



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2024 – 07:10 pm BST

PDB ID : 4BGY
Title : H5 (VN1194) Influenza Virus Haemagglutinin in Complex with Avian Receptor Analogue 3'-SLN
Authors : Xiong, X.; Coombs, P.; Martin, S.R.; Liu, J.; Xiao, H.; McCauley, J.W.; Locher, K.; Walker, P.A.; Collins, P.J.; Kawaoka, Y.; Skehel, J.J.; Gamblin, S.J.
Deposited on : 2013-03-29
Resolution : 2.68 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

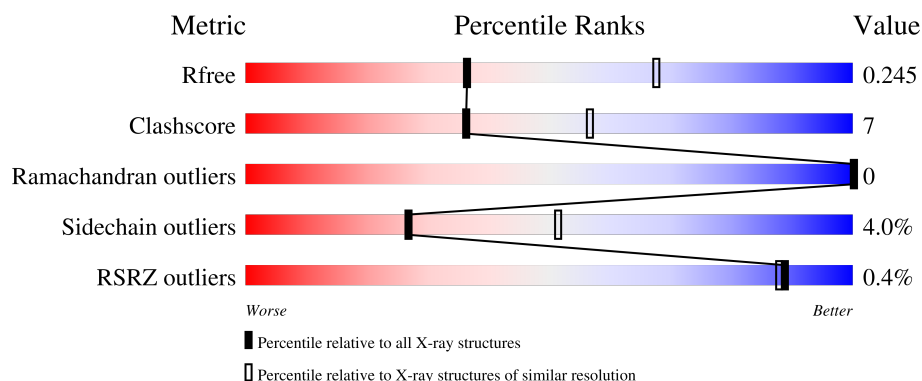
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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.




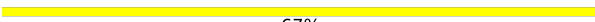

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	4708 (2.70-2.66)
Clashscore	180529	5138 (2.70-2.66)
Ramachandran outliers	177936	5071 (2.70-2.66)
Sidechain outliers	177891	5071 (2.70-2.66)
RSRZ outliers	164620	4708 (2.70-2.66)

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	326	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>..</div> </div> </div>
2	B	166	<div> <div></div> <div>85%</div> <div>12%</div> <div>..</div> </div>
3	C	2	<div> <div></div> <div>100%</div> </div>
3	D	2	<div> <div></div> <div>100%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	2	 100%
4	E	3	 67%  33%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	F	1	X	-	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4137 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 HEMAGGLUTININ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2547	1609	439	484	15			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	325	THR	ARG	conflict	UNP Q6DQ34

- Molecule 2 is a protein called HEMAGGLUTININ.

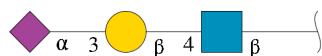
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	162	Total	C	N	O	S	0	0	0
			1312	819	226	259	8			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



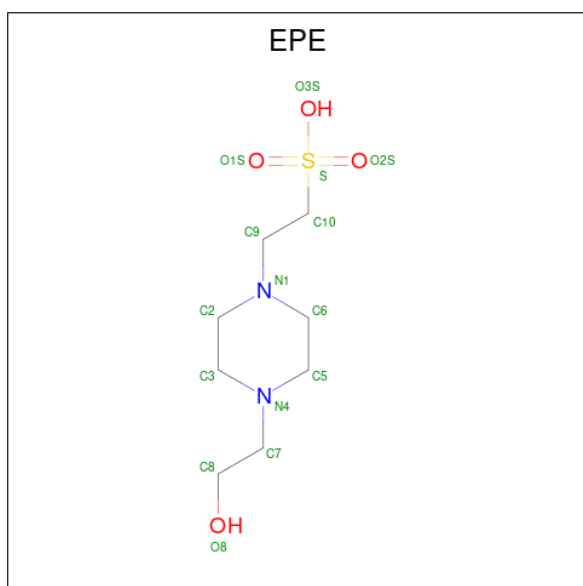
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

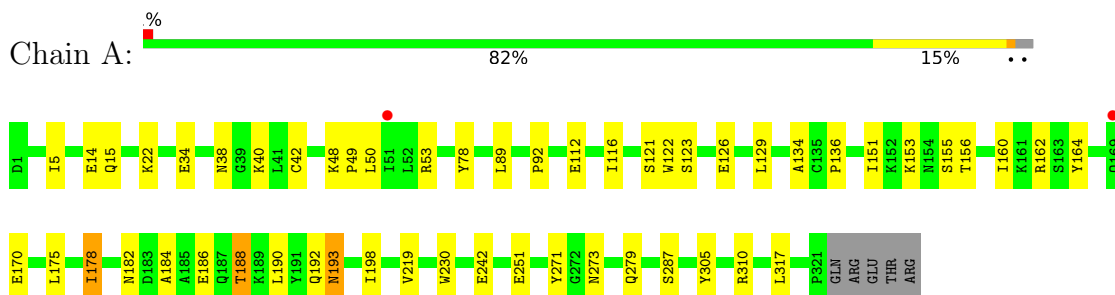
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	73	Total	O	0	0
			73	73		
6	B	60	Total	O	0	0
			60	60		

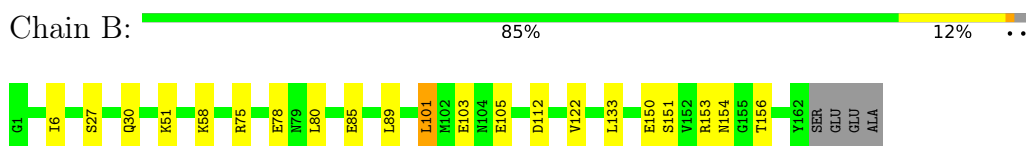
3 Residue-property plots

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.

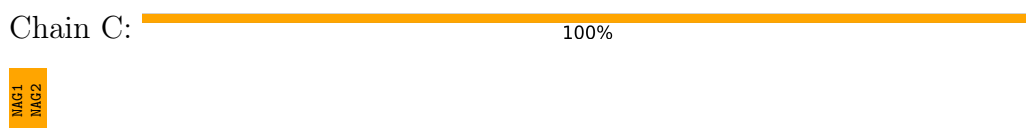
- Molecule 1: HEMAGGLUTININ



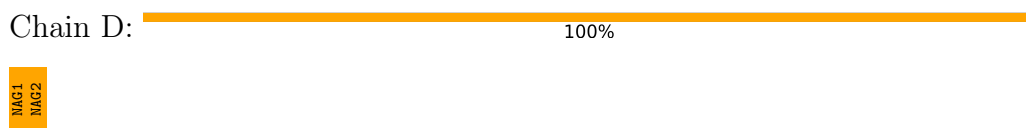
- Molecule 2: HEMAGGLUTININ



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1
MAG2

- Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E: 

MAG1
GAL2
STA3

4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	101.40Å 101.40Å 450.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.94 – 2.68 40.94 – 2.68	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.94-2.68) 99.9 (40.94-2.68)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.202 , 0.248 0.204 , 0.245	Depositor DCC
R_{free} test set	1307 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	52.6	Xtriage
Anisotropy	0.712	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4137	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% 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: EPE, GAL, NAG, SIA

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	0.35	0/2609	0.56	0/3545
2	B	0.40	0/1339	0.57	0/1802
All	All	0.37	0/3948	0.56	0/5347

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

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	2547	0	2489	39	1
2	B	1312	0	1218	13	0
3	C	28	0	25	2	0
3	D	28	0	25	1	0
3	F	28	0	25	1	0
4	E	46	0	40	3	0
5	B	15	0	18	1	0
6	A	73	0	0	1	0
6	B	60	0	0	2	0
All	All	4137	0	3840	56	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 7.

All (56) 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:B:150:GLU:OE1	2:B:153:ARG:NH1	2.08	0.86
1:A:126:GLU:HG2	1:A:129:LEU:CD1	2.11	0.81
3:C:1:NAG:O4	3:C:2:NAG:H83	1.87	0.75
1:A:129:LEU:HD12	1:A:151:ILE:HD11	1.70	0.71
1:A:126:GLU:HG2	1:A:129:LEU:HD12	1.71	0.70
1:A:193:ASN:H	1:A:193:ASN:HD22	1.42	0.66
1:A:122:TRP:CH2	1:A:162:ARG:HG3	2.32	0.64
2:B:156:THR:O	2:B:156:THR:HG22	1.98	0.63
1:A:121:SER:HB3	1:A:162:ARG:HH22	1.65	0.62
1:A:193:ASN:H	1:A:193:ASN:ND2	1.98	0.62
1:A:126:GLU:CG	1:A:129:LEU:HG	2.31	0.60
2:B:51:LYS:HE3	2:B:103:GLU:OE1	2.01	0.60
1:A:190:LEU:HD21	4:E:3:SIA:O10	2.02	0.58
1:A:34:GLU:OE1	1:A:287:SER:HB3	2.03	0.58
1:A:15:GLN:O	1:A:310:ARG:NH2	2.37	0.57
2:B:80:LEU:C	2:B:80:LEU:HD23	2.26	0.56
1:A:5:ILE:HD11	2:B:122:VAL:HG21	1.88	0.54
1:A:182:ASN:ND2	1:A:186:GLU:OE1	2.40	0.54
4:E:3:SIA:H6	4:E:3:SIA:O1B	2.07	0.53
1:A:48:LYS:HG2	1:A:49:PRO:HD2	1.91	0.53
1:A:14:GLU:HG2	6:A:2016:HOH:O	2.09	0.52
1:A:116:ILE:HG22	1:A:164:TYR:CE1	2.45	0.51
3:D:1:NAG:O4	3:D:2:NAG:O7	2.29	0.51
1:A:134:ALA:O	1:A:136:PRO:HD3	2.10	0.51
2:B:75:ARG:NH1	2:B:78:GLU:OE1	2.41	0.50
4:E:3:SIA:O10	4:E:3:SIA:O7	2.17	0.50
2:B:156:THR:O	2:B:156:THR:CG2	2.60	0.49
1:A:122:TRP:CZ3	1:A:162:ARG:NH1	2.80	0.49
2:B:80:LEU:HD23	2:B:80:LEU:O	2.12	0.48
1:A:121:SER:HB3	1:A:162:ARG:NH2	2.29	0.47
1:A:42:CYS:HB2	1:A:279:GLN:OE1	2.14	0.47
1:A:178:ILE:HD12	1:A:198:ILE:CD1	2.44	0.47
1:A:92:PRO:HB3	1:A:219:VAL:CG1	2.44	0.47
1:A:184:ALA:O	1:A:188:THR:CG2	2.63	0.46
1:A:160:ILE:O	1:A:242:GLU:HA	2.14	0.46
3:F:1:NAG:O4	3:F:2:NAG:O7	2.33	0.46
1:A:134:ALA:C	1:A:136:PRO:HD3	2.37	0.45
1:A:271:TYR:CE2	1:A:273:ASN:HA	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:LEU:HD23	1:A:230:TRP:HB3	1.97	0.45
1:A:116:ILE:HG22	1:A:164:TYR:CZ	2.52	0.44
1:A:126:GLU:HG2	1:A:129:LEU:CG	2.47	0.43
2:B:101:LEU:HD13	6:B:2035:HOH:O	2.18	0.43
1:A:122:TRP:HH2	1:A:162:ARG:HG3	1.79	0.43
1:A:22:LYS:HD3	3:C:1:NAG:H82	2.00	0.43
1:A:305:TYR:HD2	2:B:89:LEU:HD22	1.82	0.43
1:A:123:SER:O	1:A:153:LYS:NZ	2.41	0.43
2:B:151:SER:HA	2:B:154:ASN:HB2	2.01	0.42
1:A:34:GLU:OE1	1:A:287:SER:CB	2.67	0.42
1:A:184:ALA:O	1:A:188:THR:HG23	2.20	0.42
1:A:178:ILE:HD12	1:A:198:ILE:HD12	2.02	0.41
2:B:105:GLU:HG3	6:B:2036:HOH:O	2.19	0.41
1:A:38:ASN:OD1	1:A:40:LYS:N	2.47	0.41
5:B:1165:EPE:H62	5:B:1165:EPE:O2S	2.20	0.41
1:A:50:LEU:HD12	1:A:50:LEU:C	2.40	0.41
1:A:49:PRO:HB3	1:A:78:TYR:CZ	2.55	0.41
2:B:6:ILE:HD12	2:B:112:ASP:HA	2.02	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:GLU:OE2	1:A:170:GLU:OE2[11_565]	2.07	0.13

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	319/326 (98%)	304 (95%)	15 (5%)	0	100	100
2	B	160/166 (96%)	153 (96%)	7 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	479/492 (97%)	457 (95%)	22 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	289/294 (98%)	278 (96%)	11 (4%)	28	53
2	B	138/141 (98%)	132 (96%)	6 (4%)	25	49
All	All	427/435 (98%)	410 (96%)	17 (4%)	27	51

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	ARG
1	A	89	LEU
1	A	112	GLU
1	A	155	SER
1	A	156	THR
1	A	178	ILE
1	A	188	THR
1	A	192	GLN
1	A	193	ASN
1	A	251	GLU
1	A	317	LEU
2	B	27	SER
2	B	30	GLN
2	B	58	LYS
2	B	85	GLU
2	B	101	LEU
2	B	133	LEU

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

Mol	Chain	Res	Type
1	A	15	GLN
1	A	193	ASN
2	B	79	ASN

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 ⓘ

9 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	C	1	3,1	14,14,15	0.48	0	17,19,21	1.19	1 (5%)
3	NAG	C	2	3	14,14,15	0.63	0	17,19,21	1.55	4 (23%)
3	NAG	D	1	3,1	14,14,15	0.51	0	17,19,21	1.60	3 (17%)
3	NAG	D	2	3	14,14,15	0.67	0	17,19,21	1.41	2 (11%)
4	NAG	E	1	4	15,15,15	0.41	0	21,21,21	1.42	1 (4%)
4	GAL	E	2	4	11,11,12	0.28	0	15,15,17	0.89	1 (6%)
4	SIA	E	3	4	20,20,21	0.78	0	24,28,31	1.22	3 (12%)
3	NAG	F	1	2,3	14,14,15	0.66	0	17,19,21	1.26	2 (11%)
3	NAG	F	2	3	14,14,15	0.55	0	17,19,21	1.24	1 (5%)

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	NAG	C	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	C	2	3	-	4/6/23/26	0/1/1/1
3	NAG	D	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	3/6/23/26	0/1/1/1
4	NAG	E	1	4	-	0/6/26/26	0/1/1/1
4	GAL	E	2	4	-	2/2/19/22	0/1/1/1
4	SIA	E	3	4	-	1/18/34/38	0/1/1/1
3	NAG	F	1	2,3	1/1/5/7	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	1	NAG	C1-C2-N2	-5.30	104.58	110.73
3	D	1	NAG	C1-O5-C5	4.63	118.46	112.19
3	C	2	NAG	O5-C1-C2	-3.79	105.30	111.29
3	C	1	NAG	O5-C5-C6	3.71	113.02	107.20
3	F	2	NAG	C1-C2-N2	-3.55	104.42	110.49
3	D	2	NAG	C1-C2-N2	-3.54	104.45	110.49
3	C	2	NAG	C8-C7-N2	3.37	121.80	116.10
4	E	3	SIA	C4-C5-N5	-3.15	104.14	110.38
3	F	1	NAG	O5-C5-C4	-2.90	103.78	110.83
4	E	3	SIA	C3-C4-C5	-2.44	108.51	111.46
3	D	2	NAG	C4-C3-C2	2.34	114.45	111.02
3	D	1	NAG	O7-C7-N2	2.29	126.16	121.95
3	C	2	NAG	C3-C4-C5	2.17	114.11	110.24
3	C	2	NAG	O7-C7-C8	-2.16	118.05	122.06
3	F	1	NAG	O5-C5-C6	2.14	110.55	107.20
4	E	3	SIA	C6-C5-N5	-2.13	107.38	110.91
3	D	1	NAG	O7-C7-C8	-2.05	118.26	122.06
4	E	2	GAL	O3-C3-C2	-2.02	106.12	109.99

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	F	1	NAG	C1

All (16) torsion outliers are listed below:

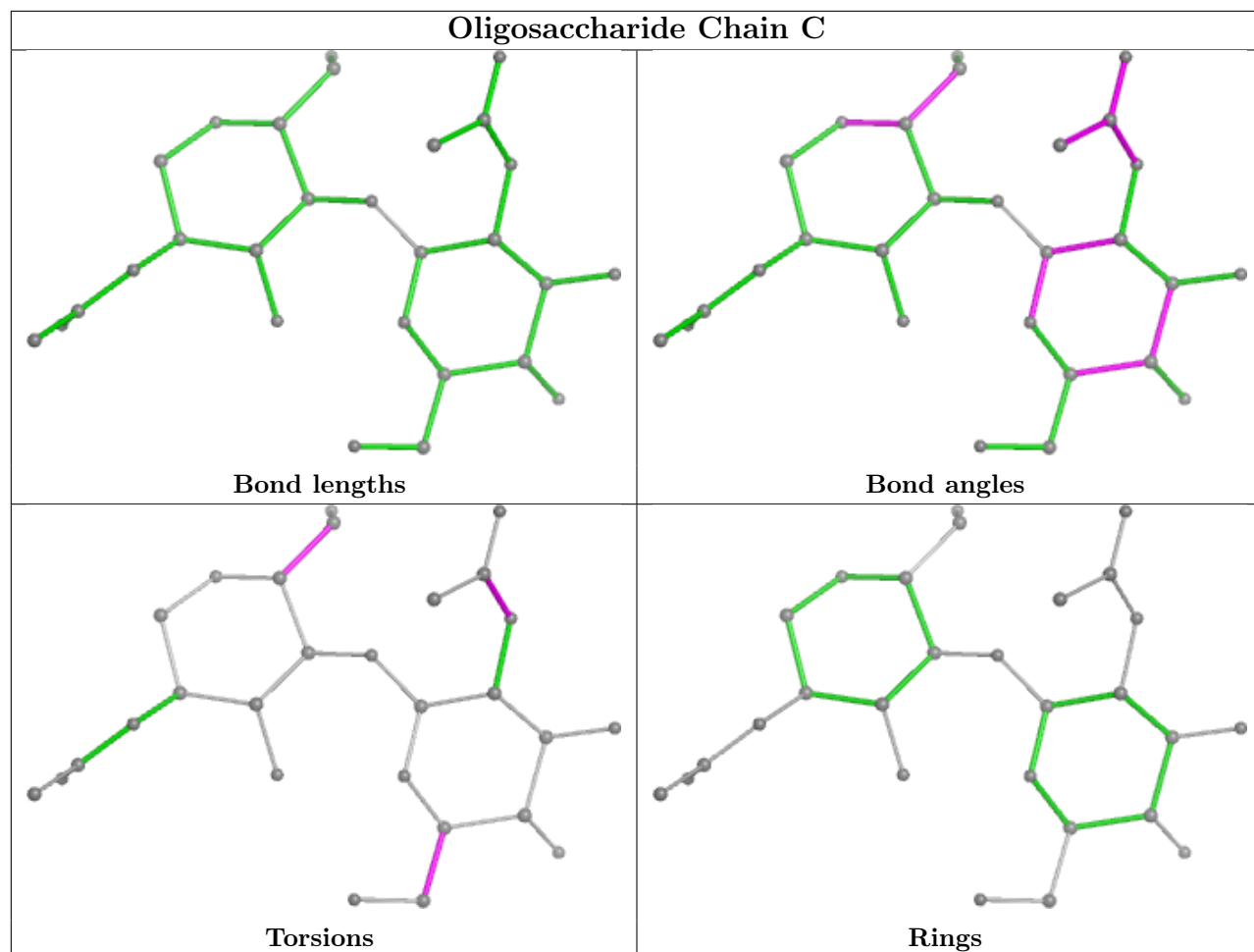
Mol	Chain	Res	Type	Atoms
3	D	1	NAG	O5-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
4	E	2	GAL	O5-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
4	E	2	GAL	C4-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6
3	F	1	NAG	C4-C5-C6-O6
3	C	2	NAG	C8-C7-N2-C2
3	C	2	NAG	O7-C7-N2-C2
3	C	1	NAG	O5-C5-C6-O6
3	C	1	NAG	C4-C5-C6-O6
3	C	2	NAG	C4-C5-C6-O6
3	D	2	NAG	C3-C2-N2-C7
4	E	3	SIA	C6-C5-N5-C10

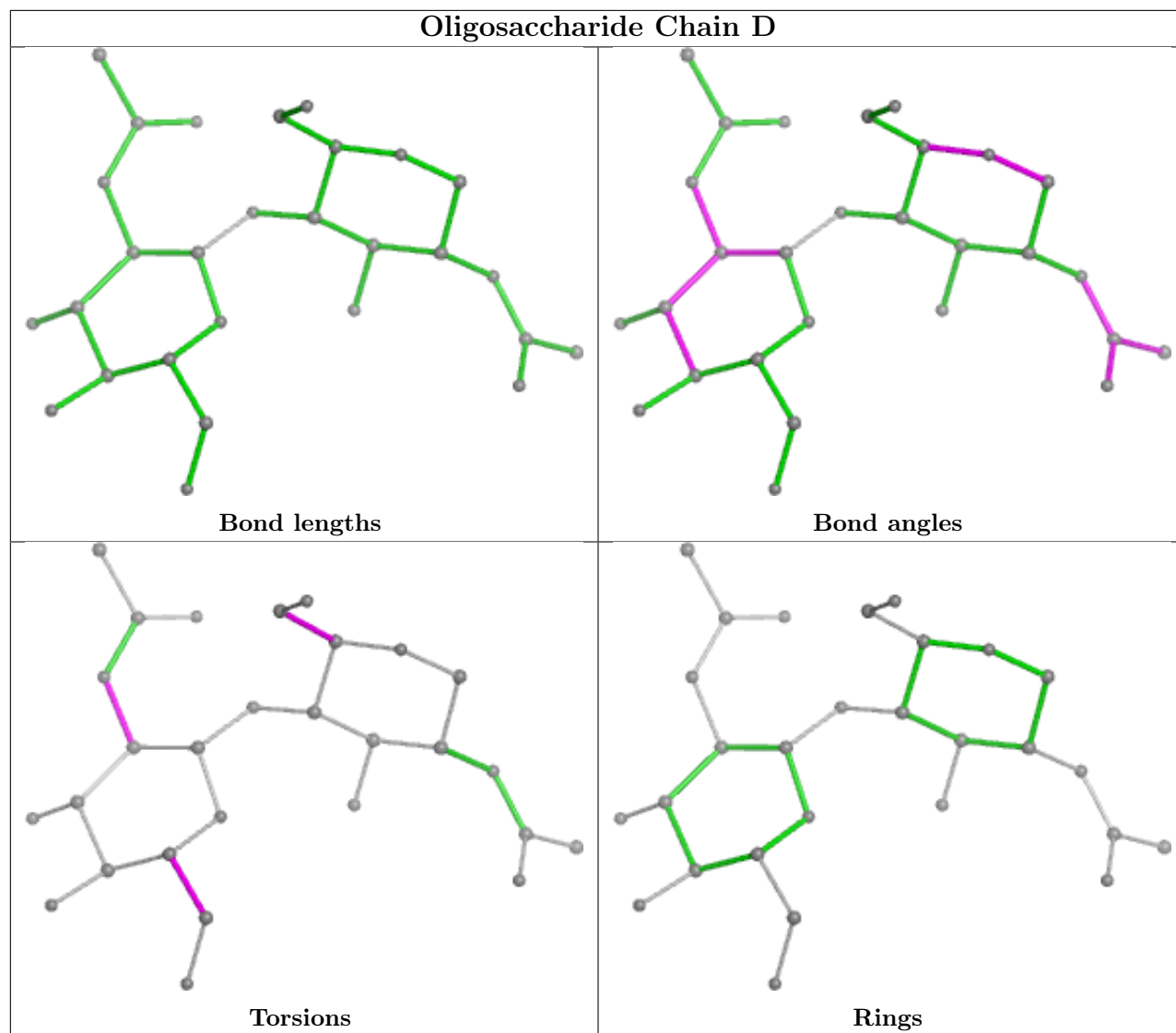
There are no ring outliers.

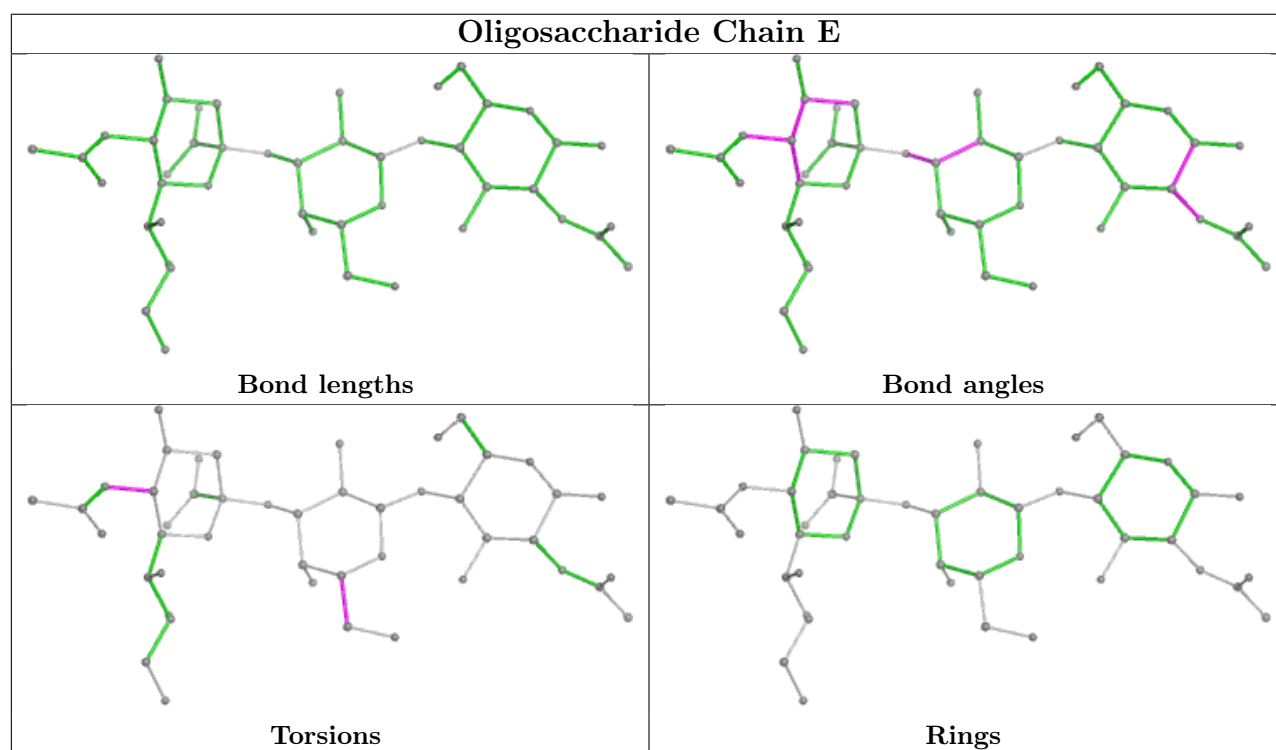
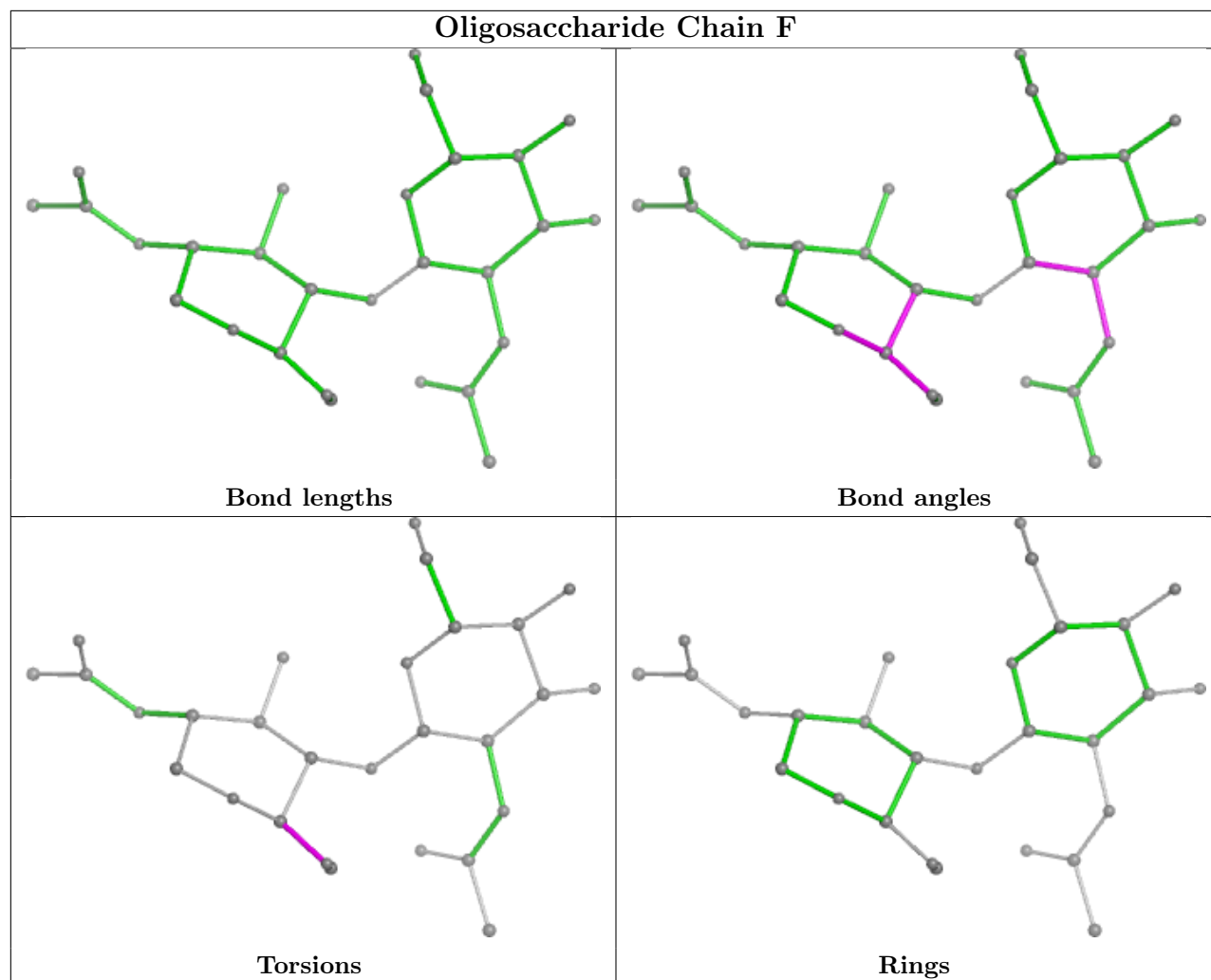
7 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	2	NAG	1	0
3	D	2	NAG	1	0
3	D	1	NAG	1	0
3	F	1	NAG	1	0
3	C	2	NAG	1	0
4	E	3	SIA	3	0
3	C	1	NAG	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

1 ligand is 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
5	EPE	B	1165	-	15,15,15	2.03	1 (6%)	18,20,20	2.86	9 (50%)

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
5	EPE	B	1165	-	-	3/9/19/19	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1165	EPE	C10-S	-7.40	1.67	1.77

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1165	EPE	C6-N1-C2	5.10	120.30	108.83
5	B	1165	EPE	C5-N4-C3	4.54	119.06	108.83
5	B	1165	EPE	O2S-S-C10	4.33	112.13	106.92
5	B	1165	EPE	C9-N1-C6	3.66	120.60	111.23
5	B	1165	EPE	C7-N4-C5	3.62	120.48	111.23
5	B	1165	EPE	C7-N4-C3	3.60	120.45	111.23
5	B	1165	EPE	C3-C2-N1	-3.51	103.44	110.64
5	B	1165	EPE	C9-N1-C2	3.07	119.09	111.23
5	B	1165	EPE	O1S-S-C10	2.68	110.15	106.92

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1165	EPE	C10-C9-N1-C6
5	B	1165	EPE	N4-C7-C8-O8
5	B	1165	EPE	C8-C7-N4-C5

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1165	EPE	1	0

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, 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	321/326 (98%)	-0.02	2 (0%) 85 84	44, 83, 107, 133	0
2	B	162/166 (97%)	-0.25	0 100 100	36, 60, 104, 134	0
All	All	483/492 (98%)	-0.10	2 (0%) 89 88	36, 77, 107, 134	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	51	ILE	2.0
1	A	169	GLN	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.

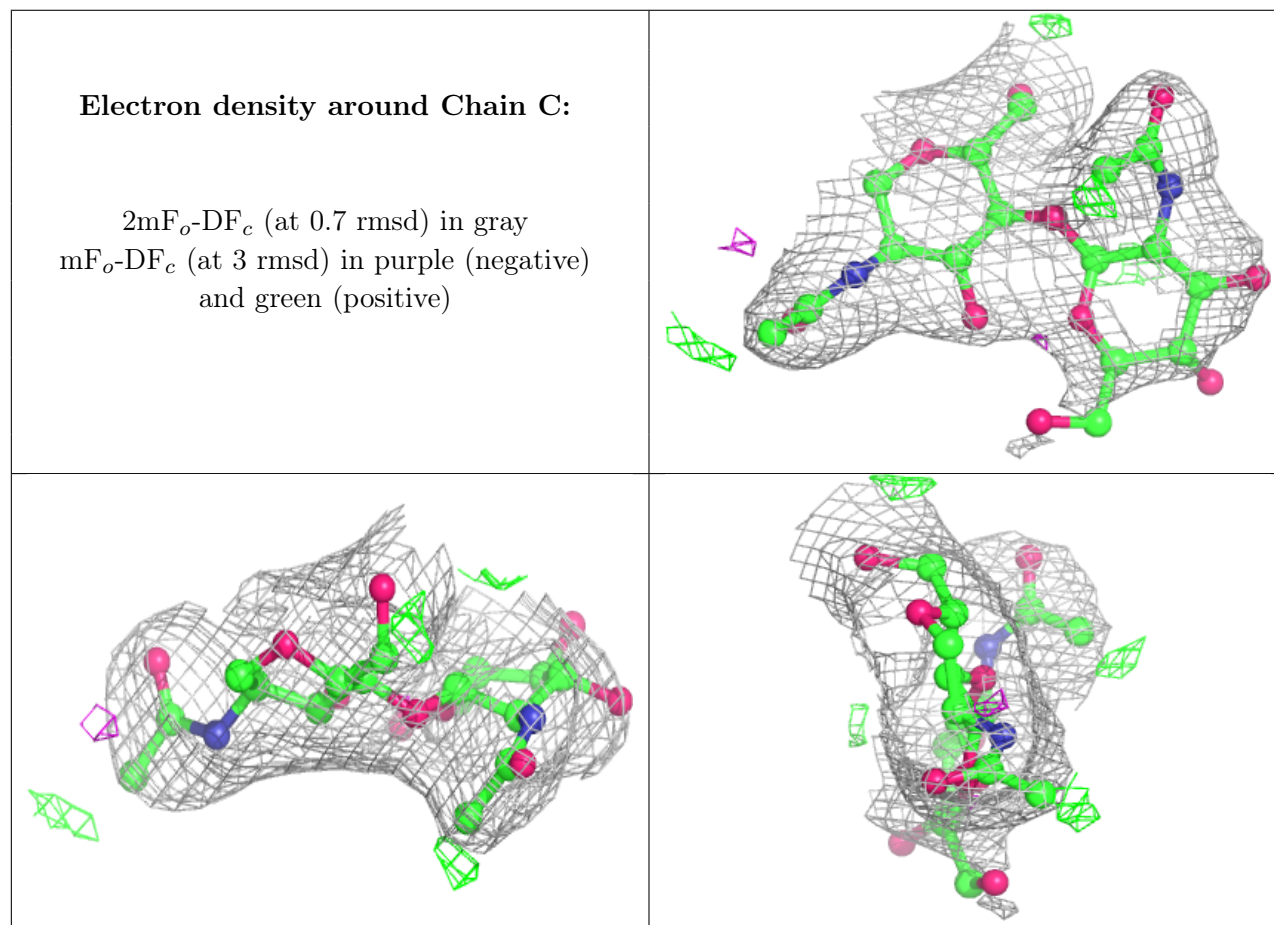
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	C	2	14/15	0.52	0.13	100,121,127,127	0
4	NAG	E	1	15/15	0.56	0.14	112,118,121,121	0
3	NAG	F	2	14/15	0.63	0.15	108,122,130,132	0
3	NAG	D	2	14/15	0.64	0.16	112,118,123,126	0
4	GAL	E	2	11/12	0.81	0.11	89,97,106,108	0
3	NAG	C	1	14/15	0.83	0.09	75,85,93,109	0
3	NAG	D	1	14/15	0.87	0.10	76,85,94,101	0

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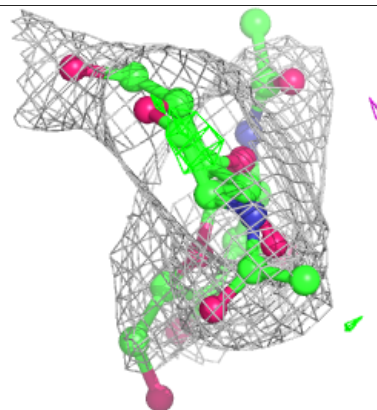
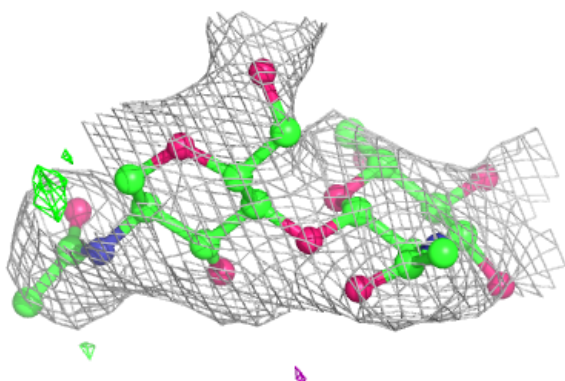
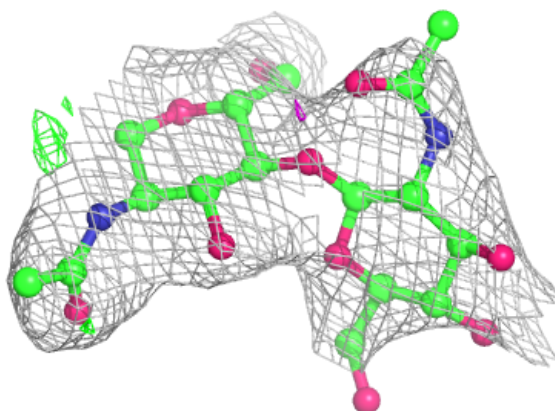
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	F	1	14/15	0.88	0.10	83,94,109,119	0
4	SIA	E	3	20/21	0.91	0.09	70,80,87,89	0

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.

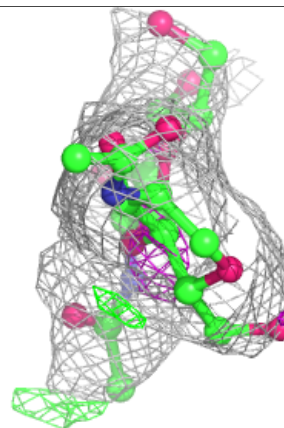
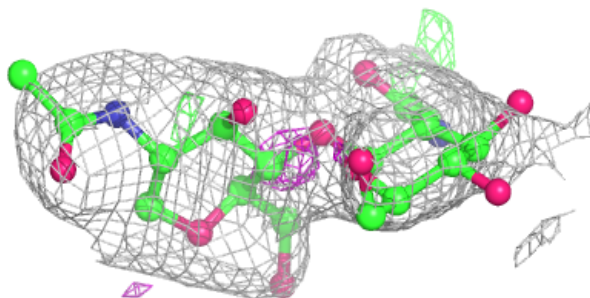
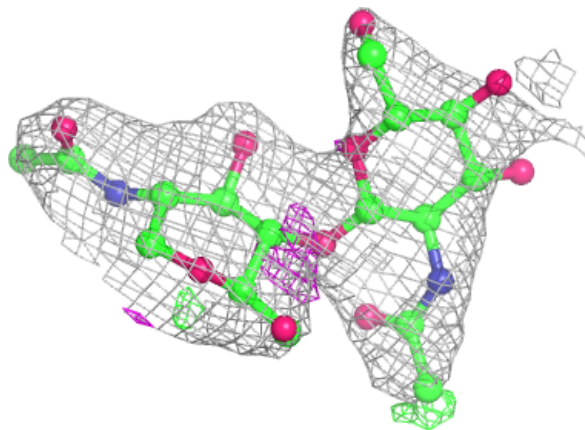


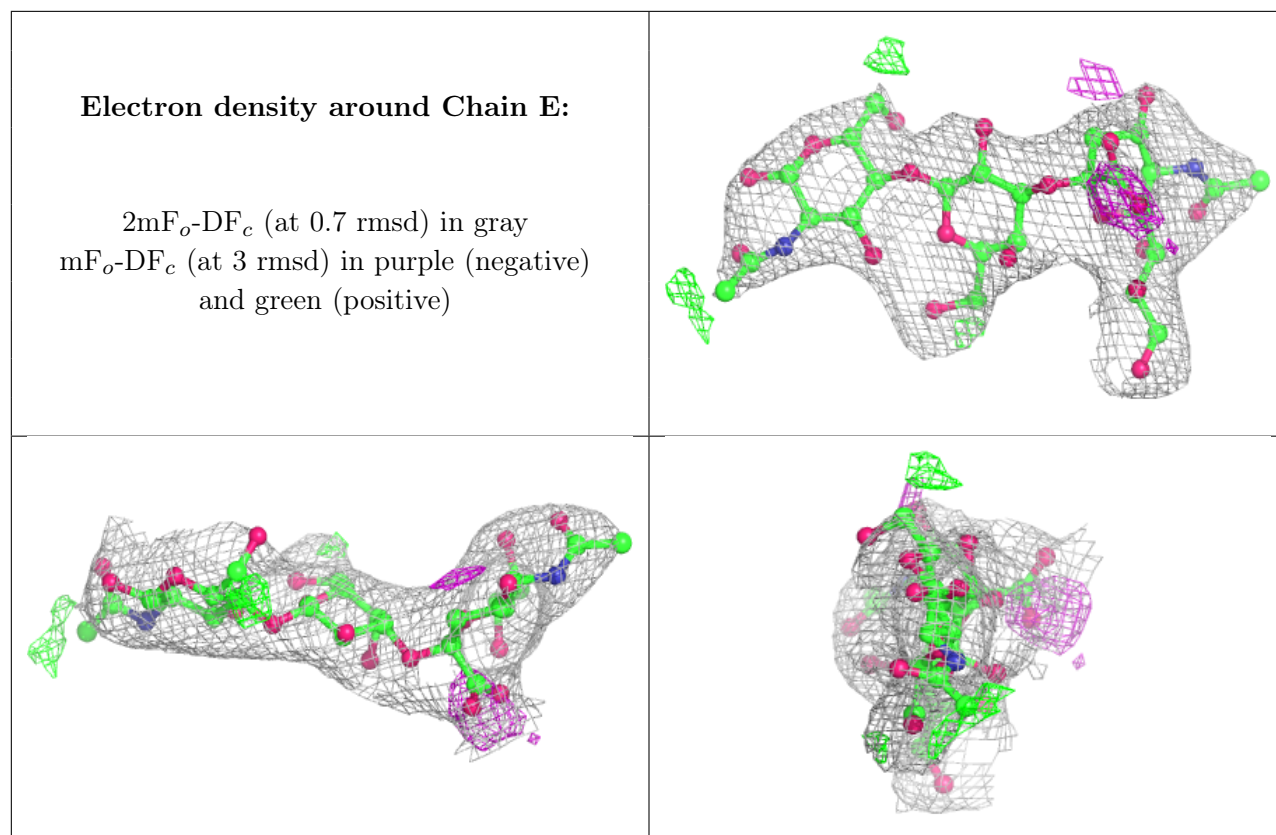
Electron density around Chain D:

$2mF_o-DF_c$ (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 [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
5	EPE	B	1165	15/15	0.86	0.16	79,82,114,117	0

6.5 Other polymers [i](#)

There are no such residues in this entry.