

wwPDB X-ray Structure Validation Summary Report (i)

May 15, 2023 – 02:27 PM EDT

PDB ID : 8CWI

Title: Fab arm of antibody 10G4 bound to CoV-2 receptor binding domain (RBD)

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Deposited on : 2022-05-19

Resolution : 1.87 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: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.32.2

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

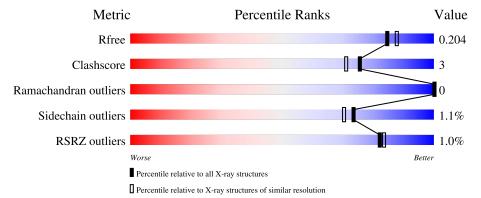
 $\begin{tabular}{lll} Validation Pipeline (wwPDB-VP) & : & 2.32.2 \end{tabular}$

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 1.87 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	Similar resolution $(\# \text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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	A	204	88%	6% • •
2	Н	228	89%	6% •
3	L	215	93%	6%
4	В	6	50% 50%	
5	С	3	100%	



2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 5666 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 Spike protein S1.

Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf	Trace	
1	Δ	195	Total	С	N	О	S	0	7	0	
1	Λ	130	1579	1009	264	298	8		'		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	529	GLY	-	expression tag	UNP P0DTC2
A	530	SER	-	expression tag	UNP P0DTC2
A	531	HIS	-	expression tag	UNP P0DTC2
A	532	HIS	-	expression tag	UNP P0DTC2
A	533	HIS	-	expression tag	UNP P0DTC2
A	534	HIS	-	expression tag	UNP P0DTC2
A	535	HIS	-	expression tag	UNP P0DTC2
A	536	HIS	-	expression tag	UNP P0DTC2

• Molecule 2 is a protein called Heavy chain of Fab arm of antibody 10G4.

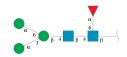
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	Н	218	Total 1650	C 1042	N 279	O 323	S 6	0	3	0

• Molecule 3 is a protein called Light chain of Fab arm of antibody 10G4.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
3	L	214	Total 1656	C 1035	N 282	O 332	S 7	0	2	0

• Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
4	В	6	Total 71	C 40	N 2	O 29	0	0	0

• Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[al pha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
5	С	3	Total C N O 38 22 2 14	0	0	0

• Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Mg 1 1	0	0
6	L	1	Total Mg 1 1	0	0

• Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

ſ	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	7	Н	2	Total Cl 2 2	0	0
	7	L	1	Total Cl 1 1	0	0

• Molecule 8 is water.

Mo	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	230	Total O 230 230	0	0
8	Н	249	Total O 249 249	0	0



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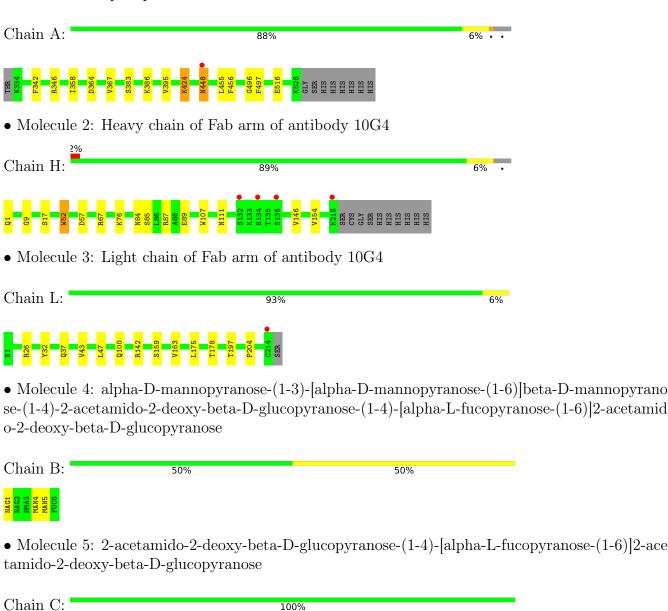
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	188	Total 188	O 188	0	0



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.







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	53.69Å 88.64Å 81.40Å	Donositor
a, b, c, α , β , γ	90.00° 95.03° 90.00°	Depositor
Resolution (Å)	46.56 - 1.87	Depositor
Resolution (A)	46.56 - 1.87	EDS
% Data completeness	98.4 (46.56-1.87)	Depositor
(in resolution range)	98.4 (46.56-1.87)	EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.41 (at 1.87Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
D D.	0.170 , 0.205	Depositor
R, R_{free}	0.170 , 0.204	DCC
R_{free} test set	3038 reflections $(4.94%)$	wwPDB-VP
Wilson B-factor (Å ²)	27.2	Xtriage
Anisotropy	0.379	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.31 , 43.1	EDS
L-test for twinning ²	$ < L > = 0.50, < L^2> = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5666	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.26% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ 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: NAG, FUC, CL, MAN, BMA, MG

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		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.36	0/1623	0.54	1/2211 (0.0%)	
2	Н	0.35	0/1691	0.57	0/2306	
3	L	0.34	0/1693	0.56	0/2301	
All	All	0.35	0/5007	0.56	1/6818 (0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	455	LEU	CA-CB-CG	5.11	127.05	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	A	1579	0	1472	11	1
2	Н	1650	0	1589	11	0
3	L	1656	0	1591	11	0
4	В	71	0	61	1	0
5	С	38	0	34	0	0
6	A	1	0	0	0	0
6	L	1	0	0	0	0



Continued	trom	mmoninonic	maaa
COHABABACA		DIEUIUU	DUIUE
0 0 1000100000			

Mol	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
7	Н	2	0	0	0	0
7	L	1	0	0	1	0
8	A	230	0	0	2	0
8	Н	249	0	0	2	1
8	L	188	0	0	1	0
All	All	5666	0	4747	30	1

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 3.

The worst 5 of 30 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
2:H:9:GLY:H	2:H:111:ASN:ND2	1.83	0.77
2:H:76:LYS:NZ	8:H:401:HOH:O	2.20	0.75
2:H:9:GLY:H	2:H:111:ASN:HD22	1.41	0.69
2:H:87:ARG:HG3	2:H:89:GLU:HG2	1.82	0.61
1:A:364:ASP:O	1:A:367:VAL:HG12	2.02	0.59

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:516:GLU:OE2	8:H:601:HOH:O[2_546]	2.12	0.08

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	Perce	ntiles
1	A	$200/204\ (98\%)$	193 (96%)	7 (4%)	0	100	100
2	Н	$219/228\ (96\%)$	215 (98%)	4 (2%)	0	100	100



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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
3	L	214/215 (100%)	207 (97%)	7 (3%)	0	100	100
All	All	633/647 (98%)	615 (97%)	18 (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	A	170/176 (97%)	166 (98%)	4 (2%)	49 39
2	Н	180/190 (95%)	176 (98%)	4 (2%)	52 43
3	L	187/189 (99%)	186 (100%)	1 (0%)	88 88
All	All	537/555 (97%)	528 (98%)	9 (2%)	73 54

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	Н	85	SER
3	L	100	GLN
1	A	448[B]	ASN
2	Н	17[A]	SER
2	Н	17[B]	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	498	GLN
2	Н	111	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)

9 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		Во	ond leng	ths	В	ond ang	les
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	В	1	1,4	14,14,15	0.39	0	17,19,21	0.67	0
4	NAG	В	2	4	14,14,15	0.39	0	17,19,21	0.55	0
4	BMA	В	3	4	11,11,12	0.66	0	15,15,17	0.74	0
4	MAN	В	4	4	11,11,12	0.66	0	15,15,17	1.08	2 (13%)
4	MAN	В	5	4	11,11,12	0.80	0	15,15,17	1.03	1 (6%)
4	FUC	В	6	4	10,10,11	0.63	0	14,14,16	0.80	0
5	NAG	С	1	3,5	14,14,15	0.30	0	17,19,21	0.50	0
5	NAG	С	2	5	14,14,15	0.31	0	17,19,21	0.50	0
5	FUC	С	3	5	10,10,11	0.86	0	14,14,16	0.85	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
4	NAG	В	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	В	2	4	-	0/6/23/26	0/1/1/1
4	BMA	В	3	4	-	0/2/19/22	0/1/1/1
4	MAN	В	4	4	-	0/2/19/22	0/1/1/1
4	MAN	В	5	4	-	0/2/19/22	0/1/1/1
4	FUC	В	6	4	-	-	0/1/1/1
5	NAG	С	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	С	2	5	-	0/6/23/26	0/1/1/1
5	FUC	С	3	5	-	-	0/1/1/1



There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
4	В	4	MAN	C1-O5-C5	3.09	116.38	112.19
4	В	5	MAN	C1-O5-C5	2.55	115.65	112.19
4	В	4	MAN	O2-C2-C3	-2.03	106.07	110.14

There are no chirality outliers.

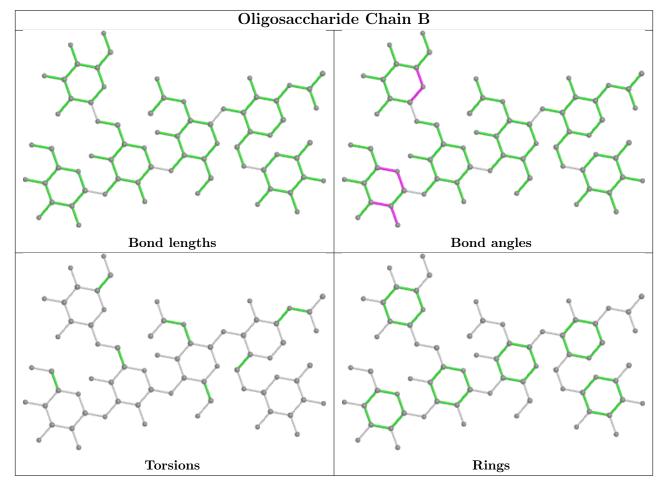
There are no torsion outliers.

There are no ring outliers.

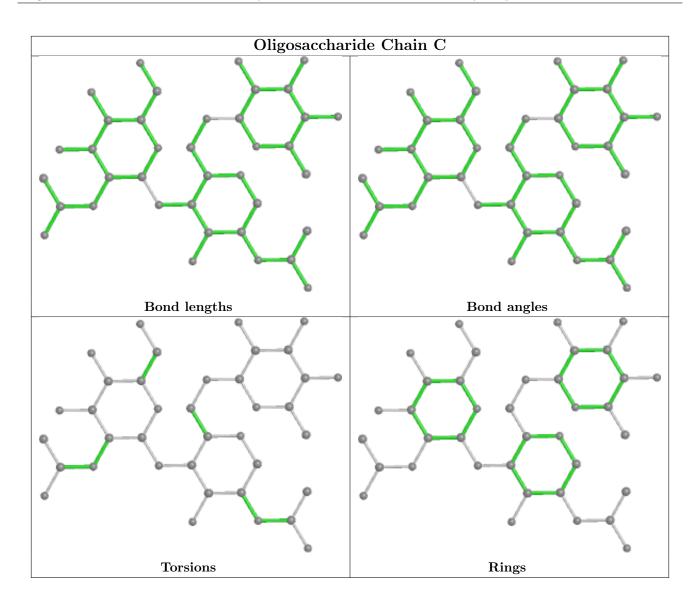
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	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)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

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>	# RSRZ > 2	$OWAB(A^2)$	Q<0.9
1	A	195/204~(95%)	-0.39	1 (0%) 91 91	21, 29, 45, 63	0
2	Н	218/228 (95%)	-0.26	4 (1%) 68 70	21, 31, 59, 86	0
3	L	214/215 (99%)	-0.29	1 (0%) 91 91	22, 33, 50, 76	0
All	All	627/647 (96%)	-0.31	6 (0%) 82 83	21, 31, 52, 86	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	Н	134	SER	3.0
2	Н	132	SER	2.8
3	L	214	CYS	2.6
2	Н	218	LYS	2.3
1	A	448[A]	ASN	2.1

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.

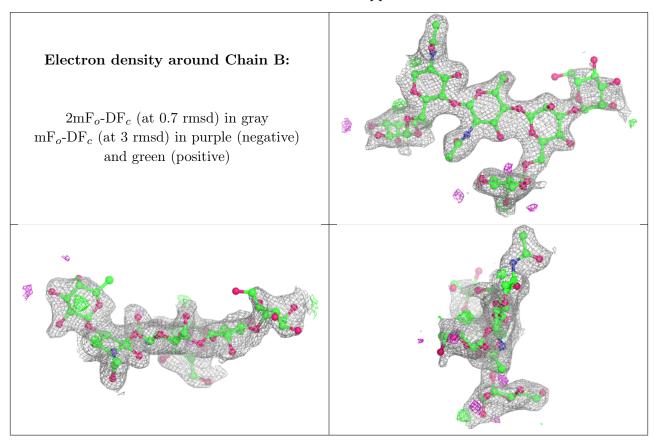
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	MAN	В	4	11/12	0.63	0.29	100,105,106,106	0
5	NAG	С	2	14/15	0.74	0.22	77,84,87,90	0
4	MAN	В	5	11/12	0.78	0.17	74,77,79,80	0



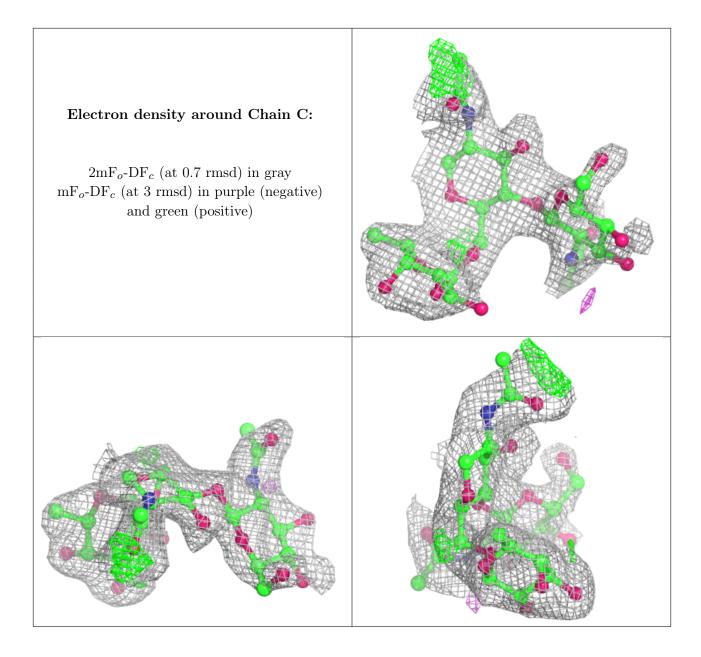
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	BMA	В	3	11/12	0.83	0.14	76,78,84,92	0
5	NAG	С	1	14/15	0.88	0.11	44,55,62,66	0
4	FUC	В	6	10/11	0.90	0.20	61,70,78,78	0
5	FUC	С	3	10/11	0.90	0.14	53,56,58,60	0
4	NAG	В	2	14/15	0.92	0.13	48,53,64,70	0
4	NAG	В	1	14/15	0.96	0.07	29,33,43,46	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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
6	MG	A	601	1/1	0.96	0.10	40,40,40,40	0
7	CL	L	302	1/1	0.96	0.08	52,52,52,52	0
7	CL	Н	302	1/1	0.97	0.06	41,41,41,41	0
6	MG	L	301	1/1	0.97	0.06	41,41,41,41	0
7	CL	Н	301	1/1	1.00	0.07	27,27,27,27	0



6.5 Other polymers (i)

There are no such residues in this entry.

