

Full wwPDB X-ray Structure Validation Report (i)

Sep 25, 2023 – 04:51 PM EDT

PDB ID	:	6ANY
Title	:	Structure of BmVAL-1
Authors	:	Asojo, O.A.
Deposited on	:	2017-08-14
Resolution	:	2.25 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:

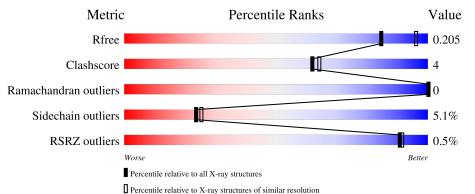
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	206	86%		12%	•		
2	В	2	50%	50%				
3	С	4	50%	50%				

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
2	FUC	В	2	-	-	-	Х
3	BMA	С	3	-	-	-	Х



6ANY

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 1883 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 Bm4233, isoform b.

Mol	Chain	Residues		Atoms			ZeroOcc	AltConf	Trace	
1	А	206	Total 1630	C 1012	N 307	0 297	S 14	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

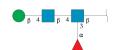
Chair	n Residue	Modelled	Actual	Comment	Reference
A	221	GLU	-	expression tag	UNP O44932
А	222	LEU	-	expression tag	UNP O44932

• Molecule 2 is an oligosaccharide called alpha-L-fucopyranose-(1-3)-2-acetamido-2-deoxy-bet a-D-glucopyranose.



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
2	В	2	Total 24	C 14	N 1	O 9	0	0	0

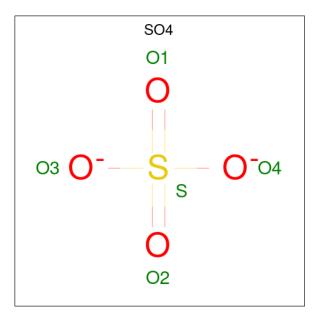
• Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-b eta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopy ranose.



Mo	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
3	С	4	Total 49	C 28	N 2	0 19	0	0	0



• Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



\mathbf{M}	ol	Chain	Residues	Ato	\mathbf{ms}		ZeroOcc	AltConf
4	:	А	1	Total 5	O 4	S 1	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	А	175	Total (175 17) 75	0	0



NAG1 FUC2

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% •	
F17 28 K30 K30 F39 F39 K47	R51 154 154 189 190 190 191 8126 8126 8126 8126 8126 1131 7131 7131 7131 8128	L139 R145 H149 H150 F151 F152 F151 F152 F154 F154 F154 F154 F154 F154 F156 F154 F156 F156 F156 F156 F156 F156 F156 F156	
• Molecule 2: alp	pha-L-fucopyranose-(1-3)	-2-acetamido-2-deoxy-beta-D-gluo	copyranose
Chain B:	50%	50%	•

• Molecule 1: Bm4233, isoform b

 • Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alp ha-L-fucopyranose-(1-3)] 2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:	50%	50%
NAG1 NAG2 BMA3 FUC4		



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants	85.79Å 85.79Å 66.67Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.87 - 2.25	Depositor
Resolution (A)	44.87 - 2.25	EDS
% Data completeness	95.8 (44.87-2.25)	Depositor
(in resolution range)	95.8 (44.87-2.25)	EDS
R _{merge}	0.03	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$3.58 (at 2.24 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
D D.	0.180 , 0.214	Depositor
R, R_{free}	0.181 , 0.205	DCC
R_{free} test set	977 reflections (8.29%)	wwPDB-VP
Wilson B-factor $(Å^2)$	20.5	Xtriage
Anisotropy	0.089	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 35.1	EDS
L-test for twinning ²	$ \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	0.94	EDS
Total number of atoms	1883	wwPDB-VP
Average B, all atoms $(Å^2)$	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.81% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities 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, FUC, SO4, 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		lengths	Bond angles		
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.93	0/1664	0.91	5/2238~(0.2%)	

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	145	ARG	NE-CZ-NH2	-6.84	116.88	120.30
1	А	174	ARG	NE-CZ-NH1	6.56	123.58	120.30
1	А	222	LEU	CB-CG-CD1	-6.23	100.40	111.00
1	А	205	ASP	CB-CG-OD2	-5.47	113.38	118.30
1	А	219	ARG	NE-CZ-NH2	5.05	122.82	120.30

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	1630	0	1609	15	0
2	В	24	0	22	1	0
3	С	49	0	43	0	0
4	А	5	0	0	0	0
5	А	175	0	0	2	0
All	All	1883	0	1674	15	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 4.

All (15) 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:90:ASN:ND2	5:A:401:HOH:O	2.13	0.81
1:A:133:GLN:HG2	2:B:1:NAG:H82	1.86	0.57
1:A:222:LEU:HD23	1:A:222:LEU:N	2.24	0.52
1:A:133:GLN:HG3	1:A:135:PRO:HD3	1.91	0.51
1:A:183:SER:HA	1:A:184:PRO:C	2.35	0.46
1:A:159:THR:HA	1:A:184:PRO:HD2	1.98	0.45
1:A:139:LEU:CD2	1:A:187:ASN:HA	2.47	0.45
1:A:17:PHE:CD2	1:A:30:LYS:HG2	2.52	0.45
1:A:17:PHE:N	5:A:411:HOH:O	2.51	0.43
1:A:149:LEU:HA	1:A:152:THR:OG1	2.18	0.43
1:A:39:PHE:HB3	1:A:124:TRP:CZ2	2.54	0.42
1:A:39:PHE:HB3	1:A:124:TRP:CH2	2.55	0.42
1:A:126:GLU:OE2	1:A:150:HIS:ND1	2.42	0.42
1:A:17:PHE:CE2	1:A:30:LYS:HG2	2.54	0.42
1:A:51:ARG:HB2	1:A:131:TYR:O	2.21	0.40

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	204/206~(99%)	203 (100%)	1 (0%)	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	176/176~(100%)	167~(95%)	9~(5%)	24 25	

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	28	GLN
1	А	36	ASN
1	А	47	LYS
1	А	51	ARG
1	А	54	THR
1	А	89	ARG
1	А	91	GLN
1	А	133	GLN
1	А	145	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type	
1	А	90	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)

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

5.5 Carbohydrates (i)

6 monosaccharides are modelled in this entry.



6ANY

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		
Mol Type	Chain	lain nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
2	NAG	В	1	1,2	14,14,15	0.29	0	17,19,21	0.61	0
2	FUC	В	2	2	10,10,11	0.30	0	14,14,16	0.72	0
3	NAG	С	1	1,3	14,14,15	0.55	0	17,19,21	0.48	0
3	NAG	С	2	3	14,14,15	0.37	0	17,19,21	0.99	1 (5%)
3	BMA	С	3	3	11,11,12	0.23	0	$15,\!15,\!17$	0.67	0
3	FUC	С	4	3	10,10,11	0.29	0	$14,\!14,\!16$	0.93	1 (7%)

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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
3	С	4	FUC	C2-C3-C4	-2.39	106.76	110.89
3	С	2	NAG	C4-C3-C2	-2.05	108.01	111.02

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	С	3	BMA	C4-C5-C6-O6

Continued on next page...



Mol	Chain	Res	Type	Atoms
3	С	2	NAG	C8-C7-N2-C2
3	С	3	BMA	O5-C5-C6-O6
3	С	2	NAG	O7-C7-N2-C2
2	В	1	NAG	C4-C5-C6-O6
2	В	1	NAG	O5-C5-C6-O6
3	С	2	NAG	O5-C5-C6-O6

Continued from previous page...

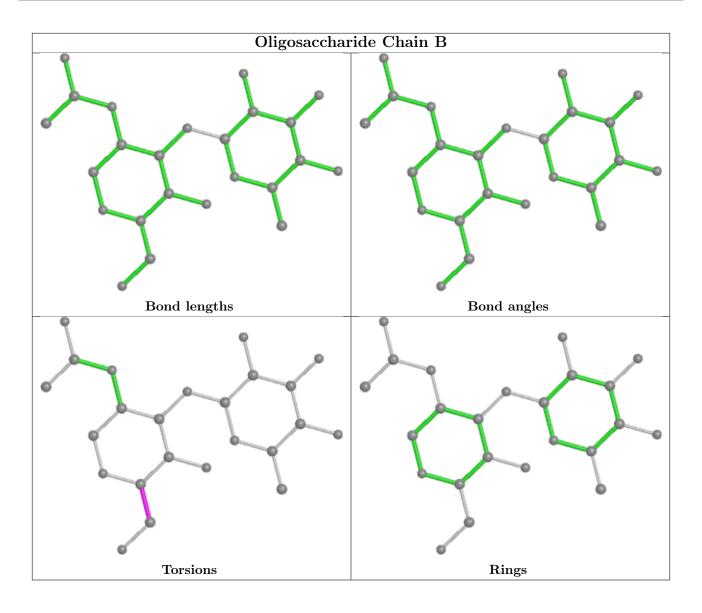
There are no ring outliers.

1 monomer is involved in 1 short contact:

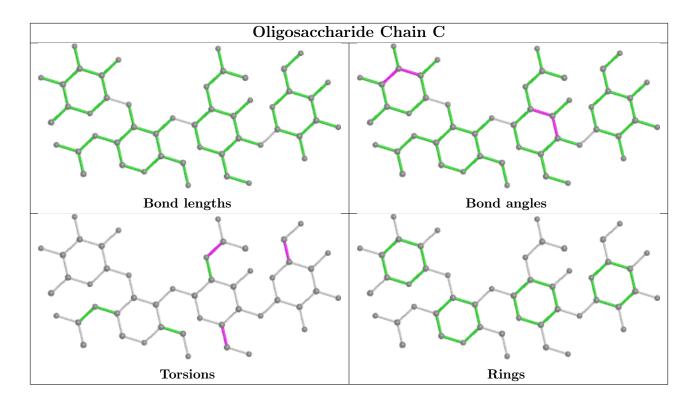
M	bl	Chain	Res	Type	Clashes	Symm-Clashes
2		В	1	NAG	1	0

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	Mol Type Chain	Chain	Res	Link	Bond lengths			Bond angles		
IVIOI		Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
4	SO4	А	307	-	4,4,4	0.20	0	$6,\!6,\!6$	0.79	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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	<RSRZ $>$	#RSRZ>2	$OWAB(Å^2)$	Q<0.9
1	А	206/206~(100%)	-0.36	1 (0%) 91 91	11, 18, 32, 49	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ	
1	А	207	ASN	2.7	

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

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

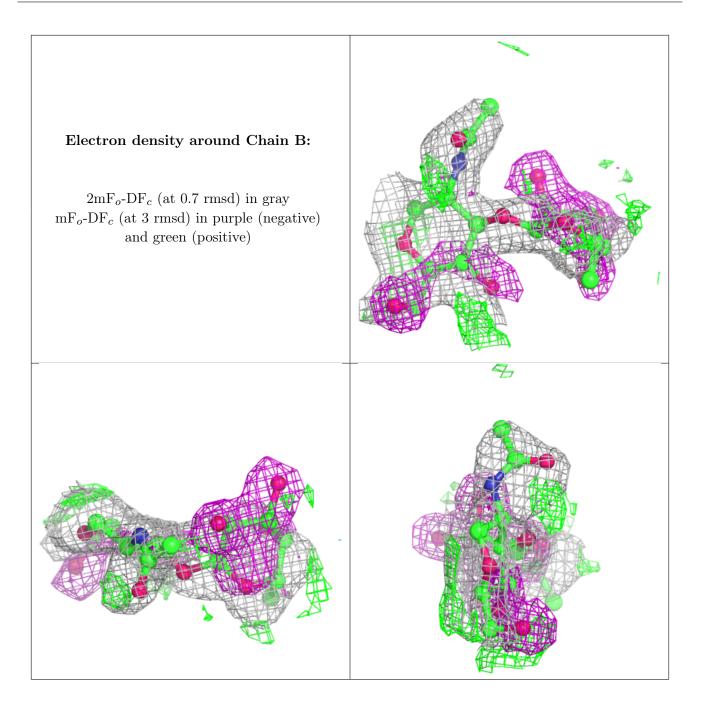
6.3 Carbohydrates (i)

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.

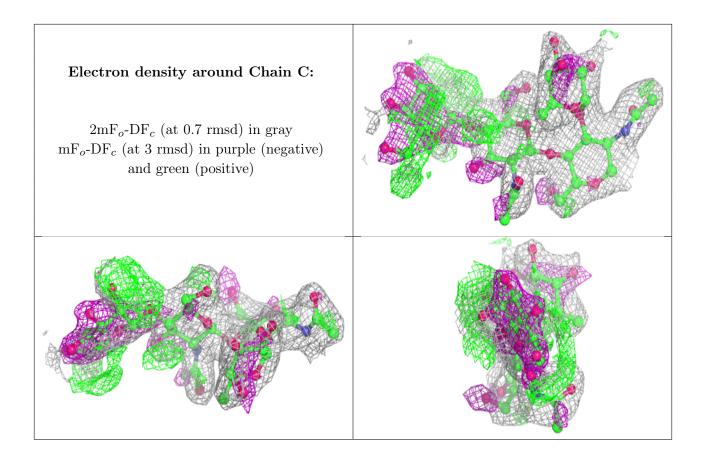
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	BMA	С	3	11/12	0.29	0.56	$30,\!30,\!30,\!30$	0
2	FUC	В	2	10/11	0.65	0.61	30,30,30,30	0
2	NAG	В	1	14/15	0.72	0.30	30,30,30,30	0
3	NAG	С	2	14/15	0.75	0.35	30,30,30,30	0
3	FUC	С	4	10/11	0.81	0.34	30,30,30,30	0
3	NAG	\mathbf{C}	1	14/15	0.83	0.21	30,30,30,30	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	SO4	А	307	5/5	0.98	0.13	$32,\!35,\!38,\!39$	0

6.5 Other polymers (i)

There are no such residues in this entry.

