



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 30, 2021 – 12:56 PM EDT

PDB ID : 6VSL
Title : Crystal structure of a human fucosylated IgG1 Fc expressed in tobacco plants
(*Nicotiana benthamiana*)
Authors : Tolbert, W.D.; Pazgier, M.
Deposited on : 2020-02-11
Resolution : 2.10 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 : 2.23.1
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) : 2.23.1

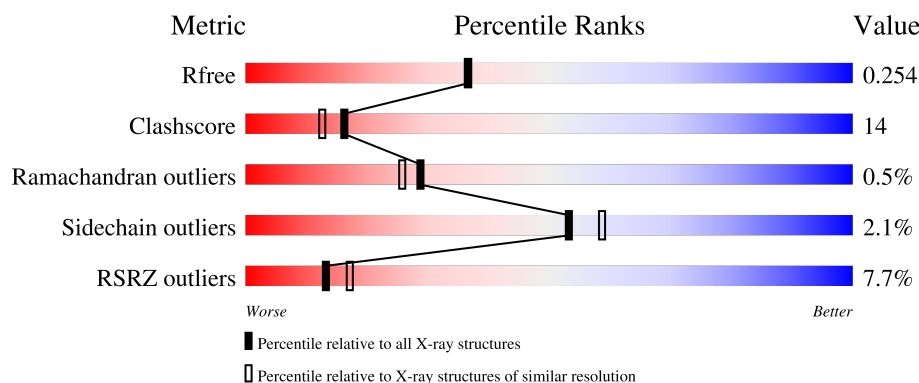
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	211	<div> <div></div> <div>85%</div> <div>14%</div> </div>
1	B	211	<div> <div>14%</div> <div>66%</div> <div>31%</div> <div>..</div> </div>
2	C	9	<div> <div>44%</div> <div>44%</div> <div>11%</div> </div>
2	D	9	<div> <div>44%</div> <div>56%</div> </div>

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
2	NAG	D	1	-	-	-	X
2	NAG	D	8	-	-	-	X
2	FUC	D	9	-	-	-	X

2 Entry composition [i](#)

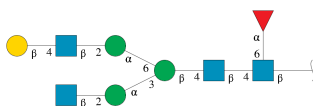
There are 4 unique types of molecules in this entry. The entry contains 3784 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Immunoglobulin gamma-1 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	0	1	0
			1684	1072	283	322	7			
1	B	208	Total	C	N	O	S	0	1	0
			1675	1065	284	320	6			

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	9	Total	C	N	O	0	0	0
			110	62	4	44			
2	D	9	Total	C	N	O	0	0	0
			110	62	4	44			

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			8	6	2		
3	B	1	Total	C	O	0	0
			8	6	2		

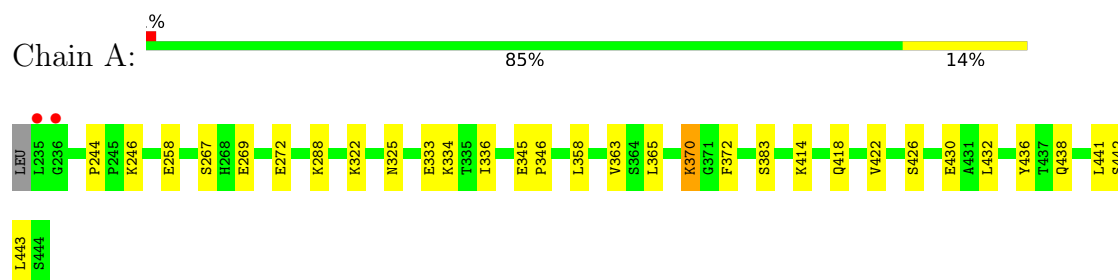
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	119	Total	O	0	0
			119	119		
4	B	70	Total	O	0	0
			70	70		

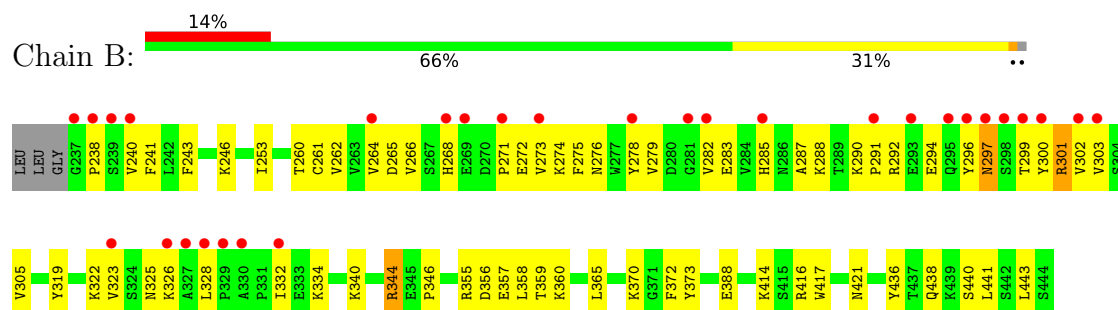
3 Residue-property plots

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.

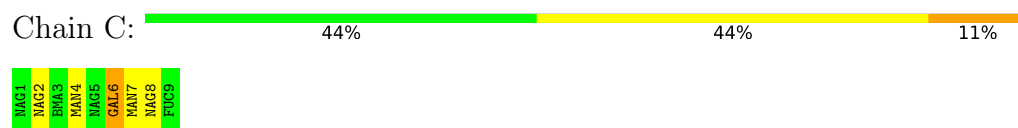
- Molecule 1: Immunoglobulin gamma-1 heavy chain



- Molecule 1: Immunoglobulin gamma-1 heavy chain



- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1
MAG2
BMA3
MAN4
NAG5
GAL6
MAN7
MAG8
FUC9

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.86Å 79.92Å 138.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.91 – 2.10 46.91 – 2.10	Depositor EDS
% Data completeness (in resolution range)	93.7 (46.91-2.10) 98.2 (46.91-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.17.1_3660: ???), REFMAC 5.8.0258	Depositor
R, R_{free}	0.207 , 0.250 0.208 , 0.254	Depositor DCC
R_{free} test set	1618 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.298	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3784	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.27% 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: MAN, MPD, FUC, GAL, 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	0.42	0/1730	0.57	0/2356
1	B	0.46	0/1721	0.64	0/2344
All	All	0.44	0/3451	0.60	0/4700

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1684	0	1652	21	0
1	B	1675	0	1642	81	0
2	C	110	0	94	1	0
2	D	110	0	94	7	0
3	A	8	0	14	1	0
3	B	8	0	14	5	0
4	A	119	0	0	2	0
4	B	70	0	0	4	0
All	All	3784	0	3510	102	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 14.

All (102) 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:B:332:ILE:HB	3:B:510:MPD:H52	1.30	1.09
1:B:268:HIS:CE1	1:B:300:TYR:HE2	1.74	1.03
1:B:292:ARG:HG3	1:B:302:VAL:HG22	1.46	0.97
1:B:268:HIS:CE1	1:B:300:TYR:CE2	2.56	0.94
1:B:292:ARG:HG2	1:B:300:TYR:HB3	1.58	0.83
1:B:275:PHE:CD1	1:B:323:VAL:HG12	2.18	0.79
1:B:240:VAL:HG11	1:B:332:ILE:HD11	1.65	0.78
1:B:292:ARG:CG	1:B:300:TYR:HB3	2.17	0.75
1:B:332:ILE:HB	3:B:510:MPD:C5	2.15	0.74
1:B:272:GLU:O	1:B:325:ASN:ND2	2.20	0.74
1:B:294:GLU:HG3	1:B:300:TYR:CE2	2.25	0.71
1:B:276:ASN:HB2	1:B:322:LYS:HE3	1.72	0.71
1:B:290:LYS:HB3	1:B:291:PRO:HD2	1.72	0.71
1:B:271:PRO:HD2	1:B:272:GLU:OE1	1.91	0.71
1:A:258:GLU:OE1	4:A:601:HOH:O	2.09	0.69
1:B:358:LEU:O	1:B:414:LYS:NZ	2.22	0.67
1:B:357:GLU:OE2	4:B:601:HOH:O	2.12	0.66
1:B:278:TYR:CD1	1:B:283:GLU:HG3	2.31	0.66
1:B:268:HIS:HA	1:B:271:PRO:HG3	1.78	0.65
1:B:240:VAL:HB	1:B:332:ILE:HD13	1.80	0.62
1:B:417:TRP:NE1	1:B:443:LEU:HD13	2.14	0.62
1:B:273:VAL:HG21	1:B:302:VAL:HG11	1.81	0.62
1:B:334:LYS:NZ	3:B:510:MPD:H32	2.16	0.61
1:B:370:LYS:NZ	4:B:602:HOH:O	2.27	0.61
1:B:265:ASP:OD1	2:D:1:NAG:N2	2.25	0.59
1:B:268:HIS:HE1	1:B:300:TYR:CE2	2.18	0.59
1:B:290:LYS:HE3	1:B:305:VAL:HG21	1.87	0.57
1:A:358:LEU:O	1:A:414:LYS:NZ	2.26	0.56
1:A:334:LYS:NZ	3:A:510:MPD:H12	2.21	0.56
1:B:246:LYS:NZ	2:D:5:NAG:O3	2.26	0.56
1:B:262:VAL:HG22	1:B:303:VAL:HG22	1.87	0.56
1:B:325:ASN:HB3	1:B:328:LEU:HD13	1.89	0.55
1:A:244:PRO:HB3	1:A:336:ILE:HD11	1.88	0.55
1:B:292:ARG:NH1	1:B:292:ARG:HB3	2.21	0.55
1:B:264:VAL:HG12	1:B:265:ASP:N	2.22	0.55
1:B:264:VAL:HG11	2:D:1:NAG:O4	2.07	0.54
1:B:275:PHE:CE1	1:B:323:VAL:HG12	2.42	0.54
1:B:294:GLU:HB2	1:B:300:TYR:CE1	2.42	0.54
1:B:264:VAL:HG12	1:B:265:ASP:H	1.73	0.54
1:A:422:VAL:HA	1:A:442:SER:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:290:LYS:HB3	1:B:291:PRO:CD	2.38	0.54
1:B:275:PHE:CE1	1:B:323:VAL:CG1	2.91	0.53
1:B:238:PRO:HB3	1:B:266:VAL:HG22	1.90	0.52
1:B:238:PRO:HA	1:B:264:VAL:O	2.10	0.52
1:A:322:LYS:HG3	1:A:333:GLU:HG2	1.92	0.51
1:A:358:LEU:HD23	1:A:363:VAL:HG11	1.92	0.51
1:B:246:LYS:HZ3	2:D:6:GAL:H2	1.75	0.51
1:A:426:SER:HB3	1:A:438:GLN:HG2	1.92	0.50
1:A:430:GLU:OE2	4:A:602:HOH:O	2.19	0.50
1:B:240:VAL:CG1	1:B:332:ILE:HD11	2.37	0.50
1:B:297:ASN:OD1	1:B:299:THR:OG1	2.18	0.50
1:B:346:PRO:HB3	1:B:372:PHE:HB3	1.93	0.49
1:B:416:ARG:O	1:B:421:ASN:ND2	2.39	0.49
1:B:278:TYR:CE1	1:B:283:GLU:HG3	2.48	0.49
1:B:266:VAL:CG1	1:B:271:PRO:HA	2.43	0.48
1:B:290:LYS:HE3	1:B:305:VAL:CG2	2.45	0.47
1:B:268:HIS:CA	1:B:271:PRO:HG3	2.42	0.47
1:A:345:GLU:HG3	1:A:432:LEU:HD23	1.96	0.47
1:B:301:ARG:HD3	1:B:303:VAL:HG23	1.96	0.47
1:A:370:LYS:HE3	4:B:610:HOH:O	2.14	0.46
1:B:287:ALA:O	1:B:288:LYS:HG3	2.15	0.46
1:B:334:LYS:HZ1	3:B:510:MPD:H32	1.79	0.46
1:B:296:TYR:O	1:B:296:TYR:CG	2.69	0.46
1:B:240:VAL:HB	1:B:332:ILE:CD1	2.44	0.46
1:B:325:ASN:CG	1:B:326:LYS:H	2.20	0.46
1:B:292:ARG:HB3	1:B:292:ARG:HH11	1.81	0.46
1:B:279:VAL:HG22	1:B:319:TYR:CD2	2.52	0.45
1:B:325:ASN:N	1:B:328:LEU:HD22	2.32	0.45
1:A:346:PRO:HB3	1:A:372:PHE:HB3	1.99	0.45
1:A:383:SER:OG	1:A:422:VAL:O	2.21	0.45
1:B:253:ILE:HG12	4:B:616:HOH:O	2.16	0.45
1:B:325:ASN:CB	1:B:328:LEU:HD13	2.46	0.45
1:B:243:PHE:HE2	2:D:3:BMA:H62	1.81	0.44
1:A:269:GLU:CD	1:A:269:GLU:H	2.21	0.44
1:B:274:LYS:HG2	1:B:275:PHE:N	2.33	0.44
1:A:288:LYS:HA	1:A:288:LYS:HD3	1.80	0.43
1:A:365:LEU:HD13	1:A:441:LEU:HD23	2.00	0.43
1:B:294:GLU:HB2	1:B:300:TYR:CZ	2.53	0.43
1:B:241:PHE:CE1	2:D:2:NAG:H4	2.53	0.43
1:B:238:PRO:HB3	1:B:266:VAL:CG2	2.48	0.43
1:B:326:LYS:N	1:B:326:LYS:HD2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:LYS:HG2	2:C:6:GAL:O4	2.18	0.43
1:B:273:VAL:CG2	1:B:302:VAL:HG11	2.47	0.43
1:B:388:GLU:OE1	1:B:416:ARG:NH2	2.31	0.42
1:B:322:LYS:HE3	1:B:322:LYS:HB3	1.79	0.42
1:B:417:TRP:CZ2	1:B:443:LEU:HA	2.54	0.42
1:B:355:ARG:HA	1:B:358:LEU:HD12	2.01	0.42
1:B:436:TYR:CE1	1:B:438:GLN:HG3	2.55	0.42
1:A:267:SER:HB2	1:A:269:GLU:OE1	2.20	0.42
1:B:275:PHE:CE1	1:B:323:VAL:HG11	2.55	0.41
1:B:292:ARG:CB	1:B:292:ARG:CZ	2.97	0.41
1:B:268:HIS:O	1:B:271:PRO:HG3	2.21	0.41
1:B:292:ARG:CZ	1:B:292:ARG:HB2	2.50	0.41
1:A:418:GLN:HA	1:A:443:LEU:HD22	2.02	0.41
1:B:240:VAL:CB	1:B:332:ILE:CD1	2.99	0.41
1:B:359:THR:HG23	1:B:360:LYS:HD2	2.03	0.40
1:A:436:TYR:OH	1:A:438:GLN:HG3	2.21	0.40
1:B:334:LYS:HZ2	3:B:510:MPD:H32	1.85	0.40
1:B:344[B]:ARG:HG3	1:B:373:TYR:HB3	2.02	0.40
1:A:272:GLU:O	1:A:325:ASN:ND2	2.53	0.40
1:B:365:LEU:HG	1:B:441:LEU:HD23	2.04	0.40
1:B:260:THR:HG21	2:D:5:NAG:H61	2.03	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	209/211 (99%)	206 (99%)	3 (1%)	0	100	100
1	B	207/211 (98%)	196 (95%)	9 (4%)	2 (1%)	15	11
All	All	416/422 (99%)	402 (97%)	12 (3%)	2 (0%)	29	26

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	282	VAL
1	B	297	ASN

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	196/196 (100%)	195 (100%)	1 (0%)	88	92
1	B	195/196 (100%)	187 (96%)	8 (4%)	30	31
All	All	391/392 (100%)	382 (98%)	9 (2%)	53	55

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	370	LYS
1	B	261	CYS
1	B	285	HIS
1	B	301	ARG
1	B	340	LYS
1	B	344[A]	ARG
1	B	344[B]	ARG
1	B	356	ASP
1	B	440	SER

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

Mol	Chain	Res	Type
1	A	347	GLN
1	B	268	HIS
1	B	325	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

18 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	NAG	C	1	1,2	14,14,15	0.44	0	17,19,21	0.52	0
2	NAG	C	2	2	14,14,15	0.66	1 (7%)	17,19,21	0.56	0
2	BMA	C	3	2	11,11,12	0.89	0	15,15,17	0.95	0
2	MAN	C	4	2	11,11,12	0.96	1 (9%)	15,15,17	1.47	2 (13%)
2	NAG	C	5	2	14,14,15	0.19	0	17,19,21	0.53	0
2	GAL	C	6	2	11,11,12	0.96	1 (9%)	15,15,17	1.52	2 (13%)
2	MAN	C	7	2	11,11,12	1.11	1 (9%)	15,15,17	1.15	2 (13%)
2	NAG	C	8	2	14,14,15	0.41	0	17,19,21	0.87	1 (5%)
2	FUC	C	9	2	10,10,11	0.80	0	14,14,16	0.78	0
2	NAG	D	1	1,2	14,14,15	0.62	1 (7%)	17,19,21	0.93	1 (5%)
2	NAG	D	2	2	14,14,15	1.40	1 (7%)	17,19,21	1.08	1 (5%)
2	BMA	D	3	2	11,11,12	1.24	2 (18%)	15,15,17	1.02	0
2	MAN	D	4	2	11,11,12	0.62	0	15,15,17	1.15	2 (13%)
2	NAG	D	5	2	14,14,15	0.79	1 (7%)	17,19,21	0.47	0
2	GAL	D	6	2	11,11,12	0.82	0	15,15,17	0.94	1 (6%)
2	MAN	D	7	2	11,11,12	1.18	1 (9%)	15,15,17	1.37	1 (6%)
2	NAG	D	8	2	14,14,15	0.52	0	17,19,21	0.97	1 (5%)
2	FUC	D	9	2	10,10,11	0.74	0	14,14,16	1.17	1 (7%)

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	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
2	MAN	C	4	2	-	0/2/19/22	0/1/1/1
2	NAG	C	5	2	-	0/6/23/26	0/1/1/1
2	GAL	C	6	2	-	2/2/19/22	0/1/1/1
2	MAN	C	7	2	-	0/2/19/22	0/1/1/1
2	NAG	C	8	2	-	3/6/23/26	0/1/1/1
2	FUC	C	9	2	-	-	0/1/1/1
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	BMA	D	3	2	-	0/2/19/22	0/1/1/1
2	MAN	D	4	2	-	0/2/19/22	0/1/1/1
2	NAG	D	5	2	-	0/6/23/26	0/1/1/1
2	GAL	D	6	2	-	2/2/19/22	0/1/1/1
2	MAN	D	7	2	-	2/2/19/22	0/1/1/1
2	NAG	D	8	2	-	2/6/23/26	0/1/1/1
2	FUC	D	9	2	-	-	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	NAG	O5-C1	-5.11	1.35	1.43
2	D	3	BMA	C2-C3	3.27	1.57	1.52
2	D	7	MAN	O5-C1	-2.97	1.39	1.43
2	D	5	NAG	O5-C1	-2.86	1.39	1.43
2	C	7	MAN	C2-C3	-2.48	1.48	1.52
2	C	2	NAG	O5-C1	-2.36	1.39	1.43
2	C	6	GAL	C1-C2	2.28	1.57	1.52
2	D	3	BMA	O5-C1	-2.25	1.40	1.43
2	D	1	NAG	C1-C2	2.03	1.55	1.52
2	C	4	MAN	C4-C3	2.02	1.57	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4	MAN	C1-O5-C5	4.62	118.45	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	6	GAL	C1-O5-C5	3.85	117.41	112.19
2	D	7	MAN	O2-C2-C3	-3.34	103.45	110.14
2	D	4	MAN	C1-O5-C5	3.32	116.69	112.19
2	D	1	NAG	O4-C4-C5	-3.23	101.27	109.30
2	D	8	NAG	C3-C4-C5	2.93	115.47	110.24
2	D	9	FUC	C1-O5-C5	2.89	119.32	112.78
2	C	7	MAN	C1-O5-C5	2.79	115.97	112.19
2	D	2	NAG	C3-C4-C5	2.65	114.97	110.24
2	C	7	MAN	O2-C2-C3	-2.58	104.96	110.14
2	C	6	GAL	C1-C2-C3	2.57	112.82	109.67
2	C	8	NAG	C3-C4-C5	2.35	114.44	110.24
2	D	4	MAN	O2-C2-C3	-2.23	105.66	110.14
2	C	4	MAN	O2-C2-C3	-2.22	105.69	110.14
2	D	6	GAL	C1-C2-C3	2.07	112.21	109.67

There are no chirality outliers.

All (11) torsion outliers are listed below:

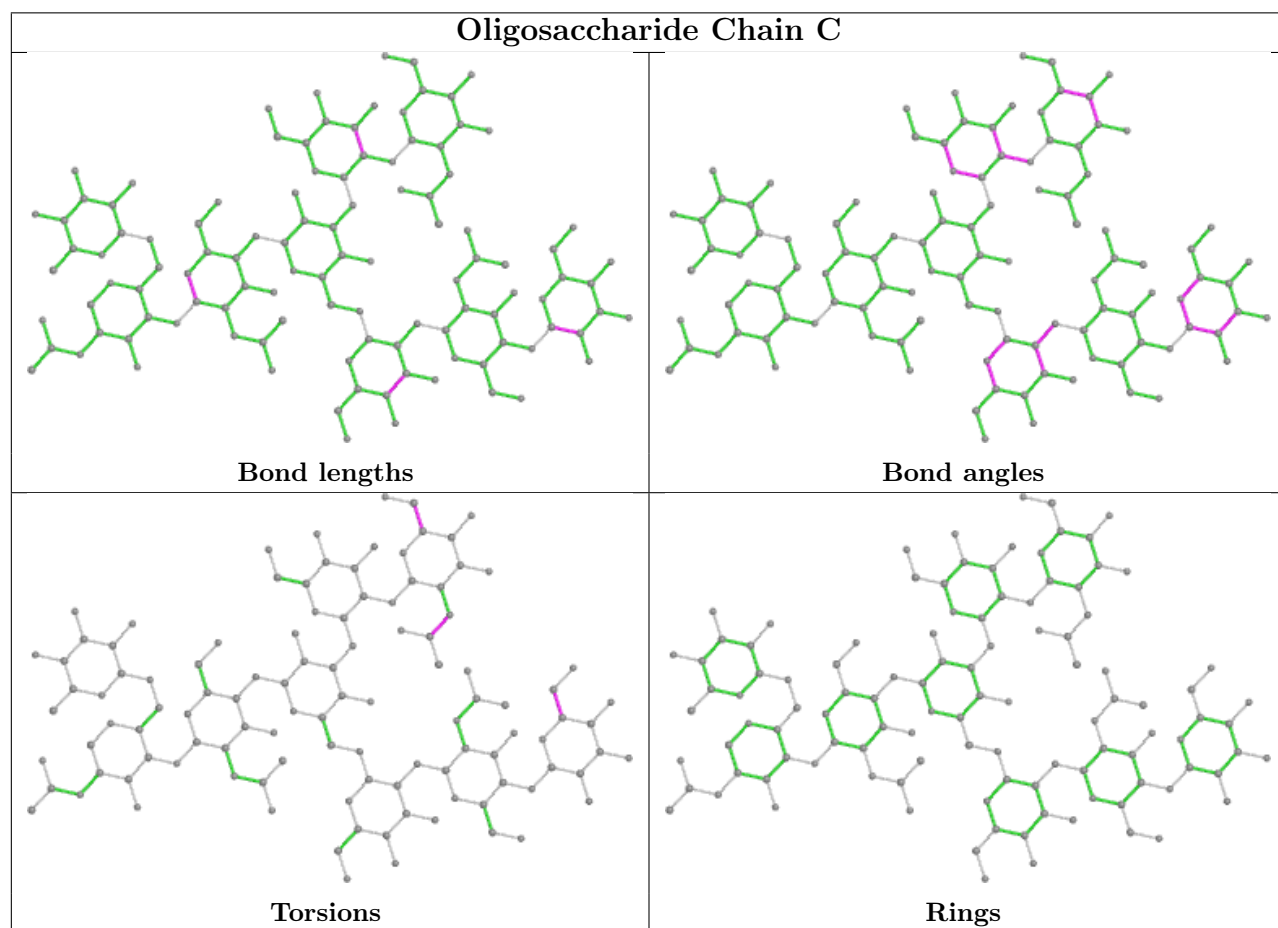
Mol	Chain	Res	Type	Atoms
2	D	6	GAL	O5-C5-C6-O6
2	D	6	GAL	C4-C5-C6-O6
2	D	7	MAN	O5-C5-C6-O6
2	C	6	GAL	C4-C5-C6-O6
2	D	7	MAN	C4-C5-C6-O6
2	C	8	NAG	C8-C7-N2-C2
2	C	8	NAG	O7-C7-N2-C2
2	D	8	NAG	C8-C7-N2-C2
2	D	8	NAG	O7-C7-N2-C2
2	C	6	GAL	O5-C5-C6-O6
2	C	8	NAG	O5-C5-C6-O6

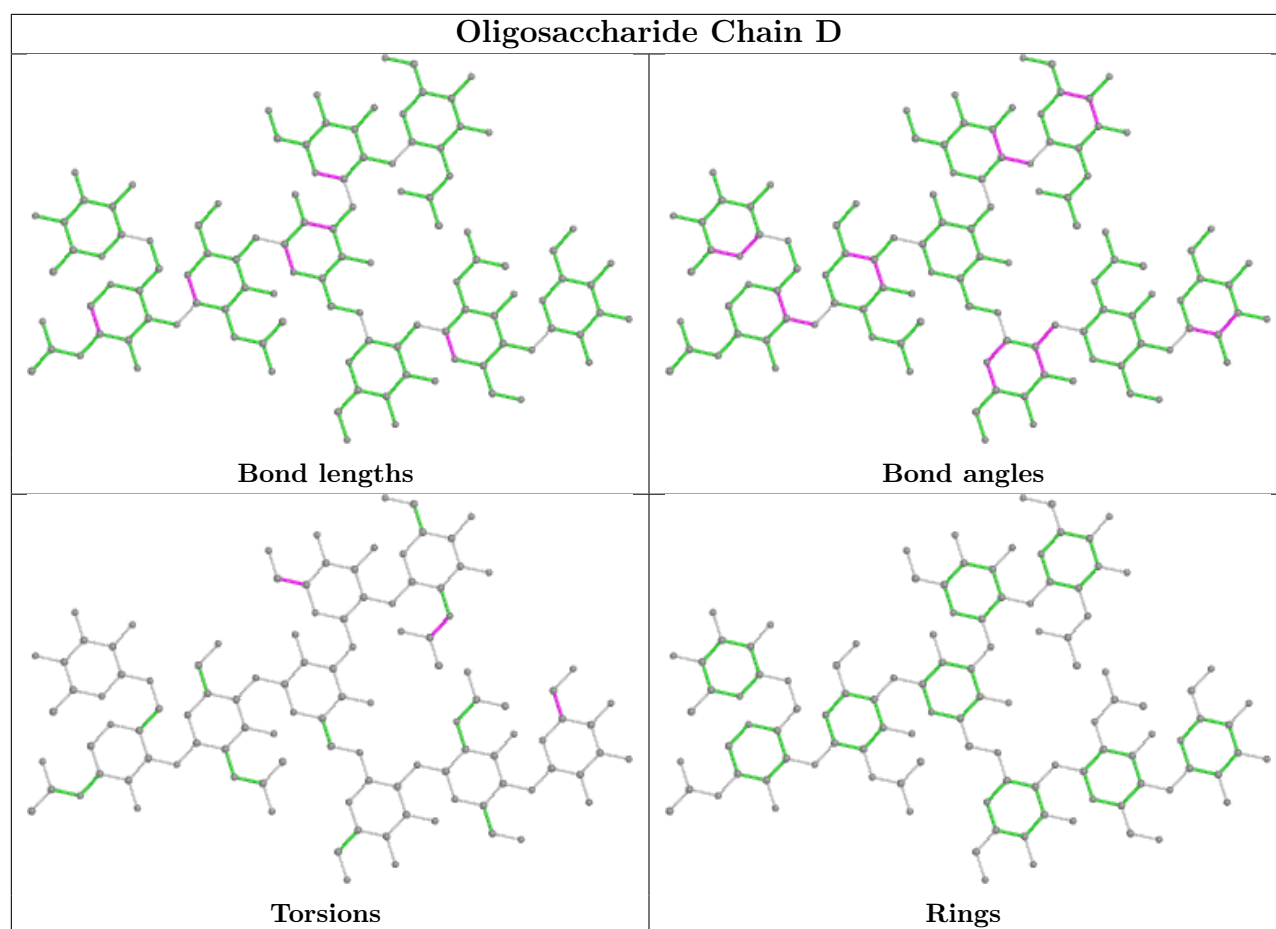
There are no ring outliers.

6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	NAG	1	0
2	D	5	NAG	2	0
2	C	6	GAL	1	0
2	D	1	NAG	2	0
2	D	3	BMA	1	0
2	D	6	GAL	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](#)

2 ligands 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	MPD	B	510	-	7,7,7	0.27	0	9,10,10	0.81	0
3	MPD	A	510	-	7,7,7	0.28	0	9,10,10	0.67	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
3	MPD	B	510	-	-	0/5/5/5	-
3	MPD	A	510	-	-	2/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	510	MPD	C2-C3-C4-O4
3	A	510	MPD	C2-C3-C4-C5

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	510	MPD	5	0
3	A	510	MPD	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	210/211 (99%)	-0.12	2 (0%) 82 85	26, 46, 77, 95	0
1	B	208/211 (98%)	0.61	30 (14%) 2 3	25, 62, 130, 166	0
All	All	418/422 (99%)	0.24	32 (7%) 13 17	25, 52, 118, 166	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	296	TYR	12.3
1	B	300	TYR	6.3
1	B	282	VAL	4.6
1	B	273	VAL	4.5
1	B	332	ILE	4.2
1	B	323	VAL	4.0
1	A	236	GLY	3.8
1	B	297	ASN	3.6
1	B	239	SER	3.6
1	B	298	SER	3.5
1	B	240	VAL	3.5
1	B	328	LEU	3.4
1	B	299	THR	3.4
1	A	235	LEU	3.3
1	B	237	GLY	3.3
1	B	291	PRO	3.0
1	B	264	VAL	3.0
1	B	295	GLN	2.9
1	B	326	LYS	2.9
1	B	281	GLY	2.8
1	B	293	GLU	2.8
1	B	327	ALA	2.7
1	B	330	ALA	2.7
1	B	303	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	271	PRO	2.6
1	B	329	PRO	2.5
1	B	285	HIS	2.4
1	B	238	PRO	2.4
1	B	268	HIS	2.3
1	B	302	VAL	2.2
1	B	269	GLU	2.1
1	B	278	TYR	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

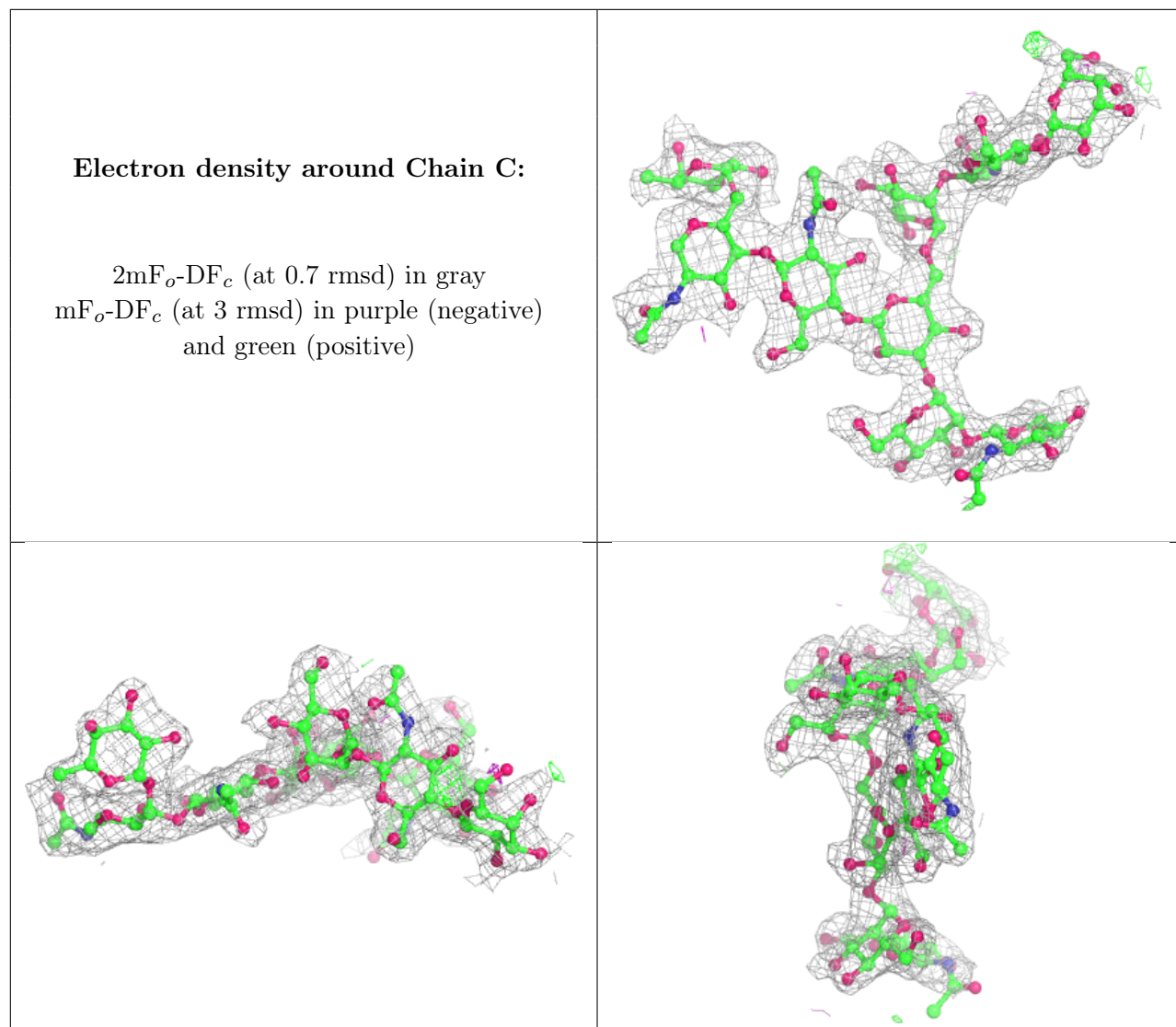
6.3 Carbohydrates ⓘ

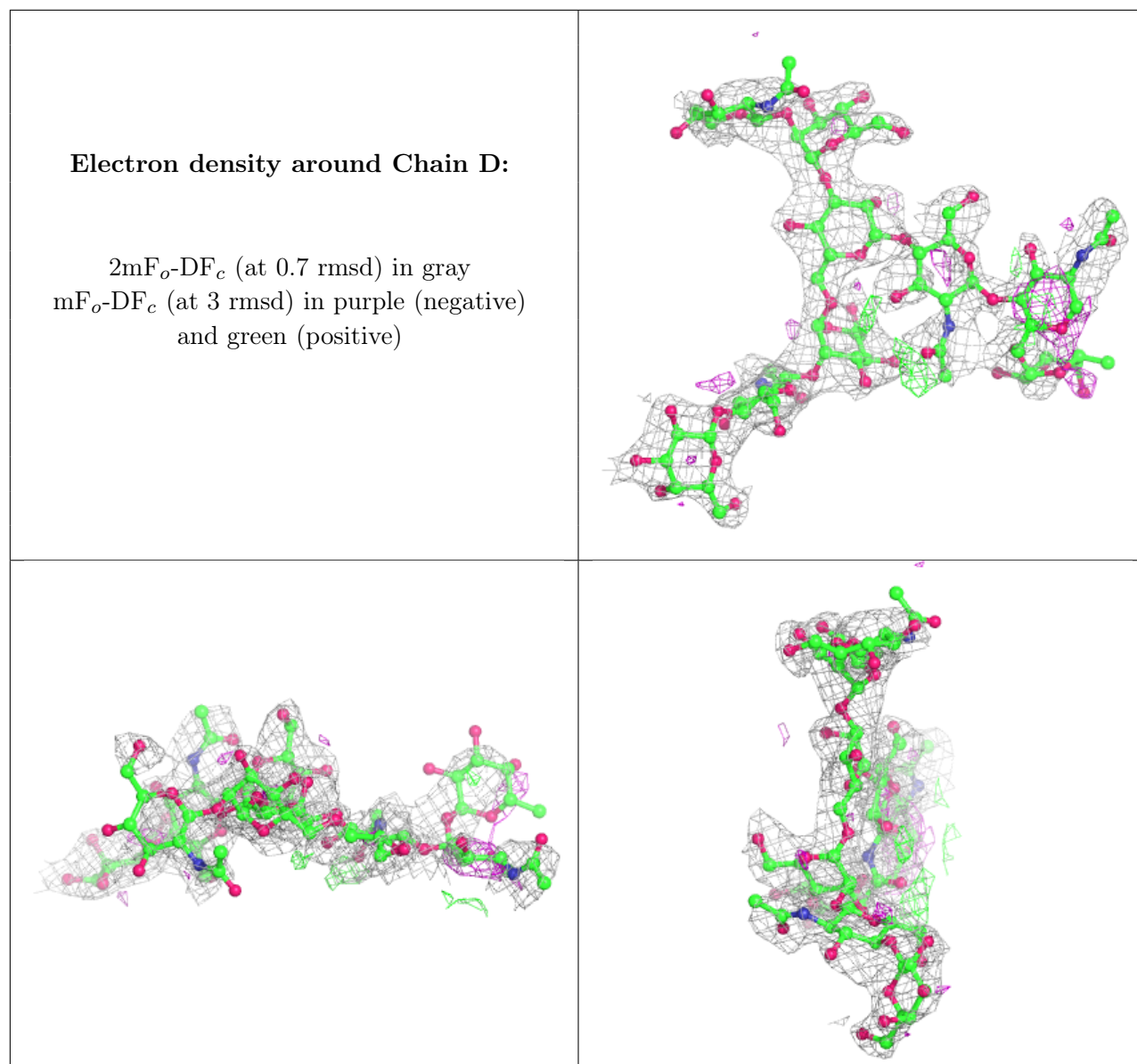
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	D	2	14/15	0.25	0.34	97,102,107,107	0
2	NAG	D	1	14/15	0.47	0.43	79,90,105,110	0
2	FUC	D	9	10/11	0.67	0.61	116,119,121,122	0
2	NAG	C	8	14/15	0.68	0.28	92,98,107,109	0
2	NAG	D	8	14/15	0.69	0.41	106,109,112,112	0
2	BMA	D	3	11/12	0.75	0.16	85,89,92,95	0
2	MAN	D	7	11/12	0.78	0.25	82,88,94,100	0
2	MAN	D	4	11/12	0.83	0.14	74,79,83,84	0
2	MAN	C	7	11/12	0.85	0.16	65,69,76,87	0
2	BMA	C	3	11/12	0.86	0.13	47,51,58,63	0
2	GAL	D	6	11/12	0.87	0.20	70,76,77,79	0
2	FUC	C	9	10/11	0.87	0.13	65,67,69,69	0
2	GAL	C	6	11/12	0.88	0.20	64,69,84,87	0
2	NAG	C	5	14/15	0.89	0.10	55,61,67,69	0
2	MAN	C	4	11/12	0.90	0.09	49,51,57,57	0
2	NAG	D	5	14/15	0.91	0.16	77,80,84,85	0
2	NAG	C	1	14/15	0.95	0.06	46,54,64,69	0
2	NAG	C	2	14/15	0.95	0.10	46,51,55,55	0

The following is a graphical depiction of the model fit to experimental electron density for oligosac-

charide. Each fit is shown from different orientation to approximate a three-dimensional view.





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
3	MPD	B	510	8/8	0.81	0.35	97,100,103,115	0
3	MPD	A	510	8/8	0.82	0.17	70,74,83,87	0

6.5 Other polymers [i](#)

There are no such residues in this entry.