



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 04:07 PM BST

PDB ID : 6D58
Title : Crystal structure of a Fc fragment of Human IgG3
Authors : Gohain, N.; Tolbert, W.D.; Pazgier, M.
Deposited on : 2018-04-19
Resolution : 2.39 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

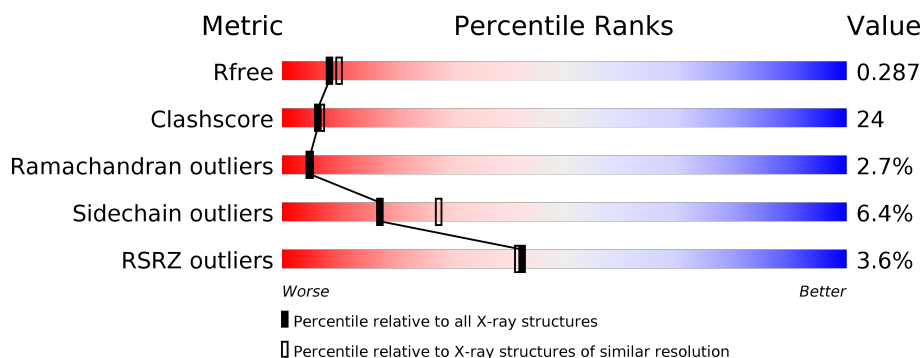
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	214	<div> <div>57%</div> <div>34%</div> <div>5%</div> </div>
1	B	214	<div> <div>7%</div> <div>52%</div> <div>38%</div> <div>6%</div> </div>
2	C	8	<div> <div>75%</div> <div>25%</div> </div>
2	D	8	<div> <div>13%</div> <div>25%</div> <div>63%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3538 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.

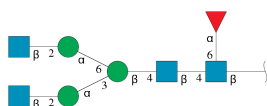
- Molecule 1 is a protein called Immunoglobulin heavy constant gamma 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1659	1056	278	317	8			
1	B	207	Total	C	N	O	S	0	0	0
			1659	1056	278	317	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	435	HIS	ARG	engineered mutation	UNP P01860
B	435	HIS	ARG	engineered mutation	UNP P01860

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	8	Total	C	N	O	0	0	0
			99	56	4	39			
2	D	8	Total	C	N	O	0	0	0
			99	56	4	39			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	12	Total	O	0	0
			12	12		

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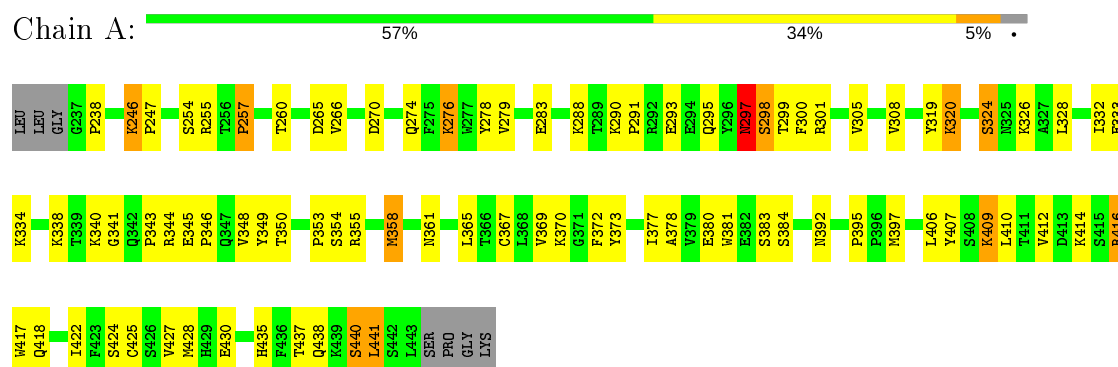
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	10	Total	O	0	0
			10	10		

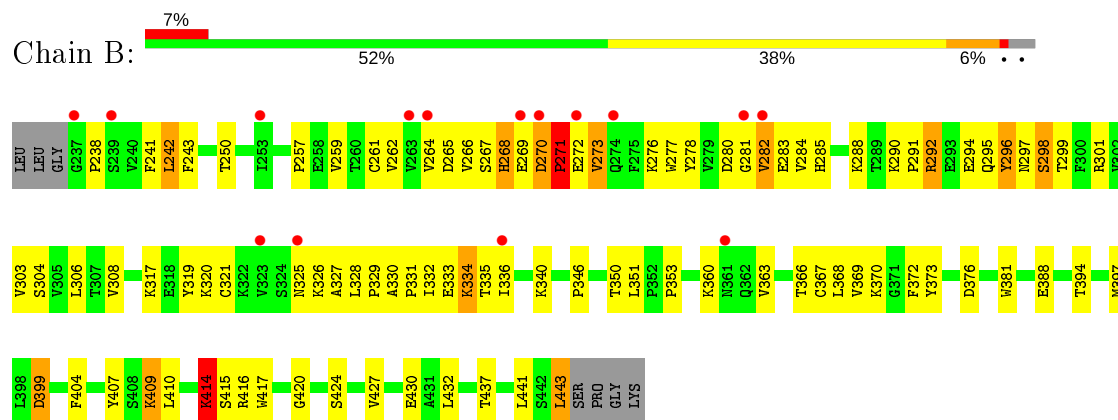
3 Residue-property plots

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.

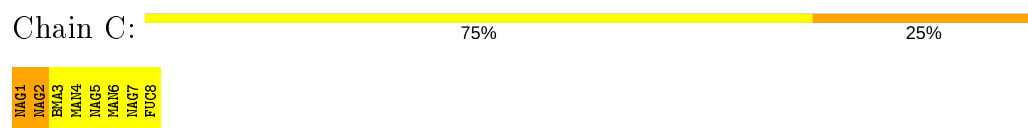
- Molecule 1: Immunoglobulin heavy constant gamma 3



- Molecule 1: Immunoglobulin heavy constant gamma 3



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



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

Chain D: 

HA61	HA62	HA63	HA64	HA65	HA66	HA67	FUC8
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.59 Å 80.68 Å 136.03 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.89 – 2.39 40.34 – 2.39	Depositor EDS
% Data completeness (in resolution range)	68.1 (35.89-2.39) 68.1 (40.34-2.39)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 2.39 Å)	Xtriage
Refinement program	PHENIX (1.13 _2998: ???)	Depositor
R, R_{free}	0.234 , 0.286 0.235 , 0.287	Depositor DCC
R_{free} test set	759 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	55.0	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 33.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3538	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.38% 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: 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	0.50	0/1705	0.67	2/2320 (0.1%)
1	B	0.59	2/1705 (0.1%)	0.76	3/2320 (0.1%)
All	All	0.55	2/3410 (0.1%)	0.72	5/4640 (0.1%)

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
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	271	PRO	N-CD	-5.55	1.40	1.47
1	B	280	ASP	CB-CG	5.04	1.62	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	242	LEU	CB-CG-CD2	-6.79	99.46	111.00
1	A	297	ASN	CB-CA-C	-6.09	98.22	110.40
1	B	280	ASP	CB-CG-OD2	-5.72	113.15	118.30
1	A	441	LEU	CA-CB-CG	5.29	127.48	115.30
1	B	280	ASP	CB-CG-OD1	5.17	122.95	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	297	ASN	Mainchain

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	1659	0	1623	77	0
1	B	1659	0	1623	96	0
2	C	99	0	85	4	0
2	D	99	0	85	10	0
3	A	12	0	0	1	0
3	B	10	0	0	0	0
All	All	3538	0	3416	168	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 24.

All (168) 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:290:LYS:HG3	1:B:291:PRO:CD	1.75	1.17
1:B:290:LYS:HG3	1:B:291:PRO:HD3	1.29	1.08
1:A:395:PRO:HD3	1:B:397:MET:HE1	1.45	0.98
1:A:290:LYS:HG2	1:A:291:PRO:HD2	1.46	0.96
1:B:333:GLU:O	1:B:334:LYS:HD2	1.72	0.90
1:A:381:TRP:CG	1:A:410:LEU:HD12	2.08	0.88
1:A:381:TRP:CD2	1:A:410:LEU:HD12	2.13	0.84
1:B:281:GLY:H	1:B:320:LYS:NZ	1.80	0.80
1:B:350:THR:HG21	1:B:441:LEU:HB2	1.64	0.79
1:B:296:TYR:HE1	2:D:8:FUC:H62	1.48	0.77
1:A:348:VAL:HG22	1:A:369:VAL:HG22	1.66	0.76
1:A:424:SER:HB3	1:A:438:GLN:HE21	1.48	0.76
1:B:328:LEU:HD23	1:B:329:PRO:HD2	1.70	0.74
1:A:297:ASN:OD1	1:A:299:THR:HG21	1.89	0.73
1:A:427:VAL:HG22	1:A:437:THR:HG22	1.70	0.73
1:B:278:TYR:HB2	1:B:320:LYS:HE2	1.71	0.72
1:B:296:TYR:CE1	2:D:8:FUC:H62	2.24	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:VAL:HG12	1:B:285:HIS:H	1.56	0.71
1:B:276:LYS:HG2	1:B:283:GLU:OE1	1.91	0.70
1:B:281:GLY:H	1:B:320:LYS:HZ1	1.40	0.70
1:B:351:LEU:HB2	1:B:366:THR:HB	1.74	0.69
1:A:320:LYS:HD3	1:A:333:GLU:OE2	1.92	0.69
1:A:290:LYS:HE3	1:A:293:GLU:OE2	1.93	0.69
1:B:268:HIS:HD2	1:B:269:GLU:N	1.91	0.68
1:B:268:HIS:CD2	1:B:269:GLU:HG3	2.27	0.68
1:B:290:LYS:HG3	1:B:291:PRO:HD2	1.71	0.67
1:B:388:GLU:HG2	1:B:410:LEU:HD11	1.75	0.67
1:B:282:VAL:HG12	1:B:283:GLU:H	1.60	0.67
1:A:367:CYS:HB2	1:A:381:TRP:CZ2	2.31	0.66
1:B:268:HIS:O	1:B:271:PRO:HG3	1.96	0.66
1:B:288:LYS:NZ	1:B:290:LYS:HD2	2.11	0.66
1:A:350:THR:HG21	1:A:441:LEU:HB2	1.78	0.66
1:A:383:SER:OG	1:A:416:ARG:NH2	2.30	0.64
1:A:265:ASP:HA	1:A:299:THR:OG1	1.96	0.64
1:A:395:PRO:HD3	1:B:397:MET:CE	2.23	0.63
1:B:320:LYS:HG3	1:B:335:THR:OG1	1.99	0.62
1:B:265:ASP:HA	1:B:299:THR:HG21	1.82	0.61
1:B:295:GLN:HE22	1:B:301:ARG:HG2	1.65	0.61
1:B:291:PRO:HA	1:B:292:ARG:HH11	1.65	0.61
1:A:407:TYR:OH	1:B:409:LYS:HG3	2.00	0.60
1:A:355:ARG:HA	1:A:358:MET:HE2	1.82	0.60
1:B:443:LEU:H	1:B:443:LEU:HD23	1.68	0.59
1:A:278:TYR:HA	1:A:283:GLU:HA	1.85	0.59
1:B:350:THR:HG22	1:B:441:LEU:HD22	1.86	0.58
1:A:320:LYS:HA	1:A:334:LYS:O	2.04	0.58
1:A:279:VAL:HG12	1:A:319:TYR:CD1	2.37	0.58
1:A:297:ASN:HB2	1:A:299:THR:HG22	1.85	0.57
1:A:290:LYS:HG2	1:A:291:PRO:CD	2.29	0.57
1:B:261:CYS:HB2	1:B:277:TRP:CH2	2.38	0.57
1:B:270:ASP:HB3	1:B:327:ALA:HB2	1.86	0.57
1:B:290:LYS:CG	1:B:291:PRO:HD3	2.19	0.57
1:B:268:HIS:CD2	1:B:269:GLU:N	2.72	0.57
1:A:343:PRO:HA	1:A:373:TYR:O	2.05	0.56
1:A:392:ASN:ND2	1:B:399:ASP:OD1	2.37	0.56
1:B:266:VAL:N	1:B:299:THR:OG1	2.38	0.56
1:B:414:LYS:HG3	1:B:415:SER:H	1.71	0.56
1:B:238:PRO:HG2	1:B:328:LEU:HG	1.88	0.55
1:A:397:MET:SD	1:B:394:THR:HA	2.46	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:TYR:OH	1:B:360:LYS:NZ	2.35	0.55
1:B:367:CYS:HB2	1:B:381:TRP:CZ2	2.42	0.55
1:A:383:SER:CB	1:A:416:ARG:NH2	2.71	0.54
1:A:384:SER:OG	1:A:384:SER:O	2.26	0.53
1:B:262:VAL:HG22	1:B:303:VAL:HG22	1.89	0.53
1:A:297:ASN:O	1:A:298:SER:HB2	2.09	0.53
1:B:272:GLU:HG2	1:B:273:VAL:N	2.24	0.53
1:A:380:GLU:OE1	1:A:380:GLU:N	2.42	0.52
1:B:242:LEU:CD2	1:B:336:ILE:HB	2.39	0.52
1:B:340:LYS:O	1:B:373:TYR:OH	2.17	0.52
1:B:328:LEU:HD13	1:B:332:ILE:HG13	1.91	0.52
1:B:350:THR:CG2	1:B:441:LEU:HB2	2.37	0.52
1:A:424:SER:HB3	1:A:438:GLN:NE2	2.21	0.51
1:A:414:LYS:HE2	1:A:418:GLN:NE2	2.25	0.51
1:B:241:PHE:CE2	2:D:2:NAG:H4	2.45	0.51
1:A:295:GLN:HG2	2:C:1:NAG:H62	1.91	0.51
1:A:266:VAL:HB	1:A:300:PHE:HB2	1.92	0.51
1:A:417:TRP:CZ3	1:A:441:LEU:HD13	2.46	0.51
1:B:257:PRO:HB2	1:B:308:VAL:HG22	1.92	0.50
1:A:238:PRO:HD2	1:A:328:LEU:HG	1.92	0.50
1:B:278:TYR:HB2	1:B:320:LYS:HB2	1.93	0.50
1:B:317:LYS:HD3	1:B:317:LYS:N	2.25	0.50
1:A:425:CYS:O	1:A:438:GLN:HG2	2.12	0.50
1:B:242:LEU:HD13	1:B:334:LYS:HB3	1.93	0.49
1:A:301:ARG:NH1	2:C:2:NAG:O7	2.44	0.49
1:B:432:LEU:HD22	1:B:437:THR:HG23	1.95	0.49
1:A:297:ASN:OD1	1:A:299:THR:CG2	2.58	0.49
1:A:424:SER:CB	1:A:438:GLN:NE2	2.75	0.49
1:A:346:PRO:HB3	1:A:372:PHE:HB3	1.94	0.49
1:B:281:GLY:H	1:B:320:LYS:HZ2	1.59	0.49
1:A:254:SER:OG	1:A:255:ARG:NH2	2.46	0.49
1:A:422:ILE:HG22	1:A:440:SER:OG	2.13	0.49
1:B:268:HIS:CD2	1:B:269:GLU:CG	2.96	0.49
1:B:259:VAL:HG23	1:B:308:VAL:HG11	1.95	0.48
1:B:290:LYS:CG	1:B:291:PRO:CD	2.69	0.48
1:B:295:GLN:C	1:B:297:ASN:H	2.16	0.48
1:B:243:PHE:HE1	2:D:3:BMA:H62	1.78	0.48
1:A:344:ARG:O	1:A:372:PHE:HA	2.14	0.48
1:A:355:ARG:HA	1:A:358:MET:CE	2.43	0.48
1:B:282:VAL:HG12	1:B:283:GLU:N	2.29	0.47
1:B:388:GLU:CD	1:B:416:ARG:HH22	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:TRP:CB	1:A:410:LEU:HD12	2.43	0.47
1:B:294:GLU:O	1:B:294:GLU:HG2	2.15	0.47
1:A:328:LEU:HD13	1:A:332:ILE:HG13	1.96	0.47
1:A:345:GLU:CD	1:A:345:GLU:H	2.18	0.47
1:B:306:LEU:HD23	1:B:308:VAL:HG12	1.96	0.47
1:B:288:LYS:HZ1	1:B:290:LYS:HD2	1.79	0.47
1:A:383:SER:CB	1:A:416:ARG:HH22	2.28	0.47
1:B:276:LYS:HE3	1:B:283:GLU:OE1	2.15	0.46
1:B:257:PRO:HB2	1:B:308:VAL:CG2	2.44	0.46
1:B:427:VAL:HG23	1:B:432:LEU:HD11	1.96	0.46
1:A:397:MET:HB3	1:A:397:MET:HE2	1.91	0.46
1:B:242:LEU:HD12	1:B:321:CYS:HB3	1.97	0.46
1:A:276:LYS:NZ	1:A:283:GLU:OE1	2.46	0.46
1:B:328:LEU:HD21	1:B:332:ILE:HD11	1.96	0.46
1:B:241:PHE:HZ	2:D:2:NAG:H61	1.81	0.46
2:D:4:MAN:H4	2:D:5:NAG:H83	1.98	0.46
1:A:341:GLY:HA3	1:A:373:TYR:CE2	2.52	0.45
1:A:260:THR:HG23	1:A:305:VAL:HG22	1.98	0.45
1:B:325:ASN:OD1	1:B:326:LYS:N	2.49	0.45
1:A:350:THR:HG22	1:A:441:LEU:HG	1.99	0.45
1:B:268:HIS:HD2	1:B:269:GLU:HG3	1.79	0.45
1:B:278:TYR:CD1	1:B:283:GLU:HG3	2.52	0.45
1:B:340:LYS:HA	1:B:340:LYS:HD2	1.72	0.45
1:A:428:MET:HA	1:A:435:HIS:O	2.17	0.44
1:B:250:THR:HG22	1:B:257:PRO:HB3	1.98	0.44
1:B:350:THR:HG22	1:B:441:LEU:HD13	1.99	0.44
1:A:246:LYS:NZ	2:C:7:NAG:O4	2.47	0.44
1:B:243:PHE:CE1	2:D:3:BMA:H62	2.53	0.44
1:A:350:THR:CG2	1:A:441:LEU:HB2	2.46	0.44
1:B:325:ASN:HB3	1:B:328:LEU:HB2	1.99	0.44
1:A:361:ASN:O	1:A:414:LYS:HB2	2.18	0.44
1:A:409:LYS:NZ	1:B:399:ASP:OD2	2.49	0.44
1:B:308:VAL:HG12	1:B:319:TYR:HE2	1.83	0.44
1:B:242:LEU:HD22	1:B:336:ILE:HB	1.99	0.44
1:B:278:TYR:CB	1:B:320:LYS:HE2	2.46	0.43
1:B:360:LYS:HB2	1:B:360:LYS:HE2	1.60	0.43
1:B:264:VAL:HG21	2:D:1:NAG:O4	2.18	0.43
2:D:2:NAG:H83	2:D:6:MAN:O4	2.18	0.43
1:B:368:LEU:HD13	1:B:407:TYR:CZ	2.54	0.43
1:B:284:VAL:HG12	1:B:285:HIS:N	2.30	0.43
1:A:354:SER:OG	1:B:351:LEU:HD23	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:LYS:HB3	1:A:276:LYS:NZ	2.34	0.42
1:A:424:SER:OG	1:A:438:GLN:NE2	2.52	0.42
1:B:346:PRO:HB3	1:B:372:PHE:HB3	2.00	0.42
1:A:370:LYS:HE3	1:A:370:LYS:HB3	1.80	0.42
1:A:257:PRO:HB2	1:A:308:VAL:HG22	2.00	0.42
1:B:296:TYR:CE1	2:D:8:FUC:C6	3.01	0.42
2:C:1:NAG:O5	2:C:8:FUC:H5	2.20	0.42
1:A:409:LYS:O	3:A:601:HOH:O	2.22	0.42
1:B:353:PRO:HB3	1:B:363:VAL:HG22	2.01	0.42
1:A:355:ARG:HG2	1:A:355:ARG:H	1.48	0.42
1:A:424:SER:CB	1:A:438:GLN:HE21	2.22	0.42
1:A:274:GLN:HB3	1:A:324:SER:HB2	2.02	0.41
1:A:340:LYS:O	1:A:373:TYR:OH	2.22	0.41
1:A:410:LEU:HD22	1:A:412:VAL:HG13	2.01	0.41
1:B:295:GLN:HE22	1:B:301:ARG:CG	2.31	0.41
1:B:330:ALA:HB1	1:B:331:PRO:HD2	2.02	0.41
1:A:290:LYS:CG	1:A:291:PRO:HD2	2.33	0.41
1:A:338:LYS:NZ	1:A:430:GLU:OE2	2.53	0.41
1:B:369:VAL:HG12	1:B:372:PHE:CD2	2.55	0.41
1:A:355:ARG:O	1:A:358:MET:HB2	2.20	0.41
1:A:353:PRO:HG3	1:A:365:LEU:HD12	2.03	0.41
1:B:238:PRO:HB2	1:B:328:LEU:HD11	2.02	0.41
1:A:378:ALA:O	1:A:427:VAL:HA	2.20	0.41
1:A:383:SER:HB3	1:A:416:ARG:NH2	2.36	0.41
1:B:295:GLN:HE22	1:B:301:ARG:CD	2.34	0.41
1:B:397:MET:O	1:B:404:PHE:HA	2.20	0.41
1:A:377:ILE:HG12	1:A:406:LEU:HD21	2.04	0.40
1:B:298:SER:O	1:B:298:SER:OG	2.39	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	205/214 (96%)	196 (96%)	8 (4%)	1 (0%)	29	41
1	B	205/214 (96%)	180 (88%)	15 (7%)	10 (5%)	2	1
All	All	410/428 (96%)	376 (92%)	23 (6%)	11 (3%)	5	5

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	298	SER
1	B	271	PRO
1	B	273	VAL
1	B	282	VAL
1	B	414	LYS
1	B	296	TYR
1	B	417	TRP
1	B	399	ASP
1	B	268	HIS
1	B	430	GLU
1	B	420	GLY

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	194/199 (98%)	181 (93%)	13 (7%)	16	26
1	B	194/199 (98%)	182 (94%)	12 (6%)	18	29
All	All	388/398 (98%)	363 (94%)	25 (6%)	17	28

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	246	LYS
1	A	247	PRO
1	A	257	PRO
1	A	270	ASP
1	A	276	LYS

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Mol	Chain	Res	Type
1	A	288	LYS
1	A	320	LYS
1	A	324	SER
1	A	326	LYS
1	A	358	MET
1	A	409	LYS
1	A	416	ARG
1	A	440	SER
1	B	267	SER
1	B	270	ASP
1	B	292	ARG
1	B	298	SER
1	B	304	SER
1	B	334	LYS
1	B	370	LYS
1	B	376	ASP
1	B	409	LYS
1	B	414	LYS
1	B	424	SER
1	B	443	LEU

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

Mol	Chain	Res	Type
1	A	438	GLN
1	B	268	HIS
1	B	315	ASN

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	$\# Z > 2$
2	NAG	C	1	1,2	14,14,15	1.58	3 (21%)	17,19,21	1.13	1 (5%)
2	NAG	C	2	2	14,14,15	0.77	1 (7%)	17,19,21	0.54	0
2	BMA	C	3	2	11,11,12	1.56	3 (27%)	15,15,17	1.24	3 (20%)
2	MAN	C	4	2	11,11,12	1.03	0	15,15,17	1.38	2 (13%)
2	NAG	C	5	2	14,14,15	0.69	1 (7%)	17,19,21	0.44	0
2	MAN	C	6	2	11,11,12	0.69	0	15,15,17	1.25	2 (13%)
2	NAG	C	7	2	14,14,15	0.21	0	17,19,21	0.41	0
2	FUC	C	8	2	10,10,11	0.73	0	14,14,16	1.03	0
2	NAG	D	1	1,2	14,14,15	0.48	0	17,19,21	1.17	1 (5%)
2	NAG	D	2	2	14,14,15	0.32	0	17,19,21	0.56	0
2	BMA	D	3	2	11,11,12	1.09	0	15,15,17	1.06	1 (6%)
2	MAN	D	4	2	11,11,12	1.10	1 (9%)	15,15,17	1.41	3 (20%)
2	NAG	D	5	2	14,14,15	0.45	0	17,19,21	0.47	0
2	MAN	D	6	2	11,11,12	0.95	0	15,15,17	1.17	2 (13%)
2	NAG	D	7	2	14,14,15	0.30	0	17,19,21	0.64	0
2	FUC	D	8	2	10,10,11	2.88	4 (40%)	14,14,16	2.10	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
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	-	1/2/19/22	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	MAN	C	6	2	-	2/2/19/22	0/1/1/1
2	NAG	C	7	2	-	0/6/23/26	0/1/1/1
2	FUC	C	8	2	-	-	0/1/1/1
2	NAG	D	1	1,2	-	2/6/23/26	0/1/1/1

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

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	8	FUC	O5-C1	-6.24	1.33	1.43
2	D	8	FUC	C2-C3	4.99	1.59	1.52
2	C	1	NAG	O5-C1	-4.15	1.37	1.43
2	D	8	FUC	C1-C2	3.27	1.59	1.52
2	D	4	MAN	O5-C1	-2.49	1.39	1.43
2	C	2	NAG	O5-C1	-2.49	1.39	1.43
2	D	8	FUC	C6-C5	-2.45	1.45	1.51
2	C	3	BMA	C4-C5	2.43	1.58	1.53
2	C	3	BMA	O5-C1	-2.41	1.39	1.43
2	C	1	NAG	C1-C2	2.34	1.55	1.52
2	C	5	NAG	O5-C1	-2.27	1.40	1.43
2	C	3	BMA	O3-C3	2.14	1.48	1.43
2	C	1	NAG	C3-C2	2.11	1.57	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	8	FUC	C1-C2-C3	4.67	115.41	109.67
2	D	1	NAG	C1-O5-C5	4.47	118.24	112.19
2	D	4	MAN	O2-C2-C3	-3.93	102.27	110.14
2	D	8	FUC	O5-C5-C6	-3.55	99.69	107.33
2	C	4	MAN	O2-C2-C3	-3.16	103.80	110.14
2	C	1	NAG	C3-C4-C5	-3.13	104.65	110.24
2	C	4	MAN	C1-O5-C5	3.10	116.39	112.19
2	C	6	MAN	O2-C2-C3	-3.07	103.98	110.14
2	D	8	FUC	C6-C5-C4	3.04	118.68	113.07
2	D	8	FUC	C3-C4-C5	-2.65	105.64	109.77
2	C	6	MAN	C1-O5-C5	2.65	115.78	112.19
2	C	3	BMA	C1-C2-C3	2.56	112.82	109.67
2	D	6	MAN	O2-C2-C1	2.56	114.39	109.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3	BMA	C2-C3-C4	2.35	114.96	110.89
2	C	3	BMA	C3-C4-C5	2.33	114.40	110.24
2	D	8	FUC	O5-C1-C2	-2.30	107.22	110.77
2	D	4	MAN	C1-O5-C5	2.30	115.30	112.19
2	C	3	BMA	O2-C2-C3	-2.19	105.75	110.14
2	D	4	MAN	C1-C2-C3	-2.15	107.03	109.67
2	D	6	MAN	O2-C2-C3	-2.14	105.84	110.14

There are no chirality outliers.

All (15) torsion outliers are listed below:

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

There are no ring outliers.

11 monomers are involved in 14 short contacts:

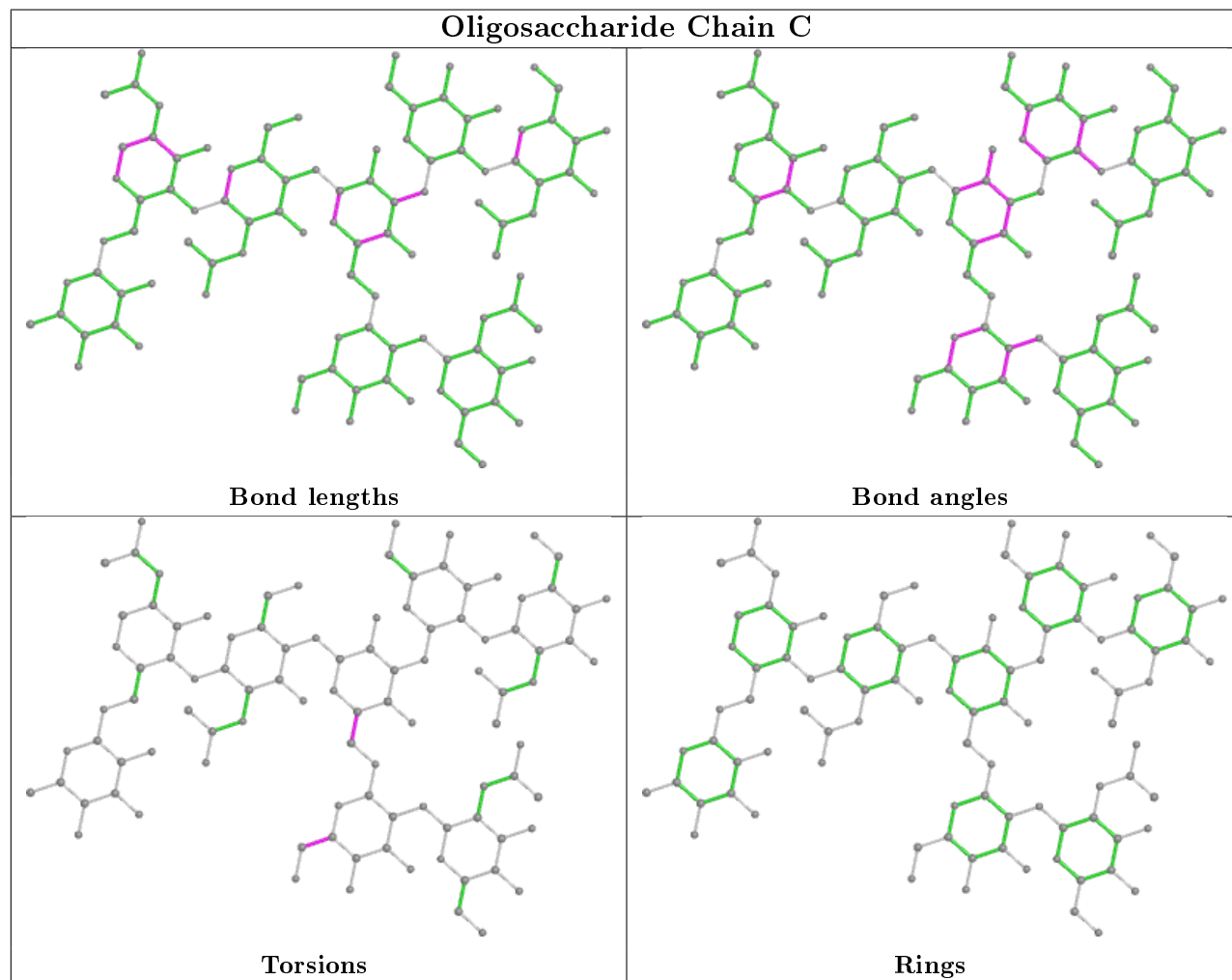
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	NAG	3	0
2	D	6	MAN	1	0
2	C	7	NAG	1	0
2	D	3	BMA	2	0
2	C	2	NAG	1	0
2	D	4	MAN	1	0
2	D	8	FUC	3	0
2	D	5	NAG	1	0
2	C	1	NAG	2	0

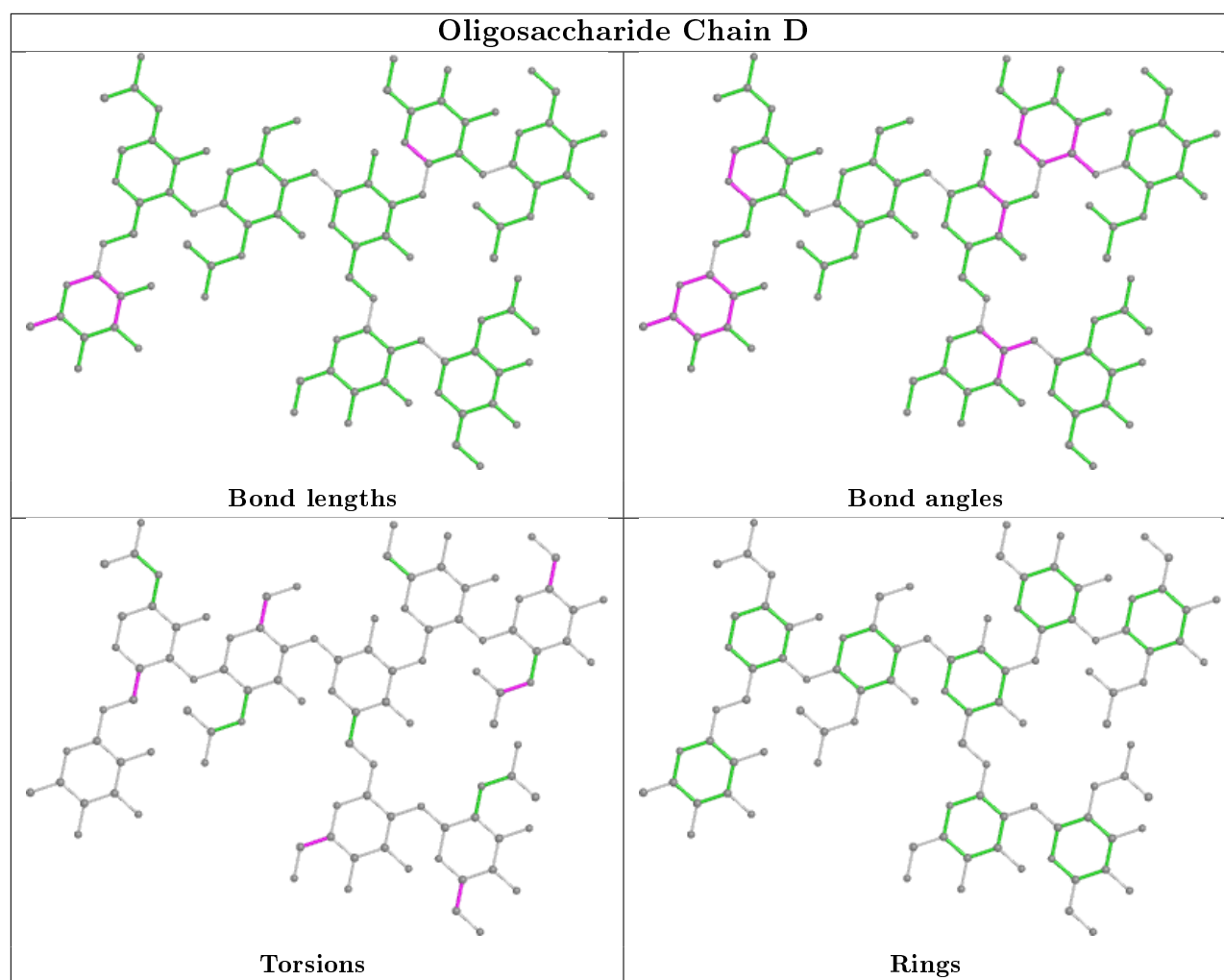
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	8	FUC	1	0
2	D	1	NAG	1	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	207/214 (96%)	0.04	0 100 100	32, 56, 84, 119	0
1	B	207/214 (96%)	0.34	15 (7%) 15 14	34, 70, 134, 171	0
All	All	414/428 (96%)	0.19	15 (3%) 42 42	32, 62, 119, 171	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	282	VAL	5.7
1	B	237	GLY	5.2
1	B	270	ASP	4.0
1	B	264	VAL	4.0
1	B	323	VAL	3.7
1	B	263	VAL	3.3
1	B	239	SER	3.1
1	B	281	GLY	3.1
1	B	253	ILE	3.0
1	B	325	ASN	2.7
1	B	361	ASN	2.5
1	B	272	GLU	2.3
1	B	269	GLU	2.3
1	B	274	GLN	2.1
1	B	336	ILE	2.1

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 ⓘ

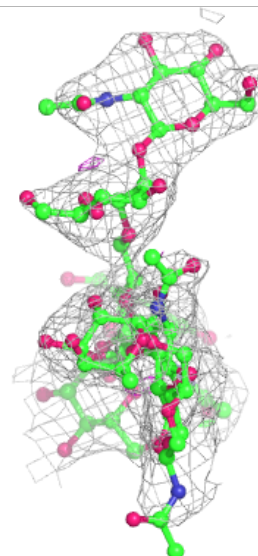
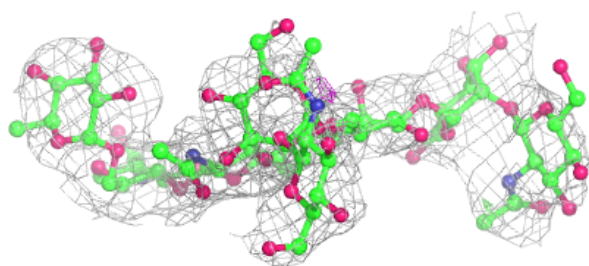
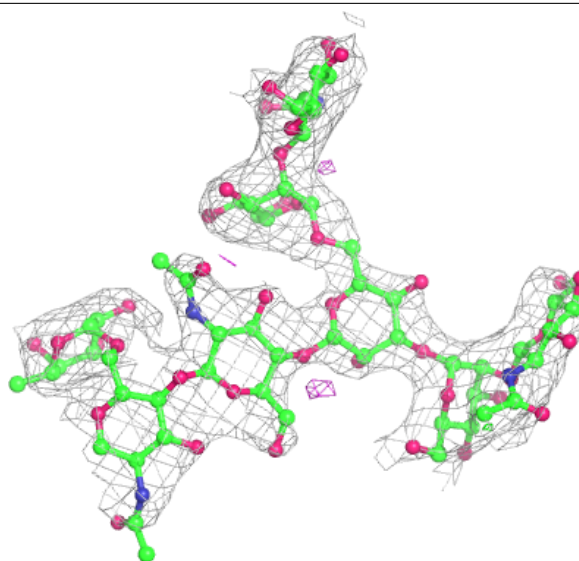
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

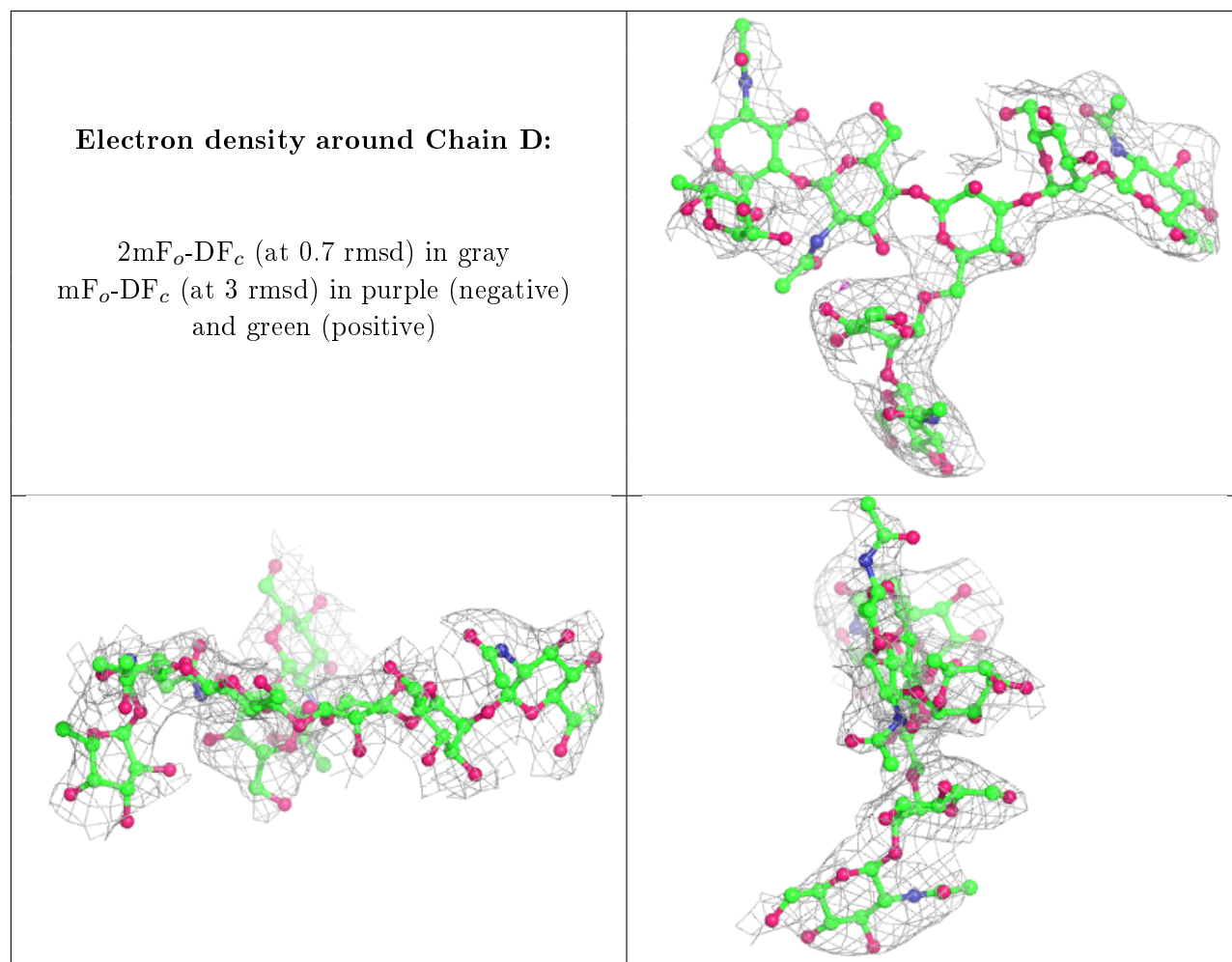
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	D	2	14/15	0.76	0.28	152,155,159,159	0
2	NAG	D	5	14/15	0.76	0.24	150,152,154,154	0
2	NAG	D	1	14/15	0.79	0.15	118,129,148,151	0
2	BMA	D	3	11/12	0.82	0.14	145,149,151,153	0
2	NAG	C	2	14/15	0.82	0.22	125,143,150,150	0
2	FUC	D	8	10/11	0.83	0.20	156,159,162,165	0
2	BMA	C	3	11/12	0.87	0.18	135,147,150,154	0
2	NAG	C	1	14/15	0.88	0.15	86,112,132,149	0
2	MAN	C	4	11/12	0.89	0.18	157,160,166,166	0
2	NAG	D	7	14/15	0.89	0.20	121,126,150,162	0
2	NAG	C	5	14/15	0.89	0.18	155,159,161,161	0
2	MAN	D	6	11/12	0.91	0.19	136,143,150,152	0
2	MAN	D	4	11/12	0.93	0.20	147,150,155,160	0
2	FUC	C	8	10/11	0.93	0.17	119,123,129,129	0
2	MAN	C	6	11/12	0.93	0.23	126,133,138,140	0
2	NAG	C	7	14/15	0.93	0.21	107,116,144,158	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.