



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:15 AM UTC

PDB ID : 7CRF / pdb_00007crf
Title : Crystal structure of human TLR8 in complex with CU-CPD107
Authors : Sakaniwa, K.; Tanji, H.; Ohto, U.; Shimizu, T.
Deposited on : 2020-08-13
Resolution : 2.89 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:

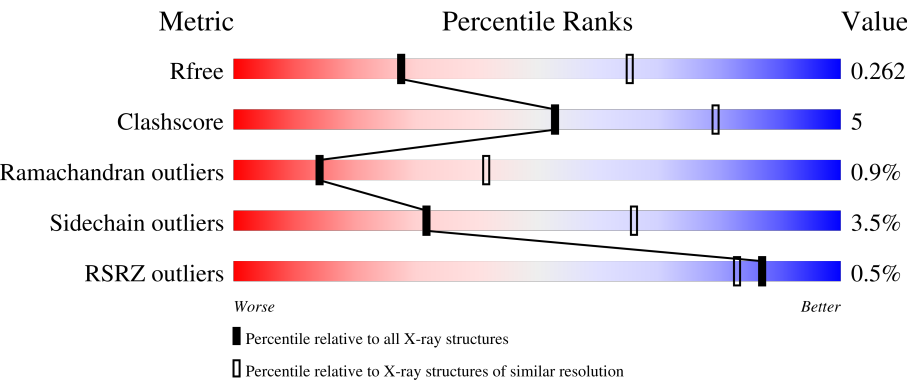
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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.



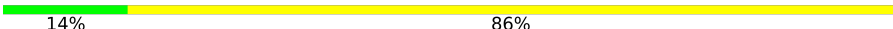
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	811	79%12%• 8%
1	B	811	78%13%• 7%
2	C	6	50%50%
3	D	3	100%
4	E	5	20%80%

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Mol	Chain	Length	Quality of chain
5	F	7	 71% 29%
6	G	4	 100%
7	H	7	 14% 86%

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 12776 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.

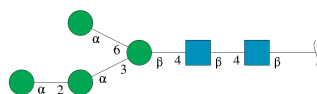
- Molecule 1 is a protein called Toll-like receptor 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	744	Total	C	N	O	S	0	0	0
			5998	3836	1019	1124	19			
1	B	751	Total	C	N	O	S	0	0	0
			6038	3859	1025	1134	20			

There are 20 discrepancies between the modelled and reference sequences:

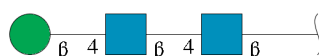
Chain	Residue	Modelled	Actual	Comment	Reference
A	23	ARG	-	expression tag	UNP Q9NR97
A	24	SER	-	expression tag	UNP Q9NR97
A	25	PRO	-	expression tag	UNP Q9NR97
A	26	TRP	-	expression tag	UNP Q9NR97
A	828	GLU	-	expression tag	UNP Q9NR97
A	829	PHE	-	expression tag	UNP Q9NR97
A	830	LEU	-	expression tag	UNP Q9NR97
A	831	VAL	-	expression tag	UNP Q9NR97
A	832	PRO	-	expression tag	UNP Q9NR97
A	833	ARG	-	expression tag	UNP Q9NR97
B	23	ARG	-	expression tag	UNP Q9NR97
B	24	SER	-	expression tag	UNP Q9NR97
B	25	PRO	-	expression tag	UNP Q9NR97
B	26	TRP	-	expression tag	UNP Q9NR97
B	828	GLU	-	expression tag	UNP Q9NR97
B	829	PHE	-	expression tag	UNP Q9NR97
B	830	LEU	-	expression tag	UNP Q9NR97
B	831	VAL	-	expression tag	UNP Q9NR97
B	832	PRO	-	expression tag	UNP Q9NR97
B	833	ARG	-	expression tag	UNP Q9NR97

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-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.



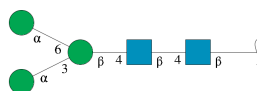
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	0	0
			72	40	2	30			

- 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.



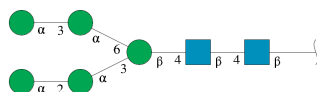
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	3	Total	C	N	O	0	0	0
			39	22	2	15			

- 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)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[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.



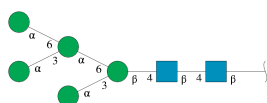
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	F	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 6 is an oligosaccharide called alpha-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.



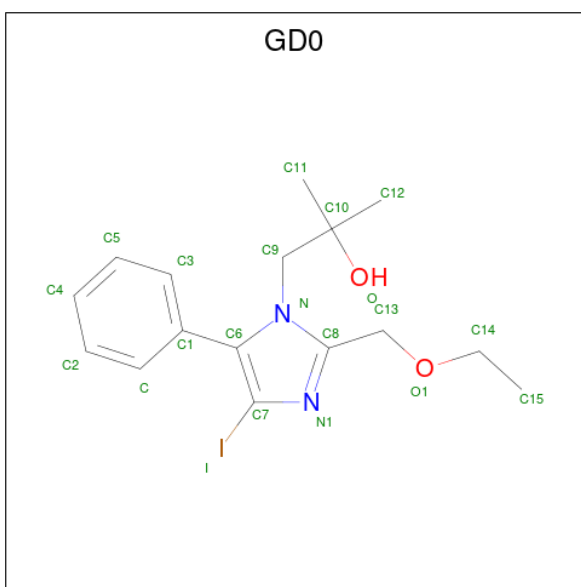
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	G	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-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.



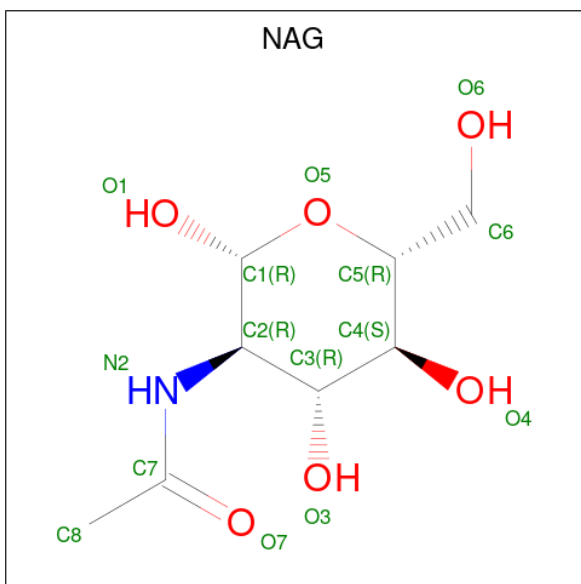
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	H	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 8 is 1-[2-(ethoxymethyl)-4-iodanyl-5-phenyl-imidazol-1-yl]-2-methyl-propan-2-ol (CCD ID: GD0) (formula: C₁₆H₂₁IN₂O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	I	N	O	0	0
			21	16	1	2	2		
8	B	1	Total	C	I	N	O	0	0
			21	16	1	2	2		

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

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

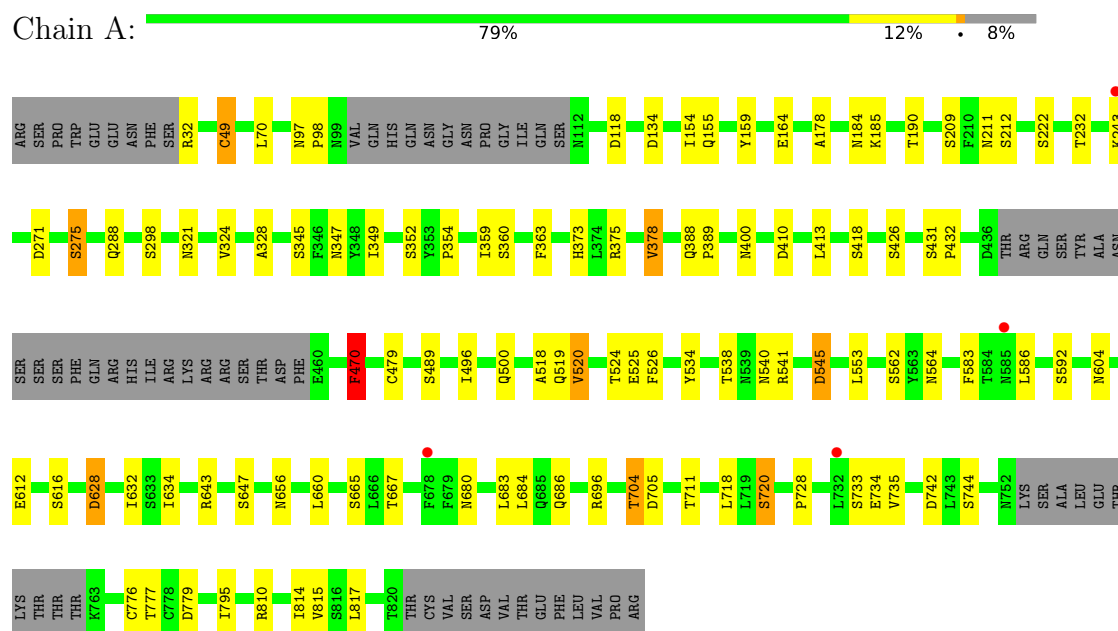
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	9	Total 9	O 9	0	0
10	B	7	Total 7	O 7	0	0

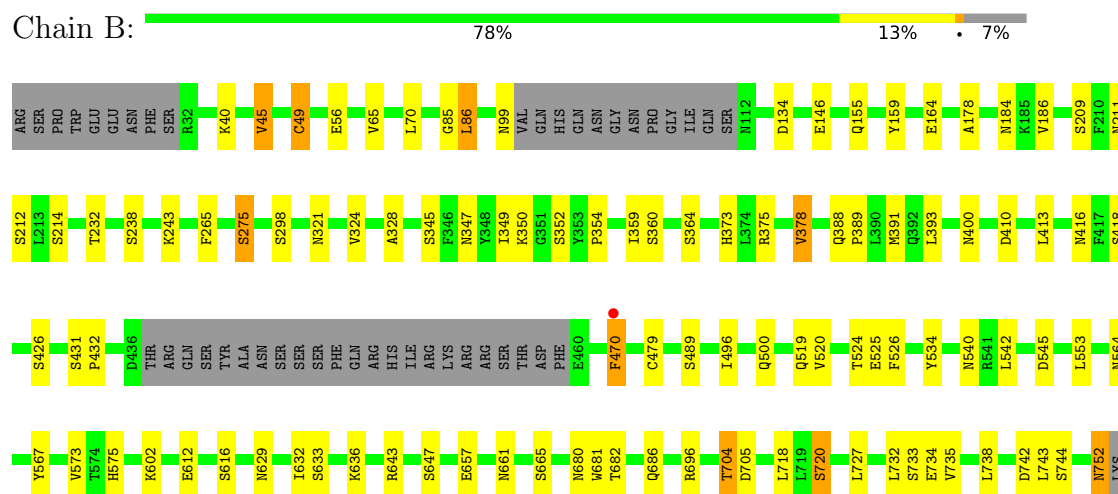
3 Residue-property plots

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: Toll-like receptor 8



• Molecule 1: Toll-like receptor 8



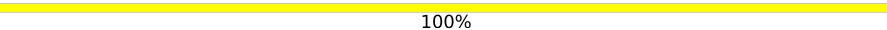


- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  50%

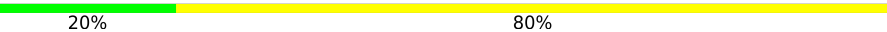


- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  100%

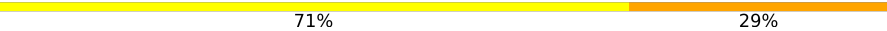


- Molecule 4: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  20%

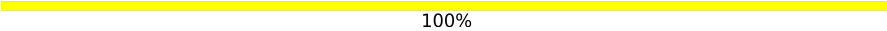


- Molecule 5: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  71%

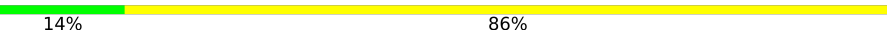


- Molecule 6: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  100%



- Molecule 7: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-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

Chain H:  14%

MAG1	MAG2	MAG3	MAG4	MAG5	MAG6	MAG7
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	144.34Å 99.97Å 140.92Å 90.00° 105.62° 90.00°	Depositor
Resolution (Å)	40.71 – 2.89 40.71 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.71-2.89) 99.8 (40.71-2.89)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.210 , 0.268 0.210 , 0.262	Depositor DCC
R_{free} test set	2216 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	75.2	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12776	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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: BMA, NAG, GD0, MAN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.02	0/6121	1.29	2/8300 (0.0%)
1	B	1.01	1/6161 (0.0%)	1.29	0/8358
All	All	1.01	1/12282 (0.0%)	1.29	2/16658 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	265	PHE	N-CA	5.19	1.50	1.45

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	545	ASP	CA-CB-CG	6.44	119.04	112.60
1	A	470	PHE	CA-CB-CG	6.04	119.84	113.80

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	5998	0	5975	61	0
1	B	6038	0	6005	65	0
2	C	72	0	61	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	39	0	34	0	0
4	E	61	0	52	0	0
5	F	83	0	70	1	0
6	G	50	0	43	0	0
7	H	83	0	70	0	0
8	A	21	0	0	6	0
8	B	21	0	0	3	0
9	A	154	0	143	0	0
9	B	140	0	130	0	0
10	A	9	0	0	0	0
10	B	7	0	0	0	0
All	All	12776	0	12583	126	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 5.

All (126) 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:660:LEU:HD21	1:A:683:LEU:HD22	1.46	0.97
1:B:45:VAL:HG13	1:B:65:VAL:HA	1.63	0.79
1:B:681:TRP:CG	1:B:704:THR:HG22	2.19	0.78
1:B:321:ASN:C	1:B:347:ASN:ND2	2.48	0.71
1:A:656:ASN:O	1:A:660:LEU:HD23	1.93	0.69
1:B:321:ASN:HB2	1:B:347:ASN:HD21	1.59	0.67
1:B:540:ASN:HB2	1:B:564:ASN:HD21	1.60	0.67
1:A:540:ASN:HB2	1:A:564:ASN:HD21	1.61	0.66
1:B:324:VAL:HG21	1:B:349:ILE:HG12	1.77	0.65
1:B:321:ASN:C	1:B:347:ASN:HD22	2.09	0.60
1:A:413:LEU:C	1:A:413:LEU:HD12	2.27	0.59
1:B:413:LEU:C	1:B:413:LEU:HD12	2.28	0.58
1:A:185:LYS:N	1:A:185:LYS:HD2	2.20	0.56
1:A:520:VAL:HG22	8:A:1001:GD0:N1	2.21	0.56
1:B:211:ASN:O	1:B:232:THR:HA	2.05	0.56
1:B:777:THR:O	1:B:779:ASP:N	2.39	0.56
8:A:1001:GD0:C4	1:B:378:VAL:HG11	2.37	0.55
1:A:211:ASN:O	1:A:232:THR:HA	2.06	0.55
1:B:777:THR:O	1:B:777:THR:OG1	2.26	0.54
1:B:720:SER:OG	1:B:742:ASP:CG	2.51	0.54
1:A:520:VAL:CG2	8:A:1001:GD0:N1	2.71	0.54
1:B:752:ASN:OD1	1:B:752:ASN:C	2.50	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:567:TYR:O	1:B:575:HIS:HE1	1.91	0.53
1:A:656:ASN:ND2	1:A:680:ASN:HD22	2.06	0.53
1:A:159:TYR:OH	1:A:184:ASN:ND2	2.43	0.52
1:A:520:VAL:HG21	1:B:350:LYS:HE3	1.91	0.52
1:A:519:GLN:HA	8:A:1001:GD0:C13	2.39	0.52
1:B:159:TYR:OH	1:B:184:ASN:ND2	2.42	0.52
1:A:720:SER:OG	1:A:742:ASP:CG	2.53	0.52
1:B:321:ASN:O	1:B:347:ASN:ND2	2.43	0.52
1:A:718:LEU:HA	1:A:742:ASP:HB3	1.92	0.51
1:B:815:VAL:O	1:B:815:VAL:HG12	2.11	0.51
1:A:815:VAL:HG12	1:A:815:VAL:O	2.11	0.51
1:B:804:ALA:O	1:B:810:ARG:NH2	2.42	0.51
1:B:799:VAL:O	1:B:802:ILE:HD11	2.10	0.50
1:A:134:ASP:HA	1:A:155:GLN:O	2.11	0.50
1:A:616:SER:HA	1:A:647:SER:O	2.11	0.50
1:B:134:ASP:HA	1:B:155:GLN:O	2.10	0.50
1:A:777:THR:O	1:A:779:ASP:N	2.41	0.50
1:B:391:MET:SD	1:B:416:ASN:HB3	2.52	0.49
1:A:328:ALA:O	1:A:360:SER:HB2	2.13	0.49
1:A:178:ALA:HA	1:A:209:SER:O	2.13	0.49
1:B:616:SER:HA	1:B:647:SER:O	2.13	0.49
1:B:718:LEU:HA	1:B:742:ASP:HB3	1.95	0.49
1:B:733:SER:O	1:B:734:GLU:HG2	2.12	0.49
1:A:526:PHE:HB3	1:A:553:LEU:HD21	1.94	0.48
1:B:720:SER:OG	1:B:742:ASP:OD2	2.31	0.48
1:B:178:ALA:HA	1:B:209:SER:O	2.13	0.48
1:B:496:ILE:HG21	1:B:525:GLU:HB3	1.96	0.48
1:A:720:SER:HA	1:A:744:SER:O	2.14	0.48
1:B:328:ALA:O	1:B:360:SER:HB2	2.14	0.47
1:B:776:CYS:HB3	1:B:817:LEU:HD21	1.96	0.47
1:A:496:ILE:HG21	1:A:525:GLU:HB3	1.96	0.47
1:B:352:SER:OG	1:B:354:PRO:HD3	2.14	0.47
1:A:776:CYS:HB3	1:A:817:LEU:HD21	1.96	0.47
1:A:720:SER:OG	1:A:742:ASP:OD2	2.32	0.47
1:B:345:SER:HA	1:B:375:ARG:O	2.15	0.47
1:B:567:TYR:O	1:B:575:HIS:CE1	2.68	0.47
1:A:352:SER:OG	1:A:354:PRO:HD3	2.15	0.46
1:B:388:GLN:N	1:B:389:PRO:CD	2.78	0.46
1:A:32:ARG:NH1	1:A:795:ILE:HD12	2.31	0.46
1:A:345:SER:HA	1:A:375:ARG:O	2.15	0.46
1:A:520:VAL:HG22	8:A:1001:GD0:C8	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:388:GLN:N	1:A:389:PRO:CD	2.78	0.46
1:B:720:SER:HA	1:B:744:SER:O	2.15	0.46
1:A:373:HIS:HA	1:A:400:ASN:HB3	1.98	0.46
1:A:520:VAL:HG22	8:A:1001:GD0:C13	2.46	0.46
1:A:656:ASN:HD22	1:A:683:LEU:HD11	1.80	0.46
1:A:378:VAL:HG11	8:B:1001:GD0:C4	2.47	0.45
1:A:684:LEU:HB3	1:A:711:THR:HG23	1.98	0.45
1:B:391:MET:SD	1:B:416:ASN:CB	3.04	0.45
1:A:359:ILE:HG23	1:A:363:PHE:CD2	2.51	0.45
1:B:526:PHE:HB3	1:B:553:LEU:HD21	1.98	0.45
1:B:542:LEU:H	1:B:564:ASN:HD22	1.65	0.45
1:B:798:LEU:HD12	1:B:798:LEU:HA	1.86	0.45
1:B:364:SER:HA	1:B:393:LEU:HD21	1.99	0.44
1:B:479:CYS:SG	1:B:534:TYR:HB3	2.57	0.44
1:A:583:PHE:HB3	1:A:586:LEU:HB2	2.00	0.44
1:A:660:LEU:CD2	1:A:683:LEU:HD13	2.47	0.44
1:B:681:TRP:CD2	1:B:704:THR:HG22	2.53	0.44
1:A:628:ASP:N	1:A:628:ASP:OD1	2.48	0.44
1:B:359:ILE:O	1:B:359:ILE:HG13	2.18	0.44
1:B:520:VAL:HG13	1:B:545:ASP:OD1	2.18	0.44
1:A:154:ILE:HG12	1:A:178:ALA:HB3	2.00	0.43
1:A:733:SER:O	1:A:734:GLU:HB2	2.19	0.43
1:B:373:HIS:HA	1:B:400:ASN:HB3	1.99	0.43
1:B:321:ASN:CB	1:B:347:ASN:HD21	2.30	0.43
1:A:705:ASP:CB	1:A:728:PRO:HB3	2.49	0.43
1:A:324:VAL:HG21	1:A:349:ILE:HG13	2.01	0.43
1:B:184:ASN:ND2	1:B:186:VAL:HB	2.34	0.43
5:F:4:MAN:HO3	5:F:5:MAN:C1	2.25	0.42
1:A:704:THR:O	1:A:705:ASP:CG	2.62	0.42
1:B:612:GLU:HG3	1:B:643:ARG:CG	2.49	0.42
1:B:636:LYS:HE3	1:B:661:ASN:HA	2.00	0.42
1:A:777:THR:O	1:A:777:THR:OG1	2.27	0.42
1:B:49:CYS:HB2	1:B:70:LEU:HD23	2.01	0.42
1:B:519:GLN:HA	8:B:1001:GD0:C13	2.50	0.42
1:A:288:GLN:OE1	1:A:288:GLN:HA	2.20	0.42
1:A:612:GLU:OE2	1:A:643:ARG:NH1	2.53	0.42
1:A:520:VAL:HG12	1:A:545:ASP:HB3	2.02	0.42
1:A:720:SER:HG	1:A:742:ASP:CG	2.27	0.42
1:B:432:PRO:HA	1:B:500:GLN:OE1	2.20	0.42
1:A:426:SER:HA	1:A:489:SER:O	2.20	0.41
1:A:518:ALA:HB2	1:A:541:ARG:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:592:SER:HA	1:A:616:SER:O	2.20	0.41
1:B:704:THR:O	1:B:705:ASP:CG	2.63	0.41
1:A:432:PRO:HA	1:A:500:GLN:OE1	2.21	0.41
1:A:696:ARG:HA	1:A:720:SER:O	2.20	0.41
1:B:520:VAL:HG23	8:B:1001:GD0:C14	2.51	0.41
1:A:479:CYS:SG	1:A:534:TYR:HB3	2.60	0.41
1:B:629:ASN:HB3	1:B:632:ILE:HG13	2.03	0.41
1:B:732:LEU:C	1:B:732:LEU:HD23	2.46	0.41
1:B:727:LEU:HD11	1:B:743:LEU:HD13	2.02	0.41
1:A:97:ASN:HA	1:A:98:PRO:HA	1.87	0.41
1:A:275:SER:HA	1:A:298:SER:HB2	2.03	0.41
1:B:275:SER:HA	1:B:298:SER:HB2	2.03	0.41
1:B:632:ILE:HG22	1:B:661:ASN:ND2	2.36	0.41
1:B:426:SER:HA	1:B:489:SER:O	2.21	0.40
1:A:604:ASN:HD22	1:A:634:ILE:HA	1.85	0.40
1:A:49:CYS:HB2	1:A:70:LEU:HD23	2.03	0.40
1:A:321:ASN:O	1:A:347:ASN:HA	2.22	0.40
1:B:85:GLY:O	1:B:86:LEU:C	2.64	0.40
1:B:680:ASN:OD1	1:B:682:THR:OG1	2.28	0.40
1:B:696:ARG:HA	1:B:720:SER:O	2.22	0.40
1:A:538:THR:HA	1:A:562:SER:O	2.22	0.40
1:B:540:ASN:CB	1:B:564:ASN:HD21	2.32	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	736/811 (91%)	667 (91%)	64 (9%)	5 (1%)	18	48
1	B	743/811 (92%)	673 (91%)	62 (8%)	8 (1%)	11	36
All	All	1479/1622 (91%)	1340 (91%)	126 (8%)	13 (1%)	14	41

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	810	ARG
1	B	810	ARG
1	B	86	LEU
1	A	470	PHE
1	B	410	ASP
1	B	470	PHE
1	A	410	ASP
1	A	378	VAL
1	B	378	VAL
1	B	214	SER
1	B	735	VAL
1	A	735	VAL
1	B	826	VAL

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	692/755 (92%)	670 (97%)	22 (3%)	34	68
1	B	697/755 (92%)	670 (96%)	27 (4%)	28	63
All	All	1389/1510 (92%)	1340 (96%)	49 (4%)	32	66

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	CYS
1	A	118	ASP
1	A	164	GLU
1	A	190	THR
1	A	212	SER
1	A	222	SER
1	A	243	LYS
1	A	271	ASP
1	A	275	SER
1	A	418	SER

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Mol	Chain	Res	Type
1	A	431	SER
1	A	470	PHE
1	A	520	VAL
1	A	524	THR
1	A	628	ASP
1	A	632	ILE
1	A	665	SER
1	A	667	THR
1	A	686	GLN
1	A	704	THR
1	A	720	SER
1	A	814	ILE
1	B	40	LYS
1	B	45	VAL
1	B	49	CYS
1	B	56	GLU
1	B	99	ASN
1	B	146	GLU
1	B	164	GLU
1	B	212	SER
1	B	238	SER
1	B	243	LYS
1	B	275	SER
1	B	418	SER
1	B	431	SER
1	B	470	PHE
1	B	524	THR
1	B	573	VAL
1	B	602	LYS
1	B	633	SER
1	B	657	GLU
1	B	665	SER
1	B	686	GLN
1	B	704	THR
1	B	720	SER
1	B	738	LEU
1	B	752	ASN
1	B	782	ASP
1	B	798	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	96	HIS
1	A	123	ASN
1	A	136	GLN
1	A	155	GLN
1	A	184	ASN
1	A	233	GLN
1	A	277	ASN
1	A	309	ASN
1	A	564	ASN
1	A	593	HIS
1	A	604	ASN
1	A	629	ASN
1	A	656	ASN
1	A	661	ASN
1	A	790	HIS
1	B	77	HIS
1	B	96	HIS
1	B	99	ASN
1	B	136	GLN
1	B	184	ASN
1	B	233	GLN
1	B	309	ASN
1	B	347	ASN
1	B	564	ASN
1	B	575	HIS
1	B	581	GLN
1	B	593	HIS
1	B	604	ASN
1	B	790	HIS
1	B	809	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 ⓘ

32 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	0.56	0	17,19,21	0.96	0
2	NAG	C	2	2	14,14,15	0.64	0	17,19,21	1.05	1 (5%)
2	BMA	C	3	2	11,11,12	0.41	0	15,15,17	1.69	4 (26%)
2	MAN	C	4	2	11,11,12	0.32	0	15,15,17	0.85	0
2	MAN	C	5	2	11,11,12	0.24	0	15,15,17	0.75	0
2	MAN	C	6	2	11,11,12	0.60	0	15,15,17	1.07	1 (6%)
3	NAG	D	1	3,1	14,14,15	0.46	0	17,19,21	1.65	3 (17%)
3	NAG	D	2	3	14,14,15	0.69	1 (7%)	17,19,21	1.22	2 (11%)
3	BMA	D	3	3	11,11,12	0.44	0	15,15,17	1.06	1 (6%)
4	NAG	E	1	4,1	14,14,15	0.42	0	17,19,21	1.36	3 (17%)
4	NAG	E	2	4	14,14,15	0.70	0	17,19,21	1.29	2 (11%)
4	BMA	E	3	4	11,11,12	0.60	0	15,15,17	1.69	3 (20%)
4	MAN	E	4	4	11,11,12	0.43	0	15,15,17	1.07	1 (6%)
4	MAN	E	5	4	11,11,12	0.64	0	15,15,17	0.68	0
5	NAG	F	1	1,5	14,14,15	0.45	0	17,19,21	1.48	2 (11%)
5	NAG	F	2	5	14,14,15	0.73	0	17,19,21	1.21	1 (5%)
5	BMA	F	3	5	11,11,12	0.36	0	15,15,17	1.73	4 (26%)
5	MAN	F	4	5	11,11,12	0.61	0	15,15,17	1.30	1 (6%)
5	MAN	F	5	5	11,11,12	0.46	0	15,15,17	1.56	2 (13%)
5	MAN	F	6	5	11,11,12	0.54	0	15,15,17	1.11	1 (6%)
5	MAN	F	7	5	11,11,12	0.86	0	15,15,17	1.49	2 (13%)
6	NAG	G	1	6,1	14,14,15	0.50	0	17,19,21	1.15	1 (5%)
6	NAG	G	2	6	14,14,15	0.47	0	17,19,21	1.15	2 (11%)
6	BMA	G	3	6	11,11,12	0.92	0	15,15,17	1.73	3 (20%)
6	MAN	G	4	6	11,11,12	0.54	0	15,15,17	1.15	2 (13%)
7	NAG	H	1	7,1	14,14,15	0.54	0	17,19,21	1.41	3 (17%)
7	NAG	H	2	7	14,14,15	0.66	0	17,19,21	1.10	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	BMA	H	3	7	11,11,12	0.48	0	15,15,17	1.13	1 (6%)
7	MAN	H	4	7	11,11,12	0.89	0	15,15,17	0.92	0
7	MAN	H	5	7	11,11,12	0.51	0	15,15,17	2.14	2 (13%)
7	MAN	H	6	7	11,11,12	0.52	0	15,15,17	1.11	1 (6%)
7	MAN	H	7	7	11,11,12	0.51	0	15,15,17	1.17	1 (6%)

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	-	0/2/19/22	0/1/1/1
2	MAN	C	5	2	-	1/2/19/22	0/1/1/1
2	MAN	C	6	2	-	0/2/19/22	0/1/1/1
3	NAG	D	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
3	BMA	D	3	3	-	2/2/19/22	0/1/1/1
4	NAG	E	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	0/2/19/22	0/1/1/1
4	MAN	E	4	4	-	2/2/19/22	0/1/1/1
4	MAN	E	5	4	-	0/2/19/22	0/1/1/1
5	NAG	F	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	F	2	5	-	0/6/23/26	0/1/1/1
5	BMA	F	3	5	-	0/2/19/22	0/1/1/1
5	MAN	F	4	5	-	2/2/19/22	0/1/1/1
5	MAN	F	5	5	-	2/2/19/22	0/1/1/1
5	MAN	F	6	5	-	1/2/19/22	0/1/1/1
5	MAN	F	7	5	-	0/2/19/22	1/1/1/1
6	NAG	G	1	6,1	-	2/6/23/26	0/1/1/1
6	NAG	G	2	6	-	1/6/23/26	0/1/1/1
6	BMA	G	3	6	-	1/2/19/22	0/1/1/1
6	MAN	G	4	6	-	1/2/19/22	0/1/1/1
7	NAG	H	1	7,1	-	2/6/23/26	0/1/1/1
7	NAG	H	2	7	-	2/6/23/26	0/1/1/1
7	BMA	H	3	7	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	MAN	H	4	7	-	2/2/19/22	0/1/1/1
7	MAN	H	5	7	-	2/2/19/22	0/1/1/1
7	MAN	H	6	7	-	2/2/19/22	0/1/1/1
7	MAN	H	7	7	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	2	NAG	O4-C4	2.02	1.48	1.43

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	H	5	MAN	C1-O5-C5	6.67	121.12	112.19
5	F	5	MAN	C1-O5-C5	4.89	118.75	112.19
6	G	3	BMA	C1-C2-C3	4.43	116.10	109.64
3	D	1	NAG	C1-O5-C5	4.38	118.05	112.19
5	F	1	NAG	C1-O5-C5	4.09	117.67	112.19
5	F	7	MAN	C1-O5-C5	4.03	117.59	112.19
6	G	1	NAG	C1-C2-N2	-3.95	104.21	110.43
4	E	3	BMA	O3-C3-C4	3.91	119.59	110.38
5	F	3	BMA	C1-O5-C5	3.76	117.23	112.19
7	H	5	MAN	O5-C1-C2	3.72	119.66	110.79
5	F	3	BMA	O5-C1-C2	3.67	119.54	110.79
5	F	1	NAG	O5-C1-C2	-3.65	105.64	111.29
2	C	3	BMA	O3-C3-C4	-3.48	102.16	110.38
6	G	3	BMA	O5-C1-C2	3.45	119.02	110.79
4	E	3	BMA	C1-O5-C5	3.22	116.50	112.19
7	H	6	MAN	C1-O5-C5	3.22	116.50	112.19
2	C	3	BMA	C3-C4-C5	3.12	115.88	110.23
4	E	2	NAG	O5-C1-C2	3.02	115.97	111.29
6	G	4	MAN	O5-C1-C2	2.97	117.87	110.79
5	F	5	MAN	O5-C1-C2	2.97	117.87	110.79
7	H	1	NAG	O4-C4-C3	-2.96	103.40	110.38
4	E	3	BMA	O5-C1-C2	2.89	117.67	110.79
7	H	2	NAG	C1-C2-N2	2.81	114.87	110.43
7	H	3	BMA	O5-C1-C2	2.79	117.44	110.79
4	E	4	MAN	C1-O5-C5	2.79	115.92	112.19
4	E	1	NAG	O5-C1-C2	-2.75	107.04	111.29
2	C	3	BMA	O4-C4-C3	-2.62	104.20	110.38
3	D	2	NAG	O5-C1-C2	2.62	115.34	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	7	MAN	O5-C1-C2	2.61	117.03	110.79
7	H	7	MAN	C3-C4-C5	2.61	114.97	110.23
6	G	4	MAN	C1-O5-C5	2.61	115.69	112.19
3	D	1	NAG	C1-C2-N2	-2.52	106.47	110.43
7	H	1	NAG	O5-C5-C6	2.48	112.48	107.66
5	F	4	MAN	C2-C3-C4	2.47	115.21	110.86
3	D	1	NAG	O7-C7-N2	2.46	126.33	121.98
3	D	3	BMA	C3-C4-C5	2.40	114.59	110.23
7	H	1	NAG	O3-C3-C4	-2.40	104.72	110.38
2	C	6	MAN	C3-C4-C5	2.37	114.53	110.23
6	G	2	NAG	C1-O5-C5	2.36	115.35	112.19
6	G	2	NAG	O5-C1-C2	2.35	114.93	111.29
5	F	2	NAG	C1-C2-N2	2.29	114.05	110.43
5	F	3	BMA	O3-C3-C4	-2.29	104.99	110.38
4	E	1	NAG	O4-C4-C3	-2.23	105.11	110.38
6	G	3	BMA	C2-C3-C4	2.22	114.77	110.86
2	C	3	BMA	C2-C3-C4	2.16	114.67	110.86
5	F	6	MAN	O5-C1-C2	2.15	115.93	110.79
4	E	2	NAG	C1-C2-N2	2.14	113.81	110.43
2	C	2	NAG	C1-C2-N2	2.12	113.78	110.43
4	E	1	NAG	O6-C6-C5	-2.12	104.12	111.33
5	F	3	BMA	C1-C2-C3	2.09	112.69	109.64
3	D	2	NAG	O7-C7-N2	2.03	125.56	121.98

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	H	6	MAN	O5-C5-C6-O6
5	F	5	MAN	O5-C5-C6-O6
7	H	4	MAN	O5-C5-C6-O6
3	D	3	BMA	C4-C5-C6-O6
7	H	6	MAN	C4-C5-C6-O6
7	H	5	MAN	O5-C5-C6-O6
5	F	4	MAN	O5-C5-C6-O6
5	F	5	MAN	C4-C5-C6-O6
5	F	4	MAN	C4-C5-C6-O6
7	H	1	NAG	C4-C5-C6-O6
4	E	4	MAN	C4-C5-C6-O6
7	H	4	MAN	C4-C5-C6-O6
3	D	3	BMA	O5-C5-C6-O6
7	H	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	H	2	NAG	O5-C5-C6-O6
7	H	5	MAN	C4-C5-C6-O6
4	E	4	MAN	O5-C5-C6-O6
6	G	2	NAG	O5-C5-C6-O6
6	G	4	MAN	O5-C5-C6-O6
6	G	3	BMA	O5-C5-C6-O6
6	G	1	NAG	C4-C5-C6-O6
2	C	5	MAN	O5-C5-C6-O6
7	H	2	NAG	C4-C5-C6-O6
5	F	6	MAN	C4-C5-C6-O6
6	G	1	NAG	O5-C5-C6-O6

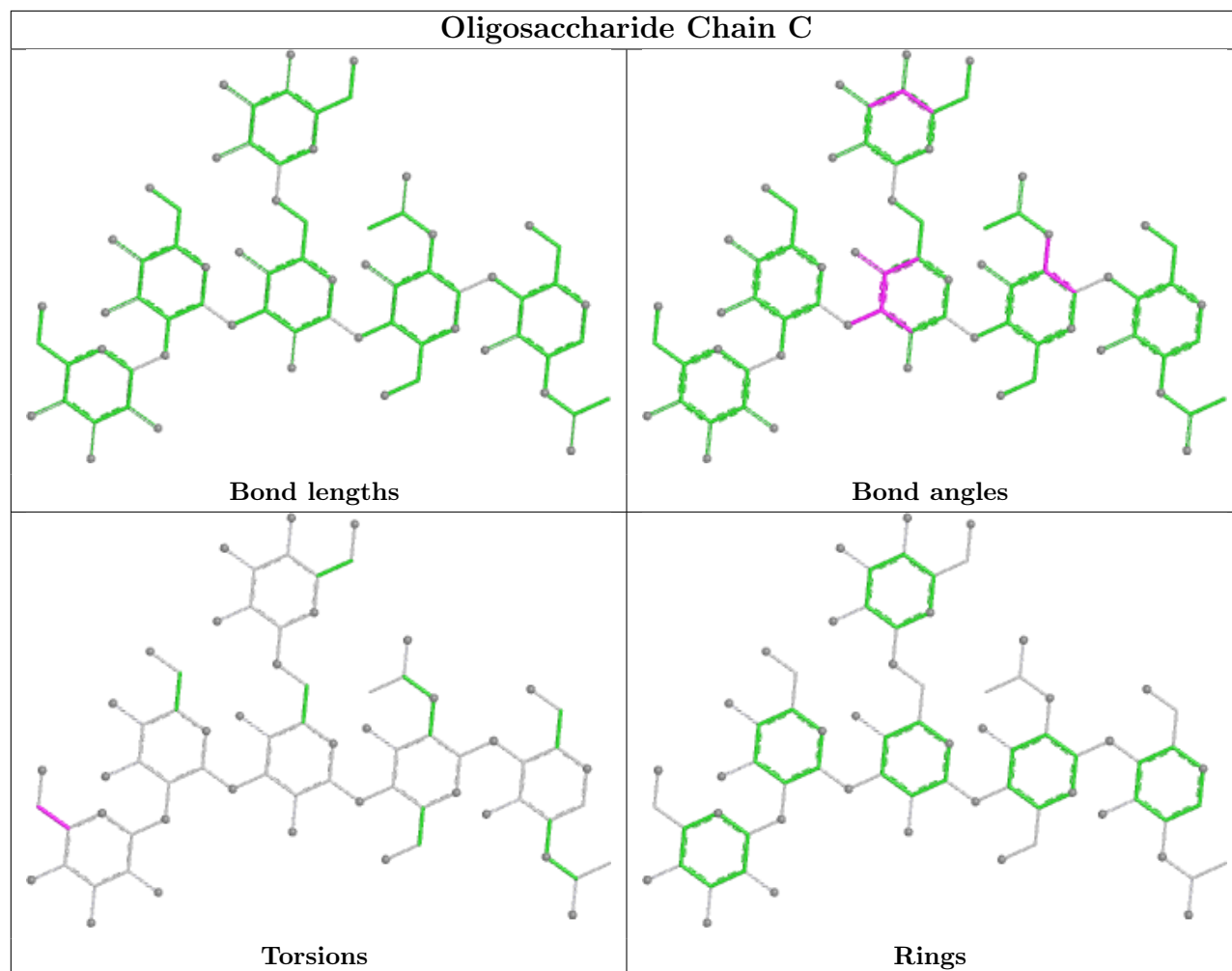
All (1) ring outliers are listed below:

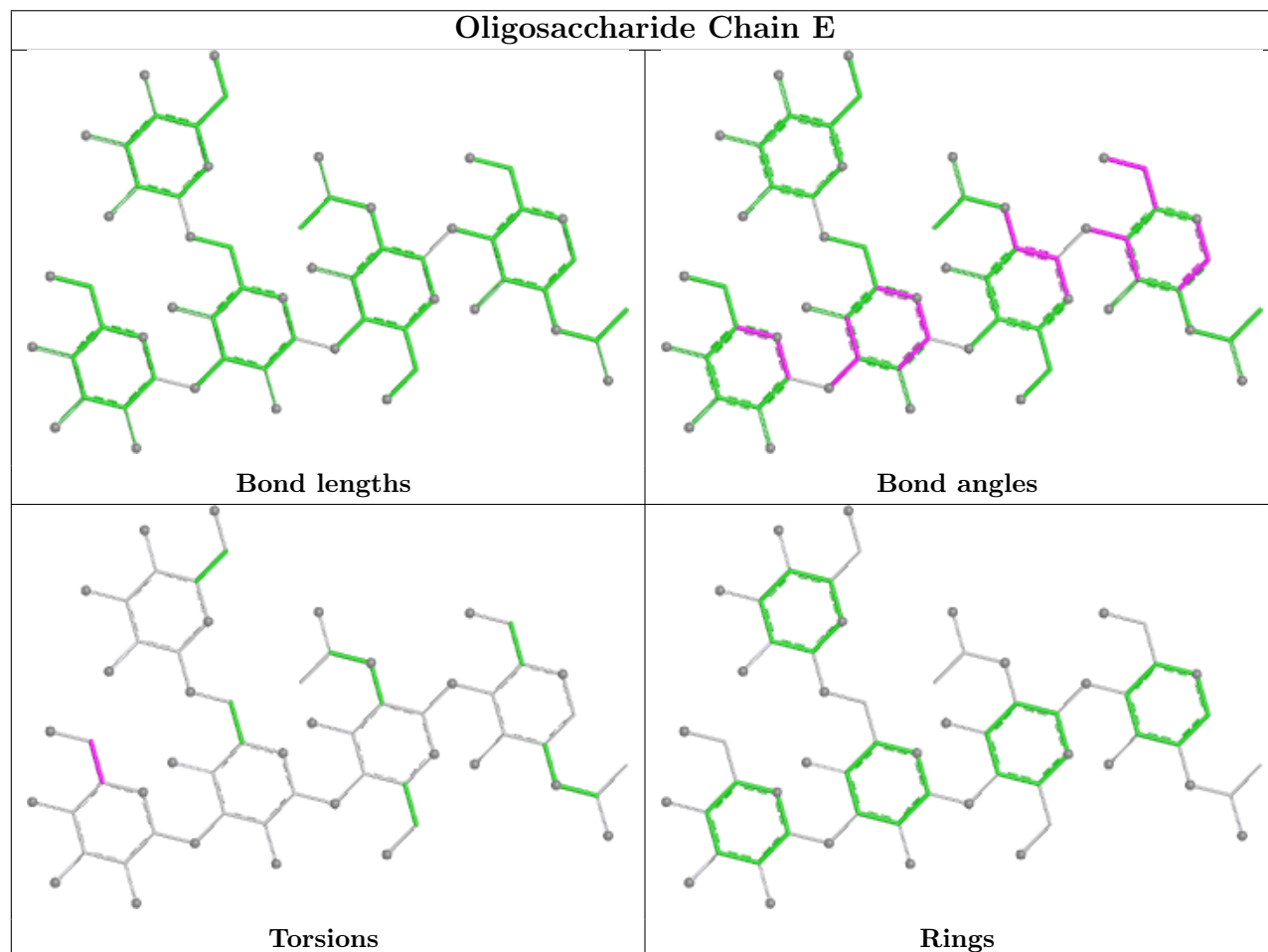
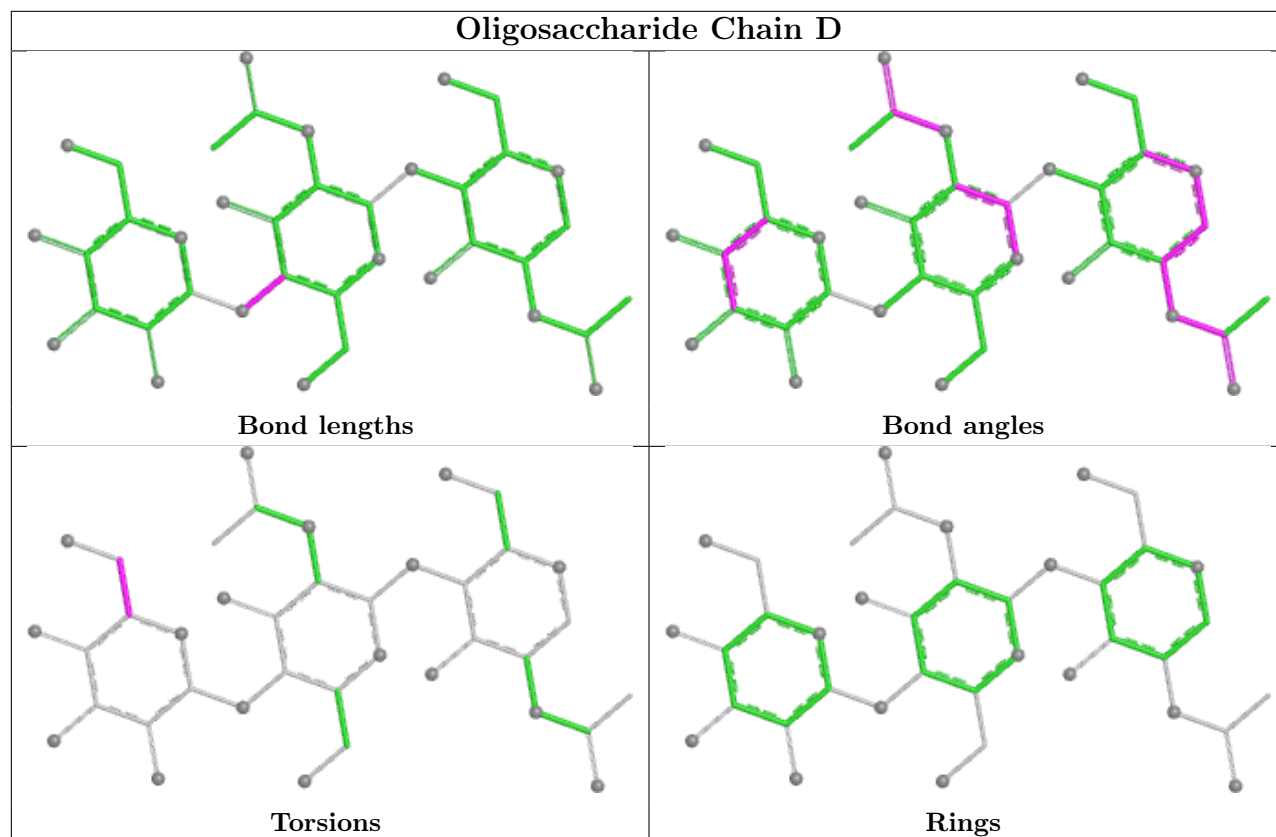
Mol	Chain	Res	Type	Atoms
5	F	7	MAN	C1-C2-C3-C4-C5-O5

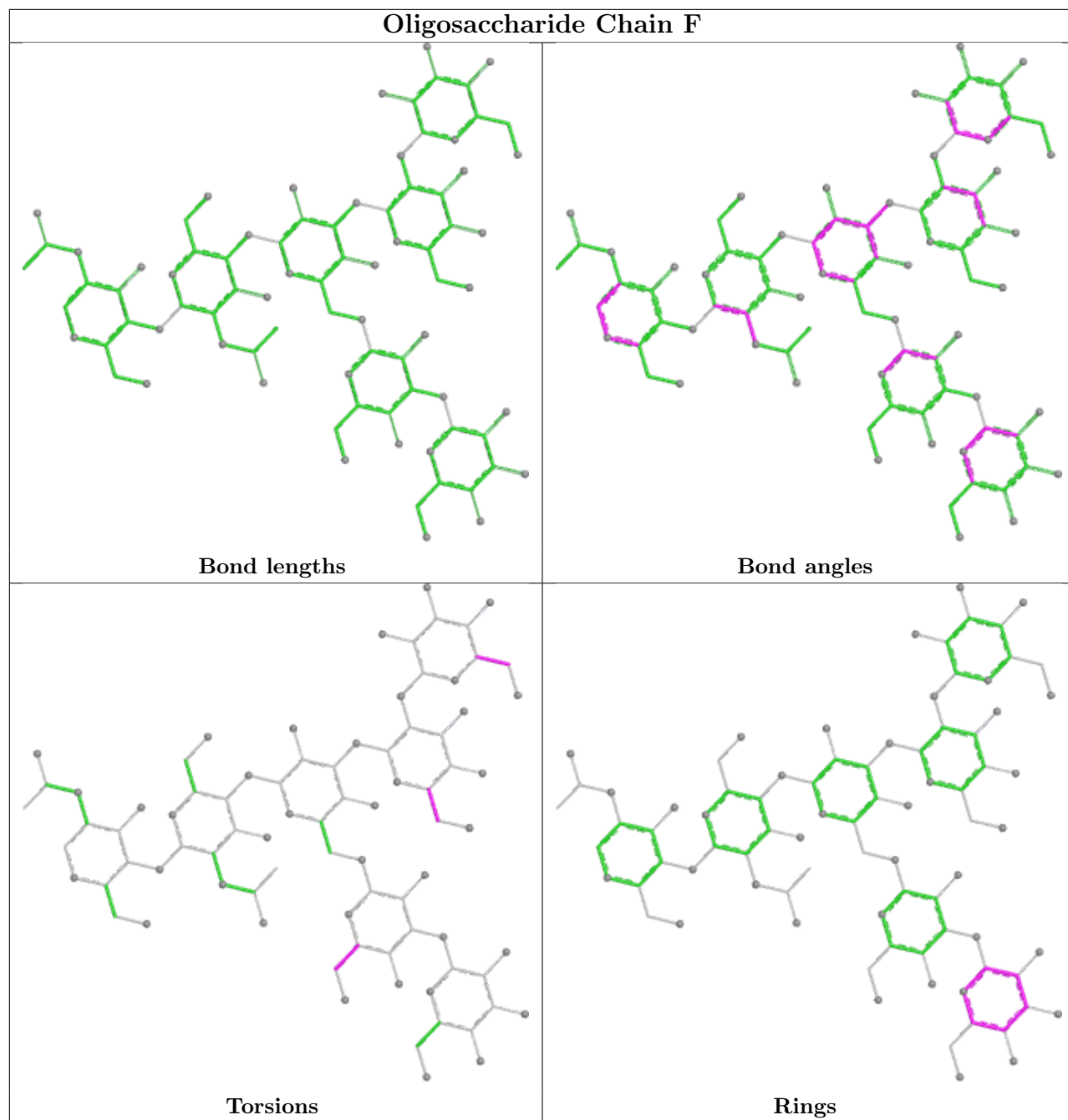
2 monomers are involved in 1 short contact:

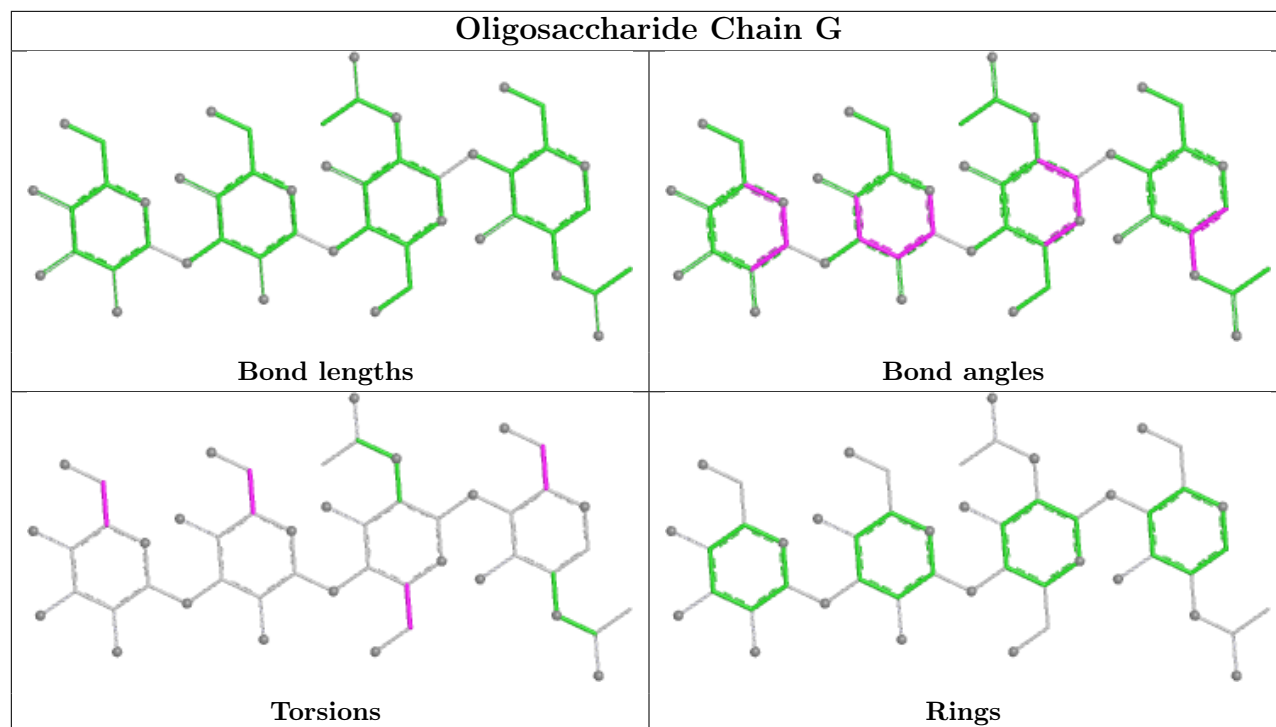
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	4	MAN	1	0
5	F	5	MAN	1	0

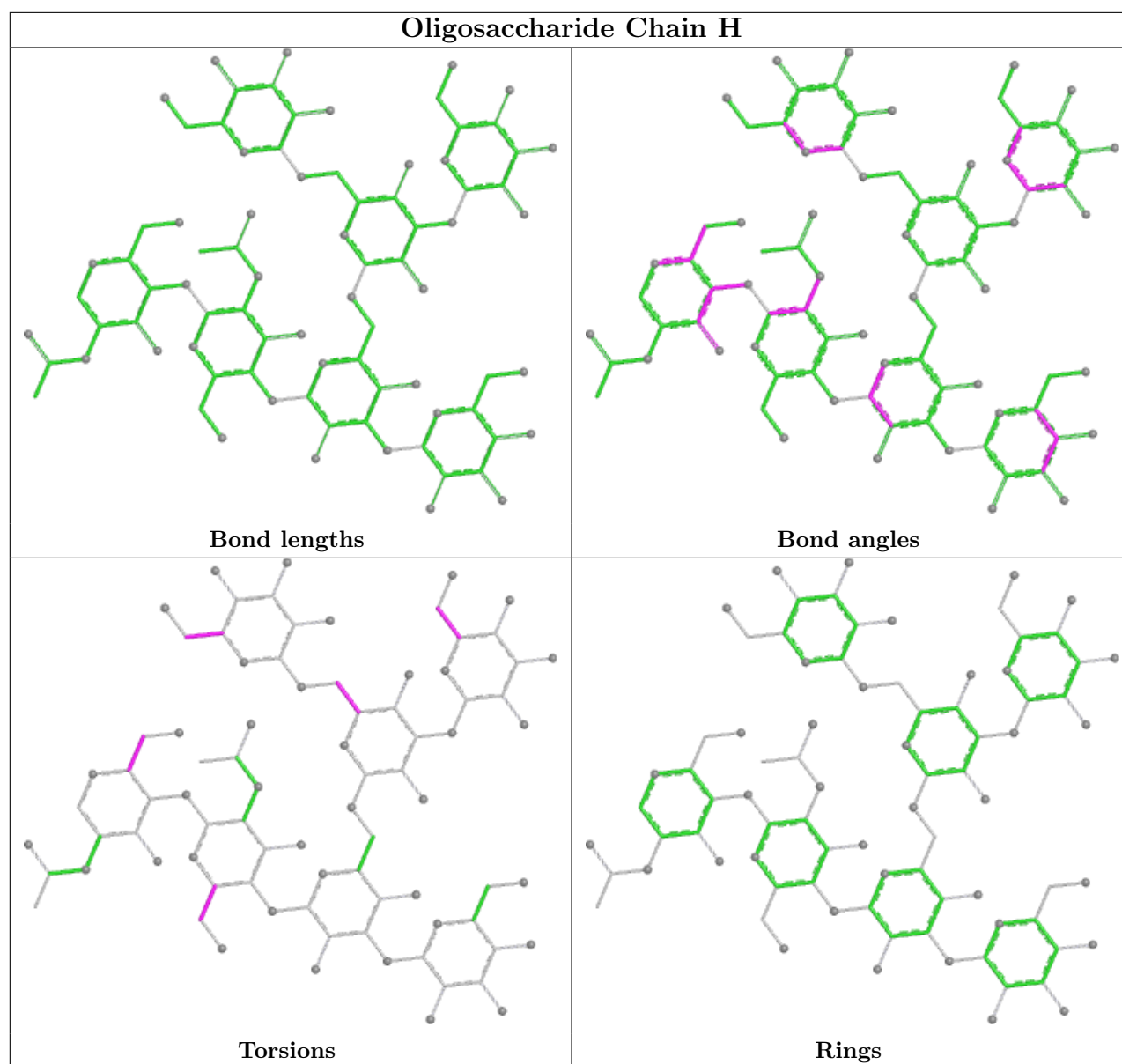
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

23 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
9	NAG	A	1006	1	14,14,15	0.41	0	17,19,21	1.37	3 (17%)
9	NAG	B	1005	1	14,14,15	0.32	0	17,19,21	1.06	2 (11%)
8	GD0	B	1001	-	21,22,22	1.61	3 (14%)	26,31,31	1.76	5 (19%)
8	GD0	A	1001	-	21,22,22	1.67	4 (19%)	26,31,31	1.78	5 (19%)
9	NAG	B	1002	1	14,14,15	0.49	0	17,19,21	0.94	0
9	NAG	B	1009	1	14,14,15	0.72	1 (7%)	17,19,21	2.02	2 (11%)
9	NAG	A	1009	1	14,14,15	0.28	0	17,19,21	0.77	0
9	NAG	B	1003	1	14,14,15	0.59	0	17,19,21	1.97	3 (17%)
9	NAG	B	1008	1	14,14,15	0.37	0	17,19,21	0.95	0
9	NAG	A	1008	1	14,14,15	0.28	0	17,19,21	0.95	0
9	NAG	A	1010	1	14,14,15	0.35	0	17,19,21	0.62	0
9	NAG	B	1011	1	14,14,15	0.51	0	17,19,21	2.58	3 (17%)
9	NAG	A	1012	1	14,14,15	0.35	0	17,19,21	1.02	1 (5%)
9	NAG	A	1002	1	14,14,15	0.45	0	17,19,21	0.75	0
9	NAG	A	1005	1	14,14,15	0.54	0	17,19,21	1.01	0
9	NAG	A	1007	1	14,14,15	0.60	0	17,19,21	1.78	3 (17%)
9	NAG	B	1010	1	14,14,15	0.20	0	17,19,21	0.95	0
9	NAG	B	1007	1	14,14,15	0.45	0	17,19,21	1.28	1 (5%)
9	NAG	A	1011	1	14,14,15	0.26	0	17,19,21	0.85	0
9	NAG	A	1003	1	14,14,15	0.50	0	17,19,21	1.69	2 (11%)
9	NAG	B	1004	1	14,14,15	0.49	0	17,19,21	1.74	3 (17%)
9	NAG	B	1006	1	14,14,15	0.60	0	17,19,21	1.15	2 (11%)
9	NAG	A	1004	1	14,14,15	0.52	0	17,19,21	0.99	1 (5%)

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
9	NAG	A	1006	1	-	1/6/23/26	0/1/1/1
9	NAG	B	1005	1	-	0/6/23/26	0/1/1/1
8	GD0	B	1001	-	-	0/13/13/13	0/2/2/2
8	GD0	A	1001	-	-	0/13/13/13	0/2/2/2
9	NAG	B	1002	1	-	2/6/23/26	0/1/1/1
9	NAG	B	1009	1	-	3/6/23/26	0/1/1/1
9	NAG	A	1009	1	-	0/6/23/26	0/1/1/1
9	NAG	B	1003	1	-	1/6/23/26	0/1/1/1
9	NAG	B	1008	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	A	1008	1	-	2/6/23/26	0/1/1/1
9	NAG	A	1010	1	-	2/6/23/26	0/1/1/1
9	NAG	B	1011	1	-	2/6/23/26	0/1/1/1
9	NAG	A	1012	1	-	0/6/23/26	0/1/1/1
9	NAG	A	1002	1	-	1/6/23/26	0/1/1/1
9	NAG	A	1005	1	-	2/6/23/26	0/1/1/1
9	NAG	A	1007	1	-	1/6/23/26	0/1/1/1
9	NAG	B	1010	1	-	0/6/23/26	0/1/1/1
9	NAG	B	1007	1	-	2/6/23/26	0/1/1/1
9	NAG	A	1011	1	-	0/6/23/26	0/1/1/1
9	NAG	A	1003	1	-	2/6/23/26	0/1/1/1
9	NAG	B	1004	1	-	0/6/23/26	0/1/1/1
9	NAG	B	1006	1	-	0/6/23/26	0/1/1/1
9	NAG	A	1004	1	-	0/6/23/26	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	1001	GD0	C7-I	5.12	2.19	2.07
8	B	1001	GD0	C7-I	4.94	2.18	2.07
8	A	1001	GD0	C6-N	-3.12	1.34	1.39
8	B	1001	GD0	C8-N1	3.08	1.37	1.32
8	B	1001	GD0	C6-N	-3.07	1.34	1.39
8	A	1001	GD0	C8-N1	3.03	1.37	1.32
8	A	1001	GD0	C8-N	-2.58	1.33	1.37
9	B	1009	NAG	C1-C2	2.26	1.55	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	1011	NAG	C1-O5-C5	9.18	124.49	112.19
9	B	1009	NAG	C1-O5-C5	7.48	122.21	112.19
9	B	1003	NAG	C1-O5-C5	6.82	121.32	112.19
8	B	1001	GD0	C9-N-C8	-6.06	122.73	127.87
9	A	1003	NAG	C1-O5-C5	5.61	119.71	112.19
9	A	1007	NAG	C2-N2-C7	5.42	130.17	122.90
8	A	1001	GD0	C9-N-C8	-5.41	123.29	127.87
9	B	1004	NAG	C3-C4-C5	3.86	117.22	110.23
9	B	1007	NAG	O5-C5-C6	3.85	115.15	107.66
9	B	1004	NAG	C4-C3-C2	3.75	116.52	111.02
8	A	1001	GD0	C10-C9-N	3.50	120.60	113.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	1004	NAG	O5-C1-C2	-3.40	106.04	111.29
8	B	1001	GD0	C10-C9-N	3.20	120.03	113.96
9	A	1006	NAG	C4-C3-C2	3.13	115.61	111.02
8	A	1001	GD0	C13-C8-N1	3.02	126.94	123.55
8	A	1001	GD0	C13-O1-C14	-3.02	106.72	112.97
9	B	1006	NAG	C3-C4-C5	2.96	115.59	110.23
9	A	1007	NAG	C3-C4-C5	2.86	115.41	110.23
9	B	1011	NAG	O5-C1-C2	2.76	115.56	111.29
9	A	1003	NAG	C4-C3-C2	-2.70	107.06	111.02
9	B	1011	NAG	C4-C3-C2	-2.68	107.09	111.02
8	A	1001	GD0	C13-C8-N	-2.68	119.94	123.34
9	A	1006	NAG	C3-C4-C5	2.54	114.83	110.23
8	B	1001	GD0	C13-O1-C14	-2.47	107.86	112.97
9	A	1006	NAG	O5-C5-C6	2.45	112.44	107.66
8	B	1001	GD0	C-C1-C3	2.42	121.65	118.57
9	B	1005	NAG	O5-C1-C2	-2.34	107.67	111.29
9	B	1005	NAG	C1-O5-C5	2.30	115.27	112.19
9	B	1003	NAG	C1-C2-N2	2.29	114.04	110.43
9	B	1009	NAG	O5-C1-C2	2.25	114.77	111.29
9	B	1003	NAG	O5-C1-C2	-2.19	107.91	111.29
9	B	1006	NAG	O4-C4-C3	-2.15	105.30	110.38
9	A	1007	NAG	O7-C7-N2	2.15	125.78	121.98
9	A	1004	NAG	O5-C5-C6	2.07	111.70	107.66
9	A	1012	NAG	O5-C1-C2	-2.07	108.08	111.29
8	B	1001	GD0	C9-N-C6	2.06	129.18	126.46

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	1007	NAG	C3-C2-N2-C7
9	A	1005	NAG	C4-C5-C6-O6
9	B	1008	NAG	O5-C5-C6-O6
9	A	1008	NAG	C4-C5-C6-O6
9	B	1011	NAG	C4-C5-C6-O6
9	B	1009	NAG	O5-C5-C6-O6
9	A	1003	NAG	C4-C5-C6-O6
9	B	1008	NAG	C4-C5-C6-O6
9	B	1011	NAG	O5-C5-C6-O6
9	A	1005	NAG	O5-C5-C6-O6
9	B	1009	NAG	C4-C5-C6-O6
9	A	1003	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
9	A	1008	NAG	O5-C5-C6-O6
9	B	1007	NAG	O5-C5-C6-O6
9	B	1002	NAG	O5-C5-C6-O6
9	A	1010	NAG	O5-C5-C6-O6
9	B	1003	NAG	C4-C5-C6-O6
9	B	1007	NAG	C4-C5-C6-O6
9	B	1002	NAG	C4-C5-C6-O6
9	B	1009	NAG	C3-C2-N2-C7
9	A	1006	NAG	O5-C5-C6-O6
9	A	1002	NAG	C4-C5-C6-O6
9	A	1010	NAG	C4-C5-C6-O6

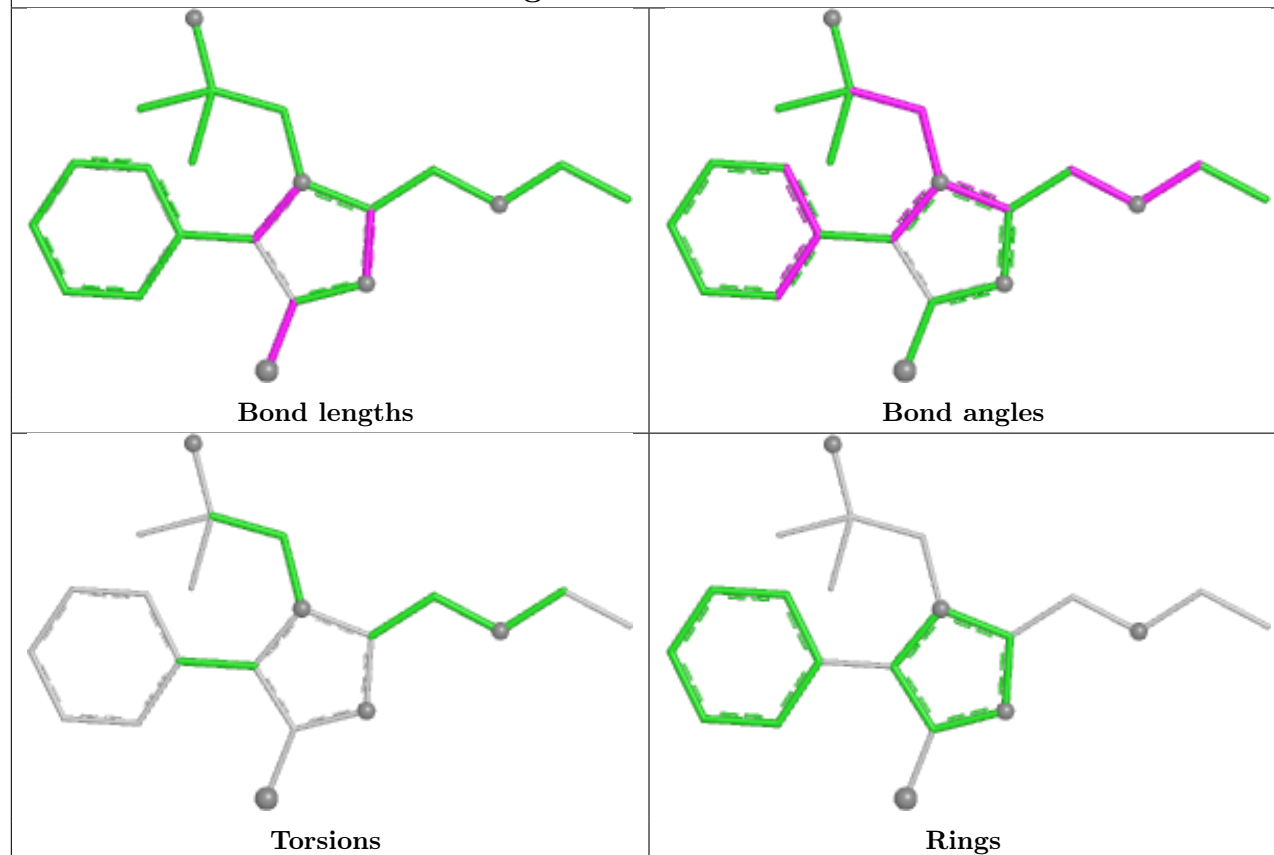
There are no ring outliers.

2 monomers are involved in 9 short contacts:

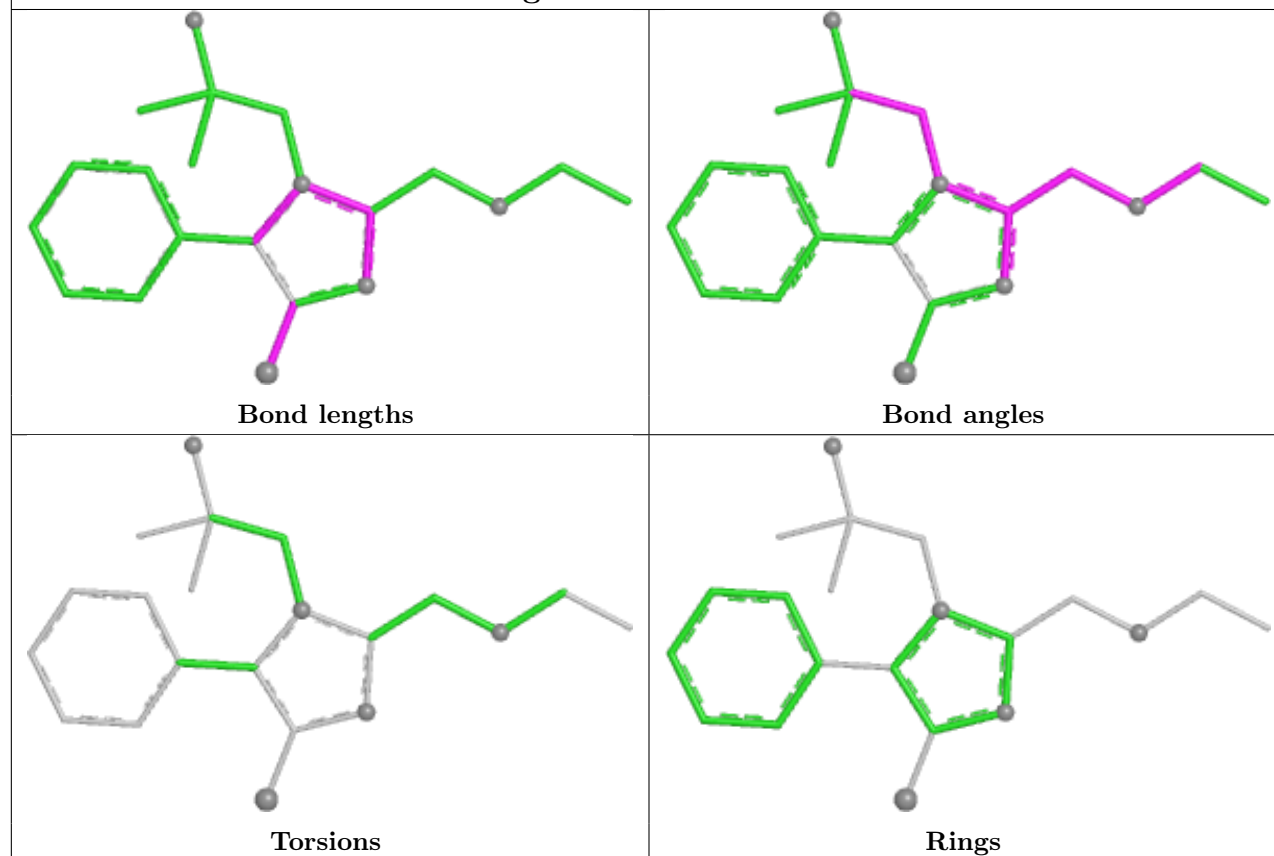
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	1001	GD0	3	0
8	A	1001	GD0	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand GD0 B 1001



Ligand GD0 A 1001



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	744/811 (91%)	-0.13	4 (0%) 87 83	44, 80, 120, 155	0
1	B	751/811 (92%)	-0.17	3 (0%) 88 85	42, 73, 113, 156	0
All	All	1495/1622 (92%)	-0.15	7 (0%) 87 83	42, 77, 117, 156	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	826	VAL	2.3
1	A	732	LEU	2.3
1	B	823	VAL	2.3
1	B	470	PHE	2.3
1	A	678	PHE	2.2
1	A	243	LYS	2.1
1	A	585	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
7	MAN	H	7	11/12	0.25	0.15	92,112,123,124	0
2	MAN	C	5	11/12	0.35	0.11	111,119,125,133	0

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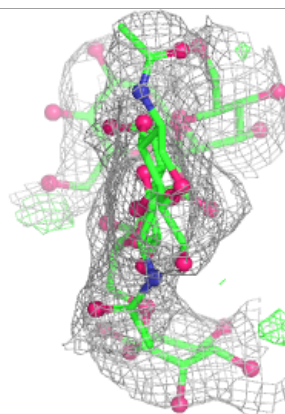
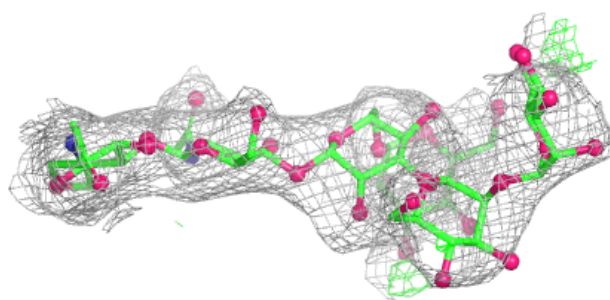
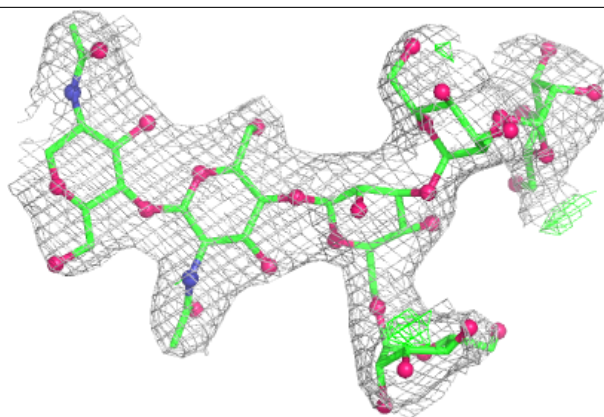
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	MAN	G	4	11/12	0.39	0.13	114,132,144,145	0
5	MAN	F	5	11/12	0.39	0.12	120,132,138,139	0
6	BMA	G	3	11/12	0.53	0.12	114,137,143,150	0
7	MAN	H	5	11/12	0.55	0.10	100,109,115,117	0
4	MAN	E	4	11/12	0.58	0.12	95,111,118,120	0
5	MAN	F	7	11/12	0.59	0.11	101,117,125,129	0
3	BMA	D	3	11/12	0.61	0.08	103,112,119,120	0
2	MAN	C	6	11/12	0.64	0.10	108,117,126,128	0
5	MAN	F	6	11/12	0.65	0.10	101,118,122,123	0
7	MAN	H	6	11/12	0.69	0.08	112,126,136,138	0
7	MAN	H	4	11/12	0.73	0.09	100,111,121,128	0
2	MAN	C	4	11/12	0.78	0.10	98,111,114,115	0
4	MAN	E	5	11/12	0.79	0.08	85,94,101,103	0
5	MAN	F	4	11/12	0.81	0.08	103,111,129,133	0
5	BMA	F	3	11/12	0.81	0.09	76,94,108,116	0
2	BMA	C	3	11/12	0.83	0.09	82,92,102,103	0
4	BMA	E	3	11/12	0.89	0.07	74,81,91,101	0
3	NAG	D	2	14/15	0.91	0.09	75,85,96,110	0
6	NAG	G	2	14/15	0.91	0.09	82,92,116,124	0
7	NAG	H	2	14/15	0.92	0.08	57,62,74,77	0
7	NAG	H	1	14/15	0.93	0.08	40,48,57,58	0
5	NAG	F	2	14/15	0.94	0.07	46,54,58,66	0
7	BMA	H	3	11/12	0.94	0.05	68,78,93,106	0
2	NAG	C	2	14/15	0.94	0.08	52,59,66,77	0
3	NAG	D	1	14/15	0.95	0.07	48,57,62,64	0
4	NAG	E	1	14/15	0.95	0.07	49,54,57,58	0
4	NAG	E	2	14/15	0.95	0.06	49,55,63,68	0
5	NAG	F	1	14/15	0.95	0.08	50,59,66,78	0
2	NAG	C	1	14/15	0.97	0.07	54,57,63,76	0
6	NAG	G	1	14/15	0.97	0.05	49,53,55,62	0

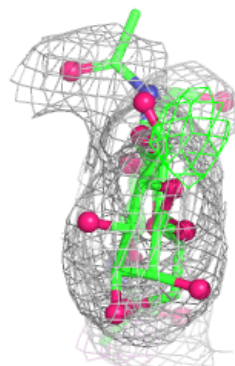
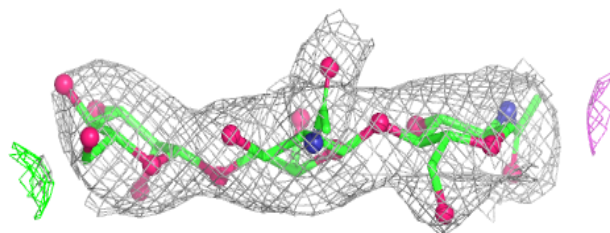
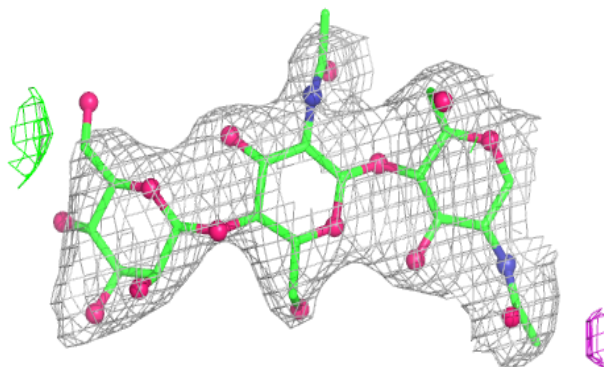
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

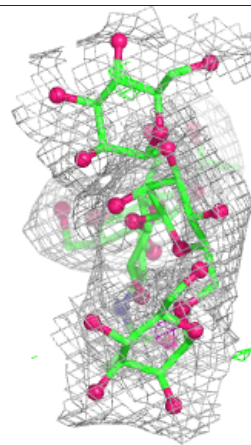
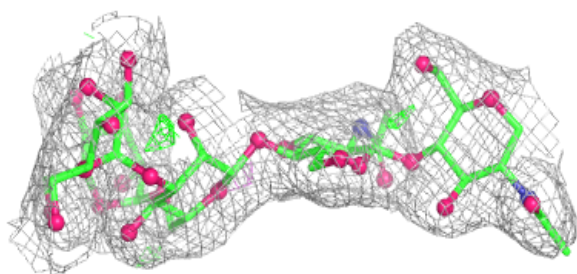
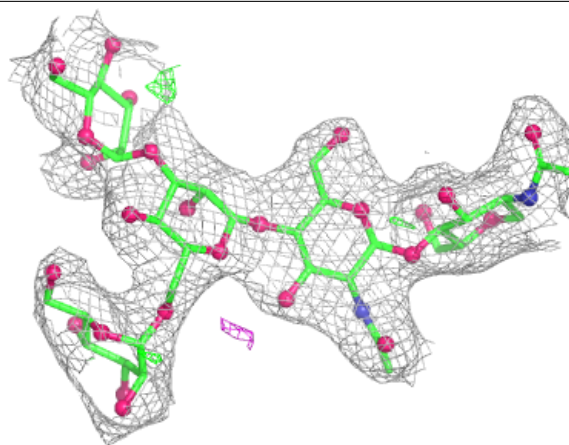
**Electron density around Chain D:**

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



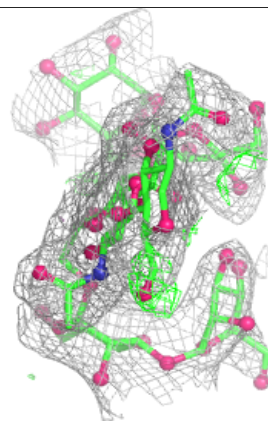
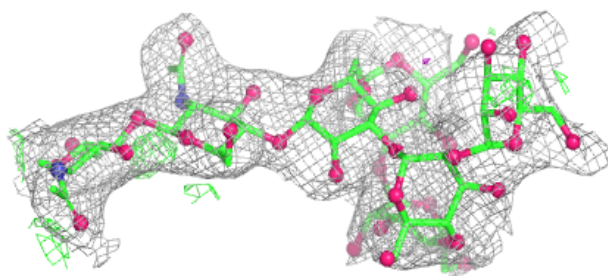
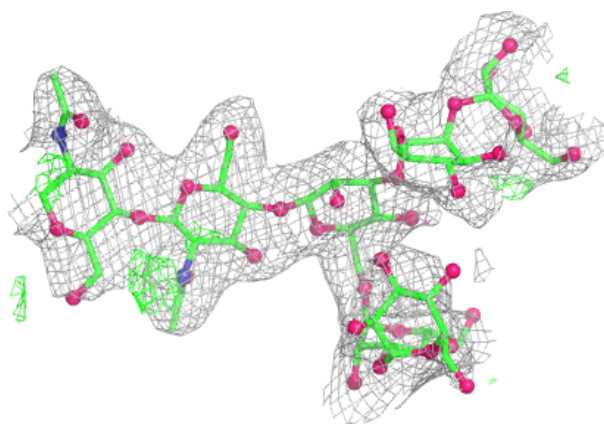
Electron density around Chain E:

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

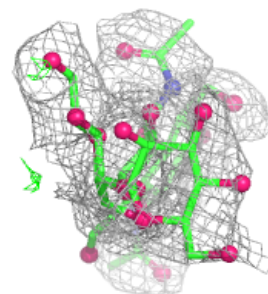
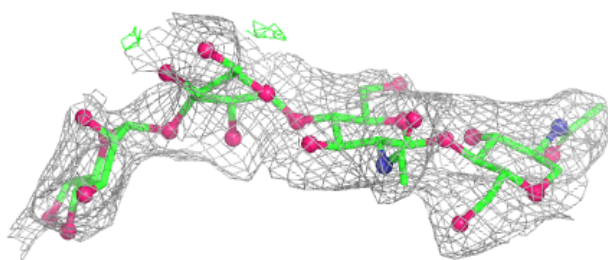
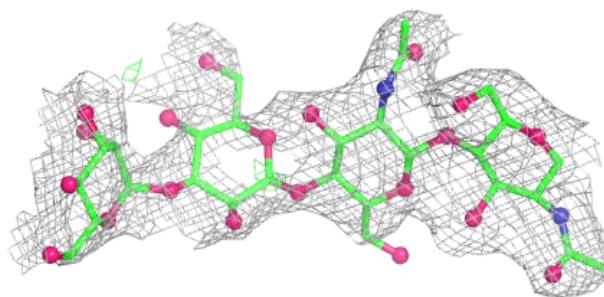


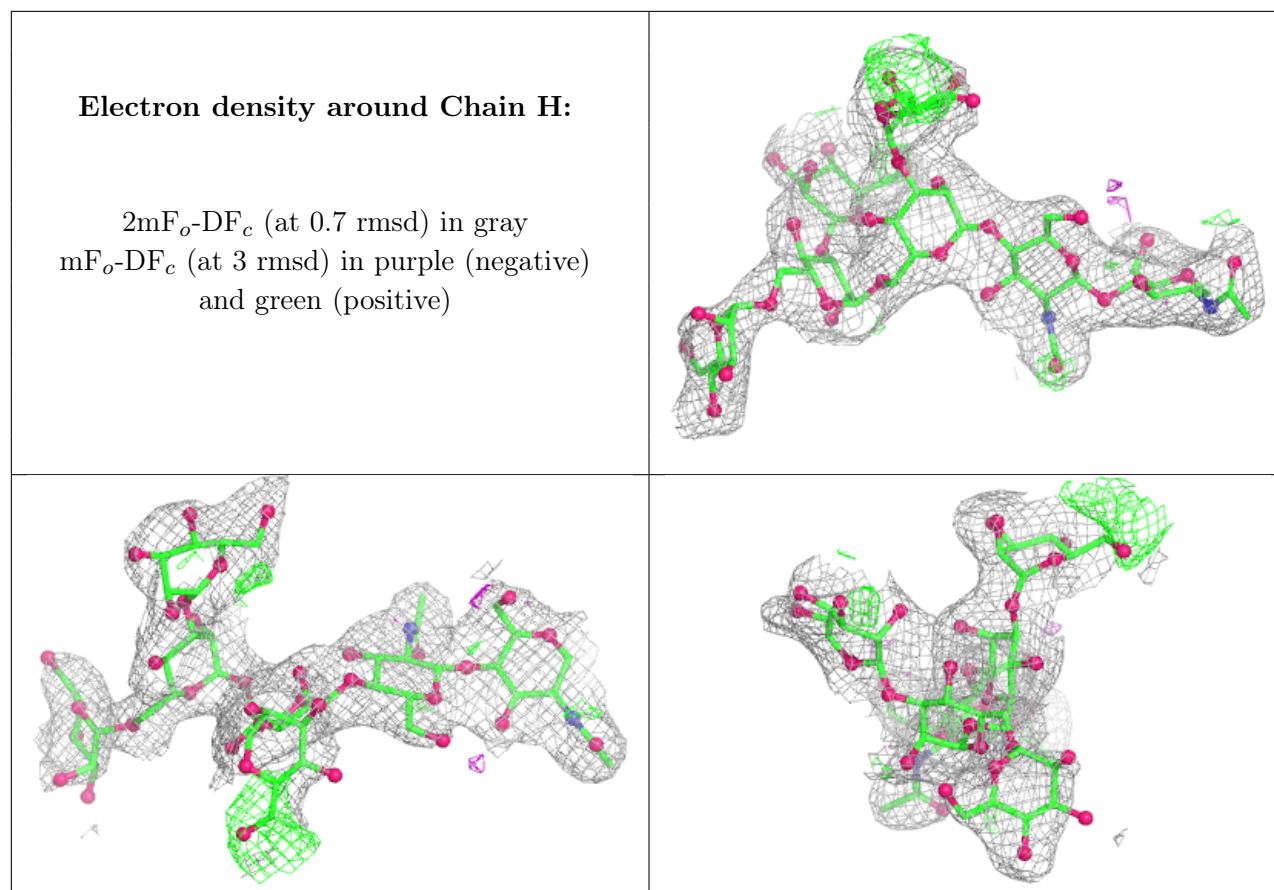
Electron density around Chain F:

$2mF_o-DF_c$ (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)





6.4 Ligands ⓘ

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
9	NAG	A	1003	14/15	0.60	0.12	104,131,139,142	0
9	NAG	A	1005	14/15	0.60	0.14	114,132,145,150	0
9	NAG	B	1002	14/15	0.61	0.12	136,141,148,149	0
9	NAG	A	1004	14/15	0.64	0.11	107,128,135,136	0
9	NAG	A	1007	14/15	0.65	0.15	82,111,121,130	0
9	NAG	A	1002	14/15	0.67	0.12	108,127,131,131	0
9	NAG	A	1006	14/15	0.68	0.10	90,107,111,112	0
9	NAG	B	1004	14/15	0.69	0.13	107,120,129,132	0
9	NAG	A	1008	14/15	0.73	0.10	110,115,121,125	0
9	NAG	A	1012	14/15	0.73	0.10	105,116,123,127	0
9	NAG	B	1009	14/15	0.75	0.09	101,115,119,128	0
9	NAG	B	1003	14/15	0.77	0.14	98,113,121,125	0
9	NAG	B	1011	14/15	0.77	0.10	103,114,118,120	0

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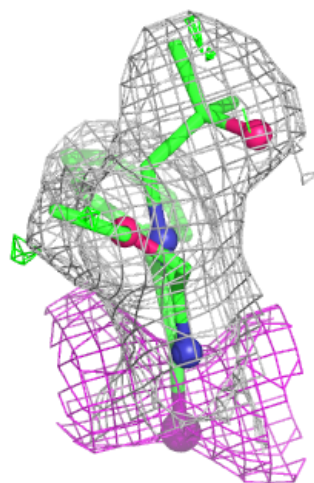
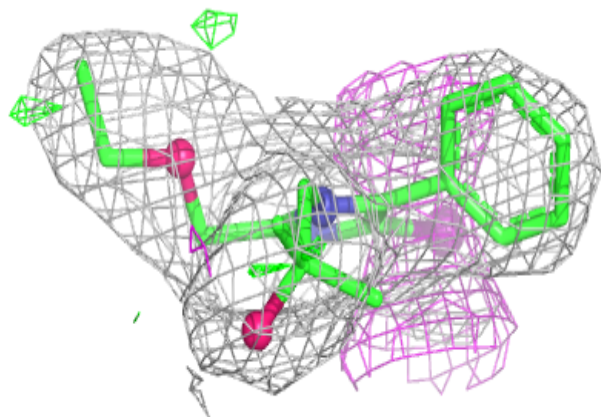
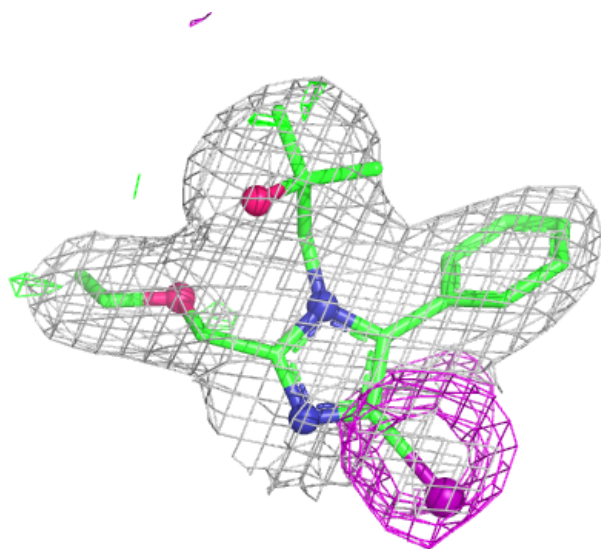
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	NAG	A	1009	14/15	0.83	0.08	86,97,102,104	0
9	NAG	B	1005	14/15	0.84	0.09	85,94,101,101	0
9	NAG	B	1010	14/15	0.86	0.10	83,92,97,98	0
9	NAG	B	1007	14/15	0.86	0.08	82,88,92,95	0
9	NAG	A	1010	14/15	0.89	0.07	92,96,104,111	0
9	NAG	B	1006	14/15	0.89	0.11	67,73,82,93	0
9	NAG	A	1011	14/15	0.90	0.07	87,93,95,97	0
9	NAG	B	1008	14/15	0.91	0.07	89,99,109,112	0
8	GD0	A	1001	21/21	0.92	0.13	62,68,82,123	0
8	GD0	B	1001	21/21	0.96	0.10	54,64,76,111	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

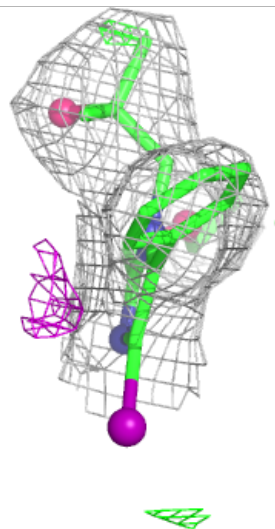
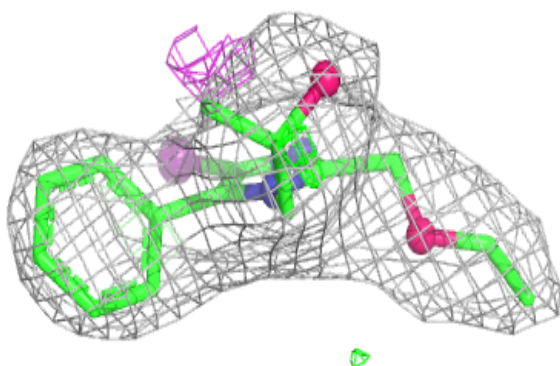
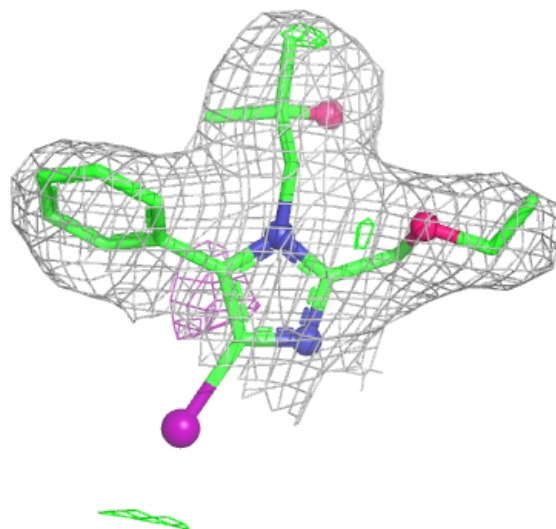
Electron density around GD0 A 1001:

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



Electron density around GD0 B 1001:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.