



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2021 – 10:03 AM EDT

PDB ID : 6XQ4
Title : Human antibody S8V2-47 in complex with the influenza hemagglutinin head domain of A/Beijing/262/1995(H1N1)
Authors : McCarthy, K.R.; Harrison, S.C.
Deposited on : 2020-07-09
Resolution : 3.35 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 : 2.18
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) : 2.18

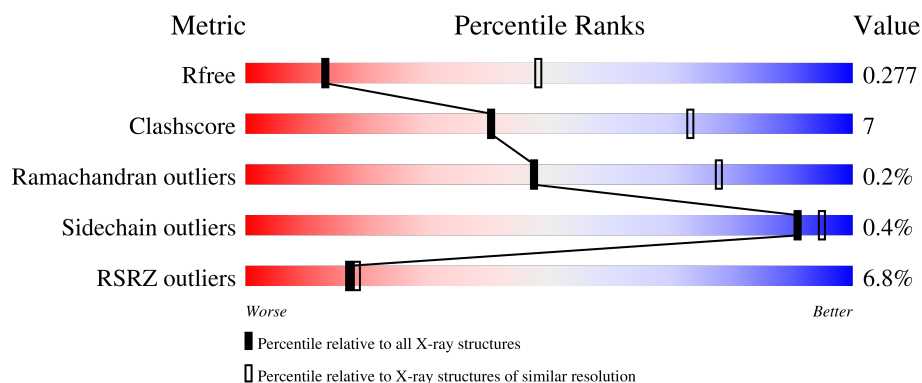
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.35 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	1060 (3.38-3.30)
Clashscore	141614	1111 (3.38-3.30)
Ramachandran outliers	138981	1090 (3.38-3.30)
Sidechain outliers	138945	1089 (3.38-3.30)
RSRZ outliers	127900	1028 (3.38-3.30)

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	223	<div> <div>4%</div> <div>78%</div> <div>16%</div> <div>6%</div> </div>
1	D	223	<div> <div>%</div> <div>80%</div> <div>15%</div> <div>5%</div> </div>
2	B	220	<div> <div>15%</div> <div>81%</div> <div>18%</div> <div>.</div> </div>
2	F	220	<div> <div>4%</div> <div>84%</div> <div>15%</div> <div>.</div> </div>
3	C	233	<div> <div>10%</div> <div>74%</div> <div>21%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
3	E	233	
4	G	2	
5	H	3	

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
5	NAG	H	2	-	-	-	X
5	BMA	H	3	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10189 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	209	Total	C	N	O	S	0	0	0
			1687	1077	289	317	4			
1	D	212	Total	C	N	O	S	0	0	0
			1711	1091	293	323	4			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	268	ALA	-	expression tag	UNP B4UPF7
A	269	LEU	-	expression tag	UNP B4UPF7
A	270	GLU	-	expression tag	UNP B4UPF7
A	271	VAL	-	expression tag	UNP B4UPF7
A	272	LEU	-	expression tag	UNP B4UPF7
A	273	PHE	-	expression tag	UNP B4UPF7
A	274	GLN	-	expression tag	UNP B4UPF7
D	268	ALA	-	expression tag	UNP B4UPF7
D	269	LEU	-	expression tag	UNP B4UPF7
D	270	GLU	-	expression tag	UNP B4UPF7
D	271	VAL	-	expression tag	UNP B4UPF7
D	272	LEU	-	expression tag	UNP B4UPF7
D	273	PHE	-	expression tag	UNP B4UPF7
D	274	GLN	-	expression tag	UNP B4UPF7

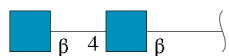
- Molecule 2 is a protein called antibody S8V2-47 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	0	0	0
			1685	1058	282	340	5			
2	F	218	Total	C	N	O	S	0	0	0
			1685	1058	282	340	5			

- Molecule 3 is a protein called antibody S8V2-47 heavy chain.

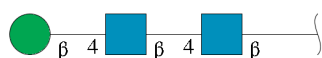
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	223	Total	C	N	O	S	0	0	0
			1677	1058	280	332	7			
3	E	223	Total	C	N	O	S	0	0	0
			1677	1058	280	332	7			

- Molecule 4 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
4	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

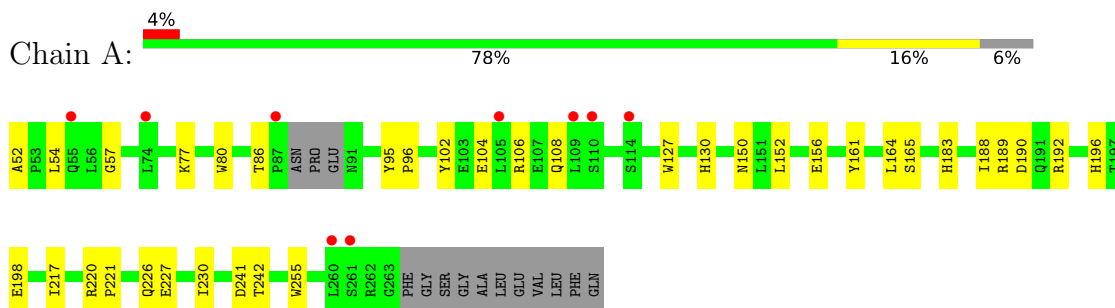


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	H	3	Total	C	N	O	0	0	0
			39	22	2	15			

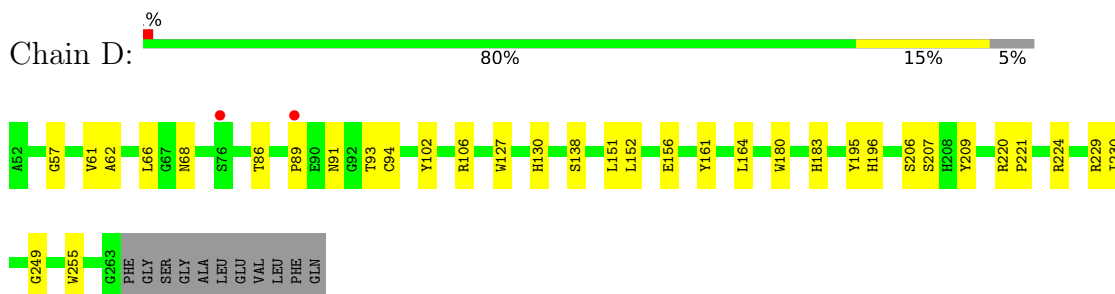
3 Residue-property plots [i](#)

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.

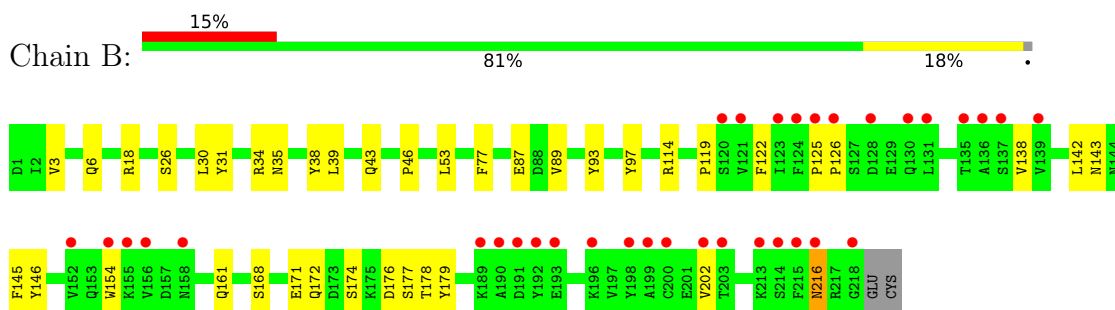
- Molecule 1: Hemagglutinin



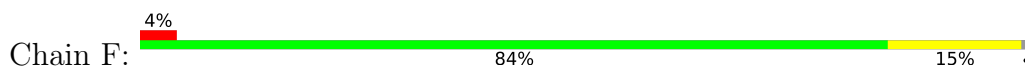
- Molecule 1: Hemagglutinin

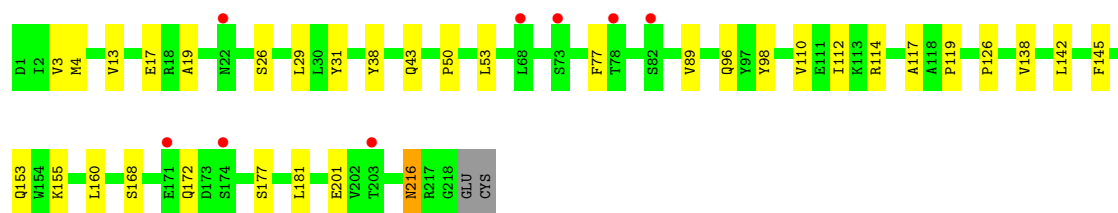


- Molecule 2: antibody S8V2-47 light chain

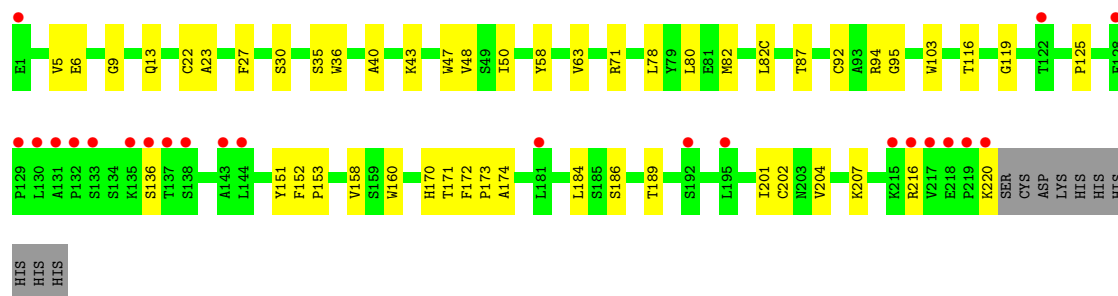
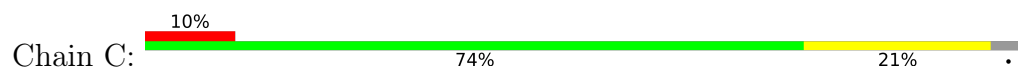


- Molecule 2: antibody S8V2-47 light chain

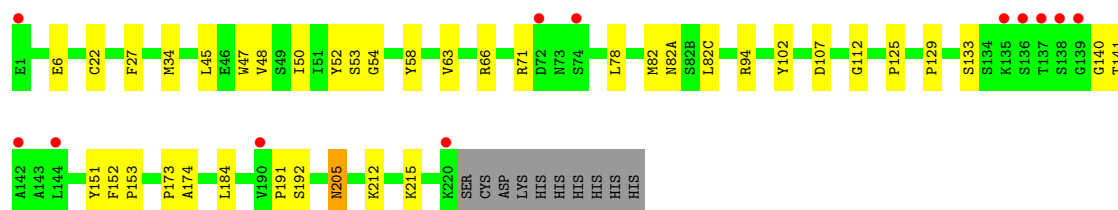
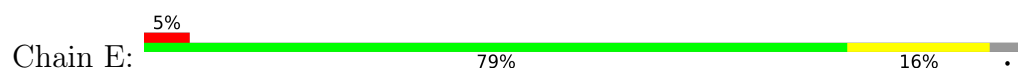




• Molecule 3: antibody S8V2-47 heavy chain



• Molecule 3: antibody S8V2-47 heavy chain



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



• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	96.09Å 66.47Å 135.18Å 90.00° 101.96° 90.00°	Depositor
Resolution (Å)	49.48 – 3.35 49.47 – 3.35	Depositor EDS
% Data completeness (in resolution range)	96.6 (49.48-3.35) 96.7 (49.47-3.35)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.238 , 0.279 0.237 , 0.277	Depositor DCC
R_{free} test set	1142 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	98.6	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	10189	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.58% 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: BMA, NAG

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	0.25	0/1738	0.47	0/2367
1	D	0.26	0/1764	0.47	0/2405
2	B	0.25	0/1721	0.45	0/2338
2	F	0.26	0/1721	0.49	0/2338
3	C	0.25	0/1719	0.46	0/2342
3	E	0.25	0/1719	0.46	0/2342
All	All	0.25	0/10382	0.47	0/14132

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	1687	0	1600	25	0
1	D	1711	0	1620	23	0
2	B	1685	0	1649	28	0
2	F	1685	0	1649	24	0
3	C	1677	0	1623	33	0
3	E	1677	0	1623	23	0
4	G	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	H	39	0	34	1	0
All	All	10189	0	9823	144	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 7.

All (144) 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:D:207:SER:HA	2:B:18:ARG:HH22	1.42	0.84
1:A:95:TYR:HE1	1:A:226:GLN:HG2	1.46	0.80
1:A:95:TYR:CE1	1:A:226:GLN:HG2	2.19	0.78
1:D:94:CYS:HA	1:D:224:ARG:HH12	1.47	0.78
1:A:220:ARG:HE	2:B:31:TYR:HE1	1.32	0.78
2:F:31:TYR:OH	2:F:98:TYR:O	2.02	0.76
1:D:68:ASN:ND2	1:D:91:ASN:O	2.20	0.75
2:F:114:ARG:HD2	2:F:177:SER:HB2	1.69	0.75
3:C:50:ILE:HD11	3:C:58:TYR:HD2	1.52	0.74
3:E:6:GLU:OE2	3:E:112:GLY:N	2.23	0.71
1:A:95:TYR:HD1	1:A:96:PRO:HD2	1.55	0.71
2:B:114:ARG:NH1	2:B:176:ASP:O	2.25	0.70
1:D:156:GLU:OE1	1:D:196:HIS:NE2	2.26	0.68
2:B:43:GLN:HB2	2:B:53:LEU:HD11	1.76	0.68
2:F:119:PRO:HB3	2:F:145:PHE:HD1	1.59	0.67
1:D:161:TYR:CE2	1:D:249:GLY:HA2	2.29	0.67
1:D:161:TYR:HE2	1:D:195:TYR:HD1	1.43	0.65
2:F:4:MET:HE2	2:F:96:GLN:HB3	1.79	0.65
2:B:89:VAL:HG11	2:B:172:GLN:HB3	1.79	0.64
1:A:57:GLY:O	1:A:86:THR:OG1	2.14	0.63
3:C:22:CYS:HB3	3:C:78:LEU:HB3	1.79	0.62
1:A:54:LEU:HG	1:A:80:TRP:CG	2.35	0.62
2:B:34:ARG:NH2	2:B:38:TYR:OH	2.28	0.61
2:B:114:ARG:HH22	2:B:178:THR:HG22	1.66	0.61
3:E:34:MET:HB3	3:E:78:LEU:HD22	1.83	0.61
3:E:125:PRO:HB3	3:E:151:TYR:HB3	1.83	0.60
2:B:125:PRO:HG2	3:C:220:LYS:HE3	1.83	0.60
2:F:168:SER:HB2	3:E:173:PRO:HG2	1.83	0.60
1:A:192:ARG:NH2	1:A:198:GLU:OE2	2.35	0.59
2:B:168:SER:HB2	3:C:173:PRO:HG2	1.83	0.58
3:E:174:ALA:HB2	3:E:184:LEU:HD23	1.84	0.58
1:D:61:VAL:HG11	1:D:106:ARG:HG2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:LEU:HG	1:A:255:TRP:HB2	1.84	0.57
3:C:9:GLY:O	3:C:207:LYS:NZ	2.35	0.57
2:B:126:PRO:HD3	2:B:138:VAL:HG22	1.87	0.57
3:C:201:ILE:HG12	3:C:216:ARG:HG2	1.86	0.57
2:F:172:GLN:HE21	2:F:177:SER:HB3	1.70	0.56
1:A:127:TRP:CZ3	1:A:164:LEU:HD21	2.42	0.55
3:E:141:THR:HA	3:E:191:PRO:HA	1.89	0.55
2:F:153:GLN:HB3	2:F:201:GLU:HB3	1.89	0.55
3:E:53:SER:O	3:E:71:ARG:NH2	2.39	0.55
1:D:161:TYR:HD2	1:D:195:TYR:HA	1.72	0.55
1:D:152:LEU:HG	1:D:255:TRP:HB2	1.87	0.54
3:C:13:GLN:HG3	3:C:119:GLY:HA2	1.88	0.54
1:D:161:TYR:CE2	1:D:195:TYR:HD1	2.24	0.54
1:D:180:TRP:HH2	1:D:209:TYR:HE2	1.56	0.54
1:A:164:LEU:HD22	1:A:165:SER:H	1.73	0.53
2:B:119:PRO:HB3	2:B:145:PHE:HD1	1.72	0.53
3:C:50:ILE:HD11	3:C:58:TYR:CD2	2.39	0.52
2:B:216:ASN:O	2:B:216:ASN:ND2	2.41	0.52
3:E:205:ASN:OD1	3:E:212:LYS:NZ	2.43	0.52
2:B:172:GLN:HE21	2:B:177:SER:HB3	1.75	0.52
3:C:158:VAL:HG22	3:C:204:VAL:HG22	1.91	0.51
3:E:129:PRO:HD3	3:E:215:LYS:HE2	1.92	0.51
2:F:126:PRO:HD3	2:F:138:VAL:HG22	1.91	0.51
3:E:94:ARG:O	3:E:107:ASP:N	2.43	0.51
3:C:87:THR:HG23	3:C:116:THR:HA	1.92	0.51
3:E:53:SER:OG	3:E:54:GLY:N	2.42	0.51
2:B:122:PHE:HB3	3:C:136:SER:HA	1.93	0.51
2:B:6:GLN:HE22	2:B:93:TYR:HA	1.74	0.50
3:C:6:GLU:OE2	3:C:92:CYS:N	2.43	0.50
1:D:130:HIS:CD2	1:D:161:TYR:CD1	3.00	0.50
3:C:82:MET:HB3	3:C:82(C):LEU:HD21	1.93	0.50
3:C:125:PRO:HB3	3:C:151:TYR:HB3	1.92	0.50
2:F:13:VAL:HG21	2:F:19:ALA:HB2	1.93	0.50
3:E:48:VAL:HG13	3:E:63:VAL:HG21	1.93	0.50
1:A:221:PRO:HD3	3:C:103:TRP:CE2	2.46	0.50
1:D:220:ARG:HH11	1:D:229:ARG:HG3	1.76	0.50
1:A:156:GLU:OE1	1:A:196:HIS:NE2	2.44	0.49
2:B:142:LEU:HD21	2:B:202:VAL:HG21	1.94	0.49
3:E:140:GLY:HA2	3:E:192:SER:OG	2.11	0.49
2:B:30:LEU:HD11	2:B:35:ASN:HA	1.94	0.49
3:C:170:HIS:HB3	3:C:172:PHE:HE1	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:174:ALA:HA	3:C:184:LEU:HB3	1.93	0.49
1:D:180:TRP:CH2	1:D:209:TYR:HE2	2.30	0.49
3:C:35:SER:OG	3:C:50:ILE:HG22	2.13	0.48
3:E:82:MET:HB3	3:E:82(C):LEU:HD21	1.94	0.48
2:F:50:PRO:HG2	3:E:45:LEU:HD21	1.96	0.47
3:E:47:TRP:HZ2	3:E:50:ILE:HG22	1.79	0.47
2:B:46:PRO:HB3	2:B:171:GLU:OE1	2.15	0.47
2:F:29:LEU:HD12	2:F:77:PHE:CE2	2.50	0.47
1:A:189:ARG:NH1	1:A:190:ASP:OD1	2.48	0.47
3:C:48:VAL:HG13	3:C:63:VAL:HG21	1.96	0.46
2:F:4:MET:CE	2:F:96:GLN:HB3	2.45	0.46
2:B:154:TRP:HB2	2:B:161:GLN:HB2	1.97	0.46
1:A:188:ILE:HA	1:A:217:ILE:HD13	1.99	0.45
1:A:220:ARG:HB3	1:A:221:PRO:HD2	1.98	0.45
2:F:114:ARG:HH22	2:F:117:ALA:HB2	1.81	0.45
1:A:227:GLU:HG2	3:C:58:TYR:OH	2.17	0.45
1:D:127:TRP:HZ3	1:D:164:LEU:HD11	1.82	0.45
3:C:172:PHE:N	3:C:172:PHE:CD1	2.85	0.45
2:F:119:PRO:HB3	2:F:145:PHE:CD1	2.46	0.45
3:C:94:ARG:HG2	3:C:95:GLY:N	2.32	0.45
2:B:114:ARG:HH12	2:B:178:THR:HG22	1.82	0.45
2:F:13:VAL:HB	2:F:17:GLU:HG2	1.99	0.44
3:E:6:GLU:HA	3:E:22:CYS:HA	1.99	0.44
3:E:66:ARG:HB3	3:E:82(A):ASN:O	2.17	0.44
1:A:130:HIS:CE1	1:A:164:LEU:HD23	2.53	0.44
2:B:172:GLN:HB2	2:B:179:TYR:CZ	2.52	0.44
2:F:114:ARG:O	2:F:114:ARG:HG3	2.17	0.44
2:F:155:LYS:HG2	2:F:160:LEU:HD23	2.00	0.44
3:C:47:TRP:CZ2	3:C:50:ILE:HG23	2.53	0.44
1:D:206:SER:O	2:B:18:ARG:NH2	2.51	0.43
2:B:143:ASN:ND2	3:C:189:THR:HG21	2.34	0.43
3:C:36:TRP:NE1	3:C:80:LEU:HB2	2.33	0.43
2:B:39:LEU:HD13	2:B:77:PHE:CG	2.53	0.43
2:B:114:ARG:HH21	2:B:146:TYR:HB2	1.83	0.43
2:F:112:ILE:H	2:F:172:GLN:HE22	1.67	0.43
1:A:52:ALA:HB1	1:A:77:LYS:NZ	2.33	0.43
3:E:152:PHE:HA	3:E:153:PRO:HA	1.84	0.43
1:D:57:GLY:O	1:D:86:THR:HB	2.19	0.43
1:D:183:HIS:HA	1:D:230:ILE:HG12	2.00	0.43
2:B:3:VAL:H	2:B:26:SER:HB3	1.82	0.43
3:E:52:TYR:OH	3:E:58:TYR:HE2	2.02	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:1:NAG:H4	5:H:2:NAG:N2	2.32	0.43
2:B:97:TYR:O	3:C:103:TRP:HB2	2.18	0.43
3:E:27:PHE:CZ	3:E:94:ARG:HG3	2.54	0.43
3:C:5:VAL:O	3:C:23:ALA:N	2.47	0.42
3:C:30:SER:HA	3:C:71:ARG:CZ	2.50	0.42
1:A:102:TYR:CE2	1:A:106:ARG:HD2	2.53	0.42
1:D:86:THR:HG21	1:D:89:PRO:HB3	2.00	0.42
1:A:164:LEU:HD22	1:A:165:SER:N	2.34	0.42
2:F:38:TYR:HE2	3:E:102:TYR:HB3	1.84	0.42
3:C:40:ALA:HB3	3:C:43:LYS:HB2	2.01	0.42
1:A:54:LEU:HG	1:A:80:TRP:CD2	2.55	0.42
1:D:66:LEU:HD22	1:D:151:LEU:HD11	2.02	0.42
3:C:171:THR:HG23	3:C:186:SER:HB2	2.02	0.42
1:A:241:ASP:OD1	1:A:242:THR:N	2.53	0.41
3:C:27:PHE:CE1	3:C:94:ARG:HD2	2.55	0.41
3:C:160:TRP:CH2	3:C:202:CYS:HB3	2.55	0.41
3:C:152:PHE:HA	3:C:153:PRO:HA	1.80	0.41
2:F:3:VAL:H	2:F:26:SER:HB2	1.86	0.41
2:F:89:VAL:HA	2:F:110:VAL:HG23	2.03	0.41
1:A:183:HIS:HA	1:A:230:ILE:HG12	2.01	0.41
2:F:142:LEU:HB2	2:F:181:LEU:HB3	2.01	0.41
1:A:104:GLU:O	1:A:108:GLN:HG2	2.21	0.41
1:D:62:ALA:HB2	1:D:102:TYR:CE1	2.56	0.41
2:F:43:GLN:HB2	2:F:53:LEU:HD11	2.02	0.41
1:A:150:ASN:C	1:A:255:TRP:HB3	2.40	0.40
1:D:220:ARG:HB3	1:D:221:PRO:HD2	2.03	0.40
3:E:174:ALA:HA	3:E:184:LEU:HB3	2.02	0.40
1:D:138:SER:O	1:D:224:ARG:NH2	2.54	0.40
2:B:87:GLU:HA	2:B:174:SER:HB3	2.03	0.40
2:F:216:ASN:O	2:F:216:ASN:ND2	2.54	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	205/223 (92%)	195 (95%)	10 (5%)	0	100	100
1	D	210/223 (94%)	200 (95%)	9 (4%)	1 (0%)	29	63
2	B	216/220 (98%)	203 (94%)	13 (6%)	0	100	100
2	F	216/220 (98%)	204 (94%)	12 (6%)	0	100	100
3	C	221/233 (95%)	207 (94%)	14 (6%)	0	100	100
3	E	221/233 (95%)	205 (93%)	15 (7%)	1 (0%)	29	63
All	All	1289/1352 (95%)	1214 (94%)	73 (6%)	2 (0%)	47	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	133	SER
1	D	93	THR

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	184/195 (94%)	183 (100%)	1 (0%)	88	93
1	D	187/195 (96%)	187 (100%)	0	100	100
2	B	192/194 (99%)	191 (100%)	1 (0%)	88	93
2	F	192/194 (99%)	191 (100%)	1 (0%)	88	93
3	C	185/195 (95%)	185 (100%)	0	100	100
3	E	185/195 (95%)	184 (100%)	1 (0%)	88	93
All	All	1125/1168 (96%)	1121 (100%)	4 (0%)	91	95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	TYR
2	B	216	ASN
2	F	216	ASN
3	E	205	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	130	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 ⓘ

5 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	$\# Z > 2$
4	NAG	G	1	1,4	14,14,15	1.18	1 (7%)	17,19,21	1.57	1 (5%)
4	NAG	G	2	4	14,14,15	0.45	0	17,19,21	0.50	0
5	NAG	H	1	5,1	14,14,15	0.32	0	17,19,21	0.49	0
5	NAG	H	2	5	14,14,15	0.50	0	17,19,21	0.51	0
5	BMA	H	3	5	11,11,12	0.61	0	15,15,17	0.82	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
4	NAG	G	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
5	NAG	H	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	H	2	5	-	0/6/23/26	0/1/1/1
5	BMA	H	3	5	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	1	NAG	O5-C1	3.97	1.50	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	1	NAG	C1-O5-C5	6.03	120.36	112.19

There are no chirality outliers.

All (7) torsion outliers are listed below:

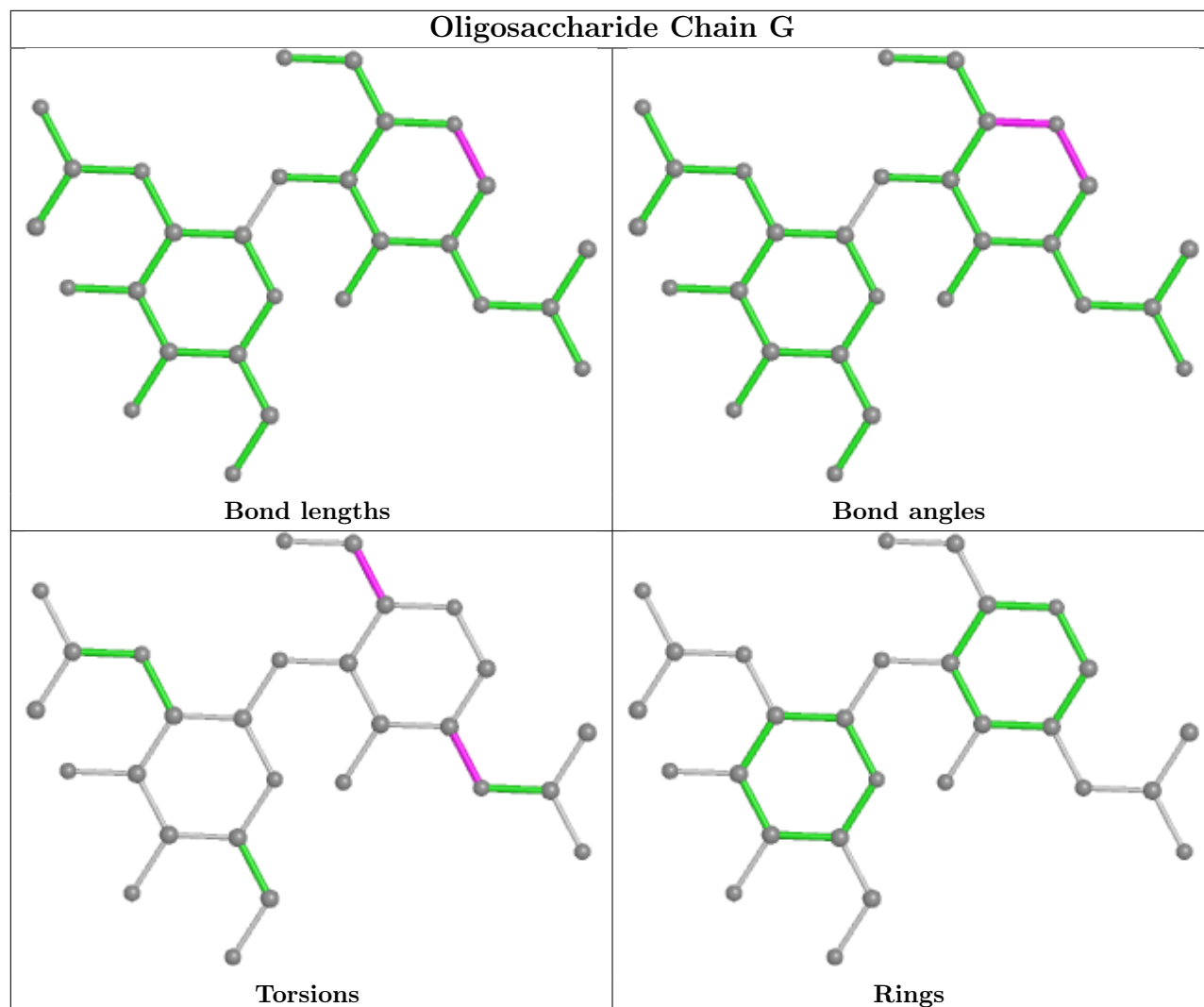
Mol	Chain	Res	Type	Atoms
5	H	1	NAG	O5-C5-C6-O6
5	H	1	NAG	C8-C7-N2-C2
5	H	1	NAG	O7-C7-N2-C2
4	G	1	NAG	C1-C2-N2-C7
5	H	1	NAG	C4-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	G	1	NAG	C3-C2-N2-C7

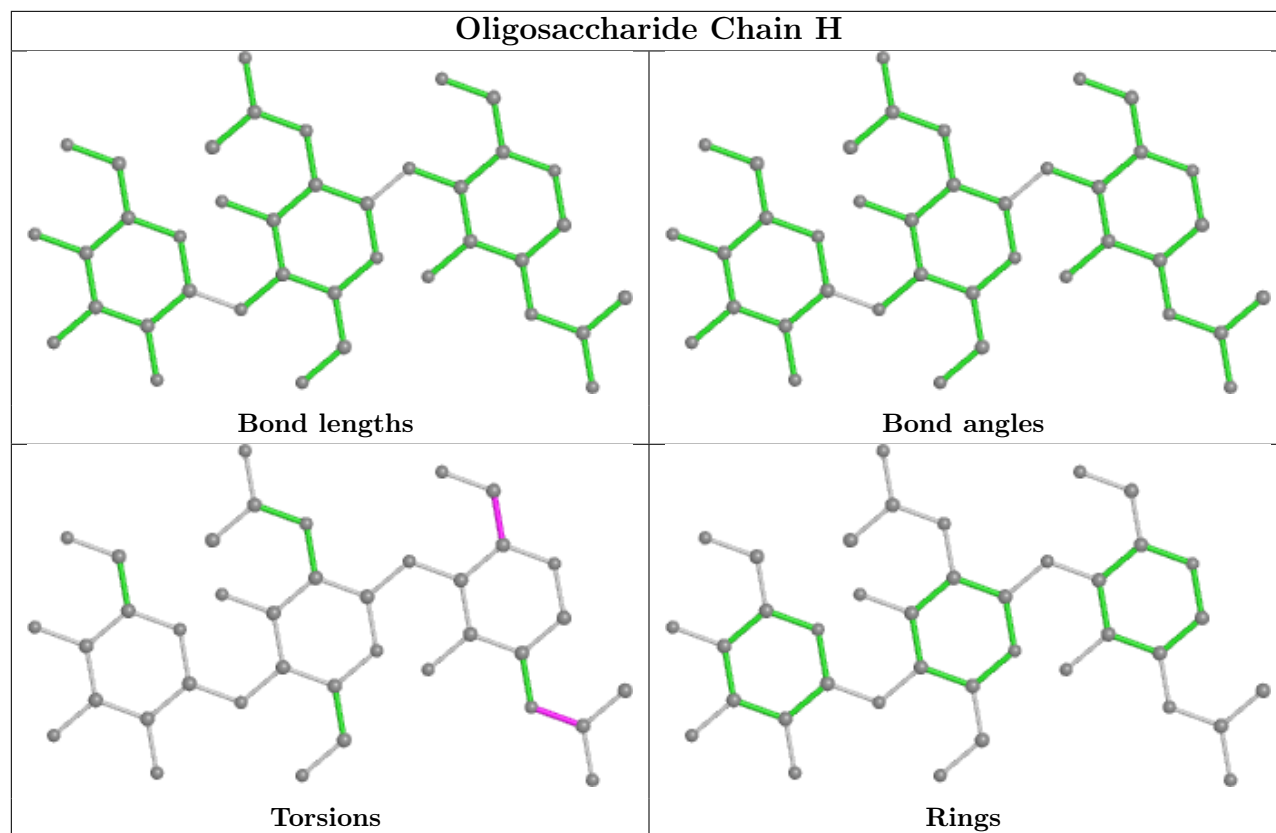
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	H	1	NAG	1	0
5	H	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](#)

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	209/223 (93%)	0.42	9 (4%) 35 36	67, 92, 138, 168	0
1	D	212/223 (95%)	0.33	2 (0%) 84 85	61, 88, 128, 141	0
2	B	218/220 (99%)	0.87	34 (15%) 2 1	61, 115, 188, 207	0
2	F	218/220 (99%)	0.39	8 (3%) 41 40	64, 96, 124, 173	0
3	C	223/233 (95%)	0.87	23 (10%) 6 6	59, 103, 184, 226	0
3	E	223/233 (95%)	0.36	12 (5%) 25 26	61, 92, 122, 157	0
All	All	1303/1352 (96%)	0.54	88 (6%) 17 18	59, 95, 164, 226	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	138	SER	10.2
3	C	137	THR	6.7
3	C	136	SER	6.3
2	B	198	TYR	5.9
3	C	132	PRO	4.9
2	B	191	ASP	4.8
3	C	131	ALA	4.8
3	E	135	LYS	4.8
2	B	158	ASN	4.7
2	B	135	THR	4.6
3	C	129	PRO	4.5
3	C	144	LEU	4.4
2	B	189	LYS	4.4
3	C	220	LYS	4.1
3	C	135	LYS	4.0
2	B	192	TYR	4.0
2	B	193	GLU	4.0
2	B	120	SER	3.8
2	B	190	ALA	3.8

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Mol	Chain	Res	Type	RSRZ
2	B	124	PHE	3.8
2	B	214	SER	3.8
2	B	131	LEU	3.8
2	B	199	ALA	3.7
3	C	1	GLU	3.7
2	F	22	ASN	3.7
3	C	218	GLU	3.6
1	A	261	SER	3.6
2	B	123	ILE	3.5
3	C	219	PRO	3.5
2	B	218	GLY	3.4
2	B	200	CYS	3.4
2	F	174	SER	3.4
3	E	137	THR	3.4
3	C	130	LEU	3.3
1	A	110	SER	3.2
3	E	136	SER	3.2
3	C	217	VAL	3.2
3	E	74	SER	3.1
1	A	114	SER	3.1
2	B	202	VAL	3.1
2	B	216	ASN	3.0
1	A	55	GLN	3.0
1	D	76	SER	3.0
3	C	128	PHE	3.0
3	C	143	ALA	2.9
2	B	215	PHE	2.9
2	B	196	LYS	2.8
2	B	125	PRO	2.8
3	C	192	SER	2.8
2	B	156	VAL	2.7
1	A	105	LEU	2.7
1	A	74	LEU	2.7
2	B	152	VAL	2.6
2	B	203	THR	2.6
3	C	215	LYS	2.6
3	C	195	LEU	2.5
1	A	109	LEU	2.5
3	C	133	SER	2.5
1	A	87	PRO	2.5
3	C	181	LEU	2.4
2	B	130	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	154	TRP	2.4
3	E	139	GLY	2.4
2	F	203	THR	2.4
2	B	137	SER	2.3
2	F	73	SER	2.3
3	C	216	ARG	2.3
1	D	89	PRO	2.3
2	F	171	GLU	2.3
2	F	82	SER	2.3
2	B	139	VAL	2.2
3	E	190	VAL	2.2
3	C	122	THR	2.2
2	B	121	VAL	2.2
2	B	136	ALA	2.1
2	F	78	THR	2.1
3	E	1	GLU	2.1
3	E	220	LYS	2.1
3	E	144	LEU	2.1
2	B	155	LYS	2.1
3	E	142	ALA	2.1
2	B	126	PRO	2.1
2	B	213	LYS	2.0
1	A	260	LEU	2.0
2	F	68	LEU	2.0
3	E	138	SER	2.0
2	B	128	ASP	2.0
3	E	72	ASP	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.

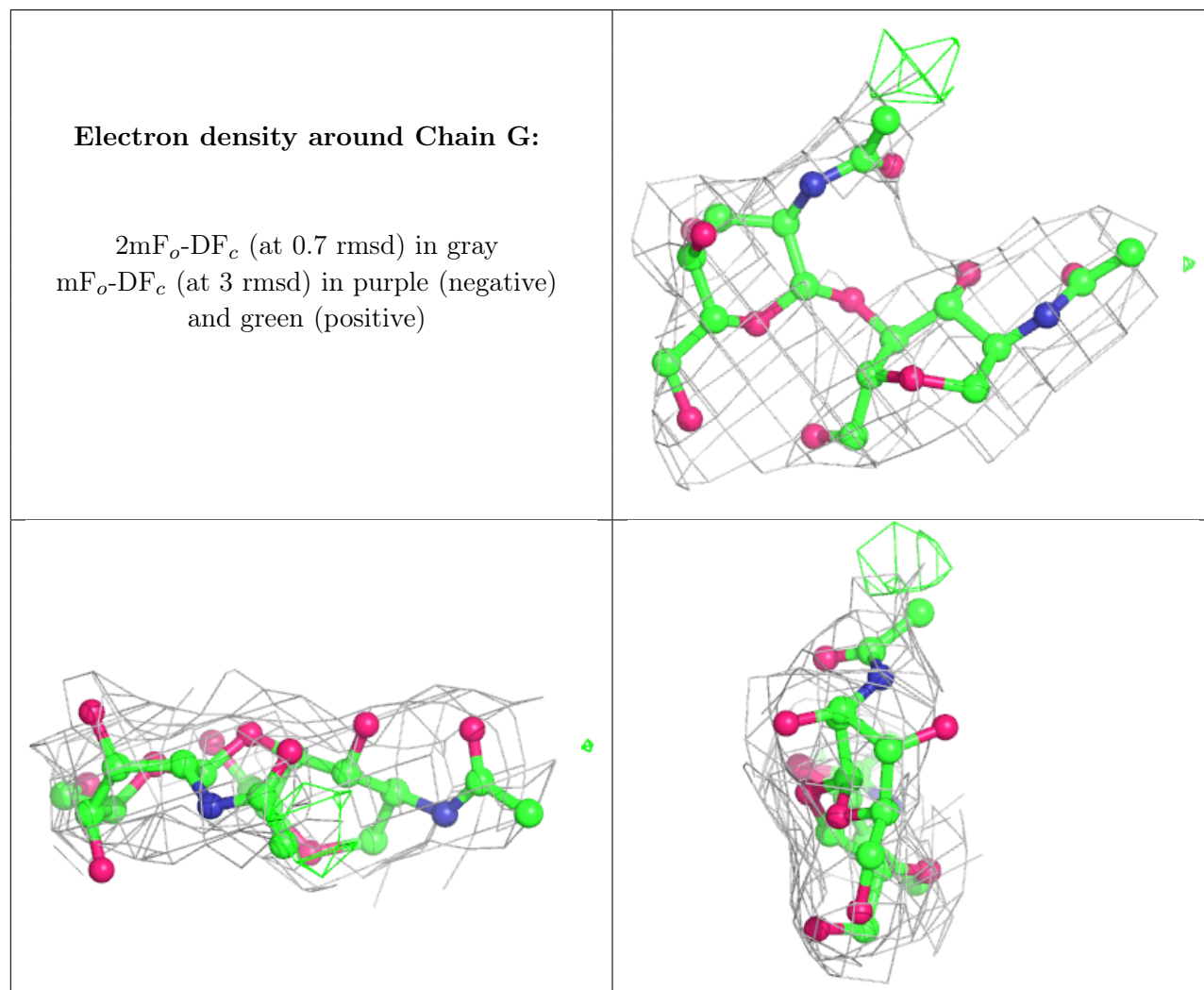
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	NAG	H	1	14/15	0.63	0.37	112,160,170,195	0

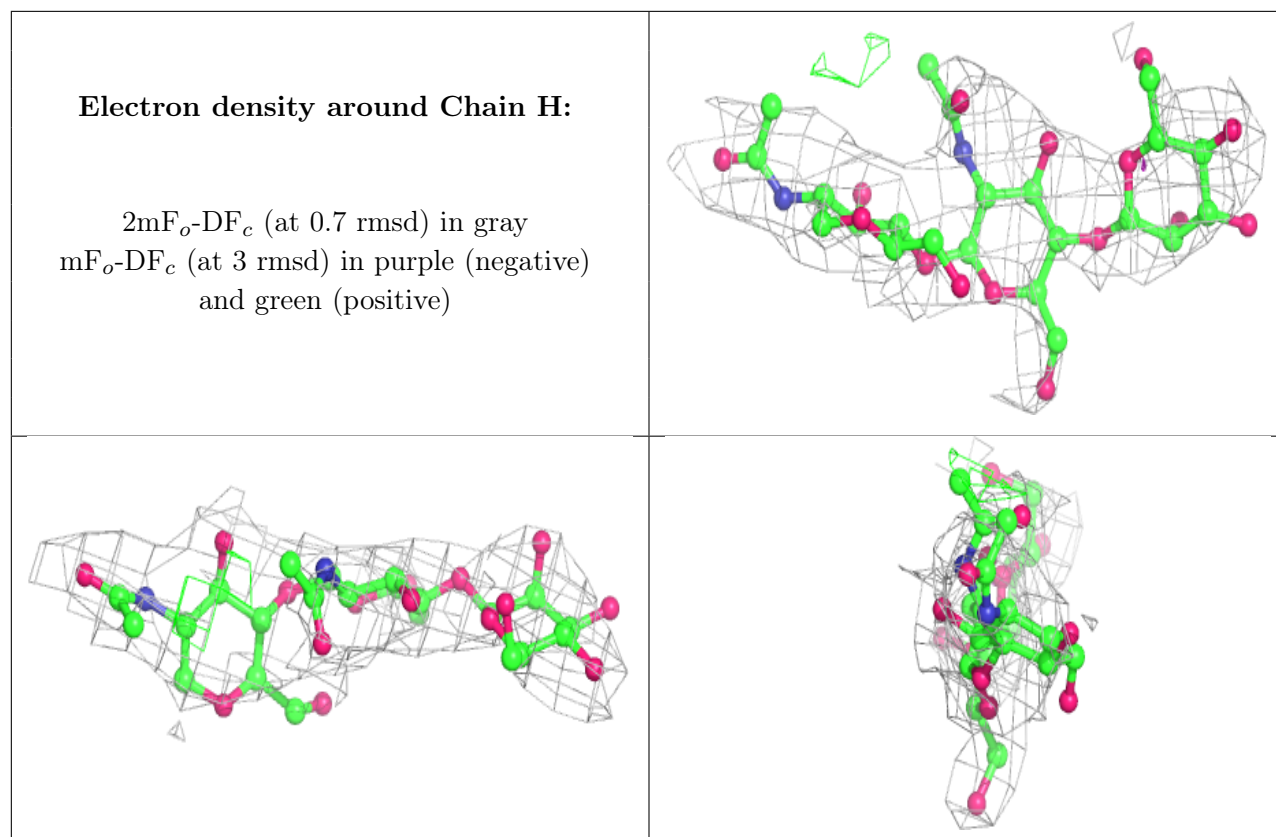
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	BMA	H	3	11/12	0.72	0.59	104,138,151,160	0
5	NAG	H	2	14/15	0.76	0.41	132,163,171,172	0
4	NAG	G	2	14/15	0.79	0.25	121,157,162,168	0
4	NAG	G	1	14/15	0.84	0.16	114,143,156,160	0

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.





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.