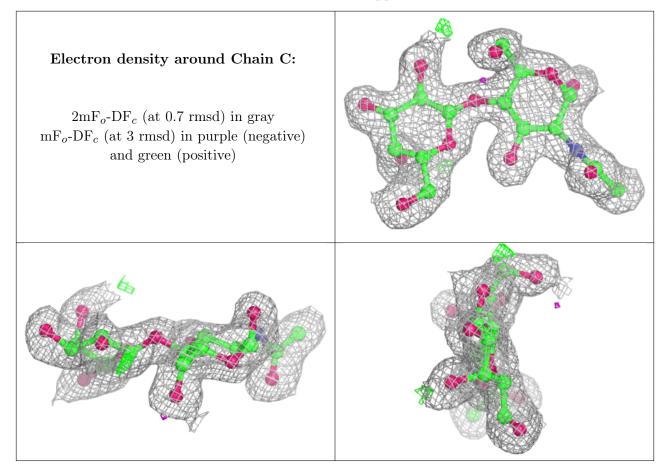
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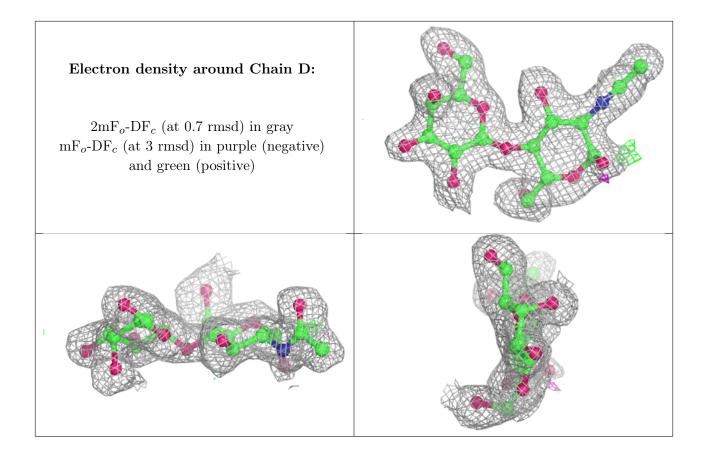
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(Å^2)$	Q<0.9
2	NDG	D	1	15/15	0.86	0.17	$19,\!22,\!27,\!27$	0
2	NDG	С	1	15/15	0.90	0.14	15,18,21,22	0

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,  $95^{th}$  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	$B-factors(A^2)$	Q < 0.9
3	EDO	A	301	4/4	0.93	0.15	$19,\!19,\!21,\!22$	0

## 6.5 Other polymers (i)

There are no such residues in this entry.

