



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 29, 2022 – 12:10 AM EST

PDB ID : 8DCY
Title : CCHFV GP38 Hoti/Kosovo bound with 13G8 Fab
Authors : Durie, I.A.; Bergeron, E.; Pegan, S.D.; McGuire, J.
Deposited on : 2022-06-17
Resolution : 3.62 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
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) : 2.31.2

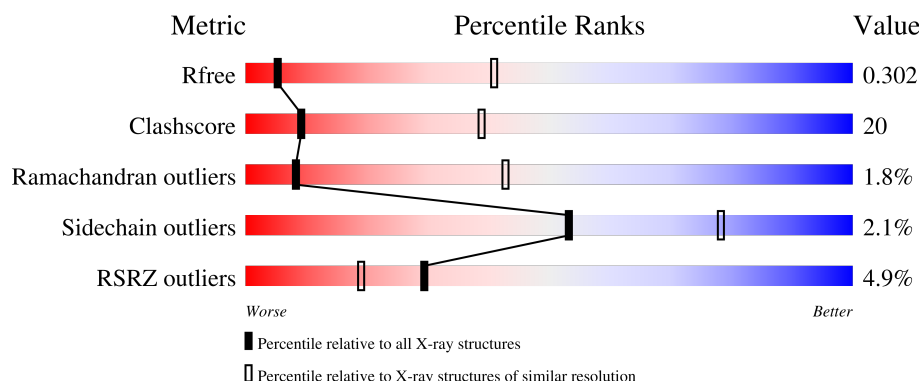
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.62 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	1290 (3.74-3.50)
Clashscore	141614	1387 (3.74-3.50)
Ramachandran outliers	138981	1339 (3.74-3.50)
Sidechain outliers	138945	1339 (3.74-3.50)
RSRZ outliers	127900	1191 (3.74-3.50)

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	233	
2	B	213	
3	C	274	
4	D	2	
4	G	2	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5109 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 13G8 Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	0	0
			1525	963	257	297	8			

- Molecule 2 is a protein called 13G8 Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	208	Total	C	N	O	S	0	0	0
			1595	1004	265	319	7			

- Molecule 3 is a protein called Envelopment polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	241	Total	C	N	O	S	0	0	0
			1930	1234	332	354	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	520	LEU	-	expression tag	UNP A0A7T6Y557
C	521	GLU	-	expression tag	UNP A0A7T6Y557
C	522	VAL	-	expression tag	UNP A0A7T6Y557
C	523	LEU	-	expression tag	UNP A0A7T6Y557
C	524	PHE	-	expression tag	UNP A0A7T6Y557
C	525	GLN	-	expression tag	UNP A0A7T6Y557

- 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			
4	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

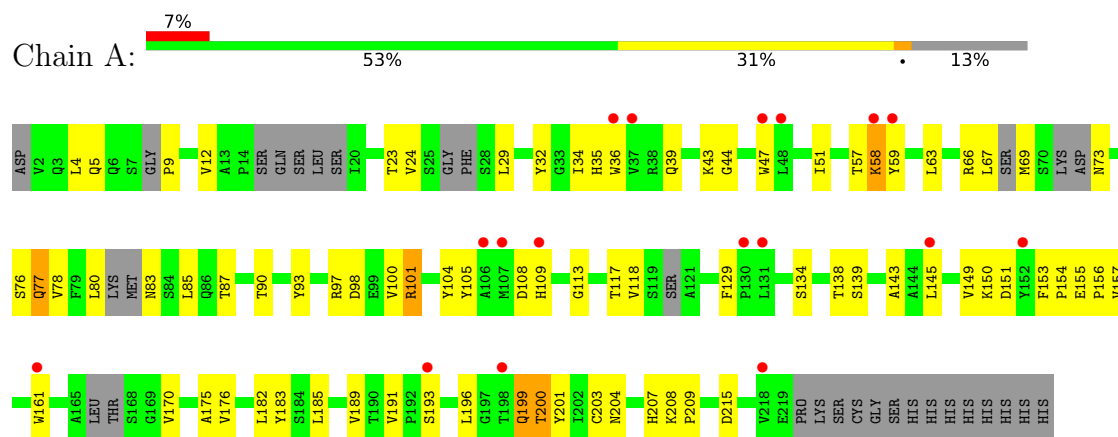
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	O	0	0
			1	1		
5	B	1	Total	O	0	0
			1	1		
5	C	1	Total	O	0	0
			1	1		

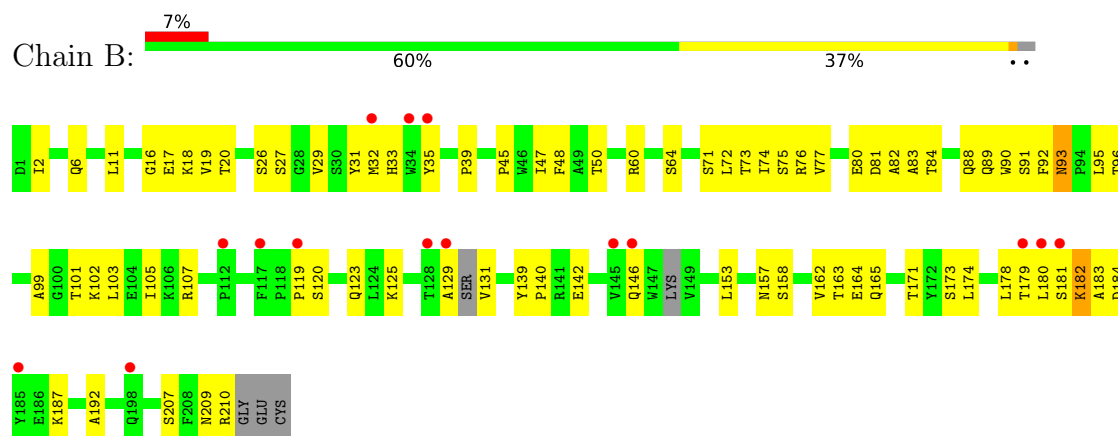
3 Residue-property plots

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.

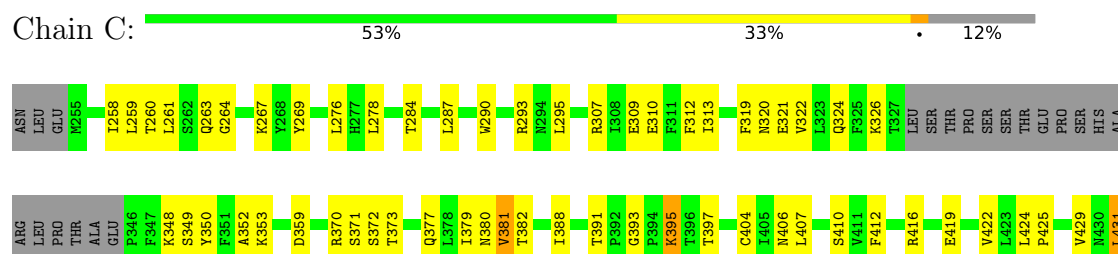
• Molecule 1: 13G8 Heavy Chain

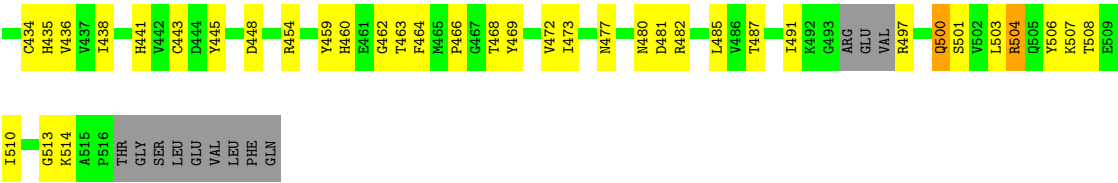


• Molecule 2: 13G8 Light Chain



• Molecule 3: Envelopment polypeptide





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



• Molecule 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 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	175.92Å 175.92Å 93.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.06 – 3.62 49.64 – 3.62	Depositor EDS
% Data completeness (in resolution range)	98.3 (45.06-3.62) 99.7 (49.64-3.62)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 3.57Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.281 , 0.299 0.282 , 0.302	Depositor DCC
R_{free} test set	818 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	138.7	Xtriage
Anisotropy	0.561	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 134.1	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	0.89	EDS
Total number of atoms	5109	wwPDB-VP
Average B, all atoms (Å ²)	184.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.55% 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: 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	0.37	0/1556	0.82	1/2111 (0.0%)
2	B	0.33	0/1634	0.71	0/2221
3	C	0.35	0/1967	0.75	1/2650 (0.0%)
All	All	0.35	0/5157	0.76	2/6982 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	145	LEU	CA-CB-CG	9.36	136.83	115.30
3	C	504	ARG	CB-CG-CD	6.38	128.19	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	93	ASN	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1525	0	1485	64	0
2	B	1595	0	1533	73	0
3	C	1930	0	1954	77	0
4	D	28	0	25	3	0
4	G	28	0	25	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
All	All	5109	0	5022	203	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 20.

All (203) 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:157:ASN:HD21	2:B:179:THR:H	1.16	0.91
1:A:150:LYS:NZ	1:A:151:ASP:OD1	2.08	0.85
3:C:259:LEU:HD11	3:C:267:LYS:HE3	1.59	0.84
1:A:35:HIS:HD1	1:A:47:TRP:HH2	1.25	0.81
2:B:6:GLN:HE22	2:B:99:ALA:HB3	1.45	0.81
1:A:93:TYR:O	1:A:113:GLY:HA2	1.82	0.80
3:C:307:ARG:NH1	3:C:310:GLU:OE1	2.14	0.80
2:B:60:ARG:NH2	2:B:80:GLU:OE2	2.18	0.77
3:C:391:THR:HB	3:C:504:ARG:HD3	1.70	0.74
2:B:157:ASN:HD21	2:B:179:THR:N	1.86	0.74
2:B:32:MET:HA	2:B:89:GLN:NE2	2.04	0.73
3:C:259:LEU:HD22	3:C:263:GLN:HB3	1.68	0.73
2:B:33:HIS:HB3	2:B:48:PHE:HA	1.71	0.72
3:C:382:THR:HG22	3:C:463:THR:H	1.53	0.72
2:B:119:PRO:HD3	2:B:131:VAL:HG22	1.70	0.71
3:C:395:LYS:HD3	3:C:395:LYS:H	1.57	0.69
3:C:377:GLN:HB3	3:C:441:HIS:CE1	2.27	0.69
2:B:18:LYS:HG2	2:B:75:SER:HA	1.75	0.69
2:B:89:GLN:HB3	2:B:92:PHE:CZ	2.27	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:ASP:HA	2:B:45:PRO:HG3	1.76	0.68
2:B:120:SER:OG	2:B:123:GLN:HG3	1.94	0.68
2:B:35:TYR:HD1	2:B:45:PRO:HA	1.59	0.67
3:C:422:VAL:O	3:C:468:THR:OG1	2.08	0.67
1:A:59:TYR:HB3	1:A:63:LEU:HD21	1.76	0.66
1:A:36:TRP:NE1	1:A:69:MET:SD	2.68	0.66
1:A:29:LEU:HD12	1:A:73:ASN:HA	1.77	0.65
3:C:406:ASN:HA	3:C:424:LEU:HD23	1.79	0.65
3:C:473:ILE:HG23	3:C:480:ASN:HD22	1.63	0.63
2:B:142:GLU:OE1	2:B:142:GLU:N	2.26	0.63
2:B:179:THR:HG23	2:B:180:LEU:HG	1.79	0.63
3:C:380:ASN:O	3:C:381:VAL:HB	1.98	0.61
3:C:429:VAL:HG23	3:C:487:THR:HB	1.80	0.61
3:C:406:ASN:OD1	3:C:424:LEU:HB2	2.01	0.61
2:B:18:LYS:HA	2:B:74:ILE:O	2.01	0.61
3:C:464:PHE:CE2	4:D:1:NAG:H82	2.36	0.61
1:A:51:ILE:CD1	1:A:57:THR:HG22	2.31	0.61
2:B:35:TYR:CD1	2:B:45:PRO:HA	2.37	0.60
2:B:91:SER:H	3:C:258:ILE:HG22	1.66	0.60
2:B:95:LEU:HD12	2:B:96:THR:H	1.66	0.60
3:C:416:ARG:HE	3:C:514:LYS:HE2	1.66	0.60
2:B:84:THR:HA	2:B:101:THR:O	2.02	0.59
2:B:17:GLU:O	2:B:77:VAL:HG23	2.02	0.59
3:C:435:HIS:O	3:C:480:ASN:ND2	2.27	0.59
2:B:82:ALA:HA	2:B:103:LEU:O	2.03	0.59
3:C:473:ILE:HG23	3:C:480:ASN:ND2	2.17	0.58
2:B:2:ILE:HG12	2:B:92:PHE:CE1	2.38	0.58
1:A:51:ILE:HD11	1:A:69:MET:HG2	1.86	0.58
2:B:182:LYS:HD3	2:B:182:LYS:N	2.19	0.58
1:A:129:PHE:CG	2:B:123:GLN:HG2	2.39	0.57
1:A:153:PHE:HB2	1:A:182:LEU:HD23	1.87	0.57
3:C:438:ILE:HD11	3:C:445:TYR:CZ	2.40	0.57
1:A:100:VAL:HG13	1:A:104:TYR:HE1	1.70	0.56
1:A:117:THR:N	1:A:155:GLU:OE2	2.37	0.56
2:B:33:HIS:CE1	2:B:88:GLN:HG2	2.41	0.56
2:B:64:SER:HG	2:B:71:SER:HG	1.52	0.56
1:A:51:ILE:HD12	1:A:57:THR:HG22	1.88	0.56
2:B:91:SER:H	3:C:258:ILE:CG2	2.19	0.56
2:B:2:ILE:HG21	2:B:92:PHE:CG	2.40	0.55
3:C:431:LEU:HD12	3:C:434:CYS:HB3	1.88	0.55
3:C:321:GLU:OE2	3:C:349:SER:OG	2.23	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:481:ASP:O	3:C:513:GLY:HA2	2.07	0.55
2:B:32:MET:SD	2:B:89:GLN:NE2	2.78	0.55
3:C:393:GLY:HA3	3:C:506:TYR:CD2	2.42	0.55
3:C:445:TYR:OH	3:C:454:ARG:O	2.17	0.55
1:A:149:VAL:HG11	1:A:157:VAL:HG21	1.89	0.54
3:C:435:HIS:N	3:C:480:ASN:OD1	2.40	0.54
1:A:24:VAL:HG22	1:A:76:SER:O	2.08	0.54
3:C:422:VAL:HG21	3:C:469:TYR:CZ	2.43	0.53
1:A:39:GLN:HG3	1:A:44:GLY:O	2.09	0.53
3:C:377:GLN:HG2	3:C:441:HIS:CG	2.44	0.53
3:C:436:VAL:HG12	3:C:473:ILE:HG12	1.89	0.53
4:D:1:NAG:O3	4:D:2:NAG:N2	2.42	0.53
1:A:12:VAL:HG21	1:A:85:LEU:HD13	1.90	0.53
1:A:73:ASN:N	1:A:77:GLN:H	2.07	0.52
1:A:150:LYS:HG2	1:A:151:ASP:H	1.74	0.52
2:B:2:ILE:HD11	2:B:27:SER:H	1.74	0.52
2:B:91:SER:HA	3:C:258:ILE:HG22	1.91	0.52
1:A:170:VAL:HG22	1:A:189:VAL:HG22	1.91	0.52
1:A:138:THR:O	1:A:139:SER:OG	2.27	0.51
3:C:312:PHE:O	3:C:313:ILE:HD13	2.10	0.51
2:B:2:ILE:HD12	2:B:26:SER:OG	2.10	0.51
2:B:20:THR:HA	2:B:72:LEU:O	2.11	0.51
2:B:32:MET:HA	2:B:89:GLN:HE22	1.76	0.51
3:C:410:SER:HG	3:C:412:PHE:HE1	1.57	0.51
2:B:31:TYR:O	2:B:89:GLN:CD	2.50	0.51
2:B:183:ALA:O	2:B:187:LYS:HG2	2.11	0.50
1:A:153:PHE:N	1:A:154:PRO:CD	2.73	0.50
1:A:35:HIS:ND1	1:A:47:TRP:HH2	2.03	0.50
2:B:182:LYS:HD3	2:B:182:LYS:H	1.76	0.50
1:A:196:LEU:HA	1:A:201:TYR:HE2	1.76	0.50
1:A:154:PRO:HG2	1:A:207:HIS:HE1	1.77	0.50
1:A:35:HIS:NE2	1:A:98:ASP:HB3	2.27	0.49
1:A:154:PRO:HB2	1:A:207:HIS:NE2	2.27	0.49
3:C:259:LEU:HD21	3:C:267:LYS:HE3	1.95	0.49
1:A:104:TYR:HE2	3:C:259:LEU:O	1.95	0.49
1:A:105:TYR:CE2	2:B:90:TRP:HA	2.48	0.49
3:C:307:ARG:CZ	3:C:324:GLN:HG2	2.42	0.49
1:A:5:GLN:O	1:A:23:THR:HB	2.13	0.49
2:B:157:ASN:CG	2:B:158:SER:H	2.15	0.49
3:C:264:GLY:HA3	3:C:290:TRP:CE2	2.47	0.49
3:C:260:THR:HG22	3:C:263:GLN:HB2	1.95	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:ILE:HG13	1:A:78:VAL:HG21	1.96	0.48
1:A:87:THR:HA	1:A:118:VAL:HB	1.94	0.48
2:B:29:VAL:HG21	2:B:89:GLN:HG2	1.96	0.48
3:C:379:ILE:HG13	3:C:381:VAL:HG23	1.95	0.48
3:C:464:PHE:CZ	4:D:1:NAG:H82	2.49	0.48
3:C:404:CYS:HB2	3:C:407:LEU:HB2	1.95	0.48
3:C:284:THR:HG23	3:C:287:LEU:HD13	1.94	0.48
2:B:33:HIS:O	2:B:88:GLN:NE2	2.46	0.48
2:B:16:GLY:HA2	2:B:76:ARG:HG3	1.95	0.48
2:B:158:SER:HA	2:B:178:LEU:HD22	1.96	0.48
3:C:295:LEU:HD21	3:C:322:VAL:HG21	1.96	0.47
3:C:419:GLU:HG2	3:C:472:VAL:HG12	1.95	0.47
2:B:39:PRO:HB3	2:B:164:GLU:HG3	1.96	0.47
3:C:264:GLY:HA3	3:C:290:TRP:CD1	2.50	0.47
3:C:320:ASN:HB3	3:C:350:TYR:CD1	2.49	0.47
2:B:47:ILE:HD12	2:B:72:LEU:HD13	1.96	0.47
2:B:60:ARG:HD2	2:B:76:ARG:O	2.15	0.47
1:A:203:CYS:O	1:A:215:ASP:HA	2.15	0.47
1:A:73:ASN:N	1:A:77:GLN:O	2.48	0.47
3:C:391:THR:HG21	3:C:491:ILE:HD11	1.97	0.46
2:B:90:TRP:CE3	2:B:91:SER:HB3	2.50	0.46
3:C:371:SER:O	3:C:373:THR:N	2.48	0.46
2:B:84:THR:OG1	2:B:102:LYS:HG3	2.14	0.46
1:A:58:LYS:HA	1:A:58:LYS:HD3	1.55	0.46
2:B:129:ALA:HB3	2:B:181:SER:O	2.15	0.46
3:C:259:LEU:O	3:C:293:ARG:NH2	2.48	0.46
2:B:81:ASP:O	2:B:103:LEU:HD22	2.16	0.46
3:C:429:VAL:HG22	3:C:485:LEU:HD12	1.97	0.46
1:A:34:ILE:HG21	1:A:78:VAL:HG11	1.97	0.46
2:B:139:TYR:CD1	2:B:140:PRO:HA	2.51	0.46
3:C:382:THR:CG2	3:C:463:THR:H	2.25	0.46
2:B:209:ASN:O	2:B:210:ARG:HB2	2.15	0.46
1:A:161:TRP:CZ3	1:A:203:CYS:HB3	2.51	0.45
2:B:163:THR:OG1	2:B:173:SER:N	2.40	0.45
1:A:150:LYS:HG2	1:A:151:ASP:N	2.31	0.45
3:C:397:THR:HB	3:C:508:THR:HG23	1.98	0.45
1:A:199:GLN:O	1:A:200:THR:OG1	2.32	0.45
2:B:181:SER:HB3	2:B:184:ASP:HB2	1.98	0.45
2:B:182:LYS:HB2	2:B:183:ALA:H	1.59	0.45
3:C:319:PHE:HE1	3:C:353:LYS:HA	1.82	0.45
3:C:429:VAL:O	3:C:429:VAL:HG13	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:491:ILE:O	3:C:504:ARG:HB2	2.17	0.45
1:A:32:TYR:HB3	1:A:97:ARG:HD2	1.99	0.44
2:B:105:ILE:H	2:B:165:GLN:HE22	1.65	0.44
2:B:95:LEU:HD12	2:B:96:THR:N	2.30	0.44
3:C:431:LEU:HA	3:C:485:LEU:HD13	1.99	0.44
3:C:459:TYR:HA	3:C:462:GLY:O	2.18	0.44
1:A:105:TYR:HE2	2:B:90:TRP:CD1	2.36	0.44
1:A:66:ARG:HG3	1:A:67:LEU:HD12	2.00	0.44
2:B:89:GLN:NE2	2:B:89:GLN:HA	2.29	0.44
3:C:422:VAL:HG22	3:C:469:TYR:O	2.18	0.43
1:A:4:LEU:HD22	1:A:24:VAL:HB	2.00	0.43
1:A:97:ARG:HB3	1:A:109:HIS:CE1	2.53	0.43
2:B:89:GLN:HB3	2:B:92:PHE:CE2	2.52	0.43
2:B:157:ASN:ND2	2:B:179:THR:H	1.99	0.43
3:C:276:LEU:HB3	3:C:278:LEU:HG	2.00	0.43
3:C:388:ILE:HB	3:C:503:LEU:HB3	2.01	0.43
1:A:154:PRO:HG2	1:A:207:HIS:CE1	2.54	0.43
2:B:91:SER:N	3:C:258:ILE:HG22	2.31	0.43
2:B:192:ALA:HB2	2:B:207:SER:HB3	2.00	0.43
1:A:176:VAL:O	1:A:183:TYR:HA	2.19	0.43
2:B:19:VAL:O	2:B:73:THR:HA	2.18	0.43
1:A:108:ASP:OD1	1:A:108:ASP:N	2.42	0.43
2:B:125:LYS:O	2:B:125:LYS:HG3	2.19	0.43
1:A:193:SER:O	1:A:196:LEU:HB3	2.19	0.42
3:C:269:TYR:HD2	3:C:359:ASP:H	1.67	0.42
3:C:377:GLN:HB3	3:C:441:HIS:NE2	2.33	0.42
3:C:380:ASN:OD1	3:C:382:THR:HG23	2.19	0.42
1:A:32:TYR:CG	1:A:97:ARG:HD2	2.55	0.42
1:A:101:ARG:O	3:C:261:LEU:HB2	2.18	0.42
2:B:2:ILE:HG12	2:B:92:PHE:CZ	2.54	0.42
3:C:503:LEU:HD12	3:C:503:LEU:HA	1.65	0.42
2:B:139:TYR:CG	2:B:140:PRO:HA	2.54	0.42
1:A:9:PRO:HG2	1:A:208:LYS:HD2	2.02	0.42
2:B:146:GLN:HG2	2:B:153:LEU:HD22	2.02	0.42
1:A:87:THR:O	1:A:90:THR:HG22	2.19	0.42
3:C:259:LEU:HD13	3:C:264:GLY:HA2	2.02	0.42
1:A:105:TYR:CE1	2:B:95:LEU:HD22	2.55	0.42
3:C:500:GLN:HB2	3:C:501:SER:H	1.56	0.42
1:A:43:LYS:HE2	1:A:43:LYS:HB2	1.95	0.42
3:C:307:ARG:HD3	3:C:307:ARG:HA	1.70	0.42
3:C:313:ILE:HG23	3:C:370:ARG:NH2	2.35	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:438:ILE:HD11	3:C:445:TYR:OH	2.19	0.42
1:A:32:TYR:CB	1:A:97:ARG:HD2	2.50	0.41
3:C:259:LEU:CD1	3:C:267:LYS:HE3	2.42	0.41
1:A:66:ARG:NH1	1:A:83:ASN:O	2.45	0.41
2:B:32:MET:O	2:B:50:THR:HA	2.20	0.41
2:B:107:ARG:HH11	2:B:171:THR:HG22	1.85	0.41
1:A:36:TRP:CD1	1:A:69:MET:SD	3.13	0.41
3:C:309:GLU:HG2	3:C:460:HIS:HB2	2.02	0.41
1:A:175:ALA:HA	1:A:185:LEU:HB3	2.03	0.41
2:B:33:HIS:CB	2:B:48:PHE:HA	2.45	0.41
3:C:326:LYS:CE	3:C:326:LYS:HA	2.51	0.41
1:A:69:MET:HA	1:A:80:LEU:C	2.41	0.41
3:C:510:ILE:HD13	3:C:510:ILE:HA	1.85	0.41
1:A:104:TYR:CE2	3:C:260:THR:HA	2.56	0.40
1:A:143:ALA:HB3	1:A:191:VAL:O	2.21	0.40
2:B:162:VAL:HG22	2:B:174:LEU:HD12	2.03	0.40
3:C:425:PRO:HG2	3:C:466:PRO:O	2.22	0.40
3:C:435:HIS:ND1	3:C:448:ASP:HB3	2.37	0.40
1:A:204:ASN:HD22	1:A:215:ASP:CG	2.25	0.40
1:A:207:HIS:CD2	1:A:209:PRO:HD2	2.57	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	184/233 (79%)	166 (90%)	14 (8%)	4 (2%)	6	39
2	B	202/213 (95%)	186 (92%)	13 (6%)	3 (2%)	10	47
3	C	235/274 (86%)	209 (89%)	22 (9%)	4 (2%)	9	44
All	All	621/720 (86%)	561 (90%)	49 (8%)	11 (2%)	8	43

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	134	SER
2	B	93	ASN
2	B	182	LYS
3	C	352	ALA
3	C	372	SER
3	C	381	VAL
1	A	156	PRO
1	A	101	ARG
3	C	500	GLN
1	A	200	THR
2	B	83	ALA

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	173/201 (86%)	170 (98%)	3 (2%)	60	82
2	B	179/183 (98%)	178 (99%)	1 (1%)	86	94
3	C	218/248 (88%)	210 (96%)	8 (4%)	34	66
All	All	570/632 (90%)	558 (98%)	12 (2%)	53	78

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	LYS
1	A	77	GLN
1	A	199	GLN
2	B	11	LEU
3	C	348	LYS
3	C	395	LYS
3	C	431	LEU
3	C	443	CYS
3	C	477	ASN
3	C	482	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	C	497	ARG
3	C	507	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	162	ASN
1	A	199	GLN
2	B	123	GLN
2	B	157	ASN
2	B	165	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 ⓘ

4 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	D	1	4,3	14,14,15	0.46	0	17,19,21	0.81	1 (5%)
4	NAG	D	2	4	14,14,15	0.33	0	17,19,21	0.64	1 (5%)
4	NAG	G	1	4,3	14,14,15	0.27	0	17,19,21	0.41	0
4	NAG	G	2	4	14,14,15	0.36	0	17,19,21	0.62	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
4	NAG	D	1	4,3	-	2/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1
4	NAG	G	1	4,3	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1	NAG	C1-O5-C5	2.81	116.00	112.19
4	D	2	NAG	C1-O5-C5	2.19	115.17	112.19
4	G	2	NAG	C1-O5-C5	2.09	115.02	112.19

There are no chirality outliers.

All (5) torsion outliers are listed below:

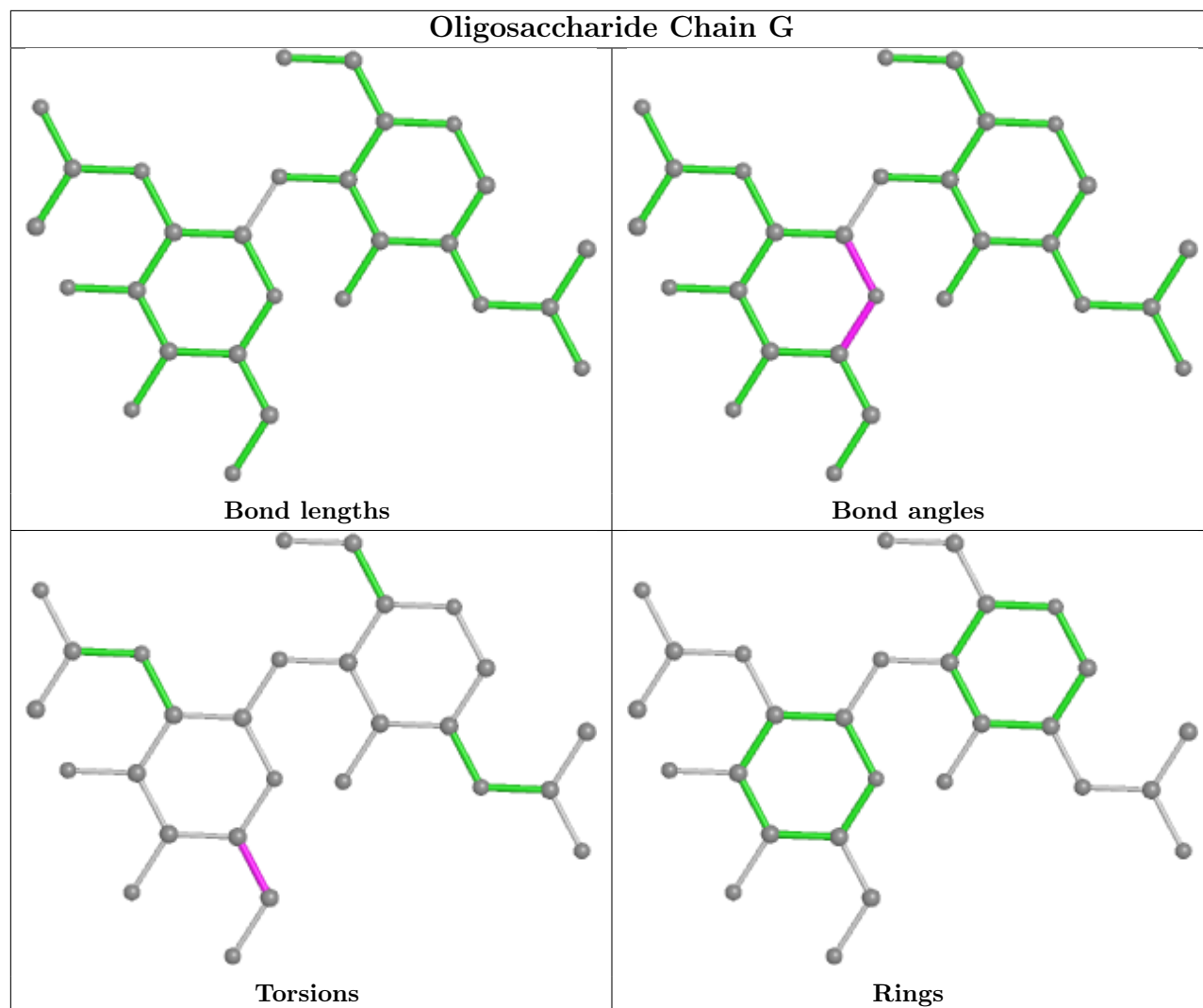
Mol	Chain	Res	Type	Atoms
4	D	1	NAG	O5-C5-C6-O6
4	D	2	NAG	C8-C7-N2-C2
4	D	2	NAG	O7-C7-N2-C2
4	D	1	NAG	C4-C5-C6-O6
4	G	2	NAG	C4-C5-C6-O6

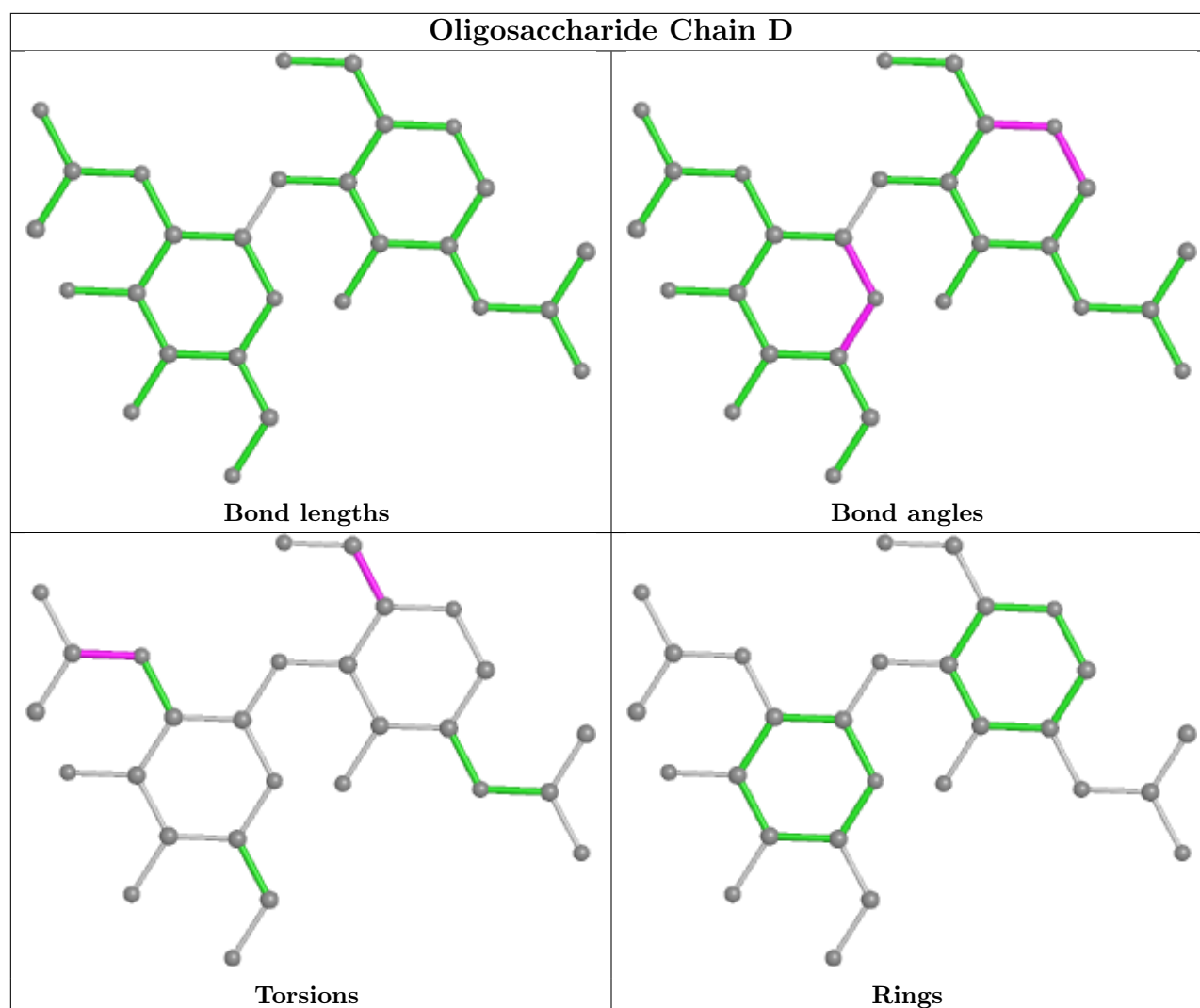
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1	NAG	3	0
4	D	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	202/233 (86%)	0.39	17 (8%)	11 7	149, 225, 302, 360	0
2	B	208/213 (97%)	0.45	15 (7%)	15 9	137, 181, 263, 297	0
3	C	241/274 (87%)	0.08	0	100 100	94, 143, 199, 307	0
All	All	651/720 (90%)	0.29	32 (4%)	29 19	94, 176, 277, 360	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	218	VAL	6.6
2	B	128	THR	4.8
2	B	129	ALA	4.3
1	A	36	TRP	4.3
2	B	119	PRO	3.8
2	B	181	SER	3.2
1	A	107	MET	3.1
1	A	47	TRP	3.1
2	B	117	PHE	3.0
1	A	131	LEU	2.9
2	B	180	LEU	2.9
2	B	185	TYR	2.8
2	B	198	GLN	2.8
2	B	146	GLN	2.7
1	A	198	THR	2.7
2	B	32	MET	2.6
1	A	48	LEU	2.6
1	A	145	LEU	2.5
2	B	179	THR	2.5
1	A	106	ALA	2.5
1	A	130	PRO	2.5
2	B	34	TRP	2.5
1	A	193	SER	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	35	TYR	2.4
1	A	37	VAL	2.3
1	A	161	TRP	2.3
2	B	112	PRO	2.3
1	A	58	LYS	2.3
2	B	145	VAL	2.3
1	A	109	HIS	2.3
1	A	59	TYR	2.2
1	A	152	TYR	2.2

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](#)

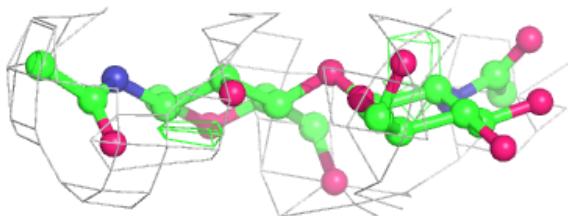
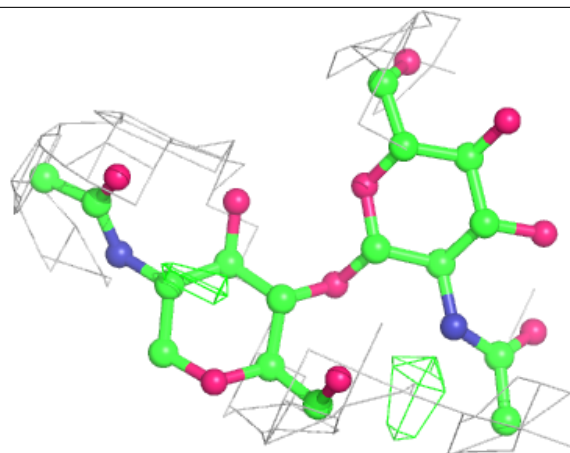
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

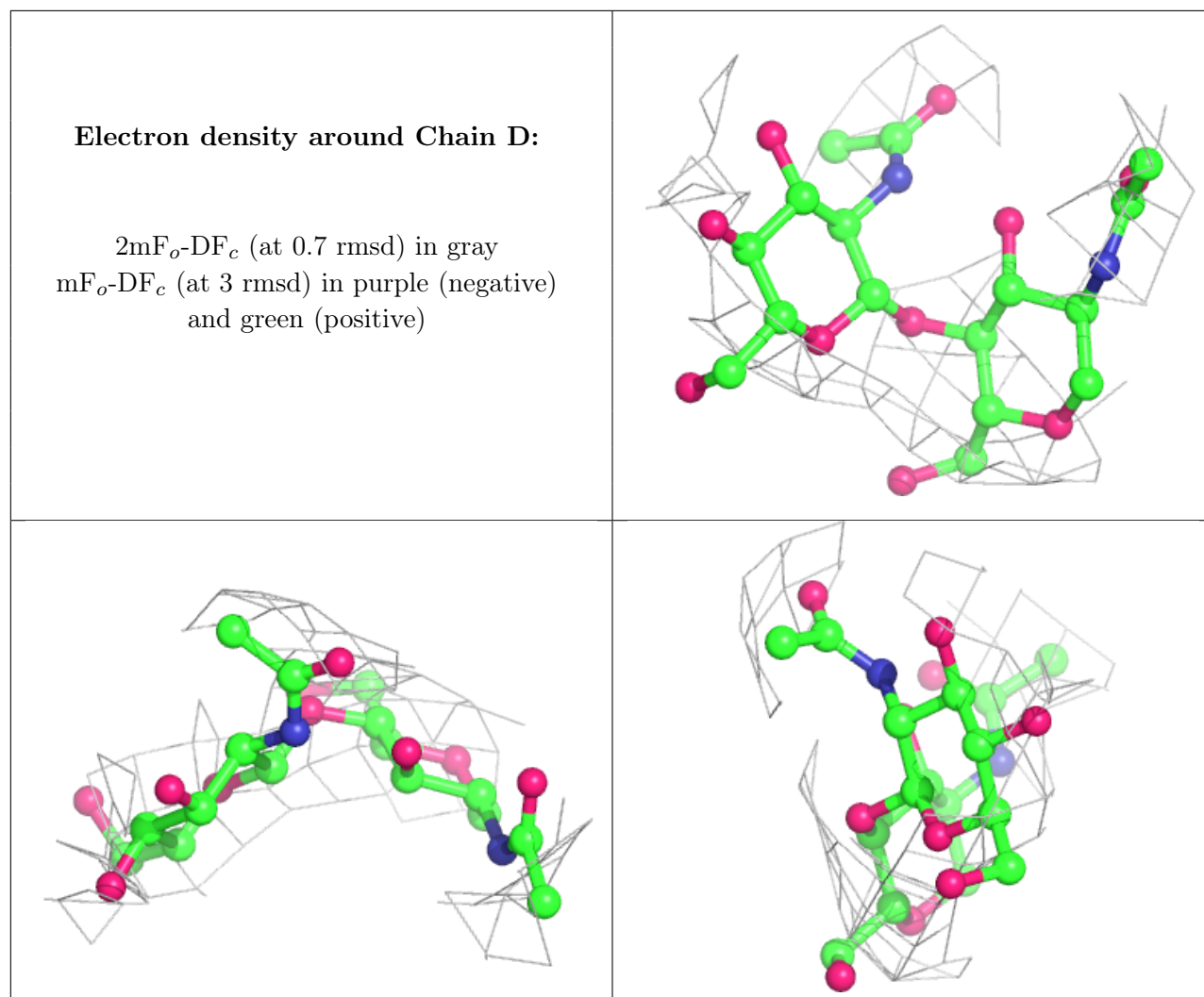
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	D	1	14/15	0.84	0.27	134,170,193,194	0
4	NAG	D	2	14/15	0.84	0.21	182,201,215,225	0
4	NAG	G	1	14/15	0.93	0.30	130,151,170,170	0
4	NAG	G	2	14/15	0.95	0.42	157,168,184,188	0

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 G:

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





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.