



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:51 PM UTC

PDB ID : 4W4N / pdb_00004w4n
Title : Crystal structure of human Fc at 1.80 Å
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Deposited on : 2014-08-15
Resolution : 1.80 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

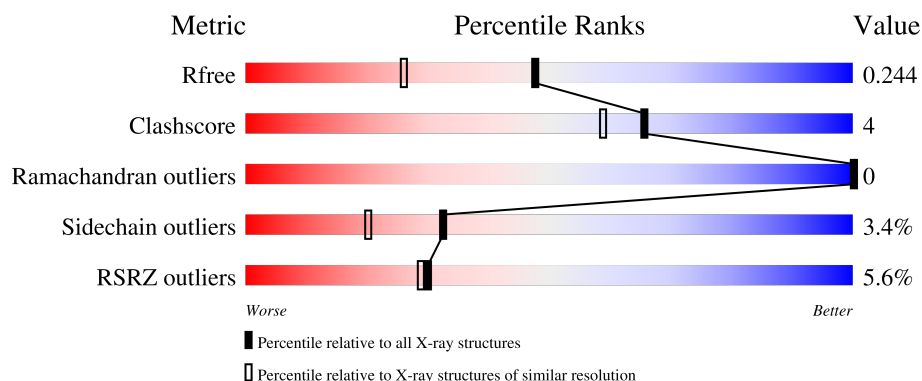
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	10% 81% 10% 8%
1	B	223	81% 9% 7%
2	C	7	100%
3	D	8	100%

2 Entry composition [i](#)

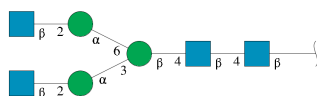
There are 5 unique types of molecules in this entry. The entry contains 3931 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ig gamma-1 chain C region.

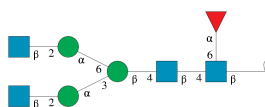
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	0	3	0
			1666	1062	279	319	6			
1	B	207	Total	C	N	O	S	0	5	0
			1688	1077	283	322	6			

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]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
2	C	7	Total	C	N	O	0	0	0
			89	50	4	35			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]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
3	D	8	Total	C	N	O	0	0	0
			99	56	4	39			

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	1
			12	6	6		

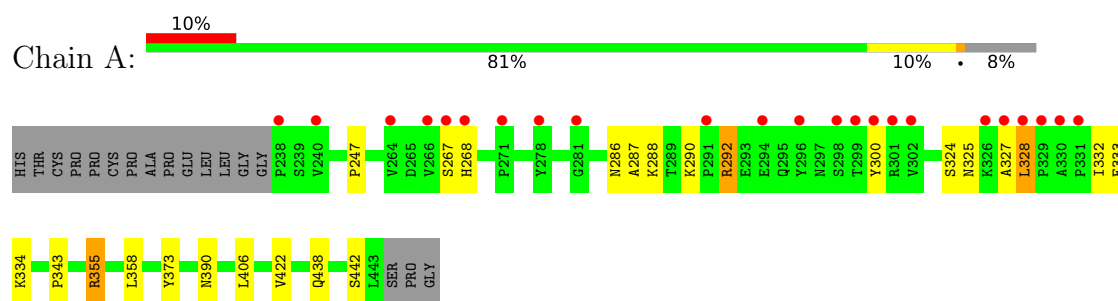
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	137	Total	O	0	0
			137	137		
5	B	237	Total	O	0	3
			240	240		

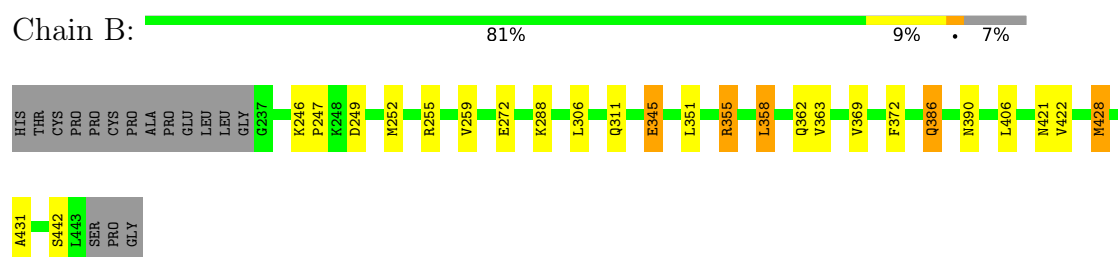
3 Residue-property plots

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: Ig gamma-1 chain C region



- Molecule 1: Ig gamma-1 chain C region



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



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]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
MAN6
MAG7
FUC8

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.71Å 79.29Å 138.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.19 – 1.80 22.19 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.9 (22.19-1.80) 98.9 (22.19-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.200 , 0.244 0.200 , 0.244	Depositor DCC
R_{free} test set	2087 reflections (4.04%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3931	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.21% 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, GOL, FUC, MAN, BMA

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	1.23	2/1721 (0.1%)	1.16	4/2344 (0.2%)
1	B	1.24	2/1749 (0.1%)	1.21	8/2383 (0.3%)
All	All	1.24	4/3470 (0.1%)	1.19	12/4727 (0.3%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	267	SER	CB-OG	5.88	1.53	1.42
1	B	369	VAL	C-O	-5.74	1.17	1.24
1	A	390	ASN	CA-C	5.44	1.60	1.52
1	B	372	PHE	N-CA	5.41	1.52	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	286	ASN	N-CA-C	7.92	122.94	113.28
1	B	345[A]	GLU	CA-C-N	6.49	126.45	119.76
1	B	345[A]	GLU	C-N-CA	6.49	126.45	119.76
1	B	345[B]	GLU	CA-C-N	6.49	126.45	119.76
1	B	345[B]	GLU	C-N-CA	6.49	126.45	119.76
1	A	328	LEU	CA-C-N	5.72	125.19	119.24
1	A	328	LEU	C-N-CA	5.72	125.19	119.24
1	B	428	MET	CG-SD-CE	-5.62	88.53	100.90
1	B	386	GLN	CA-C-N	-5.18	114.62	119.85
1	B	386	GLN	C-N-CA	-5.18	114.62	119.85
1	A	287	ALA	N-CA-C	5.09	117.67	110.50
1	B	390	ASN	N-CA-C	5.03	117.16	109.62

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	1666	0	1641	15	0
1	B	1688	0	1667	12	0
2	C	89	0	76	0	0
3	D	99	0	85	0	0
4	B	12	0	16	0	0
5	A	137	0	0	3	0
5	B	240	0	0	1	0
All	All	3931	0	3485	27	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 4.

All (27) 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:A:333:GLU:O	1:A:334:LYS:HD2	1.89	0.72
1:A:438:GLN:NE2	5:A:601:HOH:O	2.27	0.65
1:B:252:MET:SD	1:B:428:MET:HE1	2.38	0.64
1:A:328:LEU:HD22	1:A:332:ILE:HD11	1.80	0.64
1:A:292:ARG:H	1:A:292:ARG:HH11	1.47	0.60
1:A:406:LEU:C	1:A:406:LEU:HD12	2.32	0.54
1:B:422:VAL:HA	1:B:442:SER:HB3	1.89	0.54
1:B:406:LEU:HD12	1:B:406:LEU:C	2.33	0.54
1:A:438:GLN:HG3	5:A:639:HOH:O	2.12	0.50
1:A:325:ASN:ND2	1:A:327:ALA:H	2.10	0.49
1:A:343:PRO:HA	1:A:373:TYR:O	2.13	0.49
1:A:268:HIS:CD2	1:A:300:TYR:CE1	3.01	0.48
1:B:421:ASN:ND2	5:B:601:HOH:O	2.42	0.48
1:B:246:LYS:HB2	1:B:249:ASP:OD2	2.14	0.47
1:B:259[B]:VAL:HG23	1:B:306:LEU:HB3	1.97	0.47
1:A:328:LEU:HD13	1:A:332:ILE:HD12	1.98	0.46
1:A:333:GLU:C	5:A:632:HOH:O	2.60	0.45
1:A:355:ARG:HA	1:A:358:LEU:HD12	1.99	0.44
1:B:358:LEU:HD13	1:B:363:VAL:HG11	1.99	0.44
1:B:249:ASP:OD1	1:B:255[A]:ARG:NE	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:345[A]:GLU:HG3	1:B:431:ALA:O	2.18	0.43
1:B:358:LEU:HD13	1:B:358:LEU:HA	1.77	0.43
1:A:355:ARG:HE	1:A:355:ARG:HB2	1.40	0.43
1:A:290:LYS:HA	1:A:290:LYS:HD2	1.92	0.42
1:B:355:ARG:HA	1:B:358:LEU:HD23	2.02	0.42
1:A:422:VAL:HA	1:A:442:SER:HB3	2.02	0.42
1:B:272[A]:GLU:CD	1:B:272[A]:GLU:H	2.28	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	207/223 (93%)	202 (98%)	5 (2%)	0	100	100
1	B	210/223 (94%)	208 (99%)	2 (1%)	0	100	100
All	All	417/446 (94%)	410 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	196/206 (95%)	191 (97%)	5 (3%)	40	28

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	198/206 (96%)	190 (96%)	8 (4%)	28	15
All	All	394/412 (96%)	381 (97%)	13 (3%)	32	21

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	247	PRO
1	A	288	LYS
1	A	292	ARG
1	A	324	SER
1	A	355	ARG
1	B	247	PRO
1	B	288	LYS
1	B	311	GLN
1	B	351	LEU
1	B	355	ARG
1	B	358	LEU
1	B	362	GLN
1	B	386	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	HIS
1	A	276	ASN
1	A	285	HIS
1	A	325	ASN
1	A	362	GLN
1	A	390	ASN
1	A	419	GLN
1	A	421	ASN
1	A	434	ASN
1	B	285	HIS
1	B	315	ASN
1	B	361	ASN
1	B	386	GLN
1	B	390	ASN
1	B	433	HIS
1	B	434	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 ⓘ

15 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	2,1	14,14,15	0.83	0	17,19,21	3.13	7 (41%)
2	NAG	C	2	2	14,14,15	0.79	0	17,19,21	1.30	2 (11%)
2	BMA	C	3	2	11,11,12	1.08	1 (9%)	15,15,17	1.60	3 (20%)
2	MAN	C	4	2	11,11,12	1.00	0	15,15,17	1.90	4 (26%)
2	NAG	C	5	2	14,14,15	0.50	0	17,19,21	2.26	3 (17%)
2	MAN	C	6	2	11,11,12	0.68	0	15,15,17	1.18	1 (6%)
2	NAG	C	7	2	14,14,15	0.51	0	17,19,21	2.81	6 (35%)
3	NAG	D	1	3,1	14,14,15	1.15	2 (14%)	17,19,21	0.93	0
3	NAG	D	2	3	14,14,15	1.16	2 (14%)	17,19,21	1.81	5 (29%)
3	BMA	D	3	3	11,11,12	1.04	1 (9%)	15,15,17	1.36	2 (13%)
3	MAN	D	4	3	11,11,12	1.22	2 (18%)	15,15,17	1.88	4 (26%)
3	NAG	D	5	3	14,14,15	0.60	0	17,19,21	2.06	7 (41%)
3	MAN	D	6	3	11,11,12	0.86	0	15,15,17	1.86	4 (26%)
3	NAG	D	7	3	14,14,15	1.48	2 (14%)	17,19,21	1.63	3 (17%)
3	FUC	D	8	3	10,10,11	1.24	2 (20%)	14,14,16	1.10	2 (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
2	NAG	C	1	2,1	-	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	-	1/2/19/22	0/1/1/1
2	NAG	C	5	2	-	3/6/23/26	0/1/1/1
2	MAN	C	6	2	-	0/2/19/22	0/1/1/1
2	NAG	C	7	2	-	2/6/23/26	0/1/1/1
3	NAG	D	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
3	BMA	D	3	3	-	0/2/19/22	0/1/1/1
3	MAN	D	4	3	-	0/2/19/22	0/1/1/1
3	NAG	D	5	3	-	2/6/23/26	0/1/1/1
3	MAN	D	6	3	-	0/2/19/22	0/1/1/1
3	NAG	D	7	3	-	0/6/23/26	0/1/1/1
3	FUC	D	8	3	-	-	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	7	NAG	C1-C2	4.08	1.57	1.52
3	D	2	NAG	C3-C2	3.10	1.59	1.52
3	D	8	FUC	C6-C5	2.71	1.58	1.51
3	D	1	NAG	C1-C2	2.40	1.55	1.52
2	C	3	BMA	O5-C1	-2.31	1.39	1.43
3	D	8	FUC	C1-C2	2.26	1.57	1.52
3	D	1	NAG	O5-C5	-2.21	1.39	1.43
3	D	4	MAN	C4-C5	2.17	1.57	1.53
3	D	2	NAG	O7-C7	2.13	1.28	1.23
3	D	7	NAG	C4-C5	2.07	1.57	1.53
3	D	3	BMA	C2-C3	2.06	1.55	1.52
3	D	4	MAN	O5-C1	-2.02	1.40	1.43

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	7	NAG	C1-O5-C5	8.47	123.54	112.19
2	C	1	NAG	C2-N2-C7	7.40	132.81	122.90
2	C	5	NAG	C1-O5-C5	6.76	121.24	112.19
2	C	1	NAG	C1-C2-N2	6.23	120.25	110.43
2	C	4	MAN	C1-O5-C5	5.12	119.05	112.19
2	C	7	NAG	C6-C5-C4	-4.86	101.09	113.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C1-O5-C5	4.81	118.63	112.19
3	D	5	NAG	C2-N2-C7	4.80	129.33	122.90
3	D	6	MAN	C1-O5-C5	4.70	118.49	112.19
3	D	4	MAN	O2-C2-C3	-4.32	101.20	110.15
2	C	1	NAG	O7-C7-C8	-3.92	115.07	122.05
3	D	7	NAG	C1-O5-C5	-3.85	107.03	112.19
3	D	2	NAG	O7-C7-N2	3.80	128.69	121.98
2	C	5	NAG	C4-C3-C2	-3.76	105.50	111.02
3	D	2	NAG	C8-C7-N2	-3.68	110.01	116.12
3	D	5	NAG	C1-O5-C5	3.51	116.88	112.19
3	D	3	BMA	O3-C3-C2	-3.35	103.22	110.05
2	C	3	BMA	O6-C6-C5	-3.29	100.12	111.33
2	C	4	MAN	O2-C2-C3	-3.20	103.52	110.15
2	C	6	MAN	C1-O5-C5	3.20	116.47	112.19
2	C	5	NAG	C8-C7-N2	3.18	121.40	116.12
2	C	3	BMA	O4-C4-C3	-3.12	103.02	110.38
2	C	7	NAG	O5-C5-C4	3.11	118.40	110.83
2	C	7	NAG	O3-C3-C4	-3.02	103.25	110.38
3	D	4	MAN	C1-C2-C3	2.96	113.95	109.64
3	D	6	MAN	O5-C5-C6	2.90	113.32	107.66
3	D	4	MAN	C1-O5-C5	2.90	116.07	112.19
2	C	1	NAG	O7-C7-N2	2.86	127.04	121.98
2	C	1	NAG	O5-C5-C6	2.77	113.06	107.66
3	D	7	NAG	C1-C2-N2	-2.60	106.33	110.43
3	D	2	NAG	O5-C1-C2	-2.57	107.32	111.29
3	D	5	NAG	O7-C7-N2	2.49	126.39	121.98
3	D	8	FUC	O5-C1-C2	-2.48	104.88	110.79
3	D	6	MAN	C6-C5-C4	-2.47	106.97	113.02
3	D	4	MAN	O4-C4-C3	-2.42	104.67	110.38
3	D	6	MAN	O6-C6-C5	-2.40	103.15	111.33
3	D	3	BMA	O2-C2-C3	2.39	115.11	110.15
3	D	7	NAG	O5-C1-C2	-2.36	107.64	111.29
3	D	5	NAG	C6-C5-C4	-2.35	107.24	113.02
2	C	3	BMA	C1-C2-C3	2.34	113.05	109.64
2	C	7	NAG	O6-C6-C5	-2.32	103.44	111.33
3	D	2	NAG	O4-C4-C5	-2.29	103.70	109.32
2	C	4	MAN	O3-C3-C4	-2.22	105.14	110.38
3	D	2	NAG	C2-N2-C7	-2.21	119.94	122.90
2	C	2	NAG	C1-O5-C5	2.20	115.13	112.19
2	C	2	NAG	O7-C7-N2	2.18	125.83	121.98
3	D	5	NAG	O3-C3-C2	-2.16	104.91	109.40
3	D	8	FUC	C3-C4-C5	-2.11	106.61	109.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	7	NAG	C3-C4-C5	2.10	114.04	110.23
2	C	1	NAG	C4-C3-C2	2.06	114.04	111.02
3	D	5	NAG	O7-C7-C8	-2.04	118.42	122.05
3	D	5	NAG	O6-C6-C5	-2.04	104.38	111.33
2	C	4	MAN	O4-C4-C5	2.04	114.34	109.32

There are no chirality outliers.

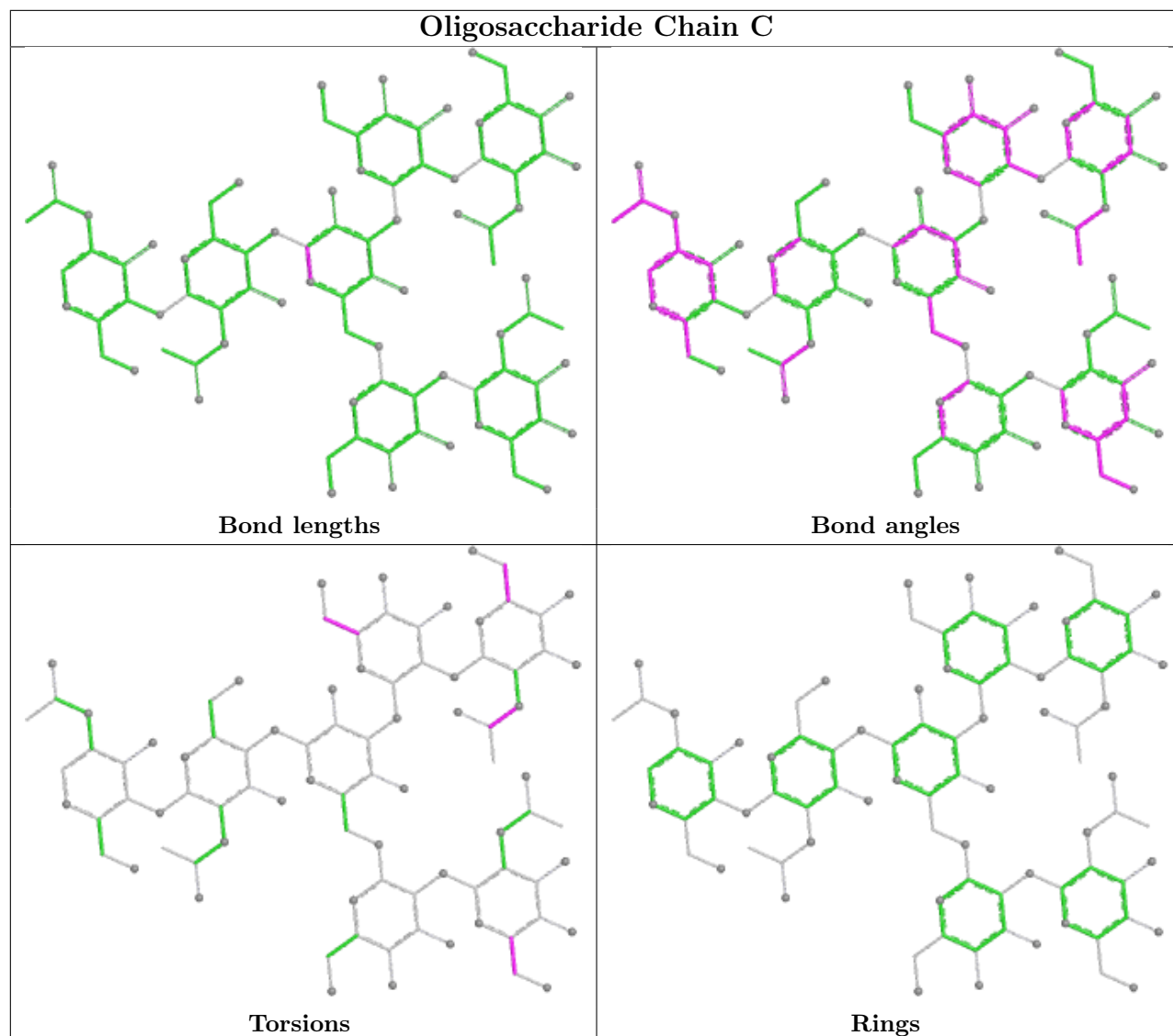
All (8) torsion outliers are listed below:

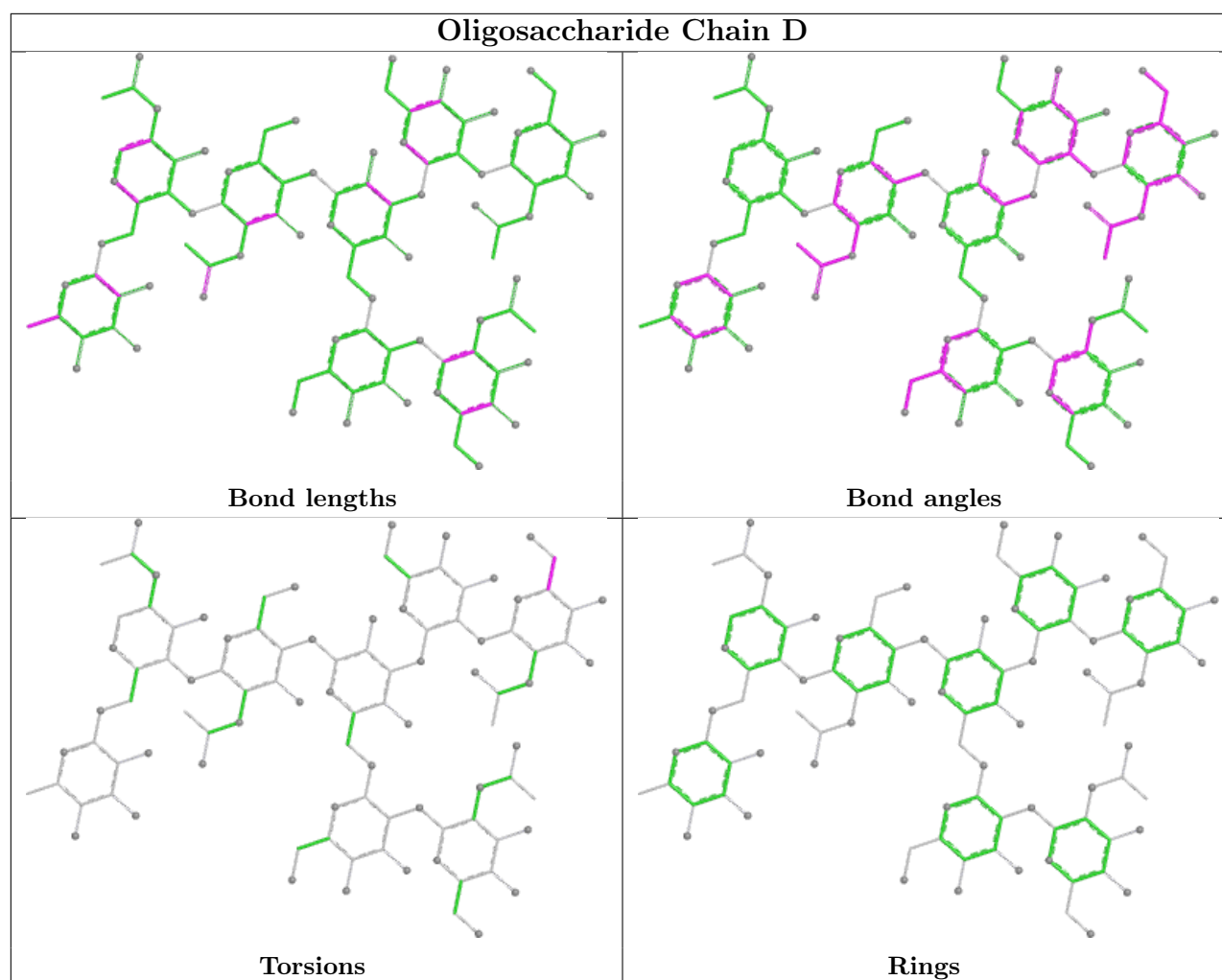
Mol	Chain	Res	Type	Atoms
2	C	7	NAG	O5-C5-C6-O6
2	C	7	NAG	C4-C5-C6-O6
2	C	5	NAG	C8-C7-N2-C2
2	C	5	NAG	O7-C7-N2-C2
3	D	5	NAG	C4-C5-C6-O6
2	C	5	NAG	O5-C5-C6-O6
3	D	5	NAG	O5-C5-C6-O6
2	C	4	MAN	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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	$\# Z > 2$
4	GOL	B	509[A]	-	5,5,5	0.17	0	5,5,5	0.64	0
4	GOL	B	509[B]	-	5,5,5	0.23	0	5,5,5	0.73	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	GOL	B	509[A]	-	-	0/4/4/4	-
4	GOL	B	509[B]	-	-	2/4/4/4	-

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
4	B	509[B]	GOL	O1-C1-C2-O2
4	B	509[B]	GOL	O1-C1-C2-C3

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	206/223 (92%)	0.41	23 (11%) 10 8	17, 45, 104, 135	3 (1%)
1	B	207/223 (92%)	-0.32	0 100 100	11, 30, 63, 85	5 (2%)
All	All	413/446 (92%)	0.04	23 (5%) 30 29	11, 35, 94, 135	8 (1%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	296	TYR	5.5
1	A	300	TYR	5.2
1	A	271	PRO	4.1
1	A	328	LEU	3.6
1	A	291	PRO	3.6
1	A	327	ALA	3.4
1	A	238	PRO	3.3
1	A	266	VAL	3.1
1	A	240	VAL	2.9
1	A	302	VAL	2.8
1	A	268	HIS	2.7
1	A	278	TYR	2.5
1	A	301	ARG	2.5
1	A	267	SER	2.4
1	A	331	PRO	2.4
1	A	264	VAL	2.4
1	A	330	ALA	2.3
1	A	299	THR	2.3
1	A	281	GLY	2.3
1	A	294	GLU	2.2
1	A	326	LYS	2.2
1	A	298	SER	2.2
1	A	329	PRO	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 ⓘ

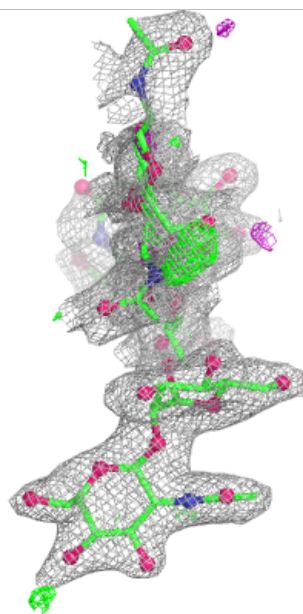
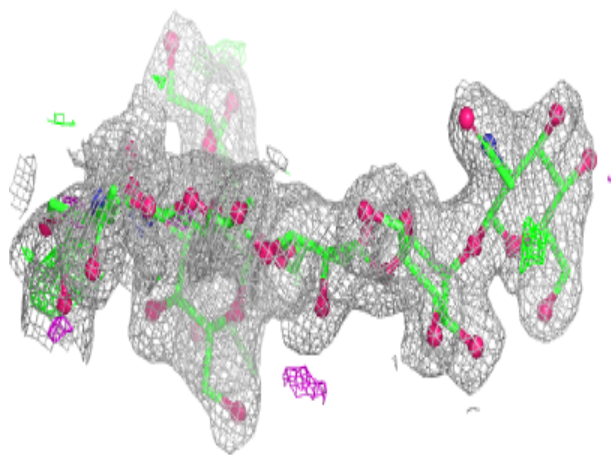
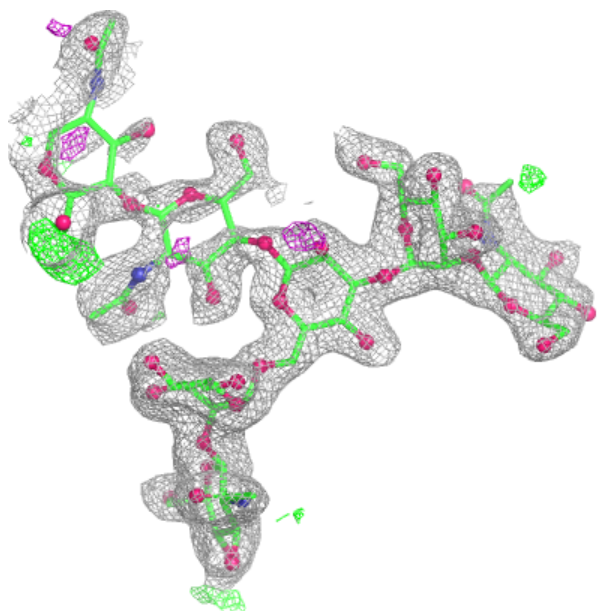
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

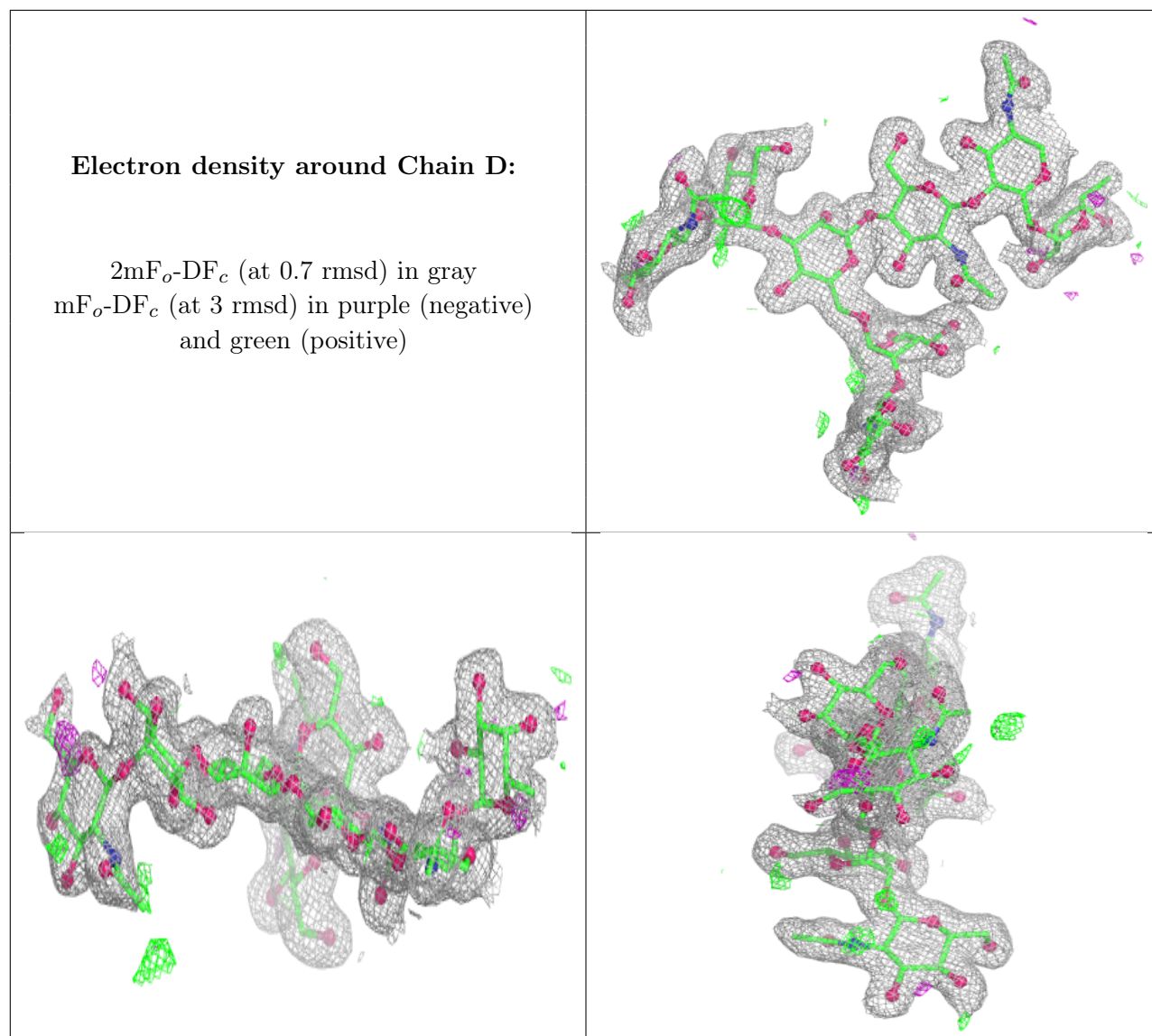
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	C	1	14/15	0.70	0.17	73,81,89,90	0
2	NAG	C	5	14/15	0.76	0.12	61,71,80,82	0
3	NAG	D	5	14/15	0.76	0.13	45,48,53,54	0
2	NAG	C	2	14/15	0.85	0.13	54,61,69,70	0
2	NAG	C	7	14/15	0.88	0.10	54,57,60,62	0
3	FUC	D	8	10/11	0.92	0.08	30,34,37,42	0
2	MAN	C	4	11/12	0.93	0.08	45,47,53,58	0
2	BMA	C	3	11/12	0.93	0.08	46,47,51,52	0
2	MAN	C	6	11/12	0.93	0.08	51,57,61,65	0
3	NAG	D	7	14/15	0.94	0.07	24,26,30,33	0
3	MAN	D	4	11/12	0.95	0.06	28,32,39,42	0
3	NAG	D	1	14/15	0.96	0.06	27,30,34,34	0
3	MAN	D	6	11/12	0.97	0.05	25,27,31,32	0
3	NAG	D	2	14/15	0.97	0.05	24,25,27,28	0
3	BMA	D	3	11/12	0.97	0.05	23,25,26,26	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 C:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	509[A]	6/6	0.83	0.12	38,49,52,52	6
4	GOL	B	509[B]	6/6	0.83	0.12	31,47,48,51	6

6.5 Other polymers [i](#)

There are no such residues in this entry.