



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 07:09 PM BST

PDB ID : 4P3I
Title : Structure of the P domain from a GI.7 Norovirus variant in complex with LeA HBGA.
Authors : Shanker, S.; Czako, R.; Sankaran, B.; Atmar, R.; Estes, M.; Prasad, B.V.V.
Deposited on : 2014-03-07
Resolution : 1.69 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

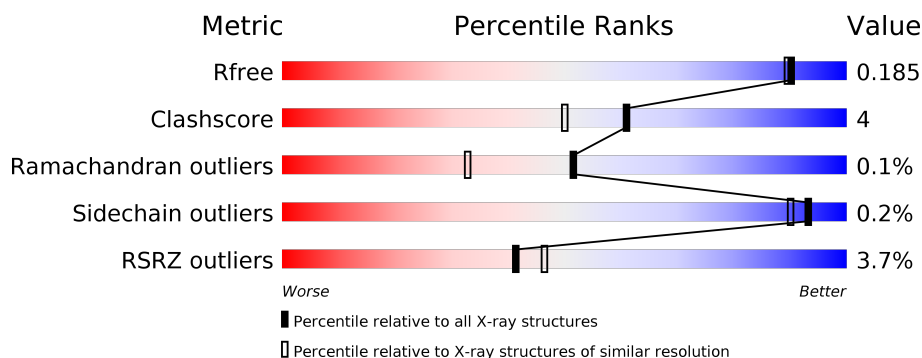
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.69 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	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 on 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	301	<div> <div>4%</div> <div>88%</div> <div>8%</div> <div>•</div> </div>
1	B	301	<div> <div>3%</div> <div>87%</div> <div>7%</div> <div>6%</div> </div>
1	C	301	<div> <div>4%</div> <div>84%</div> <div>9%</div> <div>•</div> <div>6%</div> </div>
1	D	301	<div> <div>3%</div> <div>90%</div> <div>6%</div> <div>•</div> </div>
2	E	3	<div> <div>33%</div> <div>33%</div> <div>33%</div> </div>
2	F	3	<div> <div>67%</div> <div>33%</div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	G	3	<div><div></div><div>33%</div><div>33%</div><div>33%</div></div>
2	H	3	<div><div></div><div>33%</div><div>67%</div></div>

2 Entry composition [i](#)

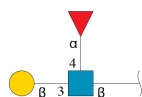
There are 3 unique types of molecules in this entry. The entry contains 10067 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 P domain of VP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	0	0	0
			2205	1407	373	417	8			
1	B	282	Total	C	N	O	S	0	0	0
			2147	1377	362	400	8			
1	C	282	Total	C	N	O	S	0	0	0
			2142	1369	361	404	8			
1	D	289	Total	C	N	O	S	0	0	0
			2192	1399	367	418	8			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	0	0	0
			36	20	1	15			
2	F	3	Total	C	N	O	0	0	0
			36	20	1	15			
2	G	3	Total	C	N	O	0	0	0
			36	20	1	15			
2	H	3	Total	C	N	O	0	0	0
			36	20	1	15			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	297	Total	O	0	0
			297	297		

Continued on next page...

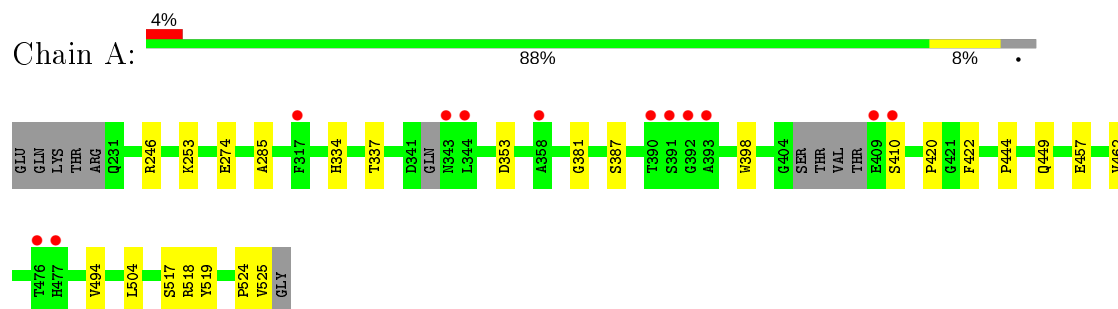
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	305	Total 305	O 305	0	0
3	C	310	Total 310	O 310	0	0
3	D	325	Total 325	O 325	0	0

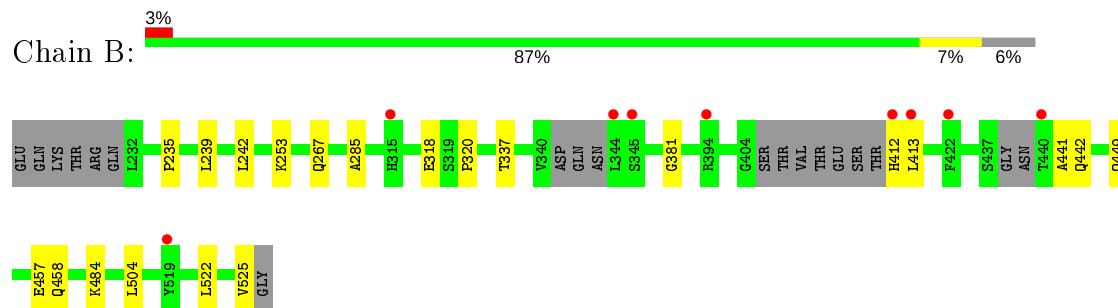
3 Residue-property plots [i](#)

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.

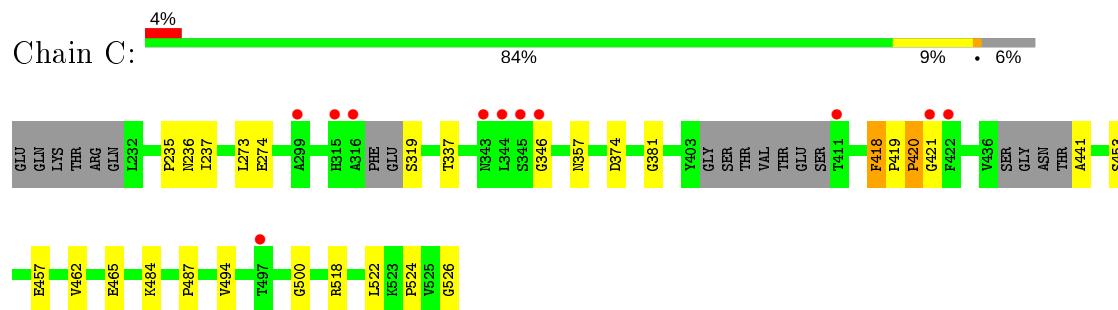
- Molecule 1: P domain of VP1



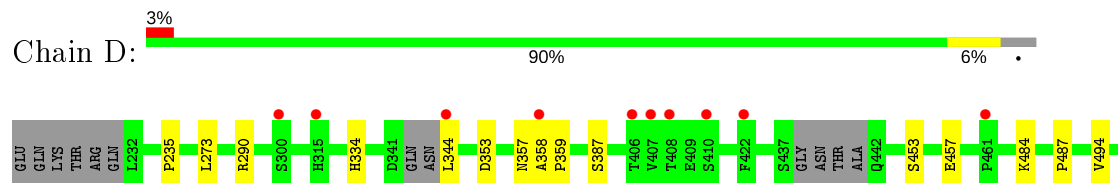
- Molecule 1: P domain of VP1

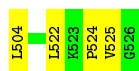


- Molecule 1: P domain of VP1

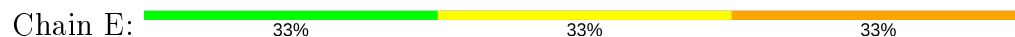


- Molecule 1: P domain of VP1





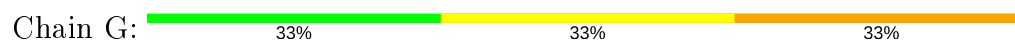
- Molecule 2: beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose



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



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.17Å 63.28Å 90.40Å 72.77° 82.20° 60.31°	Depositor
Resolution (Å)	38.92 – 1.69 38.92 – 1.69	Depositor EDS
% Data completeness (in resolution range)	96.2 (38.92-1.69) 96.2 (38.92-1.69)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 1.70Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.155 , 0.177 0.166 , 0.185	Depositor DCC
R_{free} test set	6146 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	15.8	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.008 for h-k,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10067	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% 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: GAL, NAG, 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	0.34	0/2270	0.55	0/3107
1	B	0.33	0/2211	0.55	0/3023
1	C	0.42	1/2204 (0.0%)	0.67	6/3015 (0.2%)
1	D	0.38	0/2257	0.57	0/3089
All	All	0.37	1/8942 (0.0%)	0.59	6/12234 (0.0%)

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	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	421	GLY	C-N	8.56	1.53	1.34

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	421	GLY	O-C-N	11.61	141.28	122.70
1	C	421	GLY	CA-C-N	-9.95	95.31	117.20
1	C	421	GLY	C-N-CA	-7.01	104.18	121.70
1	C	346	GLY	N-CA-C	6.52	129.41	113.10
1	C	419	PRO	C-N-CD	-5.86	107.70	120.60
1	C	418	PHE	O-C-N	-5.17	111.28	121.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	344	LEU	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	2205	0	2125	14	0
1	B	2147	0	2078	16	0
1	C	2142	0	2067	17	0
1	D	2192	0	2105	22	0
2	E	36	0	33	2	0
2	F	36	0	33	3	0
2	G	36	0	33	6	0
2	H	36	0	33	3	0
3	A	297	0	0	0	0
3	B	305	0	0	2	0
3	C	310	0	0	5	0
3	D	325	0	0	4	0
All	All	10067	0	8507	78	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 (78) 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)
2:G:1:NAG:O3	2:G:3:FUC:H5	1.56	1.04
1:B:318:GLU:O	1:B:412:HIS:N	1.97	0.97
1:D:357:ASN:C	1:D:359:PRO:HD2	1.86	0.95
1:B:320:PRO:CG	1:B:413:LEU:HD23	2.11	0.80
2:F:1:NAG:O3	2:F:3:FUC:H5	1.83	0.78
1:D:358:ALA:N	1:D:359:PRO:HD2	2.00	0.76
1:D:358:ALA:N	1:D:359:PRO:CD	2.50	0.74
1:B:253:LYS:HE2	1:B:442:GLN:OE1	1.89	0.73
1:B:457:GLU:OE1	1:B:458:GLN:NE2	2.22	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:357:ASN:C	1:D:359:PRO:CD	2.61	0.68
1:D:290:ARG:HD2	3:D:991:HOH:O	1.95	0.67
1:B:320:PRO:HG3	1:B:413:LEU:HD23	1.77	0.65
1:D:334:HIS:HE1	1:D:387:SER:OG	1.79	0.65
1:B:320:PRO:HG2	1:B:413:LEU:HD23	1.80	0.64
1:D:290:ARG:CD	3:D:991:HOH:O	2.45	0.64
1:B:267:GLN:CD	1:B:413:LEU:HD22	2.18	0.64
1:A:420:PRO:O	1:A:518:ARG:NH2	2.19	0.62
2:E:1:NAG:H61	2:E:3:FUC:C1	2.30	0.62
1:A:504:LEU:HD11	1:A:525:VAL:HG11	1.81	0.61
1:A:334:HIS:HE1	1:A:387:SER:OG	1.84	0.60
1:C:319:SER:N	3:C:938:HOH:O	2.34	0.59
1:C:457:GLU:OE2	1:D:235:PRO:HA	2.03	0.58
1:B:337:THR:HG22	1:B:381:GLY:HA3	1.87	0.57
1:C:484:LYS:HG3	1:C:522:LEU:HD21	1.87	0.56
3:C:1000:HOH:O	2:G:1:NAG:H81	2.05	0.56
1:C:274:GLU:HG3	1:C:462:VAL:HG23	1.88	0.55
1:C:273:LEU:HG	1:C:487:PRO:HA	1.88	0.55
2:G:1:NAG:C4	2:G:3:FUC:H5	2.38	0.54
1:C:441:ALA:N	3:C:967:HOH:O	2.42	0.52
3:D:703:HOH:O	2:H:1:NAG:H83	2.10	0.52
1:D:290:ARG:CG	3:D:991:HOH:O	2.58	0.51
1:D:504:LEU:HD11	1:D:525:VAL:HG21	1.94	0.50
1:B:239:LEU:HD12	1:B:242:LEU:HD12	1.94	0.50
1:D:334:HIS:HD2	1:D:353:ASP:OD1	1.95	0.50
1:C:235:PRO:HA	1:D:457:GLU:OE2	2.12	0.49
2:G:1:NAG:C3	2:G:3:FUC:H5	2.39	0.49
1:A:334:HIS:HD2	1:A:353:ASP:OD1	1.95	0.48
1:A:337:THR:HG22	1:A:381:GLY:HA3	1.96	0.48
1:A:457:GLU:OE1	1:B:235:PRO:HA	2.14	0.48
1:C:494:VAL:O	1:C:524:PRO:HA	2.14	0.48
1:A:274:GLU:HG3	1:A:462:VAL:HG23	1.94	0.47
1:C:236:ASN:HB3	1:D:457:GLU:HG2	1.96	0.47
1:C:237:ILE:HG13	1:D:457:GLU:HG3	1.96	0.47
1:C:337:THR:HG22	1:C:381:GLY:HA3	1.96	0.47
2:H:1:NAG:H61	2:H:3:FUC:C1	2.45	0.47
2:H:2:GAL:O5	2:H:3:FUC:H5	2.16	0.46
1:C:465:GLU:HB3	1:C:518:ARG:HB3	1.98	0.45
1:A:253:LYS:HE3	1:A:253:LYS:HB2	1.69	0.45
2:G:1:NAG:H61	2:G:3:FUC:C1	2.46	0.45
1:D:504:LEU:HD11	1:D:525:VAL:CG2	2.48	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:484:LYS:HG3	1:D:522:LEU:HD21	1.98	0.44
1:C:357:ASN:HB2	3:C:892:HOH:O	2.17	0.44
1:D:357:ASN:O	1:D:358:ALA:C	2.56	0.43
1:C:453:SER:O	1:C:457:GLU:HB2	2.18	0.43
1:D:273:LEU:HD13	1:D:487:PRO:HA	2.00	0.43
1:A:420:PRO:HB2	1:A:518:ARG:NH2	2.33	0.43
2:G:1:NAG:C4	2:G:3:FUC:C5	2.96	0.43
1:A:494:VAL:O	1:A:524:PRO:HA	2.18	0.43
1:C:418:PHE:O	1:C:420:PRO:HD3	2.18	0.43
1:B:412:HIS:N	3:B:969:HOH:O	2.51	0.43
1:D:494:VAL:O	1:D:524:PRO:HA	2.18	0.43
1:C:500:GLY:O	1:C:526:GLY:HA3	2.18	0.42
1:B:484:LYS:HG3	1:B:522:LEU:HD21	2.02	0.42
1:A:422:PHE:CZ	1:A:519:TYR:HD1	2.37	0.42
2:F:1:NAG:H61	2:F:3:FUC:C1	2.49	0.42
2:F:2:GAL:C1	2:F:3:FUC:H5	2.49	0.42
2:E:1:NAG:C6	2:E:3:FUC:C1	2.96	0.42
1:B:504:LEU:HD11	1:B:525:VAL:HG21	2.02	0.42
1:A:285:ALA:HA	1:A:449:GLN:HG2	2.02	0.41
1:B:441:ALA:HB1	3:B:973:HOH:O	2.20	0.41
1:B:285:ALA:HA	1:B:449:GLN:HG2	2.03	0.41
1:B:267:GLN:OE1	1:B:413:LEU:CD2	2.69	0.41
1:A:398:TRP:CE2	1:A:444:PRO:HG3	2.55	0.41
1:A:246:ARG:HD3	1:A:444:PRO:O	2.21	0.41
1:C:374:ASP:N	3:C:943:HOH:O	2.41	0.41
1:D:334:HIS:CE1	1:D:387:SER:OG	2.67	0.40
1:D:357:ASN:O	1:D:359:PRO:N	2.54	0.40
1:D:453:SER:O	1:D:457:GLU:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	284/301 (94%)	273 (96%)	11 (4%)	0	100	100
1	B	274/301 (91%)	264 (96%)	10 (4%)	0	100	100
1	C	274/301 (91%)	263 (96%)	10 (4%)	1 (0%)	34	18
1	D	283/301 (94%)	276 (98%)	7 (2%)	0	100	100
All	All	1115/1204 (93%)	1076 (96%)	38 (3%)	1 (0%)	51	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	420	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	244/257 (95%)	242 (99%)	2 (1%)	81	74
1	B	239/257 (93%)	239 (100%)	0	100	100
1	C	236/257 (92%)	236 (100%)	0	100	100
1	D	243/257 (95%)	243 (100%)	0	100	100
All	All	962/1028 (94%)	960 (100%)	2 (0%)	93	90

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	410	SER
1	A	517	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	334	HIS
1	D	334	HIS
1	D	470	HIS

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$
2	NAG	E	1	2	15,15,15	0.64	0	21,21,21	0.82	1 (4%)
2	GAL	E	2	2	11,11,12	0.28	0	15,15,17	0.84	0
2	FUC	E	3	2	10,10,11	0.33	0	14,14,16	0.79	0
2	NAG	F	1	2	15,15,15	0.64	0	21,21,21	0.83	1 (4%)
2	GAL	F	2	2	11,11,12	0.30	0	15,15,17	0.85	0
2	FUC	F	3	2	10,10,11	0.33	0	14,14,16	0.79	0
2	NAG	G	1	2	15,15,15	0.64	0	21,21,21	0.82	1 (4%)
2	GAL	G	2	2	11,11,12	0.28	0	15,15,17	0.85	0
2	FUC	G	3	2	10,10,11	0.34	0	14,14,16	0.79	0
2	NAG	H	1	2	15,15,15	0.63	0	21,21,21	0.82	1 (4%)
2	GAL	H	2	2	11,11,12	0.29	0	15,15,17	0.85	1 (6%)
2	FUC	H	3	2	10,10,11	0.34	0	14,14,16	0.79	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	E	1	2	-	0/6/26/26	0/1/1/1
2	GAL	E	2	2	-	0/2/19/22	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FUC	E	3	2	-	-	0/1/1/1
2	NAG	F	1	2	-	3/6/26/26	0/1/1/1
2	GAL	F	2	2	-	0/2/19/22	0/1/1/1
2	FUC	F	3	2	-	-	0/1/1/1
2	NAG	G	1	2	-	0/6/26/26	0/1/1/1
2	GAL	G	2	2	-	0/2/19/22	0/1/1/1
2	FUC	G	3	2	-	-	0/1/1/1
2	NAG	H	1	2	-	1/6/26/26	0/1/1/1
2	GAL	H	2	2	-	0/2/19/22	0/1/1/1
2	FUC	H	3	2	-	-	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	NAG	C1-O5-C5	-2.23	109.45	113.66
2	E	1	NAG	C1-O5-C5	-2.22	109.47	113.66
2	G	1	NAG	C1-O5-C5	-2.21	109.50	113.66
2	H	1	NAG	C1-O5-C5	-2.17	109.56	113.66
2	H	2	GAL	C1-O5-C5	-2.00	109.48	112.19

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	1	NAG	O5-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	F	1	NAG	C3-C2-N2-C7
2	H	1	NAG	O5-C5-C6-O6

There are no ring outliers.

10 monomers are involved in 14 short contacts:

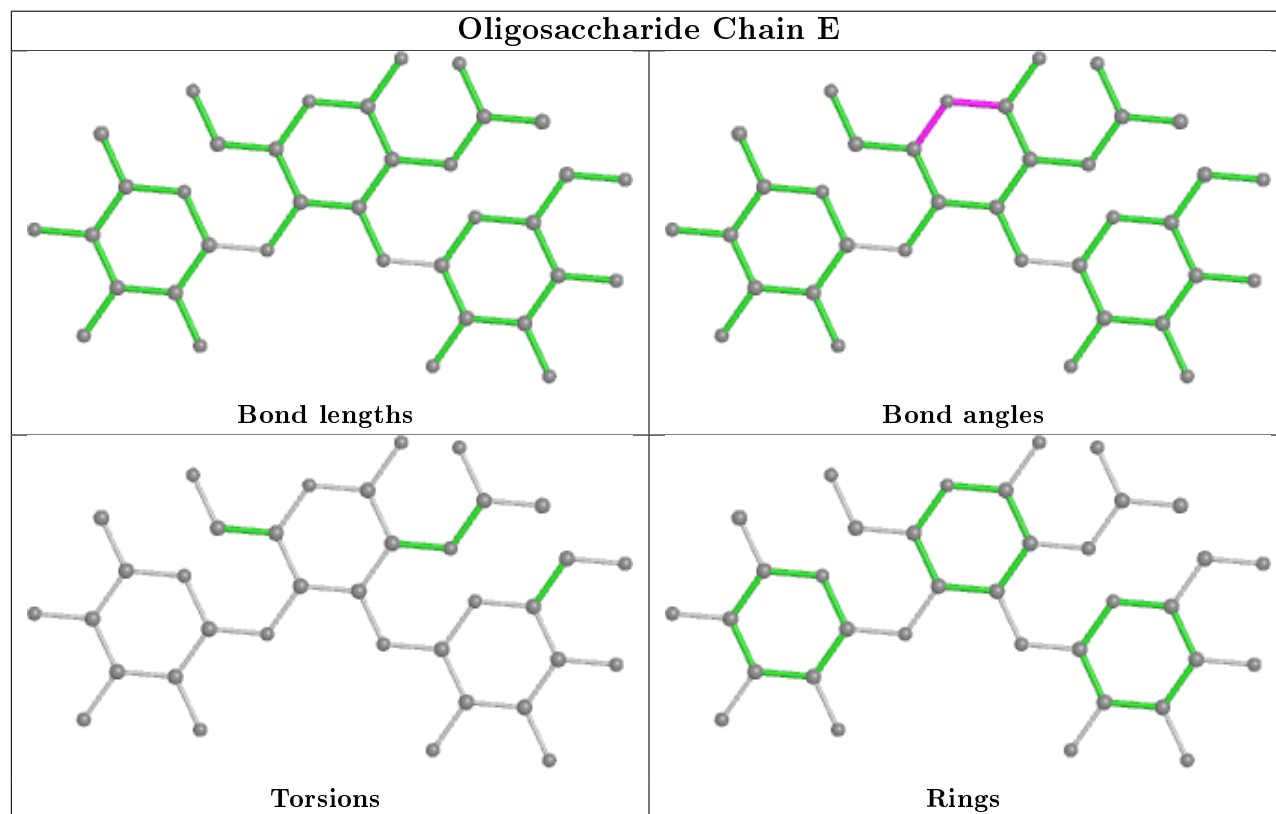
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	2	GAL	1	0
2	F	3	FUC	3	0
2	F	1	NAG	2	0
2	E	1	NAG	2	0
2	H	2	GAL	1	0
2	H	1	NAG	2	0

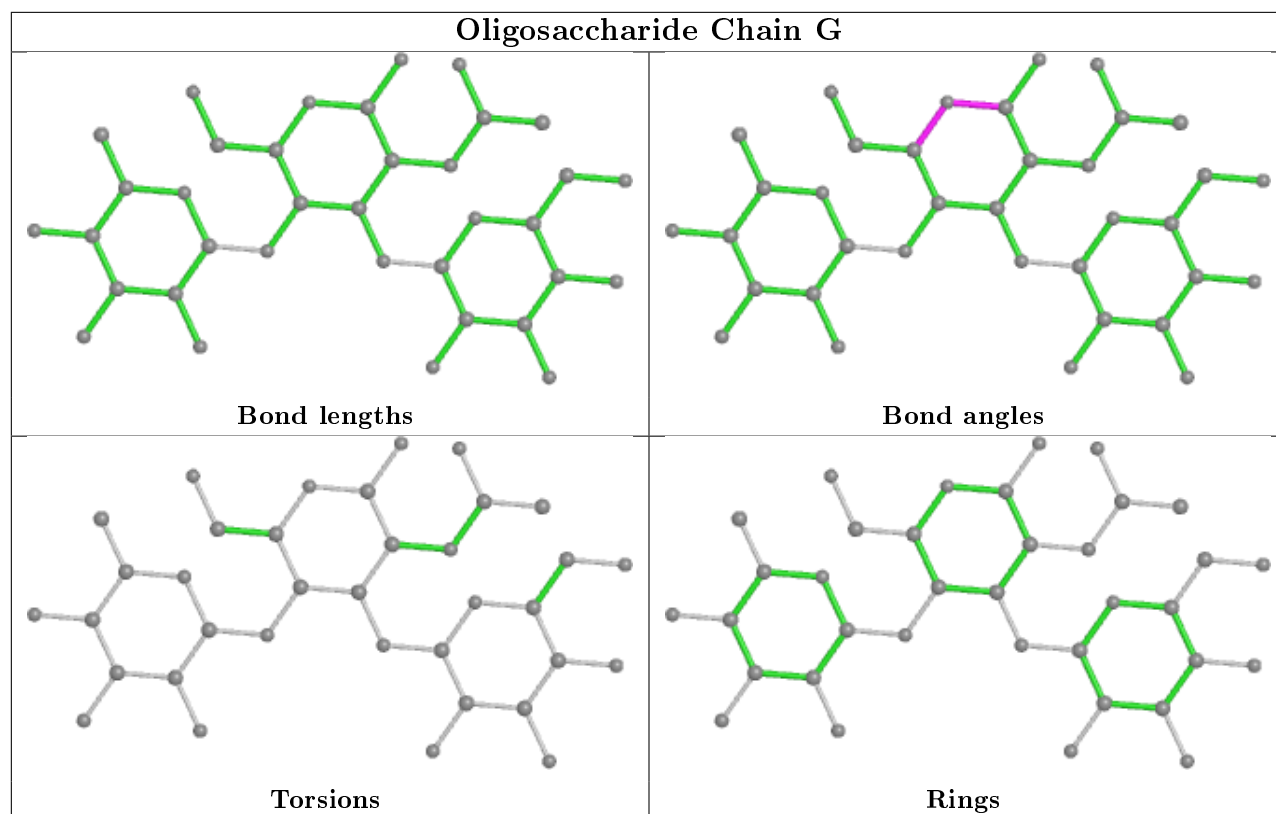
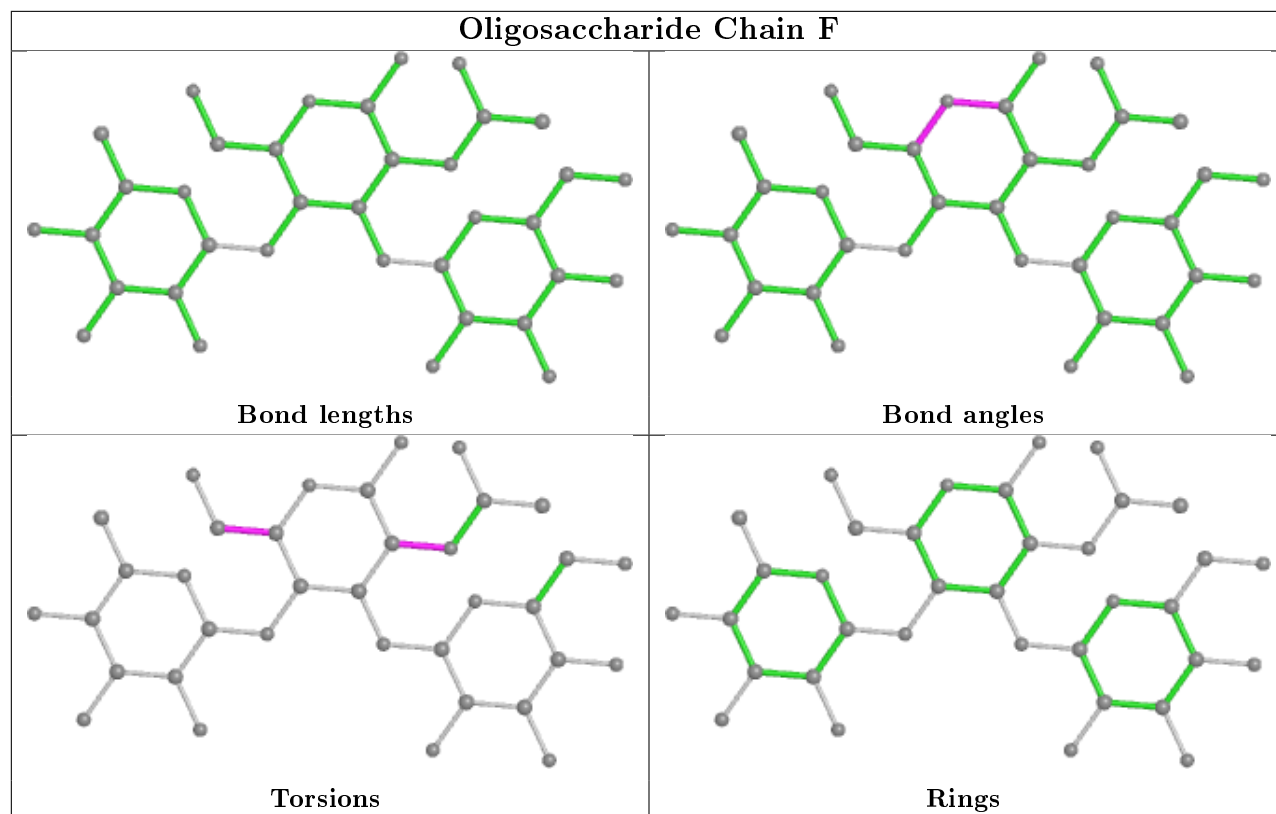
Continued on next page...

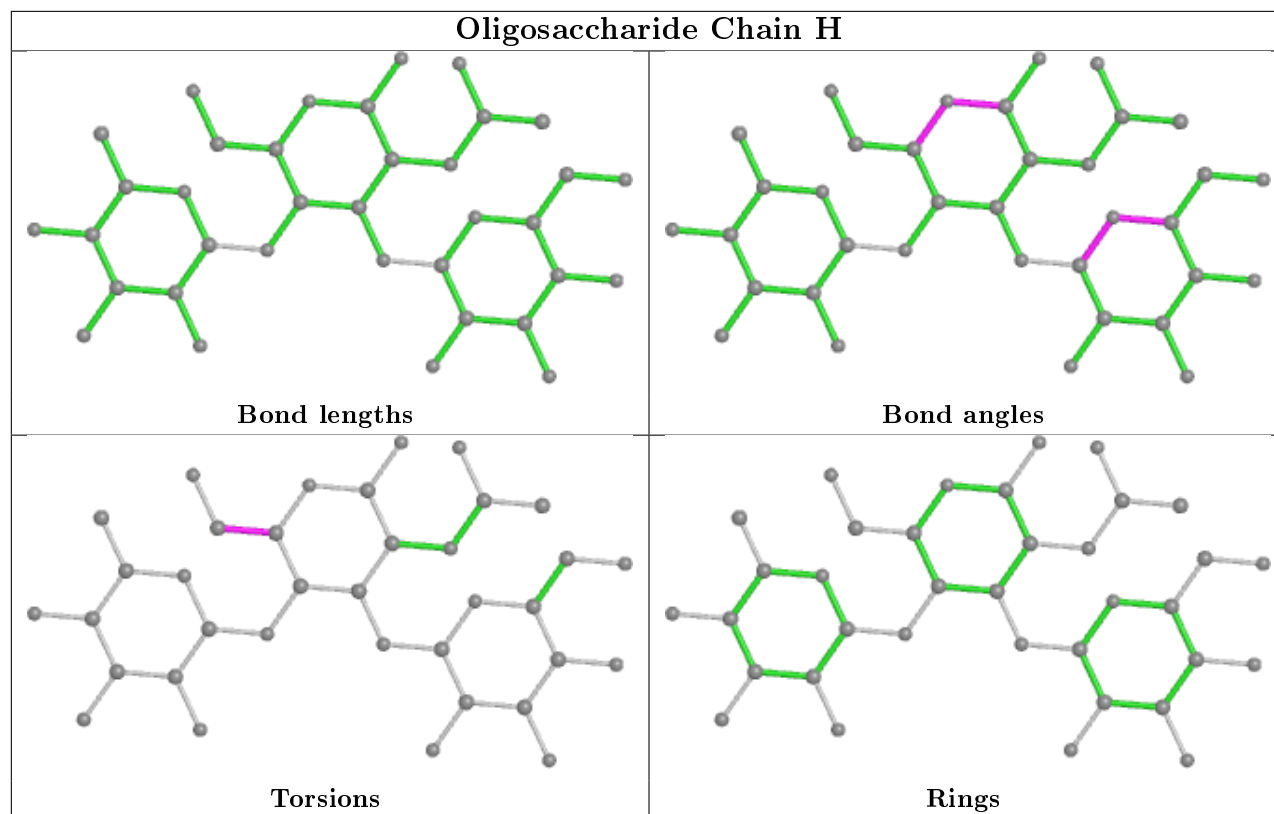
Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	1	NAG	6	0
2	H	3	FUC	2	0
2	G	3	FUC	5	0
2	E	3	FUC	2	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](#)

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 ⓘ

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	290/301 (96%)	0.02	12 (4%) 37 41	8, 16, 37, 54	0
1	B	282/301 (93%)	-0.14	9 (3%) 47 52	8, 15, 29, 66	0
1	C	282/301 (93%)	-0.16	11 (3%) 39 44	7, 14, 35, 51	0
1	D	289/301 (96%)	-0.14	10 (3%) 44 49	7, 13, 31, 49	0
All	All	1143/1204 (94%)	-0.11	42 (3%) 41 46	7, 15, 33, 66	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	407	VAL	6.9
1	D	358	ALA	6.8
1	B	440	THR	6.3
1	D	422	PHE	5.8
1	C	344	LEU	5.4
1	A	391	SER	5.2
1	C	422	PHE	5.0
1	B	412	HIS	4.9
1	A	392	GLY	4.7
1	B	422	PHE	4.6
1	A	410	SER	4.1
1	A	409	GLU	3.9
1	C	421	GLY	3.7
1	A	477	HIS	3.7
1	D	344	LEU	3.6
1	C	343	ASN	3.4
1	B	315	HIS	3.2
1	A	317	PHE	3.2
1	B	519	TYR	3.1
1	B	413	LEU	3.1
1	C	497	THR	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	406	THR	2.9
1	B	344	LEU	2.9
1	C	315	HIS	2.7
1	D	408	THR	2.6
1	C	345	SER	2.6
1	A	393	ALA	2.6
1	C	299	ALA	2.5
1	A	344	LEU	2.4
1	B	345	SER	2.4
1	A	390	THR	2.4
1	C	411	THR	2.4
1	A	343	ASN	2.3
1	D	461	PRO	2.3
1	A	358	ALA	2.3
1	A	476	THR	2.3
1	D	410	SER	2.1
1	B	394	ARG	2.1
1	D	300	SER	2.1
1	C	316	ALA	2.1
1	D	315	HIS	2.1
1	C	346	GLY	2.0

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, 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(\AA^2)	Q<0.9
2	FUC	F	3	10/11	0.47	0.31	50,75,77,78	10
2	FUC	E	3	10/11	0.50	0.35	38,45,51,52	10
2	FUC	G	3	10/11	0.55	0.32	43,50,63,64	10
2	NAG	G	1	15/15	0.64	0.26	26,40,46,47	15
2	NAG	F	1	15/15	0.68	0.27	31,48,51,51	15
2	NAG	H	1	15/15	0.77	0.25	26,29,33,42	15
2	NAG	E	1	15/15	0.81	0.24	29,40,45,46	15

Continued on next page...

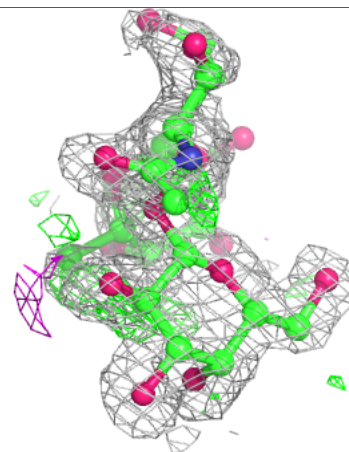
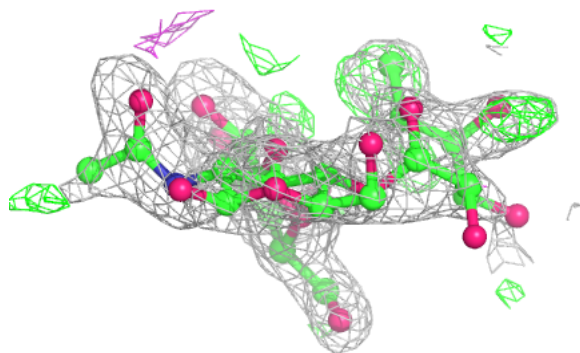
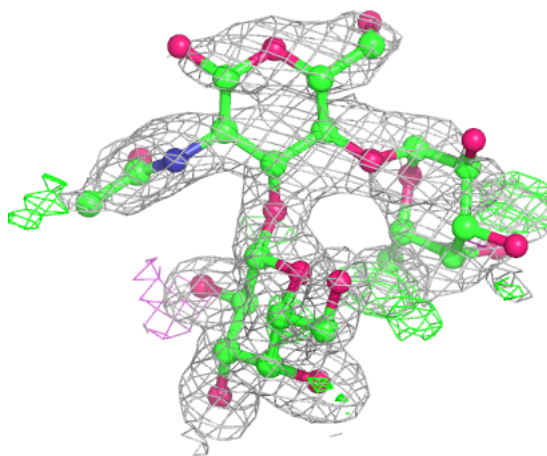
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	FUC	H	3	10/11	0.82	0.16	11,26,32,34	10
2	GAL	H	2	11/12	0.87	0.13	12,16,25,31	11
2	GAL	F	2	11/12	0.87	0.17	13,17,29,31	11
2	GAL	E	2	11/12	0.88	0.14	15,20,24,29	11
2	GAL	G	2	11/12	0.93	0.10	9,14,21,25	11

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.

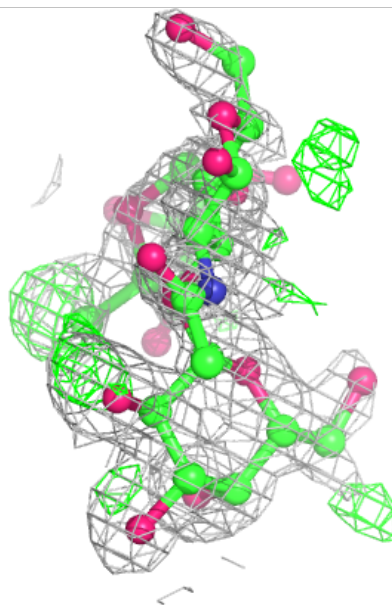
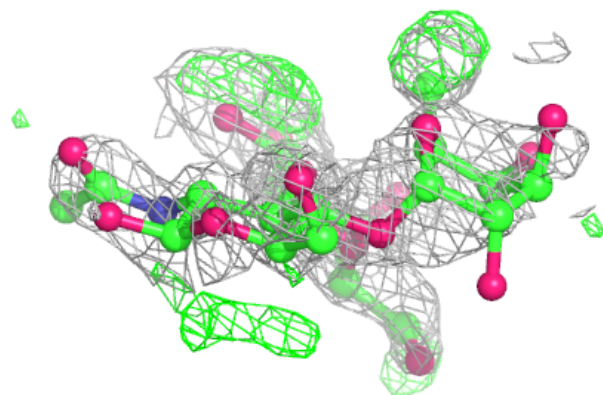
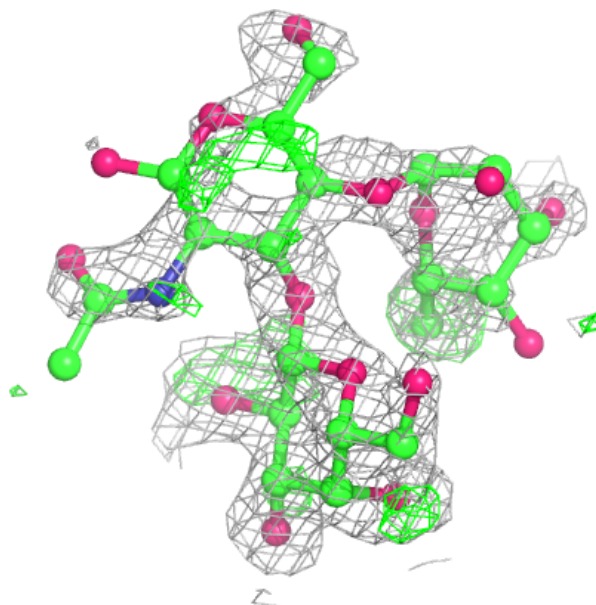
Electron density around Chain E:

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



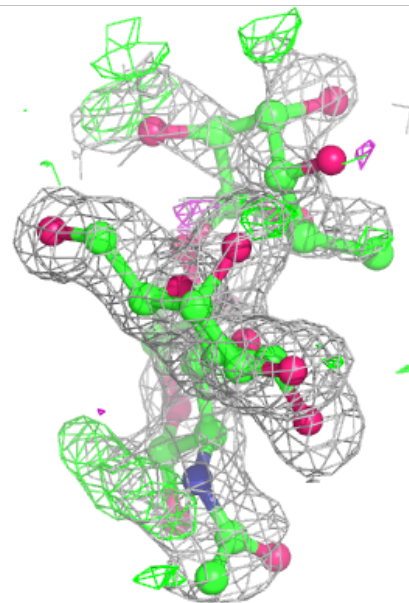
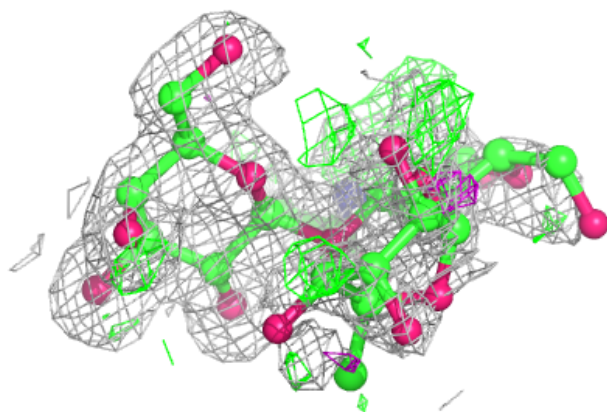
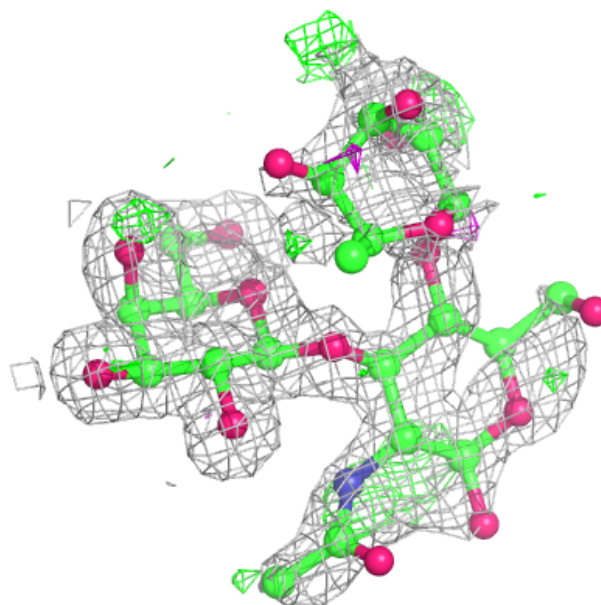
Electron density around Chain F:

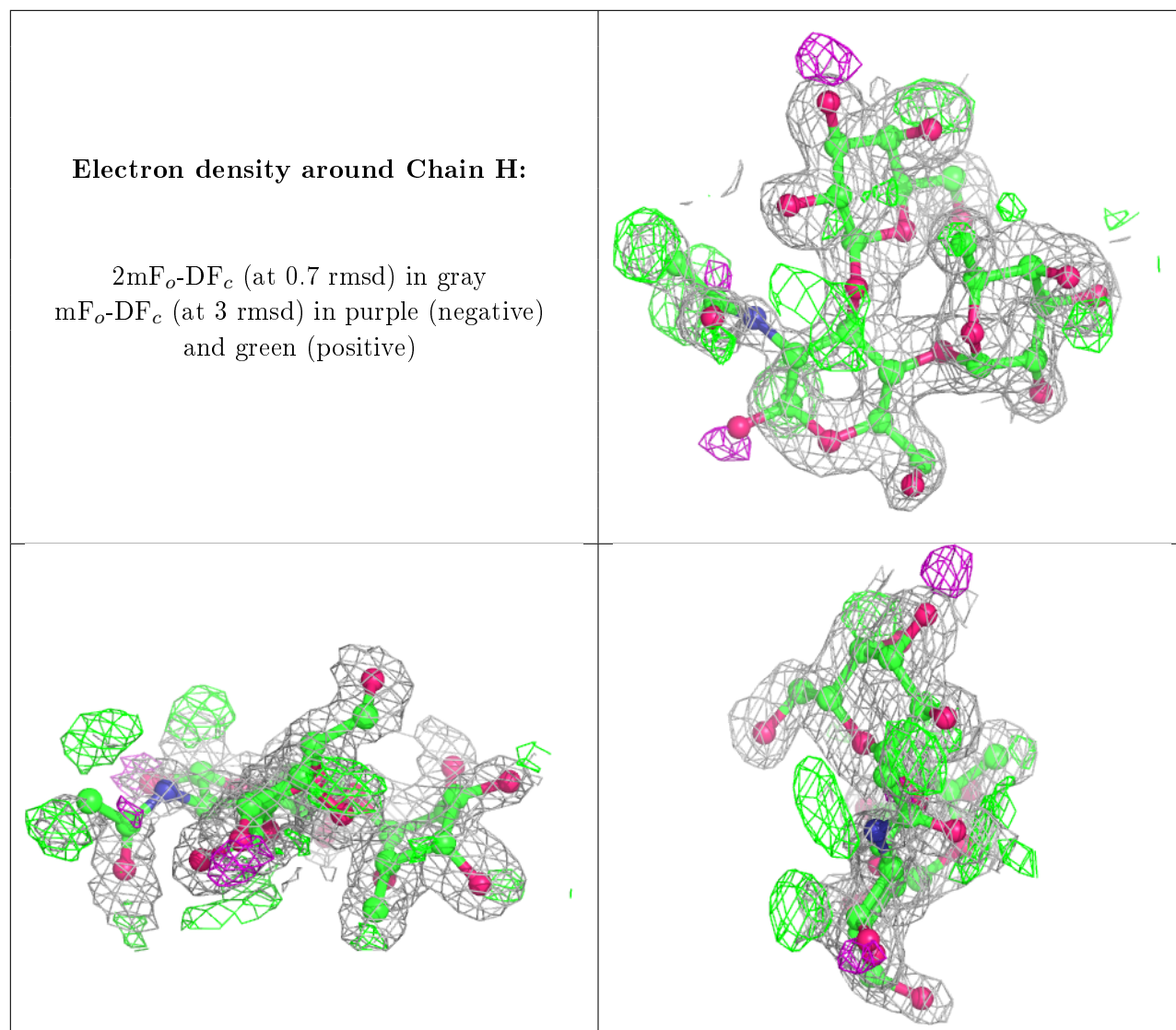
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain G:

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





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.