



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 2, 2024 – 08:10 am GMT

PDB ID : 5F8R  
Title : Blood group antigen binding adhesin BabA of Helicobacter pylori strain S831 in complex with blood group H Lewis b hexasaccharide  
Authors : Moonens, K.; Gideonsson, P.; Subedi, S.; Romao, E.; Oscarson, S.; Muyldermans, S.; Boren, T.; Remaut, H.  
Deposited on : 2015-12-09  
Resolution : 2.44 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](mailto: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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
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

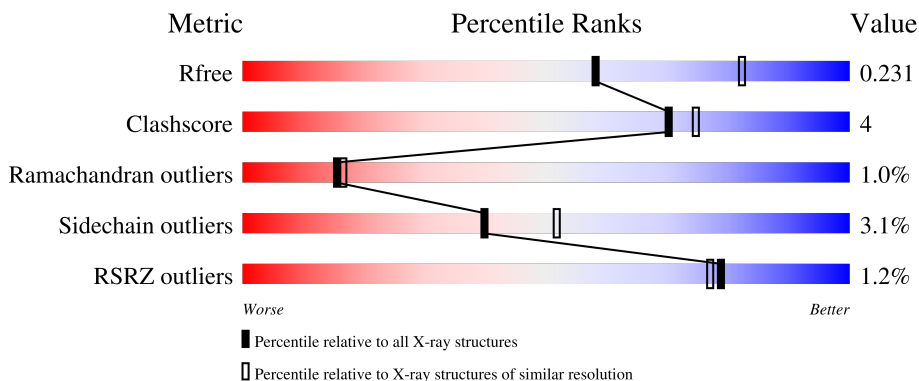
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.44 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1564 (2.46-2.42)
Clashscore	141614	1631 (2.46-2.42)
Ramachandran outliers	138981	1617 (2.46-2.42)
Sidechain outliers	138945	1617 (2.46-2.42)
RSRZ outliers	127900	1547 (2.46-2.42)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	466	 2% 80% 8% • 11%
1	B	466	 2% 80% 8% • 10%
2	C	120	 2% 85% 8% • 6%
2	D	120	 % 92% •• 5%
3	E	6	 100%

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Mol	Chain	Length	Quality of chain
3	F	6	 17% 83%

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 8224 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 Adhesin binding fucosylated histo-blood group antigen, Adhesin, Adhesin binding fucosylated histo-blood group antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	417	3111	1920	536	642	13	0	1	0
1	B	421	3139	1939	539	648	13	0	0	0

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	ALA	-	expression tag	UNP O52269
A	4	SER	-	expression tag	UNP O52269
A	5	TRP	-	expression tag	UNP O52269
A	6	SER	-	expression tag	UNP O52269
A	7	HIS	-	expression tag	UNP O52269
A	8	PRO	-	expression tag	UNP O52269
A	9	GLN	-	expression tag	UNP O52269
A	10	PHE	-	expression tag	UNP O52269
A	11	GLU	-	expression tag	UNP O52269
A	12	LYS	-	expression tag	UNP O52269
A	13	SER	-	expression tag	UNP O52269
A	14	GLY	-	expression tag	UNP O52269
A	15	GLY	-	expression tag	UNP O52269
A	16	GLY	-	expression tag	UNP O52269
A	17	GLY	-	expression tag	UNP O52269
A	18	GLY	-	expression tag	UNP O52269
A	19	LEU	-	expression tag	UNP O52269
A	20	VAL	-	expression tag	UNP O52269
A	21	PRO	-	expression tag	UNP O52269
A	22	ARG	-	expression tag	UNP O52269
A	23	GLY	-	expression tag	UNP O52269
A	24	SER	-	expression tag	UNP O52269
A	461	GLY	-	expression tag	UNP O52269
A	462	SER	-	expression tag	UNP O52269

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Chain	Residue	Modelled	Actual	Comment	Reference
A	463	HIS	-	expression tag	UNP O52269
A	464	HIS	-	expression tag	UNP O52269
A	465	HIS	-	expression tag	UNP O52269
A	466	HIS	-	expression tag	UNP O52269
A	467	HIS	-	expression tag	UNP O52269
A	468	HIS	-	expression tag	UNP O52269
B	3	ALA	-	expression tag	UNP O52269
B	4	SER	-	expression tag	UNP O52269
B	5	TRP	-	expression tag	UNP O52269
B	6	SER	-	expression tag	UNP O52269
B	7	HIS	-	expression tag	UNP O52269
B	8	PRO	-	expression tag	UNP O52269
B	9	GLN	-	expression tag	UNP O52269
B	10	PHE	-	expression tag	UNP O52269
B	11	GLU	-	expression tag	UNP O52269
B	12	LYS	-	expression tag	UNP O52269
B	13	SER	-	expression tag	UNP O52269
B	14	GLY	-	expression tag	UNP O52269
B	15	GLY	-	expression tag	UNP O52269
B	16	GLY	-	expression tag	UNP O52269
B	17	GLY	-	expression tag	UNP O52269
B	18	GLY	-	expression tag	UNP O52269
B	19	LEU	-	expression tag	UNP O52269
B	20	VAL	-	expression tag	UNP O52269
B	21	PRO	-	expression tag	UNP O52269
B	22	ARG	-	expression tag	UNP O52269
B	23	GLY	-	expression tag	UNP O52269
B	24	SER	-	expression tag	UNP O52269
B	461	GLY	-	expression tag	UNP O52269
B	462	SER	-	expression tag	UNP O52269
B	463	HIS	-	expression tag	UNP O52269
B	464	HIS	-	expression tag	UNP O52269
B	465	HIS	-	expression tag	UNP O52269
B	466	HIS	-	expression tag	UNP O52269
B	467	HIS	-	expression tag	UNP O52269
B	468	HIS	-	expression tag	UNP O52269

- Molecule 2 is a protein called Nanobody Nb-ER19.

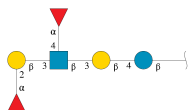
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	113	863	540	155	163	5	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	D	114	873	546	158	164	5	0	0	0

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	E	6	68	38	1	29	0	0	0
3	F	6	68	38	1	29	0	0	0

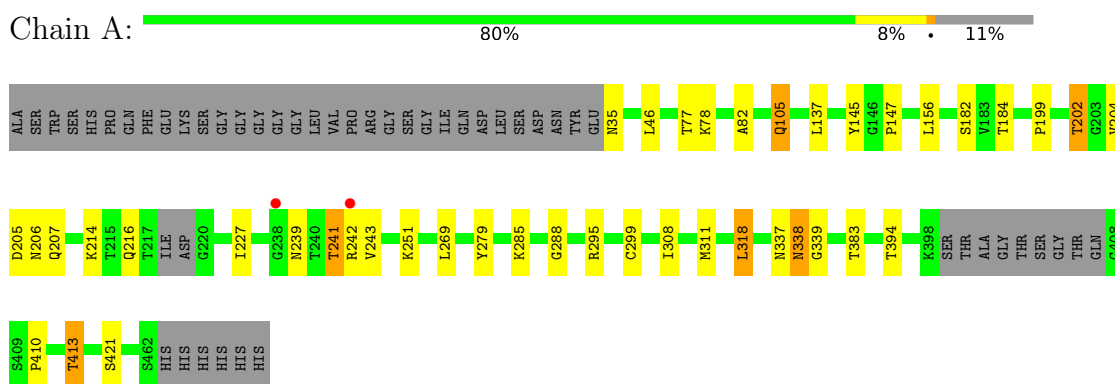
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	38	Total	O	0	0
			38	38		
4	B	28	Total	O	0	0
			28	28		
4	C	12	Total	O	0	0
			12	12		
4	D	24	Total	O	0	0
			24	24		

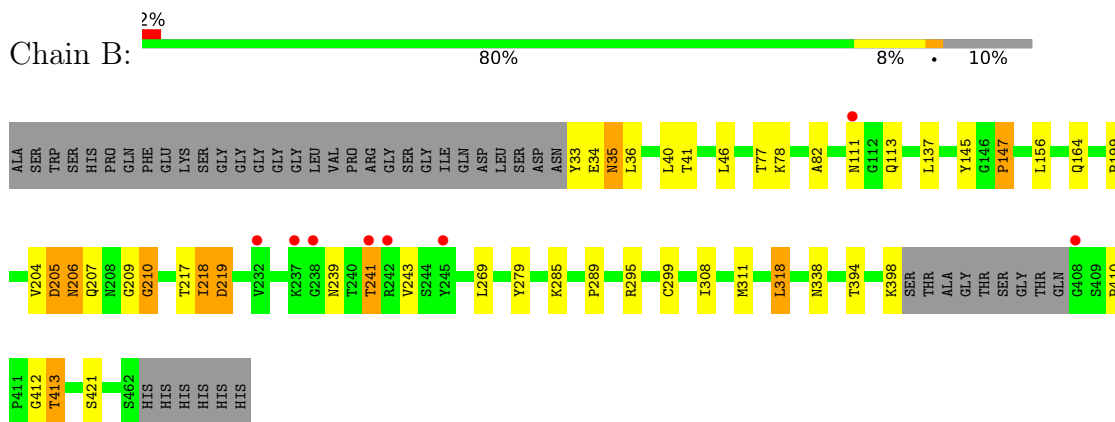
### 3 Residue-property plots [i](#)

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.

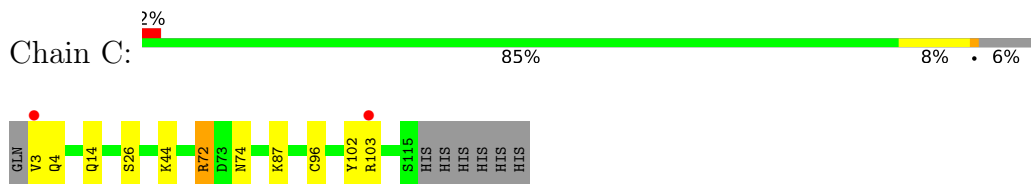
- Molecule 1: Adhesin binding fucosylated histo-blood group antigen,Adhesin,Adhesin binding fucosylated histo-blood group antigen



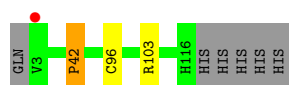
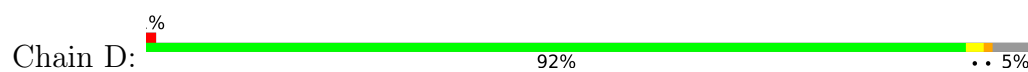
- Molecule 1: Adhesin binding fucosylated histo-blood group antigen,Adhesin,Adhesin binding fucosylated histo-blood group antigen



- Molecule 2: Nanobody Nb-ER19



- Molecule 2: Nanobody Nb-ER19



- Molecule 3: alpha-L-fucopyranose-(1-2)-beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose



- Molecule 3: alpha-L-fucopyranose-(1-2)-beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	50.79Å 131.91Å 125.39Å 90.00° 94.88° 90.00°	Depositor
Resolution (Å)	48.40 – 2.44 48.36 – 2.44	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.40-2.44) 99.5 (48.36-2.44)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.34 (at 2.45Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.189 , 0.219 0.203 , 0.231	Depositor DCC
$R_{free}$ test set	3049 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	61.0	Xtrriage
Anisotropy	0.285	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 37.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8224	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<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.

## 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: BGC, FUC, NAG, GAL

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	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.81	1/3157 (0.0%)	0.83	3/4292 (0.1%)
1	B	0.76	0/3187	0.81	2/4335 (0.0%)
2	C	0.81	0/882	0.89	2/1195 (0.2%)
2	D	0.94	0/893	0.96	1/1210 (0.1%)
All	All	0.81	1/8119 (0.0%)	0.85	8/11032 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	105	GLN	CG-CD	5.12	1.62	1.51

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	299	CYS	CA-CB-SG	11.40	134.52	114.00
1	A	299	CYS	CB-CA-C	-8.42	93.56	110.40
2	C	72	ARG	NE-CZ-NH1	7.71	124.16	120.30
2	D	42	PRO	C-N-CA	-6.47	108.71	122.30
2	C	72	ARG	NE-CZ-NH2	-6.36	117.12	120.30
1	B	318	LEU	CA-CB-CG	5.99	129.08	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	318	LEU	CA-CB-CG	5.88	128.83	115.30
1	B	299	CYS	CA-CB-SG	5.08	123.15	114.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	242	ARG	Peptide
1	B	33	TYR	Peptide

## 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	3111	0	3038	30	0
1	B	3139	0	3064	22	0
2	C	863	0	845	6	0
2	D	873	0	852	0	0
3	E	68	0	60	0	0
3	F	68	0	60	0	0
4	A	38	0	0	0	0
4	B	28	0	0	0	0
4	C	12	0	0	0	0
4	D	24	0	0	0	0
All	All	8224	0	7919	58	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 (58) 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:279:TYR:OH	1:A:295:ARG:CZ	1.86	1.22
2:C:3:VAL:N	2:C:26:SER:HG	1.51	1.08
1:A:279:TYR:HH	1:A:295:ARG:NE	1.68	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:TYR:HH	1:A:295:ARG:CZ	1.85	0.88
1:A:279:TYR:OH	1:A:295:ARG:NH1	2.09	0.84
1:A:279:TYR:OH	1:A:295:ARG:NE	2.13	0.80
2:C:3:VAL:HG13	2:C:102:TYR:OH	1.86	0.75
1:A:337:ASN:O	1:A:339:GLY:N	2.21	0.74
1:B:410:PRO:O	1:B:413:THR:HG23	1.94	0.67
2:C:3:VAL:HG13	2:C:102:TYR:CZ	2.31	0.65
2:C:72:ARG:HD3	2:C:74:ASN:OD1	2.03	0.59
1:B:241:THR:HG22	1:B:243:VAL:H	1.69	0.57
1:B:205:ASP:O	1:B:206:ASN:HB2	2.05	0.56
1:A:410:PRO:O	1:A:413:THR:CG2	2.55	0.55
1:B:308:ILE:HA	1:B:311:MET:HE3	1.88	0.55
1:B:209:GLY:O	1:B:210:GLY:O	2.25	0.55
1:A:105:GLN:CD	1:A:279:TYR:HD2	2.11	0.55
1:B:218:ILE:HG22	1:B:219:ASP:N	2.22	0.54
1:A:239:ASN:OD1	1:A:241:THR:HG22	2.07	0.53
1:A:199:PRO:HD3	1:A:207:GLN:HG3	1.90	0.53
1:B:199:PRO:HD3	1:B:207:GLN:HG3	1.91	0.52
1:A:410:PRO:O	1:A:413:THR:HG23	2.09	0.52
1:A:308:ILE:HA	1:A:311:MET:HE3	1.91	0.52
1:A:394:THR:HG22	1:A:421:SER:HB2	1.94	0.50
1:A:35:ASN:OD1	1:A:35:ASN:C	2.49	0.50
2:C:3:VAL:N	2:C:26:SER:OG	2.32	0.50
1:B:218:ILE:CG2	1:B:219:ASP:N	2.76	0.49
1:B:35:ASN:OD1	1:B:35:ASN:C	2.51	0.48
1:A:202:THR:HG22	1:A:204:VAL:HB	1.95	0.48
1:A:279:TYR:OH	1:A:295:ARG:CG	2.62	0.48
1:B:156:LEU:C	1:B:156:LEU:HD23	2.34	0.48
1:A:202:THR:HG22	1:A:204:VAL:H	1.79	0.48
1:B:137:LEU:HD23	1:B:147:PRO:HG3	1.97	0.47
1:A:214:LYS:HE3	1:A:227:ILE:HD11	1.97	0.47
1:A:156:LEU:C	1:A:156:LEU:HD23	2.35	0.47
1:B:217:THR:HG23	1:B:217:THR:O	2.15	0.47
1:A:105:GLN:CD	1:A:279:TYR:CD2	2.89	0.46
1:A:156:LEU:HG	1:A:269:LEU:HD12	1.97	0.46
1:A:137:LEU:HD23	1:A:147:PRO:HG3	1.97	0.45
1:B:239:ASN:OD1	1:B:241:THR:HB	2.16	0.45
1:B:111:ASN:HB2	1:B:113:GLN:HG2	1.98	0.45
1:B:394:THR:HG22	1:B:421:SER:HB2	1.99	0.44
1:A:337:ASN:O	1:A:338:ASN:C	2.55	0.44
1:B:156:LEU:HG	1:B:269:LEU:HD12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:164:GLN:OE1	1:B:412:GLY:HA2	2.18	0.43
1:A:202:THR:CG2	1:A:204:VAL:HB	2.49	0.43
1:B:77:THR:O	1:B:78:LYS:C	2.57	0.42
2:C:3:VAL:CG1	2:C:102:TYR:OH	2.61	0.42
1:B:36:LEU:O	1:B:40:LEU:HG	2.20	0.42
1:A:205:ASP:N	1:A:205:ASP:OD1	2.54	0.41
1:B:279:TYR:OH	1:B:295:ARG:HD2	2.20	0.41
1:A:77:THR:O	1:A:78:LYS:C	2.58	0.41
1:A:241:THR:HG23	1:A:243:VAL:H	1.86	0.41
1:B:204:VAL:HG12	1:B:206:ASN:H	1.84	0.41
1:A:82:ALA:HB1	1:A:318:LEU:HD21	2.03	0.41
1:A:184:THR:OG1	1:A:251:LYS:HD2	2.21	0.40
1:A:206:ASN:OD1	1:A:206:ASN:N	2.55	0.40
1:B:82:ALA:HB1	1:B:318:LEU:HD21	2.03	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	A	412/466 (88%)	396 (96%)	14 (3%)	2 (0%)	29	34
1	B	417/466 (90%)	397 (95%)	13 (3%)	7 (2%)	9	7
2	C	111/120 (92%)	108 (97%)	2 (2%)	1 (1%)	17	20
2	D	112/120 (93%)	109 (97%)	3 (3%)	0	100	100
All	All	1052/1172 (90%)	1010 (96%)	32 (3%)	10 (1%)	15	16

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	338	ASN

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Mol	Chain	Res	Type
1	B	34	GLU
1	B	206	ASN
1	B	210	GLY
1	B	338	ASN
1	B	289	PRO
1	B	241	THR
2	C	4	GLN
1	B	218	ILE
1	A	288	GLY

### 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	345/382 (90%)	336 (97%)	9 (3%)	46	58
1	B	348/382 (91%)	338 (97%)	10 (3%)	42	54
2	C	92/99 (93%)	87 (95%)	5 (5%)	22	29
2	D	93/99 (94%)	90 (97%)	3 (3%)	39	50
All	All	878/962 (91%)	851 (97%)	27 (3%)	40	52

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU
1	A	145	TYR
1	A	182	SER
1	A	202	THR
1	A	216	GLN
1	A	241	THR
1	A	285	LYS
1	A	383	THR
1	A	413	THR
1	B	35	ASN
1	B	41	THR
1	B	46	LEU

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Mol	Chain	Res	Type
1	B	145	TYR
1	B	147	PRO
1	B	205	ASP
1	B	219	ASP
1	B	285	LYS
1	B	398	LYS
1	B	413	THR
2	C	14	GLN
2	C	44	LYS
2	C	87	LYS
2	C	96	CYS
2	C	103	ARG
2	D	42	PRO
2	D	96	CYS
2	D	103	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	A	105	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](#)

12 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	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	BGC	E	1	3	12,12,12	0.59	0	17,17,17	1.18	1 (5%)
3	GAL	E	2	3	11,11,12	0.58	0	15,15,17	1.67	5 (33%)
3	NAG	E	3	3	14,14,15	1.21	2 (14%)	17,19,21	1.72	4 (23%)
3	GAL	E	4	3	11,11,12	0.63	0	15,15,17	1.09	1 (6%)
3	FUC	E	5	3	10,10,11	1.05	0	14,14,16	1.32	1 (7%)
3	FUC	E	6	3	10,10,11	0.90	1 (10%)	14,14,16	1.10	0
3	BGC	F	1	3	12,12,12	0.63	0	17,17,17	1.71	3 (17%)
3	GAL	F	2	3	11,11,12	0.63	0	15,15,17	2.06	4 (26%)
3	NAG	F	3	3	14,14,15	0.84	1 (7%)	17,19,21	1.33	3 (17%)
3	GAL	F	4	3	11,11,12	0.90	0	15,15,17	1.06	0
3	FUC	F	5	3	10,10,11	0.71	0	14,14,16	1.44	1 (7%)
3	FUC	F	6	3	10,10,11	0.59	0	14,14,16	1.63	4 (28%)

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
3	BGC	E	1	3	-	1/2/22/22	0/1/1/1
3	GAL	E	2	3	-	0/2/19/22	0/1/1/1
3	NAG	E	3	3	-	0/6/23/26	0/1/1/1
3	GAL	E	4	3	-	0/2/19/22	0/1/1/1
3	FUC	E	5	3	-	-	0/1/1/1
3	FUC	E	6	3	-	-	0/1/1/1
3	BGC	F	1	3	-	2/2/22/22	0/1/1/1
3	GAL	F	2	3	-	2/2/19/22	0/1/1/1
3	NAG	F	3	3	-	0/6/23/26	0/1/1/1
3	GAL	F	4	3	-	0/2/19/22	0/1/1/1
3	FUC	F	5	3	-	-	0/1/1/1
3	FUC	F	6	3	-	-	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	3	NAG	O5-C1	-2.96	1.39	1.43
3	E	3	NAG	O5-C5	-2.58	1.38	1.43
3	E	6	FUC	O5-C1	-2.43	1.39	1.43
3	F	3	NAG	O5-C1	-2.19	1.40	1.43



All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	2	GAL	O5-C5-C6	5.47	115.79	107.20
3	E	3	NAG	O5-C5-C6	-4.02	100.90	107.20
3	F	1	BGC	C3-C4-C5	3.94	117.27	110.24
3	F	5	FUC	O5-C1-C2	-3.51	105.35	110.77
3	F	1	BGC	O5-C5-C6	3.17	114.31	106.44
3	F	2	GAL	C3-C4-C5	-2.99	104.90	110.24
3	E	3	NAG	C4-C3-C2	-2.95	106.69	111.02
3	E	5	FUC	C1-C2-C3	2.92	113.25	109.67
3	E	2	GAL	C1-O5-C5	2.82	116.01	112.19
3	F	6	FUC	C1-C2-C3	2.81	113.12	109.67
3	F	2	GAL	O2-C2-C3	-2.76	104.62	110.14
3	F	6	FUC	O5-C1-C2	2.66	114.88	110.77
3	F	3	NAG	C2-N2-C7	2.54	126.52	122.90
3	E	2	GAL	O5-C1-C2	2.54	114.70	110.77
3	F	6	FUC	C6-C5-C4	-2.50	108.45	113.07
3	F	1	BGC	C1-C2-C3	-2.50	105.14	110.31
3	F	6	FUC	O5-C5-C6	2.38	112.45	107.33
3	E	2	GAL	C1-C2-C3	2.35	112.56	109.67
3	E	3	NAG	O7-C7-C8	-2.34	117.70	122.06
3	E	2	GAL	O3-C3-C2	-2.27	105.64	109.99
3	E	4	GAL	C1-O5-C5	2.25	115.25	112.19
3	F	3	NAG	O7-C7-N2	2.24	126.06	121.95
3	E	1	BGC	O5-C5-C6	2.23	111.98	106.44
3	F	3	NAG	C1-O5-C5	2.17	115.13	112.19
3	F	2	GAL	O3-C3-C4	2.11	115.22	110.35
3	E	2	GAL	O2-C2-C3	-2.06	106.01	110.14
3	E	3	NAG	O3-C3-C4	-2.02	105.69	110.35

There are no chirality outliers.

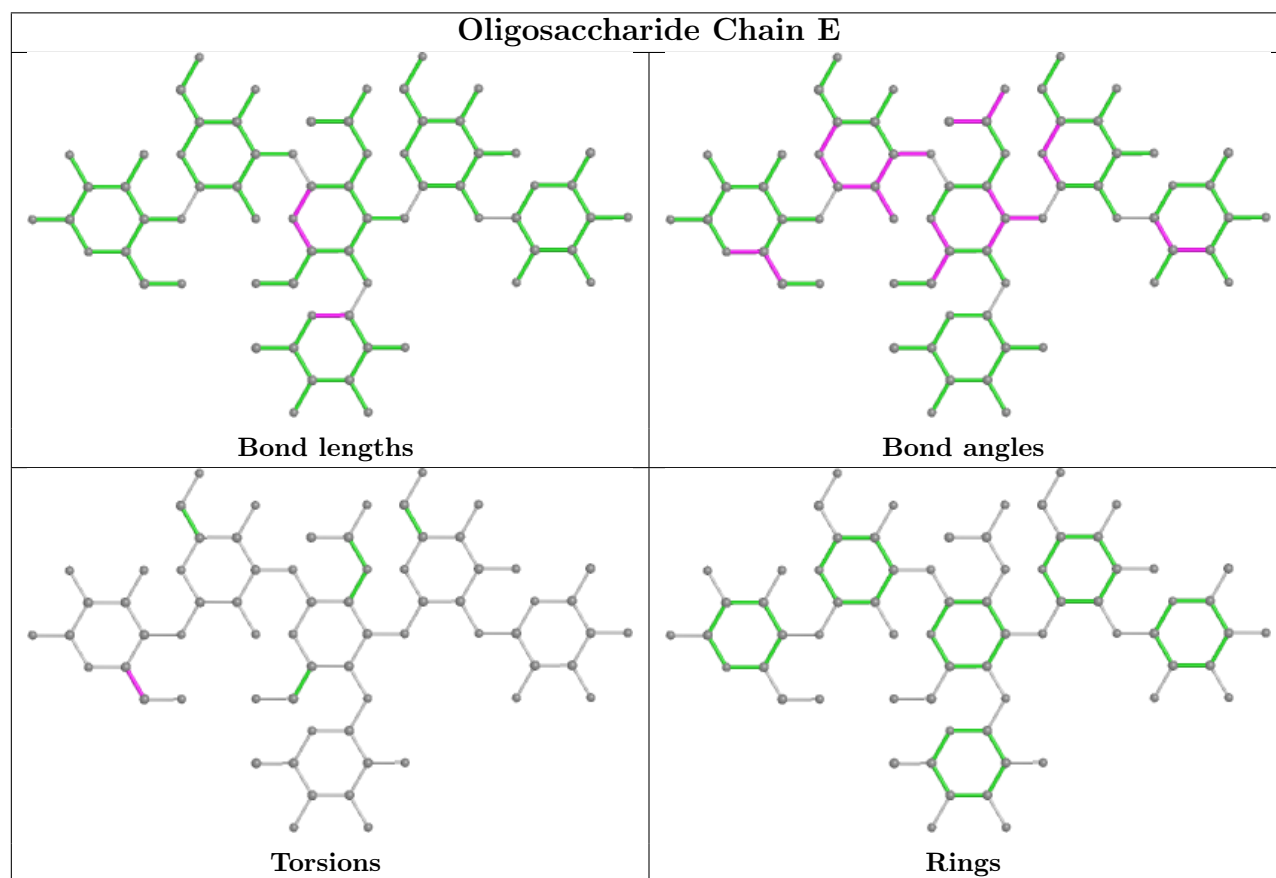
All (5) torsion outliers are listed below:

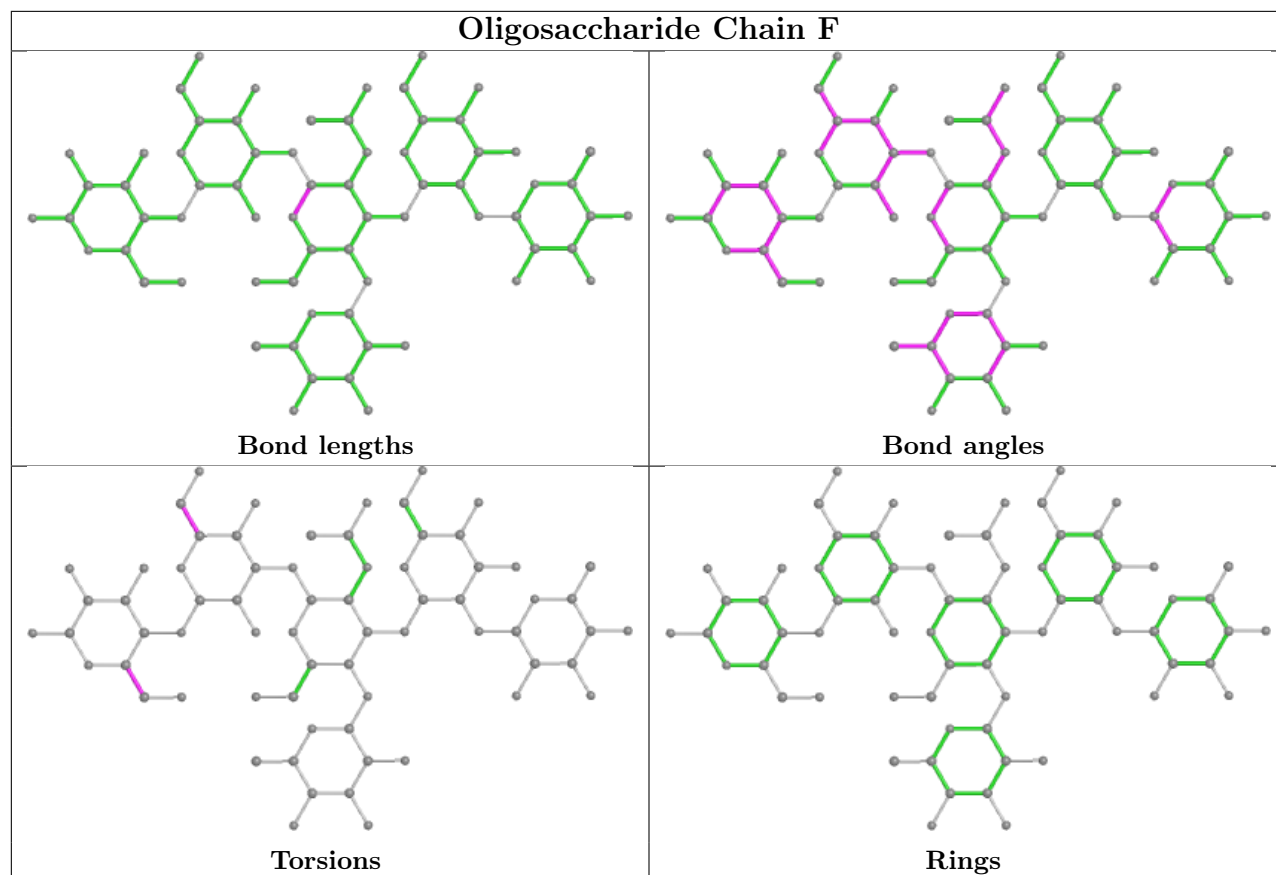
Mol	Chain	Res	Type	Atoms
3	F	2	GAL	O5-C5-C6-O6
3	F	1	BGC	O5-C5-C6-O6
3	F	2	GAL	C4-C5-C6-O6
3	F	1	BGC	C4-C5-C6-O6
3	E	1	BGC	O5-C5-C6-O6

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](#)

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	417/466 (89%)	-0.09	2 (0%) 91 91	44, 61, 105, 142	0
1	B	421/466 (90%)	-0.01	8 (1%) 66 63	45, 66, 112, 135	0
2	C	113/120 (94%)	-0.13	2 (1%) 68 64	47, 62, 80, 103	0
2	D	114/120 (95%)	-0.02	1 (0%) 84 83	43, 54, 73, 107	0
All	All	1065/1172 (90%)	-0.05	13 (1%) 79 77	43, 62, 107, 142	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	3	VAL	5.1
2	D	3	VAL	3.3
1	B	242	ARG	3.2
1	B	238	GLY	3.0
1	A	238	GLY	3.0
1	B	241	THR	2.9
1	B	111	ASN	2.8
1	B	237	LYS	2.8
1	B	408	GLY	2.7
1	A	242	ARG	2.6
1	B	245	TYR	2.5
1	B	232	VAL	2.2
2	C	103	ARG	2.1

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

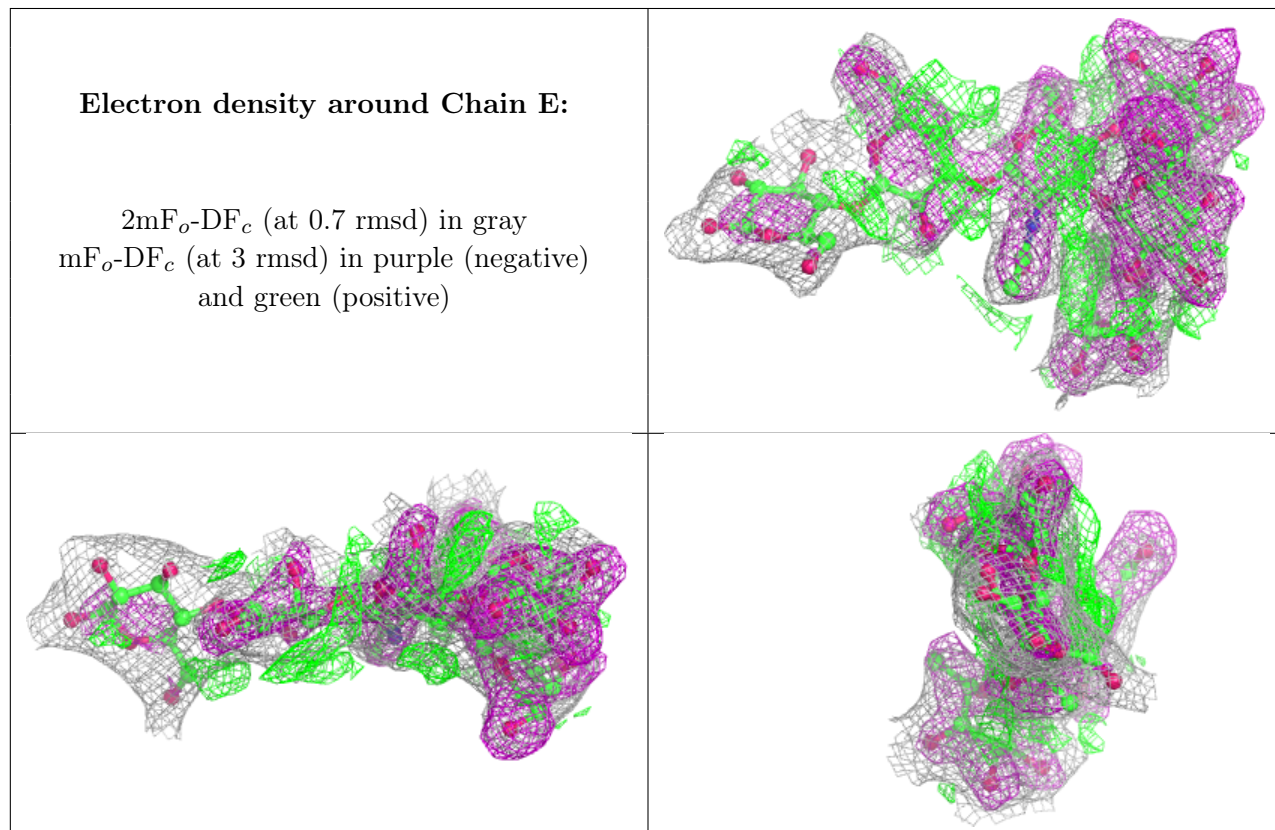
There are no non-standard protein/DNA/RNA residues in this entry.

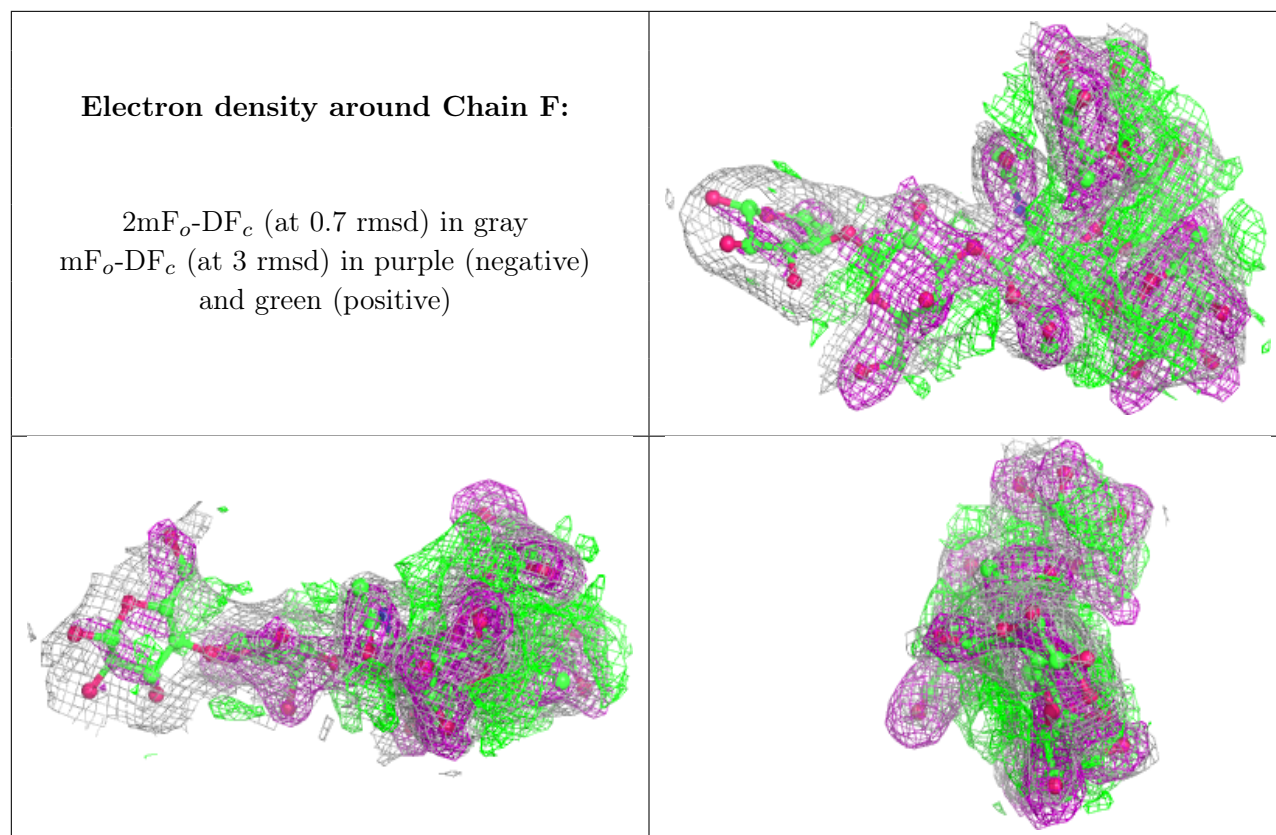
### 6.3 Carbohydrates [i](#)

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	GAL	F	2	11/12	0.86	0.31	43,48,52,56	0
3	GAL	F	4	11/12	0.86	0.30	26,29,31,31	0
3	BGC	E	1	12/12	0.91	0.28	57,68,76,80	0
3	NAG	F	3	14/15	0.91	0.27	34,38,40,43	0
3	BGC	F	1	12/12	0.91	0.27	54,61,65,66	0
3	FUC	F	6	10/11	0.91	0.37	32,35,37,37	0
3	FUC	E	6	10/11	0.92	0.28	34,36,36,37	0
3	GAL	E	2	11/12	0.92	0.27	42,46,50,51	0
3	NAG	E	3	14/15	0.94	0.19	32,35,37,39	0
3	FUC	F	5	10/11	0.94	0.26	29,31,31,32	0
3	GAL	E	4	11/12	0.94	0.22	29,31,33,33	0
3	FUC	E	5	10/11	0.95	0.18	24,27,30,30	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.