



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:22 PM UTC

PDB ID : 6D58 / pdb_00006d58
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 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

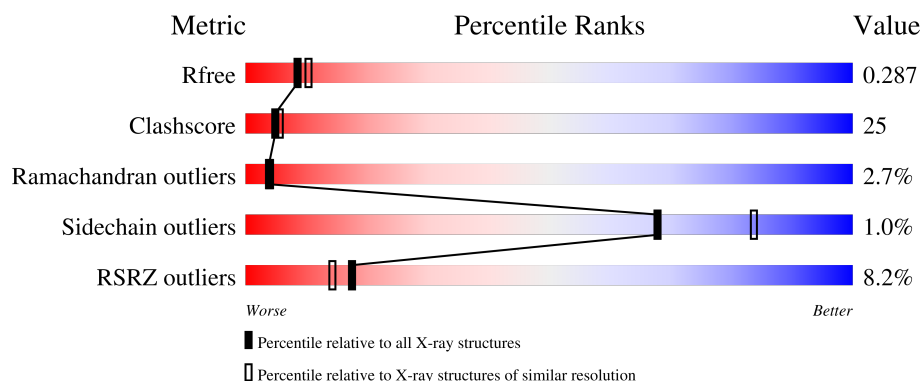
1 Overall quality at a glance

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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (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 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	214	7% 58% 37% ..
1	B	214	9% 54% 38% 5% .
2	C	8	75% 25%
2	D	8	12% 25% 62%

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.

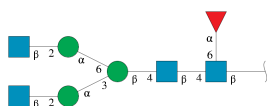
- Molecule 1 is a protein called Immunoglobulin heavy constant gamma 3.

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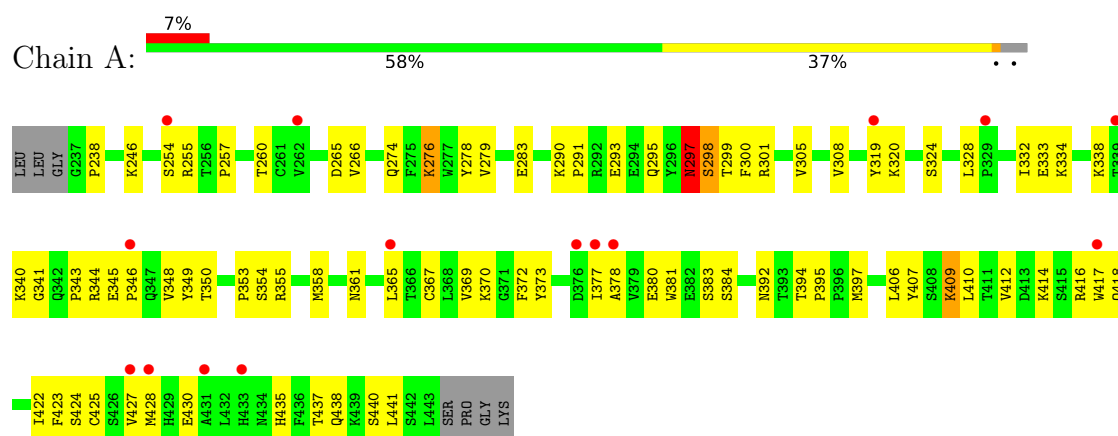
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	10	Total	O	0	0
			10	10		

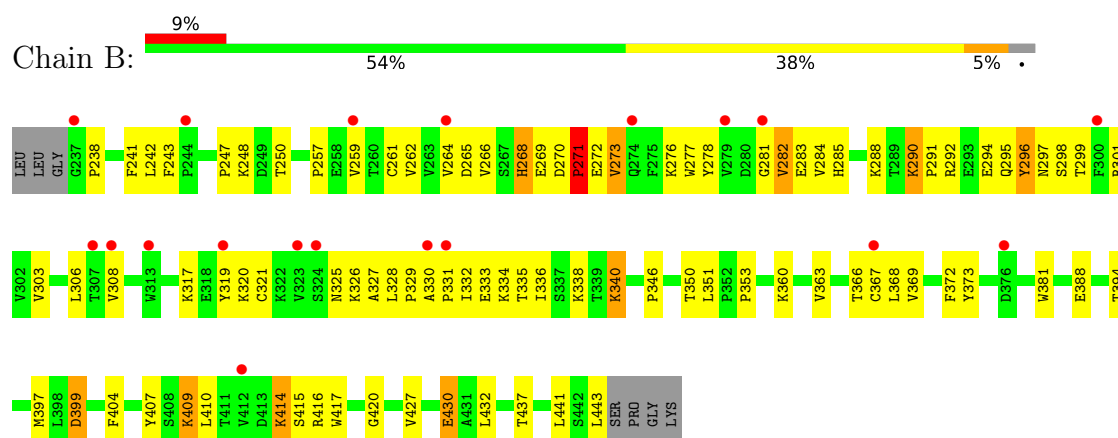
3 Residue-property plots [i](#)

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.

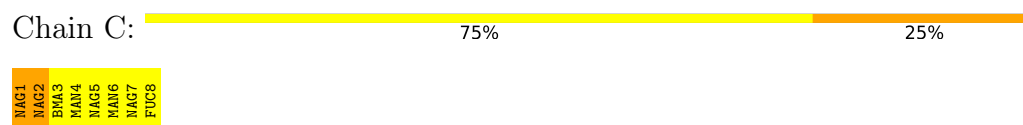
- Molecule 1: Immunoglobulin heavy constant gamma 3



- 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:  12% 25% 62%

NAG1	NAG2	BR43	MAN4	NA65	MAN6	NAG7	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	49.59Å 80.68Å 136.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.89 – 2.39 35.89 – 2.39	Depositor EDS
% Data completeness (in resolution range)	68.1 (35.89-2.39) 68.0 (35.89-2.39)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.234 , 0.286 0.236 , 0.287	Depositor DCC
R_{free} test set	759 reflections (3.40%)	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 , 32.7	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: NAG, MAN, FUC, 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	0.45	0/1705	0.72	2/2320 (0.1%)
1	B	0.53	1/1705 (0.1%)	0.80	2/2320 (0.1%)
All	All	0.49	1/3410 (0.0%)	0.76	4/4640 (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 (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	271	PRO	N-CD	-5.48	1.40	1.47

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	290	LYS	CA-C-N	-8.28	109.49	119.84
1	B	290	LYS	C-N-CA	-8.28	109.49	119.84
1	A	297	ASN	CB-CA-C	-7.33	98.22	110.68
1	A	297	ASN	CA-CB-CG	5.02	117.62	112.60

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	81	0
1	B	1659	0	1623	100	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	175	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 25.

All (175) 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: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:395:PRO:HD3	1:B:397:MET:HE1	1.53	0.89
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:A:355:ARG:HA	1:A:358:MET:HE2	1.65	0.78
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:296:TYR:CE1	2:D:8:FUC:H62	2.24	0.72
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: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:395:PRO:HD3	1:B:397:MET:CE	2.23	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:281:GLY:H	1:B:320:LYS:HZ2	1.40	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:B:340:LYS:O	1:B:373:TYR:OH	2.17	0.62
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:397:MET:SD	1:B:394:THR:HA	2.46	0.55
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: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:B:295:GLN:C	1:B:297:ASN:H	2.16	0.53
1:A:380:GLU:N	1:A:380:GLU:OE1	2.42	0.52
1:A:394:THR:HA	1:B:397:MET:HE3	1.91	0.52
1:B:242:LEU:CD2	1:B:336:ILE:HB	2.39	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:345:GLU:H	1:A:345:GLU:CD	2.18	0.52
1:A:424:SER:HB3	1:A:438:GLN:NE2	2.21	0.51
1:B:388:GLU:CD	1:B:416:ARG:HH22	2.18	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:288:LYS:HZ1	1:B:290:LYS:HD2	1.75	0.50
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:N	1:B:317:LYS:HD3	2.25	0.50
1:A:425:CYS:O	1:A:438:GLN:HG2	2.12	0.50
1:A:301:ARG:NH1	2:C:2:NAG:O7	2.44	0.49
1:B:242:LEU:HD13	1:B:334:LYS:HB3	1.93	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: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:243:PHE:HE1	2:D:3:BMA:H62	1.78	0.48
1:B:290:LYS:CG	1:B:291:PRO:CD	2.69	0.48
1:A:344:ARG:O	1:A:372:PHE:HA	2.14	0.48
1:B:282:VAL:HG12	1:B:283:GLU:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:LEU:HD13	1:A:332:ILE:HG13	1.96	0.47
1:A:381:TRP:CB	1:A:410:LEU:HD12	2.43	0.47
1:B:294:GLU:HG2	1:B:294:GLU:O	2.15	0.47
1:B:306:LEU:HD23	1:B:308:VAL:HG12	1.96	0.47
1:A:384:SER:O	1:A:384:SER:OG	2.26	0.47
1:B:281:GLY:N	1:B:320:LYS:HZ2	2.10	0.47
1:A:383:SER:CB	1:A:416:ARG:HH22	2.28	0.47
1:B:257:PRO:HB2	1:B:308:VAL:CG2	2.44	0.46
1:B:276:LYS:HE3	1:B:283:GLU:OE1	2.15	0.46
1:B:427:VAL:HG23	1:B:432:LEU:HD11	1.96	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:397:MET:HB3	1:A:397:MET:HE2	1.60	0.45
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:B:338:LYS:NZ	1:B:430:GLU:OE1	2.40	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:246:LYS:NZ	2:C:7:NAG:O4	2.47	0.44
1:A:428:MET:HA	1:A:435:HIS:O	2.17	0.44
1:B:243:PHE:CE1	2:D:3:BMA:H62	2.53	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: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:B:281:GLY:H	1:B:320:LYS:HZ1	1.59	0.44
1:A:409:LYS:NZ	1:B:399:ASP:OD2	2.49	0.44
1:A:423:PHE:O	1:A:440:SER:OG	2.26	0.44
1:B:242:LEU:HD22	1:B:336:ILE:HB	1.99	0.44
1:B:308:VAL:HG12	1:B:319:TYR:HE2	1.83	0.44
1:A:428:MET:HE2	1:A:428:MET:HB3	1.91	0.43
1:B:278:TYR:CB	1:B:320:LYS:HE2	2.46	0.43
1:B:264:VAL:HG21	2:D:1:NAG:O4	2.18	0.43
1:B:360:LYS:HB2	1:B:360:LYS:HE2	1.60	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:276:LYS:NZ	1:A:276:LYS:HB3	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:HB3	1:A:370:LYS:HE3	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:H	1:A:355:ARG:HG2	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:A:290:LYS:HE3	1:A:293:GLU:CD	2.45	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:353:PRO:HG3	1:A:365:LEU:HD12	2.03	0.41
1:A:355:ARG:O	1:A:358:MET:HB2	2.20	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:B:247:PRO:O	1:B:248:LYS:C	2.65	0.40
1:A:345:GLU:CD	1:A:345:GLU:N	2.79	0.40
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

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	205/214 (96%)	196 (96%)	8 (4%)	1 (0%)	24	37
1	B	205/214 (96%)	180 (88%)	15 (7%)	10 (5%)	1	1
All	All	410/428 (96%)	376 (92%)	23 (6%)	11 (3%)	4	4

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%)	192 (99%)	2 (1%)	68	84
1	B	194/199 (98%)	192 (99%)	2 (1%)	68	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	388/398 (98%)	384 (99%)	4 (1%)	68 84

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	276	LYS
1	A	409	LYS
1	B	340	LYS
1	B	409	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	362	GLN
1	A	434	ASN
1	A	438	GLN
1	B	268	HIS
1	B	295	GLN
1	B	342	GLN
1	B	433	HIS

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 ⓘ

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	1.57	3 (21%)	17,19,21	1.14	1 (5%)
2	NAG	C	2	2	14,14,15	0.75	1 (7%)	17,19,21	0.55	0
2	BMA	C	3	2	11,11,12	1.51	3 (27%)	15,15,17	1.16	3 (20%)
2	MAN	C	4	2	11,11,12	0.99	0	15,15,17	1.34	2 (13%)
2	NAG	C	5	2	14,14,15	0.67	1 (7%)	17,19,21	0.46	0
2	MAN	C	6	2	11,11,12	0.66	0	15,15,17	1.21	2 (13%)
2	NAG	C	7	2	14,14,15	0.21	0	17,19,21	0.43	0
2	FUC	C	8	2	10,10,11	0.72	0	14,14,16	0.96	0
2	NAG	D	1	1,2	14,14,15	0.47	0	17,19,21	1.19	1 (5%)
2	NAG	D	2	2	14,14,15	0.34	0	17,19,21	0.57	0
2	BMA	D	3	2	11,11,12	1.07	0	15,15,17	1.04	1 (6%)
2	MAN	D	4	2	11,11,12	1.06	1 (9%)	15,15,17	1.35	2 (13%)
2	NAG	D	5	2	14,14,15	0.43	0	17,19,21	0.48	0
2	MAN	D	6	2	11,11,12	0.91	0	15,15,17	1.06	2 (13%)
2	NAG	D	7	2	14,14,15	0.29	0	17,19,21	0.66	0
2	FUC	D	8	2	10,10,11	2.77	4 (40%)	14,14,16	1.93	4 (28%)

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	-	1/6/23/26	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
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	-	3/6/23/26	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	-5.92	1.33	1.43
2	D	8	FUC	C2-C3	4.81	1.59	1.52
2	C	1	NAG	O5-C1	-3.94	1.37	1.43
2	D	8	FUC	C1-C2	3.12	1.59	1.52
2	C	1	NAG	C1-C2	2.56	1.55	1.52
2	D	8	FUC	C6-C5	-2.44	1.45	1.51
2	C	3	BMA	C4-C5	2.40	1.58	1.53
2	D	4	MAN	O5-C1	-2.36	1.39	1.43
2	C	2	NAG	O5-C1	-2.35	1.39	1.43
2	C	3	BMA	O5-C1	-2.28	1.39	1.43
2	C	5	NAG	O5-C1	-2.14	1.40	1.43
2	C	1	NAG	C3-C2	2.13	1.57	1.52
2	C	3	BMA	O3-C3	2.04	1.48	1.43

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	NAG	C1-O5-C5	4.52	118.24	112.19
2	D	8	FUC	C1-C2-C3	3.96	115.41	109.64
2	D	4	MAN	O2-C2-C3	-3.81	102.27	110.15
2	D	8	FUC	O5-C5-C6	-3.56	99.69	107.40
2	C	4	MAN	C1-O5-C5	3.14	116.39	112.19
2	C	1	NAG	C3-C4-C5	-3.08	104.65	110.23
2	C	4	MAN	O2-C2-C3	-3.07	103.80	110.15
2	D	8	FUC	C6-C5-C4	3.06	118.68	113.08
2	C	6	MAN	O2-C2-C3	-2.98	103.98	110.15
2	D	8	FUC	C3-C4-C5	-2.74	105.64	109.81
2	C	6	MAN	C1-O5-C5	2.68	115.78	112.19
2	D	3	BMA	C2-C3-C4	2.33	114.96	110.86
2	D	4	MAN	C1-O5-C5	2.33	115.30	112.19
2	C	3	BMA	C3-C4-C5	2.30	114.40	110.23
2	D	6	MAN	O2-C2-C1	2.26	114.39	109.22
2	C	3	BMA	C1-C2-C3	2.18	112.82	109.64
2	C	3	BMA	O2-C2-C3	-2.13	105.75	110.15
2	D	6	MAN	O2-C2-C3	-2.08	105.84	110.15

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	6	MAN	O5-C5-C6-O6

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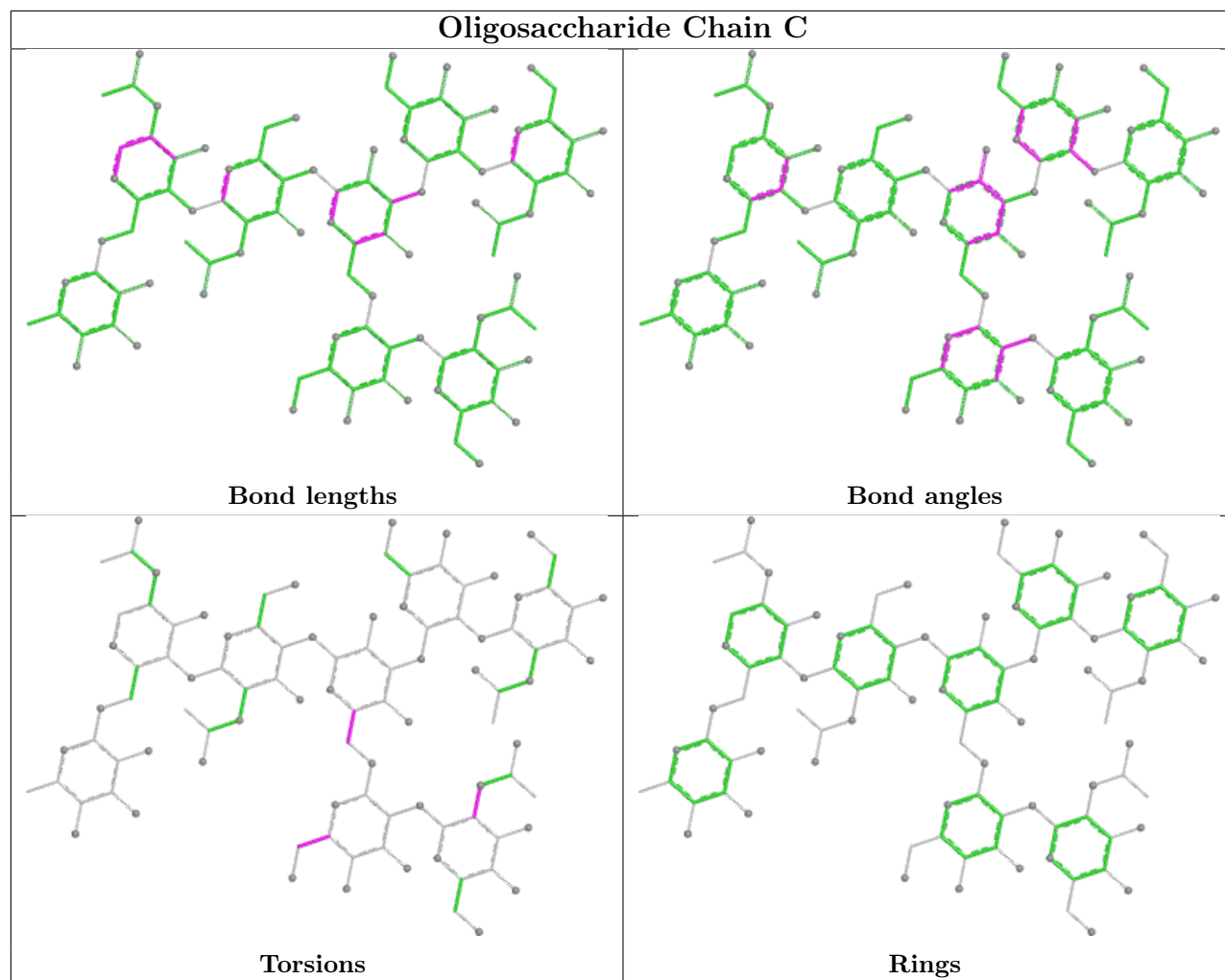
Mol	Chain	Res	Type	Atoms
2	D	5	NAG	O5-C5-C6-O6
2	C	6	MAN	C4-C5-C6-O6
2	D	5	NAG	C4-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	D	6	MAN	O5-C5-C6-O6
2	D	5	NAG	C8-C7-N2-C2
2	D	5	NAG	O7-C7-N2-C2
2	D	7	NAG	O5-C5-C6-O6
2	D	6	MAN	C4-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
2	D	7	NAG	C4-C5-C6-O6
2	C	7	NAG	C1-C2-N2-C7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

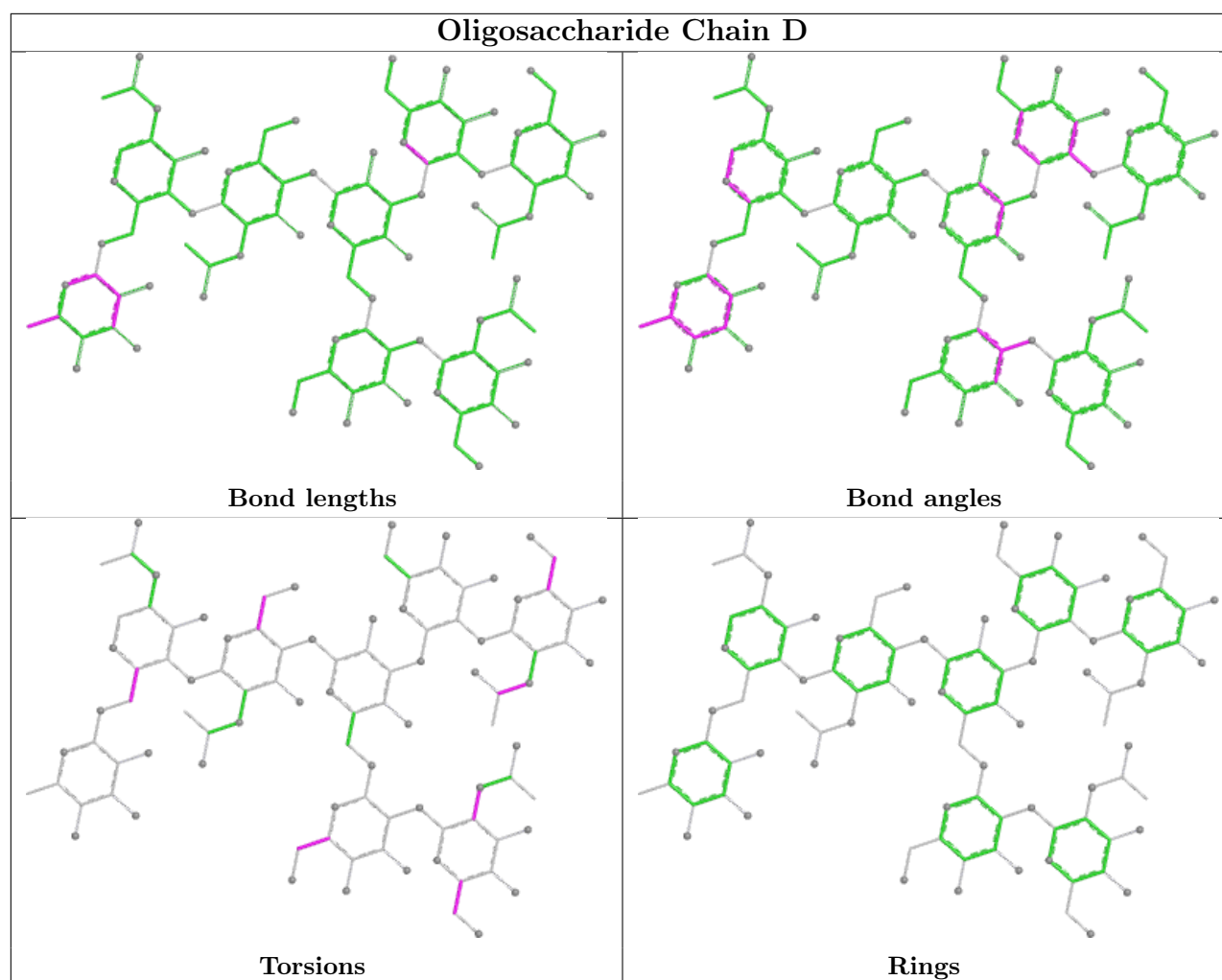
There are no ring outliers.

11 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	NAG	1	0
2	C	8	FUC	1	0
2	D	3	BMA	2	0
2	D	6	MAN	1	0
2	D	1	NAG	1	0
2	D	4	MAN	1	0
2	D	2	NAG	3	0
2	C	7	NAG	1	0
2	C	1	NAG	2	0
2	D	5	NAG	1	0
2	D	8	FUC	3	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	207/214 (96%)	0.50	15 (7%) 21 18	32, 56, 84, 119	0
1	B	207/214 (96%)	0.71	19 (9%) 14 12	34, 70, 134, 171	0
All	All	414/428 (96%)	0.61	34 (8%) 17 14	32, 62, 119, 171	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	323	VAL	5.7
1	B	319	TYR	4.2
1	A	377	ILE	3.7
1	B	324	SER	3.5
1	A	254	SER	3.4
1	B	259	VAL	3.4
1	B	237	GLY	3.1
1	B	376	ASP	2.9
1	B	307	THR	2.8
1	B	331	PRO	2.8
1	A	428	MET	2.8
1	B	279	VAL	2.8
1	B	412	VAL	2.7
1	B	313	TRP	2.6
1	B	264	VAL	2.6
1	A	417	TRP	2.5
1	B	308	VAL	2.4
1	A	431	ALA	2.4
1	A	339	THR	2.4
1	A	365	LEU	2.4
1	A	262	VAL	2.3
1	B	300	PHE	2.3
1	A	433	HIS	2.2
1	A	427	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	367	CYS	2.2
1	B	330	ALA	2.2
1	A	319	TYR	2.2
1	B	244	PRO	2.1
1	B	281	GLY	2.1
1	B	274	GLN	2.1
1	A	376	ASP	2.1
1	A	346	PRO	2.1
1	A	378	ALA	2.0
1	A	329	PRO	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.

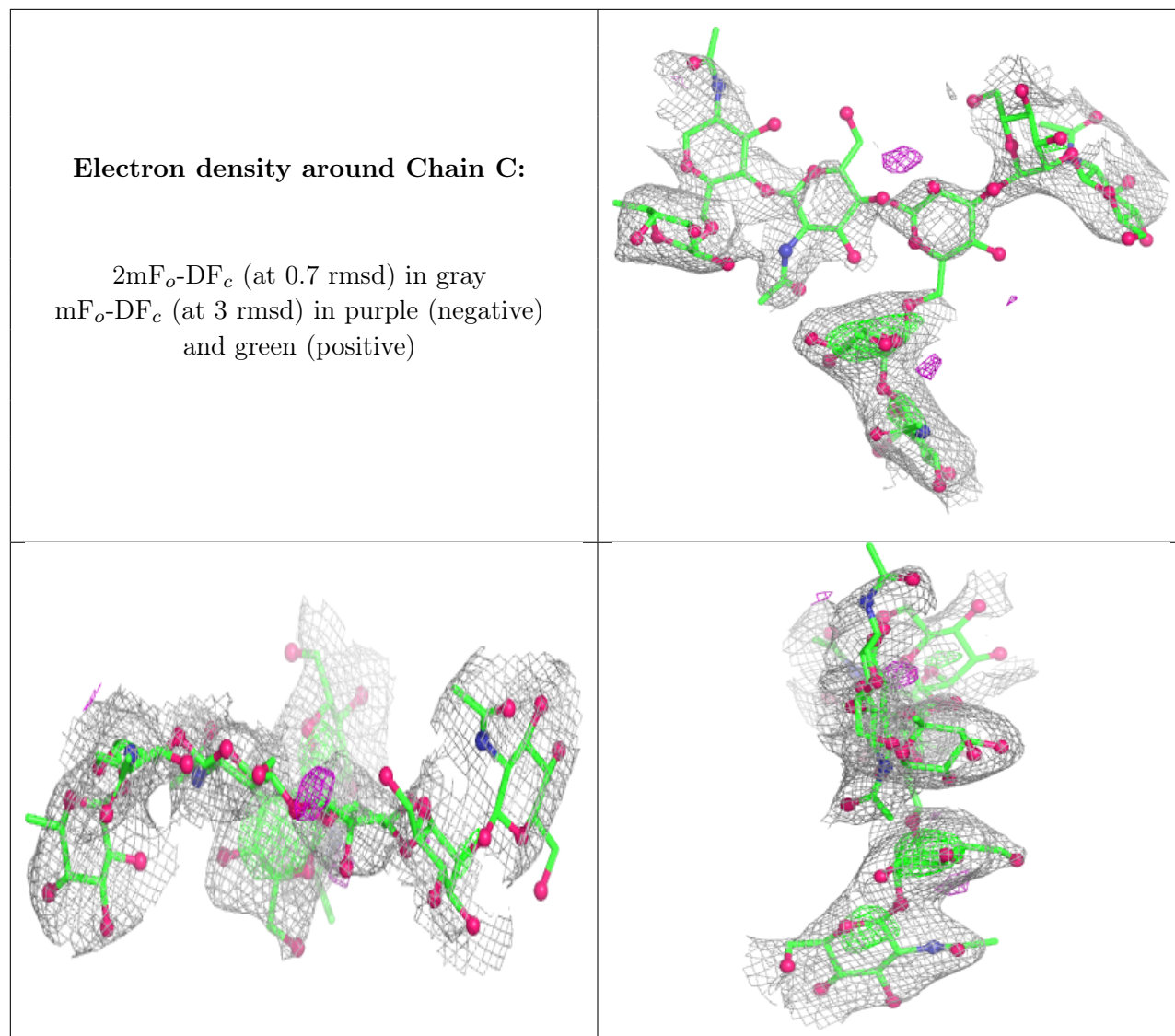
6.3 Carbohydrates [i](#)

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(Å ²)	Q<0.9
2	NAG	D	2	14/15	0.72	0.19	152,155,159,159	0
2	NAG	D	5	14/15	0.75	0.13	150,152,154,154	0
2	NAG	C	5	14/15	0.77	0.16	155,159,161,161	0
2	NAG	C	1	14/15	0.82	0.17	86,112,132,149	0
2	NAG	D	1	14/15	0.82	0.10	118,129,148,151	0
2	FUC	D	8	10/11	0.82	0.13	156,159,162,165	0
2	BMA	C	3	11/12	0.84	0.23	135,147,150,154	0
2	NAG	C	2	14/15	0.85	0.21	125,143,150,150	0
2	MAN	D	6	11/12	0.86	0.16	136,143,150,152	0
2	MAN	D	4	11/12	0.87	0.12	147,150,155,160	0
2	MAN	C	6	11/12	0.87	0.22	126,133,138,140	0
2	FUC	C	8	10/11	0.87	0.13	119,123,129,129	0
2	BMA	D	3	11/12	0.87	0.14	145,149,151,153	0
2	MAN	C	4	11/12	0.89	0.14	157,160,166,166	0
2	NAG	D	7	14/15	0.92	0.14	121,126,150,162	0
2	NAG	C	7	14/15	0.92	0.20	107,116,144,158	0

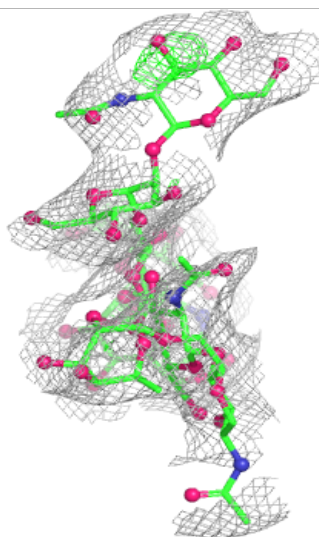
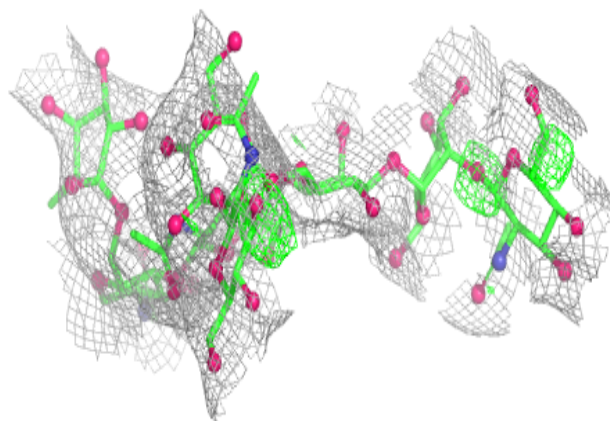
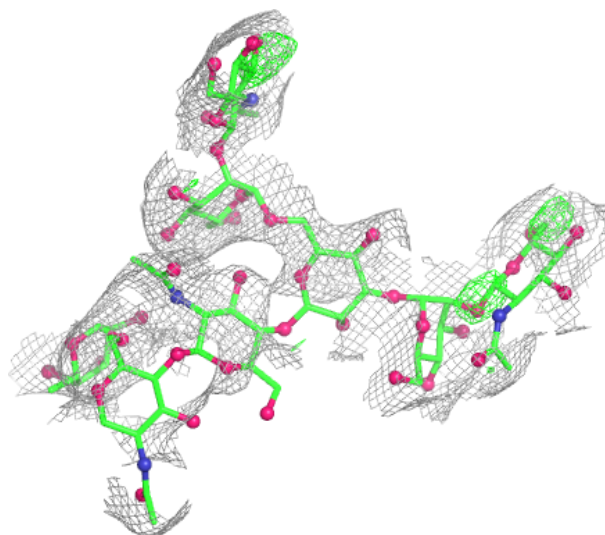
The following is a graphical depiction of the model fit to experimental electron density for oligosac-

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



Electron density around Chain D:

$2mF_o-DF_c$ (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.