



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:04 PM UTC

PDB ID : 4FOM / pdb_00004fom
Title : Crystal structure of human nectin-3 full ectodomain (D1-D3)
Authors : Harrison, O.J.; Jin, X.; Brasch, J.; Shapiro, L.
Deposited on : 2012-06-20
Resolution : 3.93 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

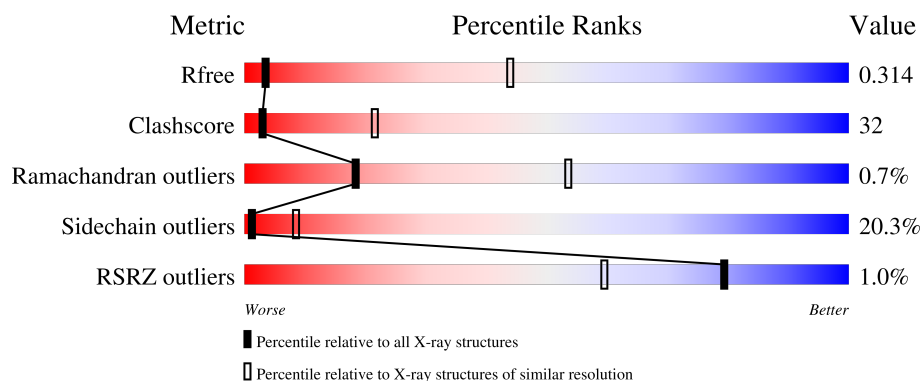
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.93 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	1022 (4.12-3.76)
Clashscore	190562	1000 (4.10-3.78)
Ramachandran outliers	187476	1009 (4.12-3.76)
Sidechain outliers	187428	1002 (4.12-3.76)
RSRZ outliers	180081	1022 (4.12-3.76)

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	308	50% 36% 12%
2	B	5	20% 80%
2	E	5	20% 60% 20%
3	C	3	67% 33%
4	D	6	17% 67% 17%

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 criteria:

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

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2588 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.

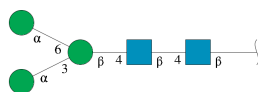
- Molecule 1 is a protein called Poliovirus receptor-related protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	0	0
			2342	1496	396	443	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	360	HIS	-	expression tag	UNP Q9NQS3
A	361	HIS	-	expression tag	UNP Q9NQS3
A	362	HIS	-	expression tag	UNP Q9NQS3
A	363	HIS	-	expression tag	UNP Q9NQS3
A	364	HIS	-	expression tag	UNP Q9NQS3
A	365	HIS	-	expression tag	UNP Q9NQS3

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



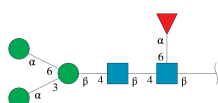
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	E	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



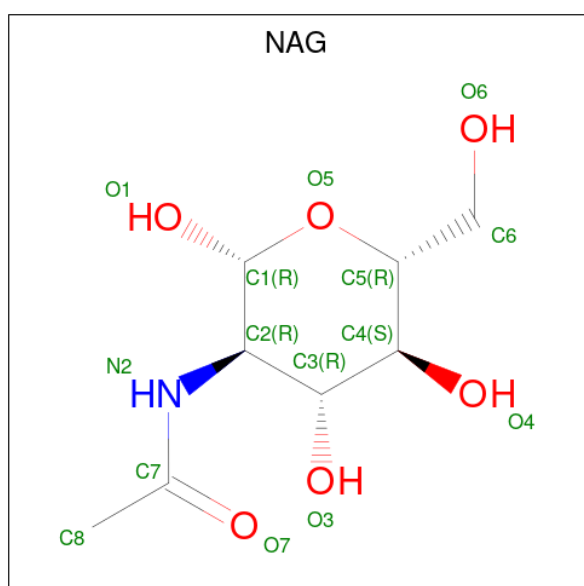
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[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
4	D	6	Total	C	N	O	0	0	0
			71	40	2	29			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

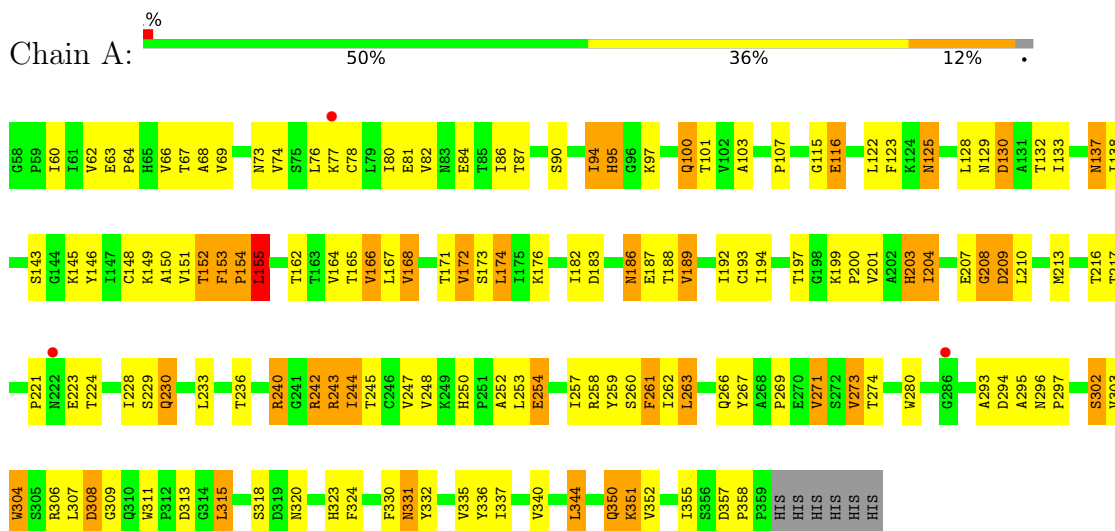


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		

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: Poliovirus receptor-related protein 3



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



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



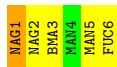
- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 4: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-[α -L-fucopyranose-(1-6)]2-acetamido-2-deoxy- β -D-glucopyranose

Chain D: 17% 67% 17%



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	131.87Å 131.87Å 247.54Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.93 20.00 – 3.93	Depositor EDS
% Data completeness (in resolution range)	98.4 (20.00-3.93) 97.8 (20.00-3.93)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.72 (at 3.98Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.247 , 0.279 0.261 , 0.314	Depositor DCC
R_{free} test set	559 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	142.2	Xtriage
Anisotropy	0.280	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 92.2	EDS
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
F_o, F_c correlation	0.92	EDS
Total number of atoms	2588	wwPDB-VP
Average B, all atoms (Å ²)	170.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.48% 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, NAG, MAN, BMA

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.13	4/2403 (0.2%)	0.97	9/3279 (0.3%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	186	ASN	CG-ND2	-7.36	1.17	1.33
1	A	137	ASN	N-CA	5.41	1.53	1.46
1	A	320	ASN	N-CA	5.02	1.51	1.46
1	A	203	HIS	CE1-NE2	5.02	1.37	1.32

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	154	PRO	N-CA-CB	-15.11	85.98	102.60
1	A	154	PRO	N-CD-CG	-13.47	87.64	103.80
1	A	308	ASP	N-CA-C	-10.01	94.88	109.59
1	A	153	PHE	CB-CA-C	-9.14	97.34	110.13
1	A	154	PRO	N-CA-C	8.39	133.91	112.10
1	A	95	HIS	N-CA-C	-7.71	100.92	110.41
1	A	153	PHE	N-CA-C	6.15	122.21	109.11
1	A	208	GLY	N-CA-C	6.10	122.22	114.66
1	A	207	GLU	N-CA-C	5.99	120.59	113.16

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	2342	0	2300	140	0
2	B	61	0	52	13	0
2	E	61	0	52	4	0
3	C	39	0	34	4	0
4	D	71	0	61	9	0
5	A	14	0	13	0	0
All	All	2588	0	2512	162	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 32.

All (162) 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:186:ASN:CG	4:D:1:NAG:C1	1.93	1.41
1:A:186:ASN:OD1	4:D:1:NAG:C1	1.84	1.23
3:C:2:NAG:C6	3:C:3:BMA:H2	1.80	1.09
1:A:243:ARG:HH11	1:A:243:ARG:HG3	1.07	1.08
1:A:186:ASN:ND2	4:D:1:NAG:O5	1.92	1.03
3:C:2:NAG:H61	3:C:3:BMA:C2	1.90	1.02
1:A:69:VAL:HG22	1:A:167:LEU:HD11	1.03	1.02
1:A:176:LYS:HG2	1:A:261:PHE:HZ	1.26	1.01
1:A:69:VAL:CG2	1:A:167:LEU:HD11	1.91	1.00
1:A:63:GLU:O	1:A:162:THR:HG22	1.65	0.97
3:C:2:NAG:H61	3:C:3:BMA:H2	0.97	0.95
2:B:1:NAG:C6	2:B:2:NAG:N2	2.30	0.95
2:B:1:NAG:C6	2:B:2:NAG:HN2	1.79	0.95
1:A:167:LEU:HD12	1:A:167:LEU:O	1.66	0.93
1:A:67:THR:HB	1:A:167:LEU:HD23	1.51	0.92
2:B:1:NAG:H61	2:B:2:NAG:N2	1.85	0.90
2:B:3:BMA:H61	2:B:5:MAN:O2	1.67	0.89
1:A:176:LYS:HG2	1:A:261:PHE:CZ	2.08	0.89
1:A:186:ASN:ND2	4:D:1:NAG:C2	2.35	0.89
1:A:69:VAL:HG22	1:A:167:LEU:CD1	1.98	0.88
1:A:242:ARG:HG3	1:A:242:ARG:HH11	1.38	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:ARG:HH11	1:A:240:ARG:HG3	1.41	0.86
1:A:244:ILE:HG22	1:A:261:PHE:HB3	1.59	0.83
1:A:172:VAL:HG11	1:A:248:VAL:HG11	1.57	0.83
1:A:63:GLU:O	1:A:162:THR:CG2	2.27	0.81
2:B:1:NAG:H62	2:B:2:NAG:HN2	1.46	0.80
1:A:266:GLN:HG2	1:A:297:PRO:HD3	1.64	0.80
1:A:243:ARG:HG3	1:A:243:ARG:NH1	1.88	0.79
2:B:1:NAG:H61	2:B:2:NAG:C1	2.13	0.79
2:B:1:NAG:H61	2:B:2:NAG:HN2	1.42	0.78
1:A:269:PRO:HA	1:A:295:ALA:HB2	1.66	0.78
1:A:189:VAL:CG1	1:A:230:GLN:HG2	2.14	0.78
1:A:331:ASN:O	1:A:332:TYR:HB2	1.83	0.77
1:A:240:ARG:HG3	1:A:240:ARG:NH1	1.95	0.77
1:A:87:THR:HG23	1:A:151:VAL:HG12	1.65	0.76
1:A:168:VAL:HG22	1:A:199:LYS:HB3	1.67	0.76
2:B:1:NAG:H62	2:B:2:NAG:N2	1.98	0.73
1:A:243:ARG:HH11	1:A:243:ARG:CG	1.93	0.73
1:A:318:SER:HB3	1:A:323:HIS:CE1	2.24	0.72
4:D:3:BMA:H62	4:D:5:MAN:H5	1.72	0.71
1:A:248:VAL:CG1	1:A:257:ILE:HG23	2.21	0.71
1:A:67:THR:HB	1:A:167:LEU:CD2	2.20	0.71
1:A:296:ASN:HA	1:A:297:PRO:C	2.16	0.70
1:A:242:ARG:HG3	1:A:242:ARG:NH1	2.03	0.69
1:A:94:ILE:HG12	1:A:94:ILE:O	1.93	0.68
2:B:3:BMA:C6	2:B:5:MAN:O2	2.42	0.68
1:A:74:VAL:HG12	1:A:138:ILE:HD11	1.77	0.67
1:A:186:ASN:ND2	4:D:1:NAG:H2	2.09	0.67
1:A:137:ASN:CG	1:A:137:ASN:O	2.37	0.66
1:A:271:VAL:HG12	1:A:293:ALA:HB2	1.77	0.66
1:A:248:VAL:HG13	1:A:257:ILE:HG23	1.79	0.65
1:A:273:VAL:HG11	1:A:304:TRP:CH2	2.32	0.65
1:A:244:ILE:HG22	1:A:261:PHE:CB	2.25	0.65
1:A:240:ARG:HH11	1:A:240:ARG:CG	2.09	0.65
1:A:189:VAL:HG11	1:A:230:GLN:HG2	1.77	0.65
1:A:60:ILE:HG12	1:A:80:ILE:HG22	1.78	0.64
1:A:78:CYS:HG	1:A:148:CYS:HG	0.66	0.64
1:A:125:ASN:HB3	1:A:130:ASP:HB2	1.80	0.64
1:A:294:ASP:OD2	1:A:294:ASP:C	2.39	0.64
1:A:247:VAL:HG12	1:A:258:ARG:HG2	1.80	0.62
1:A:125:ASN:CB	1:A:130:ASP:HB2	2.30	0.61
1:A:153:PHE:CD2	1:A:153:PHE:C	2.78	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ILE:HG21	1:A:229:SER:HB2	1.82	0.60
1:A:62:VAL:HG12	1:A:162:THR:CG2	2.31	0.60
1:A:68:ALA:O	1:A:167:LEU:HG	2.01	0.60
1:A:97:LYS:HG2	1:A:97:LYS:O	2.02	0.59
1:A:280:TRP:HZ3	1:A:324:PHE:HE2	1.47	0.59
2:E:3:BMA:H62	2:E:5:MAN:H5	1.84	0.59
1:A:302:SER:HB3	1:A:340:VAL:HG22	1.83	0.59
2:B:1:NAG:H61	2:B:2:NAG:C2	2.33	0.58
1:A:192:ILE:HG12	1:A:230:GLN:HG3	1.84	0.57
1:A:174:LEU:HD12	1:A:259:TYR:HB3	1.87	0.57
1:A:197:THR:CG2	2:E:2:NAG:H83	2.35	0.56
2:B:2:NAG:H61	2:B:3:BMA:H2	1.85	0.56
1:A:280:TRP:HZ3	1:A:324:PHE:CE2	2.23	0.56
1:A:351:LYS:HA	1:A:351:LYS:HE2	1.87	0.56
1:A:269:PRO:HA	1:A:295:ALA:CB	2.34	0.55
1:A:189:VAL:HG22	4:D:6:FUC:O4	2.06	0.55
1:A:233:LEU:HD12	1:A:233:LEU:O	2.06	0.55
1:A:100:GLN:HE21	1:A:100:GLN:HA	1.71	0.55
1:A:167:LEU:HD13	1:A:252:ALA:HB2	1.88	0.55
1:A:69:VAL:HA	1:A:167:LEU:CD1	2.38	0.54
1:A:248:VAL:HG13	1:A:257:ILE:CG2	2.37	0.54
1:A:143:SER:HB2	1:A:166:VAL:HG22	1.89	0.54
2:B:5:MAN:H62	2:B:5:MAN:O3	2.08	0.53
1:A:250:HIS:HB3	1:A:253:LEU:HD12	1.90	0.53
1:A:90:SER:HB2	1:A:149:LYS:HB3	1.91	0.53
2:B:1:NAG:H62	2:B:2:NAG:C7	2.40	0.51
2:E:3:BMA:H61	2:E:5:MAN:H3	1.93	0.51
1:A:103:ALA:HB1	1:A:123:PHE:CE2	2.46	0.51
1:A:82:VAL:C	1:A:84:GLU:H	2.19	0.51
1:A:152:THR:OG1	1:A:155:LEU:HG	2.11	0.51
1:A:331:ASN:O	1:A:332:TYR:CB	2.54	0.51
1:A:62:VAL:HG22	1:A:78:CYS:HA	1.93	0.50
1:A:315:LEU:CD2	1:A:332:TYR:HE1	2.24	0.50
1:A:167:LEU:HD12	1:A:167:LEU:C	2.34	0.50
1:A:66:VAL:HB	1:A:164:VAL:HG12	1.93	0.49
1:A:94:ILE:O	1:A:94:ILE:CG1	2.60	0.49
1:A:199:LYS:HB2	1:A:224:THR:HG22	1.93	0.49
1:A:280:TRP:CZ3	1:A:324:PHE:HE2	2.28	0.49
1:A:221:PRO:C	1:A:223:GLU:H	2.20	0.49
1:A:267:TYR:HE2	1:A:295:ALA:HA	1.78	0.49
1:A:130:ASP:C	1:A:132:THR:H	2.19	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:GLU:N	1:A:254:GLU:OE2	2.45	0.48
1:A:62:VAL:CG1	1:A:162:THR:CG2	2.91	0.48
1:A:173:SER:OG	1:A:194:ILE:HB	2.13	0.48
1:A:62:VAL:HG12	1:A:162:THR:HG21	1.94	0.47
1:A:233:LEU:HD21	1:A:263:LEU:HD21	1.97	0.47
1:A:107:PRO:HB3	1:A:128:LEU:HD11	1.95	0.47
1:A:273:VAL:HG11	1:A:304:TRP:HH2	1.76	0.47
4:D:3:BMA:C6	4:D:5:MAN:H3	2.44	0.47
1:A:245:THR:HG22	1:A:260:SER:HA	1.97	0.47
1:A:62:VAL:CG1	1:A:162:THR:HG21	2.45	0.46
1:A:69:VAL:HA	1:A:167:LEU:HD12	1.96	0.46
1:A:208:GLY:O	1:A:209:ASP:C	2.58	0.46
3:C:2:NAG:C6	3:C:3:BMA:C2	2.68	0.46
1:A:183:ASP:OD1	1:A:236:THR:HA	2.16	0.46
1:A:74:VAL:CG1	1:A:138:ILE:HD11	2.44	0.45
1:A:296:ASN:ND2	1:A:297:PRO:HA	2.31	0.45
1:A:197:THR:HG21	2:E:2:NAG:H83	1.98	0.45
1:A:307:LEU:HB3	1:A:335:VAL:HG13	1.99	0.45
1:A:66:VAL:O	1:A:164:VAL:HA	2.17	0.45
4:D:3:BMA:H61	4:D:5:MAN:H3	1.99	0.45
1:A:82:VAL:C	1:A:84:GLU:N	2.75	0.44
1:A:228:ILE:HG13	1:A:228:ILE:O	2.17	0.44
1:A:267:TYR:CE2	1:A:295:ALA:HA	2.53	0.44
1:A:153:PHE:O	1:A:153:PHE:HD2	2.02	0.43
1:A:77:LYS:HA	1:A:132:THR:HG23	2.00	0.43
1:A:315:LEU:CD2	1:A:332:TYR:CE1	3.01	0.43
1:A:199:LYS:CB	1:A:224:THR:HG22	2.49	0.43
1:A:172:VAL:CG1	1:A:248:VAL:HG11	2.40	0.43
1:A:311:TRP:CE2	1:A:315:LEU:HD12	2.54	0.43
1:A:194:ILE:HG12	1:A:228:ILE:HG22	2.01	0.42
1:A:143:SER:CB	1:A:166:VAL:HG22	2.50	0.42
1:A:154:PRO:O	1:A:155:LEU:C	2.62	0.42
1:A:186:ASN:O	1:A:187:GLU:C	2.61	0.42
1:A:318:SER:HB3	1:A:323:HIS:HE1	1.81	0.42
1:A:337:ILE:HB	1:A:350:GLN:HB3	2.01	0.42
1:A:344:LEU:HD13	1:A:344:LEU:HA	1.85	0.42
1:A:94:ILE:C	1:A:95:HIS:O	2.58	0.42
1:A:76:LEU:HB2	1:A:133:ILE:HG13	2.00	0.42
1:A:97:LYS:HE2	1:A:97:LYS:HB3	1.90	0.41
1:A:115:GLY:HA2	1:A:116:GLU:HA	1.78	0.41
1:A:60:ILE:HD11	1:A:150:ALA:HB3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:PRO:O	1:A:201:VAL:C	2.63	0.41
1:A:357:ASP:HA	1:A:358:PRO:HD3	1.82	0.41
1:A:193:CYS:SG	1:A:204:ILE:HD11	2.60	0.41
1:A:306:ARG:HD2	1:A:336:TYR:CE1	2.55	0.41
1:A:94:ILE:HD13	1:A:145:LYS:O	2.21	0.41
1:A:130:ASP:C	1:A:132:THR:N	2.79	0.41
1:A:243:ARG:NH1	1:A:243:ARG:CG	2.63	0.41
1:A:306:ARG:NH1	1:A:308:ASP:HB2	2.36	0.41
1:A:308:ASP:O	1:A:309:GLY:C	2.64	0.41
1:A:153:PHE:HA	1:A:154:PRO:HA	1.74	0.41
1:A:63:GLU:O	1:A:162:THR:HG21	2.15	0.40
1:A:69:VAL:HA	1:A:167:LEU:O	2.21	0.40
1:A:63:GLU:HA	1:A:64:PRO:HD3	1.99	0.40
1:A:66:VAL:HG12	1:A:67:THR:H	1.86	0.40
1:A:153:PHE:CD2	1:A:153:PHE:O	2.73	0.40
1:A:221:PRO:C	1:A:223:GLU:N	2.79	0.40
1:A:245:THR:HG22	1:A:260:SER:CB	2.52	0.40
1:A:271:VAL:HG12	1:A:293:ALA:CB	2.49	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	300/308 (97%)	262 (87%)	36 (12%)	2 (1%)	18 53

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	155	LEU
1	A	254	GLU

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	261/267 (98%)	208 (80%)	53 (20%)	1 8

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	81	GLU
1	A	86	ILE
1	A	94	ILE
1	A	100	GLN
1	A	101	THR
1	A	116	GLU
1	A	122	LEU
1	A	125	ASN
1	A	129	ASN
1	A	130	ASP
1	A	146	TYR
1	A	152	THR
1	A	155	LEU
1	A	165	THR
1	A	166	VAL
1	A	168	VAL
1	A	171	THR
1	A	172	VAL
1	A	174	LEU
1	A	182	ILE
1	A	188	THR
1	A	189	VAL
1	A	203	HIS
1	A	204	ILE
1	A	209	ASP
1	A	210	LEU
1	A	213	MET
1	A	216	THR
1	A	217	THR

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Mol	Chain	Res	Type
1	A	230	GLN
1	A	240	ARG
1	A	242	ARG
1	A	243	ARG
1	A	244	ILE
1	A	261	PHE
1	A	262	ILE
1	A	263	LEU
1	A	271	VAL
1	A	273	VAL
1	A	274	THR
1	A	302	SER
1	A	303	VAL
1	A	304	TRP
1	A	313	ASP
1	A	315	LEU
1	A	330	PHE
1	A	331	ASN
1	A	344	LEU
1	A	350	GLN
1	A	351	LYS
1	A	352	VAL
1	A	355	ILE

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

Mol	Chain	Res	Type
1	A	100	GLN
1	A	230	GLN
1	A	292	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

19 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	B	1	2,1	14,14,15	0.36	0	17,19,21	0.57	0
2	NAG	B	2	2	14,14,15	0.34	0	17,19,21	0.57	0
2	BMA	B	3	2	11,11,12	0.65	0	15,15,17	0.53	0
2	MAN	B	4	2	11,11,12	0.66	0	15,15,17	0.52	0
2	MAN	B	5	2	11,11,12	0.25	0	15,15,17	0.53	0
3	NAG	C	1	3,1	14,14,15	1.03	1 (7%)	17,19,21	0.91	1 (5%)
3	NAG	C	2	3	14,14,15	1.31	2 (14%)	17,19,21	1.36	2 (11%)
3	BMA	C	3	3	11,11,12	0.26	0	15,15,17	0.53	0
4	NAG	D	1	4,1	14,14,15	1.98	3 (21%)	17,19,21	1.40	2 (11%)
4	NAG	D	2	4	14,14,15	0.75	0	17,19,21	1.47	3 (17%)
4	BMA	D	3	4	11,11,12	0.26	0	15,15,17	0.52	0
4	MAN	D	4	4	11,11,12	0.25	0	15,15,17	0.52	0
4	MAN	D	5	4	11,11,12	0.72	0	15,15,17	0.74	0
4	FUC	D	6	4	10,10,11	0.71	0	14,14,16	0.57	0
2	NAG	E	1	2,1	14,14,15	1.46	2 (14%)	17,19,21	2.15	2 (11%)
2	NAG	E	2	2	14,14,15	0.86	0	17,19,21	1.13	1 (5%)
2	BMA	E	3	2	11,11,12	0.25	0	15,15,17	0.53	0
2	MAN	E	4	2	11,11,12	0.26	0	15,15,17	0.52	0
2	MAN	E	5	2	11,11,12	0.73	0	15,15,17	0.88	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	B	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	2/6/23/26	0/1/1/1
2	BMA	B	3	2	-	0/2/19/22	0/1/1/1
2	MAN	B	4	2	-	2/2/19/22	0/1/1/1

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

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1	NAG	C1-C2	-5.51	1.44	1.52
2	E	1	NAG	C1-C2	3.40	1.57	1.52
4	D	1	NAG	O5-C1	-3.38	1.38	1.43
2	E	1	NAG	O7-C7	2.61	1.29	1.23
3	C	1	NAG	C1-C2	2.53	1.55	1.52
3	C	2	NAG	C2-N2	2.46	1.50	1.46
4	D	1	NAG	O5-C5	-2.20	1.39	1.43
3	C	2	NAG	C3-C2	2.01	1.56	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	NAG	C1-O5-C5	7.66	122.45	112.19
3	C	2	NAG	C4-C3-C2	4.18	117.14	111.02
4	D	2	NAG	C4-C3-C2	-3.43	105.99	111.02
4	D	2	NAG	O5-C1-C2	-2.84	106.90	111.29
4	D	1	NAG	C3-C4-C5	2.77	115.26	110.23
4	D	2	NAG	C1-O5-C5	2.76	115.89	112.19
2	E	1	NAG	C2-N2-C7	-2.59	119.42	122.90
4	D	1	NAG	O5-C1-C2	2.38	114.97	111.29
3	C	1	NAG	C1-O5-C5	2.37	115.36	112.19
2	E	2	NAG	O5-C1-C2	-2.13	107.99	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	NAG	C3-C4-C5	2.13	114.10	110.23

There are no chirality outliers.

All (25) torsion outliers are listed below:

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

There are no ring outliers.

13 monomers are involved in 30 short contacts:

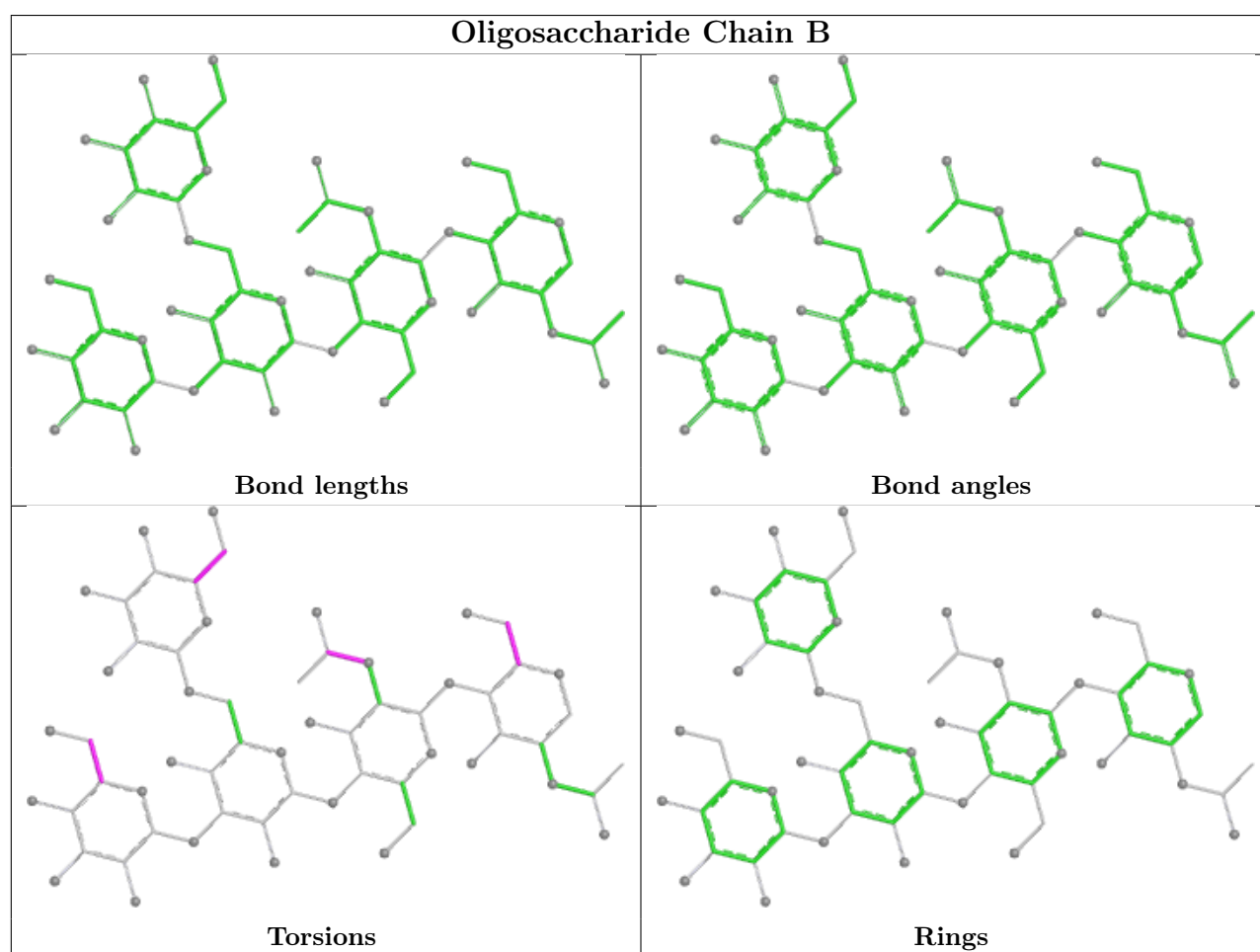
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	3	BMA	2	0
2	B	2	NAG	10	0
3	C	2	NAG	4	0
2	E	5	MAN	2	0
4	D	3	BMA	3	0

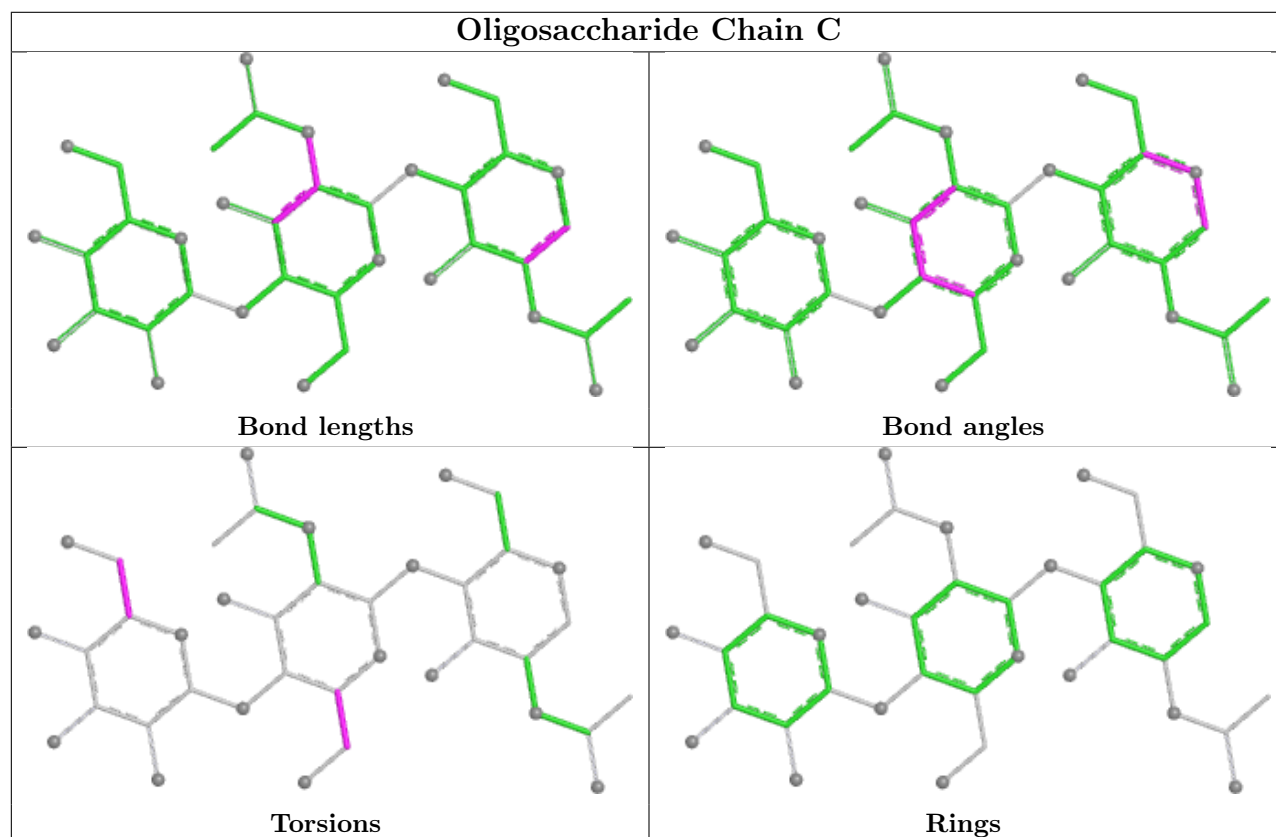
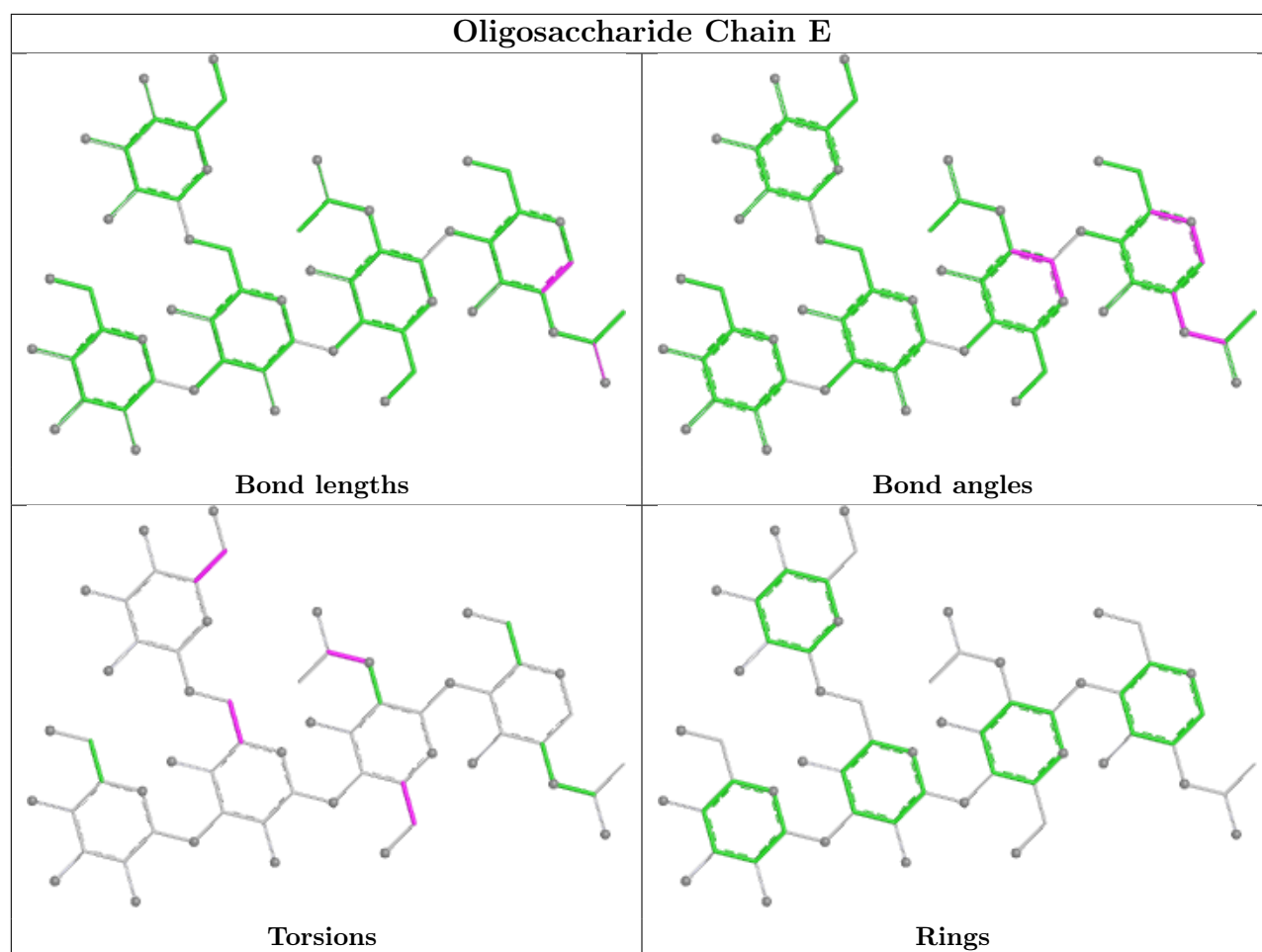
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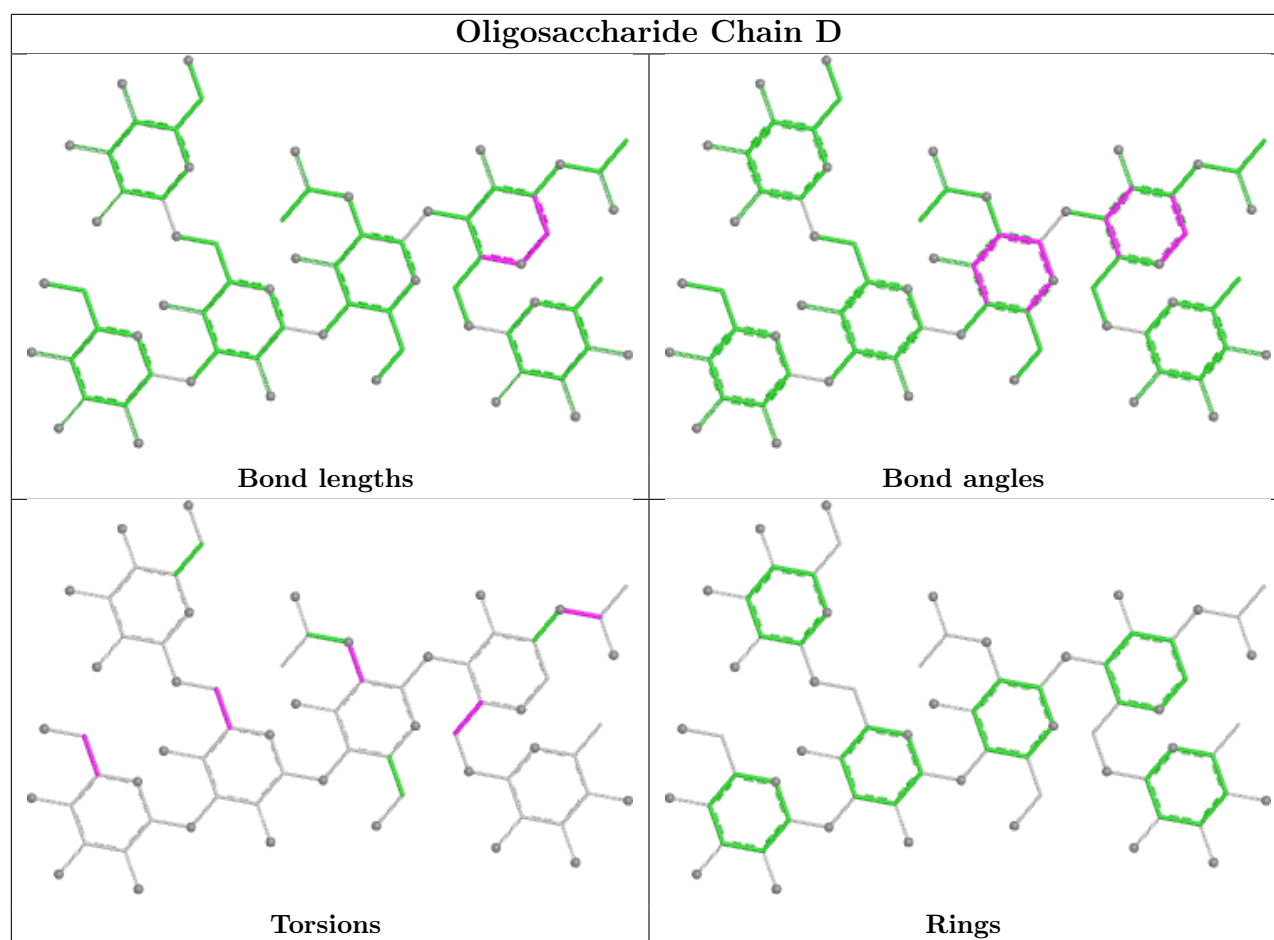
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3	BMA	3	0
4	D	5	MAN	3	0
4	D	1	NAG	5	0
3	C	3	BMA	4	0
2	B	1	NAG	9	0
2	E	2	NAG	2	0
4	D	6	FUC	1	0
2	B	5	MAN	3	0

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](#)

1 ligand is 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	$\# Z > 2$
5	NAG	A	420	1	14,14,15	1.61	5 (35%)	17,19,21	0.76	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
5	NAG	A	420	1	-	2/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	420	NAG	C1-C2	3.43	1.57	1.52
5	A	420	NAG	C3-C2	2.16	1.57	1.52
5	A	420	NAG	C4-C5	2.13	1.57	1.53
5	A	420	NAG	C2-N2	2.13	1.49	1.46
5	A	420	NAG	C4-C3	2.06	1.57	1.52

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	420	NAG	O5-C5-C6-O6
5	A	420	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	302/308 (98%)	0.15	3 (0%) 79 59	107, 169, 237, 278	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	286	GLY	3.4
1	A	77	LYS	2.4
1	A	222	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.

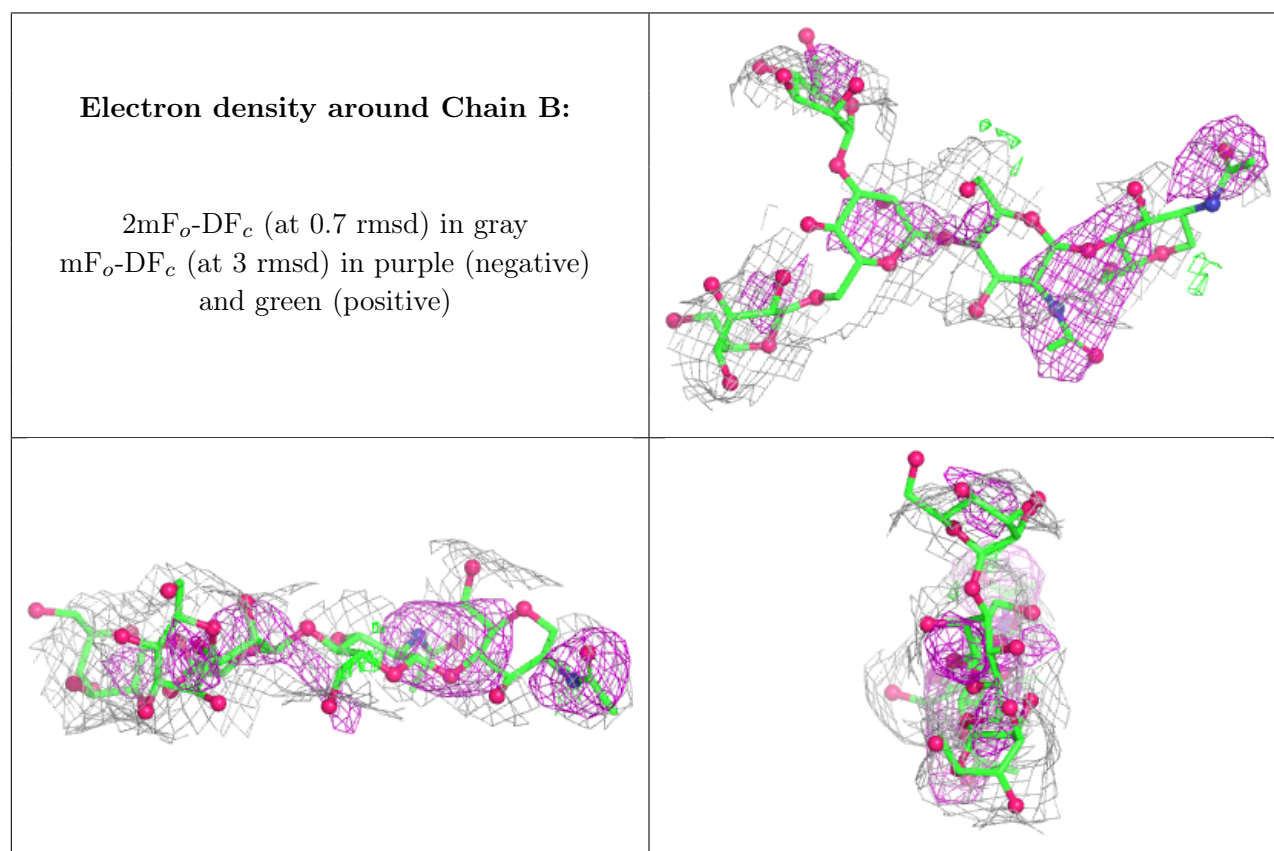
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	C	2	14/15	-0.13	0.19	174,189,197,198	0
2	MAN	B	4	11/12	-0.04	0.16	231,233,233,234	0
3	BMA	C	3	11/12	0.03	0.18	210,211,212,212	0
2	MAN	B	5	11/12	0.28	0.15	247,249,250,250	0
4	MAN	D	5	11/12	0.31	0.13	208,211,212,212	0
3	NAG	C	1	14/15	0.34	0.14	136,149,154,160	0
2	NAG	B	2	14/15	0.42	0.16	180,185,189,196	0
2	MAN	E	5	11/12	0.47	0.23	135,141,142,143	0
2	BMA	B	3	11/12	0.51	0.11	222,226,232,236	0

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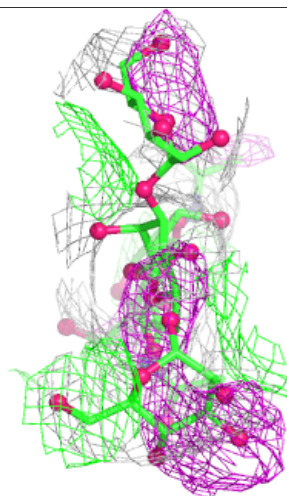
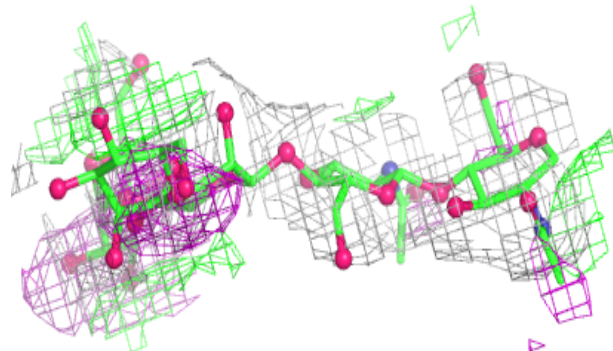
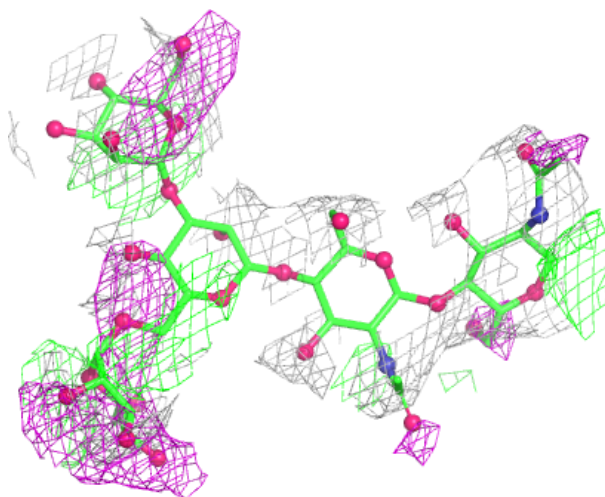
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	BMA	D	3	11/12	0.55	0.10	191,196,198,199	0
2	BMA	E	3	11/12	0.55	0.12	140,143,145,146	0
4	MAN	D	4	11/12	0.56	0.10	207,209,209,210	0
2	MAN	E	4	11/12	0.57	0.12	150,153,157,159	0
2	NAG	B	1	14/15	0.60	0.12	113,127,143,147	0
4	FUC	D	6	10/11	0.62	0.26	106,111,113,114	0
4	NAG	D	1	14/15	0.81	0.11	111,119,122,127	0
2	NAG	E	2	14/15	0.85	0.13	114,122,128,130	0
4	NAG	D	2	14/15	0.86	0.09	142,156,164,170	0
2	NAG	E	1	14/15	0.88	0.12	93,98,103,104	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.



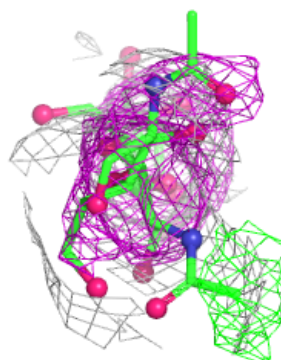
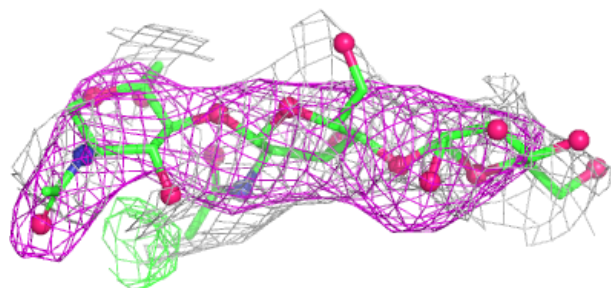
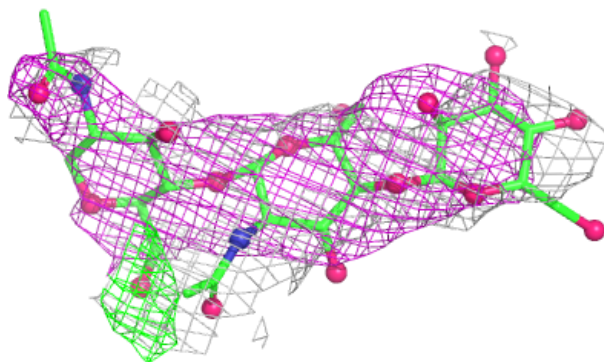
Electron density around Chain E:

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

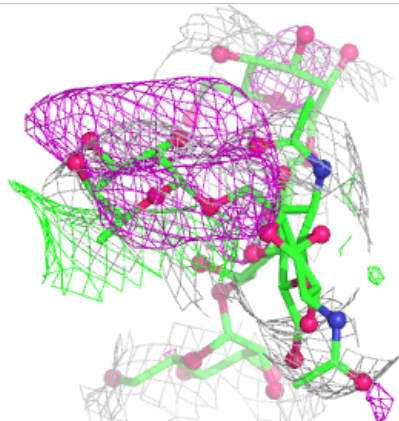
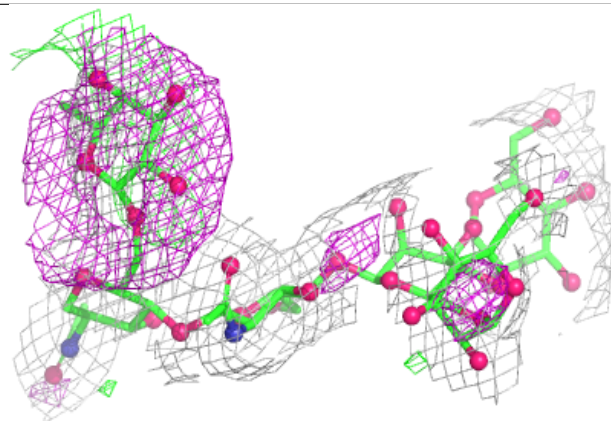
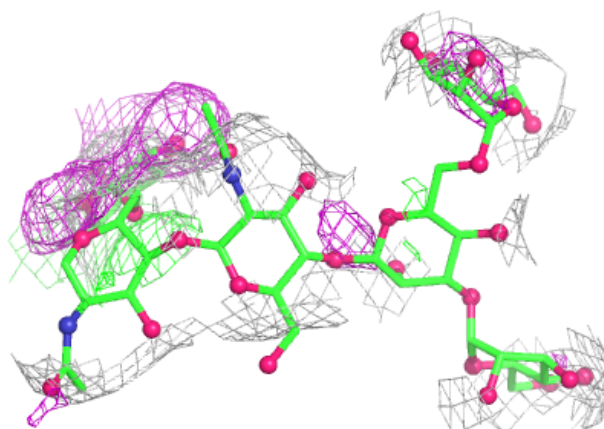


Electron density around Chain C:

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

**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](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	A	420	14/15	0.78	0.26	128,136,139,139	0

6.5 Other polymers [i](#)

There are no such residues in this entry.