

wwPDB X-ray Structure Validation Summary Report (i)

Oct 10, 2023 – 02:12 AM EDT

PDB ID : 7RGF

Title: Protocadherin gammaC4 EC1-4 crystal structure disrupted trans interface

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Deposited on : 2021-07-15

Resolution : 2.40 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 : 2.35.1

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

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

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

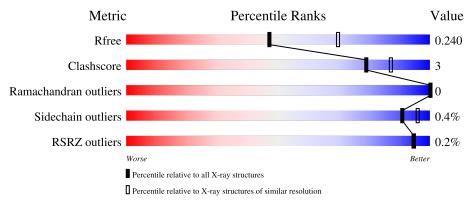
Validation Pipeline (wwPDB-VP) : 2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	Whole archive $(\# \mathrm{Entries})$	Similar resolution $(\# \text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	429	92%	6%				
1	В	429	89%	9%	-			
2	С	8	100%					
3	D	7	86%	14%				



2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 6736 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 Protocadherin gamma C4.

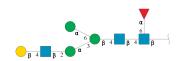
Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf	Trace
1	Λ	418	Total	С	N	О	S	0	0	0
1	Λ	410	3177	1977	559	634	7	0		
1	В	418	Total	С	N	О	S	0	0	0
1	Ъ	410	3150	1962	550	632	6	0		

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	422	HIS	-	expression tag	UNP Q91XX0
A	423	HIS	-	expression tag	UNP Q91XX0
A	424	HIS	-	expression tag	UNP Q91XX0
A	425	HIS	-	expression tag	UNP Q91XX0
A	426	HIS	-	expression tag	UNP Q91XX0
A	427	HIS	-	expression tag	UNP Q91XX0
A	428	HIS	-	expression tag	UNP Q91XX0
A	429	HIS	-	expression tag	UNP Q91XX0
В	422	HIS	_	expression tag	UNP Q91XX0
В	423	HIS	-	expression tag	UNP Q91XX0
В	424	HIS	_	expression tag	UNP Q91XX0
В	425	HIS	-	expression tag	UNP Q91XX0
В	426	HIS	-	expression tag	UNP Q91XX0
В	427	HIS	-	expression tag	UNP Q91XX0
В	428	HIS	-	expression tag	UNP Q91XX0
В	429	HIS	-	expression tag	UNP Q91XX0

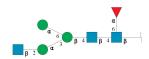
• Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-b eta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]be ta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.





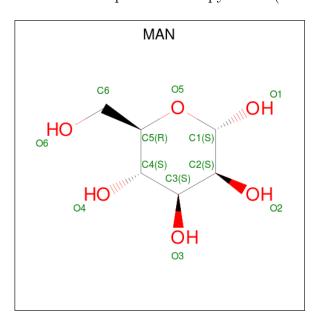
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace	
2	С	8	10001	C N 54 3	O 39	0	0	0

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



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
3	D	7	Total 85	C 48	N 3	O 34	0	0	0

• Molecule 4 is alpha-D-mannopyranose (three-letter code: MAN) (formula: $C_6H_{12}O_6$).

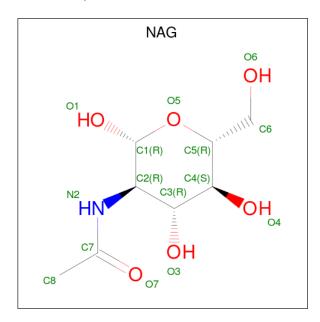


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 11 6 5	0	0
4	A	1	Total C O 11 6 5	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	В	1	Total C O 11 6 5	0	0
4	В	1	Total C O 11 6 5	0	0

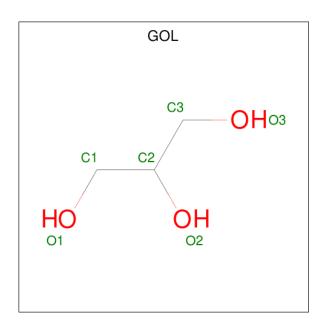
 \bullet Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $\rm C_8H_{15}NO_6).$



Mo	Chai	n Residues	A	Atoms				AltConf
5	A	1	Total 14			O 5	0	0
5	В	1	Total 14		N 1		0	0

• Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total 6	C 3	O 3	0	0

• Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	9	Total Ca 9 9	0	0
7	В	9	Total Ca 9 9	0	0

• Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	3	Total Cl 3 3	0	0
8	В	4	Total Cl 4 4	0	0

• Molecule 9 is water.

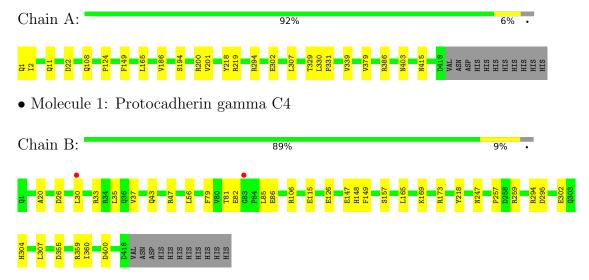
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	50	Total O 50 50	0	0
9	В	75	Total O 75 75	0	0



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: Protocadherin gamma C4



 $\bullet \ \, \text{Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-6)]2-acetamido-2-deoxy-$

Chain C: 100%

NAG1 NAG2 BMA3 MAN4 NAG5 GAL6 MAN7

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

Chain D: 86% 14%





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	51.94Å 103.95Å 200.59Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.16 - 2.40	Depositor
rtesolution (A)	38.16 - 2.40	EDS
% Data completeness	98.5 (38.16-2.40)	Depositor
(in resolution range)	94.3 (38.16-2.40)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.22 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
D D.	0.196 , 0.240	Depositor
R, R_{free}	0.196 , 0.240	DCC
R_{free} test set	2144 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	48.4	Xtriage
Anisotropy	0.683	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.28 , 34.8	EDS
L-test for twinning ²	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6736	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 17.66% 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: BMA, CL, MAN, CA, FUC, GAL, NAG, GOL

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
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.26	0/3241	0.47	0/4418
1	В	0.26	0/3214	0.48	0/4386
All	All	0.26	0/6455	0.47	0/8804

There are no bond length outliers.

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	A	3177	0	3072	14	0
1	В	3150	0	3021	25	0
2	С	96	0	82	0	0
3	D	85	0	73	0	0
4	A	22	0	20	0	0
4	В	22	0	20	0	0
5	A	14	0	13	0	0
5	В	14	0	13	0	0
6	A	6	0	8	0	0
7	A	9	0	0	0	0
7	В	9	0	0	0	0



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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	3	0	0	0	0
8	В	4	0	0	0	0
9	A	50	0	0	1	0
9	В	75	0	0	2	0
All	All	6736	0	6322	36	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 3.

The worst 5 of 36 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:B:294:ARG:HG2	1:B:302:GLU:HG2	1.74	0.69
1:B:81:THR:O	1:B:85:LEU:HD12	1.95	0.67
1:B:359:ARG:NH2	9:B:603:HOH:O	2.29	0.65
1:B:47:ARG:HH21	1:B:56:LEU:HD12	1.61	0.64
1:B:218:TYR:HB2	1:B:307:LEU:HD23	1.78	0.64

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	Perce	entiles
1	A	$416/429 \ (97\%)$	412 (99%)	4 (1%)	0	100	100
1	В	416/429 (97%)	411 (99%)	5 (1%)	0	100	100
All	All	832/858 (97%)	823 (99%)	9 (1%)	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	352/371 (95%)	349 (99%)	3 (1%)	78 90
1	В	346/371 (93%)	346 (100%)	0	100 100
All	All	698/742 (94%)	695 (100%)	3 (0%)	91 96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	201	VAL
1	A	386	ARG
1	A	403	ASN

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)

15 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	Tuna	Chain	Res	Link	Вс	Bond lengths			Bond angles			
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2		
2	NAG	С	1	2,1	14,14,15	0.82	0	17,19,21	0.63	0		
2	NAG	С	2	2	14,14,15	0.66	0	17,19,21	0.50	0		
2	BMA	С	3	2	11,11,12	0.87	0	15,15,17	1.11	0		
2	MAN	С	4	2	11,11,12	0.64	0	15,15,17	0.71	0		
2	NAG	С	5	2	14,14,15	0.73	0	17,19,21	0.64	0		
2	GAL	С	6	2	11,11,12	0.50	0	15,15,17	0.88	0		
2	MAN	С	7	2	11,11,12	0.45	0	15,15,17	0.62	0		
2	FUC	С	8	2	10,10,11	0.65	0	14,14,16	0.71	0		
3	NAG	D	1	1,3	14,14,15	0.70	0	17,19,21	0.46	0		
3	NAG	D	2	3	14,14,15	0.71	0	17,19,21	0.62	0		
3	BMA	D	3	3	11,11,12	1.04	1 (9%)	15,15,17	0.99	0		
3	MAN	D	4	3	11,11,12	0.50	0	15,15,17	0.77	0		
3	NAG	D	5	3	14,14,15	0.77	0	17,19,21	0.47	0		
3	MAN	D	6	3	11,11,12	0.51	0	15,15,17	0.63	0		
3	FUC	D	7	3	10,10,11	0.64	0	14,14,16	0.75	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	С	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	С	2	2	-	0/6/23/26	0/1/1/1
2	BMA	С	3	2	-	0/2/19/22	0/1/1/1
2	MAN	С	4	2	-	2/2/19/22	0/1/1/1
2	NAG	С	5	2	-	2/6/23/26	0/1/1/1
2	GAL	С	6	2	-	0/2/19/22	0/1/1/1
2	MAN	С	7	2	-	1/2/19/22	0/1/1/1
2	FUC	С	8	2	-	-	0/1/1/1
3	NAG	D	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	2/6/23/26	0/1/1/1
3	BMA	D	3	3	-	0/2/19/22	0/1/1/1
3	MAN	D	4	3	-	0/2/19/22	0/1/1/1
3	NAG	D	5	3	-	2/6/23/26	0/1/1/1
3	MAN	D	6	3	-	0/2/19/22	0/1/1/1
3	FUC	D	7	3	-	-	0/1/1/1

All (1) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(ext{\AA})$
3	D	3	BMA	O5-C1	-2.10	1.40	1.43

There are no bond angle outliers.

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

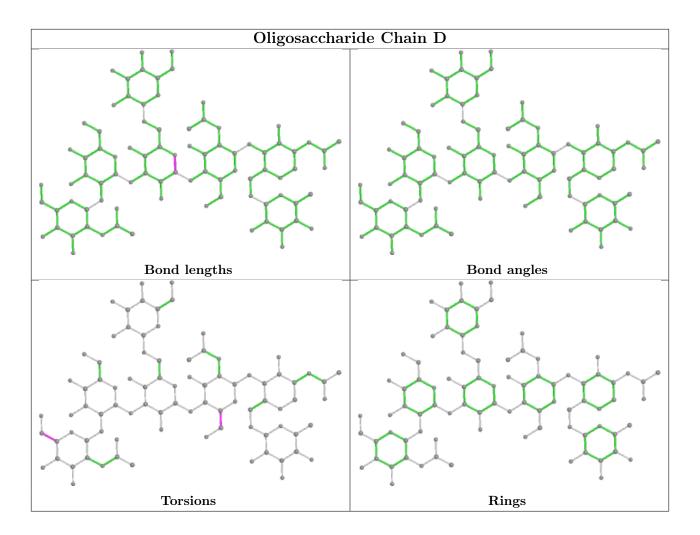
Mol	Chain	Res	Type	Atoms
2	С	4	MAN	O5-C5-C6-O6
2	С	4	MAN	C4-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
3	D	5	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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 32 ligands modelled in this entry, 25 are monoatomic - leaving 7 for Mogul analysis.

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	Tuno	Type Chain Res		Link	Во	ond leng	$ ag{ths}$	В	ond ang	les
WIOI	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GOL	A	504	-	5,5,5	0.91	0	5,5,5	1.06	0
4	MAN	В	501	1	11,11,12	0.47	0	15,15,17	0.73	0
4	MAN	A	501	1	11,11,12	0.51	0	15,15,17	0.70	0
4	MAN	В	502	1	11,11,12	0.52	0	15,15,17	0.64	0
5	NAG	В	503	1	14,14,15	0.67	0	17,19,21	0.66	1 (5%)
5	NAG	A	503	1	14,14,15	0.76	1 (7%)	17,19,21	0.55	0



Mol Type Chain	Res	Res Link	Bo	Bond lengths			Bond angles			
	Chain	rtes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2	
4	MAN	A	502	1	11,11,12	0.44	0	15,15,17	0.69	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
6	GOL	A	504	-	-	0/4/4/4	-
4	MAN	В	501	1	-	2/2/19/22	0/1/1/1
4	MAN	A	501	1	-	2/2/19/22	0/1/1/1
4	MAN	В	502	1	-	0/2/19/22	0/1/1/1
5	NAG	В	503	1	-	1/6/23/26	0/1/1/1
5	NAG	A	503	1	-	3/6/23/26	0/1/1/1
4	MAN	A	502	1	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
5	A	503	NAG	C1-C2	2.03	1.55	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
5	В	503	NAG	C2-N2-C7	2.25	126.10	122.90

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	503	NAG	O5-C5-C6-O6
4	A	501	MAN	O5-C5-C6-O6
4	A	501	MAN	C4-C5-C6-O6
4	В	501	MAN	C4-C5-C6-O6
5	A	503	NAG	C1-C2-N2-C7

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	$\langle { m RSRZ} \rangle$	# RSRZ > 2	$OWAB(Å^2)$	Q < 0.9
1	A	418/429 (97%)	-0.28	0 100 100	39, 72, 107, 118	0
1	В	418/429 (97%)	-0.26	2 (0%) 91 89	38, 65, 115, 153	0
All	All	836/858 (97%)	-0.27	2 (0%) 95 94	38, 68, 110, 153	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	30	LEU	4.3
1	В	83	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.

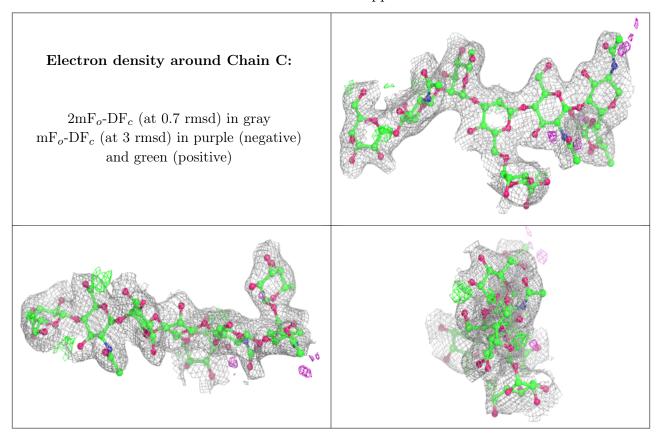
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	GAL	С	6	11/12	0.79	0.16	96,102,104,104	0
3	MAN	D	6	11/12	0.83	0.20	113,117,117,120	0
2	MAN	С	7	11/12	0.88	0.24	124,128,131,131	0
3	FUC	D	7	10/11	0.89	0.15	96,100,103,107	0
3	MAN	D	4	11/12	0.91	0.16	74,83,88,89	0
2	BMA	С	3	11/12	0.92	0.15	79,91,119,132	0
2	NAG	С	2	14/15	0.92	0.15	77,81,91,91	0



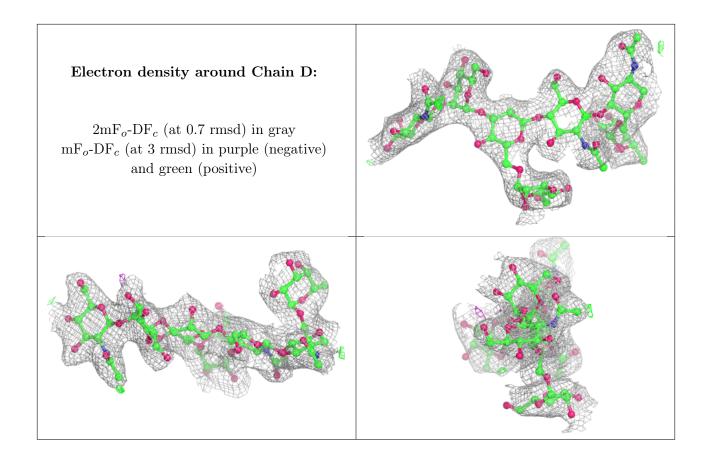
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	FUC	С	8	10/11	0.93	0.25	89,93,97,100	0
3	NAG	D	5	14/15	0.93	0.14	75,77,81,90	0
3	NAG	D	2	14/15	0.94	0.10	91,95,98,100	0
3	BMA	D	3	11/12	0.94	0.12	96,103,109,112	0
2	NAG	С	5	14/15	0.95	0.11	61,76,78,88	0
2	NAG	С	1	14/15	0.96	0.13	47,58,77,86	0
3	NAG	D	1	14/15	0.96	0.08	50,62,91,107	0
2	MAN	С	4	11/12	0.97	0.11	62,66,71,72	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	$\operatorname{B-factors}({ ext{\AA}}^2)$	Q < 0.9
4	MAN	A	501	11/12	0.80	0.16	86,93,100,103	0
4	MAN	В	501	11/12	0.88	0.17	76,93,96,97	0
5	NAG	A	503	14/15	0.90	0.15	87,102,110,119	0
5	NAG	В	503	14/15	0.90	0.20	83,105,114,118	0
4	MAN	В	502	11/12	0.91	0.18	74,92,98,101	0
6	GOL	A	504	6/6	0.91	0.21	75,85,90,94	0
4	MAN	A	502	11/12	0.92	0.16	66,81,88,91	0
7	CA	A	513	1/1	0.92	0.13	64,64,64,64	0
7	CA	A	505	1/1	0.93	0.11	76,76,76,76	0
7	CA	В	504	1/1	0.94	0.07	62,62,62,62	0
7	CA	A	512	1/1	0.95	0.09	60,60,60,60	0
8	CL	A	515	1/1	0.96	0.23	74,74,74,74	0
8	CL	В	514	1/1	0.96	0.17	64,64,64,64	0
7	CA	A	511	1/1	0.97	0.12	66,66,66,66	0



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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
8	CL	A	516	1/1	0.97	0.07	74,74,74,74	0
8	CL	В	513	1/1	0.97	0.17	48,48,48,48	0
7	CA	В	509	1/1	0.97	0.12	52,52,52,52	0
8	CL	В	515	1/1	0.97	0.07	69,69,69,69	0
8	CL	В	516	1/1	0.97	0.10	74,74,74,74	0
7	CA	В	512	1/1	0.98	0.14	38,38,38,38	0
7	CA	A	507	1/1	0.98	0.14	52,52,52,52	0
7	CA	A	508	1/1	0.98	0.07	77,77,77,77	0
7	CA	A	510	1/1	0.98	0.15	51,51,51,51	0
7	CA	В	505	1/1	0.98	0.09	63,63,63,63	0
7	CA	В	507	1/1	0.98	0.08	73,73,73,73	0
7	CA	A	506	1/1	0.98	0.10	60,60,60,60	0
7	CA	В	508	1/1	0.99	0.11	54,54,54,54	0
7	CA	В	506	1/1	0.99	0.12	55,55,55,55	0
7	CA	В	510	1/1	0.99	0.13	54,54,54,54	0
7	CA	В	511	1/1	0.99	0.14	42,42,42,42	0
7	CA	A	509	1/1	0.99	0.15	69,69,69,69	0
8	CL	A	514	1/1	0.99	0.06	56,56,56,56	0

6.5 Other polymers (i)

There are no such residues in this entry.

