



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 10:39 PM UTC

PDB ID : 5F7M / pdb_00005f7m
Title : Blood group antigen binding adhesin BabA of Helicobacter pylori strain 17875
in complex with blood group H Lewis b hexasaccharide
Authors : Moonens, K.; Gideonsson, P.; Subedi, S.; Romao, E.; Oscarson, S.; Muylder-
mans, S.; Boren, T.; Remaut, H.
Deposited on : 2015-12-08
Resolution : 2.72 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

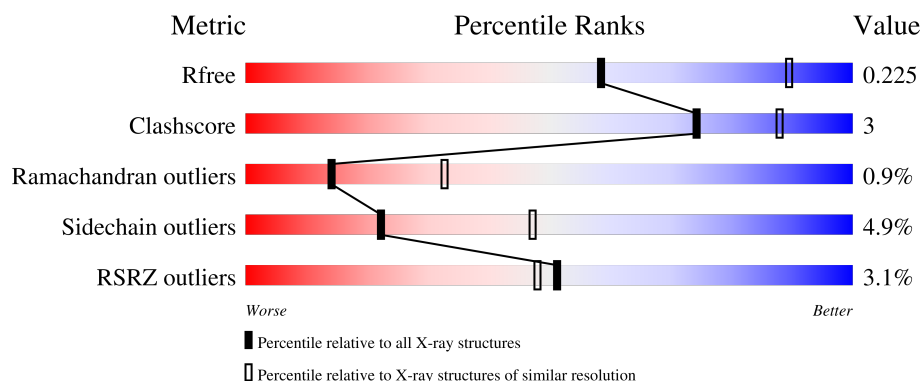
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.72 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}	180053	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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 $\leq 5\%$. 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	4% 79% 10% • 10%
1	B	466	3% 78% 11% • 10%
2	C	120	% 84% 9% • 6%
2	D	120	% 86% 8% • 5%
3	E	6	100%

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Mol	Chain	Length	Quality of chain
3	F	6	 100%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8260 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	419	Total	C	N	O	S	0	0	0
			3114	1920	537	644	13			
1	B	420	Total	C	N	O	S	0	0	0
			3124	1926	540	645	13			

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
A	463	HIS	-	expression tag	UNP O52269

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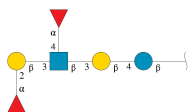
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Chain	Residue	Modelled	Actual	Comment	Reference
A	464	HIS	-	expression tag	UNP O52269
A	465	HIS	-	expression tag	UNP O52269
A	466	HIS	-	expression tag	UNP O52269
A	467	HIS	-	expression tag	UNP O52269
A	468	HIS	-	expression tag	UNP O52269
B	3	ALA	-	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
2	C	113	Total	C	N	O	S	0	0	0
			863	540	155	163	5			
2	D	114	Total	C	N	O	S	0	0	0
			873	546	158	164	5			

- 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
3	E	6	Total	C	N	O	0	0	0
			68	38	1	29			
3	F	6	Total	C	N	O	0	0	0
			68	38	1	29			

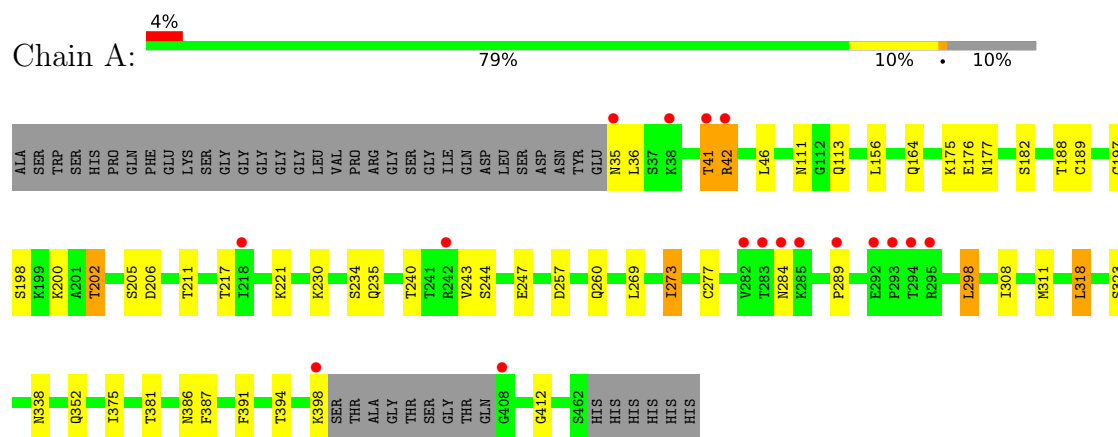
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	68	Total	O	0	0
			68	68		
4	B	40	Total	O	0	0
			40	40		
4	C	14	Total	O	0	0
			14	14		
4	D	28	Total	O	0	0
			28	28		

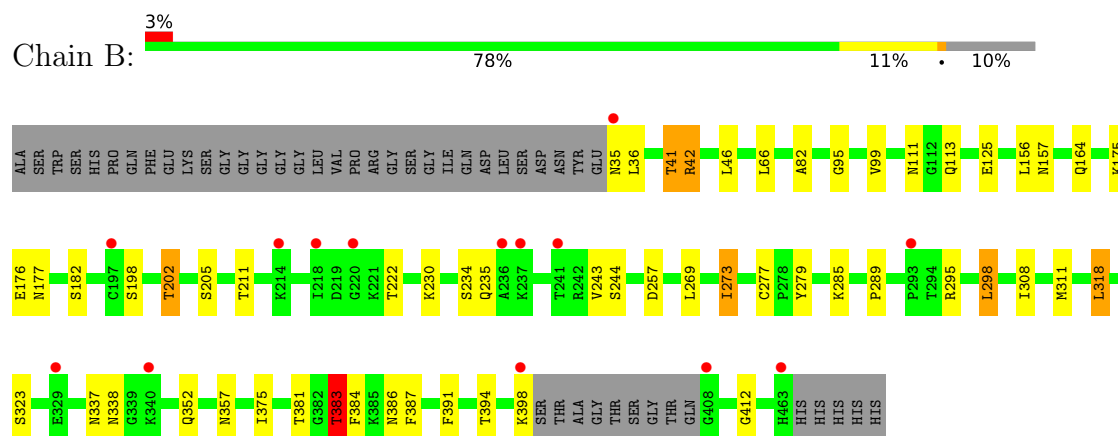
3 Residue-property plots [i](#)

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.

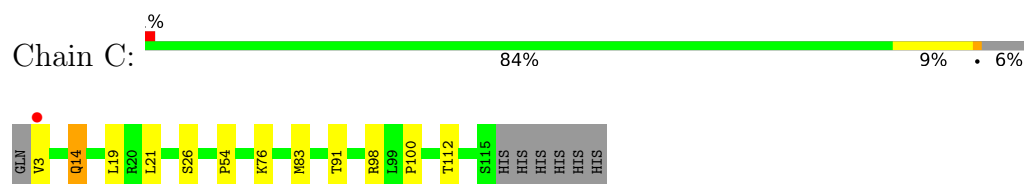
- Molecule 1: Adhesin binding fucosylated histo-blood group antigen



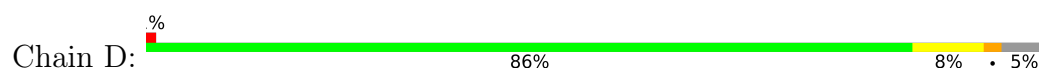
- Molecule 1: Adhesin binding fucosylated histo-blood group antigen



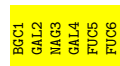
- Molecule 2: Nanobody Nb-ER19



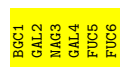
- Molecule 2: Nanobody Nb-ER19



- 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, α , β , γ	51.10Å 131.89Å 123.10Å 90.00° 94.77° 90.00°	Depositor
Resolution (Å)	45.00 – 2.72 45.00 – 2.72	Depositor EDS
% Data completeness (in resolution range)	99.5 (45.00-2.72) 99.5 (45.00-2.72)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.179 , 0.218 0.189 , 0.225	Depositor DCC
R_{free} test set	2194 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	42.5	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8260	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, GAL, BGC, FUC

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	1.08	1/3160 (0.0%)	1.02	3/4295 (0.1%)
1	B	1.02	1/3171 (0.0%)	1.01	4/4310 (0.1%)
2	C	1.02	0/882	0.98	2/1195 (0.2%)
2	D	1.11	0/893	1.01	3/1210 (0.2%)
All	All	1.05	2/8106 (0.0%)	1.01	12/11010 (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
2	D	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	277	CYS	C-O	-5.49	1.21	1.23
1	B	277	CYS	C-O	-5.02	1.21	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	14	GLN	CA-C-N	-5.91	113.67	119.76
2	C	14	GLN	C-N-CA	-5.91	113.67	119.76
2	D	115	SER	N-CA-C	5.89	118.81	110.50
2	D	14	GLN	CA-C-N	-5.59	114.00	119.76
2	D	14	GLN	C-N-CA	-5.59	114.00	119.76
1	B	383	THR	N-CA-C	5.40	117.86	111.33
1	B	222	THR	N-CA-C	5.33	116.14	107.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	260	GLN	N-CA-C	-5.31	105.49	111.28
1	A	273	ILE	CB-CA-C	-5.27	105.23	111.97
1	A	189	CYS	CB-CA-C	-5.07	102.02	110.14
1	B	273	ILE	CB-CA-C	-5.07	105.48	111.97
1	B	383	THR	N-CA-CB	-5.06	102.47	110.06

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	3	VAL	Peptide

5.2 Too-close contacts

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	3114	0	3044	20	0
1	B	3124	0	3051	23	0
2	C	863	0	845	4	0
2	D	873	0	852	3	0
3	E	68	0	60	0	0
3	F	68	0	60	0	0
4	A	68	0	0	1	0
4	B	40	0	0	0	0
4	C	14	0	0	0	0
4	D	28	0	0	0	0
All	All	8260	0	7912	50	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.

All (50) 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:B:279:TYR:OH	1:B:295:ARG:HD2	1.82	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ASN:OD1	1:A:36:LEU:N	2.22	0.72
1:A:35:ASN:OD1	1:A:35:ASN:C	2.32	0.71
1:B:35:ASN:OD1	1:B:36:LEU:N	2.24	0.71
1:B:35:ASN:OD1	1:B:35:ASN:C	2.35	0.70
1:A:217:THR:HA	1:A:221:LYS:O	1.98	0.63
2:C:3:VAL:N	2:C:26:SER:HG	2.00	0.59
1:B:391:PHE:O	1:B:394:THR:HB	2.07	0.54
1:A:391:PHE:O	1:A:394:THR:HB	2.09	0.53
1:B:318:LEU:C	1:B:318:LEU:HD23	2.34	0.53
1:A:234:SER:N	1:A:244:SER:HB3	2.25	0.51
1:A:198:SER:O	1:A:202:THR:HB	2.11	0.51
2:C:19:LEU:HB2	2:C:83:MET:HE3	1.94	0.50
1:B:198:SER:O	1:B:202:THR:HB	2.12	0.50
1:B:235:GLN:HA	1:B:235:GLN:OE1	2.13	0.48
1:A:235:GLN:HA	1:A:235:GLN:OE1	2.13	0.48
1:A:318:LEU:C	1:A:318:LEU:HD23	2.40	0.47
1:B:41:THR:O	1:B:42:ARG:C	2.58	0.47
1:A:197:CYS:HB2	4:A:636:HOH:O	2.15	0.47
2:D:19:LEU:HB2	2:D:83:MET:HE3	1.96	0.46
1:A:386:ASN:O	1:A:387:PHE:C	2.58	0.46
1:B:234:SER:N	1:B:244:SER:HB3	2.31	0.46
1:A:308:ILE:HA	1:A:311:MET:HE3	1.98	0.46
1:B:257:ASP:OD1	1:B:257:ASP:C	2.58	0.45
1:A:41:THR:O	1:A:42:ARG:C	2.58	0.45
1:A:311:MET:HE1	1:A:375:ILE:HD13	1.98	0.45
1:A:273:ILE:HG23	1:A:298:LEU:HD23	1.99	0.44
1:B:156:LEU:HG	1:B:269:LEU:HD12	2.00	0.44
1:A:257:ASP:OD1	1:A:257:ASP:C	2.60	0.44
1:A:175:LYS:HE3	1:A:176:GLU:OE2	2.18	0.43
1:B:125:GLU:OE2	1:B:157:ASN:ND2	2.38	0.43
1:B:311:MET:HE1	1:B:375:ILE:HD13	2.00	0.43
1:B:111:ASN:HB2	1:B:113:GLN:HG2	2.00	0.43
1:B:175:LYS:HE3	1:B:176:GLU:OE2	2.18	0.43
1:B:386:ASN:O	1:B:387:PHE:C	2.59	0.43
1:A:111:ASN:HB2	1:A:113:GLN:HG2	2.00	0.42
1:A:156:LEU:HG	1:A:269:LEU:HD12	2.00	0.42
2:D:91:THR:HG23	2:D:112:THR:HA	2.01	0.42
1:A:164:GLN:OE1	1:A:412:GLY:HA2	2.19	0.42
1:B:273:ILE:HG23	1:B:298:LEU:HD23	2.01	0.42
1:B:308:ILE:HA	1:B:311:MET:HE3	2.02	0.41
1:B:357:ASN:OD1	1:B:357:ASN:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:91:THR:HG23	2:C:112:THR:HA	2.01	0.41
2:C:98:ARG:HG3	2:C:100:PRO:O	2.21	0.41
1:B:82:ALA:HB1	1:B:318:LEU:HD21	2.01	0.41
1:B:164:GLN:OE1	1:B:412:GLY:HA2	2.20	0.41
1:B:383:THR:HG22	1:B:384:PHE:N	2.35	0.41
2:D:98:ARG:HG3	2:D:100:PRO:O	2.20	0.41
1:A:188:THR:HG22	1:A:247:GLU:HG3	2.03	0.41
1:B:95:GLY:O	1:B:99:VAL:HG23	2.20	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	415/466 (89%)	390 (94%)	21 (5%)	4 (1%)	12	30
1	B	416/466 (89%)	395 (95%)	16 (4%)	5 (1%)	10	25
2	C	111/120 (92%)	105 (95%)	6 (5%)	0	100	100
2	D	112/120 (93%)	107 (96%)	5 (4%)	0	100	100
All	All	1054/1172 (90%)	997 (95%)	48 (5%)	9 (1%)	14	33

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	338	ASN
1	A	41	THR
1	A	338	ASN
1	B	42	ARG
1	A	205	SER
1	B	41	THR
1	B	205	SER

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Mol	Chain	Res	Type
1	B	289	PRO
1	A	289	PRO

5.3.2 Protein sidechains ⓘ

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	345/381 (91%)	327 (95%)	18 (5%)	21	45
1	B	346/381 (91%)	329 (95%)	17 (5%)	22	47
2	C	92/99 (93%)	88 (96%)	4 (4%)	26	52
2	D	93/99 (94%)	89 (96%)	4 (4%)	26	52
All	All	876/960 (91%)	833 (95%)	43 (5%)	22	47

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ARG
1	A	46	LEU
1	A	177	ASN
1	A	182	SER
1	A	200	LYS
1	A	202	THR
1	A	206	ASP
1	A	211	THR
1	A	230	LYS
1	A	240	THR
1	A	243	VAL
1	A	284	ASN
1	A	298	LEU
1	A	318	LEU
1	A	323	SER
1	A	352	GLN
1	A	381	THR
1	A	398	LYS
1	B	46	LEU

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Mol	Chain	Res	Type
1	B	66	LEU
1	B	177	ASN
1	B	182	SER
1	B	202	THR
1	B	211	THR
1	B	230	LYS
1	B	243	VAL
1	B	285	LYS
1	B	298	LEU
1	B	318	LEU
1	B	323	SER
1	B	337	ASN
1	B	352	GLN
1	B	381	THR
1	B	383	THR
1	B	398	LYS
2	C	14	GLN
2	C	21	LEU
2	C	54	PRO
2	C	76	LYS
2	D	3	VAL
2	D	21	LEU
2	D	76	LYS
2	D	115	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	ASN
1	A	157	ASN
1	A	161	GLN
1	A	178	ASN
1	A	274	ASN
1	A	316	GLN
1	A	352	GLN
1	A	361	GLN
1	A	366	ASN
1	A	396	ASN
1	B	124	ASN
1	B	157	ASN
1	B	161	GLN
1	B	260	GLN

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Mol	Chain	Res	Type
1	B	265	GLN
1	B	274	ASN
1	B	290	GLN
1	B	321	GLN
1	B	337	ASN
1	B	366	ASN
1	B	396	ASN
2	C	40	GLN
2	C	82	GLN
2	D	40	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

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	0.57	0	17,17,17	1.53	3 (17%)
3	GAL	E	2	3	11,11,12	0.77	0	15,15,17	1.20	1 (6%)
3	NAG	E	3	3	14,14,15	0.79	1 (7%)	17,19,21	1.43	2 (11%)
3	GAL	E	4	3	11,11,12	1.00	0	15,15,17	1.38	1 (6%)
3	FUC	E	5	3	10,10,11	0.74	0	14,14,16	1.36	2 (14%)
3	FUC	E	6	3	10,10,11	0.69	0	14,14,16	1.28	1 (7%)
3	BGC	F	1	3	12,12,12	0.51	0	17,17,17	1.25	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GAL	F	2	3	11,11,12	0.64	0	15,15,17	1.73	4 (26%)
3	NAG	F	3	3	14,14,15	0.75	0	17,19,21	1.40	2 (11%)
3	GAL	F	4	3	11,11,12	0.88	1 (9%)	15,15,17	1.53	1 (6%)
3	FUC	F	5	3	10,10,11	0.89	0	14,14,16	1.19	1 (7%)
3	FUC	F	6	3	10,10,11	0.88	0	14,14,16	1.43	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
3	BGC	E	1	3	-	0/2/22/22	0/1/1/1
3	GAL	E	2	3	-	1/2/19/22	0/1/1/1
3	NAG	E	3	3	-	0/6/23/26	0/1/1/1
3	GAL	E	4	3	-	1/2/19/22	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	-	0/2/22/22	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	-	2/2/19/22	0/1/1/1
3	FUC	F	5	3	-	-	0/1/1/1
3	FUC	F	6	3	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	4	GAL	O5-C1	-2.06	1.40	1.43
3	E	3	NAG	O5-C1	-2.00	1.40	1.43

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	4	GAL	C1-O5-C5	5.22	119.18	112.19
3	F	2	GAL	C1-C2-C3	4.15	115.68	109.64
3	F	1	BGC	C4-C3-C2	-3.87	104.03	110.83
3	E	4	GAL	C1-O5-C5	-3.60	107.36	112.19
3	E	1	BGC	C3-C4-C5	3.37	116.34	110.23
3	F	6	FUC	C1-C2-C3	3.27	114.40	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	3	NAG	C1-O5-C5	3.21	116.48	112.19
3	E	2	GAL	O3-C3-C2	-3.08	103.77	110.05
3	F	3	NAG	O7-C7-N2	3.02	127.32	121.98
3	F	3	NAG	O7-C7-C8	-2.96	116.78	122.05
3	E	6	FUC	O2-C2-C3	-2.70	104.56	110.15
3	E	5	FUC	O5-C5-C6	2.61	113.06	107.40
3	F	2	GAL	O5-C5-C6	2.45	112.42	107.66
3	F	2	GAL	O3-C3-C4	2.24	115.65	110.38
3	E	3	NAG	C1-C2-N2	2.22	113.94	110.43
3	E	1	BGC	O5-C1-C2	-2.13	106.55	110.30
3	E	5	FUC	O2-C2-C3	-2.12	105.76	110.15
3	F	5	FUC	C1-C2-C3	2.12	112.73	109.64
3	E	1	BGC	O5-C5-C4	2.11	113.50	109.70
3	F	2	GAL	C3-C4-C5	-2.07	106.48	110.23

There are no chirality outliers.

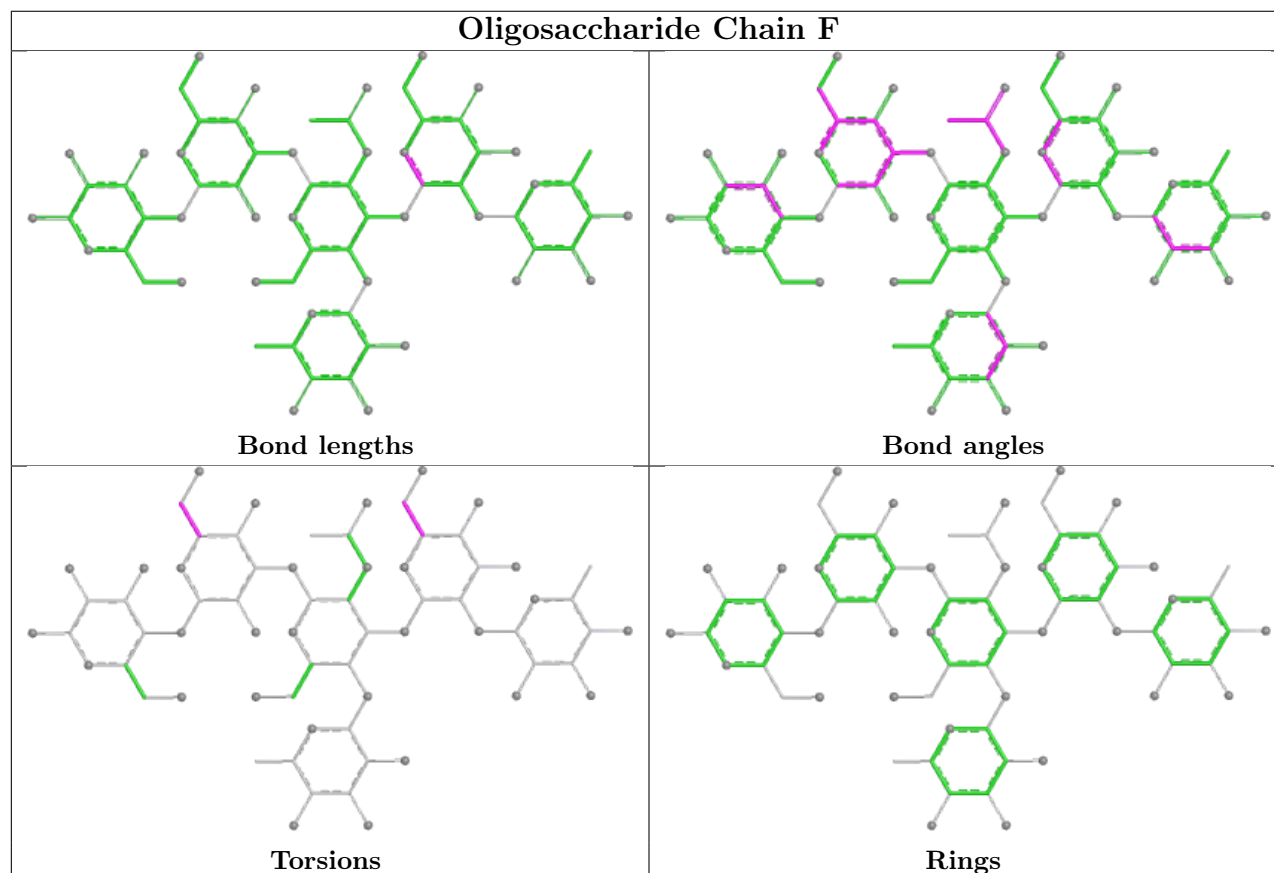
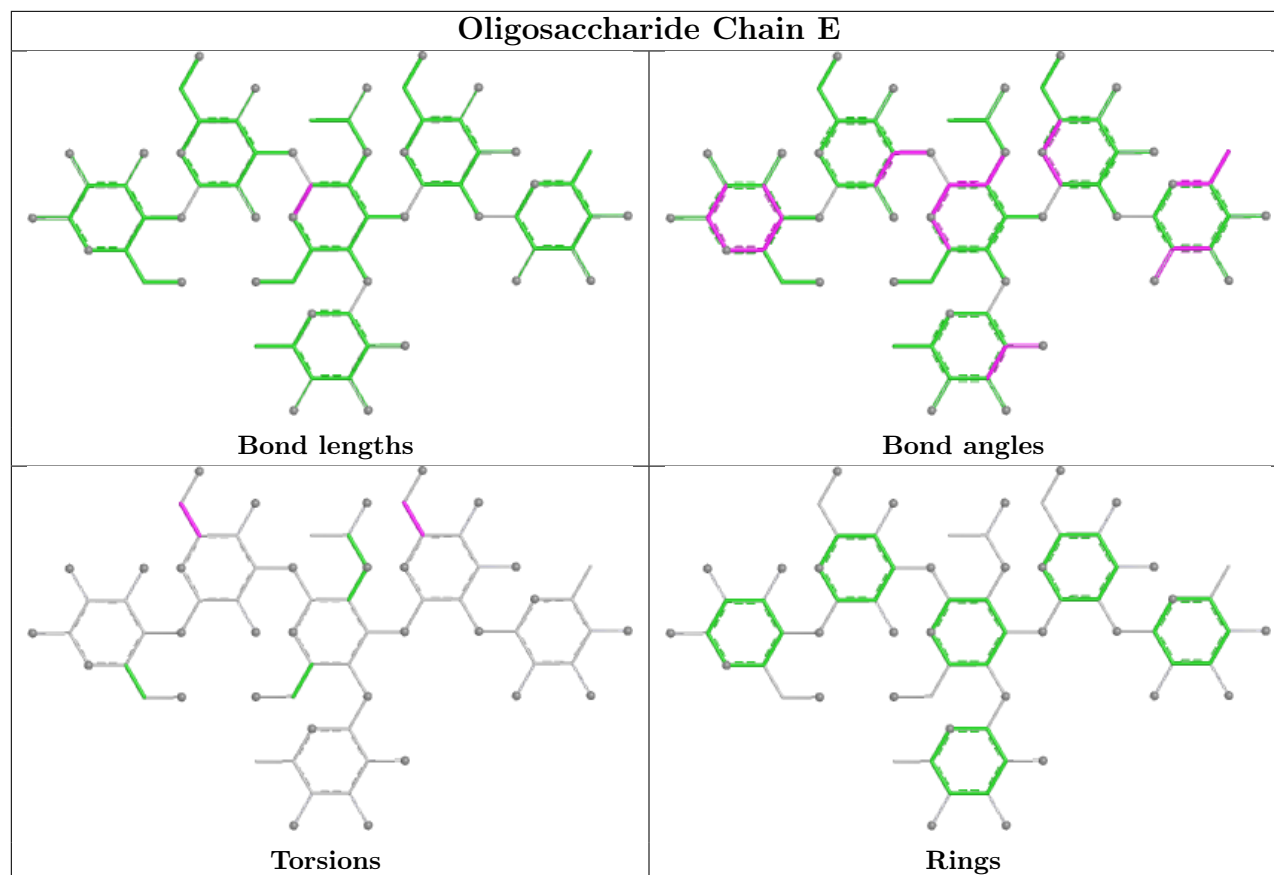
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	4	GAL	C4-C5-C6-O6
3	F	4	GAL	O5-C5-C6-O6
3	F	2	GAL	C4-C5-C6-O6
3	F	2	GAL	O5-C5-C6-O6
3	E	4	GAL	O5-C5-C6-O6
3	E	2	GAL	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

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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, 95th 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	A	419/466 (89%)	-0.11	17 (4%) 41 38	25, 44, 97, 128	0
1	B	420/466 (90%)	0.03	14 (3%) 49 45	28, 49, 96, 120	0
2	C	113/120 (94%)	-0.28	1 (0%) 81 80	28, 45, 66, 95	0
2	D	114/120 (95%)	-0.53	1 (0%) 81 80	22, 36, 57, 77	0
All	All	1066/1172 (90%)	-0.11	33 (3%) 51 48	22, 45, 93, 128	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	3	VAL	6.0
1	B	218	ILE	4.9
1	A	218	ILE	4.5
1	B	237	LYS	4.4
1	B	398	LYS	4.2
1	B	408	GLY	3.4
1	A	398	LYS	3.3
1	A	282	VAL	3.3
1	B	293	PRO	3.0
1	A	38	LYS	3.0
1	A	295	ARG	3.0
1	A	35	ASN	2.7
1	A	292	GLU	2.7
1	B	463	HIS	2.6
1	B	35	ASN	2.5
1	A	294	THR	2.5
1	B	340	LYS	2.5
1	A	293	PRO	2.4
1	B	214	LYS	2.4
1	B	197	CYS	2.4
1	A	41	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	285	LYS	2.4
1	B	220	GLY	2.4
1	B	329	GLU	2.3
1	B	241	THR	2.2
1	A	42	ARG	2.2
1	B	236	ALA	2.1
1	A	283	THR	2.1
1	A	289	PRO	2.1
1	A	408	GLY	2.1
2	D	3	VAL	2.1
1	A	242	ARG	2.1
1	A	284	ASN	2.0

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

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

6.3 Carbohydrates ⓘ

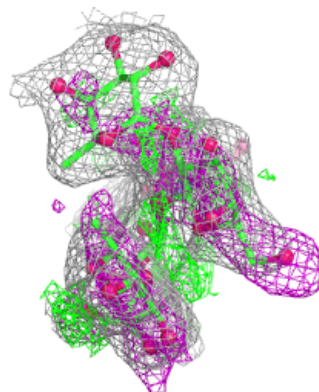
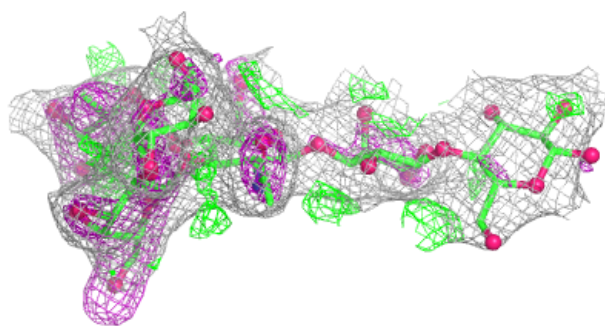
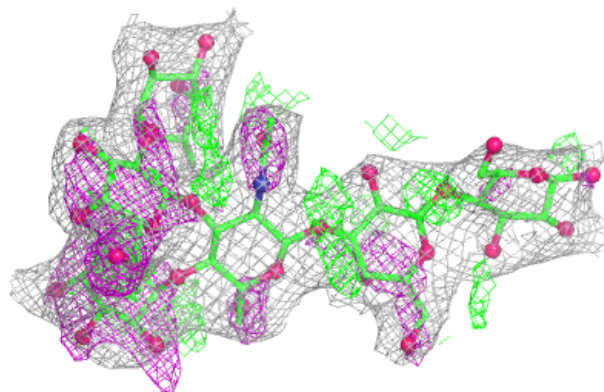
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	BGC	F	1	12/12	0.60	0.16	47,50,52,53	0
3	BGC	E	1	12/12	0.61	0.16	56,64,68,69	0
3	GAL	F	2	11/12	0.65	0.16	33,40,44,44	0
3	GAL	F	4	11/12	0.76	0.15	22,24,25,26	0
3	FUC	F	6	10/11	0.76	0.17	23,24,25,25	0
3	GAL	E	2	11/12	0.80	0.12	35,40,45,47	0
3	NAG	F	3	14/15	0.80	0.13	25,28,30,30	0
3	FUC	E	6	10/11	0.85	0.14	25,27,28,28	0
3	FUC	F	5	10/11	0.88	0.11	21,24,25,25	0
3	GAL	E	4	11/12	0.88	0.12	21,24,25,25	0
3	NAG	E	3	14/15	0.91	0.10	22,27,29,30	0
3	FUC	E	5	10/11	0.95	0.08	17,18,19,20	0

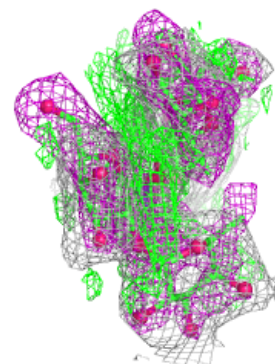
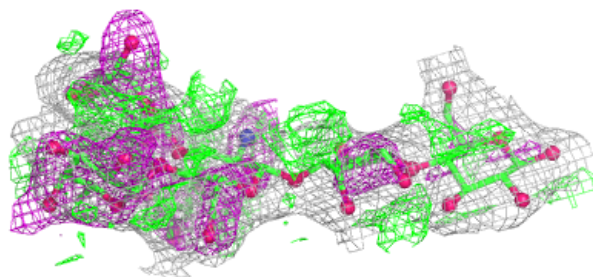
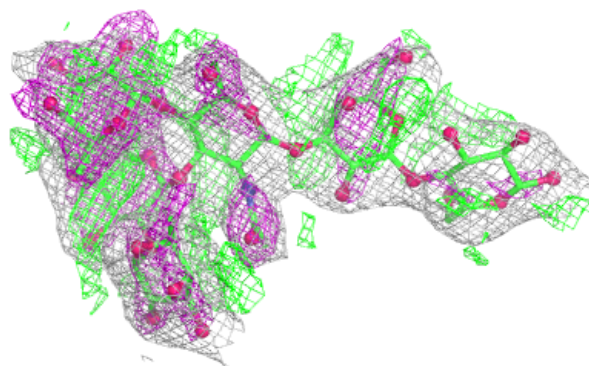
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around Chain F:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.