



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:56 AM UTC

PDB ID : 5WYX / pdb_00005wyx
Title : Crystal structure of human TLR8 in complex with CU-CPT8m
Authors : Tanji, H.; Ohto, U.; Shimizu, T.
Deposited on : 2017-01-16
Resolution : 2.40 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

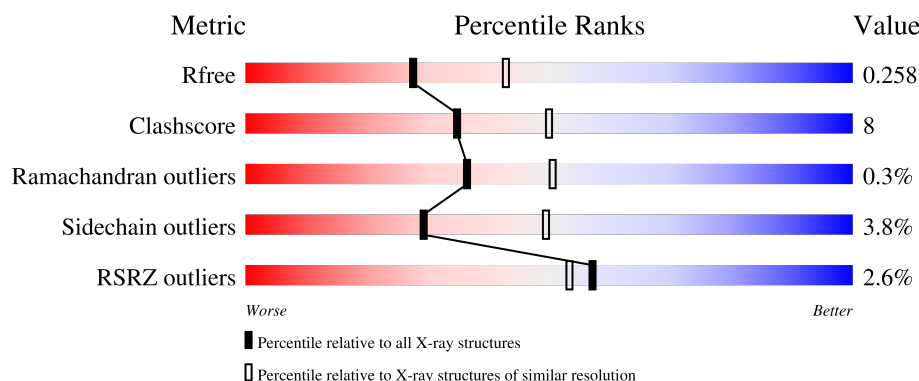
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



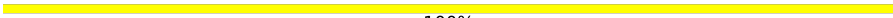
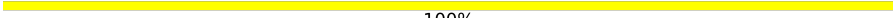
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	811	3% 76% 14% • 8%
1	B	811	2% 79% 11% • 9%
2	C	4	50% 50%
3	D	2	100%
3	G	2	100%

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Mol	Chain	Length	Quality of chain
4	E	5	 100%
5	F	4	 50%50%
5	H	4	 100%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 12621 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.

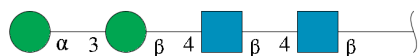
- Molecule 1 is a protein called Toll-like receptor 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	750	Total	C	N	O	S	0	0	0
			6038	3861	1026	1132	19			
1	B	739	Total	C	N	O	S	0	0	0
			5957	3810	1012	1116	19			

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



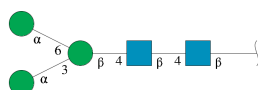
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 3 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
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- 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.



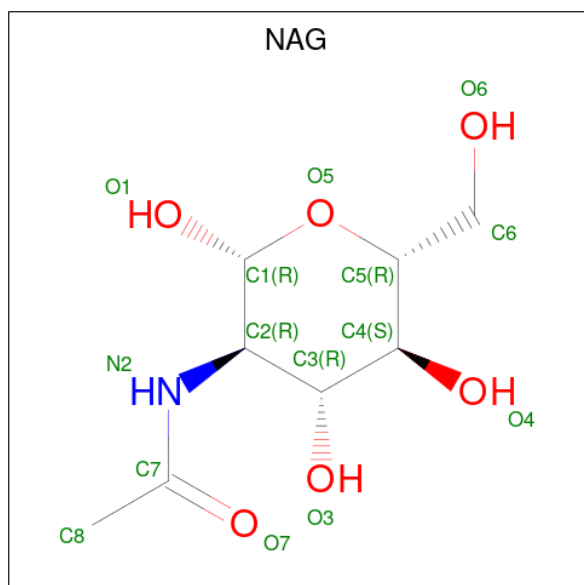
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



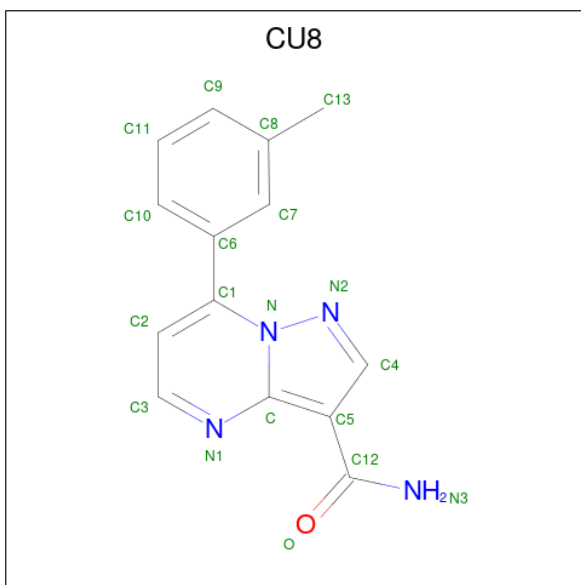
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	F	4	Total	C	N	O	0	0	0
			50	28	2	20			
5	H	4	Total	C	N	O	0	0	0
			50	28	2	20			

- 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	B	1	Total	C	N	O	0	0
			14	8	1	5		
6	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 7 is 7-(3-methylphenyl)pyrazolo[1,5-a]pyrimidine-3-carboxamide (CCD ID: CU8) (formula: $C_{14}H_{12}N_4O$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			19	14	4	1		
7	B	1	Total	C	N	O	0	0
			19	14	4	1		

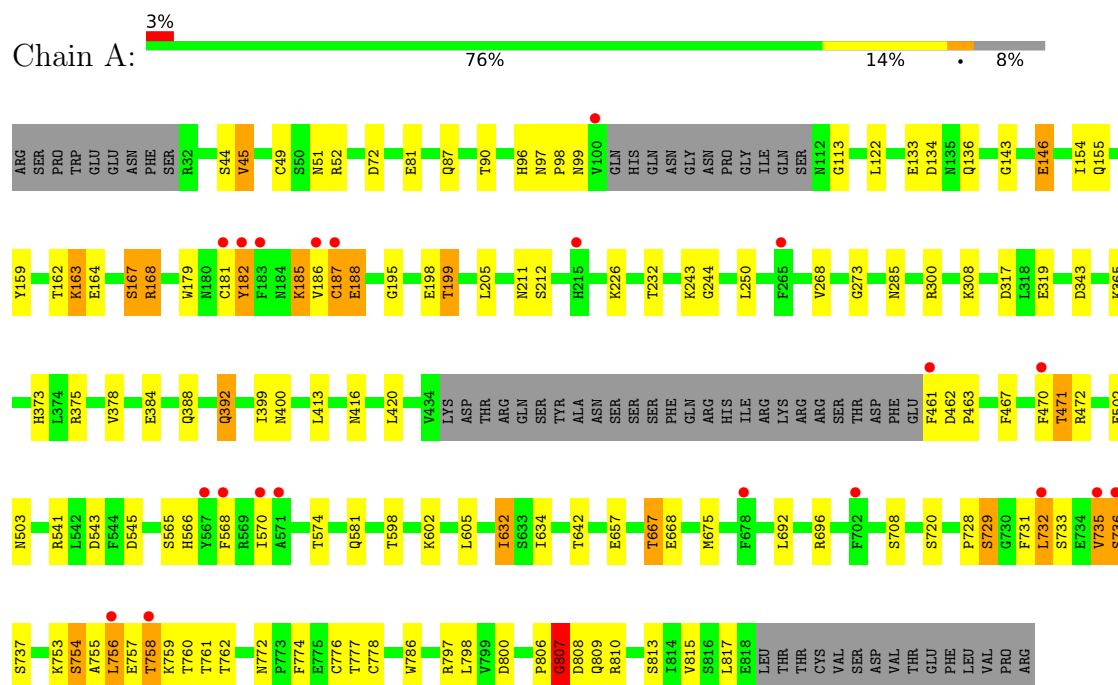
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	133	Total	O	0	0
			133	133		
8	B	104	Total	O	0	0
			104	104		

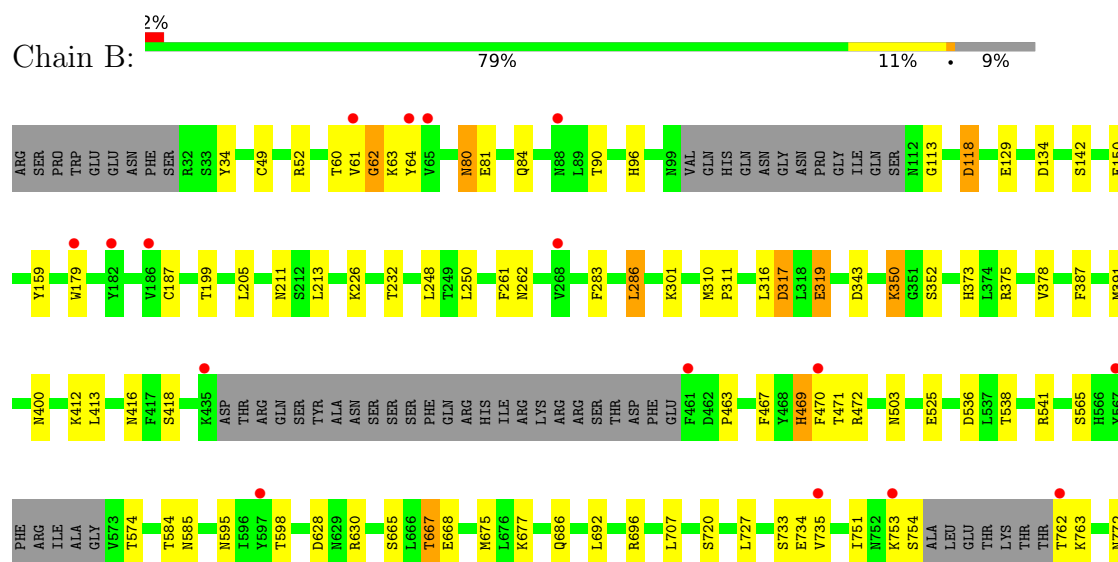
3 Residue-property plots [i](#)

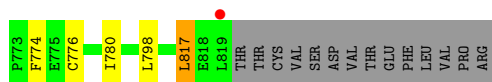
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Toll-like receptor 8

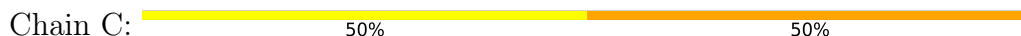


• Molecule 1: Toll-like receptor 8





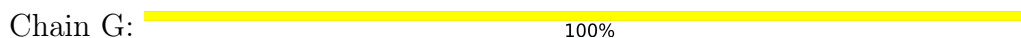
- Molecule 2: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



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



- 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



- Molecule 5: 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 5: 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



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	144.14Å 99.15Å 140.25Å 90.00° 108.07° 90.00°	Depositor
Resolution (Å)	40.16 – 2.40 40.16 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (40.16-2.40) 99.7 (40.16-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.201 , 0.256 0.206 , 0.258	Depositor DCC
R_{free} test set	3589 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	50.8	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12621	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: CU8, NAG, BMA, 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	0.92	0/6162	1.00	10/8358 (0.1%)
1	B	0.95	1/6078 (0.0%)	0.99	9/8240 (0.1%)
All	All	0.93	1/12240 (0.0%)	1.00	19/16598 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	665	SER	C-O	-5.47	1.17	1.24

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	62	GLY	N-CA-C	-10.99	97.88	110.96
1	A	163	LYS	N-CA-C	-9.99	97.44	110.33
1	A	188	GLU	N-CA-C	-8.96	97.98	110.35
1	A	182	TYR	N-CA-C	-7.56	98.00	109.79
1	A	99	ASN	N-CA-C	7.19	121.56	110.42
1	B	317	ASP	N-CA-C	6.83	120.03	108.90
1	B	472	ARG	CA-C-N	-6.41	113.38	119.85
1	B	472	ARG	C-N-CA	-6.41	113.38	119.85
1	B	316	LEU	CA-C-N	-6.31	114.10	122.99
1	B	316	LEU	C-N-CA	-6.31	114.10	122.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	471	THR	N-CA-C	6.07	121.43	113.30
1	A	167	SER	N-CA-C	-5.69	101.85	110.28
1	B	469	HIS	N-CA-C	5.43	117.61	107.99
1	A	807	GLY	N-CA-C	-5.30	100.62	113.18
1	A	268	VAL	CA-C-N	-5.23	114.86	120.03
1	A	268	VAL	C-N-CA	-5.23	114.86	120.03
1	B	113	GLY	N-CA-C	5.18	125.46	113.18
1	A	45	VAL	CB-CA-C	5.17	118.31	110.62
1	B	817	LEU	N-CA-C	5.03	116.81	110.33

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	185	LYS	Peptide
1	A	187	CYS	Peptide
1	A	807	GLY	Peptide

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	6038	0	6032	124	0
1	B	5957	0	5950	63	0
2	C	50	0	43	2	0
3	D	28	0	25	0	0
3	G	28	0	25	0	0
4	E	61	0	52	0	0
5	F	50	0	43	4	0
5	H	50	0	43	0	0
6	A	56	0	52	0	0
6	B	28	0	26	0	0
7	A	19	0	0	0	0
7	B	19	0	0	0	0
8	A	133	0	0	8	0
8	B	104	0	0	3	0
All	All	12621	0	12291	186	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 8.

All (186) 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:273:GLY:O	1:A:300:ARG:NH2	1.74	1.18
1:A:753:LYS:HA	1:A:756:LEU:HB3	1.45	0.99
1:A:167:SER:HB2	1:A:199:THR:HG21	1.44	0.98
1:A:777:THR:HB	1:A:778:CYS:HA	1.46	0.98
1:B:34:TYR:O	1:B:60:THR:HG21	1.67	0.95
1:B:469:HIS:CD2	1:B:470:PHE:CD2	2.56	0.94
1:A:735:VAL:HA	1:A:736:SER:CB	2.01	0.91
1:A:732:LEU:HD22	1:A:755:ALA:O	1.73	0.87
1:A:735:VAL:HA	1:A:736:SER:HB3	1.57	0.86
1:A:319:GLU:HG2	8:A:1055:HOH:O	1.74	0.85
1:A:167:SER:O	1:A:168:ARG:HB2	1.77	0.84
1:A:167:SER:CB	1:A:199:THR:HG21	2.07	0.83
1:B:62:GLY:HA3	1:B:63:LYS:HG3	1.60	0.82
1:B:469:HIS:CD2	1:B:470:PHE:HD2	1.99	0.81
1:A:186:VAL:HA	1:A:187:CYS:C	2.06	0.80
1:A:732:LEU:CD2	1:A:755:ALA:O	2.32	0.77
1:B:776:CYS:HB3	1:B:817:LEU:HD21	1.68	0.76
1:A:754:SER:N	1:A:756:LEU:O	2.19	0.75
1:A:187:CYS:SG	1:A:188:GLU:O	2.44	0.75
1:A:732:LEU:O	1:A:762:THR:HG21	1.88	0.73
1:B:469:HIS:NE2	1:B:470:PHE:CD2	2.57	0.72
1:B:319:GLU:OE1	1:B:469:HIS:HB2	1.88	0.72
1:A:807:GLY:HA2	1:A:810:ARG:H	1.54	0.72
1:A:777:THR:CB	1:A:778:CYS:HA	2.19	0.69
1:A:753:LYS:CA	1:A:756:LEU:HB3	2.20	0.68
1:B:226:LYS:NZ	5:F:2:NAG:H81	2.07	0.68
1:A:187:CYS:CB	1:A:188:GLU:O	2.42	0.68
1:A:733:SER:OG	1:A:757:GLU:HG2	1.94	0.67
1:A:545:ASP:OD2	8:A:1001:HOH:O	2.11	0.67
1:A:187:CYS:HB2	1:A:188:GLU:O	1.94	0.67
1:A:122:LEU:HD13	1:A:143:GLY:O	1.95	0.67
1:B:62:GLY:CA	1:B:63:LYS:HG3	2.24	0.67
1:B:63:LYS:O	1:B:64:TYR:HB2	1.95	0.66
1:A:807:GLY:HA3	1:A:810:ARG:HB3	1.75	0.65
1:A:154:ILE:HD13	1:A:463:PRO:HG2	1.79	0.65
1:A:809:GLN:NE2	1:A:817:LEU:HD11	2.11	0.65
1:A:777:THR:HB	1:A:778:CYS:CA	2.27	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:807:GLY:HA2	1:A:809:GLN:N	2.13	0.63
1:A:467:PHE:HB3	2:C:1:NAG:H81	1.81	0.63
1:B:735:VAL:HG23	8:B:1014:HOH:O	1.99	0.63
1:A:754:SER:H	1:A:756:LEU:C	2.06	0.62
1:B:375:ARG:HD2	1:B:400:ASN:HD21	1.63	0.62
1:A:632:ILE:HD12	1:A:657:GLU:HG2	1.82	0.61
1:A:735:VAL:HG13	1:A:736:SER:OG	2.01	0.60
1:A:195:GLY:O	1:A:198:GLU:HG2	2.02	0.60
1:A:732:LEU:O	1:A:762:THR:CG2	2.48	0.60
1:A:760:THR:H	1:A:761:THR:HA	1.65	0.60
1:B:584:THR:HG22	8:B:1092:HOH:O	2.02	0.59
1:A:179:TRP:CD1	1:A:463:PRO:HB3	2.38	0.58
1:A:732:LEU:C	1:A:732:LEU:HD23	2.29	0.58
1:A:807:GLY:CA	1:A:810:ARG:H	2.15	0.58
1:A:541:ARG:HA	1:A:565:SER:O	2.03	0.57
1:A:159:TYR:CD1	1:A:187:CYS:HB3	2.40	0.57
1:A:163:LYS:O	1:A:164:GLU:CB	2.52	0.56
1:B:469:HIS:NE2	1:B:470:PHE:CE2	2.73	0.56
1:B:734:GLU:HA	1:B:735:VAL:C	2.28	0.56
1:A:211:ASN:O	1:A:232:THR:HA	2.05	0.56
1:B:80:ASN:N	1:B:80:ASN:OD1	2.39	0.56
1:A:776:CYS:O	1:A:806:PRO:CD	2.55	0.55
1:A:243:LYS:HD2	1:A:285:ASN:HD22	1.72	0.55
1:A:728:PRO:O	1:A:731:PHE:HB3	2.06	0.55
1:A:776:CYS:O	1:A:806:PRO:HD3	2.06	0.55
1:B:226:LYS:HZ2	5:F:2:NAG:H81	1.70	0.55
1:A:243:LYS:HD2	1:A:285:ASN:ND2	2.23	0.54
1:B:211:ASN:O	1:B:232:THR:HA	2.07	0.54
1:B:129:GLU:HG2	1:B:150:GLU:HB3	1.89	0.54
1:A:162:THR:C	1:A:163:LYS:O	2.47	0.54
1:A:375:ARG:HD2	1:A:400:ASN:HD21	1.73	0.54
1:A:758:THR:OG1	1:A:759:LYS:HA	2.07	0.54
1:B:317:ASP:HB2	1:B:343:ASP:HB3	1.91	0.53
1:A:756:LEU:HG	1:A:786:TRP:CD1	2.44	0.53
1:B:261:PHE:CE2	1:B:350:LYS:HD3	2.43	0.53
1:B:541:ARG:HA	1:B:565:SER:O	2.08	0.53
1:A:735:VAL:HA	1:A:736:SER:OG	2.07	0.53
1:A:392:GLN:HB2	8:A:1002:HOH:O	2.08	0.52
1:B:470:PHE:O	1:B:471:THR:HB	2.09	0.52
1:A:146:GLU:H	1:A:146:GLU:CD	2.17	0.52
1:A:226:LYS:NZ	2:C:2:NAG:H81	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:729:SER:HA	1:A:754:SER:C	2.34	0.52
1:B:469:HIS:NE2	1:B:470:PHE:HD2	1.99	0.52
1:A:728:PRO:HB2	1:A:731:PHE:HB2	1.92	0.52
1:B:96:HIS:HD2	1:B:134:ASP:OD2	1.93	0.52
1:A:807:GLY:HA2	1:A:808:ASP:C	2.36	0.51
1:A:642:THR:CG2	8:A:1117:HOH:O	2.60	0.50
1:A:181:CYS:SG	1:A:187:CYS:N	2.84	0.50
1:A:605:LEU:HD13	1:A:634:ILE:HG12	1.92	0.50
1:A:777:THR:CB	1:A:778:CYS:CA	2.86	0.50
1:A:566:HIS:HE1	1:B:262:ASN:O	1.95	0.50
1:B:34:TYR:O	1:B:60:THR:CG2	2.49	0.50
1:A:566:HIS:HD2	1:A:568:PHE:H	1.59	0.49
1:A:470:PHE:CZ	1:A:472:ARG:HG2	2.48	0.49
1:A:185:LYS:N	1:A:186:VAL:O	2.45	0.49
1:B:776:CYS:HB3	1:B:817:LEU:CD2	2.41	0.49
1:A:113:GLY:HA3	1:A:136:GLN:HB2	1.95	0.49
1:B:205:LEU:C	1:B:205:LEU:HD23	2.37	0.49
1:A:44:SER:HB3	8:A:1122:HOH:O	2.12	0.49
1:A:797:ARG:NH1	1:A:800:ASP:OD2	2.46	0.49
1:A:96:HIS:HD2	1:A:134:ASP:OD2	1.96	0.49
1:A:756:LEU:C	1:A:756:LEU:HD23	2.39	0.48
1:B:373:HIS:HA	1:B:400:ASN:HB3	1.95	0.48
1:B:61:VAL:HG12	1:B:62:GLY:N	2.28	0.48
1:A:416:ASN:ND2	8:A:1008:HOH:O	2.46	0.48
1:A:753:LYS:O	1:A:754:SER:OG	2.29	0.48
1:B:261:PHE:CE2	1:B:350:LYS:CD	2.96	0.48
1:A:155:GLN:HE21	1:A:463:PRO:HG3	1.78	0.48
1:B:733:SER:O	1:B:734:GLU:C	2.57	0.48
1:A:502:GLU:O	1:A:503:ASN:HB2	2.13	0.47
1:A:185:LYS:CG	1:A:186:VAL:H	2.28	0.47
1:A:568:PHE:C	1:A:570:ILE:H	2.21	0.47
1:A:754:SER:HA	1:A:755:ALA:C	2.40	0.47
1:A:813:SER:OG	1:A:815:VAL:HG22	2.14	0.47
1:B:226:LYS:HZ3	5:F:2:NAG:H81	1.79	0.47
1:B:179:TRP:CD1	1:B:463:PRO:HB3	2.49	0.47
1:A:731:PHE:O	1:A:732:LEU:HD22	2.14	0.47
1:A:754:SER:H	1:A:756:LEU:CA	2.28	0.47
1:B:387:PHE:O	1:B:391:MET:HG3	2.15	0.47
1:B:707:LEU:HD23	1:B:735:VAL:HG11	1.96	0.46
1:B:283:PHE:HA	1:B:286:LEU:HD22	1.96	0.46
1:B:772:ASN:HB2	1:B:774:PHE:CE2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:733:SER:HB2	1:A:757:GLU:HB3	1.97	0.46
1:B:753:LYS:O	1:B:754:SER:C	2.59	0.46
1:B:467:PHE:HB3	5:F:1:NAG:H81	1.98	0.46
1:A:817:LEU:HA	8:A:1077:HOH:O	2.16	0.45
1:A:373:HIS:HA	1:A:400:ASN:HB3	1.98	0.45
1:A:243:LYS:CB	1:A:244:GLY:HA2	2.46	0.45
1:A:735:VAL:HG12	1:A:737:SER:N	2.31	0.45
1:B:667:THR:HG22	1:B:668:GLU:HG3	1.97	0.45
1:A:182:TYR:H	1:A:182:TYR:HD1	1.63	0.45
1:A:696:ARG:HG2	1:A:720:SER:OG	2.16	0.45
1:A:776:CYS:C	1:A:777:THR:HG23	2.41	0.45
1:A:753:LYS:CB	1:A:756:LEU:HD13	2.46	0.45
1:B:470:PHE:O	1:B:471:THR:CB	2.64	0.45
1:A:692:LEU:C	1:A:692:LEU:HD23	2.41	0.44
1:B:412:LYS:HB3	1:B:503:ASN:HB3	1.99	0.44
1:B:310:MET:N	1:B:311:PRO:HD3	2.32	0.44
1:A:807:GLY:HA2	1:A:810:ARG:N	2.28	0.44
1:A:186:VAL:CA	1:A:187:CYS:C	2.83	0.44
1:A:461:PHE:CG	1:A:462:ASP:N	2.83	0.44
1:B:159:TYR:CG	1:B:187:CYS:HB2	2.53	0.44
1:B:692:LEU:C	1:B:692:LEU:HD23	2.43	0.44
1:B:696:ARG:HG2	1:B:720:SER:OG	2.18	0.44
1:A:543:ASP:C	1:A:543:ASP:OD2	2.60	0.44
1:A:574:THR:HG22	1:A:598:THR:HG23	2.00	0.44
1:A:181:CYS:HB2	1:A:187:CYS:HB3	1.93	0.44
1:A:413:LEU:C	1:A:413:LEU:HD12	2.43	0.44
1:A:733:SER:CB	1:A:757:GLU:HB3	2.48	0.43
1:B:584:THR:CG2	8:B:1092:HOH:O	2.63	0.43
1:B:118:ASP:OD2	1:B:118:ASP:N	2.51	0.43
1:A:754:SER:N	1:A:756:LEU:N	2.66	0.43
1:A:185:LYS:O	1:A:187:CYS:O	2.35	0.43
1:A:667:THR:HG22	1:A:668:GLU:HG3	2.01	0.43
1:B:584:THR:C	1:B:585:ASN:HD22	2.26	0.43
1:A:97:ASN:HA	1:A:98:PRO:HA	1.81	0.43
1:A:205:LEU:C	1:A:205:LEU:HD23	2.44	0.43
1:B:471:THR:O	1:B:471:THR:HG22	2.19	0.43
1:A:122:LEU:HA	1:A:122:LEU:HD12	1.81	0.43
1:A:319:GLU:HG3	1:A:343:ASP:CG	2.44	0.43
1:A:133:GLU:CG	1:A:154:ILE:HG13	2.49	0.43
1:B:250:LEU:C	1:B:250:LEU:HD23	2.44	0.43
1:B:536:ASP:OD1	1:B:538:THR:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:733:SER:OG	1:A:757:GLU:CG	2.64	0.42
1:A:735:VAL:HG12	1:A:737:SER:H	1.84	0.42
1:B:310:MET:N	1:B:311:PRO:CD	2.83	0.42
1:A:51:ASN:HA	1:A:72:ASP:O	2.19	0.42
1:A:133:GLU:HG2	1:A:154:ILE:HG13	2.00	0.42
1:B:574:THR:HG22	1:B:598:THR:HG23	2.01	0.41
1:A:757:GLU:HG3	1:A:758:THR:N	2.34	0.41
1:B:628:ASP:HB2	1:B:630:ARG:HD3	2.02	0.41
1:B:707:LEU:HD23	1:B:735:VAL:CG1	2.50	0.41
1:A:155:GLN:HE21	1:A:463:PRO:CG	2.34	0.41
1:A:735:VAL:CA	1:A:736:SER:HB3	2.39	0.41
1:A:731:PHE:CD2	1:A:731:PHE:C	2.98	0.41
1:A:756:LEU:C	1:A:756:LEU:CD2	2.94	0.41
1:B:727:LEU:HD23	1:B:751:ILE:HG23	2.03	0.41
1:A:365:LYS:NZ	8:A:1015:HOH:O	2.53	0.41
1:A:776:CYS:HB3	1:A:777:THR:CG2	2.50	0.41
1:B:413:LEU:HD12	1:B:413:LEU:C	2.46	0.41
1:A:772:ASN:HB2	1:A:774:PHE:CE2	2.56	0.41
1:A:250:LEU:HD23	1:A:250:LEU:C	2.46	0.40
1:A:399:ILE:HG13	1:A:420:LEU:HD21	2.02	0.40
1:B:375:ARG:HD2	1:B:400:ASN:ND2	2.32	0.40
1:A:375:ARG:HD2	1:A:400:ASN:ND2	2.35	0.40
1:B:763:LYS:HD3	1:B:763:LYS:HA	1.90	0.40
1:A:753:LYS:HB3	1:A:756:LEU:HD13	2.03	0.40
1:A:798:LEU:HD23	1:A:798:LEU:HA	1.96	0.40
1:B:319:GLU:H	1:B:319:GLU:HG3	1.64	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	744/811 (92%)	692 (93%)	48 (6%)	4 (0%)	24	37
1	B	729/811 (90%)	686 (94%)	42 (6%)	1 (0%)	48	64
All	All	1473/1622 (91%)	1378 (94%)	90 (6%)	5 (0%)	36	50

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	168	ARG
1	A	736	SER
1	A	378	VAL
1	B	378	VAL
1	A	754	SER

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	697/755 (92%)	671 (96%)	26 (4%)	30	51
1	B	689/755 (91%)	662 (96%)	27 (4%)	28	48
All	All	1386/1510 (92%)	1333 (96%)	53 (4%)	29	49

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

Mol	Chain	Res	Type
1	A	45	VAL
1	A	49	CYS
1	A	52	ARG
1	A	81	GLU
1	A	87	GLN
1	A	90	THR
1	A	146	GLU
1	A	199	THR
1	A	212	SER
1	A	308	LYS
1	A	317	ASP

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Mol	Chain	Res	Type
1	A	384	GLU
1	A	388	GLN
1	A	392	GLN
1	A	471	THR
1	A	581	GLN
1	A	602	LYS
1	A	632	ILE
1	A	667	THR
1	A	675	MET
1	A	708	SER
1	A	729	SER
1	A	732	LEU
1	A	735	VAL
1	A	756	LEU
1	A	758	THR
1	B	49	CYS
1	B	52	ARG
1	B	80	ASN
1	B	81	GLU
1	B	84	GLN
1	B	90	THR
1	B	118	ASP
1	B	142	SER
1	B	199	THR
1	B	213	LEU
1	B	248	LEU
1	B	286	LEU
1	B	301	LYS
1	B	319	GLU
1	B	350	LYS
1	B	352	SER
1	B	416	ASN
1	B	418	SER
1	B	525	GLU
1	B	595	ASN
1	B	667	THR
1	B	675	MET
1	B	677	LYS
1	B	686	GLN
1	B	762	THR
1	B	780	ILE
1	B	798	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	GLN
1	A	96	HIS
1	A	155	GLN
1	A	160	ASN
1	A	191	ASN
1	A	285	ASN
1	A	415	GLN
1	A	469	HIS
1	A	566	HIS
1	A	581	GLN
1	B	96	HIS
1	B	157	ASN
1	B	191	ASN
1	B	312	HIS
1	B	499	ASN
1	B	581	GLN
1	B	582	ASN
1	B	585	ASN
1	B	604	ASN
1	B	686	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 ⓘ

21 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	1	1,2	14,14,15	0.93	1 (7%)	17,19,21	1.77	3 (17%)
2	NAG	C	2	2	14,14,15	0.58	0	17,19,21	1.50	3 (17%)
2	BMA	C	3	2	11,11,12	0.89	0	15,15,17	1.60	5 (33%)
2	MAN	C	4	2	11,11,12	0.76	0	15,15,17	1.38	3 (20%)
3	NAG	D	1	3,1	14,14,15	1.09	0	17,19,21	1.71	3 (17%)
3	NAG	D	2	3	14,14,15	1.11	0	17,19,21	1.60	6 (35%)
4	NAG	E	1	4,1	14,14,15	0.60	0	17,19,21	1.47	3 (17%)
4	NAG	E	2	4	14,14,15	1.01	1 (7%)	17,19,21	1.60	4 (23%)
4	BMA	E	3	4	11,11,12	0.63	0	15,15,17	1.51	3 (20%)
4	MAN	E	4	4	11,11,12	0.61	0	15,15,17	1.09	1 (6%)
4	MAN	E	5	4	11,11,12	0.64	0	15,15,17	1.59	3 (20%)
5	NAG	F	1	1,5	14,14,15	0.74	0	17,19,21	1.53	3 (17%)
5	NAG	F	2	5	14,14,15	1.06	1 (7%)	17,19,21	2.01	7 (41%)
5	BMA	F	3	5	11,11,12	0.90	0	15,15,17	1.15	2 (13%)
5	MAN	F	4	5	11,11,12	1.06	1 (9%)	15,15,17	2.20	4 (26%)
3	NAG	G	1	3,1	14,14,15	1.16	1 (7%)	17,19,21	1.65	4 (23%)
3	NAG	G	2	3	14,14,15	0.70	0	17,19,21	1.11	1 (5%)
5	NAG	H	1	1,5	14,14,15	0.85	0	17,19,21	1.38	2 (11%)
5	NAG	H	2	5	14,14,15	0.91	1 (7%)	17,19,21	1.59	5 (29%)
5	BMA	H	3	5	11,11,12	0.85	0	15,15,17	1.70	4 (26%)
5	MAN	H	4	5	11,11,12	0.67	0	15,15,17	1.37	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	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
2	MAN	C	4	2	-	2/2/19/22	0/1/1/1
3	NAG	D	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	2/6/23/26	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

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	4	MAN	C2-C3	3.22	1.57	1.52
5	F	2	NAG	O5-C1	-2.53	1.39	1.43
4	E	2	NAG	O5-C1	-2.42	1.39	1.43
2	C	1	NAG	C1-C2	-2.28	1.49	1.52
3	G	1	NAG	C2-N2	-2.13	1.42	1.46
5	H	2	NAG	O5-C1	-2.09	1.40	1.43

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	4	MAN	C1-C2-C3	5.23	117.27	109.64
2	C	1	NAG	O5-C1-C2	-4.57	104.22	111.29
5	H	4	MAN	O5-C5-C6	4.38	116.18	107.66
5	F	2	NAG	O5-C1-C2	-4.30	104.64	111.29
4	E	5	MAN	O5-C5-C6	3.59	114.66	107.66
5	F	1	NAG	O5-C1-C2	-3.54	105.81	111.29
4	E	2	NAG	C4-C3-C2	-3.50	105.89	111.02
5	H	3	BMA	C2-C3-C4	-3.37	104.93	110.86
3	D	2	NAG	O5-C5-C6	3.35	114.19	107.66
5	F	4	MAN	C2-C3-C4	3.33	116.72	110.86
5	F	4	MAN	O5-C5-C6	3.31	114.11	107.66
3	D	1	NAG	C1-C2-N2	-3.31	105.22	110.43
5	H	3	BMA	C1-C2-C3	3.27	114.41	109.64
5	F	2	NAG	O7-C7-C8	-3.23	116.30	122.05
5	H	2	NAG	C4-C3-C2	-3.14	106.42	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	1	NAG	C1-O5-C5	3.11	116.35	112.19
3	G	1	NAG	O3-C3-C4	3.09	117.66	110.38
3	G	1	NAG	C1-C2-N2	-3.06	105.61	110.43
4	E	1	NAG	O5-C1-C2	-3.04	106.58	111.29
2	C	1	NAG	O7-C7-C8	-3.00	116.72	122.05
2	C	2	NAG	O5-C1-C2	-2.91	106.78	111.29
2	C	3	BMA	C3-C4-C5	2.90	115.49	110.23
5	F	2	NAG	C1-C2-N2	2.88	114.97	110.43
5	H	2	NAG	O6-C6-C5	-2.87	101.55	111.33
2	C	3	BMA	O5-C5-C6	2.86	113.22	107.66
3	D	1	NAG	C1-O5-C5	-2.84	108.38	112.19
4	E	3	BMA	O3-C3-C2	2.84	115.86	110.05
3	D	1	NAG	C6-C5-C4	-2.77	106.22	113.02
5	F	2	NAG	C3-C4-C5	-2.75	105.25	110.23
5	H	3	BMA	O3-C3-C4	2.67	116.68	110.38
5	F	1	NAG	O3-C3-C2	2.64	114.89	109.40
2	C	4	MAN	C3-C4-C5	2.61	114.97	110.23
4	E	1	NAG	C8-C7-N2	2.61	120.44	116.12
5	H	2	NAG	C1-O5-C5	2.60	115.67	112.19
3	G	1	NAG	C4-C3-C2	-2.58	107.24	111.02
4	E	3	BMA	C1-O5-C5	-2.57	108.74	112.19
5	H	3	BMA	O4-C4-C5	-2.54	103.08	109.32
3	D	2	NAG	C1-O5-C5	2.51	115.56	112.19
3	D	2	NAG	O5-C1-C2	-2.47	107.46	111.29
2	C	2	NAG	C1-O5-C5	-2.41	108.95	112.19
4	E	2	NAG	C6-C5-C4	-2.41	107.11	113.02
3	G	2	NAG	O5-C1-C2	-2.38	107.61	111.29
3	G	1	NAG	O3-C3-C2	-2.37	104.48	109.40
2	C	3	BMA	C1-O5-C5	2.35	115.34	112.19
4	E	3	BMA	O4-C4-C3	-2.35	104.84	110.38
5	F	2	NAG	C1-O5-C5	-2.33	109.06	112.19
2	C	1	NAG	C4-C3-C2	2.33	114.43	111.02
3	D	2	NAG	C1-C2-N2	2.32	114.09	110.43
2	C	3	BMA	O4-C4-C3	-2.31	104.92	110.38
4	E	5	MAN	C1-O5-C5	-2.29	109.12	112.19
5	F	1	NAG	O7-C7-N2	2.25	125.96	121.98
5	H	1	NAG	C1-C2-N2	2.24	113.97	110.43
2	C	3	BMA	O6-C6-C5	2.24	118.97	111.33
2	C	2	NAG	O7-C7-C8	-2.24	118.06	122.05
4	E	4	MAN	O2-C2-C1	-2.22	104.14	109.22
5	H	2	NAG	O4-C4-C5	-2.21	103.88	109.32
5	F	2	NAG	C8-C7-N2	2.21	119.78	116.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4	MAN	C2-C3-C4	2.17	114.68	110.86
4	E	2	NAG	O4-C4-C5	-2.16	104.00	109.32
2	C	4	MAN	O3-C3-C4	-2.14	105.32	110.38
5	F	4	MAN	O2-C2-C3	2.14	114.58	110.15
4	E	1	NAG	O4-C4-C3	-2.12	105.38	110.38
5	H	2	NAG	C6-C5-C4	-2.11	107.83	113.02
5	F	3	BMA	O5-C1-C2	-2.11	105.77	110.79
3	D	2	NAG	C3-C4-C5	-2.10	106.42	110.23
4	E	5	MAN	C2-C3-C4	2.04	114.45	110.86
4	E	2	NAG	O5-C5-C4	-2.03	105.88	110.83
5	F	2	NAG	O5-C5-C6	2.03	111.62	107.66
5	F	3	BMA	C1-O5-C5	2.02	114.90	112.19
3	D	2	NAG	C4-C3-C2	-2.01	108.07	111.02

There are no chirality outliers.

All (20) torsion outliers are listed below:

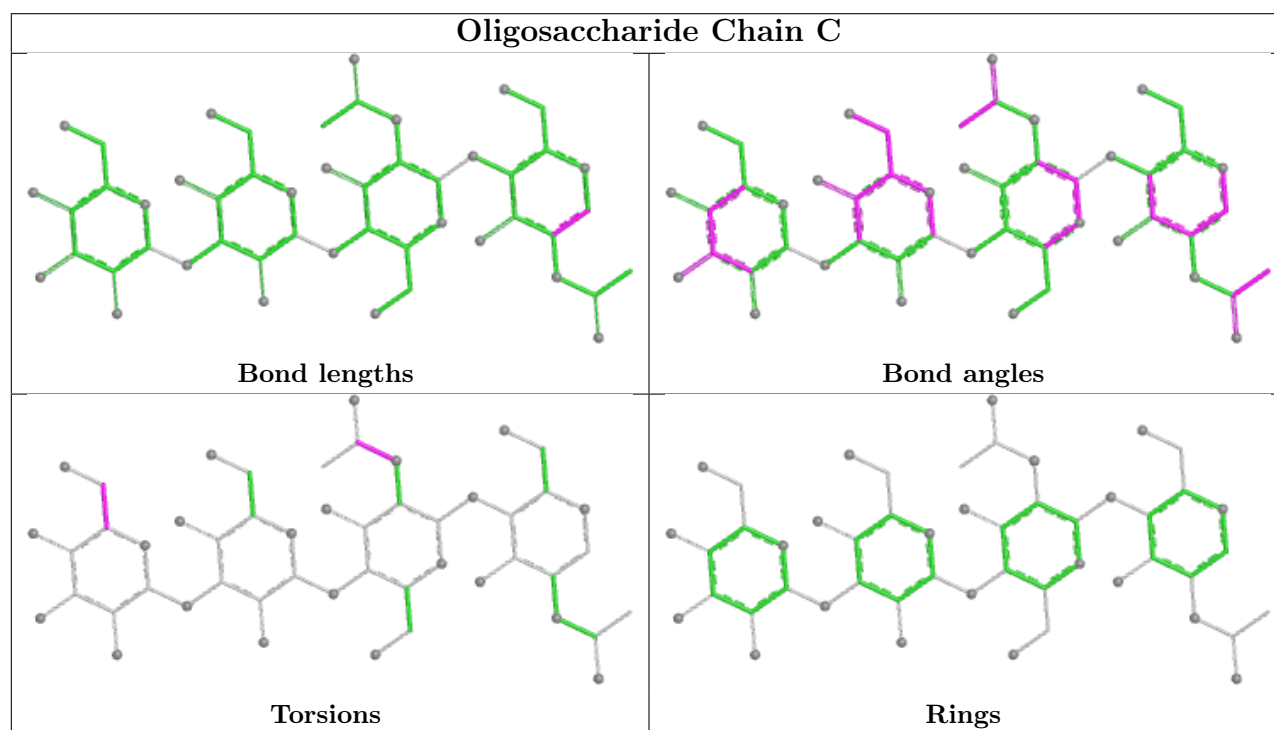
Mol	Chain	Res	Type	Atoms
2	C	4	MAN	O5-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
2	C	4	MAN	C4-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
5	F	2	NAG	C8-C7-N2-C2
5	F	2	NAG	O7-C7-N2-C2
5	F	4	MAN	O5-C5-C6-O6
4	E	4	MAN	C4-C5-C6-O6
5	F	2	NAG	C4-C5-C6-O6
5	H	4	MAN	O5-C5-C6-O6
4	E	5	MAN	O5-C5-C6-O6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

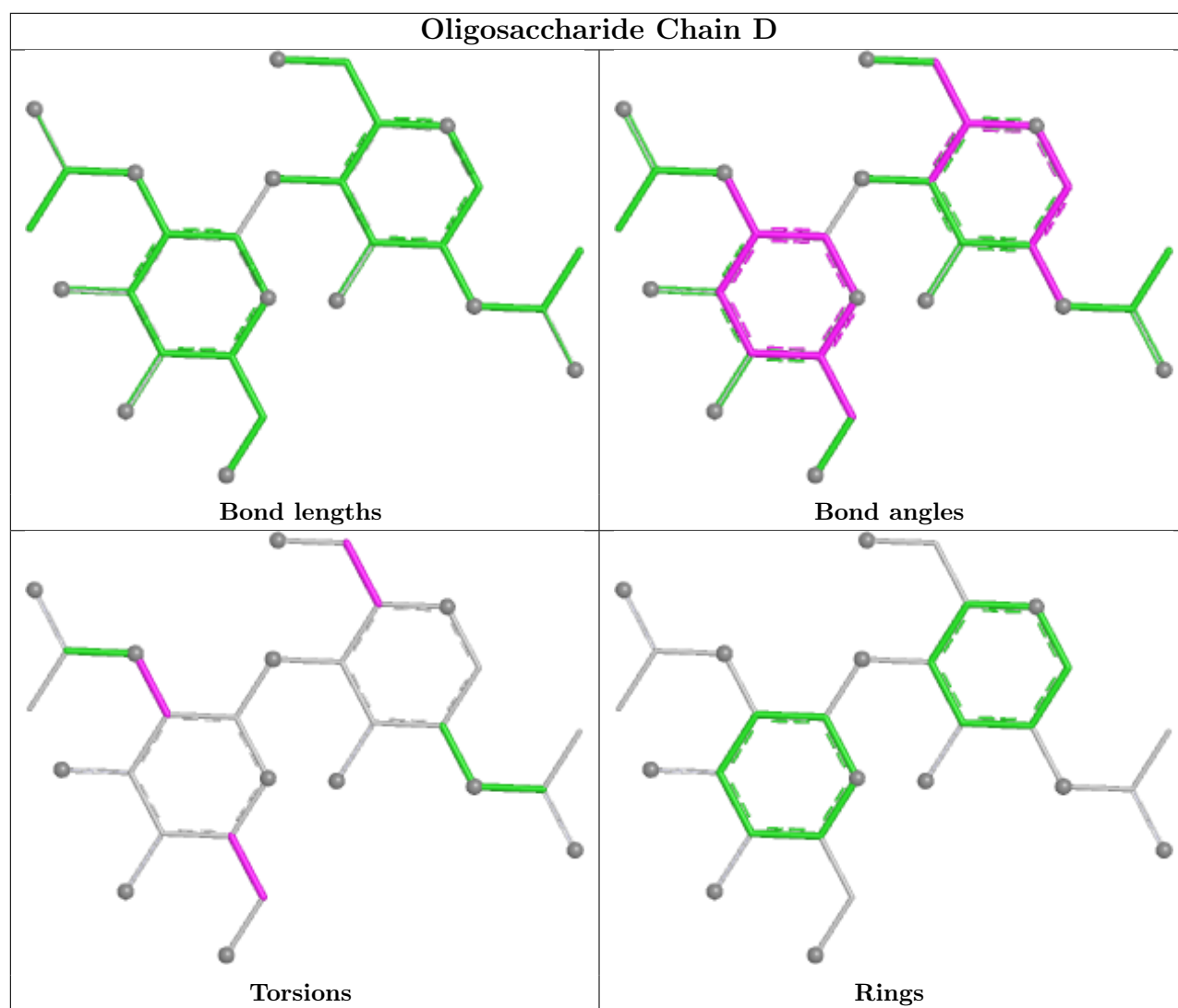
There are no ring outliers.

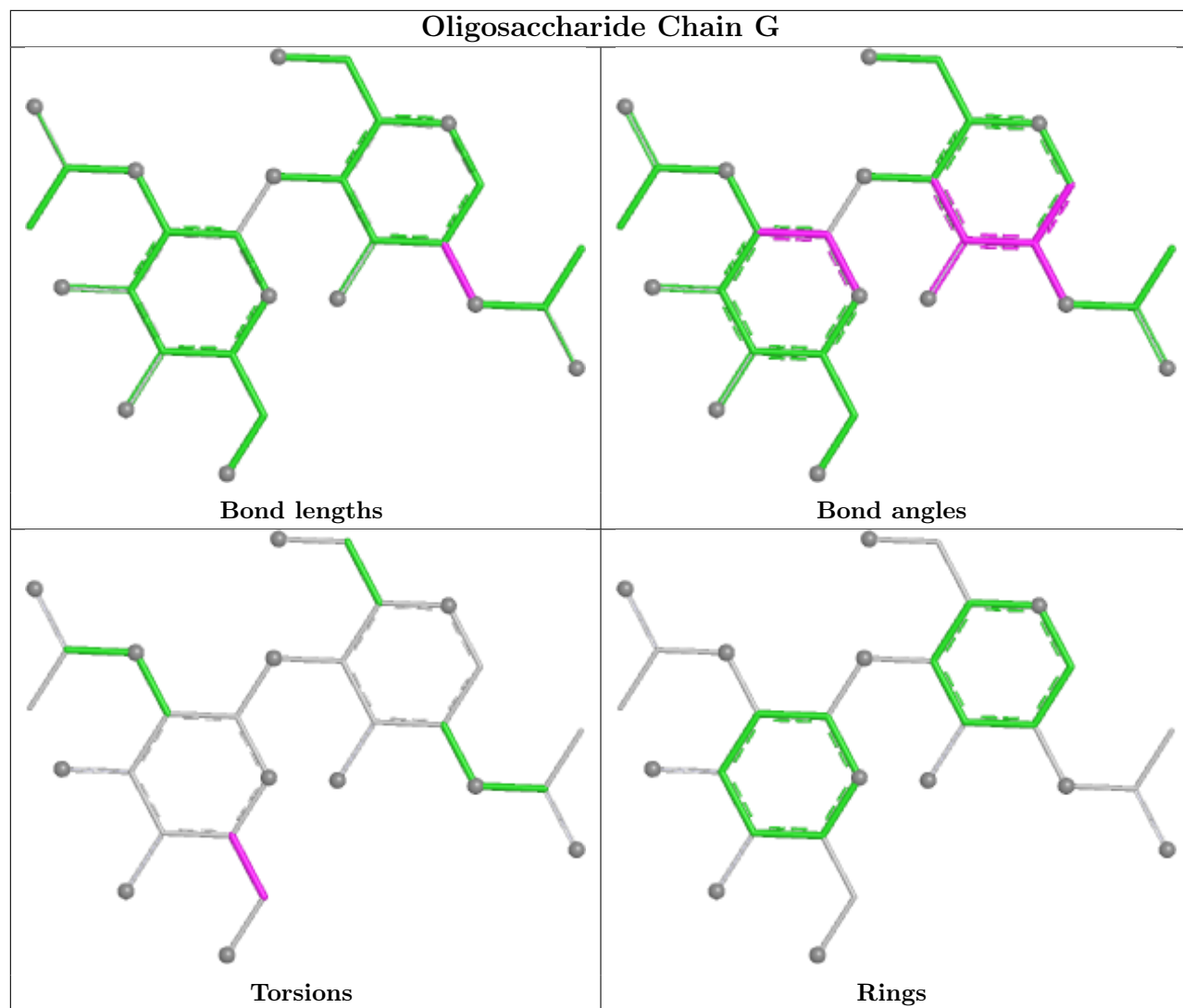
4 monomers are involved in 6 short contacts:

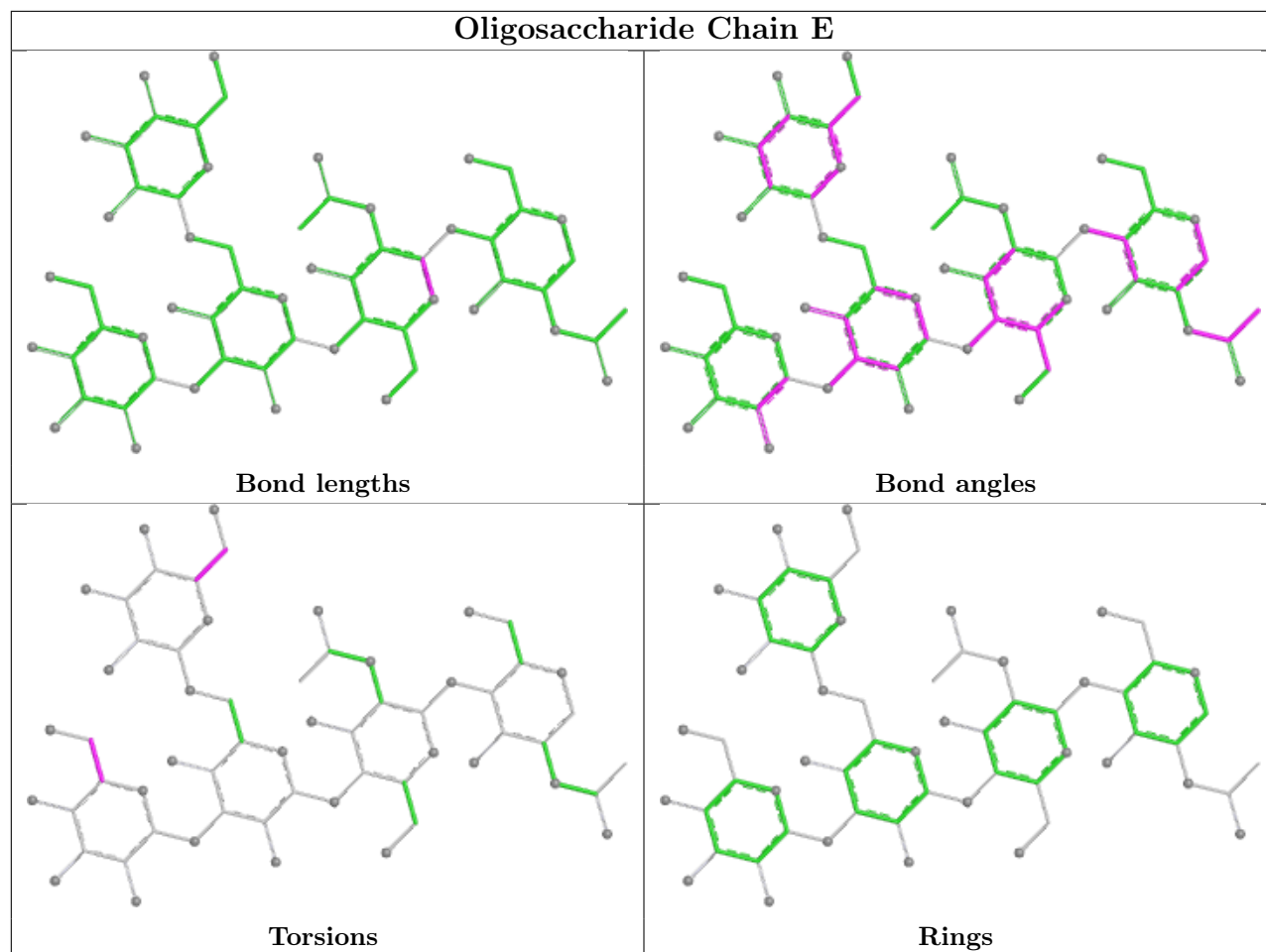
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	2	NAG	3	0
2	C	1	NAG	1	0
5	F	1	NAG	1	0
2	C	2	NAG	1	0

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

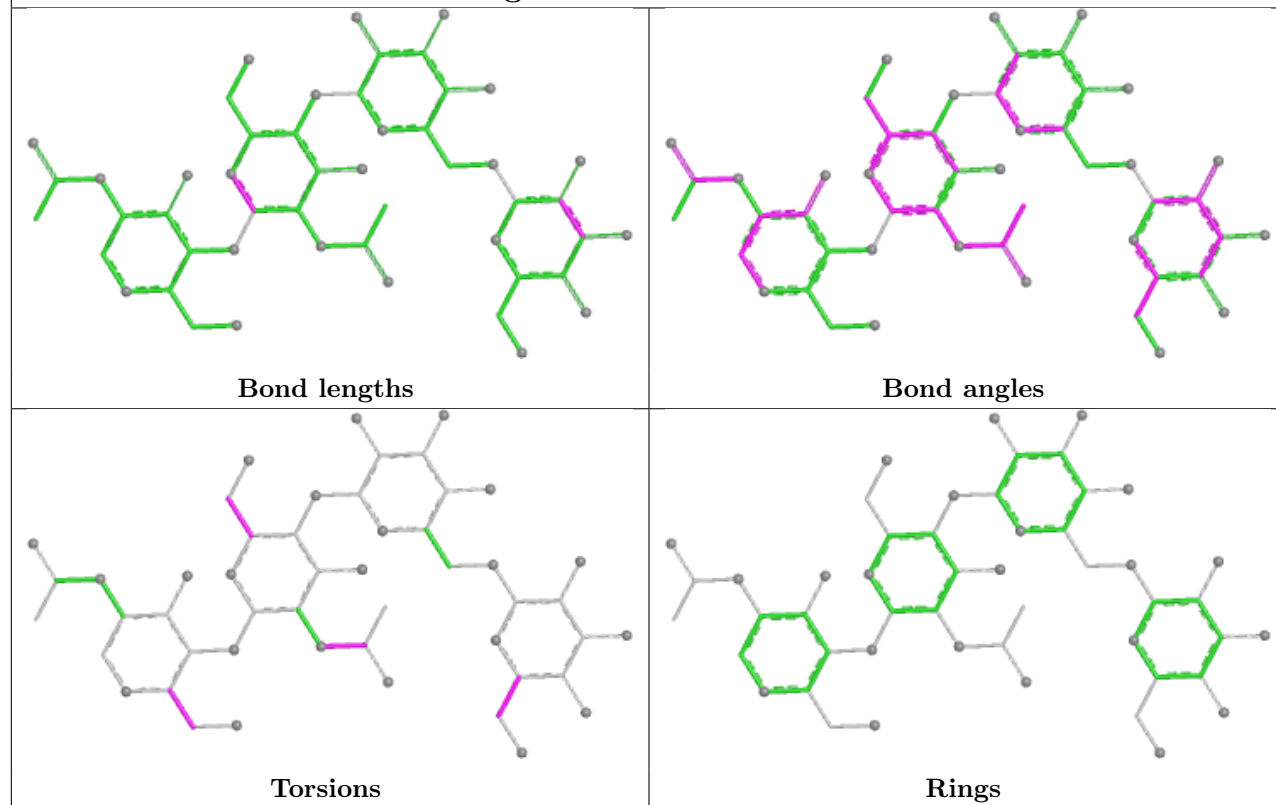




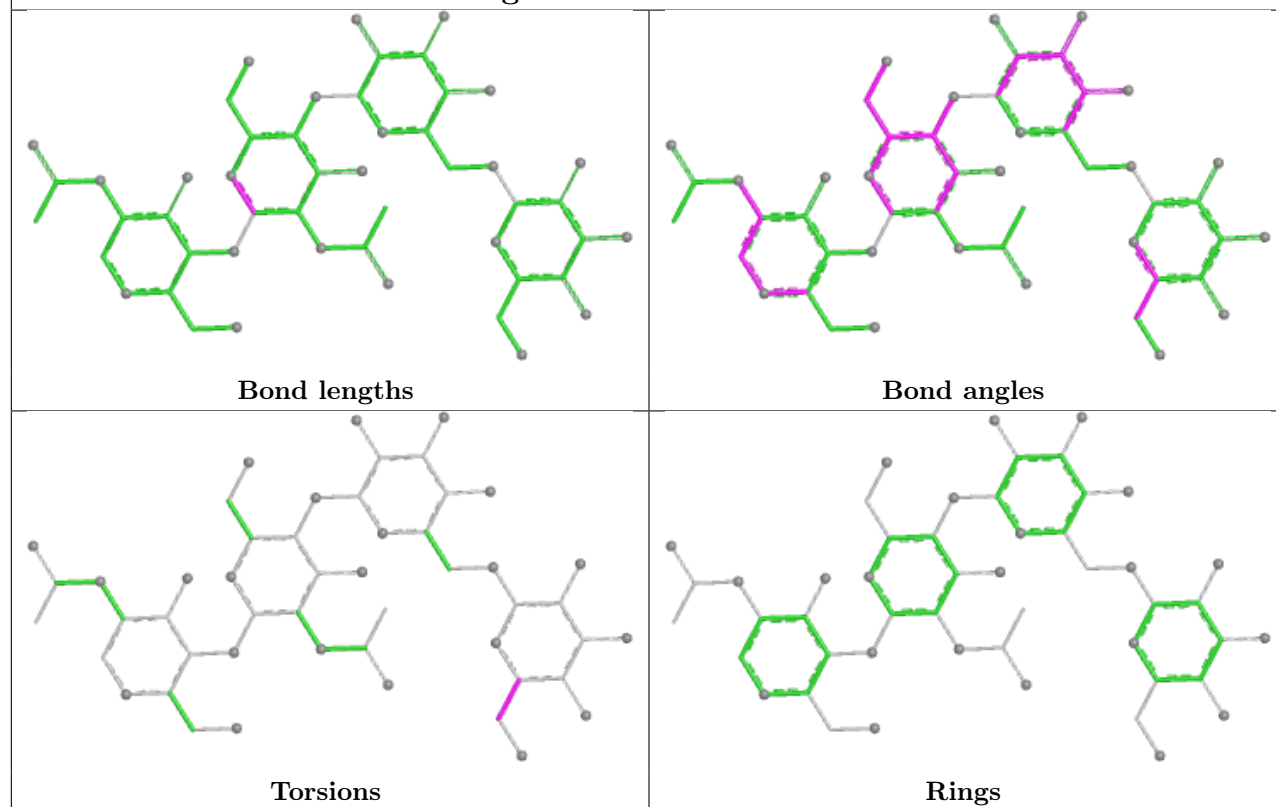




Oligosaccharide Chain F



Oligosaccharide Chain H



5.6 Ligand geometry

8 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
7	CU8	B	901	-	21,21,21	2.19	4 (19%)	29,30,30	3.14	10 (34%)
7	CU8	A	916	-	21,21,21	2.19	4 (19%)	29,30,30	3.31	12 (41%)
6	NAG	A	906	1	14,14,15	0.76	0	17,19,21	1.85	5 (29%)
6	NAG	B	913	1	14,14,15	0.93	1 (7%)	17,19,21	1.20	3 (17%)
6	NAG	A	914	1	14,14,15	0.82	0	17,19,21	1.68	3 (17%)
6	NAG	A	901	1	14,14,15	0.49	0	17,19,21	2.54	6 (35%)
6	NAG	A	915	1	14,14,15	0.66	0	17,19,21	1.59	3 (17%)
6	NAG	B	902	1	14,14,15	0.65	0	17,19,21	1.64	5 (29%)

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
7	CU8	B	901	-	-	0/7/8/8	0/3/3/3
7	CU8	A	916	-	-	0/7/8/8	0/3/3/3
6	NAG	A	906	1	-	2/6/23/26	0/1/1/1
6	NAG	B	913	1	-	0/6/23/26	0/1/1/1
6	NAG	A	914	1	-	0/6/23/26	0/1/1/1
6	NAG	A	901	1	-	2/6/23/26	0/1/1/1
6	NAG	A	915	1	-	2/6/23/26	0/1/1/1
6	NAG	B	902	1	-	2/6/23/26	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	901	CU8	C-N	-7.08	1.28	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	916	CU8	C-N	-7.01	1.28	1.38
7	B	901	CU8	C1-N	-3.73	1.32	1.38
7	A	916	CU8	C1-N	-3.65	1.32	1.38
7	A	916	CU8	C4-C5	-3.37	1.36	1.41
7	B	901	CU8	C4-C5	-3.21	1.37	1.41
7	A	916	CU8	O-C12	-2.92	1.17	1.24
6	B	913	NAG	C1-C2	2.87	1.56	1.52
7	B	901	CU8	O-C12	-2.66	1.18	1.24

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	901	CU8	C2-C3-N1	-10.13	116.73	124.75
7	A	916	CU8	C2-C3-N1	-9.88	116.93	124.75
6	A	901	NAG	C1-O5-C5	8.00	122.90	112.19
7	A	916	CU8	C5-C-N	-6.99	101.93	105.58
7	B	901	CU8	C5-C-N	-6.73	102.06	105.58
7	A	916	CU8	C4-C5-C	5.75	108.86	104.57
7	A	916	CU8	C4-N2-N	5.48	107.28	103.43
7	B	901	CU8	C3-N1-C	5.42	119.76	115.55
7	A	916	CU8	C3-N1-C	4.66	119.17	115.55
7	B	901	CU8	C4-N2-N	4.50	106.60	103.43
6	A	906	NAG	C1-O5-C5	4.46	118.16	112.19
6	A	915	NAG	C1-O5-C5	4.33	117.99	112.19
7	A	916	CU8	C5-C4-N2	-4.33	107.08	112.45
7	B	901	CU8	C4-C5-C	4.09	107.62	104.57
6	A	914	NAG	C1-O5-C5	4.04	117.60	112.19
7	A	916	CU8	N-C-N1	3.60	125.73	122.20
7	A	916	CU8	C1-N-C	-3.60	118.34	121.90
6	A	901	NAG	C8-C7-N2	3.59	122.07	116.12
6	A	906	NAG	C2-N2-C7	-3.57	118.12	122.90
7	B	901	CU8	C5-C4-N2	-3.53	108.06	112.45
6	A	906	NAG	C4-C3-C2	-3.37	106.09	111.02
7	B	901	CU8	C-N-N2	3.23	115.64	112.24
7	B	901	CU8	C2-C1-N	3.12	118.94	115.22
7	A	916	CU8	C2-C1-N	2.98	118.76	115.22
6	B	902	NAG	C1-C2-N2	-2.97	105.75	110.43
6	A	915	NAG	O5-C1-C2	-2.94	106.74	111.29
7	B	901	CU8	C1-N-C	-2.93	119.01	121.90
6	A	914	NAG	O5-C5-C6	-2.88	102.06	107.66
6	A	901	NAG	C4-C3-C2	-2.78	106.95	111.02
6	B	902	NAG	C4-C3-C2	2.78	115.09	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	901	CU8	N-C-N1	2.66	124.81	122.20
6	B	902	NAG	O5-C1-C2	2.66	115.40	111.29
6	A	915	NAG	C1-C2-N2	-2.55	106.42	110.43
6	B	902	NAG	C3-C4-C5	2.50	114.77	110.23
7	A	916	CU8	C-N-N2	2.48	114.85	112.24
6	B	913	NAG	O5-C5-C6	2.47	112.47	107.66
6	A	906	NAG	O3-C3-C2	-2.43	104.35	109.40
6	A	901	NAG	C2-N2-C7	2.22	125.88	122.90
7	A	916	CU8	C5-C12-N3	2.22	120.94	117.41
6	B	913	NAG	O7-C7-C8	-2.18	118.17	122.05
6	B	913	NAG	O5-C1-C2	-2.16	107.94	111.29
6	A	901	NAG	O5-C1-C2	2.10	114.53	111.29
6	A	914	NAG	O6-C6-C5	-2.07	104.27	111.33
7	A	916	CU8	C9-C8-C7	2.04	120.96	117.96
6	A	901	NAG	O7-C7-C8	-2.03	118.44	122.05
6	B	902	NAG	O5-C5-C6	2.02	111.58	107.66
6	A	906	NAG	O5-C1-C2	-2.01	108.18	111.29

There are no chirality outliers.

All (8) torsion outliers are listed below:

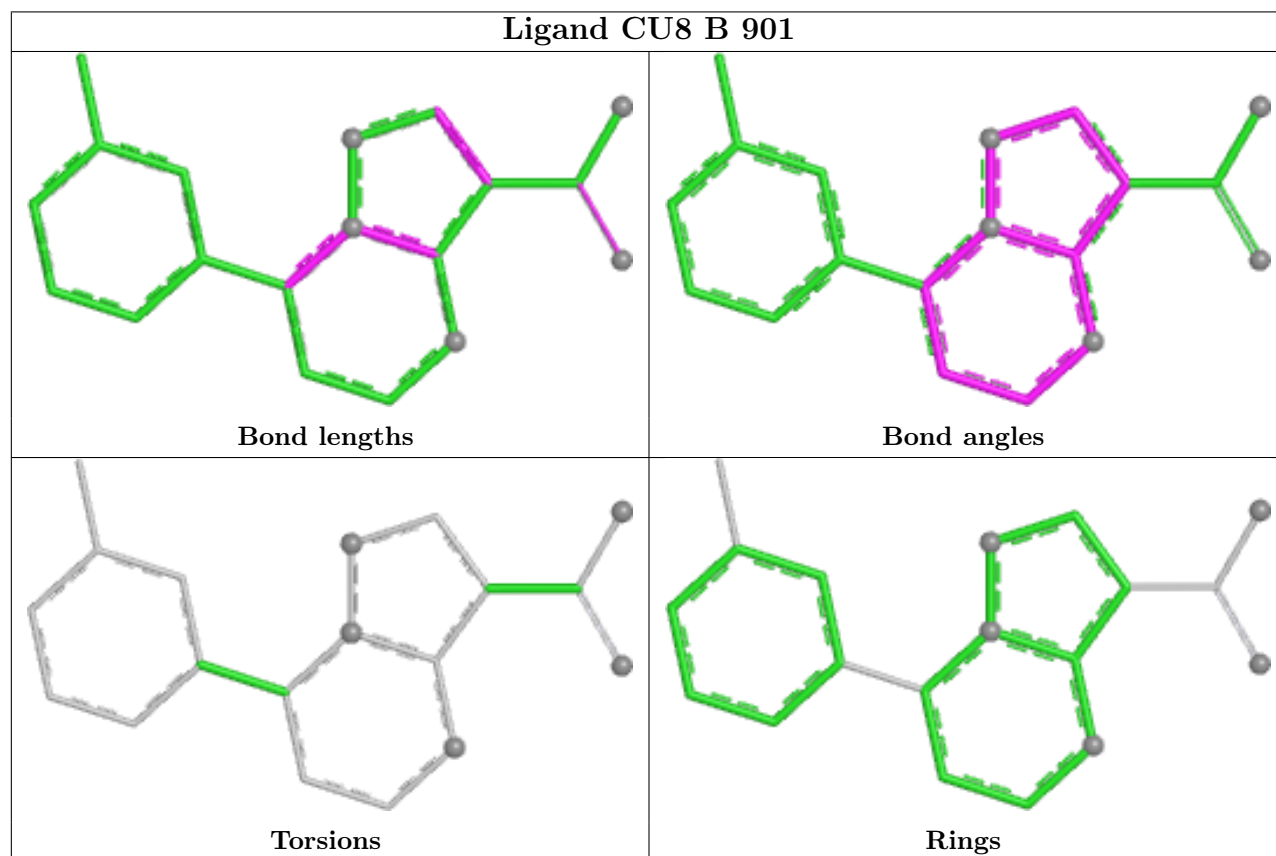
Mol	Chain	Res	Type	Atoms
6	B	902	NAG	O5-C5-C6-O6
6	A	915	NAG	O5-C5-C6-O6
6	A	915	NAG	C4-C5-C6-O6
6	B	902	NAG	C4-C5-C6-O6
6	A	906	NAG	O5-C5-C6-O6
6	A	906	NAG	C4-C5-C6-O6
6	A	901	NAG	C8-C7-N2-C2
6	A	901	NAG	O7-C7-N2-C2

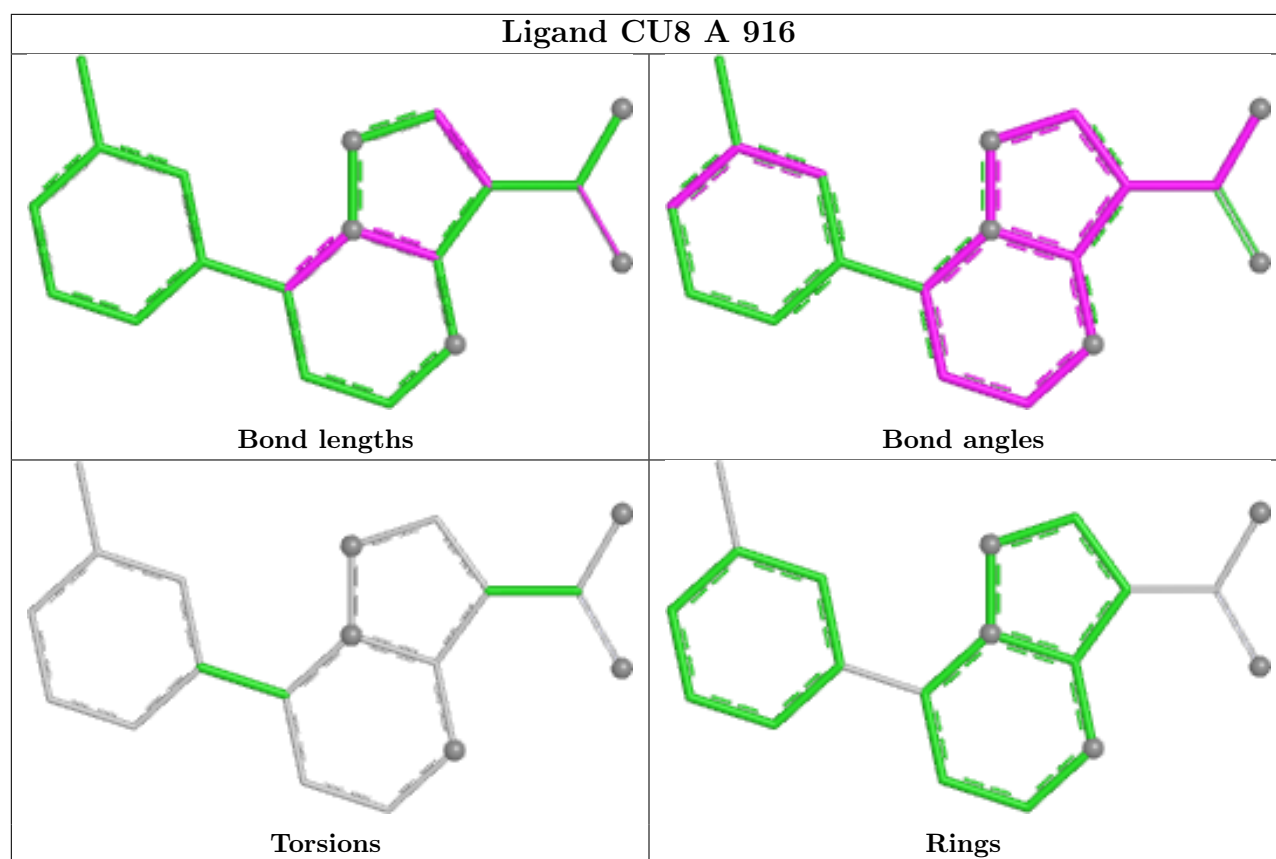
There are no ring outliers.

No monomer is involved in short contacts.

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.





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	750/811 (92%)	0.14	21 (2%) 55 51	29, 57, 103, 151	0
1	B	739/811 (91%)	0.11	17 (2%) 61 57	30, 55, 98, 134	0
All	All	1489/1622 (91%)	0.12	38 (2%) 57 53	29, 56, 99, 151	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	819	LEU	5.3
1	A	265	PHE	5.0
1	A	100	VAL	4.8
1	A	182	TYR	4.8
1	A	735	VAL	4.2
1	B	461	PHE	4.2
1	A	461	PHE	4.1
1	A	756	LEU	3.9
1	A	470	PHE	3.5
1	A	186	VAL	3.4
1	B	762	THR	3.4
1	B	435	LYS	3.2
1	A	702	PHE	3.1
1	B	182	TYR	2.9
1	A	183	PHE	2.9
1	B	64	TYR	2.9
1	B	268	VAL	2.7
1	A	187	CYS	2.6
1	A	678	PHE	2.6
1	A	215	HIS	2.6
1	B	567	TYR	2.6
1	A	758	THR	2.5
1	A	736	SER	2.5
1	A	570	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	470	PHE	2.4
1	B	61	VAL	2.4
1	B	186	VAL	2.4
1	B	179	TRP	2.4
1	A	732	LEU	2.3
1	B	88	ASN	2.3
1	A	568	PHE	2.2
1	B	65	VAL	2.2
1	B	735	VAL	2.2
1	B	597	TYR	2.1
1	B	753	LYS	2.1
1	A	181	CYS	2.1
1	A	567	TYR	2.0
1	A	571	ALA	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.

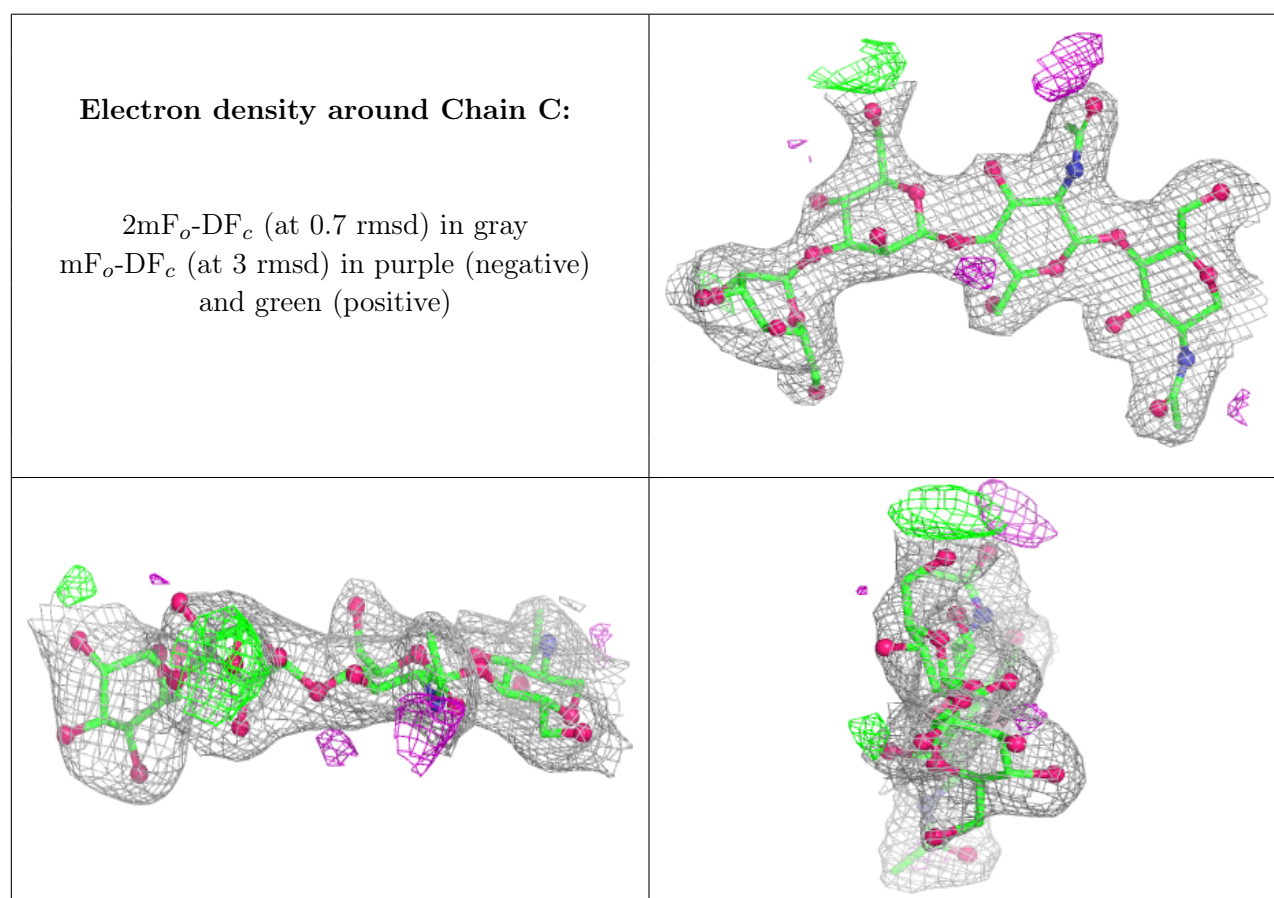
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MAN	H	4	11/12	0.56	0.15	73,96,113,114	0
5	MAN	F	4	11/12	0.64	0.14	79,91,97,98	0
4	MAN	E	4	11/12	0.69	0.11	87,95,108,110	0
2	MAN	C	4	11/12	0.75	0.12	74,84,88,91	0
4	MAN	E	5	11/12	0.76	0.10	79,84,89,96	0
5	BMA	F	3	11/12	0.78	0.12	63,68,84,85	0
5	BMA	H	3	11/12	0.83	0.09	53,59,80,103	0
2	BMA	C	3	11/12	0.84	0.10	55,67,79,82	0
3	NAG	D	2	14/15	0.84	0.10	49,62,79,81	0
3	NAG	G	2	14/15	0.86	0.11	55,71,82,95	0
4	BMA	E	3	11/12	0.90	0.07	49,59,69,73	0
2	NAG	C	2	14/15	0.93	0.07	38,46,52,61	0
5	NAG	F	2	14/15	0.94	0.07	33,45,58,64	0
5	NAG	H	2	14/15	0.94	0.07	36,41,45,50	0

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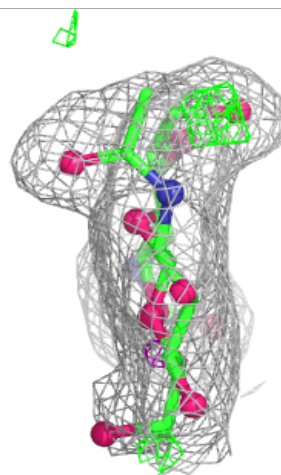
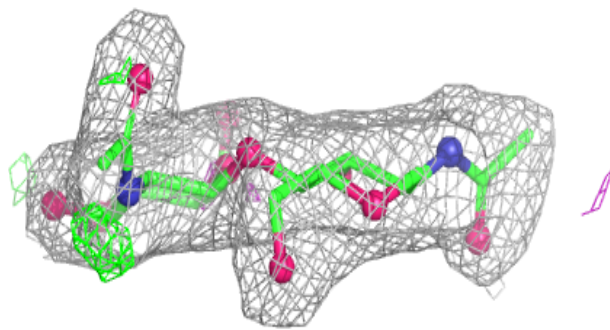
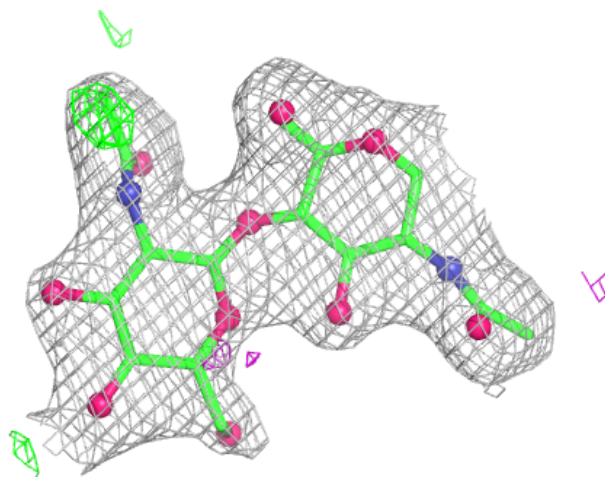
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	C	1	14/15	0.95	0.07	35,40,47,47	0
5	NAG	H	1	14/15	0.96	0.07	33,36,37,38	0
4	NAG	E	2	14/15	0.96	0.06	31,37,46,53	0
4	NAG	E	1	14/15	0.96	0.06	27,35,38,38	0
5	NAG	F	1	14/15	0.96	0.06	33,37,46,55	0
3	NAG	D	1	14/15	0.97	0.05	31,36,41,41	0
3	NAG	G	1	14/15	0.97	0.06	34,38,46,51	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.



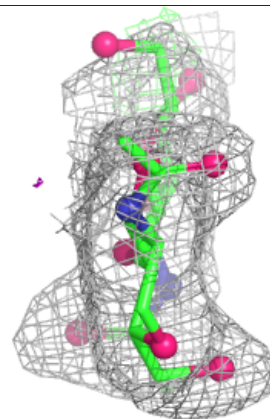
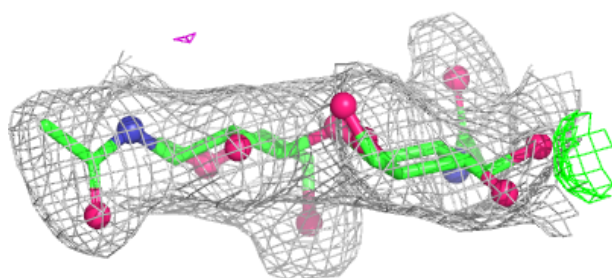
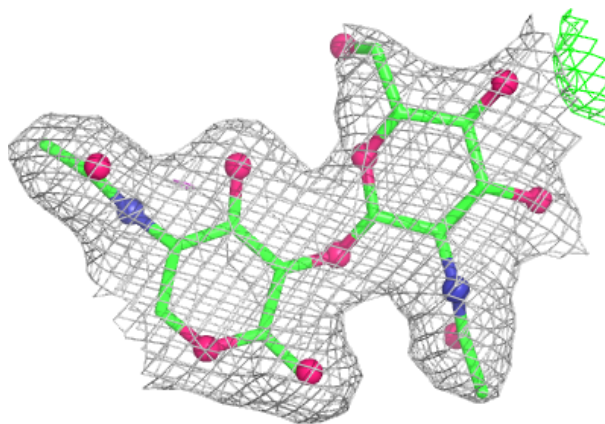
Electron density around Chain D:

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



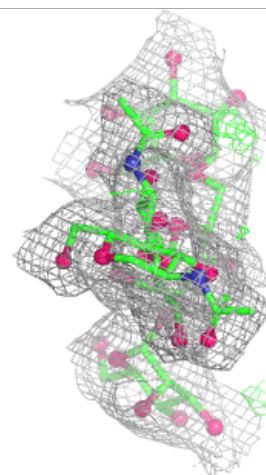
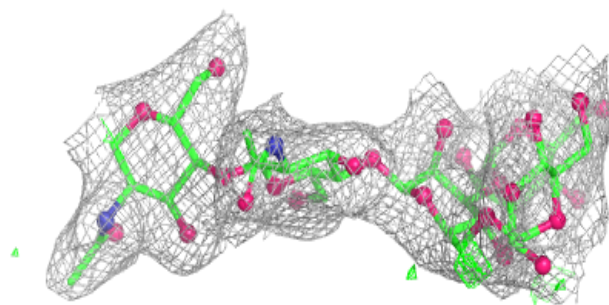
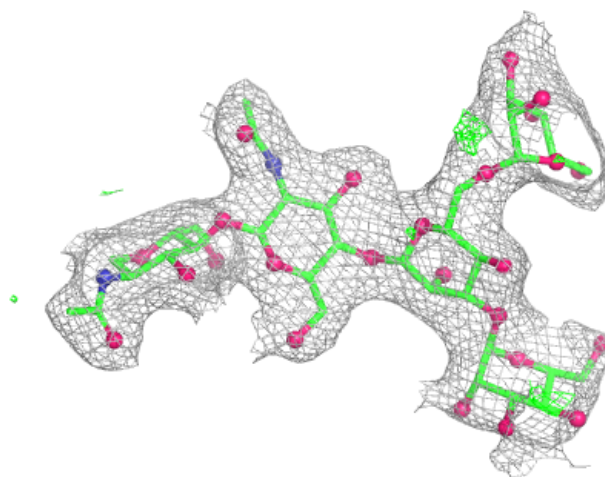
Electron density around Chain G:

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



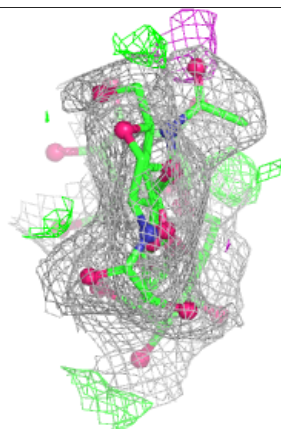
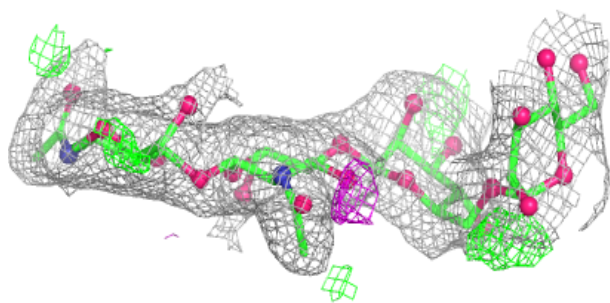
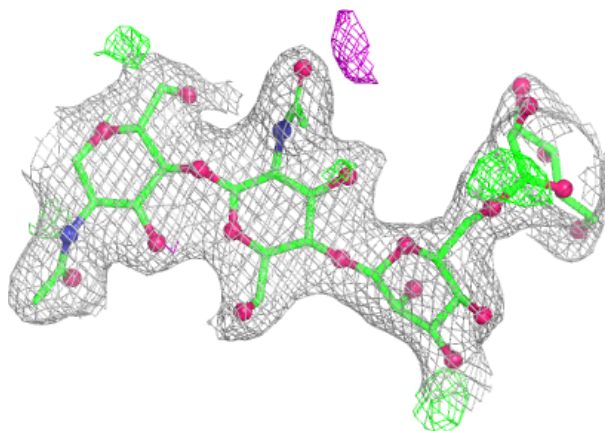
Electron density around Chain E:

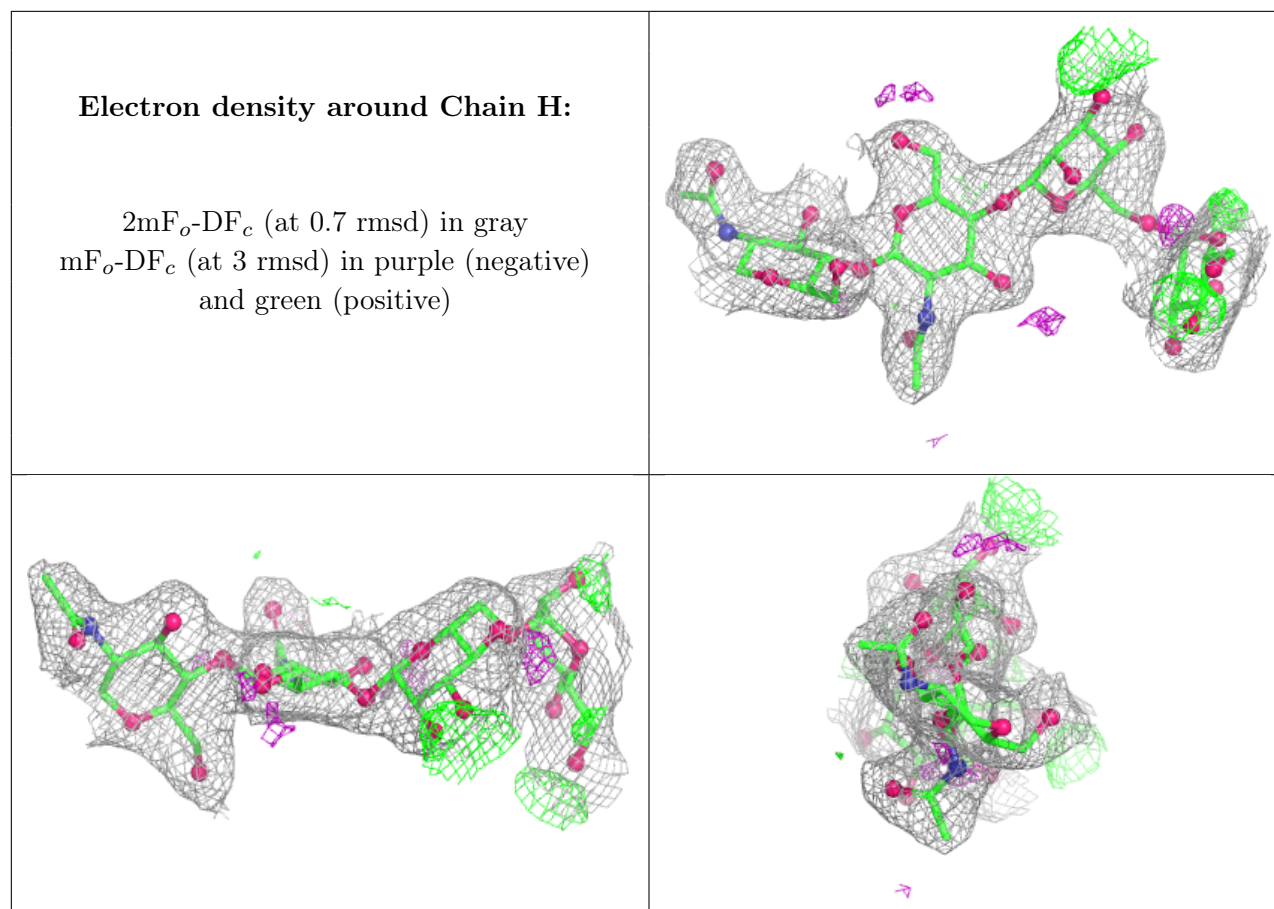
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.4 Ligands [i](#)

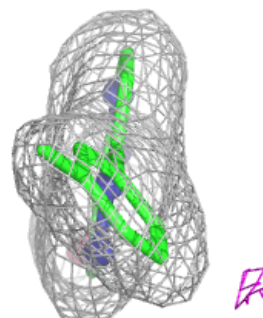
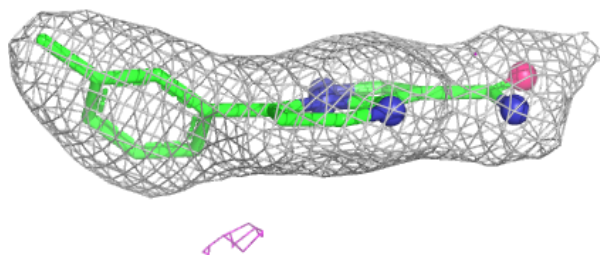
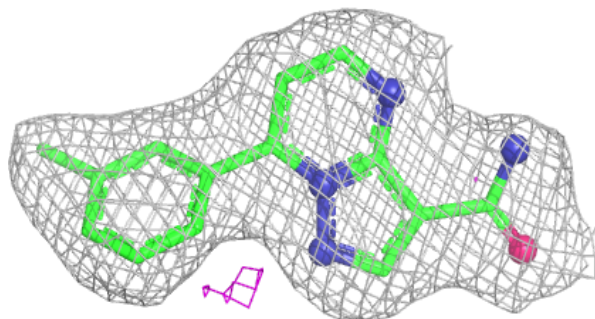
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	915	14/15	0.62	0.15	76,90,106,106	0
6	NAG	B	902	14/15	0.71	0.13	79,87,98,100	0
6	NAG	A	901	14/15	0.77	0.12	76,85,90,92	0
6	NAG	B	913	14/15	0.90	0.10	53,61,71,72	0
6	NAG	A	906	14/15	0.93	0.07	52,62,69,72	0
6	NAG	A	914	14/15	0.96	0.06	40,44,57,61	0
7	CU8	A	916	19/19	0.96	0.07	33,39,51,53	0
7	CU8	B	901	19/19	0.97	0.07	36,42,49,53	0

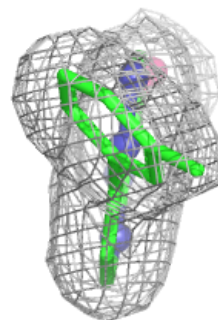
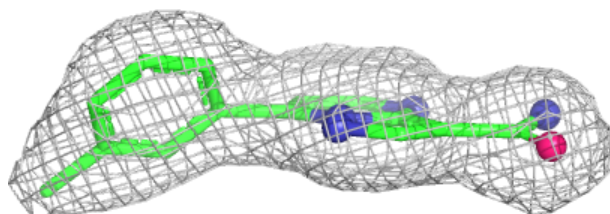
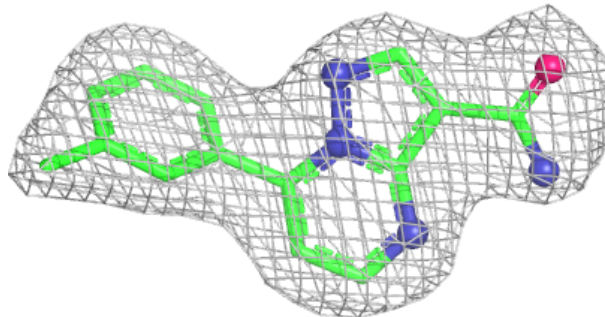
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CU8 A 916:

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

**Electron density around CU8 B 901:**

$2mF_o-DF_c$ (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.