

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

#### Jan 8, 2024 - 01:58 am GMT

PDB ID	:	5MRI
Title	:	Crystal structure of the Vps10p domain of human sortilin/NTS3 in complex
		with Triazolone 18
Authors	:	Andersen, J.L.; Strandbygaard, D.; Thirup, S.
Deposited on	:	2016-12-23
Resolution	:	2.00  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:

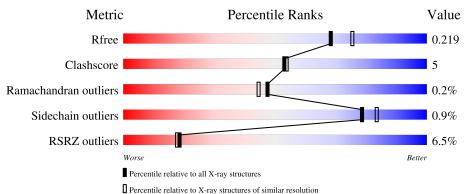
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as $541$ be (2020)
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal geometry (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 2.00 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	696	<mark>6%</mark> 83%	12%	5%
2	В	3	100%		
3	С	2	50% 50%		



# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5686 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.

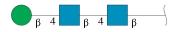
• Molecule 1 is a protein called Sortilin.

Mol	Chain	Residues		$\mathbf{A}$	toms			ZeroOcc	AltConf	Trace
1	А	662	Total 5243	C 3321	N 875	O 1017	S 30	0	7	0

Chain	Residue	Modelled	Actual	Comment	Reference
А	617	MET	VAL	conflict	UNP Q99523
А	724	GLY	-	expression tag	UNP Q99523
А	725	SER	-	expression tag	UNP Q99523
А	726	ALA	-	expression tag	UNP Q99523
А	727	MET	-	expression tag	UNP Q99523
А	728	ILE	-	expression tag	UNP Q99523
А	729	GLU	-	expression tag	UNP Q99523
А	730	GLY	-	expression tag	UNP Q99523
A	731	ARG	-	expression tag	UNP Q99523
A	732	GLY	-	expression tag	UNP Q99523
А	733	VAL	-	expression tag	UNP Q99523
А	734	GLY	-	expression tag	UNP Q99523
А	735	HIS	-	expression tag	UNP Q99523
А	736	HIS	-	expression tag	UNP Q99523
А	737	HIS	-	expression tag	UNP Q99523
А	738	HIS	-	expression tag	UNP Q99523
А	739	HIS	-	expression tag	UNP Q99523
А	740	HIS	-	expression tag	UNP Q99523

There are 18 discrepancies between the modelled and reference sequences:

• Molecule 2 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-b eta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.





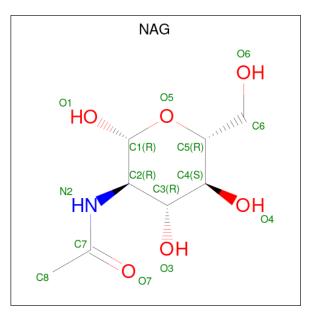
[	Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
	2	В	3	Total 39	C 22	N 2	O 15	0	0	0

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



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace	
3	С	2	Total 28	C N 16 2	0 10	0	0	0

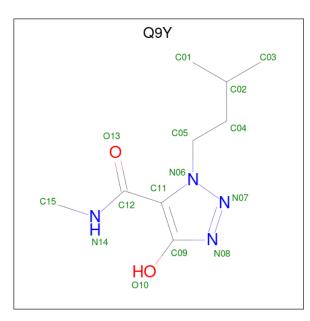
• Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



Mol	Chain	Residues	A	tor	ns		ZeroOcc	AltConf
4	А	1	Total 14	C 8	N 1	O 5	0	0

• Molecule 5 is {N}-methyl-3-(3-methylbutyl)-5-oxidanyl-1,2,3-triazole-4-carboxamide (three-letter code: Q9Y) (formula:  $C_9H_{16}N_4O_2$ ).





I	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	5	А	1	$\begin{array}{ccccc} \text{Total} & \text{C} & \text{N} & \text{O} \\ 15 & 9 & 4 & 2 \end{array}$	0	0

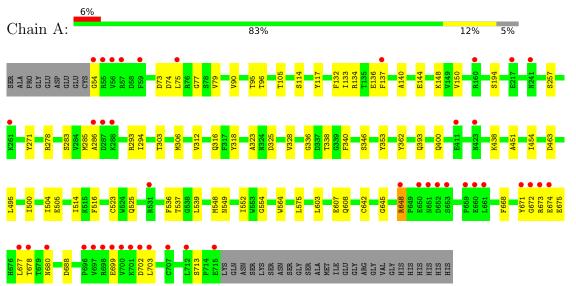
• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	347	Total         O           347         347	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: Sortilin

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

Chain B:	100%
NAG1 BMA3	

• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:	50%	50%





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	158.30Å 81.06Å 105.69Å	Deperitor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $131.10^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	43.91 - 2.00	Depositor
Resolution (A)	64.32 - 2.00	EDS
% Data completeness	98.4 (43.91-2.00)	Depositor
(in resolution range)	98.4 (64.32-2.00)	EDS
R <sub>merge</sub>	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.64 (at 2.00 \text{\AA})$	Xtriage
Refinement program	PHENIX (dev_2614: ???)	Depositor
$R, R_{free}$	0.195 , $0.219$	Depositor
II, IIfree	0.195 , $0.219$	DCC
$R_{free}$ test set	2010 reflections $(3.00%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	35.2	Xtriage
Anisotropy	0.670	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	$0.37 \;,  50.3$	EDS
L-test for twinning <sup>2</sup>	$< L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	0.023 for -h-2*l,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5686	wwPDB-VP
Average B, all atoms $(Å^2)$	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.69% 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: Q9Y, BMA, 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	Bo	ond angles
	Unain	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	А	0.26	0/5388	0.52	2/7301~(0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	278	ARG	NE-CZ-NH1	-5.86	117.37	120.30
1	А	648	ARG	C-N-CD	-5.06	109.46	120.60

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	5243	0	5061	56	0
2	В	39	0	34	0	0
3	С	28	0	25	2	0
4	А	14	0	13	1	0
5	А	15	0	0	1	0
6	А	347	0	0	13	0
All	All	5686	0	5133	57	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 5.

The worst 5 of 57 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:607:GLU:OE2	6:A:901:HOH:O	2.03	0.74
1:A:713:SER:OG	6:A:902:HOH:O	2.06	0.74
1:A:438:LYS:NZ	6:A:911:HOH:O	2.23	0.71
1:A:514:ILE:HD12	1:A:552:ILE:HD11	1.73	0.69
1:A:303:THR:HB	1:A:673:ARG:HH12	1.58	0.69

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	Percenti	les
1	А	667/696~(96%)	649~(97%)	17 (2%)	1 (0%)	51 49	9

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type	
1	А	144	GLU	

#### 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	585/605~(97%)	580~(99%)	5(1%)	78 83	

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	257	SER
1	А	325	ASP
1	А	362	TYR
1	А	393	GLN
1	А	699	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)

5 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	ype Chain Res Link		Bo	Bond lengths			Bond angles		
	Type	Unam	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
2	NAG	В	1	2,1	14,14,15	0.32	0	$17,\!19,\!21$	0.49	0
2	NAG	В	2	2	14,14,15	0.31	0	17,19,21	0.43	0
2	BMA	В	3	2	11,11,12	0.53	0	$15,\!15,\!17$	0.90	0
3	NAG	С	1	1,3	14,14,15	0.53	0	17,19,21	0.78	1 (5%)
3	NAG	С	2	3	14,14,15	0.25	0	17,19,21	0.56	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	NAG	В	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	В	2	2	-	0/6/23/26	0/1/1/1
2	BMA	В	3	2	-	2/2/19/22	0/1/1/1
3	NAG	С	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	С	2	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	С	1	NAG	C1-O5-C5	2.19	115.16	112.19

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	3	BMA	O5-C5-C6-O6
2	В	3	BMA	C4-C5-C6-O6
3	С	1	NAG	C4-C5-C6-O6
3	С	1	NAG	O5-C5-C6-O6

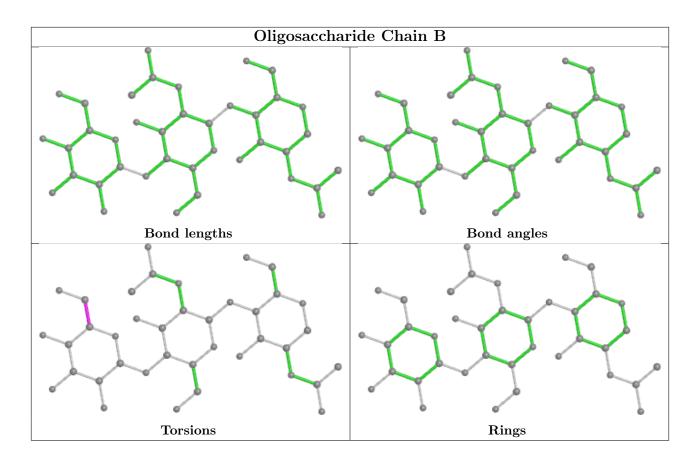
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	С	1	NAG	2	0

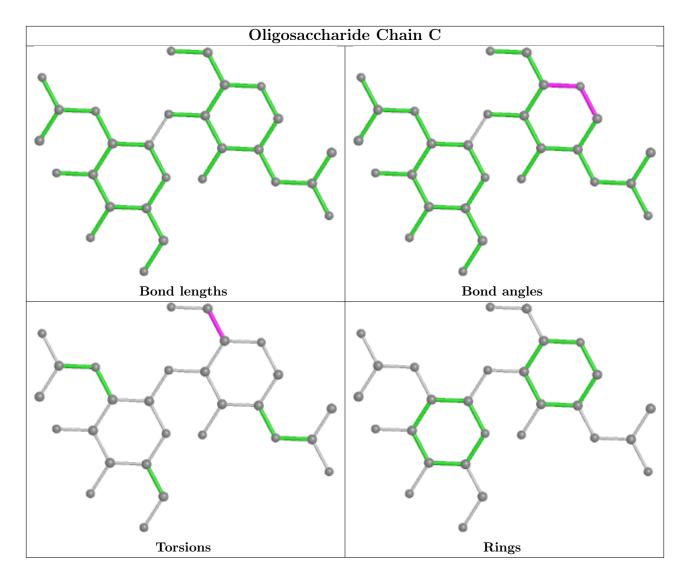
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











### 5.6 Ligand geometry (i)

2 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	Bo	Bond lengths			Bond angles		
IVIOI			nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
5	Q9Y	А	807	-	$13,\!15,\!15$	3.07	3 (23%)	10,20,20	2.16	5 (50%)	
4	NAG	А	806	1	14,14,15	0.32	0	17,19,21	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
5	Q9Y	А	807	-	-	3/7/11/11	0/1/1/1
4	NAG	А	806	1	-	1/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
5	А	807	Q9Y	C12-N14	9.59	1.45	1.33
5	А	807	Q9Y	C05-N06	-4.55	1.40	1.47
5	А	807	Q9Y	O10-C09	2.22	1.39	1.29

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
5	А	807	Q9Y	C11-N06-N07	-3.88	108.06	112.63
5	А	807	Q9Y	C15-N14-C12	-3.25	118.20	121.89
5	А	807	Q9Y	C04-C05-N06	-2.72	106.26	111.62
5	А	807	Q9Y	O13-C12-N14	-2.46	118.75	122.47
5	А	807	Q9Y	C05-N06-N07	2.09	125.68	119.93

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	$\operatorname{Res}$	Type	Atoms
5	А	807	Q9Y	C03-C02-C04-C05
5	А	807	Q9Y	C02-C04-C05-N06
5	А	807	Q9Y	C01-C02-C04-C05
4	А	806	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	А	807	Q9Y	1	0
4	А	806	NAG	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,  $95^{th}$  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	662/696~(95%)	0.24	43 (6%) 18 18	27, 46, 84, 136	0

The worst 5 of 43 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	697	VAL	9.7
1	А	701	LYS	6.3
1	А	54	GLY	6.2
1	А	55	ARG	6.2
1	А	286	ALA	6.0

## 6.2 Non-standard residues in protein, DNA, RNA chains (i)

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

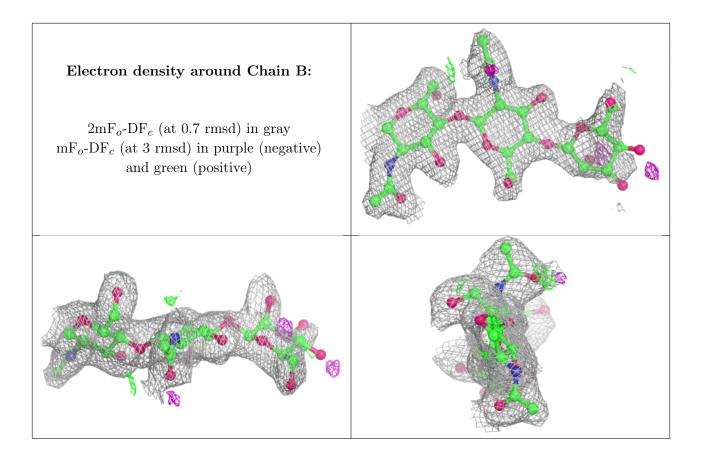
## 6.3 Carbohydrates (i)

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.

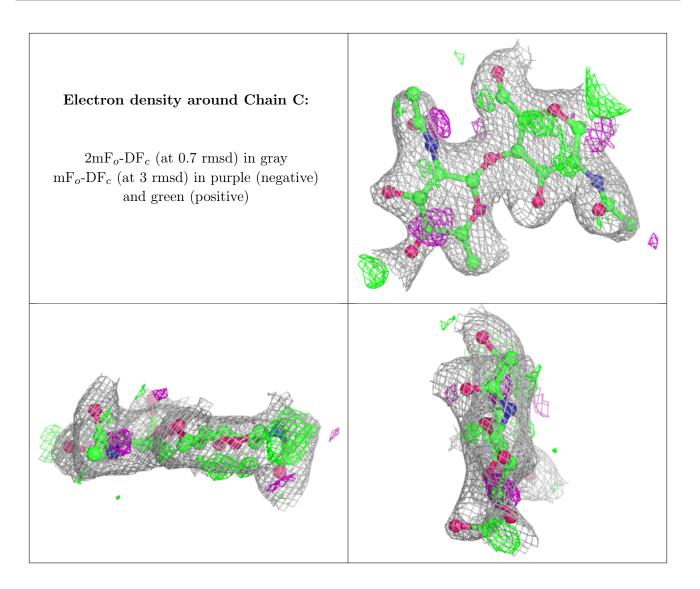
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
2	BMA	В	3	11/12	0.63	0.28	87,94,97,99	0
3	NAG	С	2	14/15	0.79	0.27	56,67,73,74	0
2	NAG	В	2	14/15	0.90	0.12	41,56,66,79	0
3	NAG	С	1	14/15	0.91	0.14	33,40,42,43	0
2	NAG	В	1	14/15	0.96	0.12	27,32,36,40	0

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	$B-factors(Å^2)$	Q<0.9
4	NAG	А	806	14/15	0.87	0.20	59,69,73,74	0
5	Q9Y	А	807	15/15	0.93	0.17	34,37,64,65	0

### 6.5 Other polymers (i)

There are no such residues in this entry.

