



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 05:22 PM UTC

PDB ID : 4MJ0 / pdb_00004mj0
Title : BK Polyomavirus VP1 pentamer in complex with GD3 oligosaccharide
Authors : Neu, U.; Stroh, L.J.; Stehle, T.
Deposited on : 2013-09-03
Resolution : 1.70 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

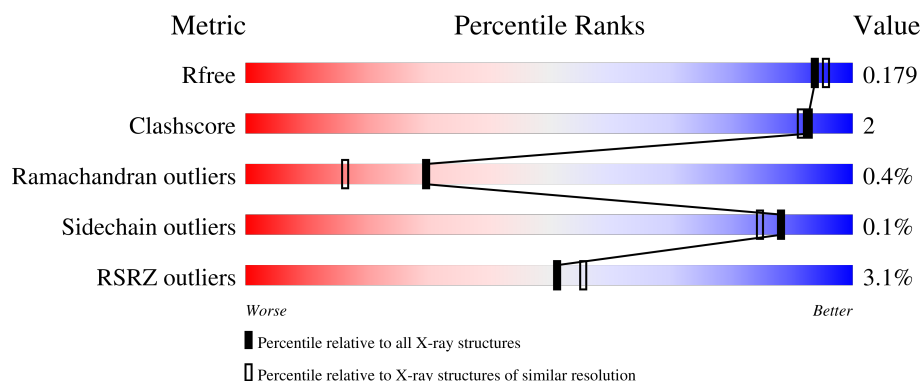
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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.



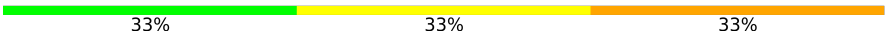
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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 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	275	% 94% . .
1	B	275	4% 93% . .
1	C	275	4% 86% . 10%
1	D	275	% 92% 5% .
1	E	275	3% 91% . .

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Mol	Chain	Length	Quality of chain
2	F	4	 25% 75%
3	G	2	 100%
3	I	2	 100%
4	H	3	 33% 33% 33%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11705 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 VP1 capsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	8	0
			2062	1300	349	399	14			
1	B	263	Total	C	N	O	S	0	7	0
			2035	1279	348	394	14			
1	C	247	Total	C	N	O	S	0	5	0
			1916	1208	329	366	13			
1	D	267	Total	C	N	O	S	0	6	0
			2073	1305	356	399	13			
1	E	264	Total	C	N	O	S	0	12	0
			2095	1320	360	401	14			

There are 20 discrepancies between the modelled and reference sequences:

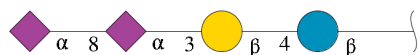
Chain	Residue	Modelled	Actual	Comment	Reference
A	26	GLY	-	expression tag	UNP Q85235
A	27	SER	-	expression tag	UNP Q85235
A	28	HIS	-	expression tag	UNP Q85235
A	29	MET	-	expression tag	UNP Q85235
B	26	GLY	-	expression tag	UNP Q85235
B	27	SER	-	expression tag	UNP Q85235
B	28	HIS	-	expression tag	UNP Q85235
B	29	MET	-	expression tag	UNP Q85235
C	26	GLY	-	expression tag	UNP Q85235
C	27	SER	-	expression tag	UNP Q85235
C	28	HIS	-	expression tag	UNP Q85235
C	29	MET	-	expression tag	UNP Q85235
D	26	GLY	-	expression tag	UNP Q85235
D	27	SER	-	expression tag	UNP Q85235
D	28	HIS	-	expression tag	UNP Q85235
D	29	MET	-	expression tag	UNP Q85235
E	26	GLY	-	expression tag	UNP Q85235
E	27	SER	-	expression tag	UNP Q85235
E	28	HIS	-	expression tag	UNP Q85235

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Chain	Residue	Modelled	Actual	Comment	Reference
E	29	MET	-	expression tag	UNP Q85235

- Molecule 2 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-8)-N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose.



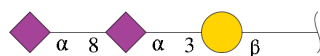
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	F	4	Total	C	N	O	0	0	0
			63	34	2	27			

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-8)-N-acetyl-alpha-neuraminic acid.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	2	Total	C	N	O	0	0	0
			41	22	2	17			
3	I	2	Total	C	N	O	0	0	0
			41	22	2	17			

- Molecule 4 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-8)-N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	3	Total	C	N	O	0	0	0
			52	28	2	22			

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cl	0	0
			1	1		
6	B	1	Total	Cl	0	0
			1	1		
6	C	1	Total	Cl	0	0
			1	1		
6	D	1	Total	Cl	0	0
			1	1		
6	E	1	Total	Cl	0	0
			1	1		

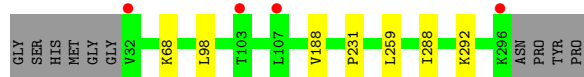
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	300	Total 300	O 300	0	0
7	B	245	Total 245	O 245	0	0
7	C	192	Total 192	O 192	0	0
7	D	266	Total 266	O 266	0	0
7	E	277	Total 277	O 277	0	0

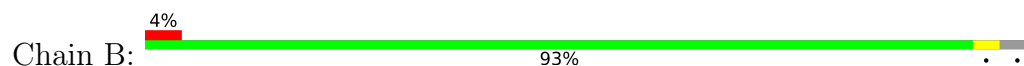
3 Residue-property plots

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.

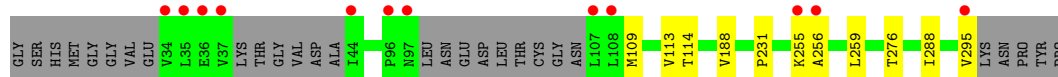
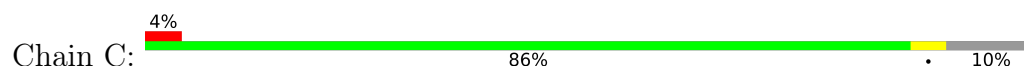
- Molecule 1: VP1 capsid protein



- Molecule 1: VP1 capsid protein



- Molecule 1: VP1 capsid protein



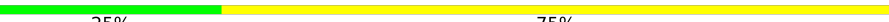
- Molecule 1: VP1 capsid protein



- Molecule 1: VP1 capsid protein

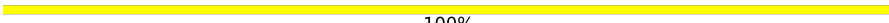


- Molecule 2: N-acetyl-alpha-neuraminic acid-(2-8)-N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain F:  25% 75%

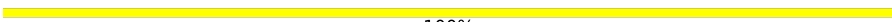
 BGL1
GAL2
SIA3
SIA4

- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-8)-N-acetyl-alpha-neuraminic acid

Chain G:  100%

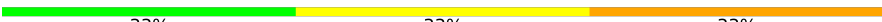
 SIA1
SIA2

- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-8)-N-acetyl-alpha-neuraminic acid

Chain I:  100%

 SIA1
SIA2

- Molecule 4: N-acetyl-alpha-neuraminic acid-(2-8)-N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose

Chain H:  33% 33% 33%

 GAL1
SIA2
SIA3

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	144.68Å 152.64Å 63.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.00 – 1.70 48.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.00-1.70) 99.9 (48.00-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.150 , 0.177 0.156 , 0.179	Depositor DCC
R_{free} test set	5395 reflections (3.42%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11705	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BGC, CL, GAL, SIA, GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.57	0/2132	0.72	0/2902
1	B	0.55	0/2101	0.71	0/2862
1	C	0.54	0/1975	0.68	0/2686
1	D	0.57	0/2138	0.71	0/2911
1	E	0.59	0/2177	0.72	0/2957
All	All	0.57	0/10523	0.71	0/14318

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2062	0	2013	5	0
1	B	2035	0	1963	4	0
1	C	1916	0	1852	10	0
1	D	2073	0	2029	8	0
1	E	2095	0	2073	7	0
2	F	63	0	53	0	0
3	G	41	0	34	0	0
3	I	41	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	52	0	44	1	0
5	A	6	0	8	0	0
5	B	12	0	16	0	0
5	C	6	0	8	0	0
5	D	6	0	8	0	0
5	E	12	0	16	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
6	E	1	0	0	0	0
7	A	300	0	0	1	0
7	B	245	0	0	0	0
7	C	192	0	0	0	0
7	D	266	0	0	0	0
7	E	277	0	0	1	0
All	All	11705	0	10151	31	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 2.

All (31) 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:C:109:MET:HE2	1:C:295:VAL:HG21	1.56	0.85
1:C:109:MET:CE	1:C:295:VAL:HG11	2.11	0.80
1:C:109:MET:HE2	1:C:295:VAL:HG11	1.78	0.65
1:C:113:VAL:HG12	1:C:114:THR:HG23	1.85	0.59
1:E:35:LEU:HD11	1:E:296:LYS:HG2	1.86	0.56
1:C:259:LEU:HD21	1:C:288:ILE:HD13	1.88	0.56
1:D:109:MET:HE1	1:D:255:LYS:HD3	1.90	0.53
1:A:259:LEU:HD21	1:A:288:ILE:HD13	1.89	0.53
1:C:109:MET:HE2	1:C:295:VAL:CG2	2.35	0.51
1:A:68:LYS:HD3	7:A:775:HOH:O	2.10	0.51
1:C:109:MET:HE3	1:C:295:VAL:HG11	1.93	0.50
1:C:255:LYS:O	1:C:256:ALA:HB3	2.14	0.48
1:E:98:LEU:HD21	1:E:110:TRP:CZ2	2.48	0.48
1:D:48:GLU:OE1	1:D:285:TYR:OH	2.28	0.48
1:E:259:LEU:HD21	1:E:288:ILE:HD13	1.96	0.47
1:D:188:VAL:HG23	1:D:189:MET:H	1.80	0.46
1:B:106:ASN:O	1:B:107:LEU:HD23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:231:PRO:HB3	1:D:226:GLY:O	2.18	0.43
1:D:98:LEU:HD21	1:D:110:TRP:CZ2	2.54	0.43
1:D:46:GLU:HG2	1:D:289:ARG:HG2	2.01	0.43
1:B:58:PRO:HA	1:B:83:LYS:HD3	2.01	0.42
1:A:231:PRO:HB3	1:B:230[A]:VAL:HG11	2.00	0.42
1:E:103:THR:O	1:E:103:THR:HG22	2.20	0.42
1:D:43:ALA:HB3	1:D:292:LYS:HE3	2.01	0.42
1:D:142:LYS:HE3	1:D:145[B]:GLN:OE1	2.20	0.42
1:A:231:PRO:HB3	1:B:226:GLY:O	2.20	0.41
1:E:188:VAL:HG23	1:E:189:MET:H	1.86	0.41
1:A:98:LEU:HD21	1:A:292:LYS:HE2	2.02	0.41
1:C:276:THR:CG2	4:H:2:SIA:H112	2.51	0.41
1:E:172:TYR:CG	1:E:178:THR:HG21	2.55	0.41
1:E:116[B]:GLN:HG3	7:E:680:HOH:O	2.21	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	271/275 (98%)	258 (95%)	12 (4%)	1 (0%)	30	16
1	B	266/275 (97%)	255 (96%)	10 (4%)	1 (0%)	30	16
1	C	246/275 (90%)	236 (96%)	9 (4%)	1 (0%)	30	16
1	D	271/275 (98%)	260 (96%)	10 (4%)	1 (0%)	30	16
1	E	272/275 (99%)	261 (96%)	10 (4%)	1 (0%)	30	16
All	All	1326/1375 (96%)	1270 (96%)	51 (4%)	5 (0%)	30	16

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	188	VAL
1	C	188	VAL
1	E	188	VAL
1	B	188	VAL
1	D	188	VAL

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	230/236 (98%)	230 (100%)	0	100	100
1	B	225/236 (95%)	225 (100%)	0	100	100
1	C	211/236 (89%)	211 (100%)	0	100	100
1	D	232/236 (98%)	232 (100%)	0	100	100
1	E	238/236 (101%)	237 (100%)	1 (0%)	84	80
All	All	1136/1180 (96%)	1135 (100%)	1 (0%)	88	85

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	247	GLU

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

Mol	Chain	Res	Type
1	E	248	GLN
1	E	272	ASN
1	E	278	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 ⓘ

11 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	BGC	F	1	2	12,12,12	0.46	0	17,17,17	0.65	0
2	GAL	F	2	2	11,11,12	0.47	0	15,15,17	0.96	1 (6%)
2	SIA	F	3	2	20,20,21	0.61	0	21,28,31	1.35	3 (14%)
2	SIA	F	4	2	20,20,21	0.62	0	21,28,31	1.21	3 (14%)
3	SIA	G	1	3	21,21,21	0.96	1 (4%)	24,31,31	1.34	5 (20%)
3	SIA	G	2	3	20,20,21	0.60	0	21,28,31	0.87	1 (4%)
4	GAL	H	1	4	12,12,12	0.49	0	17,17,17	0.74	0
4	SIA	H	2	4	20,20,21	0.63	0	21,28,31	1.39	3 (14%)
4	SIA	H	3	4	20,20,21	0.60	0	21,28,31	1.17	4 (19%)
3	SIA	I	1	3	21,21,21	0.98	1 (4%)	24,31,31	1.41	2 (8%)
3	SIA	I	2	3	20,20,21	0.68	0	21,28,31	1.04	2 (9%)

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	BGC	F	1	2	-	0/2/22/22	0/1/1/1
2	GAL	F	2	2	-	0/2/19/22	0/1/1/1
2	SIA	F	3	2	-	2/18/34/38	0/1/1/1
2	SIA	F	4	2	-	1/18/34/38	0/1/1/1
3	SIA	G	1	3	-	6/20/38/38	0/1/1/1
3	SIA	G	2	3	-	0/18/34/38	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GAL	H	1	4	-	0/2/22/22	0/1/1/1
4	SIA	H	2	4	-	2/18/34/38	0/1/1/1
4	SIA	H	3	4	-	0/18/34/38	0/1/1/1
3	SIA	I	1	3	-	6/20/38/38	0/1/1/1
3	SIA	I	2	3	-	0/18/34/38	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	1	SIA	O2-C2	3.20	1.44	1.39
3	I	1	SIA	O2-C2	2.96	1.43	1.39

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	1	SIA	O1A-C1-C2	-3.54	117.94	123.85
3	G	1	SIA	O1A-C1-C2	-3.37	118.23	123.85
2	F	3	SIA	O1B-C1-C2	3.34	121.41	112.71
2	F	3	SIA	O6-C2-C1	3.33	114.01	107.72
3	G	1	SIA	O6-C6-C5	3.09	112.67	109.84
4	H	2	SIA	O6-C2-C3	-3.00	106.52	110.56
2	F	4	SIA	O6-C2-C3	-2.89	106.67	110.56
3	I	2	SIA	O1B-C1-C2	2.63	119.56	112.71
4	H	2	SIA	O1B-C1-C2	2.59	119.44	112.71
3	I	1	SIA	O8-C8-C9	-2.56	103.20	109.03
3	I	2	SIA	O1A-C1-C2	-2.56	117.33	122.85
4	H	2	SIA	O6-C2-C1	2.50	112.44	107.72
3	G	1	SIA	O8-C8-C7	-2.48	103.44	109.25
2	F	3	SIA	O1A-C1-C2	-2.47	117.52	122.85
4	H	3	SIA	O1B-C1-C2	2.33	118.78	112.71
3	G	2	SIA	O1B-C1-C2	2.31	118.73	112.71
2	F	4	SIA	C8-C7-C6	-2.21	108.91	113.05
3	G	1	SIA	C3-C4-C5	2.16	113.06	109.72
2	F	4	SIA	O1B-C1-C2	2.11	118.19	112.71
4	H	3	SIA	C8-C7-C6	-2.09	109.13	113.05
2	F	2	GAL	O5-C5-C6	2.06	111.68	107.66
4	H	3	SIA	O6-C2-C1	2.05	111.59	107.72
4	H	3	SIA	O6-C2-C3	-2.05	107.80	110.56
3	G	1	SIA	C9-C8-C7	2.04	116.32	112.17

There are no chirality outliers.

All (17) torsion outliers are listed below:

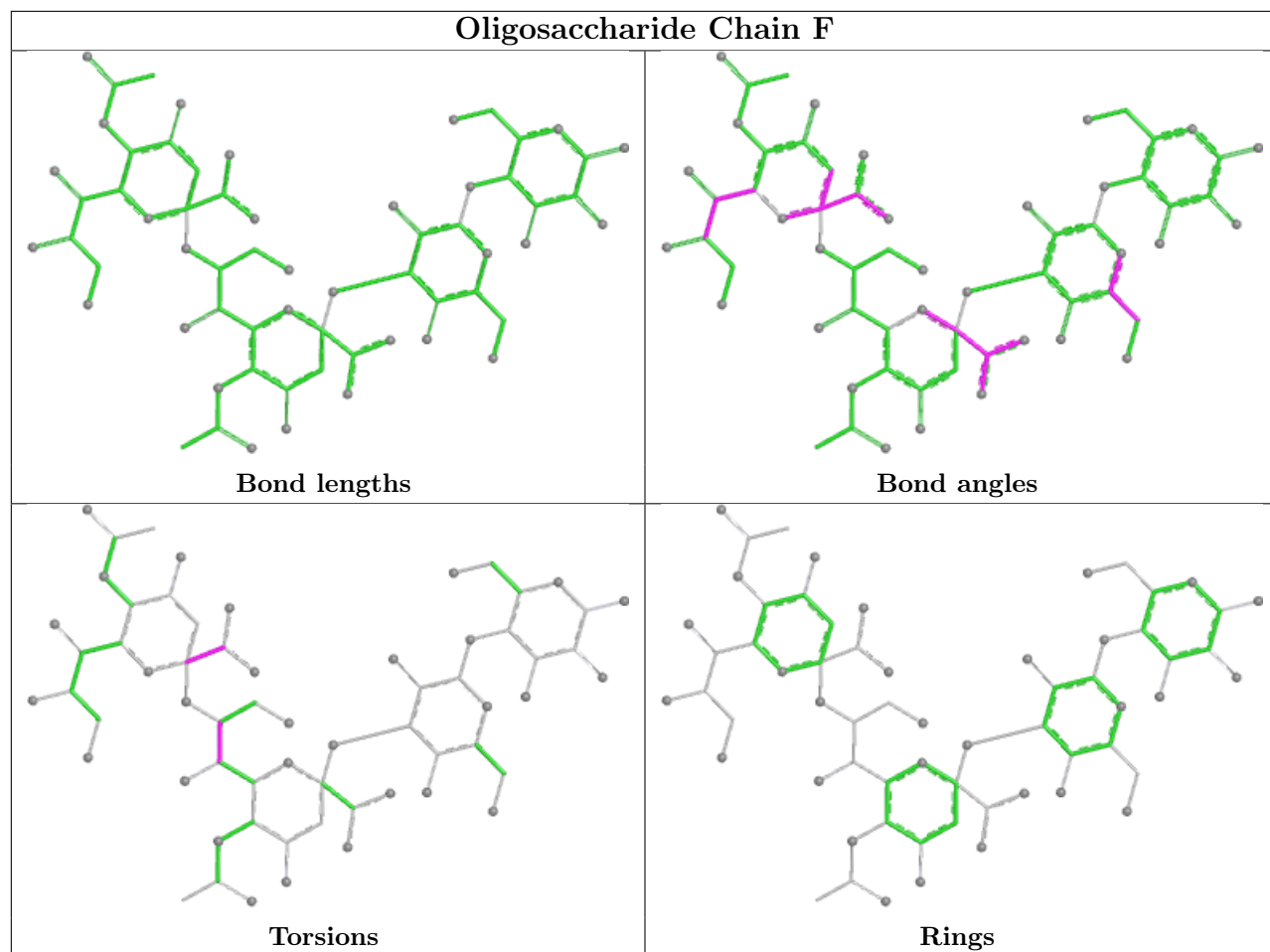
Mol	Chain	Res	Type	Atoms
3	G	1	SIA	O6-C6-C7-O7
3	I	1	SIA	C6-C7-C8-C9
3	I	1	SIA	C6-C7-C8-O8
3	I	1	SIA	O7-C7-C8-C9
3	I	1	SIA	O7-C7-C8-O8
4	H	2	SIA	C7-C8-C9-O9
3	G	1	SIA	C7-C8-C9-O9
4	H	2	SIA	O8-C8-C9-O9
3	G	1	SIA	O8-C8-C9-O9
2	F	3	SIA	C6-C7-C8-C9
3	I	1	SIA	O1B-C1-C2-C3
3	G	1	SIA	O1A-C1-C2-O6
2	F	3	SIA	C6-C7-C8-O8
2	F	4	SIA	O1A-C1-C2-O6
3	G	1	SIA	C5-C6-C7-O7
3	G	1	SIA	O1B-C1-C2-O6
3	I	1	SIA	O1A-C1-C2-C3

There are no ring outliers.

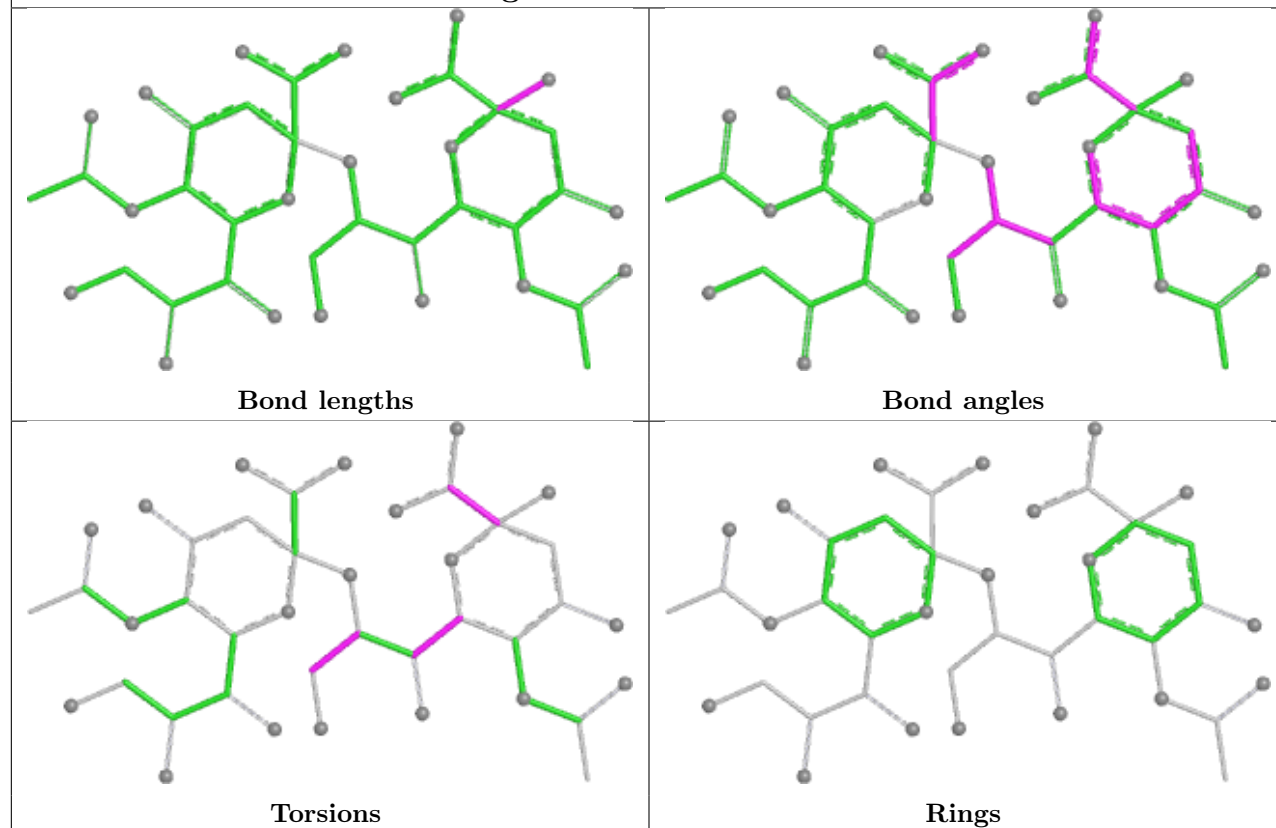
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	2	SIA	1	0

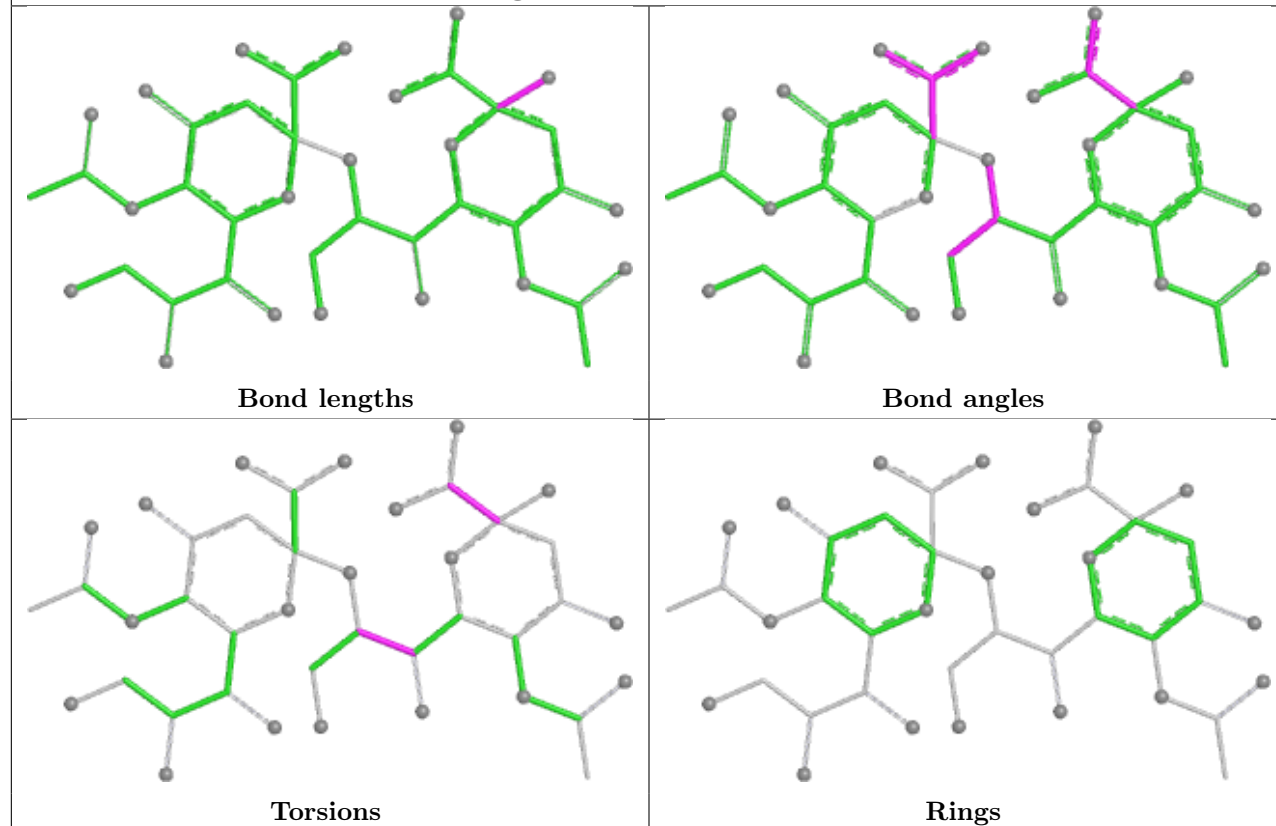
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

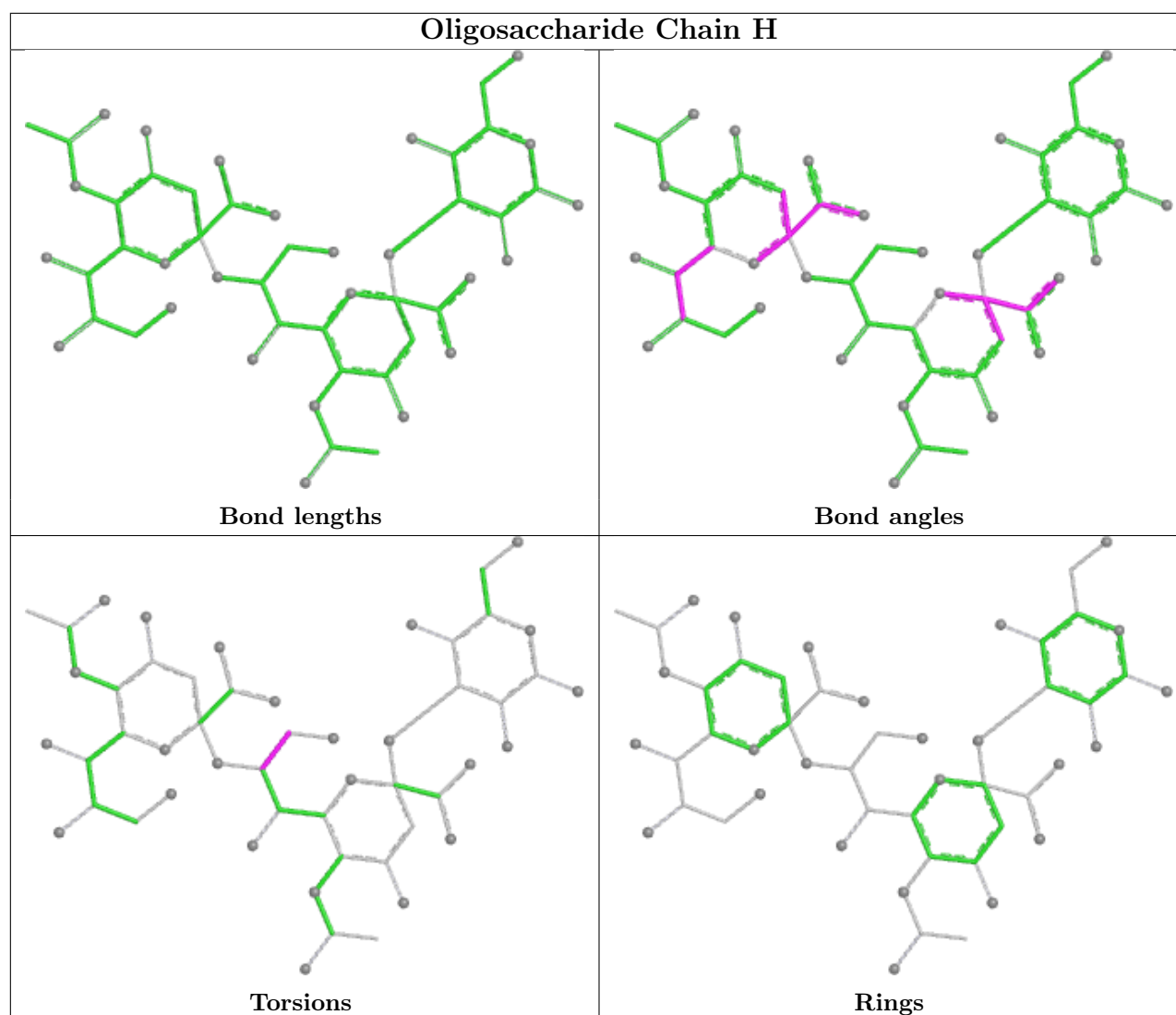


Oligosaccharide Chain G



Oligosaccharide Chain I





5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 5 are monoatomic - leaving 7 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	GOL	A	405	-	5,5,5	0.40	0	5,5,5	0.14	0
5	GOL	D	403	-	5,5,5	0.43	0	5,5,5	0.27	0
5	GOL	E	402	-	5,5,5	0.39	0	5,5,5	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GOL	B	404	-	5,5,5	0.39	0	5,5,5	0.34	0
5	GOL	C	404	-	5,5,5	0.44	0	5,5,5	0.38	0
5	GOL	E	401	-	5,5,5	0.43	0	5,5,5	0.25	0
5	GOL	B	403	-	5,5,5	0.45	0	5,5,5	0.28	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
5	GOL	A	405	-	-	4/4/4/4	-
5	GOL	D	403	-	-	2/4/4/4	-
5	GOL	E	402	-	-	2/4/4/4	-
5	GOL	B	404	-	-	4/4/4/4	-
5	GOL	C	404	-	-	4/4/4/4	-
5	GOL	E	401	-	-	4/4/4/4	-
5	GOL	B	403	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	405	GOL	O1-C1-C2-C3
5	A	405	GOL	C1-C2-C3-O3
5	B	404	GOL	O1-C1-C2-C3
5	B	404	GOL	C1-C2-C3-O3
5	C	404	GOL	O1-C1-C2-C3
5	E	401	GOL	O1-C1-C2-C3
5	C	404	GOL	O1-C1-C2-O2
5	B	403	GOL	C1-C2-C3-O3
5	C	404	GOL	C1-C2-C3-O3
5	D	403	GOL	C1-C2-C3-O3
5	E	402	GOL	C1-C2-C3-O3
5	B	403	GOL	O2-C2-C3-O3
5	B	404	GOL	O1-C1-C2-O2
5	B	404	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
5	D	403	GOL	O2-C2-C3-O3
5	A	405	GOL	O1-C1-C2-O2
5	E	402	GOL	O2-C2-C3-O3
5	A	405	GOL	O2-C2-C3-O3
5	C	404	GOL	O2-C2-C3-O3
5	E	401	GOL	O1-C1-C2-O2
5	E	401	GOL	O2-C2-C3-O3
5	E	401	GOL	C1-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	265/275 (96%)	-0.41	4 (1%) 72 76	9, 17, 33, 48	8 (3%)
1	B	263/275 (95%)	-0.22	11 (4%) 40 45	10, 20, 56, 75	7 (2%)
1	C	247/275 (89%)	-0.11	12 (4%) 35 38	12, 22, 60, 87	5 (2%)
1	D	267/275 (97%)	-0.39	4 (1%) 72 76	9, 17, 37, 60	6 (2%)
1	E	264/275 (96%)	-0.38	9 (3%) 48 52	7, 16, 44, 90	12 (4%)
All	All	1306/1375 (94%)	-0.31	40 (3%) 51 55	7, 18, 45, 90	38 (2%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	107	LEU	5.7
1	B	102	LEU	5.1
1	E	298	PRO	4.5
1	E	102	LEU	4.4
1	B	39	THR	4.3
1	E	44	ILE	4.0
1	E	103	THR	3.7
1	C	108	LEU	3.5
1	B	103	THR	3.5
1	B	105	GLY	3.5
1	C	34	VAL	3.4
1	C	44	ILE	3.3
1	B	104	CYS	3.3
1	C	35	LEU	3.2
1	B	255	LYS	3.2
1	C	255	LYS	3.2
1	E	105	GLY	3.2
1	D	298	PRO	3.1
1	E	104	CYS	3.0
1	C	37	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	E	32	VAL	2.8
1	C	36	GLU	2.7
1	E	100	GLU	2.6
1	D	104	CYS	2.6
1	A	107	LEU	2.5
1	B	44	ILE	2.5
1	C	96	PRO	2.5
1	C	97	ASN	2.5
1	A	32	VAL	2.5
1	D	105	GLY	2.4
1	E	40	GLY	2.3
1	B	42	ASP	2.2
1	C	295	VAL	2.2
1	A	103	THR	2.2
1	B	101	ASP	2.1
1	B	33	GLU	2.1
1	A	296	LYS	2.1
1	B	38	LYS	2.1
1	C	256	ALA	2.0
1	D	297	ASN	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
4	GAL	H	1	12/12	0.54	0.29	38,40,41,43	12
4	SIA	H	2	20/21	0.73	0.14	35,42,49,50	0
2	BGC	F	1	12/12	0.82	0.10	30,38,43,44	0
3	SIA	G	1	21/21	0.82	0.11	31,45,51,54	0
2	SIA	F	3	20/21	0.87	0.10	29,32,39,41	0
3	SIA	I	1	21/21	0.89	0.10	21,31,34,36	0
2	GAL	F	2	11/12	0.90	0.09	25,30,35,41	0
4	SIA	H	3	20/21	0.92	0.09	25,29,33,35	0

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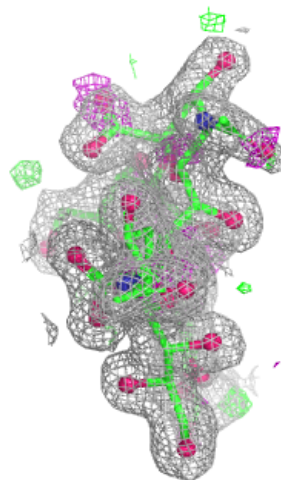
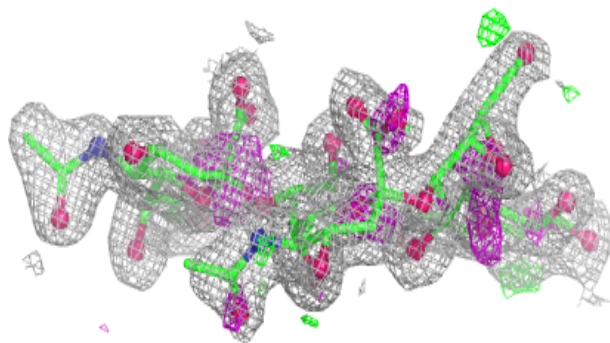
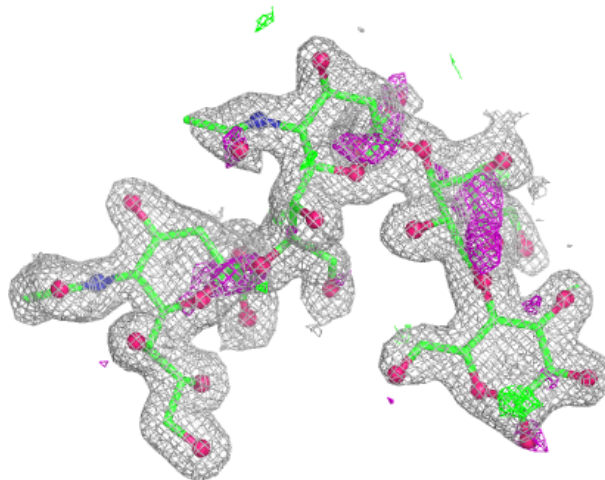
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SIA	F	4	20/21	0.93	0.08	21,27,32,32	0
3	SIA	G	2	20/21	0.94	0.07	20,25,29,34	0
3	SIA	I	2	20/21	0.96	0.06	17,19,22,22	0

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.

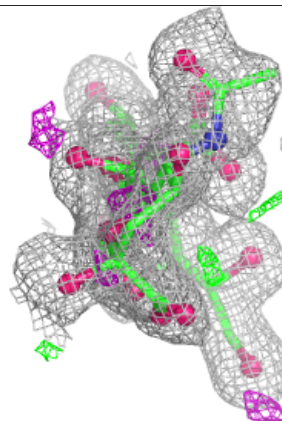
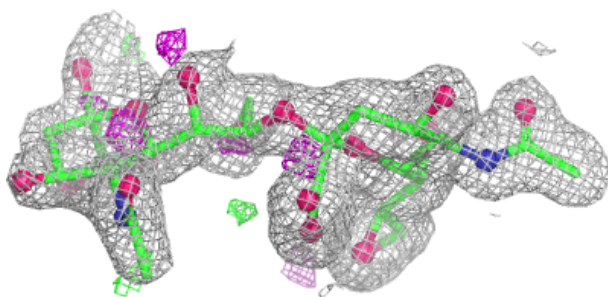
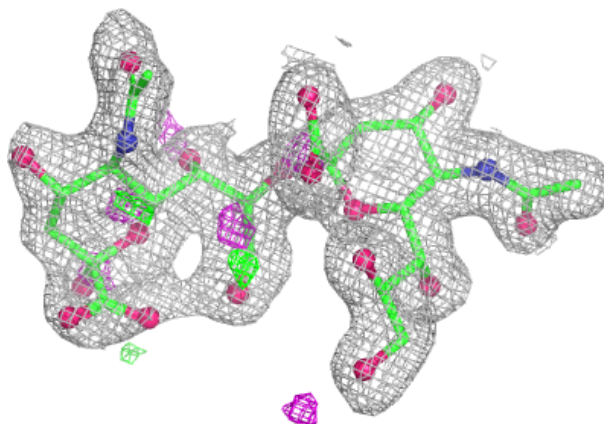
Electron density around Chain F:

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



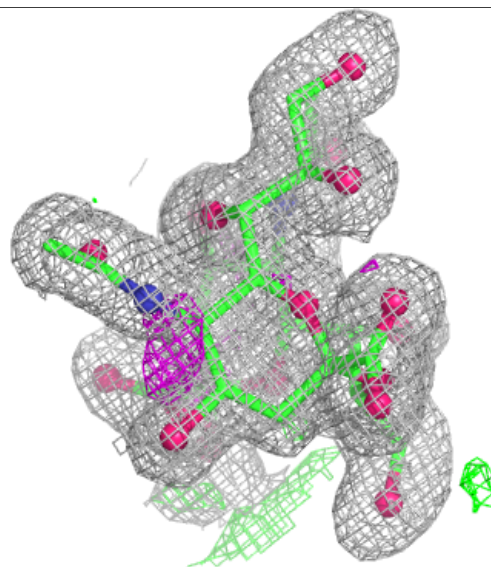
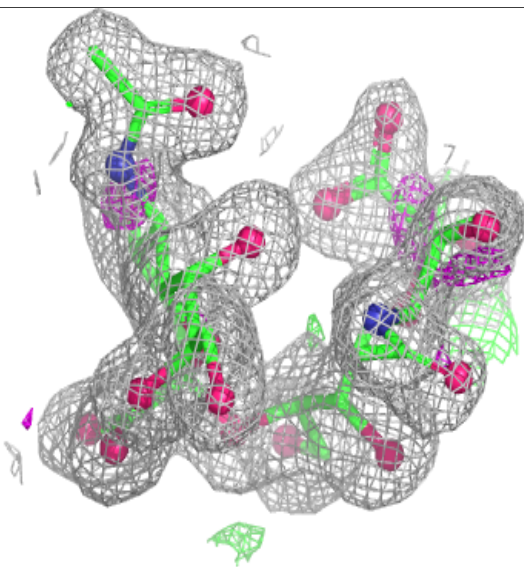
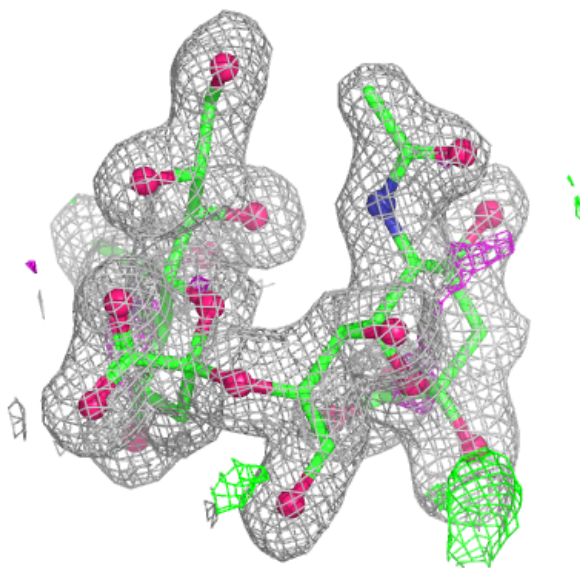
Electron density around Chain G:

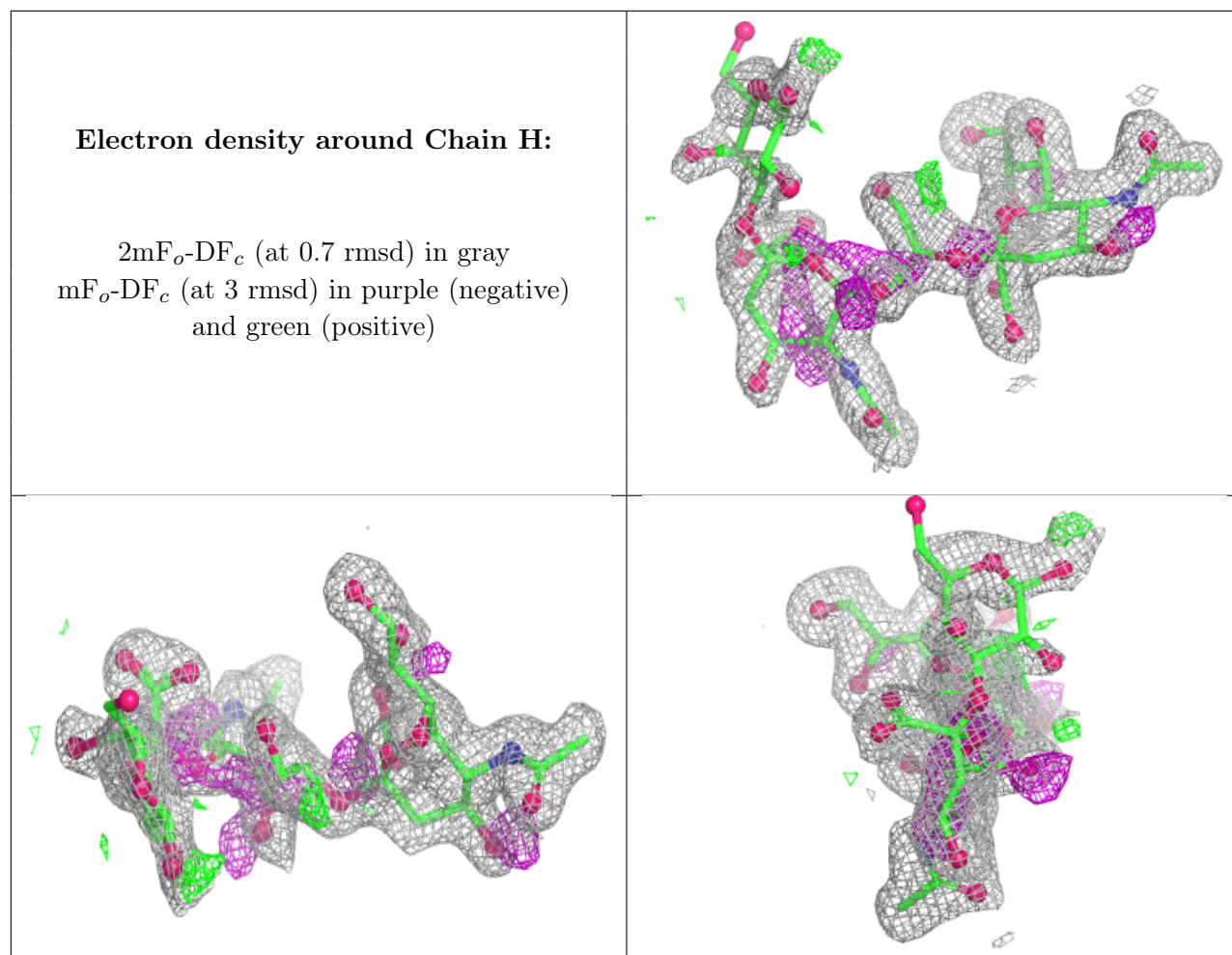
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain I:

$2mF_o-DF_c$ (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(Å ²)	Q<0.9
5	GOL	B	404	6/6	0.76	0.15	32,41,45,48	0
5	GOL	A	405	6/6	0.79	0.14	33,39,42,42	0
5	GOL	D	403	6/6	0.79	0.17	36,44,46,48	0
5	GOL	E	402	6/6	0.79	0.15	30,41,43,47	0
5	GOL	C	404	6/6	0.84	0.12	32,41,43,46	0
5	GOL	E	401	6/6	0.85	0.12	37,42,44,45	0
5	GOL	B	403	6/6	0.86	0.11	40,43,44,46	0
6	CL	C	405	1/1	0.94	0.14	35,35,35,35	0
6	CL	A	406	1/1	0.96	0.09	26,26,26,26	0
6	CL	B	405	1/1	0.97	0.09	28,28,28,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	CL	D	404	1/1	0.97	0.11	27,27,27,27	0
6	CL	E	403	1/1	0.98	0.07	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.