



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 14, 2025 – 06:16 PM EDT

PDB ID : 3DO3 / pdb_00003do3
Title : Human IgG1 Fc fragment, 2.5 Angstrom structure
Authors : Braden, B.C.
Deposited on : 2008-07-03
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 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : **FAILED**
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.42

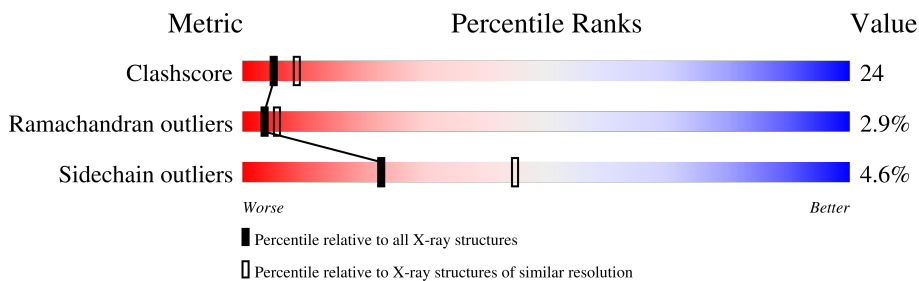
1 Overall quality at a glance

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))
Clashscore	180529	6282 (2.50-2.50)
Ramachandran outliers	177936	6191 (2.50-2.50)
Sidechain outliers	177891	6193 (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 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 $\leq 5\%$.

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	212	59% 37% ..
1	B	212	51% 42% 6% .
2	C	9	67% 33%
2	D	9	67% 22% 11%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3664 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.

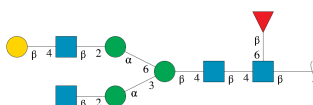
- Molecule 1 is a protein called Ig gamma-1 chain C region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	16	0	0
			1669	1061	281	320	7			
1	B	210	Total	C	N	O	S	32	0	0
			1676	1066	282	321	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	356	GLU	ASP	variant	UNP P01857
A	358	MET	LEU	variant	UNP P01857
B	356	GLU	ASP	variant	UNP P01857
B	358	MET	LEU	variant	UNP P01857

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



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

- Molecule 3 is water.

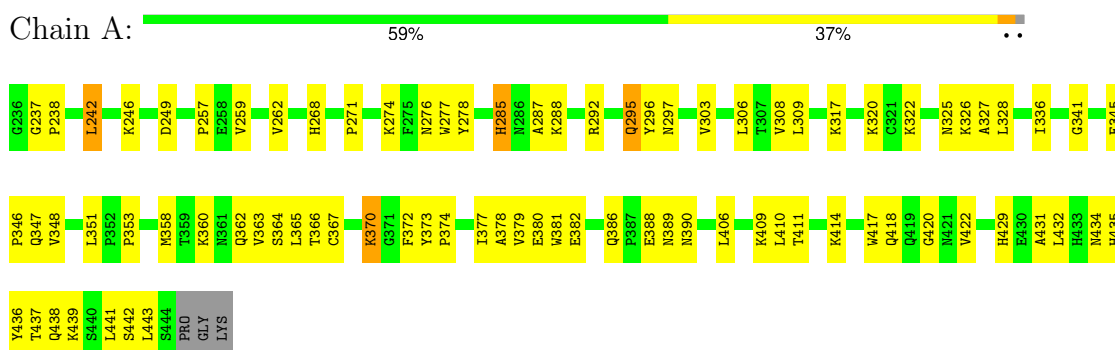
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total 54	O 54	0	0
3	B	45	Total 45	O 45	0	0

3 Residue-property plots

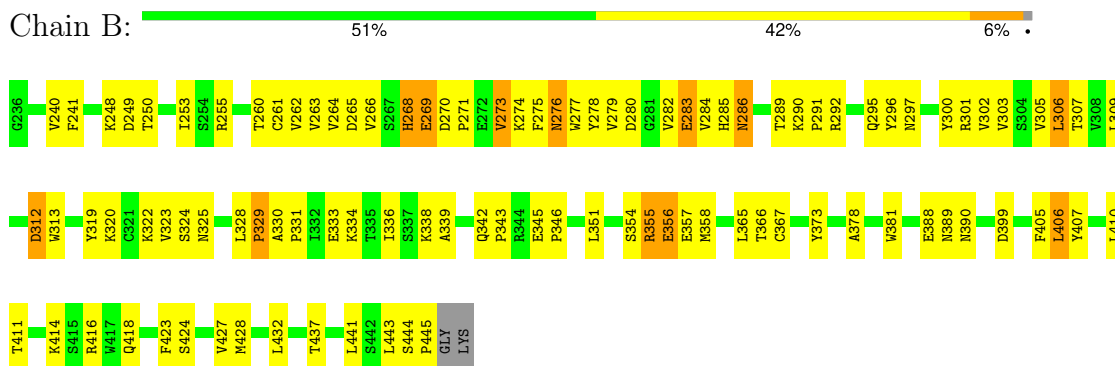
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS failed to run properly.

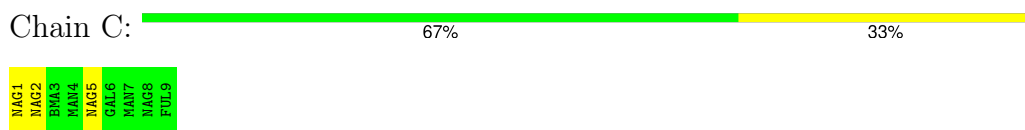
- Molecule 1: Ig gamma-1 chain C region



- Molecule 1: Ig gamma-1 chain C region



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



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

anose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[beta-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:



4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.21Å 80.22Å 138.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	17.50 – 2.50	Depositor
% Data completeness (in resolution range)	93.0 (17.50-2.50)	Depositor
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.49Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.280 , 0.330	Depositor
Wilson B-factor (Å ²)	55.0	Xtriage
Anisotropy	0.305	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3664	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.92% 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: BMA, GAL, MAN, FUL, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.41	0/1715	0.64	0/2335
1	B	0.40	0/1723	0.63	0/2347
All	All	0.41	0/3438	0.63	0/4682

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1669	0	1633	71	8
1	B	1676	0	1640	96	3
2	C	110	0	94	0	0
2	D	110	0	94	2	0
3	A	54	0	0	2	0
3	B	45	0	0	2	0
All	All	3664	0	3461	167	8

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 (167) 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:274:LYS:HE3	1:A:276:ASN:HD21	1.24	1.00
1:A:346:PRO:HB3	1:A:372:PHE:HB3	1.57	0.86
1:B:289:THR:HG22	1:B:290:LYS:H	1.40	0.84
1:A:242:LEU:HD13	1:A:336:ILE:HG13	1.60	0.83
1:A:274:LYS:HE3	1:A:276:ASN:ND2	1.95	0.81
1:A:238:PRO:HD2	1:A:328:LEU:HD13	1.64	0.79
1:B:290:LYS:HD3	1:B:303:VAL:HG23	1.68	0.76
1:B:249:ASP:OD1	1:B:255:ARG:HD3	1.87	0.75
1:A:438:GLN:O	1:A:439:LYS:HD2	1.87	0.73
1:B:289:THR:HG22	1:B:290:LYS:N	2.03	0.72
1:A:268:HIS:O	1:A:271:PRO:HD3	1.90	0.71
1:B:432:LEU:HD13	1:B:437:THR:HB	1.72	0.70
1:A:348:VAL:O	1:A:439:LYS:HG2	1.92	0.69
1:B:264:VAL:HG12	1:B:265:ASP:N	2.06	0.69
1:A:277:TRP:CE3	1:A:306:LEU:HD22	2.29	0.67
1:B:418:GLN:HA	1:B:443:LEU:HD22	1.76	0.67
1:A:295:GLN:O	1:A:297:ASN:N	2.26	0.67
1:B:279:VAL:O	1:B:280:ASP:HB2	1.93	0.67
1:A:363:VAL:HG12	1:A:364:SER:N	2.09	0.67
1:B:309:LEU:HB2	1:B:312:ASP:HB2	1.77	0.67
1:A:259:VAL:HG11	1:A:336:ILE:HD11	1.77	0.66
1:B:343:PRO:HA	1:B:373:TYR:O	1.97	0.65
1:A:360:LYS:NZ	1:A:360:LYS:HB2	2.12	0.65
1:A:379:VAL:O	1:A:380:GLU:HG3	1.97	0.65
1:A:276:ASN:HB2	1:A:322:LYS:HB3	1.79	0.65
1:A:420:GLY:HA2	1:A:443:LEU:HB3	1.78	0.64
1:B:414:LYS:HG2	1:B:418:GLN:HE21	1.65	0.62
1:A:422:VAL:HG22	1:A:442:SER:HB2	1.82	0.61
1:B:273:VAL:HG21	1:B:302:VAL:HG21	1.81	0.61
1:B:289:THR:CG2	1:B:290:LYS:H	2.13	0.61
1:B:390:ASN:ND2	1:B:411:THR:HB	2.16	0.61
1:B:264:VAL:CG1	1:B:265:ASP:N	2.64	0.60
1:B:262:VAL:HG13	1:B:303:VAL:HG12	1.84	0.59
1:B:285:HIS:O	1:B:286:ASN:HB3	2.02	0.59
1:A:432:LEU:O	1:A:435:HIS:N	2.34	0.59
1:B:444:SER:HB3	1:B:445:PRO:HD3	1.85	0.58
1:B:336:ILE:C	1:B:336:ILE:HD12	2.24	0.58
1:A:377:ILE:HG12	1:A:378:ALA:N	2.19	0.58
1:B:284:VAL:HG12	1:B:285:HIS:N	2.18	0.58
1:B:336:ILE:HD12	1:B:336:ILE:O	2.05	0.57
1:B:274:LYS:HB3	1:B:324:SER:HB2	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:417:TRP:O	1:A:443:LEU:HD22	2.05	0.56
1:B:406:LEU:HD12	1:B:406:LEU:C	2.25	0.56
1:A:257:PRO:HG2	1:A:308:VAL:O	2.06	0.56
1:B:330:ALA:HB1	1:B:331:PRO:CD	2.35	0.56
1:B:355:ARG:HA	1:B:358:MET:HG2	1.87	0.56
1:B:266:VAL:HG21	1:B:302:VAL:HG23	1.88	0.56
1:B:240:VAL:HG11	1:B:323:VAL:HG21	1.89	0.55
1:B:390:ASN:O	1:B:410:LEU:HD12	2.07	0.55
1:B:264:VAL:CG1	1:B:265:ASP:H	2.20	0.55
1:A:341:GLY:HA3	1:A:373:TYR:CE1	2.42	0.54
1:A:295:GLN:C	1:A:297:ASN:H	2.11	0.54
1:A:381:TRP:CE3	1:A:410:LEU:HD22	2.42	0.54
1:A:367:CYS:HB2	1:A:381:TRP:CZ2	2.42	0.54
1:B:268:HIS:O	1:B:271:PRO:HD3	2.07	0.54
1:A:358:MET:HE1	1:A:363:VAL:HG21	1.90	0.52
1:B:345:GLU:HG3	1:B:432:LEU:HD23	1.91	0.52
1:B:324:SER:O	1:B:325:ASN:HB2	2.08	0.52
1:B:250:THR:HG21	1:B:313:TRP:CD1	2.45	0.52
1:B:306:LEU:HD22	1:B:307:THR:N	2.25	0.52
1:B:406:LEU:HD12	1:B:406:LEU:O	2.10	0.52
1:B:289:THR:CG2	1:B:290:LYS:N	2.73	0.52
1:B:319:TYR:HB2	1:B:336:ILE:CD1	2.40	0.52
1:B:266:VAL:HB	1:B:300:TYR:HB2	1.92	0.51
1:A:363:VAL:HG12	1:A:364:SER:H	1.72	0.51
1:B:444:SER:HB3	1:B:445:PRO:CD	2.39	0.51
1:B:260:THR:HG23	1:B:305:VAL:HG12	1.93	0.51
1:B:248:LYS:HE3	1:B:255:ARG:CZ	2.41	0.51
1:B:414:LYS:HG2	1:B:418:GLN:NE2	2.25	0.51
1:B:274:LYS:HG2	1:B:275:PHE:N	2.26	0.50
1:A:358:MET:HA	1:A:358:MET:HE2	1.93	0.50
2:D:1:NAG:H4	2:D:9:FUL:H61	1.92	0.50
1:B:264:VAL:HG12	1:B:265:ASP:H	1.77	0.50
1:B:269:GLU:HA	1:B:269:GLU:OE1	2.12	0.50
1:B:278:TYR:HB2	1:B:320:LYS:HB3	1.94	0.50
1:B:325:ASN:HB3	1:B:328:LEU:HB3	1.93	0.50
1:A:325:ASN:OD1	1:A:327:ALA:HB3	2.13	0.49
1:B:273:VAL:CG2	1:B:302:VAL:HG21	2.43	0.49
1:A:363:VAL:CG1	1:A:364:SER:N	2.75	0.49
1:A:346:PRO:CB	1:A:372:PHE:HB3	2.38	0.49
1:B:313:TRP:HZ3	1:B:336:ILE:HD13	1.78	0.48
1:B:253:ILE:HG12	3:B:477:HOH:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:351:LEU:HB2	1:B:366:THR:HB	1.95	0.48
1:A:429:HIS:CD2	1:A:431:ALA:H	2.31	0.48
1:B:355:ARG:C	1:B:357:GLU:H	2.16	0.48
1:A:409:LYS:HG3	1:B:407:TYR:OH	2.14	0.48
1:B:330:ALA:HB1	1:B:331:PRO:HD2	1.96	0.48
1:A:238:PRO:HD2	1:A:328:LEU:CD1	2.41	0.48
1:A:382:GLU:HB2	1:A:386:GLN:O	2.14	0.47
1:B:284:VAL:HG12	1:B:285:HIS:H	1.77	0.47
1:A:262:VAL:HG13	1:A:303:VAL:HG22	1.96	0.47
1:A:420:GLY:H	1:A:443:LEU:HD23	1.77	0.47
1:A:436:TYR:CD1	1:A:437:THR:N	2.83	0.47
1:A:325:ASN:C	1:A:327:ALA:H	2.17	0.47
1:A:363:VAL:HG23	1:A:414:LYS:HB2	1.97	0.47
1:B:240:VAL:CG1	1:B:323:VAL:HG21	2.45	0.47
1:B:319:TYR:HB2	1:B:336:ILE:HD11	1.97	0.47
1:B:354:SER:HA	1:B:355:ARG:NH2	2.30	0.47
1:A:345:GLU:HG3	1:A:432:LEU:HD23	1.97	0.47
1:A:346:PRO:HG2	1:A:432:LEU:HD21	1.97	0.47
1:B:388:GLU:OE2	1:B:423:PHE:HE2	1.98	0.47
1:B:240:VAL:HG12	1:B:241:PHE:N	2.29	0.46
1:B:351:LEU:N	1:B:366:THR:O	2.43	0.46
1:A:365:LEU:CG	1:A:441:LEU:HD23	2.45	0.46
1:B:432:LEU:HD13	1:B:437:THR:CB	2.45	0.46
1:A:287:ALA:O	1:A:288:LYS:HG2	2.16	0.46
1:B:355:ARG:NE	1:B:355:ARG:H	2.13	0.46
1:B:261:CYS:HB2	1:B:277:TRP:CH2	2.50	0.46
1:B:390:ASN:HD22	1:B:411:THR:HB	1.78	0.46
1:A:259:VAL:CG1	1:A:336:ILE:HD11	2.43	0.46
1:B:358:MET:O	1:B:414:LYS:HD2	2.16	0.46
1:B:240:VAL:HG13	1:B:263:VAL:HG22	1.98	0.45
1:B:283:GLU:HG2	1:B:283:GLU:O	2.16	0.45
1:A:377:ILE:HG12	1:A:378:ALA:H	1.79	0.45
1:B:355:ARG:H	1:B:355:ARG:CZ	2.29	0.45
1:A:406:LEU:C	1:A:406:LEU:HD12	2.35	0.45
1:A:295:GLN:NE2	3:A:489:HOH:O	2.49	0.45
1:A:360:LYS:HB2	1:A:360:LYS:HZ2	1.80	0.45
1:B:240:VAL:HG12	1:B:241:PHE:H	1.82	0.45
1:B:282:VAL:HG12	1:B:282:VAL:O	2.15	0.44
1:A:347:GLN:HB2	1:A:370:LYS:HG2	1.99	0.44
1:B:333:GLU:C	1:B:334:LYS:HG2	2.36	0.44
1:B:365:LEU:HB3	1:B:441:LEU:HD23	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:333:GLU:O	1:B:334:LYS:HG2	2.18	0.44
1:B:406:LEU:C	1:B:406:LEU:CD1	2.85	0.44
1:B:378:ALA:HB3	1:B:428:MET:HB2	1.99	0.44
1:B:443:LEU:HG	1:B:445:PRO:HD2	1.98	0.44
1:A:358:MET:CE	1:A:363:VAL:HG21	2.48	0.44
1:A:278:TYR:CD1	1:A:278:TYR:N	2.85	0.43
1:A:351:LEU:HB2	1:A:366:THR:HB	1.99	0.43
1:A:353:PRO:CB	1:A:363:VAL:CG1	2.96	0.43
1:A:360:LYS:HG2	1:A:362:GLN:H	1.83	0.43
1:A:246:LYS:O	1:A:249:ASP:HB2	2.18	0.43
1:B:265:ASP:OD1	2:D:1:NAG:H3	2.18	0.43
1:B:306:LEU:O	1:B:306:LEU:HD13	2.17	0.43
1:B:427:VAL:HB	1:B:437:THR:HG22	1.99	0.43
1:A:358:MET:HE2	1:A:414:LYS:HD3	2.00	0.43
1:A:377:ILE:CG1	1:A:378:ALA:N	2.80	0.43
1:B:346:PRO:HG2	1:B:432:LEU:HD21	2.00	0.43
1:B:248:LYS:HG2	1:B:255:ARG:HD2	2.00	0.43
1:B:274:LYS:NZ	1:B:276:ASN:ND2	2.66	0.43
1:A:288:LYS:HG3	3:A:490:HOH:O	2.19	0.42
1:B:295:GLN:O	1:B:297:ASN:N	2.52	0.42
1:B:290:LYS:HB3	1:B:291:PRO:HD2	2.00	0.42
1:B:355:ARG:HG2	1:B:356:GLU:N	2.34	0.42
1:B:322:LYS:HE3	1:B:331:PRO:HB3	2.00	0.42
1:B:249:ASP:HA	1:B:255:ARG:HG3	2.01	0.42
1:B:365:LEU:HD13	1:B:441:LEU:HD23	2.00	0.42
1:A:336:ILE:HD12	1:A:336:ILE:O	2.20	0.42
1:B:381:TRP:HA	1:B:424:SER:O	2.20	0.42
1:A:363:VAL:CG1	1:A:364:SER:H	2.32	0.41
1:B:405:PHE:CD1	1:B:405:PHE:C	2.93	0.41
1:A:377:ILE:CG1	1:A:378:ALA:H	2.34	0.41
1:B:264:VAL:HG12	1:B:265:ASP:OD1	2.20	0.41
1:A:390:ASN:ND2	1:A:411:THR:HB	2.36	0.41
1:A:418:GLN:HA	1:A:443:LEU:HD22	2.02	0.41
1:B:365:LEU:HB3	1:B:441:LEU:CD2	2.51	0.41
1:B:367:CYS:HB2	1:B:381:TRP:CZ2	2.55	0.41
1:A:278:TYR:HB2	1:A:320:LYS:HB3	2.02	0.41
1:B:338:LYS:HG2	1:B:339:ALA:N	2.36	0.41
1:A:365:LEU:HD23	1:A:441:LEU:HD23	2.02	0.41
1:A:374:PRO:O	1:A:429:HIS:HE1	2.03	0.41
1:A:439:LYS:HG3	3:B:469:HOH:O	2.20	0.41
1:B:302:VAL:HG12	1:B:303:VAL:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:GLU:HA	1:A:388:GLU:HG2	2.03	0.40
1:A:365:LEU:HG	1:A:441:LEU:HD23	2.03	0.40
1:A:309:LEU:N	1:A:309:LEU:HD23	2.36	0.40

All (8) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:389:ASN:ND2	1:B:389:ASN:ND2[3_545]	0.91	1.29
1:A:268:HIS:CE1	1:A:285:HIS:CE1[4_455]	1.02	1.18
1:A:268:HIS:ND1	1:A:285:HIS:CE1[4_455]	1.12	1.08
1:A:268:HIS:ND1	1:A:285:HIS:ND1[4_455]	1.13	1.07
1:A:389:ASN:CG	1:B:389:ASN:ND2[3_545]	1.96	0.24
1:A:268:HIS:CE1	1:A:285:HIS:NE2[4_455]	2.07	0.13
1:A:268:HIS:CE1	1:A:285:HIS:ND1[4_455]	2.08	0.12
1:A:389:ASN:ND2	1:B:389:ASN:CG[3_545]	2.11	0.09

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	207/212 (98%)	191 (92%)	12 (6%)	4 (2%)	6	12
1	B	208/212 (98%)	174 (84%)	26 (12%)	8 (4%)	2	3
All	All	415/424 (98%)	365 (88%)	38 (9%)	12 (3%)	3	6

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	292	ARG
1	A	296	TYR
1	B	286	ASN

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Mol	Chain	Res	Type
1	B	292	ARG
1	B	329	PRO
1	A	326	LYS
1	B	296	TYR
1	B	269	GLU
1	B	283	GLU
1	B	312	ASP
1	B	356	GLU
1	A	237	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	194/196 (99%)	188 (97%)	6 (3%)	35	62
1	B	195/196 (100%)	183 (94%)	12 (6%)	15	31
All	All	389/392 (99%)	371 (95%)	18 (5%)	23	45

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	242	LEU
1	A	285	HIS
1	A	295	GLN
1	A	317	LYS
1	A	370	LYS
1	A	434	ASN
1	B	268	HIS
1	B	270	ASP
1	B	273	VAL
1	B	276	ASN
1	B	301	ARG
1	B	306	LEU
1	B	329	PRO
1	B	342	GLN

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Mol	Chain	Res	Type
1	B	355	ARG
1	B	399	ASP
1	B	406	LEU
1	B	416	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	276	ASN
1	A	315	ASN
1	A	347	GLN
1	A	362	GLN
1	A	421	ASN
1	A	429	HIS
1	A	438	GLN
1	B	276	ASN
1	B	347	GLN
1	B	361	ASN
1	B	390	ASN
1	B	418	GLN
1	B	421	ASN
1	B	433	HIS
1	B	438	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

18 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	1	2,1	14,14,15	0.62	0	17,19,21	0.85	1 (5%)
2	NAG	C	2	2	14,14,15	0.48	0	17,19,21	0.77	1 (5%)
2	BMA	C	3	2	11,11,12	0.63	0	15,15,17	0.46	0
2	MAN	C	4	2	11,11,12	0.59	0	15,15,17	0.59	0
2	NAG	C	5	2	14,14,15	0.52	0	17,19,21	0.78	1 (5%)
2	GAL	C	6	2	11,11,12	0.62	0	15,15,17	0.28	0
2	MAN	C	7	2	11,11,12	0.48	0	15,15,17	0.50	0
2	NAG	C	8	2	14,14,15	0.66	0	17,19,21	0.76	0
2	FUL	C	9	2	10,10,11	0.49	0	14,14,16	0.51	0
2	NAG	D	1	2,1	14,14,15	0.71	0	17,19,21	0.72	1 (5%)
2	NAG	D	2	2	14,14,15	0.61	0	17,19,21	0.70	0
2	BMA	D	3	2	11,11,12	0.54	0	15,15,17	0.98	1 (6%)
2	MAN	D	4	2	11,11,12	0.53	0	15,15,17	0.64	0
2	NAG	D	5	2	14,14,15	0.54	0	17,19,21	0.88	0
2	GAL	D	6	2	11,11,12	0.51	0	15,15,17	0.55	0
2	MAN	D	7	2	11,11,12	0.44	0	15,15,17	0.63	0
2	NAG	D	8	2	14,14,15	0.63	0	17,19,21	0.84	0
2	FUL	D	9	2	10,10,11	0.63	0	14,14,16	0.66	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
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	-	2/6/23/26	0/1/1/1
2	GAL	C	6	2	-	2/2/19/22	0/1/1/1
2	MAN	C	7	2	-	2/2/19/22	0/1/1/1
2	NAG	C	8	2	-	2/6/23/26	0/1/1/1
2	FUL	C	9	2	-	-	0/1/1/1
2	NAG	D	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	BMA	D	3	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAN	D	4	2	-	0/2/19/22	0/1/1/1
2	NAG	D	5	2	-	1/6/23/26	0/1/1/1
2	GAL	D	6	2	-	2/2/19/22	0/1/1/1
2	MAN	D	7	2	-	0/2/19/22	0/1/1/1
2	NAG	D	8	2	-	2/6/23/26	0/1/1/1
2	FUL	D	9	2	-	-	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3	BMA	C1-C2-C3	2.75	113.64	109.64
2	C	1	NAG	C2-N2-C7	-2.67	119.32	122.90
2	C	5	NAG	C2-N2-C7	-2.45	119.62	122.90
2	D	1	NAG	C2-N2-C7	-2.29	119.83	122.90
2	C	2	NAG	C2-N2-C7	-2.03	120.18	122.90

There are no chirality outliers.

All (16) torsion outliers are listed below:

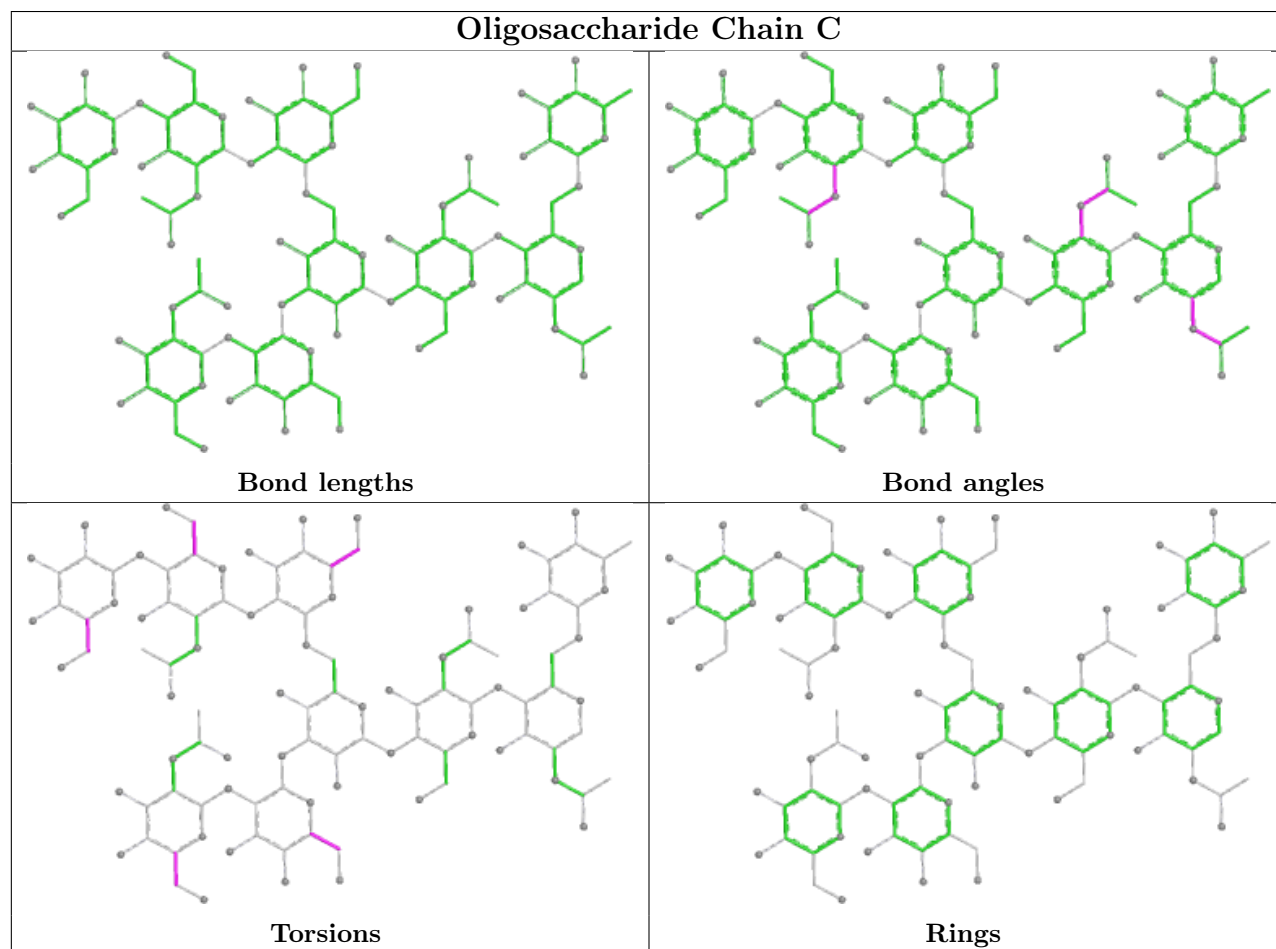
Mol	Chain	Res	Type	Atoms
2	C	6	GAL	O5-C5-C6-O6
2	C	6	GAL	C4-C5-C6-O6
2	D	6	GAL	O5-C5-C6-O6
2	D	8	NAG	O5-C5-C6-O6
2	C	5	NAG	O5-C5-C6-O6
2	C	5	NAG	C4-C5-C6-O6
2	C	7	MAN	C4-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
2	C	7	MAN	O5-C5-C6-O6
2	C	8	NAG	C4-C5-C6-O6
2	D	1	NAG	O5-C5-C6-O6
2	D	6	GAL	C4-C5-C6-O6
2	D	8	NAG	C4-C5-C6-O6
2	C	8	NAG	O5-C5-C6-O6
2	D	5	NAG	C1-C2-N2-C7
2	C	4	MAN	C4-C5-C6-O6

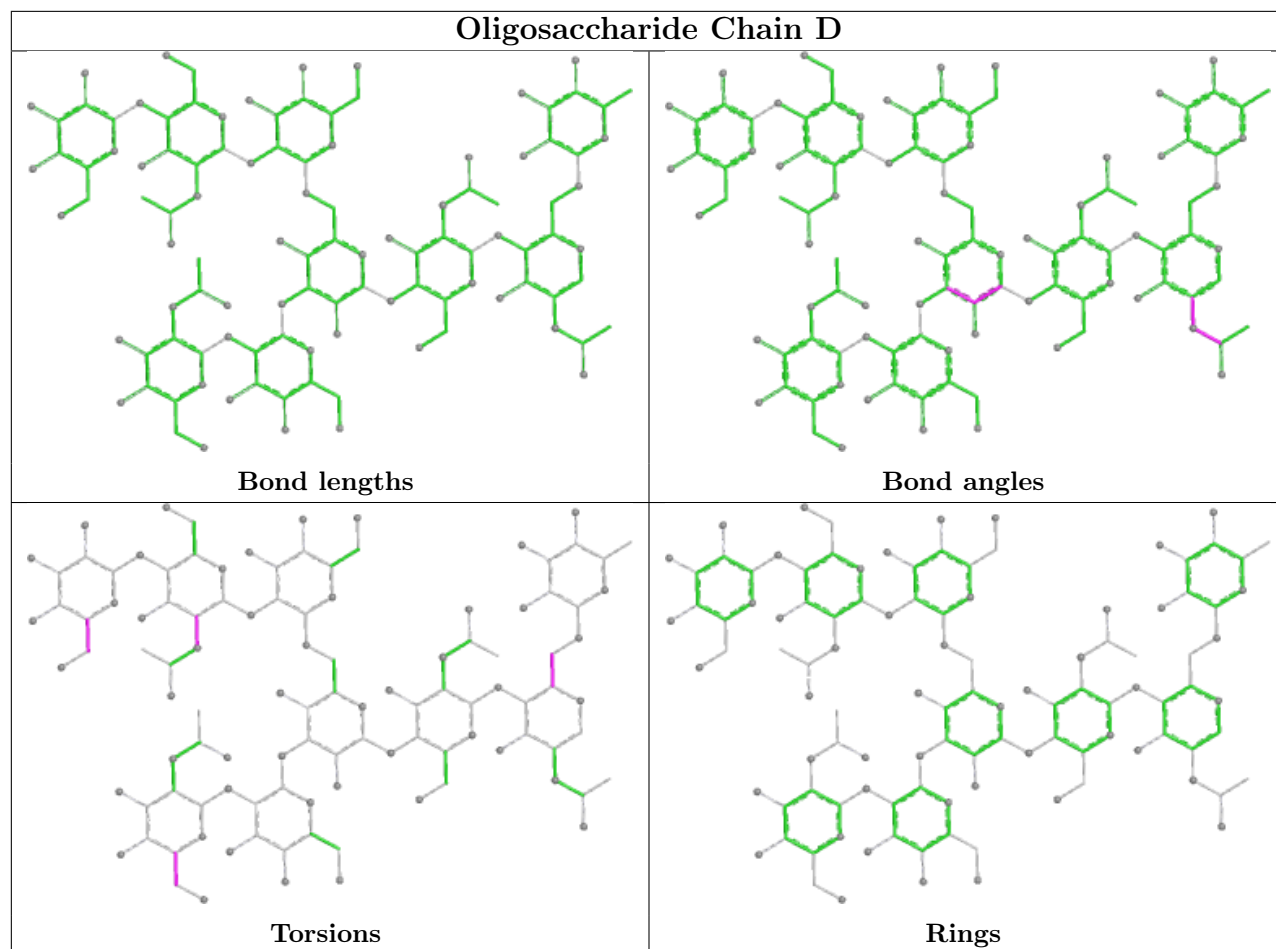
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	9	FUL	1	0
2	D	1	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 [i](#)

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

EDS failed to run properly - this section is therefore empty.

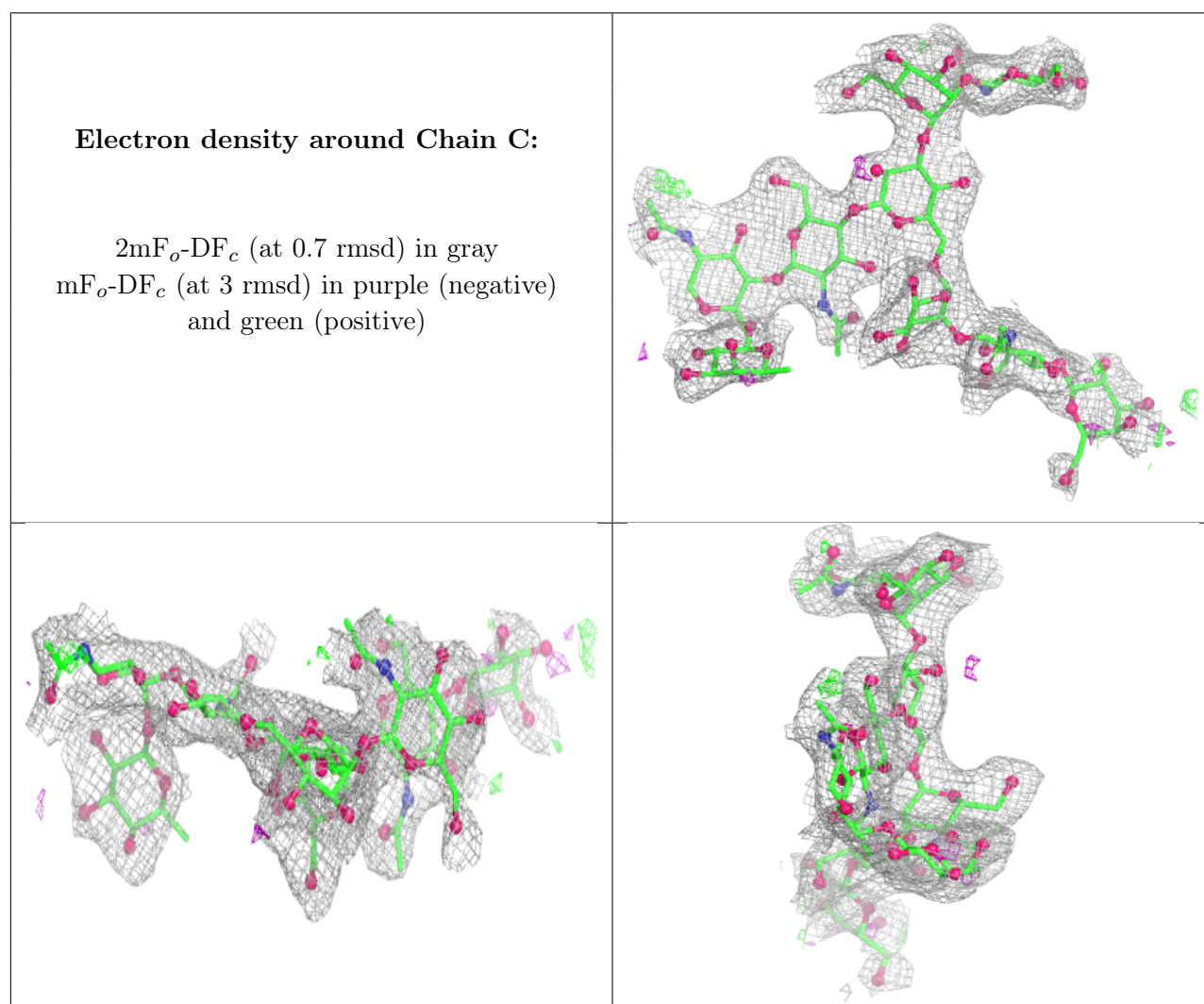
6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

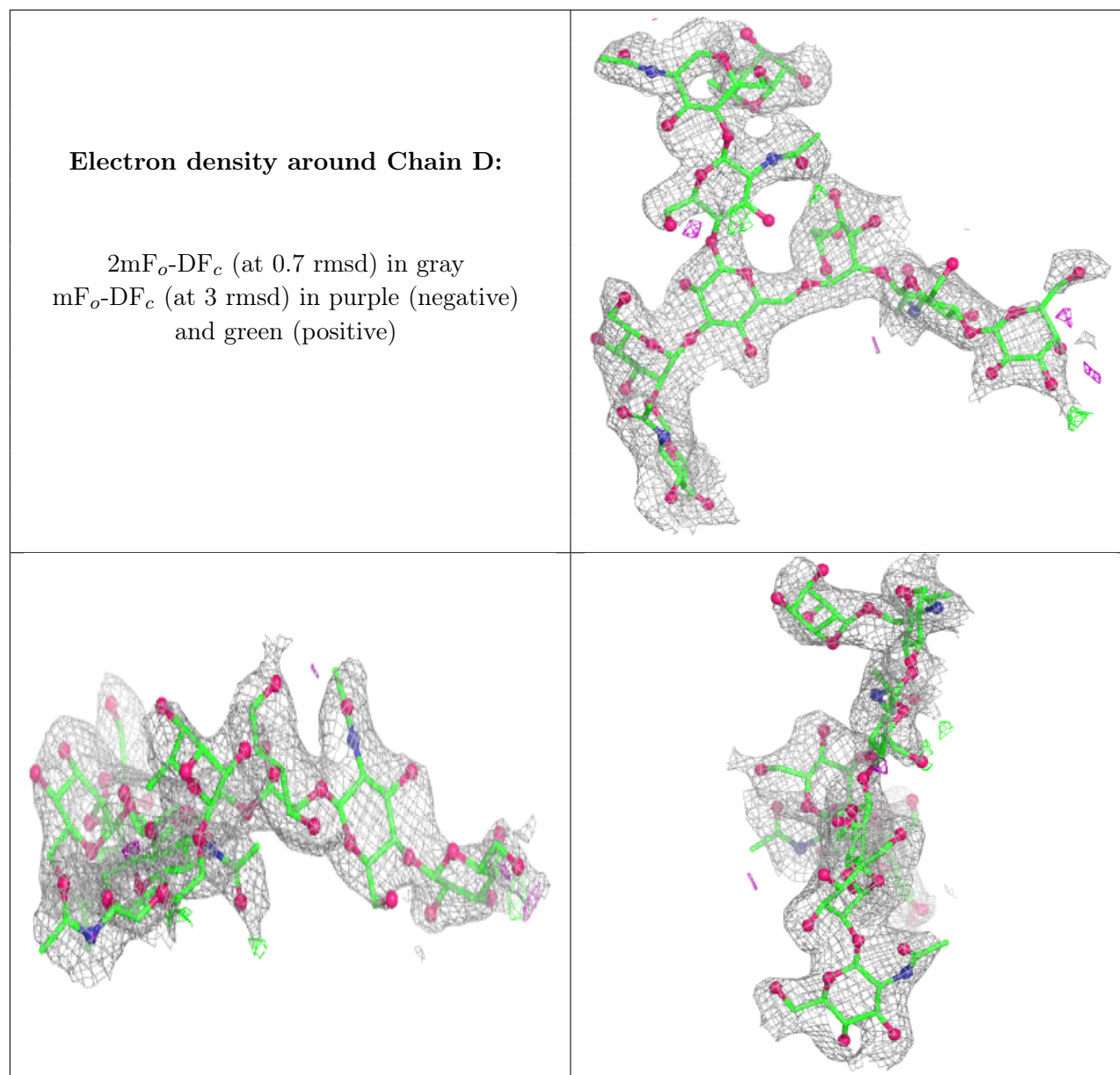
EDS failed to run properly - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS failed to run properly - this section is therefore empty.

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.





6.4 Ligands [i](#)

EDS failed to run properly - this section is therefore empty.

6.5 Other polymers [i](#)

EDS failed to run properly - this section is therefore empty.