



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:18 AM UTC

PDB ID : 4LXS / pdb_00004lxs
Title : Structure of the Toll - Spatzle complex, a molecular hub in Drosophila development and innate immunity (glycosylated form)
Authors : Stelter, M.; Parthier, C.; Breithaupt, C.; Stubbs, M.T.
Deposited on : 2013-07-30
Resolution : 3.30 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

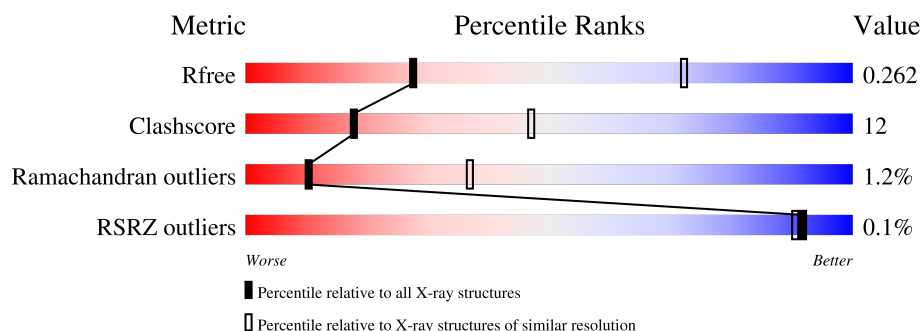
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

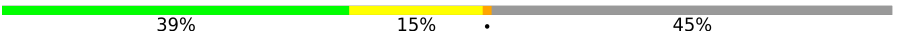
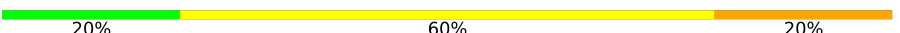
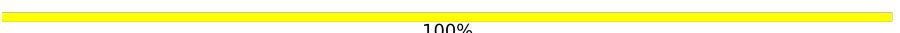
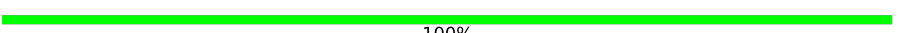
The reported resolution of this entry is 3.30 Å.

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.



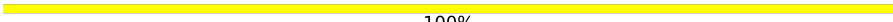
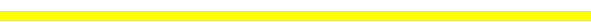
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	783	
2	J	114	
2	K	114	
3	B	5	
4	C	2	
4	E	2	

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Mol	Chain	Length	Quality of chain
4	F	2	 100%
4	G	2	 100%
4	H	2	 50%  50%
5	D	3	 33%  67%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7311 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 Protein toll.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	747	Total	C	N	O	S	0	0	0
			5977	3749	1080	1108	40			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	803	THR	-	expression tag	UNP P08953
A	804	GLY	-	expression tag	UNP P08953
A	805	HIS	-	expression tag	UNP P08953
A	806	HIS	-	expression tag	UNP P08953
A	807	HIS	-	expression tag	UNP P08953
A	808	HIS	-	expression tag	UNP P08953
A	809	HIS	-	expression tag	UNP P08953
A	810	HIS	-	expression tag	UNP P08953

- Molecule 2 is a protein called Protein spaetzle C-106.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	J	63	Total	C	N	O	S	0	0	0
			506	317	88	94	7			
2	K	65	Total	C	N	O	S	0	0	0
			503	313	89	94	7			

There are 16 discrepancies between the modelled and reference sequences:

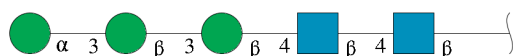
Chain	Residue	Modelled	Actual	Comment	Reference
J	107	LEU	-	expression tag	UNP P48607
J	108	GLU	-	expression tag	UNP P48607
J	109	HIS	-	expression tag	UNP P48607
J	110	HIS	-	expression tag	UNP P48607
J	111	HIS	-	expression tag	UNP P48607
J	112	HIS	-	expression tag	UNP P48607

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Chain	Residue	Modelled	Actual	Comment	Reference
J	113	HIS	-	expression tag	UNP P48607
J	114	HIS	-	expression tag	UNP P48607
K	107	LEU	-	expression tag	UNP P48607
K	108	GLU	-	expression tag	UNP P48607
K	109	HIS	-	expression tag	UNP P48607
K	110	HIS	-	expression tag	UNP P48607
K	111	HIS	-	expression tag	UNP P48607
K	112	HIS	-	expression tag	UNP P48607
K	113	HIS	-	expression tag	UNP P48607
K	114	HIS	-	expression tag	UNP P48607

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



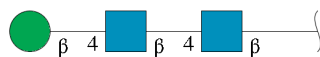
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	B	5	Total	C	N	O	0	0	0
			61	34	2	25			

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



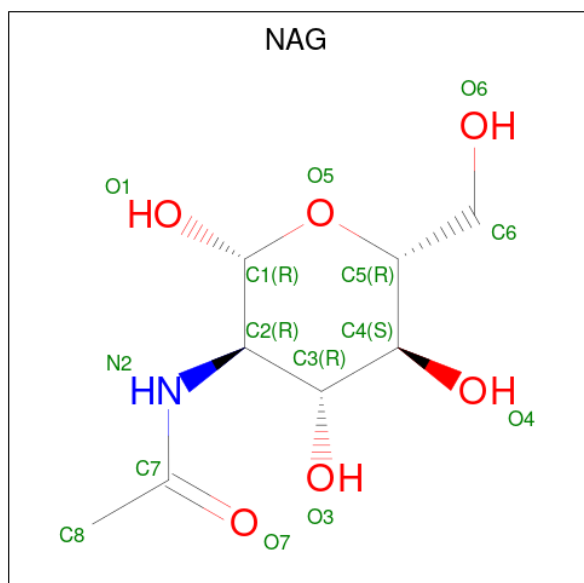
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 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.



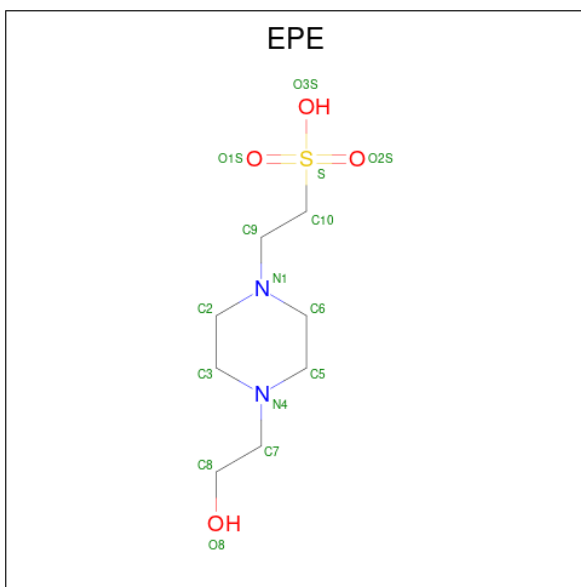
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	D	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 7 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).

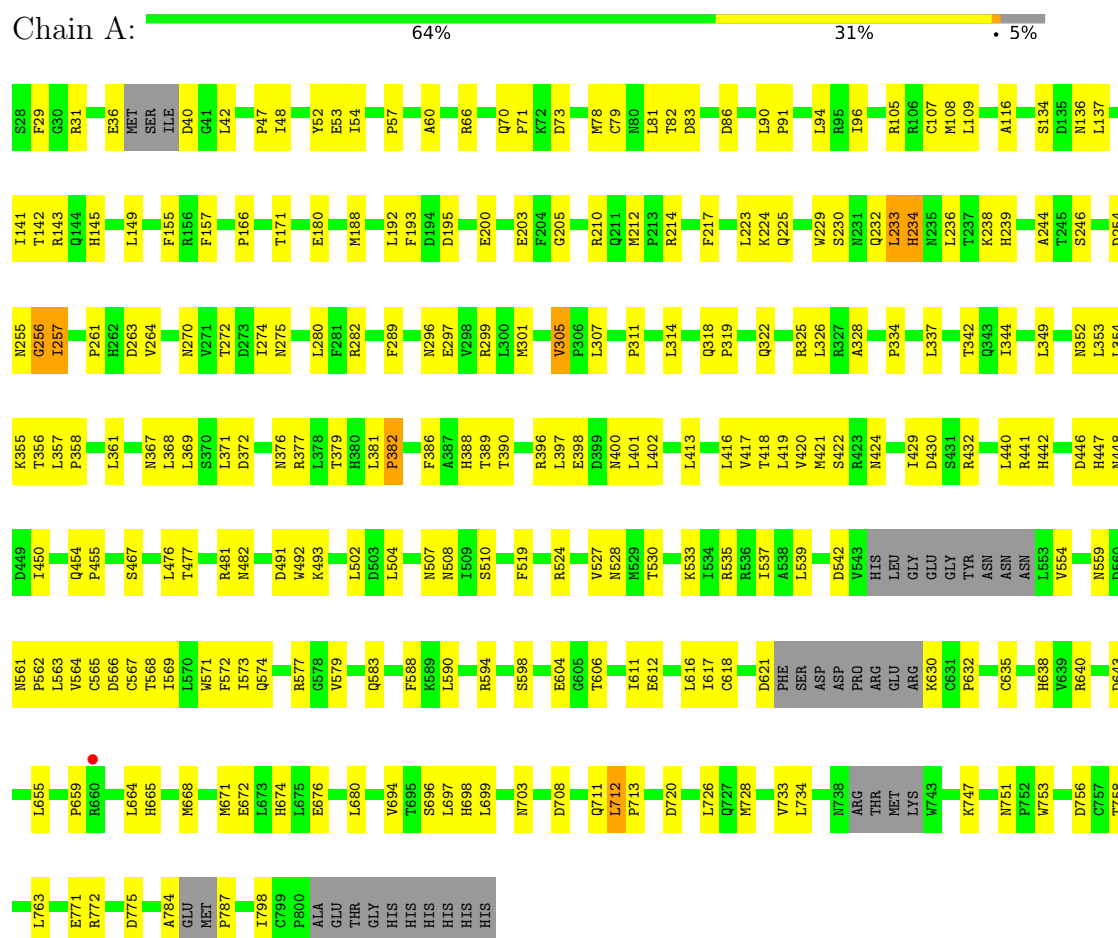


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
7	A	1	15	8	2	4	1	0	0

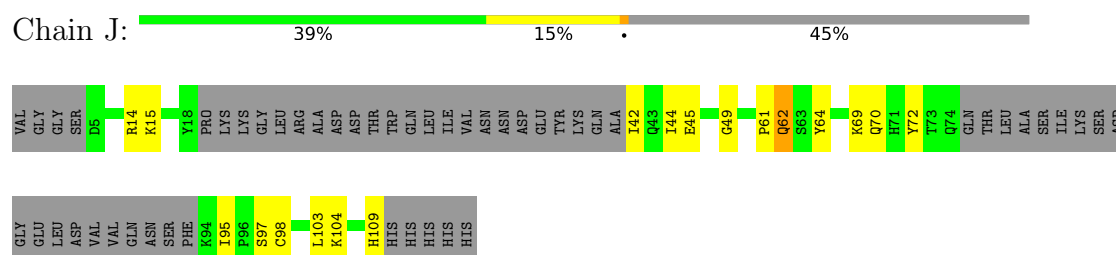
3 Residue-property plots

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 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.

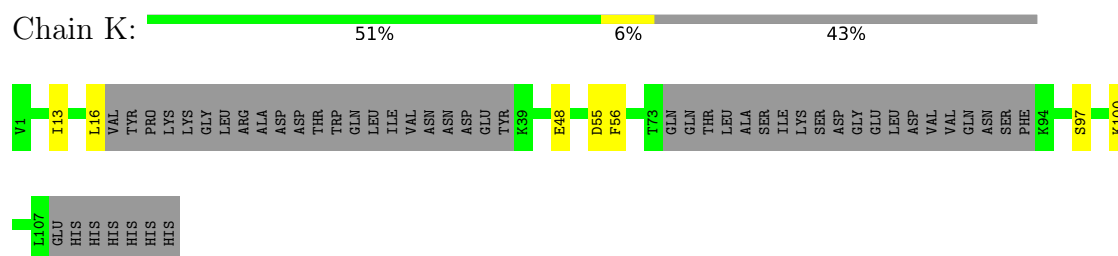
• Molecule 1: Protein toll



• Molecule 2: Protein spaetzle C-106



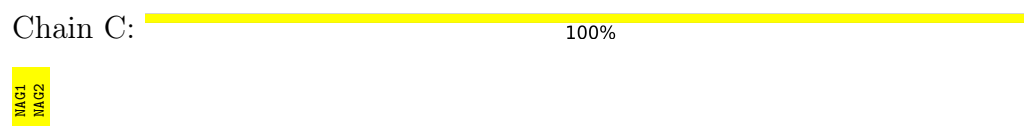
- Molecule 2: Protein spaetzle C-106



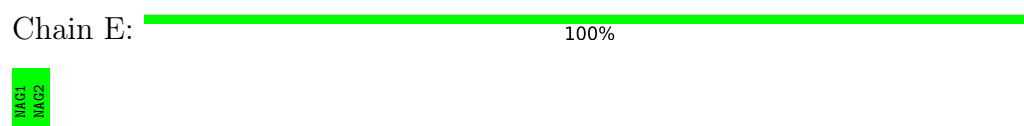
- Molecule 3: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



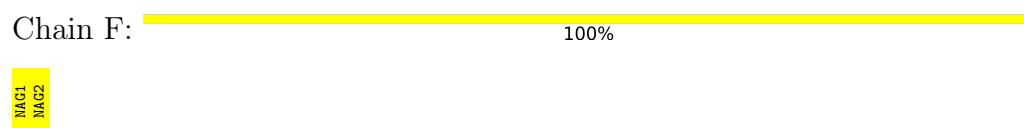
- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



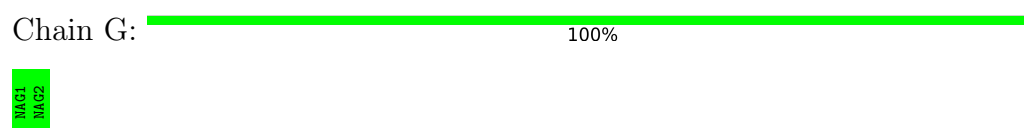
- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



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



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

Chain H:  50% 50%

MAG1
MAG2

- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  33% 67%

MAG1
MAG2
BNA3

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	171.28Å 76.82Å 123.82Å 90.00° 126.29° 90.00°	Depositor
Resolution (Å)	33.95 – 3.30 33.95 – 3.30	Depositor EDS
% Data completeness (in resolution range)	93.6 (33.95-3.30) 93.5 (33.95-3.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 3.32Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.207 , 0.264 (Not available) , 0.262	Depositor DCC
R_{free} test set	926 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	74.2	Xtriage
Anisotropy	0.517	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 7.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7311	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EPE, MAN, NAG, 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	0.36	0/6091	0.84	11/8271 (0.1%)
2	J	0.36	0/515	0.82	1/690 (0.1%)
2	K	0.32	0/510	0.75	0/681
All	All	0.36	0/7116	0.83	12/9642 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	62	GLN	N-CA-C	7.30	119.31	111.36
1	A	70	GLN	CA-C-N	6.00	127.34	119.84
1	A	70	GLN	C-N-CA	6.00	127.34	119.84
1	A	712	LEU	CA-C-N	5.67	125.57	119.85
1	A	712	LEU	C-N-CA	5.67	125.57	119.85
1	A	606	THR	CA-C-N	5.34	126.09	120.11
1	A	606	THR	C-N-CA	5.34	126.09	120.11
1	A	612	GLU	CA-C-N	5.30	126.46	119.84
1	A	612	GLU	C-N-CA	5.30	126.46	119.84
1	A	542	ASP	N-CA-C	5.10	116.05	108.86
1	A	305	VAL	CA-C-N	5.05	126.15	119.84
1	A	305	VAL	C-N-CA	5.05	126.15	119.84

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5977	0	5951	161	0
2	J	506	0	482	16	0
2	K	503	0	492	8	0
3	B	61	0	52	1	0
4	C	28	0	25	1	0
4	E	28	0	25	0	0
4	F	28	0	25	1	0
4	G	28	0	25	0	0
4	H	28	0	25	0	0
5	D	39	0	34	2	0
6	A	70	0	65	1	0
7	A	15	0	17	1	0
All	All	7311	0	7218	179	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 12.

All (179) 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:282:ARG:HH12	1:A:305:VAL:HB	1.39	0.87
1:A:640:ARG:NH2	1:A:643:ASP:OD2	2.17	0.76
1:A:270:ASN:N	1:A:270:ASN:OD1	2.25	0.70
1:A:617:ILE:HG22	1:A:640:ARG:HB3	1.73	0.70
1:A:256:GLY:HA2	1:A:280:LEU:HD13	1.74	0.70
1:A:446:ASP:OD1	1:A:447:HIS:HD2	1.75	0.69
1:A:696:SER:OG	1:A:698:HIS:NE2	2.19	0.69
1:A:573:ILE:HG12	1:A:611:ILE:HD11	1.73	0.69
1:A:441:ARG:HB3	1:A:476:LEU:HD12	1.76	0.66
1:A:567:CYS:N	1:A:618:CYS:SG	2.68	0.65
1:A:467:SER:OG	1:A:491:ASP:OD2	2.09	0.64
1:A:142:THR:H	1:A:145:HIS:HD2	1.46	0.64
1:A:230:SER:OG	1:A:254:ASP:OD2	2.15	0.63
1:A:274:ILE:HD12	1:A:289:PHE:HE1	1.64	0.63
1:A:232:GLN:O	1:A:234:HIS:N	2.32	0.63
1:A:519:PHE:HB2	1:A:554:VAL:HG21	1.81	0.62
2:J:14:ARG:HA	2:J:45:GLU:HA	1.80	0.62
1:A:655:LEU:HD23	1:A:659:PRO:HD3	1.81	0.62
1:A:210:ARG:HH11	1:A:232:GLN:HE22	1.47	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:671:MET:HB3	1:A:694:VAL:HA	1.82	0.61
1:A:422:SER:HB3	1:A:446:ASP:OD1	2.00	0.61
1:A:680:LEU:O	1:A:703:ASN:HB2	2.01	0.60
1:A:334:PRO:HG2	1:A:337:LEU:HB2	1.84	0.60
1:A:733:VAL:HG12	1:A:734:LEU:HD13	1.84	0.60
1:A:726:LEU:HB2	1:A:751:ASN:HD22	1.66	0.60
1:A:502:LEU:HB3	1:A:527:VAL:HG22	1.84	0.59
1:A:79:CYS:HB3	1:A:81:LEU:HD13	1.83	0.59
1:A:275:ASN:HB2	2:J:62:GLN:HE21	1.65	0.59
1:A:429:ILE:HD11	1:A:450:ILE:HD13	1.84	0.59
1:A:430:ASP:OD1	1:A:432:ARG:N	2.34	0.59
1:A:772:ARG:HH11	1:A:772:ARG:HG3	1.68	0.58
1:A:566:ASP:OD1	1:A:568:THR:OG1	2.22	0.58
1:A:36:GLU:C	1:A:40:ASP:HB2	2.29	0.58
1:A:528:ASN:OD1	1:A:530:THR:OG1	2.20	0.57
1:A:210:ARG:HH11	1:A:232:GLN:NE2	2.02	0.56
1:A:402:LEU:HB2	1:A:424:ASN:HD22	1.70	0.56
1:A:396:ARG:NH1	1:A:398:GLU:OE2	2.39	0.56
1:A:756:ASP:OD1	1:A:758:THR:HG22	2.05	0.56
2:K:55:ASP:HB3	2:K:56:PHE:CD2	2.41	0.56
1:A:282:ARG:NH1	1:A:305:VAL:HB	2.15	0.55
1:A:371:LEU:HD12	1:A:372:ASP:H	1.72	0.55
1:A:574:GLN:OE1	1:A:577:ARG:NH2	2.40	0.54
1:A:322:GLN:HA	1:A:344:ILE:HA	1.90	0.54
1:A:233:LEU:HB3	1:A:236:LEU:HD21	1.89	0.54
1:A:419:LEU:HD21	1:A:421:MET:HE3	1.90	0.54
1:A:142:THR:H	1:A:145:HIS:CD2	2.24	0.53
2:J:44:ILE:HG22	2:J:95:ILE:HG23	1.89	0.53
1:A:223:LEU:O	1:A:246:SER:OG	2.26	0.53
1:A:368:LEU:HD12	1:A:369:LEU:H	1.74	0.53
1:A:354:LEU:HD13	1:A:357:LEU:HD21	1.90	0.53
1:A:771:GLU:OE1	1:A:771:GLU:N	2.37	0.53
1:A:388:HIS:O	1:A:390:THR:N	2.42	0.52
1:A:708:ASP:HB3	1:A:711:GLN:HG3	1.91	0.52
2:J:61:PRO:HB2	2:J:64:TYR:CD1	2.45	0.52
1:A:442:HIS:HA	1:A:477:THR:HB	1.91	0.52
1:A:665:HIS:HB3	1:A:668:MET:HG2	1.92	0.51
1:A:108:MET:HE1	1:A:136:ASN:ND2	2.26	0.51
1:A:664:LEU:HG	1:A:668:MET:HB2	1.93	0.51
1:A:763:LEU:HD13	1:A:798:ILE:HD13	1.92	0.51
1:A:274:ILE:HD12	1:A:289:PHE:CE1	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:ALA:HB2	1:A:149:LEU:HG	1.93	0.51
1:A:418:THR:HB	1:A:442:HIS:HB2	1.93	0.51
1:A:81:LEU:HG	1:A:86:ASP:HB2	1.92	0.51
1:A:210:ARG:NH1	1:A:232:GLN:HE22	2.09	0.51
1:A:78:MET:HE3	2:K:13:ILE:HG23	1.92	0.51
1:A:143:ARG:HD3	1:A:171:THR:HB	1.93	0.50
1:A:446:ASP:O	1:A:448:ASN:ND2	2.43	0.50
2:J:42:ILE:HG21	2:J:95:ILE:HD13	1.92	0.50
1:A:632:PRO:HG2	1:A:635:CYS:HB2	1.93	0.50
2:J:97:SER:HB2	2:K:97:SER:HB2	1.94	0.50
1:A:751:ASN:HB2	1:A:753:TRP:CE2	2.47	0.50
1:A:447:HIS:HA	1:A:482:ASN:HB3	1.94	0.50
2:J:15:LYS:N	2:J:44:ILE:O	2.42	0.50
1:A:358:PRO:HG2	1:A:361:LEU:HB2	1.93	0.49
1:A:419:LEU:HD21	1:A:421:MET:CE	2.43	0.49
1:A:537:ILE:HG13	1:A:563:LEU:HD11	1.95	0.49
2:J:49:GLY:HA3	3:B:4:BMA:H61	1.95	0.49
1:A:594:ARG:NH1	5:D:2:NAG:H2	2.28	0.49
1:A:141:ILE:HG23	1:A:166:PRO:HD2	1.95	0.49
1:A:588:PHE:CD1	1:A:590:LEU:HD13	2.48	0.49
2:J:45:GLU:OE1	2:K:100:LYS:NZ	2.42	0.49
1:A:299:ARG:HB3	1:A:301:MET:HE2	1.94	0.48
1:A:398:GLU:HB3	1:A:420:VAL:HG12	1.95	0.48
1:A:356:THR:OG1	1:A:357:LEU:N	2.45	0.48
2:K:13:ILE:HD11	2:K:48:GLU:OE2	2.13	0.48
1:A:29:PHE:CZ	1:A:54:ILE:HD11	2.49	0.48
1:A:379:THR:HG22	1:A:401:LEU:HB2	1.96	0.48
1:A:371:LEU:HD23	1:A:386:PHE:HE1	1.79	0.47
1:A:107:CYS:H	1:A:134:SER:HB2	1.79	0.47
1:A:82:THR:OG1	1:A:83:ASP:N	2.47	0.47
1:A:326:LEU:HD12	1:A:349:LEU:HD21	1.96	0.47
1:A:491:ASP:OD1	1:A:491:ASP:N	2.39	0.47
1:A:299:ARG:HD2	1:A:325:ARG:HG3	1.97	0.47
1:A:712:LEU:HA	1:A:713:PRO:HD3	1.79	0.47
1:A:775:ASP:OD1	1:A:775:ASP:N	2.38	0.47
1:A:214:ARG:NH1	1:A:239:HIS:ND1	2.63	0.47
1:A:621:ASP:O	1:A:630:LYS:HB2	2.14	0.47
1:A:784:ALA:HB3	1:A:787:PRO:N	2.29	0.47
2:J:98:CYS:SG	2:K:100:LYS:HD3	2.54	0.47
1:A:217:PHE:CD2	1:A:244:ALA:HB2	2.50	0.46
1:A:232:GLN:O	1:A:234:HIS:ND1	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:MET:HE3	1:A:192:LEU:HD13	1.98	0.46
1:A:255:ASN:O	1:A:257:ILE:N	2.48	0.46
1:A:40:ASP:N	1:A:42:LEU:H	2.12	0.46
1:A:640:ARG:NH1	1:A:672:GLU:OE1	2.49	0.46
1:A:561:ASN:HA	1:A:562:PRO:HD3	1.78	0.46
1:A:416:LEU:HD21	1:A:419:LEU:HD13	1.97	0.46
1:A:272:THR:HB	1:A:296:ASN:HB2	1.99	0.45
1:A:454:GLN:HA	1:A:455:PRO:HD2	1.80	0.45
1:A:397:LEU:O	1:A:400:ASN:ND2	2.45	0.45
1:A:772:ARG:HH11	1:A:772:ARG:CG	2.29	0.45
1:A:200:GLU:O	1:A:224:LYS:N	2.28	0.45
1:A:413:LEU:HB3	1:A:416:LEU:HB2	1.99	0.45
1:A:481:ARG:HH12	5:D:1:NAG:H61	1.82	0.45
1:A:565:CYS:HA	1:A:569:ILE:HD11	1.99	0.45
1:A:91:PRO:HB2	1:A:94:LEU:HD13	1.99	0.45
1:A:78:MET:HE1	1:A:105:ARG:CZ	2.47	0.44
1:A:318:GLN:HA	1:A:319:PRO:HD2	1.89	0.44
1:A:539:LEU:O	1:A:571:TRP:HB2	2.16	0.44
2:J:70:GLN:HG3	2:J:72:TYR:CE1	2.52	0.44
2:J:104:LYS:HE2	2:J:104:LYS:HB3	1.76	0.44
1:A:109:LEU:HD12	1:A:137:LEU:HB3	2.00	0.44
1:A:539:LEU:HD12	1:A:572:PHE:HA	1.99	0.44
1:A:31:ARG:HB2	1:A:31:ARG:NH1	2.33	0.44
1:A:53:GLU:HG2	2:K:16:LEU:HD13	2.00	0.43
1:A:261:PRO:O	1:A:264:VAL:HG23	2.18	0.43
1:A:307:LEU:HB3	1:A:328:ALA:HB1	2.00	0.43
1:A:210:ARG:NH1	1:A:232:GLN:NE2	2.65	0.43
1:A:492:TRP:HZ3	1:A:504:LEU:HD11	1.83	0.43
1:A:94:LEU:HD23	1:A:96:ILE:HD11	2.01	0.43
7:A:2024:EPE:H102	7:A:2024:EPE:H62	1.65	0.43
1:A:367:ASN:O	1:A:369:LEU:HD13	2.18	0.43
1:A:71:PRO:C	1:A:73:ASP:H	2.26	0.43
1:A:225:GLN:OE1	2:J:109:HIS:NE2	2.46	0.43
1:A:355:LYS:HD2	1:A:377:ARG:HB2	2.00	0.43
1:A:270:ASN:HD22	4:F:1:NAG:C7	2.32	0.43
1:A:376:ASN:HB2	1:A:400:ASN:OD1	2.19	0.43
1:A:311:PRO:HG2	1:A:314:LEU:HD12	2.00	0.42
1:A:507:ASN:HB3	1:A:508:ASN:H	1.66	0.42
1:A:728:MET:HE3	1:A:728:MET:HB3	1.83	0.42
1:A:48:ILE:HD13	2:K:16:LEU:HD11	2.01	0.42
1:A:200:GLU:HA	1:A:223:LEU:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:617:ILE:HD12	1:A:638:HIS:CG	2.55	0.42
1:A:559:ASN:ND2	1:A:594:ARG:HB2	2.35	0.42
2:J:64:TYR:HB3	2:J:103:LEU:HB3	2.00	0.42
2:J:69:LYS:HA	2:J:69:LYS:HD3	1.64	0.42
1:A:342:THR:O	1:A:342:THR:OG1	2.37	0.42
1:A:47:PRO:HB3	1:A:52:TYR:CE1	2.55	0.42
1:A:263:ASP:N	1:A:263:ASP:OD1	2.51	0.42
4:C:1:NAG:H4	4:C:2:NAG:H2	1.81	0.42
1:A:192:LEU:HD23	1:A:193:PHE:CE1	2.55	0.42
1:A:255:ASN:C	1:A:257:ILE:H	2.28	0.42
1:A:275:ASN:HB2	2:J:62:GLN:NE2	2.33	0.42
1:A:180:GLU:HG3	1:A:203:GLU:HG3	2.02	0.41
1:A:155:PHE:CZ	1:A:157:PHE:HB2	2.55	0.41
1:A:205:GLY:HA2	1:A:229:TRP:O	2.20	0.41
1:A:674:HIS:HB3	1:A:676:GLU:OE2	2.21	0.41
1:A:355:LYS:NZ	1:A:379:THR:HG21	2.35	0.41
1:A:90:LEU:HG	1:A:91:PRO:HD2	2.03	0.41
1:A:212:MET:HE2	1:A:212:MET:HB2	1.99	0.41
1:A:381:LEU:HA	1:A:382:PRO:HD2	1.83	0.41
1:A:697:LEU:HG	1:A:699:LEU:HD21	2.02	0.41
1:A:195:ASP:OD1	1:A:195:ASP:N	2.54	0.41
1:A:441:ARG:HA	1:A:441:ARG:HD3	1.91	0.41
1:A:53:GLU:OE2	1:A:66:ARG:HD3	2.20	0.41
1:A:417:VAL:O	1:A:440:LEU:HD12	2.20	0.41
1:A:352:ASN:HB3	1:A:353:LEU:H	1.68	0.41
1:A:535:ARG:HH12	1:A:564:VAL:CG1	2.33	0.40
1:A:611:ILE:HD12	1:A:616:LEU:HD21	2.03	0.40
1:A:598:SER:O	1:A:604:GLU:HG3	2.22	0.40
1:A:510:SER:HA	1:A:533:LYS:O	2.21	0.40
1:A:559:ASN:HD22	1:A:594:ARG:HB2	1.87	0.40
6:A:2023:NAG:O7	6:A:2023:NAG:O3	2.37	0.40
1:A:57:PRO:HD2	1:A:60:ALA:CB	2.51	0.40
1:A:493:LYS:HG2	1:A:519:PHE:CE1	2.57	0.40
1:A:229:TRP:CD1	1:A:229:TRP:H	2.40	0.40
1:A:297:GLU:OE1	1:A:299:ARG:HD3	2.22	0.40
1:A:566:ASP:HB2	1:A:618:CYS:SG	2.62	0.40
1:A:720:ASP:HA	1:A:747:LYS:HB2	2.03	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	735/783 (94%)	629 (86%)	96 (13%)	10 (1%)	9	33
2	J	57/114 (50%)	49 (86%)	8 (14%)	0	100	100
2	K	59/114 (52%)	54 (92%)	5 (8%)	0	100	100
All	All	851/1011 (84%)	732 (86%)	109 (13%)	10 (1%)	10	37

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	LEU
1	A	234	HIS
1	A	238	LYS
1	A	256	GLY
1	A	389	THR
1	A	524	ARG
1	A	382	PRO
1	A	583	GLN
1	A	579	VAL
1	A	257	ILE

5.3.2 Protein sidechains [i](#)

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

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 ⓘ

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
3	NAG	B	1	3,1	14,14,15	0.48	0	17,19,21	0.49	0
3	NAG	B	2	3	14,14,15	0.66	1 (7%)	17,19,21	0.85	0
3	BMA	B	3	3	11,11,12	1.35	2 (18%)	15,15,17	1.58	4 (26%)
3	BMA	B	4	3	11,11,12	1.02	1 (9%)	15,15,17	1.10	0
3	MAN	B	5	3	11,11,12	1.03	1 (9%)	15,15,17	1.25	1 (6%)
4	NAG	C	1	1,4	14,14,15	0.39	0	17,19,21	0.69	0
4	NAG	C	2	4	14,14,15	0.33	0	17,19,21	0.41	0
5	NAG	D	1	5,1	14,14,15	0.47	0	17,19,21	0.44	0
5	NAG	D	2	5	14,14,15	0.46	0	17,19,21	0.67	0
5	BMA	D	3	5	11,11,12	0.70	0	15,15,17	0.61	0
4	NAG	E	1	1,4	14,14,15	0.31	0	17,19,21	0.54	0
4	NAG	E	2	4	14,14,15	0.41	0	17,19,21	0.52	0
4	NAG	F	1	1,4	14,14,15	0.32	0	17,19,21	0.44	0
4	NAG	F	2	4	14,14,15	0.64	0	17,19,21	0.81	1 (5%)
4	NAG	G	1	1,4	14,14,15	0.39	0	17,19,21	0.39	0
4	NAG	G	2	4	14,14,15	0.18	0	17,19,21	0.42	0
4	NAG	H	1	1,4	14,14,15	0.73	1 (7%)	17,19,21	0.69	0
4	NAG	H	2	4	14,14,15	0.41	0	17,19,21	0.54	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
3	NAG	B	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	B	2	3	-	4/6/23/26	0/1/1/1
3	BMA	B	3	3	-	1/2/19/22	0/1/1/1
3	BMA	B	4	3	-	0/2/19/22	0/1/1/1
3	MAN	B	5	3	-	2/2/19/22	0/1/1/1

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	4	BMA	C2-C3	2.64	1.56	1.52
4	H	1	NAG	C1-C2	2.48	1.55	1.52
3	B	3	BMA	O3-C3	2.35	1.48	1.43
3	B	5	MAN	C2-C3	2.20	1.55	1.52
3	B	2	NAG	O5-C1	-2.10	1.40	1.43
3	B	3	BMA	C4-C5	2.10	1.57	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	3	BMA	C1-C2-C3	-3.67	104.30	109.64
3	B	5	MAN	C1-O5-C5	3.39	116.73	112.19
4	F	2	NAG	C1-O5-C5	3.10	116.34	112.19
3	B	3	BMA	O3-C3-C2	3.08	116.34	110.05
3	B	3	BMA	O3-C3-C4	2.20	115.56	110.38
3	B	3	BMA	O2-C2-C3	-2.01	105.98	110.15

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	H	1	NAG	C1-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
4	E	1	NAG	O5-C5-C6-O6
3	B	2	NAG	O5-C5-C6-O6
5	D	1	NAG	O5-C5-C6-O6
4	E	1	NAG	C4-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
3	B	2	NAG	C8-C7-N2-C2
3	B	2	NAG	O7-C7-N2-C2
4	C	2	NAG	C8-C7-N2-C2
4	C	2	NAG	O7-C7-N2-C2
4	E	1	NAG	C8-C7-N2-C2
4	E	1	NAG	O7-C7-N2-C2
4	F	2	NAG	C8-C7-N2-C2
4	F	2	NAG	O7-C7-N2-C2
4	G	2	NAG	C8-C7-N2-C2
4	G	2	NAG	O7-C7-N2-C2
3	B	2	NAG	C4-C5-C6-O6
4	E	2	NAG	C4-C5-C6-O6
5	D	1	NAG	C4-C5-C6-O6
4	C	1	NAG	O5-C5-C6-O6
4	H	2	NAG	O5-C5-C6-O6
3	B	1	NAG	O5-C5-C6-O6
4	E	2	NAG	O5-C5-C6-O6
3	B	1	NAG	C4-C5-C6-O6
5	D	2	NAG	O5-C5-C6-O6
4	F	1	NAG	O5-C5-C6-O6
3	B	5	MAN	C4-C5-C6-O6
5	D	2	NAG	C1-C2-N2-C7
3	B	5	MAN	O5-C5-C6-O6
3	B	3	BMA	C4-C5-C6-O6
4	C	1	NAG	C4-C5-C6-O6
5	D	2	NAG	C3-C2-N2-C7

There are no ring outliers.

6 monomers are involved in 5 short contacts:

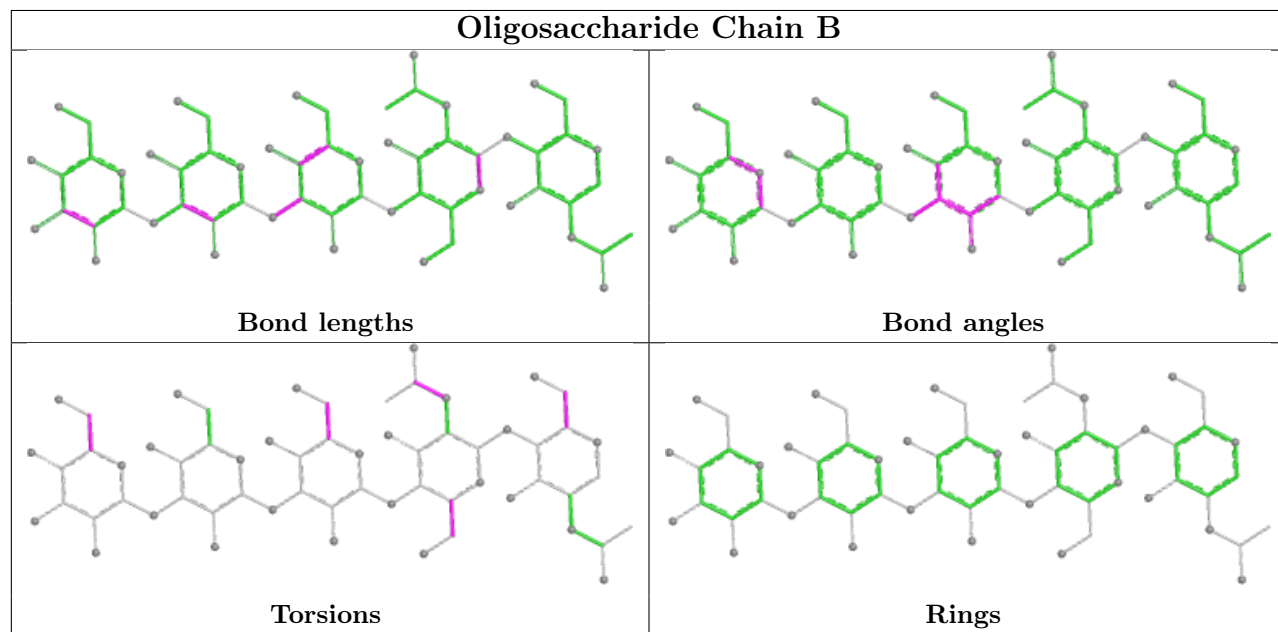
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	1	NAG	1	0
3	B	4	BMA	1	0
4	F	1	NAG	1	0
5	D	2	NAG	1	0
4	C	1	NAG	1	0

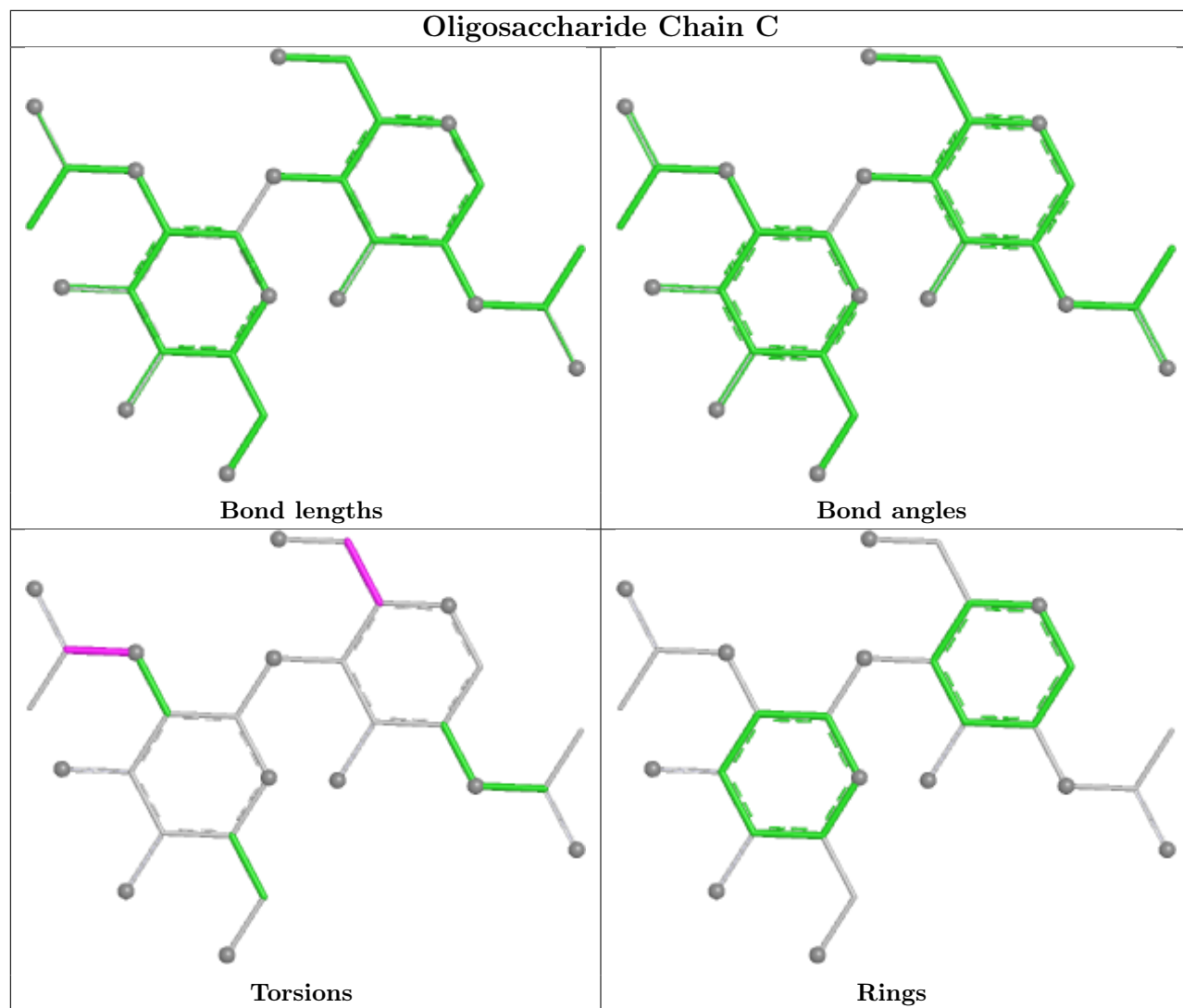
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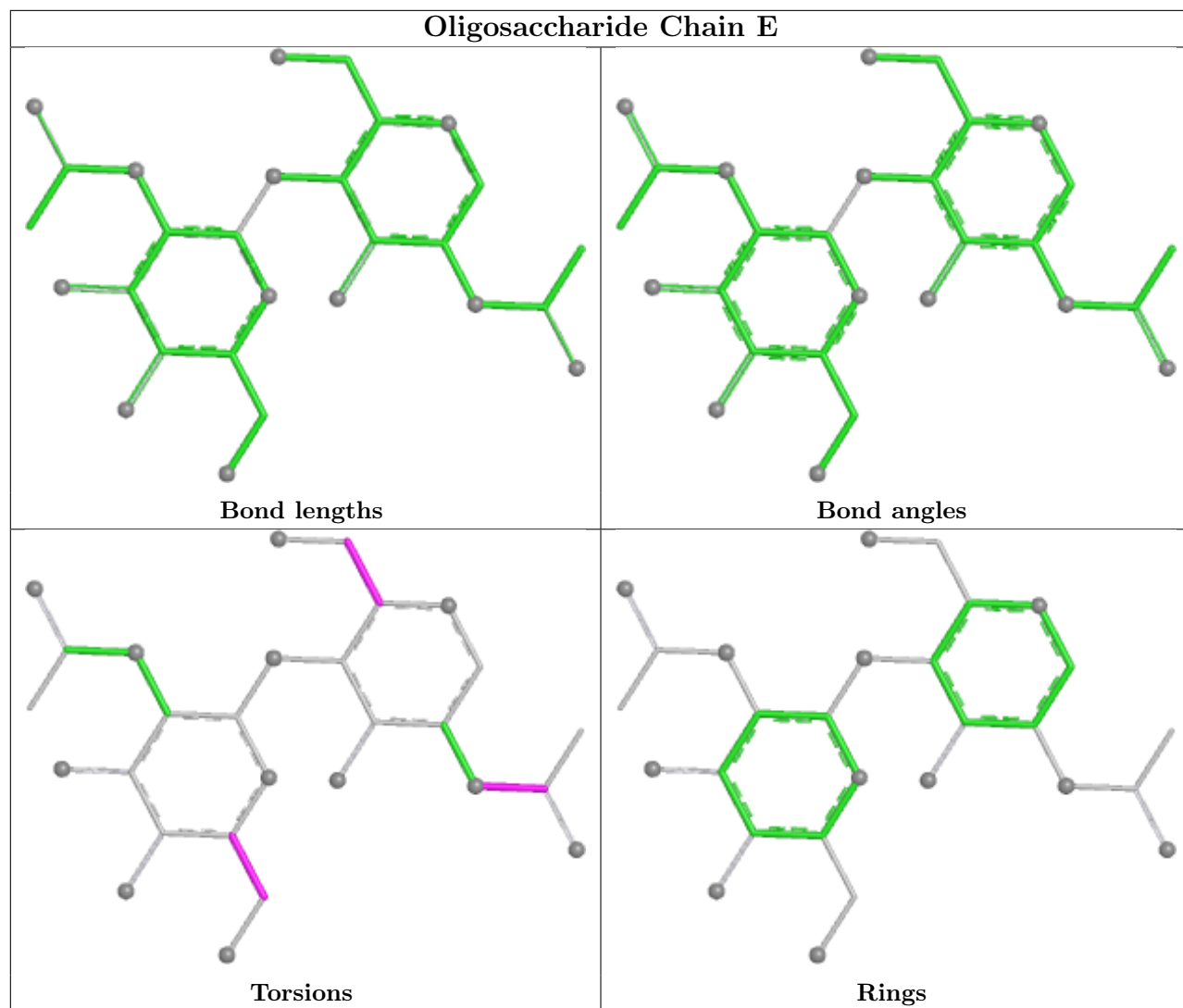
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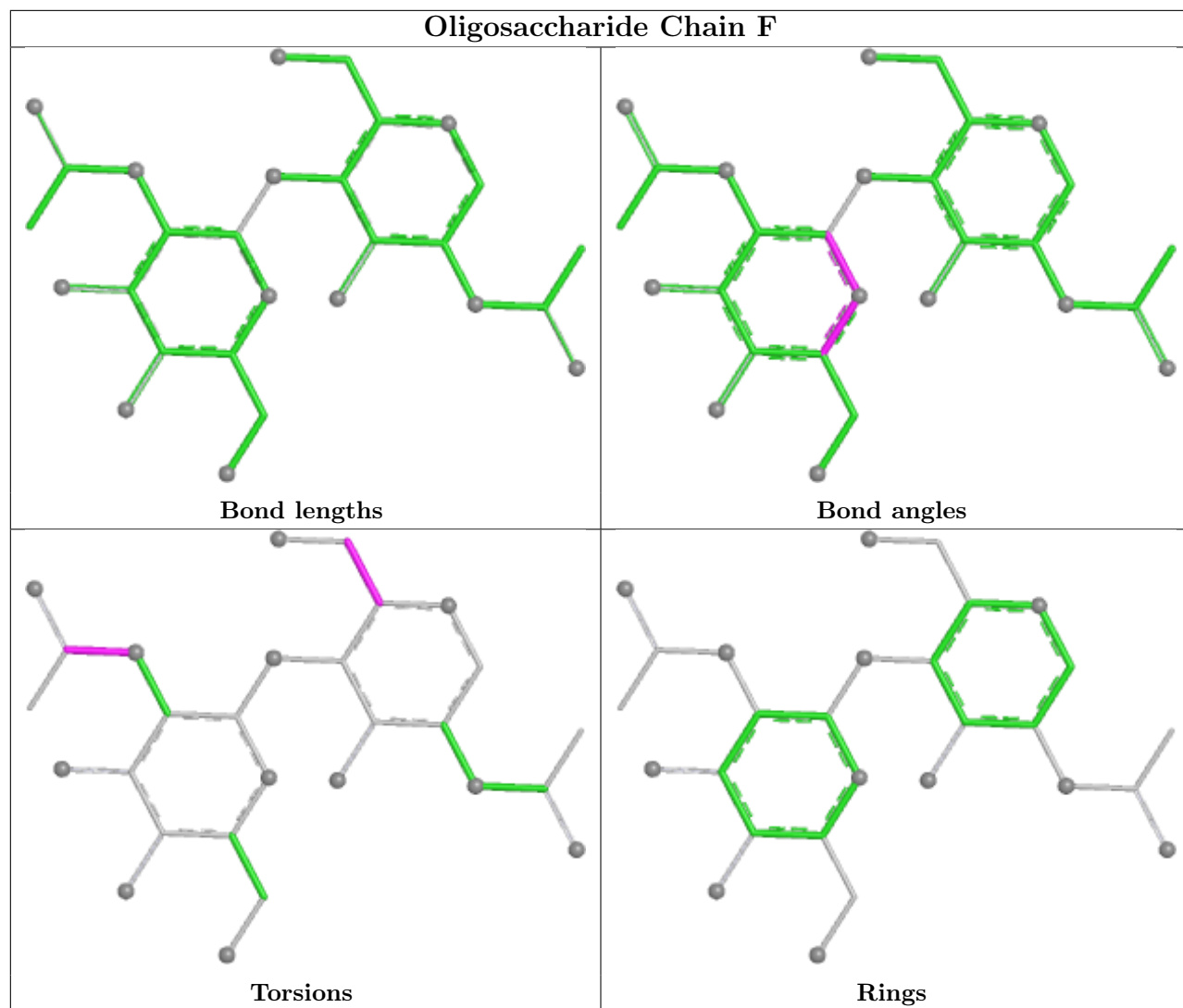
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	2	NAG	1	0

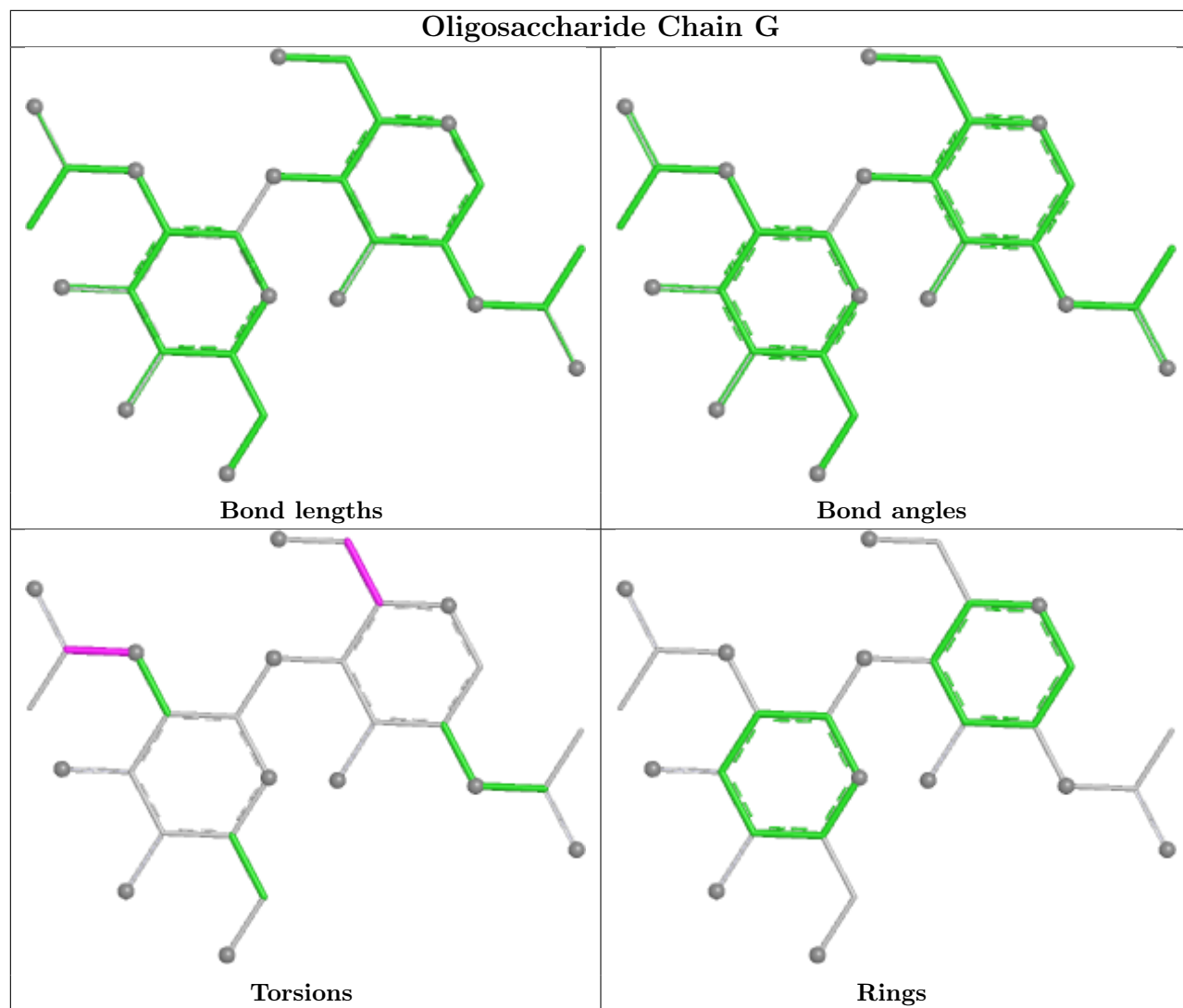
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

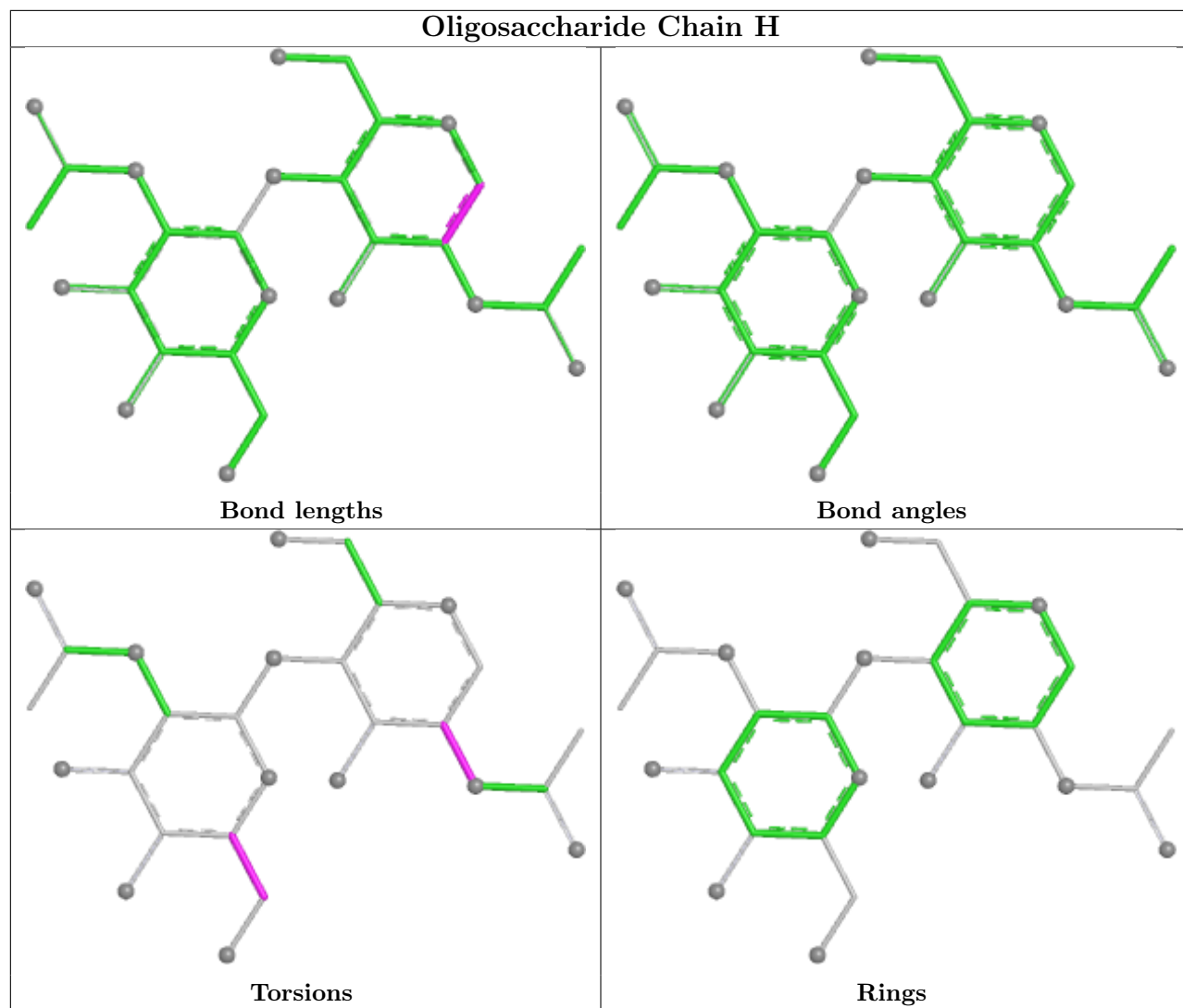


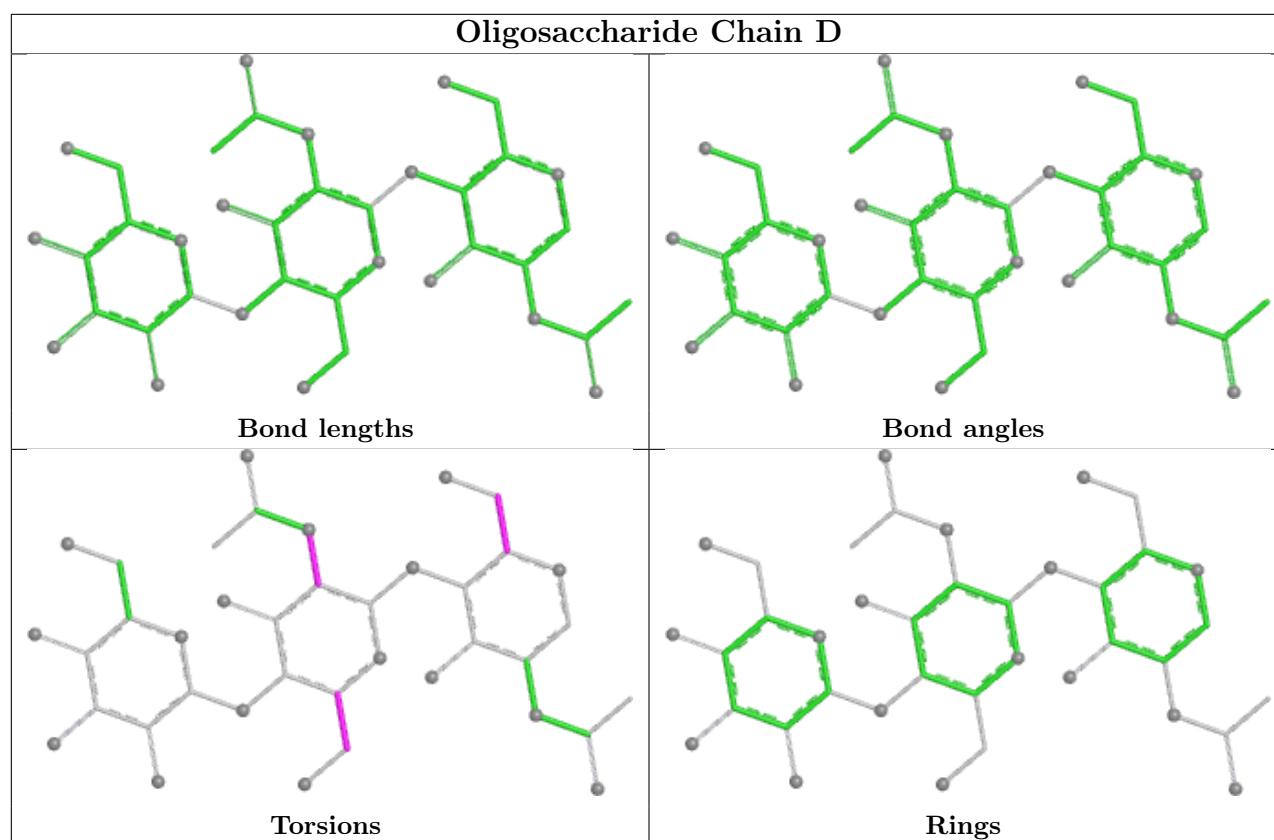












5.6 Ligand geometry [i](#)

6 ligands 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	$\# Z > 2$
6	NAG	A	2021	1	14,14,15	0.52	0	17,19,21	0.52	0
6	NAG	A	2022	1	14,14,15	1.41	2 (14%)	17,19,21	0.82	1 (5%)
7	EPE	A	2024	-	15,15,15	0.83	1 (6%)	19,20,20	1.76	4 (21%)
6	NAG	A	2023	1	14,14,15	0.39	0	17,19,21	0.49	0
6	NAG	A	2006	1	14,14,15	0.31	0	17,19,21	0.36	0
6	NAG	A	2007	1	14,14,15	0.29	0	17,19,21	0.36	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
6	NAG	A	2021	1	-	4/6/23/26	0/1/1/1
6	NAG	A	2022	1	-	3/6/23/26	0/1/1/1
7	EPE	A	2024	-	-	6/9/19/19	0/1/1/1
6	NAG	A	2023	1	-	3/6/23/26	0/1/1/1
6	NAG	A	2006	1	-	2/6/23/26	0/1/1/1
6	NAG	A	2007	1	-	1/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	2022	NAG	O5-C1	-4.51	1.36	1.43
7	A	2024	EPE	C10-S	2.84	1.81	1.77
6	A	2022	NAG	C1-C2	2.42	1.55	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	2024	EPE	C5-N4-C3	4.99	119.59	108.84
7	A	2024	EPE	C7-N4-C3	2.91	119.00	111.24
7	A	2024	EPE	C7-N4-C5	2.87	118.88	111.24
7	A	2024	EPE	O3S-S-C10	2.26	110.42	106.00
6	A	2022	NAG	C4-C3-C2	2.17	114.20	111.02

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	2024	EPE	C10-C9-N1-C6
6	A	2021	NAG	O5-C5-C6-O6
6	A	2023	NAG	O5-C5-C6-O6
6	A	2006	NAG	O5-C5-C6-O6
6	A	2022	NAG	C8-C7-N2-C2
6	A	2022	NAG	O7-C7-N2-C2
6	A	2021	NAG	C4-C5-C6-O6
6	A	2007	NAG	O5-C5-C6-O6
6	A	2023	NAG	C4-C5-C6-O6
7	A	2024	EPE	C9-C10-S-O3S
7	A	2024	EPE	C8-C7-N4-C3

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Mol	Chain	Res	Type	Atoms
7	A	2024	EPE	C9-C10-S-O2S
7	A	2024	EPE	S-C10-C9-N1
6	A	2006	NAG	C4-C5-C6-O6
6	A	2021	NAG	C3-C2-N2-C7
6	A	2021	NAG	C1-C2-N2-C7
6	A	2022	NAG	C1-C2-N2-C7
7	A	2024	EPE	C9-C10-S-O1S
6	A	2023	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	2024	EPE	1	0
6	A	2023	NAG	1	0

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	747/783 (95%)	-0.61	1 (0%) 92 90	4, 27, 58, 87	0
2	J	63/114 (55%)	-0.56	0 100 100	11, 33, 64, 71	0
2	K	65/114 (57%)	-0.52	0 100 100	10, 37, 60, 76	0
All	All	875/1011 (86%)	-0.60	1 (0%) 92 90	4, 28, 59, 87	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	660	ARG	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.

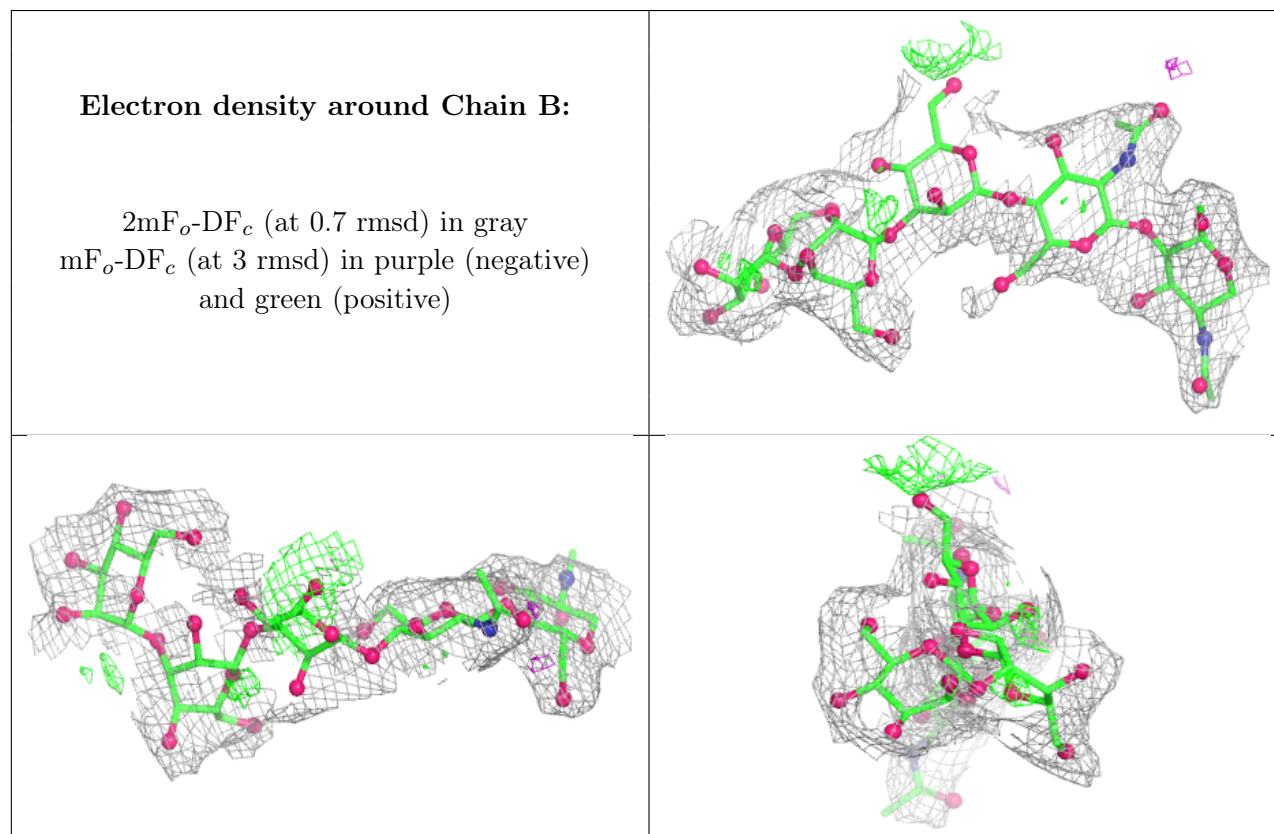
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	B	1	14/15	-	-	16,22,31,31	0
3	NAG	B	2	14/15	-	-	23,30,46,49	0
3	BMA	B	3	11/12	-	-	15,25,35,36	0
3	BMA	B	4	11/12	-	-	30,38,43,43	0
3	MAN	B	5	11/12	-	-	29,31,39,46	0
4	NAG	C	1	14/15	-	-	26,37,53,61	0
4	NAG	C	2	14/15	-	-	56,61,68,78	0

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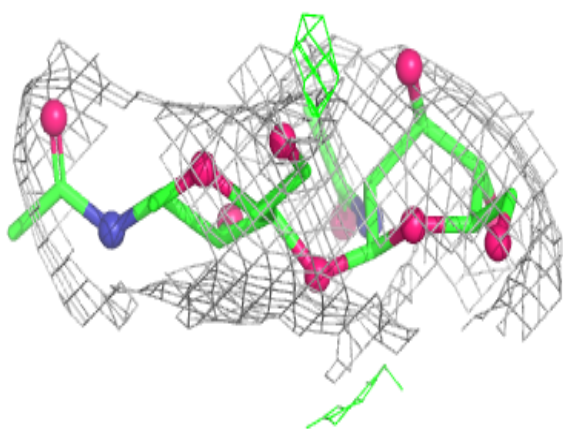
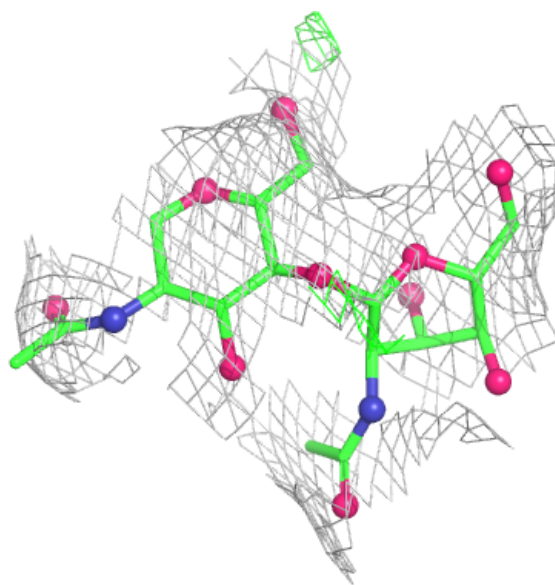
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	E	2	14/15	0.68	0.10	47,73,81,83	0
4	NAG	G	2	14/15	0.70	0.09	39,54,64,68	0
4	NAG	H	2	14/15	0.77	0.09	49,60,63,66	0
4	NAG	G	1	14/15	0.85	0.07	47,55,66,69	0
5	NAG	D	2	14/15	0.87	0.08	36,47,66,66	0
5	BMA	D	3	11/12	0.88	0.07	51,62,70,71	0
4	NAG	E	1	14/15	0.90	0.08	25,40,55,66	0
5	NAG	D	1	14/15	0.92	0.11	20,36,45,46	0
4	NAG	H	1	14/15	0.92	0.07	44,49,59,61	0
4	NAG	F	2	14/15	0.92	0.07	48,59,70,75	0
4	NAG	F	1	14/15	0.94	0.08	34,42,49,51	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.



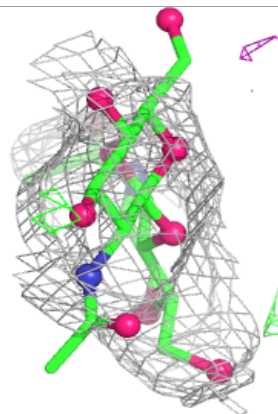
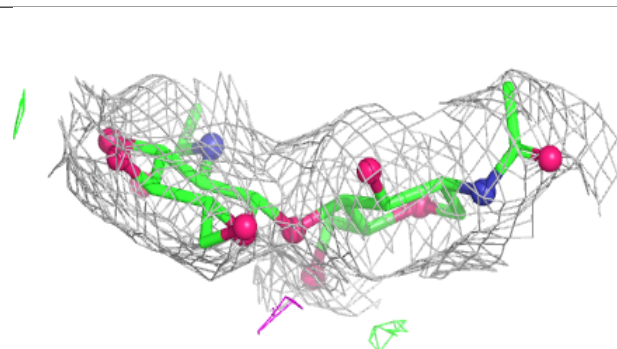
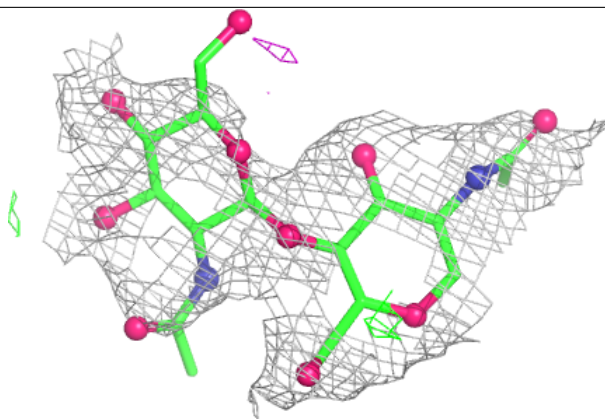
Electron density around Chain C:

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

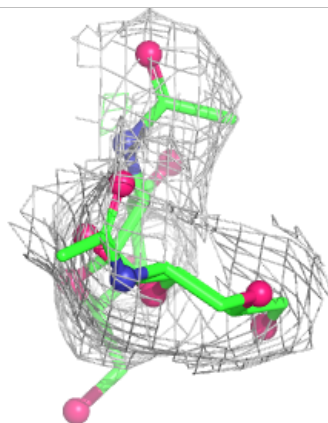
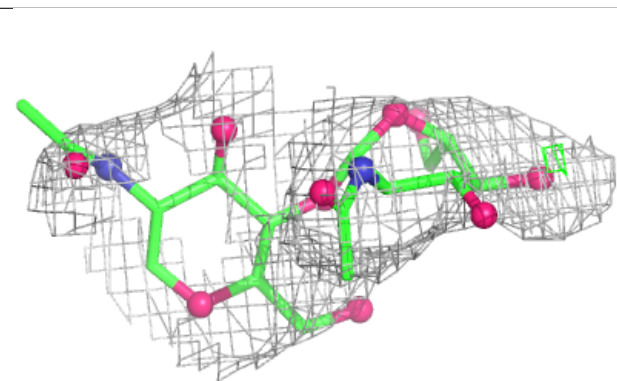
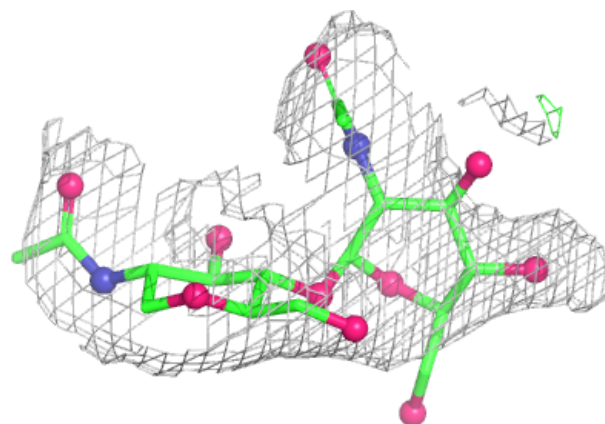


Electron density around Chain E:

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

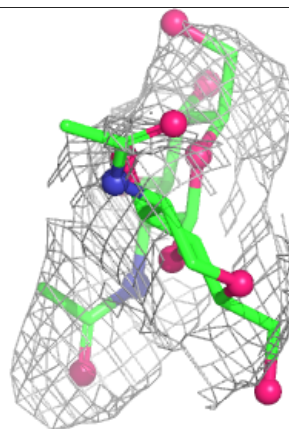
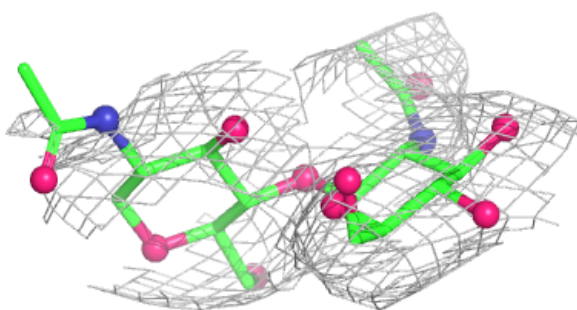
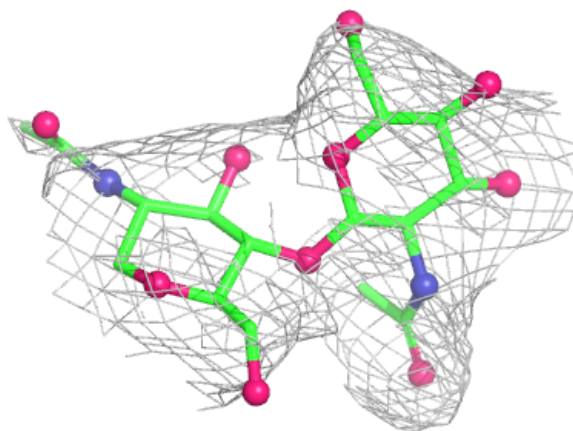
**Electron density around Chain F:**

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



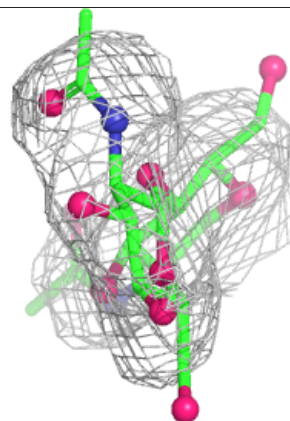
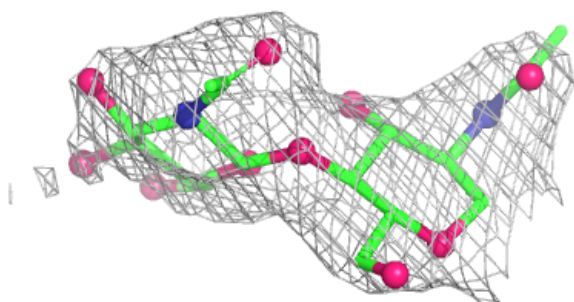
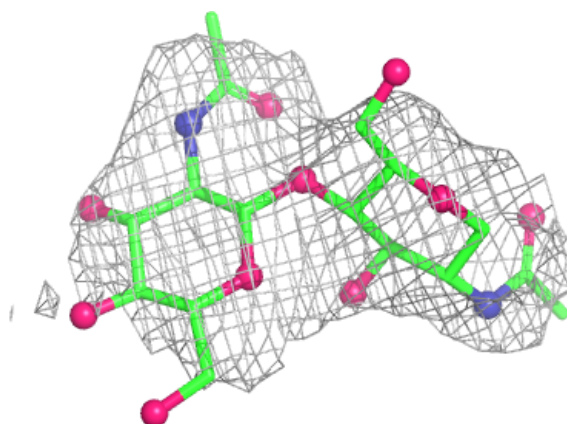
Electron density around Chain G:

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

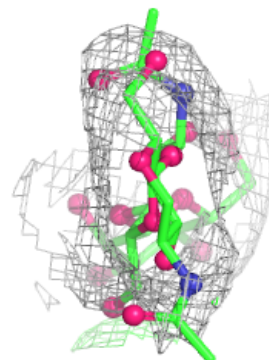
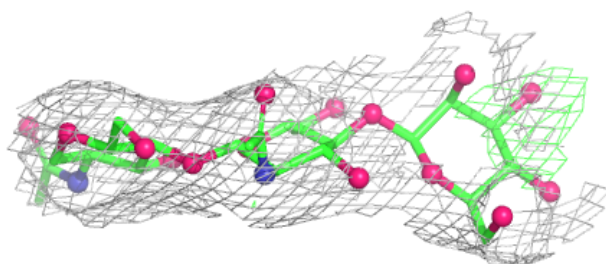
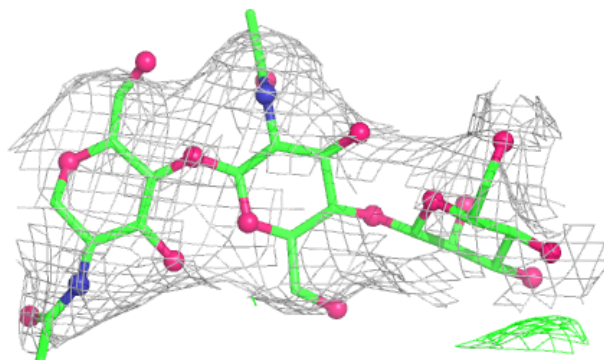


Electron density around Chain H:

$2mF_o-DF_c$ (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
6	NAG	A	2023	14/15	0.73	0.11	39,60,70,84	0
6	NAG	A	2021	14/15	0.77	0.11	55,65,80,83	0
6	NAG	A	2022	14/15	0.77	0.09	33,50,54,63	0
6	NAG	A	2006	14/15	0.77	0.08	36,59,66,75	0
7	EPE	A	2024	15/15	0.91	0.10	25,33,45,47	0
6	NAG	A	2007	14/15	0.95	0.05	28,33,43,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.