



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 01:39 AM BST

PDB ID : 4X99
Title : Immunoglobulin Fc heterodimers variant
Authors : Seok, S.H.; Choi, H.J.; Kim, Y.J.; Seo, M.D.; Kim, Y.S.
Deposited on : 2014-12-11
Resolution : 2.50 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 : 2.13.1
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) : 2.13.1

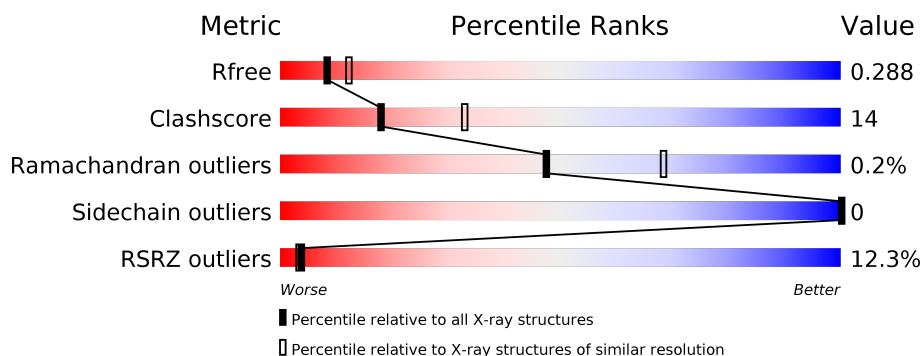
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 on 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>72%</div> <div>21%</div> <div>7%</div> </div>
2	B	223	<div> <div>22%</div> <div>68%</div> <div>23%</div> <div>8%</div> </div>
3	C	8	<div> <div>63%</div> <div>38%</div> </div>
3	D	8	<div> <div>38%</div> <div>63%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	D	1	-	-	-	X
3	NAG	D	2	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3502 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	208	Total	C	N	O	S	0	0	0
			1663	1057	279	320	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	349	CYS	TYR	engineered mutation	UNP P01857
A	360	GLU	LYS	engineered mutation	UNP P01857
A	409	TRP	LYS	engineered mutation	UNP P01857

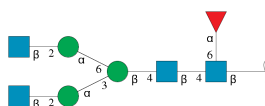
- Molecule 2 is a protein called Ig gamma-1 chain C region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	205	Total	C	N	O	S	0	0	0
			1600	1018	273	302	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	347	ARG	GLN	engineered mutation	UNP P01857
B	354	CYS	SER	engineered mutation	UNP P01857
B	399	VAL	ASP	engineered mutation	UNP P01857
B	405	THR	PHE	engineered mutation	UNP P01857

- 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	C	8	Total	C	N	O	0	0	0
			99	56	4	39			
3	D	8	Total	C	N	O	0	0	0
			99	56	4	39			

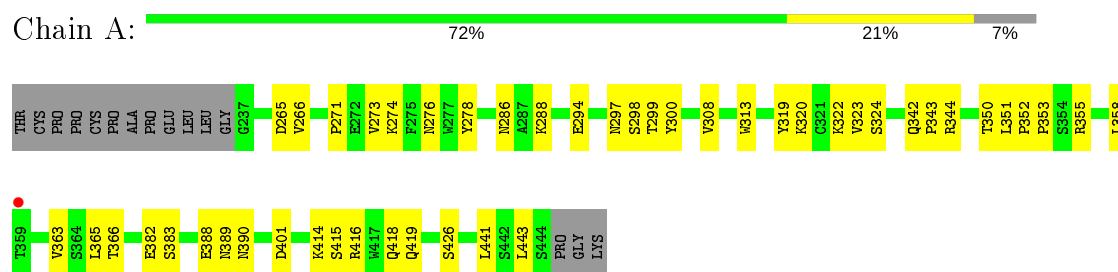
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	30	Total	O	0	0
			30	30		
4	B	11	Total	O	0	0
			11	11		

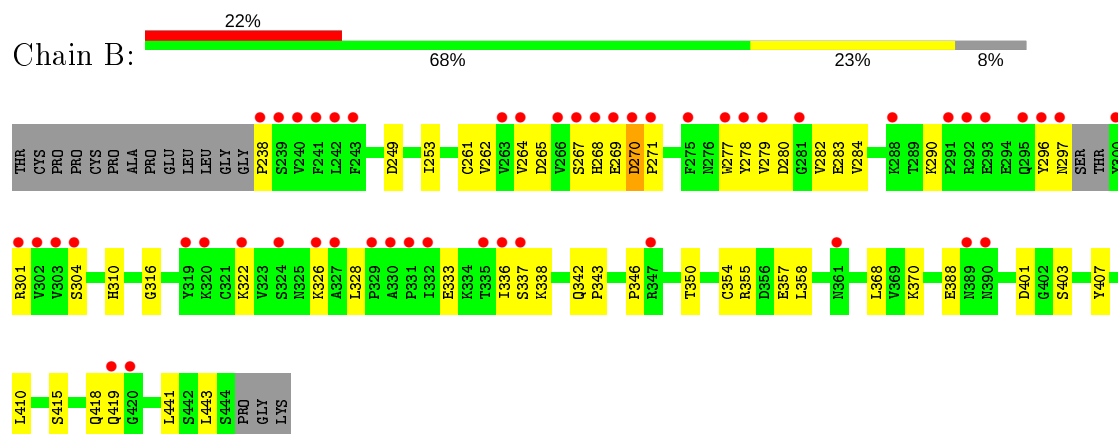
3 Residue-property plots

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 2: Ig gamma-1 chain C region



- 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



- 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-mannopyra

nose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:



MAG1
MAG2
BOA3
MAT4
MAG5
MAG6
MAG7
FUC8

4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	152.77Å 152.77Å 109.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.63 – 2.50 45.52 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (44.63-2.50) 93.4 (45.52-2.50)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.62 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.248 , 0.283 0.263 , 0.288	Depositor DCC
R_{free} test set	2000 reflections (7.49%)	wwPDB-VP
Wilson B-factor (Å ²)	51.0	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3502	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: FUC, BMA, NAG, MAN

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	0.63	0/1710	0.70	0/2333
2	B	0.63	0/1642	0.78	3/2241 (0.1%)
All	All	0.63	0/3352	0.74	3/4574 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	268	HIS	CB-CA-C	-10.77	88.87	110.40
2	B	290	LYS	C-N-CD	5.86	140.70	128.40
2	B	270	ASP	C-N-CD	5.59	140.15	128.40

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	1663	0	1614	30	0
2	B	1600	0	1555	47	0
3	C	99	0	83	7	0
3	D	99	0	85	10	0
4	A	30	0	0	1	0
4	B	11	0	0	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3502	0	3337	92	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 (92) 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:278:TYR:CE1	2:B:283:GLU:HB2	1.30	1.56
2:B:278:TYR:CE1	2:B:283:GLU:CB	2.23	1.20
2:B:267:SER:HB2	2:B:271:PRO:HG2	1.33	1.05
1:A:265:ASP:OD1	3:C:1:NAG:O7	1.83	0.96
2:B:278:TYR:HE1	2:B:283:GLU:HB2	1.28	0.94
3:C:4:MAN:H4	3:C:5:NAG:H83	1.50	0.94
2:B:278:TYR:CD1	2:B:283:GLU:CA	2.55	0.90
2:B:278:TYR:CZ	2:B:283:GLU:HB2	2.06	0.88
2:B:267:SER:O	2:B:271:PRO:HD2	1.75	0.87
2:B:278:TYR:CD1	2:B:283:GLU:HB2	2.09	0.86
3:D:1:NAG:H62	3:D:8:FUC:H5	1.59	0.83
1:A:350:THR:HB	1:A:441:LEU:HD13	1.61	0.81
1:A:294:GLU:OE2	1:A:298:SER:HA	1.84	0.77
2:B:269:GLU:C	2:B:271:PRO:HD3	2.06	0.76
2:B:278:TYR:CD1	2:B:283:GLU:N	2.55	0.75
2:B:253:ILE:H	2:B:253:ILE:HD12	1.52	0.74
2:B:278:TYR:HD1	2:B:282:VAL:C	1.90	0.74
3:C:4:MAN:O2	3:C:5:NAG:H83	1.87	0.73
2:B:278:TYR:CD1	2:B:283:GLU:HA	2.21	0.73
2:B:278:TYR:HD1	2:B:283:GLU:N	1.87	0.72
3:D:1:NAG:H62	3:D:8:FUC:C5	2.21	0.70
1:A:418:GLN:HA	1:A:443:LEU:HD22	1.75	0.69
2:B:278:TYR:CD1	2:B:283:GLU:CB	2.69	0.69
3:D:1:NAG:C6	3:D:8:FUC:H5	2.23	0.68
3:C:4:MAN:H4	3:C:5:NAG:C8	2.22	0.68
3:D:4:MAN:H4	3:D:5:NAG:H4	1.78	0.66
1:A:415:SER:O	1:A:419:GLN:HG3	1.96	0.66
3:D:4:MAN:O3	3:D:5:NAG:H61	1.96	0.65
1:A:274:LYS:HB3	1:A:324:SER:HB2	1.79	0.65
1:A:353:PRO:HB3	1:A:363:VAL:HG13	1.79	0.64
1:A:355:ARG:HA	1:A:358:LEU:HD12	1.79	0.64
2:B:265:ASP:OD1	3:D:1:NAG:H83	1.98	0.63
2:B:279:VAL:O	2:B:280:ASP:HB2	2.01	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:VAL:HG11	1:A:313:TRP:HB2	1.83	0.60
2:B:354:CYS:SG	2:B:357:GLU:HB2	2.42	0.60
3:C:4:MAN:C4	3:C:5:NAG:H83	2.30	0.59
2:B:279:VAL:O	2:B:279:VAL:HG23	2.02	0.58
2:B:262:VAL:CG1	2:B:301:ARG:HH21	2.17	0.57
1:A:308:VAL:HG22	1:A:319:TYR:CE1	2.40	0.57
2:B:267:SER:HB2	2:B:271:PRO:CG	2.21	0.57
2:B:261:CYS:HB2	2:B:277:TRP:CH2	2.41	0.55
1:A:266:VAL:HB	1:A:300:TYR:HB2	1.88	0.54
1:A:350:THR:CB	1:A:441:LEU:HD13	2.35	0.54
2:B:418:GLN:HA	2:B:443:LEU:HD22	1.90	0.54
1:A:388:GLU:OE2	1:A:416:ARG:NH1	2.40	0.54
2:B:322:LYS:HG3	2:B:333:GLU:HG2	1.90	0.54
3:C:4:MAN:O2	3:C:5:NAG:O6	2.27	0.53
3:D:7:NAG:O3	3:D:7:NAG:O7	2.26	0.53
3:D:5:NAG:H5	3:D:5:NAG:O7	2.08	0.53
2:B:316:GLY:O	2:B:337:SER:HB2	2.10	0.51
3:D:1:NAG:C6	3:D:8:FUC:C5	2.85	0.51
2:B:238:PRO:HB2	2:B:328:LEU:HD11	1.93	0.50
2:B:278:TYR:CE1	2:B:283:GLU:CA	2.84	0.50
1:A:297:ASN:OD1	1:A:299:THR:OG1	2.27	0.50
2:B:238:PRO:HA	2:B:264:VAL:O	2.14	0.48
2:B:338:LYS:HE3	4:B:606:HOH:O	2.13	0.48
1:A:351:LEU:HB2	1:A:366:THR:HB	1.95	0.48
2:B:262:VAL:HG11	2:B:301:ARG:HH21	1.77	0.48
2:B:355:ARG:HA	2:B:358:LEU:HD12	1.96	0.48
3:D:6:MAN:O3	3:D:7:NAG:N2	2.48	0.47
1:A:308:VAL:CG1	1:A:313:TRP:HB2	2.44	0.47
1:A:271:PRO:HD2	4:A:628:HOH:O	2.15	0.47
2:B:388:GLU:HG2	2:B:410:LEU:HD11	1.95	0.47
2:B:253:ILE:CD1	2:B:253:ILE:H	2.24	0.47
2:B:401:ASP:HB3	2:B:403:SER:H	1.80	0.47
1:A:358:LEU:O	1:A:414:LYS:NZ	2.23	0.47
2:B:346:PRO:HA	2:B:370:LYS:O	2.17	0.45
2:B:350:THR:OG1	2:B:441:LEU:HD22	2.17	0.45
1:A:382:GLU:OE1	1:A:426:SER:OG	2.35	0.44
2:B:262:VAL:HG13	2:B:301:ARG:HH21	1.82	0.44
1:A:286:ASN:O	1:A:288:LYS:NZ	2.37	0.44
2:B:368:LEU:HD13	2:B:407:TYR:CZ	2.53	0.44
1:A:342:GLN:HA	1:A:343:PRO:HD3	1.91	0.44
1:A:344:ARG:HH12	1:A:401:ASP:CG	2.21	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:ASN:HB2	1:A:322:LYS:HB3	2.01	0.43
2:B:296:TYR:O	2:B:297:ASN:C	2.56	0.43
2:B:283:GLU:HG2	2:B:284:VAL:N	2.34	0.43
1:A:353:PRO:HD3	1:A:365:LEU:HD23	2.01	0.43
1:A:383:SER:HB2	1:A:388:GLU:OE1	2.19	0.43
2:B:415:SER:O	2:B:419:GLN:HG3	2.19	0.42
1:A:308:VAL:HG22	1:A:319:TYR:CZ	2.54	0.42
2:B:337:SER:O	2:B:338:LYS:C	2.57	0.42
1:A:278:TYR:CD1	1:A:320:LYS:HD2	2.55	0.42
2:B:336:ILE:O	2:B:337:SER:HB3	2.20	0.41
2:B:249:ASP:O	2:B:310:HIS:HE1	2.03	0.41
1:A:389:ASN:O	1:A:390:ASN:HB2	2.20	0.41
1:A:273:VAL:HG13	1:A:323:VAL:HG13	2.03	0.41
1:A:351:LEU:HA	1:A:352:PRO:HD2	1.83	0.41
2:B:342:GLN:OE1	2:B:343:PRO:HD2	2.20	0.41
2:B:277:TRP:HZ2	2:B:304:SER:HB3	1.86	0.40
2:B:326:LYS:C	2:B:328:LEU:H	2.24	0.40
3:C:1:NAG:O7	3:C:1:NAG:C3	2.70	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	206/223 (92%)	200 (97%)	6 (3%)	0	100	100
2	B	201/223 (90%)	188 (94%)	12 (6%)	1 (0%)	29	48
All	All	407/446 (91%)	388 (95%)	18 (4%)	1 (0%)	47	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	270	ASP

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	194/206 (94%)	194 (100%)	0	100	100
2	B	183/206 (89%)	183 (100%)	0	100	100
All	All	377/412 (92%)	377 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

16 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	C	1	1,3	14,14,15	1.31	1 (7%)	17,19,21	1.82	4 (23%)
3	NAG	C	2	3	14,14,15	1.32	1 (7%)	17,19,21	1.39	3 (17%)
3	BMA	C	3	3	11,11,12	1.61	2 (18%)	15,15,17	1.62	3 (20%)
3	MAN	C	4	3	11,11,12	1.41	2 (18%)	15,15,17	2.16	7 (46%)
3	NAG	C	5	3	14,14,15	0.97	1 (7%)	17,19,21	1.96	5 (29%)
3	MAN	C	6	3	11,11,12	2.00	4 (36%)	15,15,17	2.22	4 (26%)
3	NAG	C	7	3	14,14,15	1.42	2 (14%)	17,19,21	1.55	4 (23%)
3	FUC	C	8	3	10,10,11	1.87	5 (50%)	14,14,16	1.18	0
3	NAG	D	1	3,2	14,14,15	0.79	0	17,19,21	0.80	0
3	NAG	D	2	3	14,14,15	0.91	1 (7%)	17,19,21	1.34	3 (17%)
3	BMA	D	3	3	11,11,12	1.26	1 (9%)	15,15,17	2.23	4 (26%)
3	MAN	D	4	3	11,11,12	0.80	0	15,15,17	1.96	3 (20%)
3	NAG	D	5	3	14,14,15	1.41	4 (28%)	17,19,21	1.83	4 (23%)
3	MAN	D	6	3	11,11,12	1.04	1 (9%)	15,15,17	1.90	4 (26%)
3	NAG	D	7	3	14,14,15	1.00	1 (7%)	17,19,21	1.42	3 (17%)
3	FUC	D	8	3	10,10,11	1.17	1 (10%)	14,14,16	1.55	5 (35%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	1	1,3	-	1/6/23/26	0/1/1/1
3	NAG	C	2	3	-	0/6/23/26	0/1/1/1
3	BMA	C	3	3	-	0/2/19/22	0/1/1/1
3	MAN	C	4	3	-	2/2/19/22	0/1/1/1
3	NAG	C	5	3	-	3/6/23/26	0/1/1/1
3	MAN	C	6	3	-	2/2/19/22	0/1/1/1
3	NAG	C	7	3	-	2/6/23/26	0/1/1/1
3	FUC	C	8	3	-	-	0/1/1/1
3	NAG	D	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	2/6/23/26	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	-	2/2/19/22	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	D	7	3	-	3/6/23/26	0/1/1/1
3	FUC	D	8	3	-	-	0/1/1/1

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	6	MAN	O3-C3	-3.50	1.34	1.43
3	C	8	FUC	O5-C1	-3.29	1.38	1.43
3	C	6	MAN	O5-C1	-3.20	1.38	1.43
3	C	7	NAG	O4-C4	-3.06	1.35	1.43
3	C	1	NAG	C1-C2	-3.03	1.47	1.52
3	D	5	NAG	C3-C2	2.96	1.58	1.52
3	C	2	NAG	O5-C1	-2.83	1.39	1.43
3	C	3	BMA	O2-C2	-2.64	1.37	1.43
3	D	8	FUC	C4-C5	-2.63	1.46	1.52
3	C	6	MAN	C2-C3	-2.61	1.48	1.52
3	D	3	BMA	C4-C3	-2.55	1.45	1.52
3	C	3	BMA	O3-C3	-2.47	1.37	1.43
3	D	2	NAG	C1-C2	-2.42	1.48	1.52
3	D	5	NAG	C4-C5	2.39	1.58	1.53
3	C	7	NAG	C4-C5	-2.35	1.48	1.53
3	C	8	FUC	C4-C5	-2.29	1.47	1.52
3	C	5	NAG	C3-C2	-2.28	1.47	1.52
3	C	8	FUC	O4-C4	-2.27	1.37	1.43
3	C	6	MAN	C4-C5	-2.24	1.48	1.53
3	D	6	MAN	O5-C1	2.22	1.47	1.43
3	C	8	FUC	O3-C3	-2.14	1.37	1.43
3	C	8	FUC	O2-C2	-2.13	1.38	1.43
3	D	5	NAG	C1-C2	2.09	1.55	1.52
3	D	5	NAG	O5-C5	2.05	1.47	1.43
3	C	4	MAN	O2-C2	-2.04	1.39	1.43
3	C	4	MAN	O5-C1	-2.01	1.40	1.43
3	D	7	NAG	C1-C2	-2.00	1.49	1.52

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	6	MAN	C1-O5-C5	5.99	120.30	112.19
3	D	4	MAN	O5-C5-C6	-5.42	98.70	107.20
3	D	3	BMA	C1-O5-C5	5.31	119.39	112.19
3	D	5	NAG	C1-O5-C5	5.15	119.17	112.19
3	D	3	BMA	C1-C2-C3	-5.00	103.52	109.67

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	NAG	C1-O5-C5	4.80	118.70	112.19
3	C	5	NAG	C1-O5-C5	4.70	118.56	112.19
3	D	6	MAN	O5-C5-C6	4.27	113.90	107.20
3	C	4	MAN	O5-C5-C6	-4.13	100.73	107.20
3	D	7	NAG	C1-O5-C5	4.05	117.67	112.19
3	C	6	MAN	O5-C5-C6	-3.81	101.23	107.20
3	C	5	NAG	C1-C2-N2	-3.73	104.12	110.49
3	D	6	MAN	O2-C2-C1	3.70	116.72	109.15
3	C	5	NAG	O5-C1-C2	3.49	116.81	111.29
3	D	2	NAG	O5-C1-C2	-3.31	106.06	111.29
3	C	4	MAN	C1-O5-C5	3.27	116.62	112.19
3	C	2	NAG	O5-C1-C2	-3.21	106.22	111.29
3	C	3	BMA	C1-O5-C5	3.16	116.48	112.19
3	C	7	NAG	O5-C1-C2	-3.12	106.36	111.29
3	C	1	NAG	C4-C3-C2	-3.07	106.52	111.02
3	C	1	NAG	O5-C1-C2	-3.07	106.45	111.29
3	D	4	MAN	O2-C2-C1	-2.91	103.19	109.15
3	C	7	NAG	O5-C5-C6	2.84	111.66	107.20
3	C	4	MAN	O2-C2-C3	-2.83	104.46	110.14
3	C	4	MAN	C3-C4-C5	-2.80	105.25	110.24
3	C	3	BMA	O5-C1-C2	-2.74	106.55	110.77
3	D	7	NAG	O5-C1-C2	-2.68	107.06	111.29
3	D	8	FUC	C1-C2-C3	-2.57	106.50	109.67
3	D	5	NAG	C3-C4-C5	-2.51	105.77	110.24
3	D	2	NAG	C2-N2-C7	-2.50	119.34	122.90
3	D	8	FUC	C6-C5-C4	-2.49	108.47	113.07
3	D	7	NAG	C2-N2-C7	-2.47	119.38	122.90
3	D	3	BMA	C6-C5-C4	-2.43	107.32	113.00
3	C	7	NAG	C2-N2-C7	-2.41	119.48	122.90
3	D	8	FUC	C3-C4-C5	-2.39	106.05	109.77
3	C	6	MAN	C6-C5-C4	-2.34	107.51	113.00
3	C	3	BMA	O2-C2-C1	-2.28	104.49	109.15
3	D	8	FUC	O5-C1-C2	-2.25	107.29	110.77
3	C	4	MAN	O5-C5-C4	-2.23	105.40	110.83
3	D	3	BMA	O4-C4-C3	-2.21	105.23	110.35
3	C	5	NAG	C4-C3-C2	-2.21	107.78	111.02
3	C	6	MAN	C1-C2-C3	-2.20	106.96	109.67
3	C	5	NAG	C3-C4-C5	2.20	114.16	110.24
3	D	6	MAN	C1-O5-C5	-2.19	109.22	112.19
3	D	4	MAN	C1-O5-C5	2.19	115.16	112.19
3	C	2	NAG	C1-O5-C5	2.15	115.11	112.19
3	D	6	MAN	O5-C1-C2	2.13	114.06	110.77

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	8	FUC	O4-C4-C5	-2.11	104.98	109.67
3	C	2	NAG	C6-C5-C4	-2.11	108.06	113.00
3	C	4	MAN	O4-C4-C3	-2.10	105.49	110.35
3	C	1	NAG	C3-C4-C5	-2.06	106.56	110.24
3	C	7	NAG	C1-C2-N2	-2.06	106.97	110.49
3	C	4	MAN	O5-C1-C2	-2.05	107.60	110.77
3	D	2	NAG	O5-C5-C6	-2.04	104.00	107.20
3	D	5	NAG	C2-N2-C7	2.03	125.79	122.90
3	D	5	NAG	O7-C7-N2	2.03	125.68	121.95

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1	NAG	C3-C2-N2-C7
3	D	2	NAG	O5-C5-C6-O6
3	C	6	MAN	O5-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
3	D	6	MAN	O5-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
3	C	6	MAN	C4-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
3	C	5	NAG	C8-C7-N2-C2
3	C	5	NAG	O7-C7-N2-C2
3	C	4	MAN	O5-C5-C6-O6
3	C	4	MAN	C4-C5-C6-O6
3	C	7	NAG	O5-C5-C6-O6
3	D	6	MAN	C4-C5-C6-O6
3	C	7	NAG	C4-C5-C6-O6
3	D	7	NAG	C1-C2-N2-C7
3	D	5	NAG	C4-C5-C6-O6
3	D	7	NAG	C4-C5-C6-O6
3	D	5	NAG	C1-C2-N2-C7
3	C	5	NAG	C4-C5-C6-O6
3	D	7	NAG	C3-C2-N2-C7

There are no ring outliers.

9 monomers are involved in 17 short contacts:

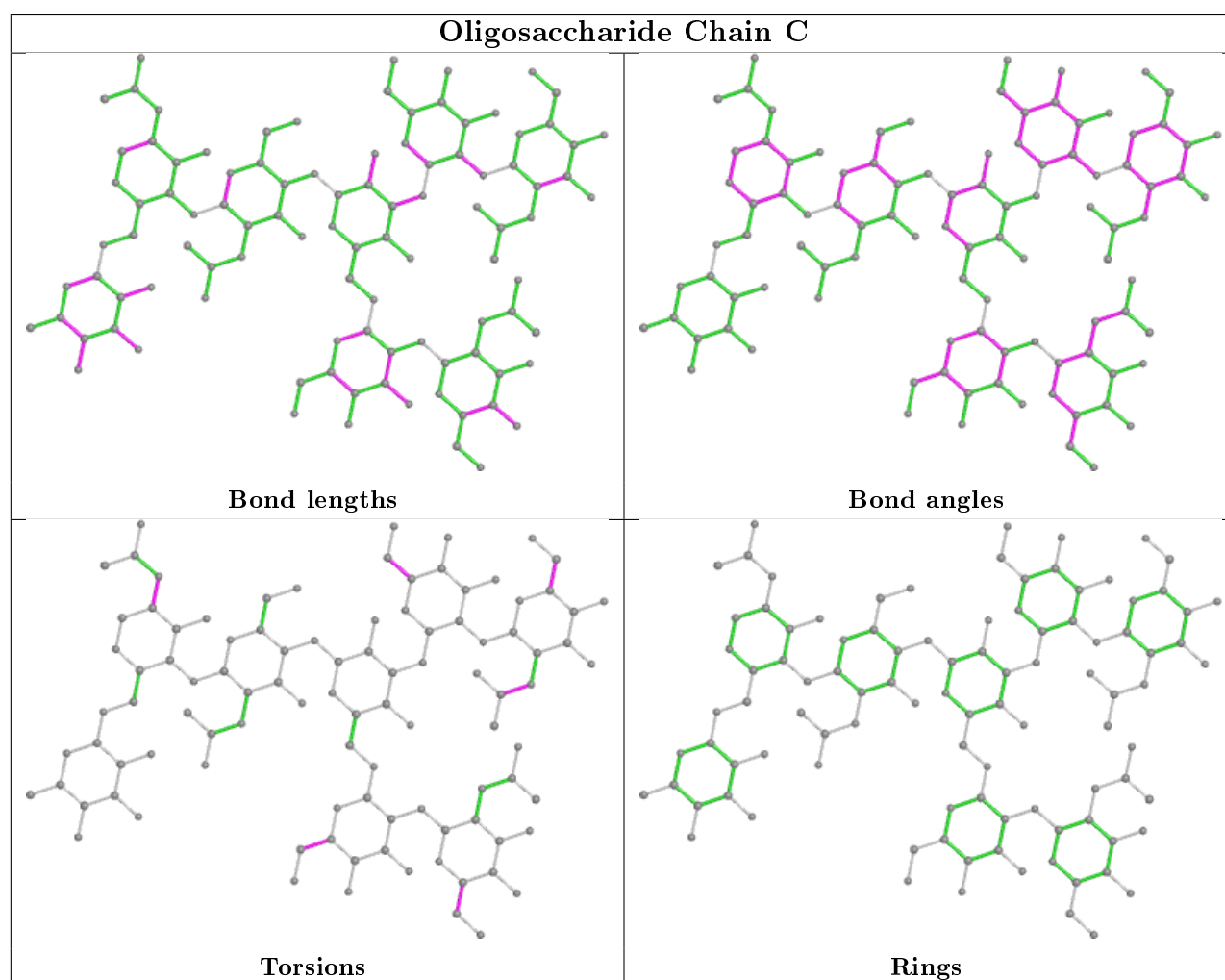
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	4	MAN	2	0

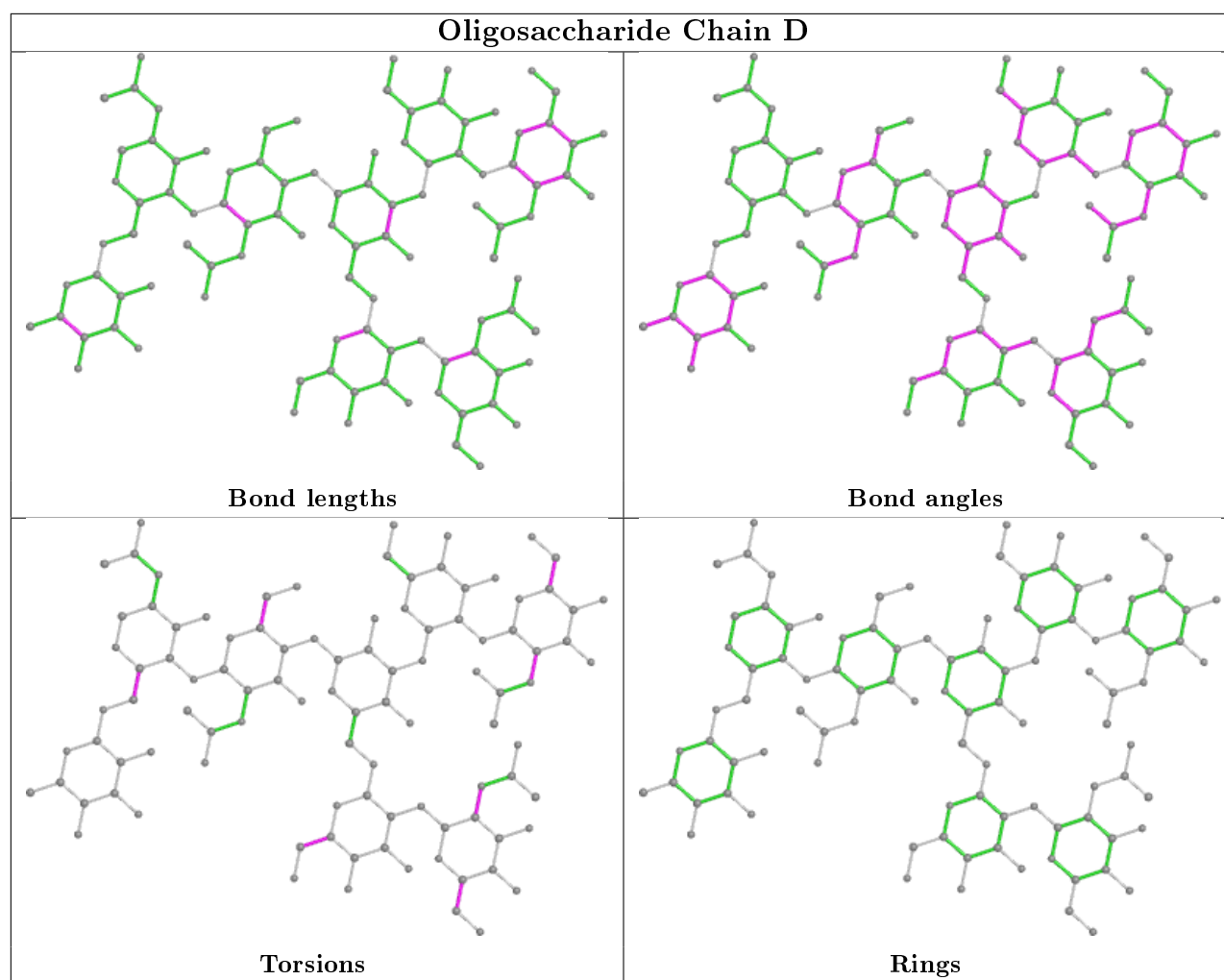
Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1	NAG	5	0
3	D	5	NAG	3	0
3	D	6	MAN	1	0
3	C	1	NAG	2	0
3	C	5	NAG	5	0
3	C	4	MAN	5	0
3	D	8	FUC	4	0
3	D	7	NAG	2	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	208/223 (93%)	0.43	1 (0%) 91 91	39, 52, 82, 95	0
2	B	205/223 (91%)	1.36	50 (24%) 0 0	37, 71, 109, 132	0
All	All	413/446 (92%)	0.89	51 (12%) 4 3	37, 59, 104, 132	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	302	VAL	21.4
2	B	270	ASP	10.1
2	B	241	PHE	7.4
2	B	263	VAL	6.9
2	B	266	VAL	5.5
2	B	296	TYR	5.4
2	B	242	LEU	5.1
2	B	268	HIS	5.0
2	B	240	VAL	4.8
2	B	264	VAL	4.8
2	B	301	ARG	4.6
2	B	291	PRO	4.6
2	B	297	ASN	4.4
2	B	326	LYS	4.4
2	B	331	PRO	4.3
2	B	303	VAL	4.3
2	B	332	ILE	4.3
2	B	281	GLY	4.1
2	B	330	ALA	4.0
2	B	267	SER	3.8
2	B	335	THR	3.8
2	B	336	ILE	3.7
2	B	319	TYR	3.6
2	B	324	SER	3.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	278	TYR	3.5
2	B	271	PRO	3.5
2	B	347	ARG	3.4
2	B	239	SER	3.3
2	B	304	SER	3.3
2	B	275	PHE	3.3
2	B	277	TRP	3.2
2	B	269	GLU	3.0
2	B	420	GLY	2.9
2	B	238	PRO	2.9
2	B	419	GLN	2.9
2	B	292	ARG	2.7
2	B	390	ASN	2.7
2	B	293	GLU	2.6
1	A	359	THR	2.5
2	B	243	PHE	2.5
2	B	329	PRO	2.5
2	B	327	ALA	2.4
2	B	322	LYS	2.3
2	B	295	GLN	2.2
2	B	361	ASN	2.2
2	B	279	VAL	2.1
2	B	320	LYS	2.1
2	B	300	TYR	2.1
2	B	288	LYS	2.1
2	B	337	SER	2.0
2	B	389	ASN	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
3	NAG	D	5	14/15	0.45	0.38	97,105,112,113	0

Continued on next page...

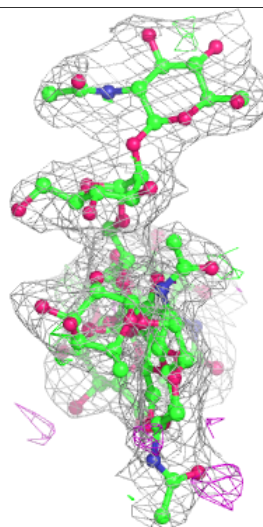
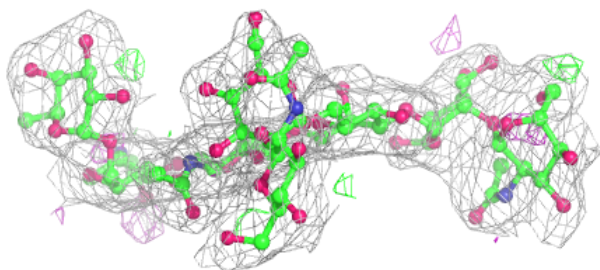
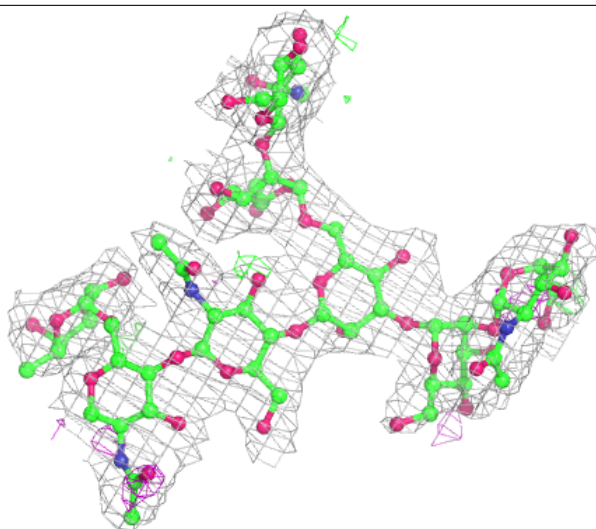
Continued from previous page...

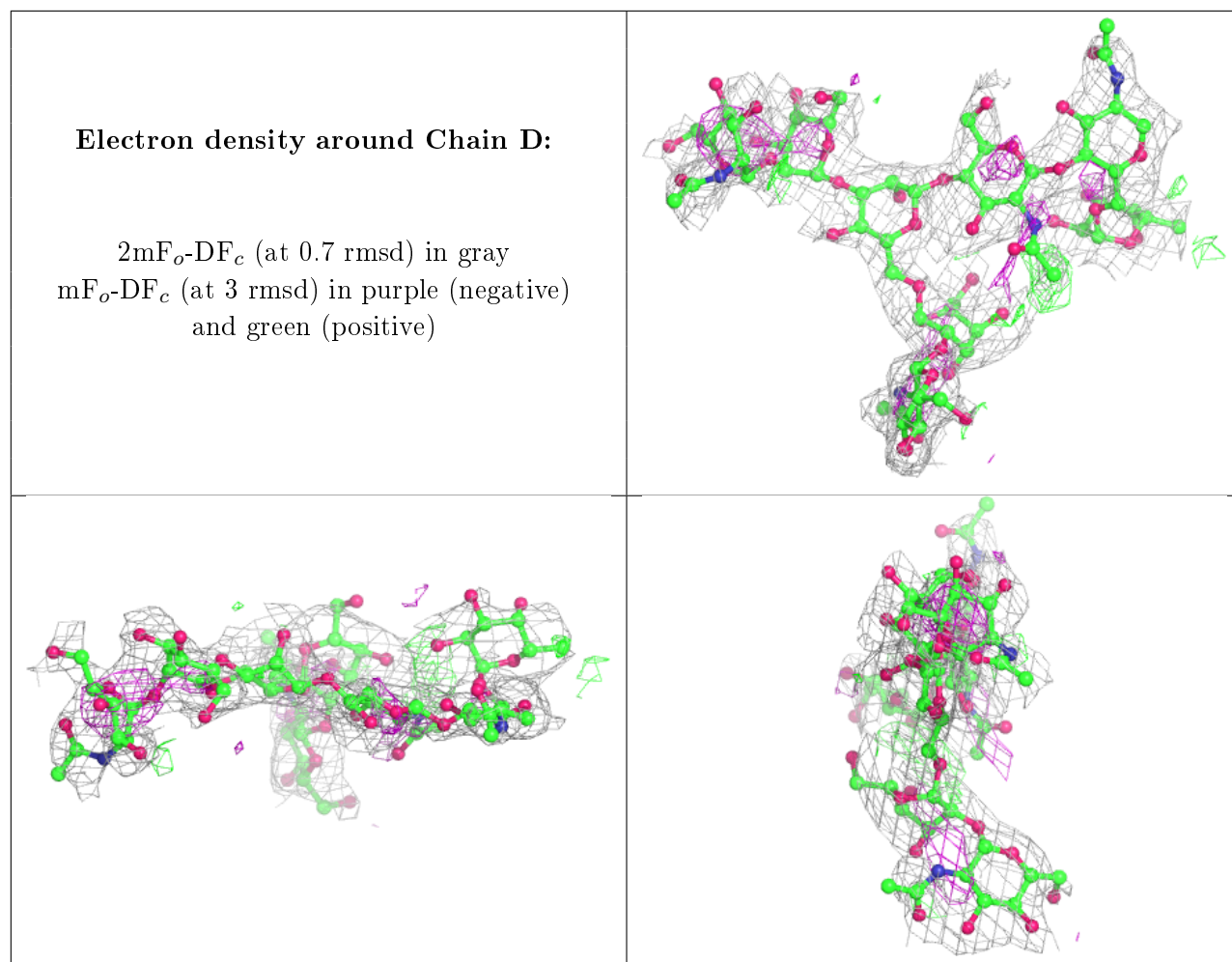
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	D	2	14/15	0.54	0.46	103,106,111,112	0
3	MAN	D	6	11/12	0.62	0.29	94,101,106,108	0
3	NAG	D	7	14/15	0.65	0.32	96,101,106,107	0
3	MAN	D	4	11/12	0.70	0.23	92,100,105,106	0
3	FUC	D	8	10/11	0.72	0.24	97,104,108,109	0
3	NAG	D	1	14/15	0.72	0.40	95,109,112,113	0
3	BMA	D	3	11/12	0.75	0.20	97,104,109,110	0
3	NAG	C	5	14/15	0.82	0.15	74,83,91,91	0
3	MAN	C	4	11/12	0.89	0.17	73,79,87,87	0
3	NAG	C	1	14/15	0.89	0.16	44,49,54,54	0
3	NAG	C	2	14/15	0.91	0.12	43,51,57,59	0
3	BMA	C	3	11/12	0.93	0.14	50,56,62,71	0
3	FUC	C	8	10/11	0.94	0.13	48,52,56,59	0
3	NAG	C	7	14/15	0.94	0.17	53,59,66,70	0
3	MAN	C	6	11/12	0.96	0.15	56,58,64,75	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.