

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

#### Oct 17, 2023 – 02:42 PM EDT

PDB ID	:	2DSU
Title	:	Binding of chitin-like polysaccharide to protective signalling factor: Crystal
		structure of the complex formed between signalling protein from sheep (SPS-
		40) with a tetrasaccharide at 2.2 A resolution
Authors	:	Srivastava, D.B.; Ethayathulla, A.S.; Kumar, J.; Singh, N.; Sharma, S.; Kaur,
		P.; Bhushan, A.; Singh, T.P.
Deposited on	:	2006-07-07
Resolution	:	2.20 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 (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:

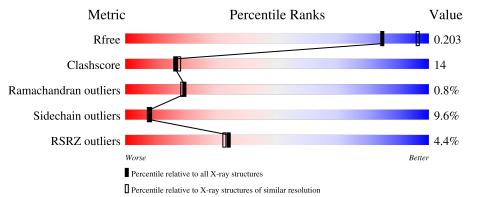
MolProbity Mogul		4.02b-467 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)		
$\mathrm{EDS}$	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal geometry (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-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	361	4%	77%	20%	•			
2	В	4	25%	75%					
3	С	4		100%					

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



Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NDG	В	2	-	-	-	Х
3	NDG	С	1	-	-	-	Х
3	NAG	С	2	-	-	Х	Х
3	NAG	С	3	-	-	-	Х

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



# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3144 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 Chitinase-3-like protein 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	361	Total 2869	C 1832	N 499	O 529	${ m S} 9$	0	0	0

• Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-beta-D-mannopyranos e-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-gluc opyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	В	4	Total 50	C 28	N 2	O 20	0	0	0

• Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-a cetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace
3	C	4	Total         C         N           57         32         4	O 21	0	0	0

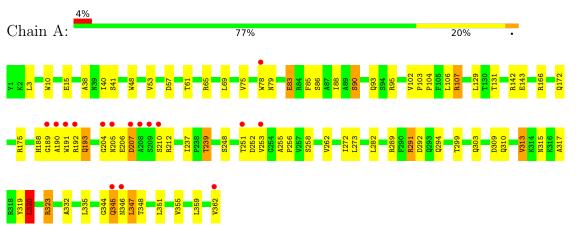
• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	168	Total         O           168         168	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.



• Molecule 1: Chitinase-3-like protein 1

• Molecule 2: alpha-D-mannopyranose-(1-4)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-a lpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain B:		25%					75%					
NAG1 NDG2 BMA3 MAN4												
	1		. 1	•	Б	,		1		 <b>a</b> 1		-

• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-gluc opyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-alpha-D-g lucopyranose

Chain C:

100%

NDG1 NAG2 NAG3 NAG3 NAG4



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	62.75Å 66.51Å 107.57Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	56.00 - 2.20	Depositor
Resolution (A)	45.64 - 2.20	EDS
% Data completeness	95.7 (56.00-2.20)	Depositor
(in resolution range)	95.9(45.64-2.20)	EDS
R <sub>merge</sub>	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$< I/\sigma(I) > 1$	$2.80 (at 2.20 \text{\AA})$	Xtriage
Refinement program	CNS 1.1	Depositor
D D.	0.187 , $0.209$	Depositor
$R, R_{free}$	0.187 , $0.203$	DCC
$R_{free}$ test set	446 reflections $(1.98\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	28.1	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.35 , $60.3$	EDS
L-test for twinning <sup>2</sup>	$ \langle L  \rangle = 0.49, \langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3144	wwPDB-VP
Average B, all atoms $(Å^2)$	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.15% 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: BMA, NDG, MAN, NAG

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		
	Unain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.38	0/2945	0.69	1/3995~(0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	320	LEU	CA-CB-CG	8.68	135.25	115.30

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	2869	0	2794	68	0
2	В	50	0	42	9	0
3	С	57	0	48	12	0
4	А	168	0	0	19	0
All	All	3144	0	2884	84	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 14.

The worst 5 of 84 close contacts within the same asymmetric unit are listed below, sorted by their



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:3:BMA:H61	2:B:4:MAN:C2	1.92	0.99
1:A:204:GLY:HA2	1:A:292:ASP:HB3	1.46	0.97
1:A:323:ARG:HH11	1:A:323:ARG:HG2	1.27	0.96
1:A:142:ARG:HD3	4:A:482:HOH:O	1.71	0.91
2:B:1:NAG:O3	2:B:2:NDG:H6C2	1.71	0.90

clash magnitude.

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	0		Outliers		
1	А	359/361~(99%)	343 (96%)	13~(4%)	3(1%)	19 19	

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	190	ALA
1	А	346	ASN
1	А	207	ASP

#### 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.

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Mol	Chain	Analysed	Analysed Rotameric Outliers		Percentiles	
Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	А	302/302~(100%)	273~(90%)	29 (10%)	8 8	

5 of 29 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	А	239	THR
1	А	351	LEU
1	А	253	VAL
1	А	323	ARG
1	А	252	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	А	345	GLN
1	А	315	ASN
1	А	294	GLN
1	А	205	ASN
1	А	303	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)

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



Mol	l Type Chain Res Link		Link	Bo	ond leng	ths	Bond angles			
MIOI	туре	Ullalli	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
2	NAG	В	1	1,2	14,14,15	0.96	1 (7%)	17,19,21	1.45	5 (29%)
2	NDG	В	2	2	14,14,15	1.16	3 (21%)	17,19,21	1.09	1 (5%)
2	BMA	В	3	2	11,11,12	1.03	2 (18%)	15,15,17	2.84	5 (33%)
2	MAN	В	4	2	11,11,12	0.73	0	15,15,17	0.75	0
3	NDG	С	1	3	15,15,15	0.54	0	21,21,21	1.14	2 (9%)
3	NAG	С	2	3	14,14,15	0.91	0	17,19,21	1.18	2 (11%)
3	NAG	С	3	3	14,14,15	1.01	0	17,19,21	1.05	1 (5%)
3	NAG	С	4	3	14,14,15	0.95	1 (7%)	17,19,21	1.50	2 (11%)

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).

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	В	1	1,2	-	3/6/23/26	0/1/1/1
2	NDG	В	2	2	-	4/6/23/26	0/1/1/1
2	BMA	В	3	2	-	2/2/19/22	0/1/1/1
2	MAN	В	4	2	-	0/2/19/22	1/1/1/1
3	NDG	С	1	3	-	4/6/26/26	0/1/1/1
3	NAG	С	2	3	-	0/6/23/26	0/1/1/1
3	NAG	С	3	3	-	4/6/23/26	0/1/1/1
3	NAG	С	4	3	_	4/6/23/26	0/1/1/1

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

Mol	Chain	Res	Type	Atoms	Z	$Observed(\text{\AA})$	$\mathrm{Ideal}(\mathrm{\AA})$
3	С	4	NAG	C1-C2	3.01	1.56	1.52
2	В	2	NDG	C1-C2	2.56	1.56	1.52
2	В	1	NAG	C4-C3	2.18	1.57	1.52
2	В	3	BMA	C1-C2	2.17	1.57	1.52
2	В	2	NDG	C4-C3	2.12	1.57	1.52

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



Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	В	3	BMA	C3-C4-C5	-6.80	98.12	110.24
2	В	3	BMA	C2-C3-C4	-5.48	101.42	110.89
3	С	4	NAG	C1-C2-N2	4.06	117.43	110.49
2	В	3	BMA	C1-C2-C3	3.95	114.53	109.67
3	С	1	NDG	C4-C3-C2	-3.19	105.67	110.34

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	1	NAG	C3-C2-N2-C7
2	В	1	NAG	C8-C7-N2-C2
2	В	1	NAG	O7-C7-N2-C2
2	В	2	NDG	C8-C7-N2-C2
2	В	2	NDG	O7-C7-N2-C2

All (1) ring outliers are listed below:

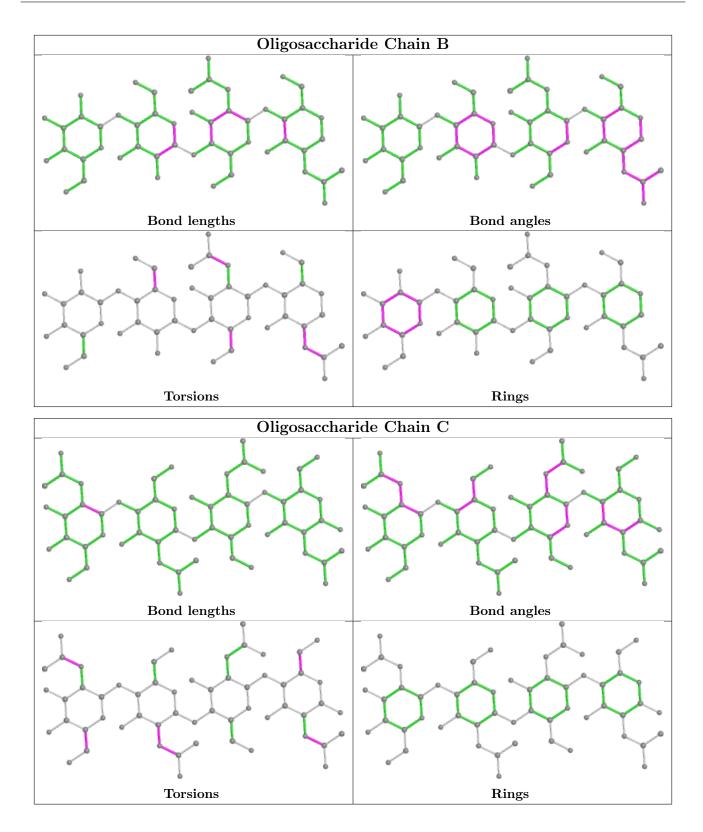
Mol	Chain	Res	Type	Atoms
2	В	4	MAN	C1-C2-C3-C4-C5-O5

8 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	С	3	NAG	3	0
2	В	3	BMA	5	0
3	С	4	NAG	2	0
2	В	4	MAN	5	0
2	В	2	NDG	1	0
3	С	2	NAG	8	0
2	В	1	NAG	4	0
3	С	1	NDG	4	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)

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,  $95^{th}$  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	<RSRZ $>$	#RSRZ>2	$OWAB(Å^2)$	Q < 0.9
1	А	361/361~(100%)	-0.13	16 (4%) 34 32	14, 28, 59, 98	0

The worst 5 of 16 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	208	ALA	15.1
1	А	191	TRP	9.2
1	А	189	GLY	6.5
1	А	207	ASP	6.5
1	А	209	SER	5.3

### 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	BMA	В	3	11/12	0.29	0.38	$55,\!56,\!56,\!56$	11
2	MAN	В	4	11/12	0.45	0.40	54,55,56,56	11
3	NAG	С	3	14/15	0.53	0.54	$53,\!55,\!57,\!57$	14
3	NDG	С	1	15/15	0.63	0.55	52,58,61,61	0
3	NAG	С	2	14/15	0.67	0.66	59,63,67,67	0
2	NDG	В	2	14/15	0.67	0.43	50,55,59,60	14
3	NAG	С	4	14/15	0.76	0.35	52,54,56,56	14

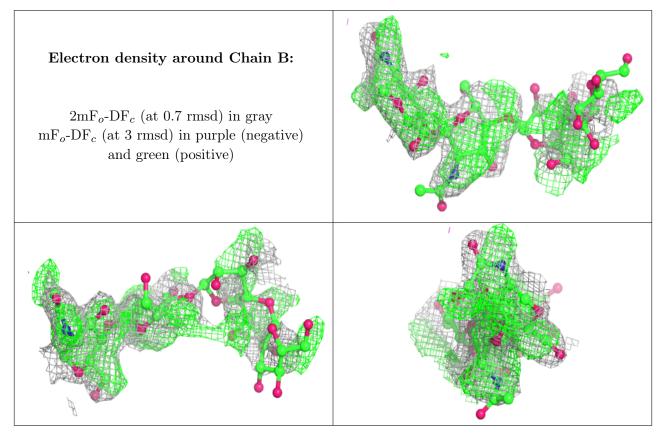
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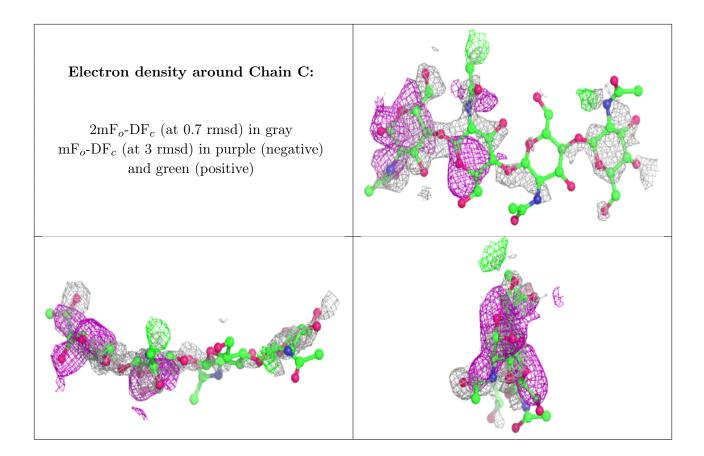
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
2	NAG	В	1	14/15	0.78	0.33	$26,\!32,\!37,\!44$	14

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.

