



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:00 AM UTC

PDB ID : 4N63 / pdb_00004n63
Title : Crystal structure of hemagglutinin from an H7N9 influenza virus in complex with an O-linked glycan receptor
Authors : Xu, R.; Wilson, I.A.
Deposited on : 2013-10-11
Resolution : 2.75 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

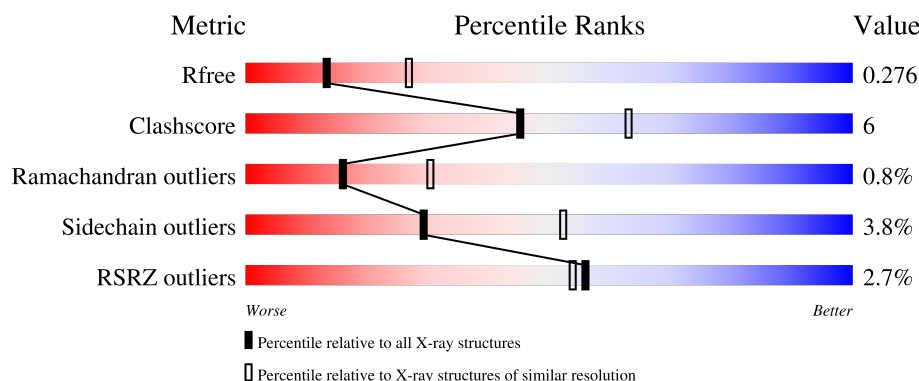
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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.



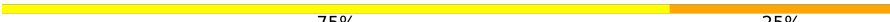
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	321	79% 19% ..
1	C	321	4% 81% 17% ..
2	B	183	2% 79% 13% 8%
2	D	183	4% 84% 8% 8%
3	E	2	100%

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Mol	Chain	Length	Quality of chain
4	F	4	 75%25%
4	G	4	 50%50%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7824 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 HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	0	0
			2427	1509	438	465	15			
1	C	318	Total	C	N	O	S	0	0	0
			2427	1509	438	465	15			

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	169	Total	C	N	O	S	0	0	0
			1373	845	241	280	7			
2	D	169	Total	C	N	O	S	0	0	0
			1373	845	241	280	7			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	177	SER	-	expression tag	UNP R4NN21
B	178	GLY	-	expression tag	UNP R4NN21
B	179	ARG	-	expression tag	UNP R4NN21
B	180	LEU	-	expression tag	UNP R4NN21
B	181	VAL	-	expression tag	UNP R4NN21
B	182	PRO	-	expression tag	UNP R4NN21
B	183	ARG	-	expression tag	UNP R4NN21
D	177	SER	-	expression tag	UNP R4NN21
D	178	GLY	-	expression tag	UNP R4NN21
D	179	ARG	-	expression tag	UNP R4NN21
D	180	LEU	-	expression tag	UNP R4NN21
D	181	VAL	-	expression tag	UNP R4NN21
D	182	PRO	-	expression tag	UNP R4NN21
D	183	ARG	-	expression tag	UNP R4NN21

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

cetamido-2-deoxy-beta-D-glucopyranose.



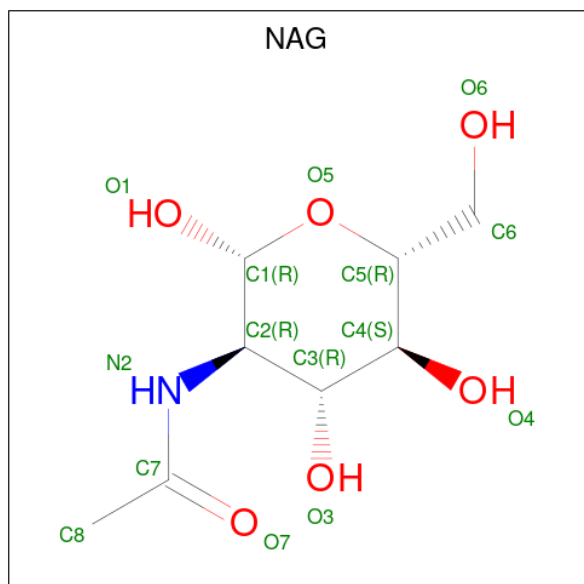
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

- 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-(1-3)-2-acetamido-2-deoxy-beta-D-galactopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	4	Total	C	N	O	0	0	0
			60	33	3	24			
4	G	4	Total	C	N	O	0	0	0
			60	33	3	24			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			14	8	1	5		

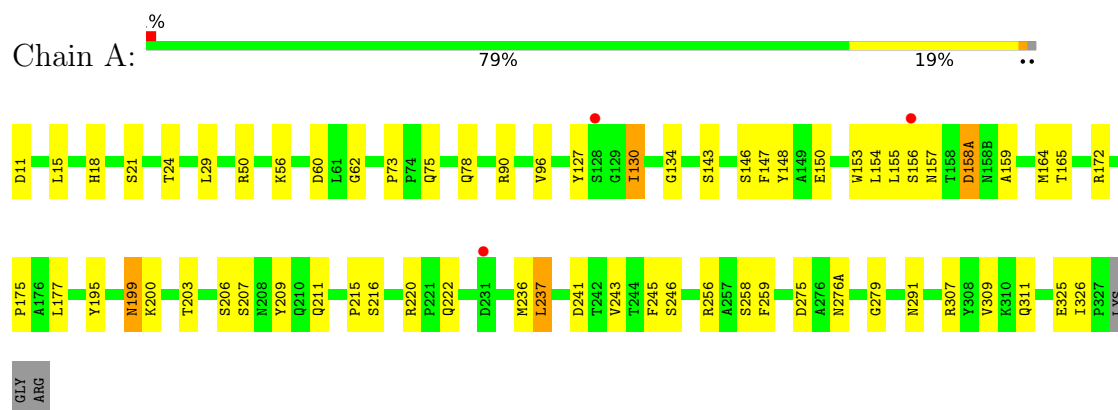
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	28	Total	O	0	0
			28	28		
6	B	24	Total	O	0	0
			24	24		
6	C	7	Total	O	0	0
			7	7		
6	D	3	Total	O	0	0
			3	3		

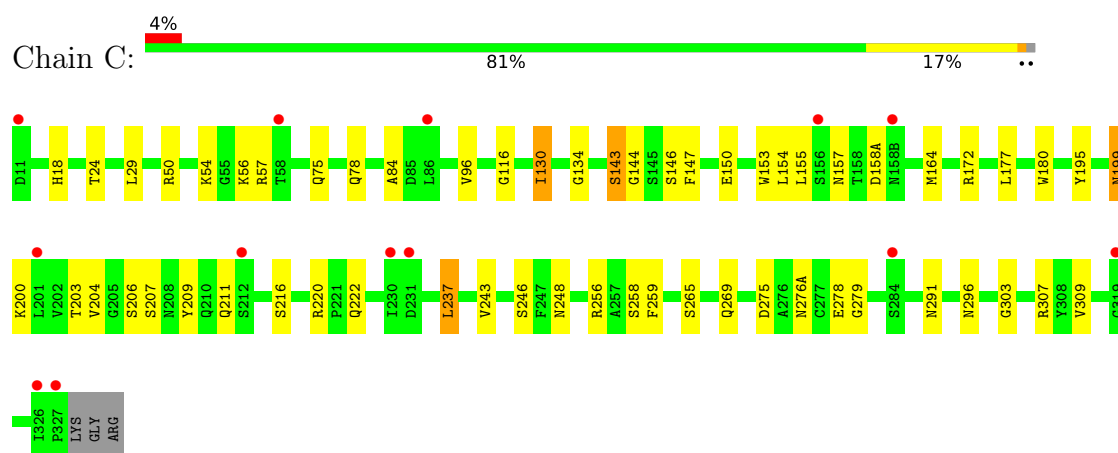
3 Residue-property plots [i](#)

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.

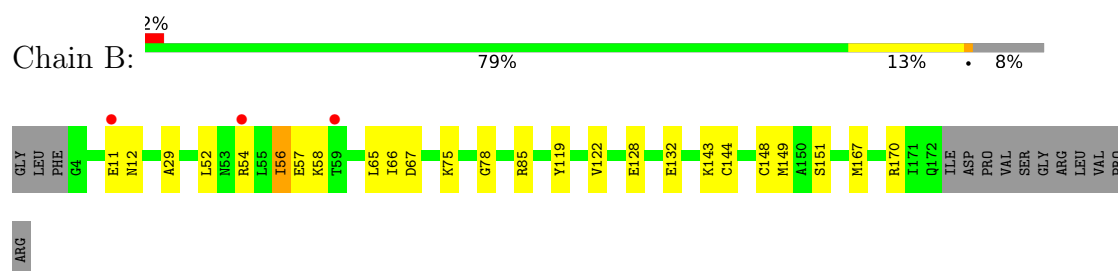
• Molecule 1: Hemagglutinin HA1



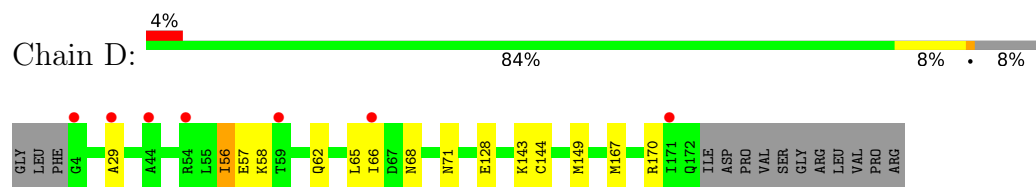
• Molecule 1: Hemagglutinin HA1



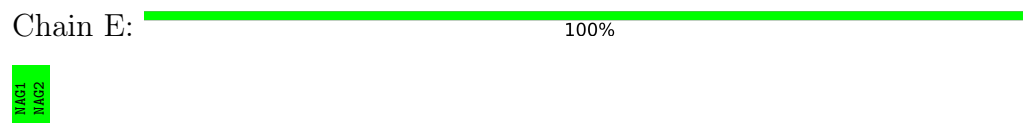
• Molecule 2: Hemagglutinin HA2



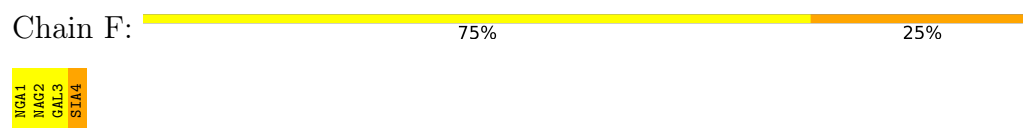
- Molecule 2: Hemagglutinin HA2



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



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	153.61Å 153.61Å 153.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.60 – 2.75 42.60 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.2 (42.60-2.75) 98.9 (42.60-2.75)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.213 , 0.269 0.217 , 0.276	Depositor DCC
R_{free} test set	1584 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	64.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 71.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.035 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7824	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.06% 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: GAL, NAG, NGA, 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	0.28	0/2474	0.71	1/3345 (0.0%)
1	C	0.27	0/2474	0.71	0/3345
2	B	0.29	0/1396	0.67	0/1881
2	D	0.28	0/1396	0.67	0/1881
All	All	0.28	0/7740	0.70	1/10452 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	SER	N-CA-C	-5.02	107.15	113.28

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	2427	0	2388	40	0
1	C	2427	0	2389	34	0
2	B	1373	0	1273	18	0
2	D	1373	0	1274	11	0
3	E	28	0	25	0	0
4	F	60	0	52	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	G	60	0	52	3	0
5	B	14	0	13	2	0
6	A	28	0	0	1	0
6	B	24	0	0	0	0
6	C	7	0	0	0	0
6	D	3	0	0	0	0
All	All	7824	0	7466	90	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 6.

All (90) 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:291:ASN:OD1	2:D:58:LYS:NZ	1.93	1.01
2:B:128:GLU:HG3	2:B:170:ARG:HH12	1.36	0.90
1:C:56:LYS:NZ	1:C:279:GLY:O	2.13	0.81
2:D:128:GLU:HG3	2:D:170:ARG:HH12	1.47	0.79
1:A:56:LYS:NZ	1:A:279:GLY:O	2.16	0.78
1:A:222:GLN:NE2	4:F:2:NAG:O7	2.17	0.77
2:B:128:GLU:HG3	2:B:170:ARG:NH1	2.05	0.71
2:D:128:GLU:HG3	2:D:170:ARG:NH1	2.05	0.71
1:C:155:LEU:HD21	4:G:4:SIA:H111	1.73	0.71
1:C:222:GLN:NE2	4:G:2:NAG:O7	2.24	0.69
2:B:29:ALA:H	2:B:143:LYS:HZ1	1.43	0.65
2:D:29:ALA:H	2:D:143:LYS:HZ1	1.42	0.64
1:C:203:THR:OG1	1:C:246:SER:OG	2.17	0.62
1:A:177:LEU:HB3	1:A:258:SER:HB2	1.80	0.62
1:A:275:ASP:OD1	1:A:276(A):ASN:ND2	2.33	0.62
1:A:195:TYR:O	1:A:200:LYS:NZ	2.28	0.62
1:C:275:ASP:OD1	1:C:276(A):ASN:ND2	2.33	0.61
1:A:155:LEU:HD21	4:F:4:SIA:H111	1.83	0.60
1:A:291:ASN:OD1	2:B:58:LYS:NZ	2.35	0.58
2:B:75:LYS:HD3	5:B:201:NAG:H81	1.84	0.58
1:C:130:ILE:HD11	1:C:154:LEU:HB3	1.84	0.58
1:C:134:GLY:HA3	1:C:153:TRP:HB3	1.85	0.58
1:A:146:SER:OG	1:A:147:PHE:N	2.34	0.58
1:A:11:ASP:N	6:A:507:HOH:O	2.37	0.57
1:A:134:GLY:HA3	1:A:153:TRP:HB3	1.85	0.57
1:C:199:ASN:OD1	1:C:199:ASN:N	2.37	0.56
1:A:50:ARG:NE	1:A:275:ASP:OD2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ILE:HD11	1:A:154:LEU:HB3	1.87	0.56
2:B:119:TYR:OH	2:B:132:GLU:OE2	2.13	0.55
1:C:206:SER:HB3	1:C:209:TYR:HB3	1.88	0.55
1:C:206:SER:OG	1:C:207:SER:N	2.39	0.55
1:A:150:GLU:OE1	1:A:256:ARG:HD3	2.08	0.54
1:C:291:ASN:HA	2:D:58:LYS:NZ	2.23	0.53
1:A:206:SER:OG	1:A:207:SER:N	2.41	0.53
1:C:195:TYR:O	1:C:200:LYS:NZ	2.35	0.52
1:A:325:GLU:HG3	2:B:12:ASN:HD22	1.75	0.52
1:A:237:LEU:HG	1:A:243:VAL:HG13	1.92	0.52
1:C:216:SER:O	1:C:220:ARG:NH2	2.38	0.52
1:A:203:THR:OG1	1:A:246:SER:OG	2.22	0.51
1:A:325:GLU:HA	2:B:12:ASN:HB2	1.92	0.51
2:B:128:GLU:O	2:B:170:ARG:NH1	2.37	0.51
1:C:303:GLY:HA2	2:D:62:GLN:HG3	1.93	0.50
1:C:150:GLU:OE1	1:C:256:ARG:HD3	2.10	0.50
1:A:326:ILE:HG12	2:B:12:ASN:HB3	1.94	0.50
2:D:144:CYS:HB3	2:D:149:MET:HE2	1.95	0.49
2:B:78:GLY:HA3	5:B:201:NAG:H82	1.95	0.48
1:A:237:LEU:HD22	1:A:241:ASP:HB3	1.96	0.48
1:C:180:TRP:CE2	1:C:204:VAL:HG21	2.49	0.48
1:A:172:ARG:HD3	1:A:259:PHE:CZ	2.49	0.48
1:C:146:SER:OG	1:C:147:PHE:N	2.36	0.48
1:C:237:LEU:HG	1:C:243:VAL:HG13	1.96	0.47
1:A:216:SER:O	1:A:220:ARG:NH2	2.37	0.47
1:C:269:GLN:HE22	2:D:68:ASN:CG	2.23	0.47
1:C:296:ASN:ND2	1:C:309:VAL:O	2.42	0.47
1:C:291:ASN:HA	2:D:58:LYS:HZ2	1.79	0.47
1:A:206:SER:HB3	1:A:209:TYR:HB3	1.97	0.46
1:A:158(A):ASP:C	1:A:159:ALA:H	2.23	0.46
1:C:200:LYS:HA	1:C:248:ASN:HD21	1.79	0.46
1:C:54:LYS:HD3	1:C:278:GLU:HB2	1.98	0.46
1:A:199:ASN:OD1	1:A:199:ASN:N	2.49	0.45
1:A:291:ASN:HA	2:B:58:LYS:NZ	2.32	0.45
1:C:57:ARG:HG2	1:C:84:ALA:HA	1.98	0.45
1:A:60:ASP:OD1	1:A:90:ARG:NE	2.34	0.45
1:C:172:ARG:HD3	1:C:259:PHE:CZ	2.52	0.44
2:B:148:CYS:O	2:B:151:SER:OG	2.28	0.44
1:C:177:LEU:HB3	1:C:258:SER:HB2	2.00	0.44
1:A:175:PRO:HB2	1:A:236:MET:HE3	1.98	0.44
4:G:2:NAG:O3	4:G:3:GAL:O5	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:307:ARG:HH11	1:C:307:ARG:HG2	1.83	0.43
1:C:116:GLY:HA2	1:C:265:SER:HB3	2.00	0.43
1:A:127:TYR:HB3	1:A:130:ILE:HG21	2.00	0.43
1:C:291:ASN:HB3	2:D:56:ILE:HG13	2.01	0.42
1:A:130:ILE:HA	1:A:156:SER:O	2.20	0.42
1:C:130:ILE:HG21	1:C:164:MET:HE1	2.01	0.42
1:A:309:VAL:HB	1:A:311:GLN:OE1	2.19	0.42
1:A:15:LEU:HD11	2:B:122:VAL:HG21	2.01	0.42
2:B:67:ASP:OD2	2:B:85:ARG:NH2	2.43	0.42
1:C:143:SER:OG	1:C:144:GLY:N	2.52	0.42
1:A:291:ASN:HA	2:B:58:LYS:HZ2	1.85	0.42
1:A:73:PRO:HG3	1:A:148:TYR:CE1	2.55	0.41
1:A:307:ARG:HH11	1:A:307:ARG:HG2	1.85	0.41
2:D:68:ASN:HB3	2:D:71:ASN:O	2.20	0.41
1:C:50:ARG:NE	1:C:275:ASP:OD2	2.53	0.41
1:A:62:GLY:O	1:A:90:ARG:HB2	2.20	0.41
1:A:291:ASN:HB3	2:B:56:ILE:HG13	2.01	0.41
2:B:144:CYS:HB3	2:B:149:MET:HE2	2.02	0.41
1:A:165:THR:HA	1:A:245:PHE:O	2.21	0.41
1:C:237:LEU:HD23	1:C:237:LEU:HA	1.89	0.40
1:A:130:ILE:HG21	1:A:164:MET:HE1	2.03	0.40
1:A:200:LYS:HB2	1:A:215:PRO:HG3	2.04	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	316/321 (98%)	293 (93%)	20 (6%)	3 (1%)	14	28
1	C	316/321 (98%)	293 (93%)	20 (6%)	3 (1%)	14	28
2	B	167/183 (91%)	158 (95%)	8 (5%)	1 (1%)	21	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	167/183 (91%)	157 (94%)	9 (5%)	1 (1%)	21	38
All	All	966/1008 (96%)	901 (93%)	57 (6%)	8 (1%)	16	30

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	157	ASN
2	B	57	GLU
1	C	157	ASN
1	A	158(A)	ASP
1	C	158(A)	ASP
2	D	57	GLU
1	A	143	SER
1	C	143	SER

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	267/269 (99%)	257 (96%)	10 (4%)	30	54
1	C	267/269 (99%)	257 (96%)	10 (4%)	30	54
2	B	145/157 (92%)	138 (95%)	7 (5%)	23	44
2	D	145/157 (92%)	141 (97%)	4 (3%)	38	61
All	All	824/852 (97%)	793 (96%)	31 (4%)	29	52

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	24	THR
1	A	29	LEU
1	A	75	GLN
1	A	78	GLN
1	A	96	VAL

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Mol	Chain	Res	Type
1	A	130	ILE
1	A	199	ASN
1	A	211	GLN
1	A	237	LEU
2	B	11	GLU
2	B	52	LEU
2	B	54	ARG
2	B	56	ILE
2	B	65	LEU
2	B	66	ILE
2	B	167	MET
1	C	18	HIS
1	C	24	THR
1	C	29	LEU
1	C	75	GLN
1	C	78	GLN
1	C	96	VAL
1	C	130	ILE
1	C	199	ASN
1	C	211	GLN
1	C	237	LEU
2	D	56	ILE
2	D	65	LEU
2	D	66	ILE
2	D	167	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	211	GLN
2	B	26	HIS
2	B	30	GLN
2	B	125	GLN
2	B	155	ASN
2	B	159	HIS
2	D	30	GLN
2	D	105	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 ⓘ

10 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	E	1	1,3	14,14,15	0.54	0	17,19,21	0.63	0
3	NAG	E	2	3	14,14,15	0.22	0	17,19,21	0.46	0
4	NGA	F	1	4	15,15,15	0.48	0	21,21,21	1.43	2 (9%)
4	NAG	F	2	4	14,14,15	0.43	0	17,19,21	0.48	0
4	GAL	F	3	4	11,11,12	1.06	2 (18%)	15,15,17	1.03	1 (6%)
4	SIA	F	4	4	20,20,21	0.72	0	21,28,31	1.18	4 (19%)
4	NGA	G	1	4	15,15,15	0.43	0	21,21,21	1.15	2 (9%)
4	NAG	G	2	4	14,14,15	0.29	0	17,19,21	0.50	0
4	GAL	G	3	4	11,11,12	1.15	2 (18%)	15,15,17	1.07	2 (13%)
4	SIA	G	4	4	20,20,21	0.73	0	21,28,31	1.17	3 (14%)

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	E	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
4	NGA	F	1	4	-	4/6/26/26	0/1/1/1
4	NAG	F	2	4	-	1/6/23/26	0/1/1/1
4	GAL	F	3	4	-	0/2/19/22	0/1/1/1
4	SIA	F	4	4	-	4/18/34/38	0/1/1/1
4	NGA	G	1	4	-	3/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
4	GAL	G	3	4	-	0/2/19/22	0/1/1/1
4	SIA	G	4	4	-	4/18/34/38	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	3	GAL	C2-C3	2.52	1.56	1.52
4	F	3	GAL	C2-C3	2.17	1.55	1.52
4	F	3	GAL	C1-C2	2.09	1.57	1.52
4	G	3	GAL	C1-C2	2.06	1.57	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	1	NGA	C1-C2-N2	-3.60	106.56	110.73
4	G	1	NGA	O5-C1-C2	3.00	112.53	109.52
4	F	4	SIA	O6-C2-C3	2.58	114.03	110.56
4	F	4	SIA	O1B-C1-C2	2.55	119.35	112.71
4	F	3	GAL	C1-C2-C3	2.54	113.34	109.64
4	G	4	SIA	O1B-C1-C2	2.52	119.27	112.71
4	G	4	SIA	O6-C2-C1	-2.45	103.09	107.72
4	F	4	SIA	O1A-C1-C2	-2.41	117.64	122.85
4	G	4	SIA	O1A-C1-C2	-2.38	117.71	122.85
4	G	3	GAL	C1-C2-C3	2.30	113.00	109.64
4	F	1	NGA	O4-C4-C3	2.24	115.66	110.38
4	F	4	SIA	O6-C2-C1	-2.20	103.55	107.72
4	G	3	GAL	C1-O5-C5	2.08	114.98	112.19
4	G	1	NGA	O3-C3-C2	2.05	113.66	109.58

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	1	NGA	C1-C2-N2-C7
4	F	1	NGA	C3-C2-N2-C7
4	F	1	NGA	C8-C7-N2-C2
4	F	1	NGA	O7-C7-N2-C2
4	F	4	SIA	C6-C7-C8-O8
4	F	4	SIA	O7-C7-C8-O8
4	G	1	NGA	C3-C2-N2-C7

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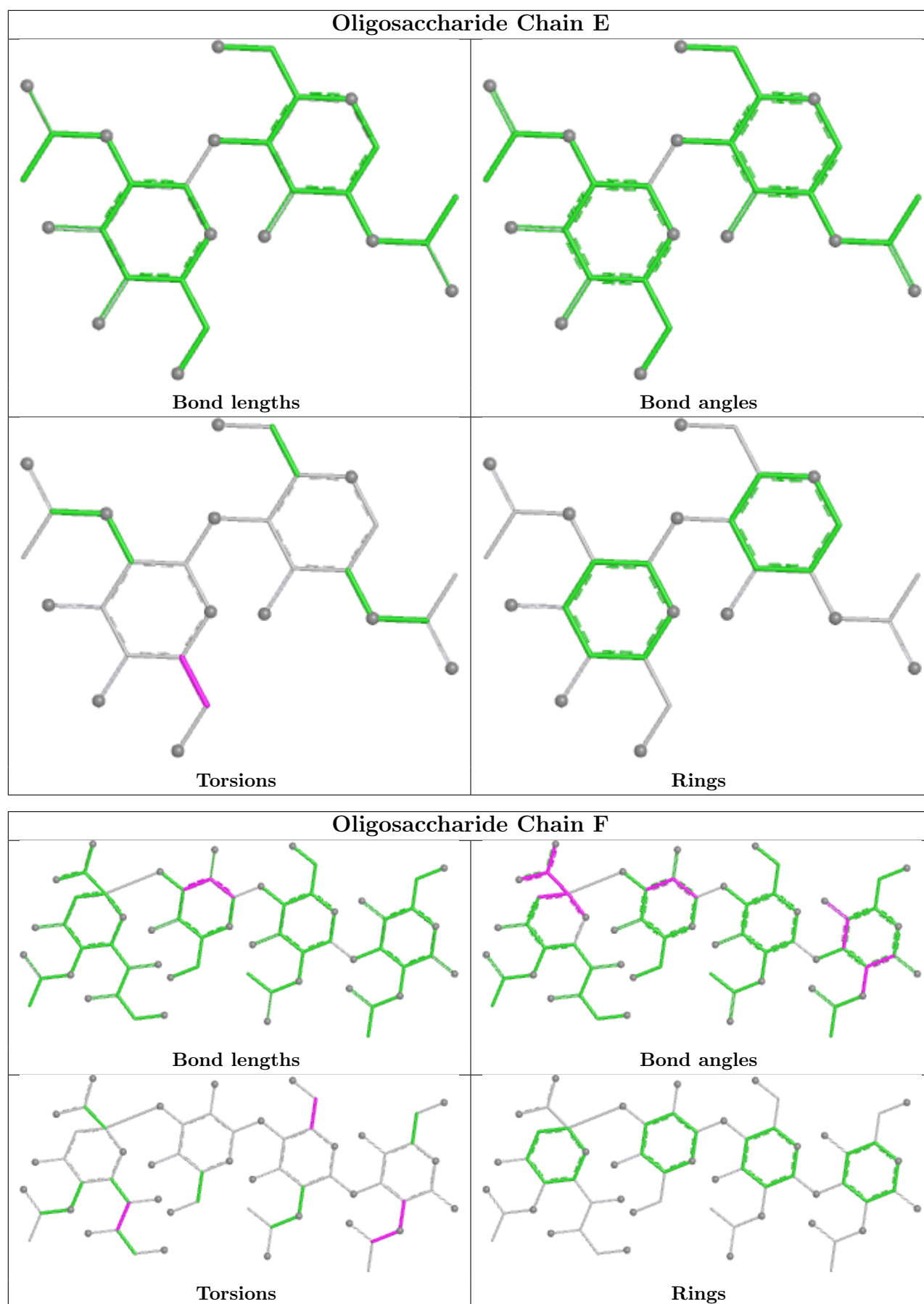
Mol	Chain	Res	Type	Atoms
4	G	1	NGA	C8-C7-N2-C2
4	G	1	NGA	O7-C7-N2-C2
4	G	4	SIA	C6-C7-C8-C9
4	G	4	SIA	C6-C7-C8-O8
4	G	4	SIA	O7-C7-C8-O8
4	G	2	NAG	C4-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
4	F	4	SIA	O7-C7-C8-C9
4	G	4	SIA	O7-C7-C8-C9
4	F	4	SIA	C6-C7-C8-C9
3	E	2	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
4	F	2	NAG	C4-C5-C6-O6

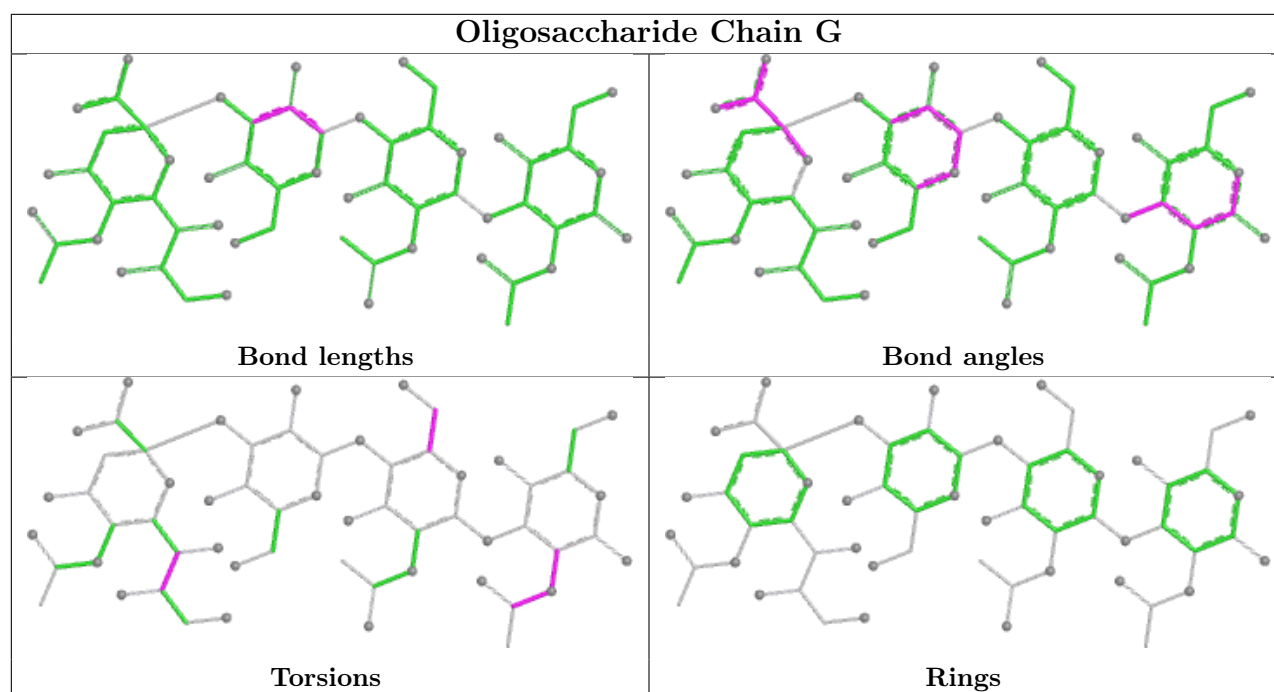
There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	4	SIA	1	0
4	G	2	NAG	2	0
4	G	4	SIA	1	0
4	G	3	GAL	1	0
4	F	2	NAG	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.





5.6 Ligand geometry [i](#)

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	NAG	B	201	2	14,14,15	0.35	0	17,19,21	0.43	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	NAG	B	201	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	201	NAG	2	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 ⓘ

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	318/321 (99%)	-0.19	3 (0%) 81 81	47, 69, 104, 172	0
1	C	318/321 (99%)	0.57	13 (4%) 41 39	70, 115, 173, 227	0
2	B	169/183 (92%)	-0.27	3 (1%) 67 65	42, 61, 106, 172	0
2	D	169/183 (92%)	0.61	7 (4%) 41 39	92, 153, 189, 252	0
All	All	974/1008 (96%)	0.18	26 (2%) 56 54	42, 91, 170, 252	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	231	ASP	3.6
1	C	156	SER	3.3
2	D	4	GLY	3.2
1	C	11	ASP	3.1
1	A	156	SER	3.1
1	C	231	ASP	3.0
1	C	326	ILE	2.8
2	D	171	ILE	2.7
2	B	11	GLU	2.6
2	D	44	ALA	2.5
1	C	230	ILE	2.4
1	C	319	GLY	2.4
2	B	54	ARG	2.4
2	D	29	ALA	2.4
2	B	59	THR	2.3
2	D	66	ILE	2.2
1	C	284	SER	2.2
1	A	128	SER	2.2
1	C	86	LEU	2.1
1	C	212	SER	2.1
2	D	54	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	58	THR	2.1
1	C	327	PRO	2.1
1	C	201	LEU	2.1
1	C	158(B)	ASN	2.0
2	D	59	THR	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](#)

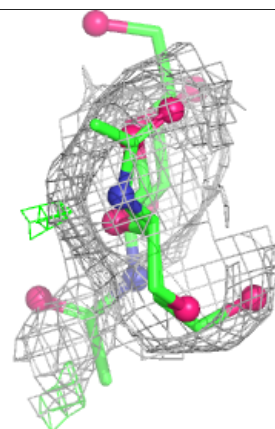
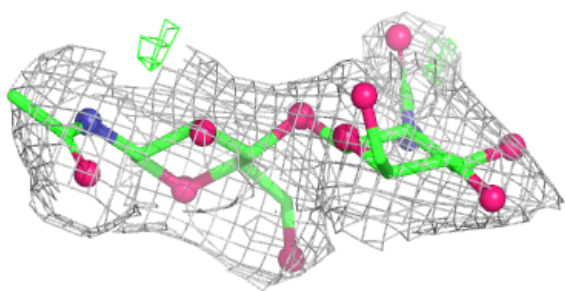
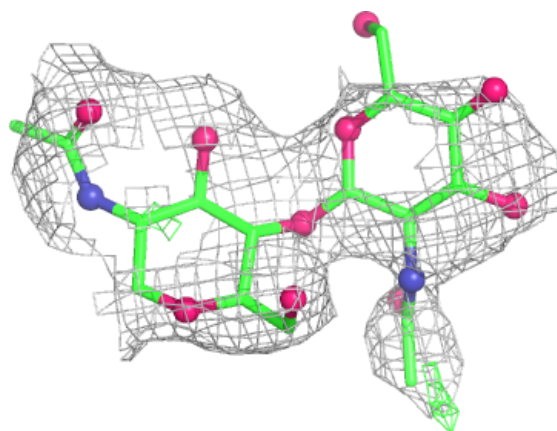
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	NGA	G	1	15/15	0.25	0.17	156,164,187,197	0
3	NAG	E	2	14/15	0.57	0.12	131,148,164,167	0
4	NAG	G	2	14/15	0.60	0.17	182,208,213,213	0
4	GAL	G	3	11/12	0.72	0.12	96,109,135,151	0
4	NGA	F	1	15/15	0.73	0.11	107,129,140,149	0
4	NAG	F	2	14/15	0.83	0.12	94,122,133,138	0
3	NAG	E	1	14/15	0.86	0.10	100,109,124,140	0
4	SIA	G	4	20/21	0.88	0.13	64,99,124,134	0
4	SIA	F	4	20/21	0.91	0.10	42,69,90,100	0
4	GAL	F	3	11/12	0.92	0.09	61,77,85,86	0

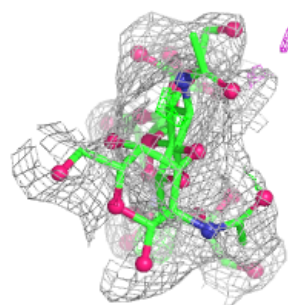
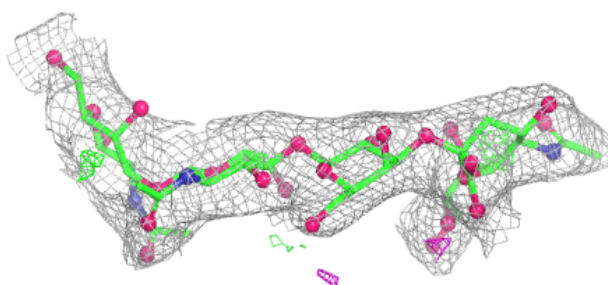
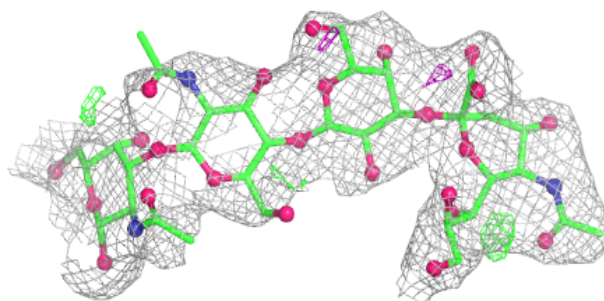
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

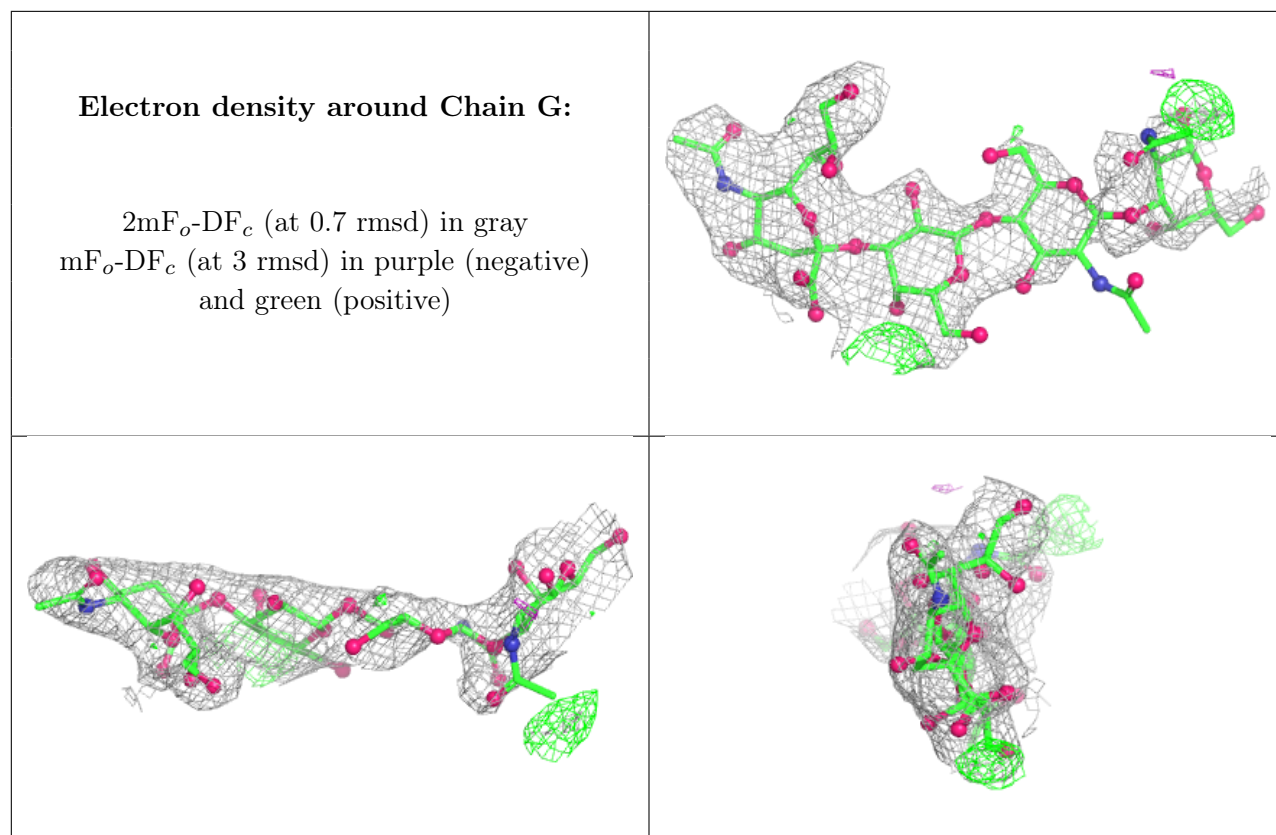
Electron density around Chain E:

$2mF_o-DF_c$ (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 [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($\text{\AA}^2$)	Q<0.9
5	NAG	B	201	14/15	0.72	0.14	86,108,112,113	0

6.5 Other polymers [i](#)

There are no such residues in this entry.