



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2020 – 08:08 PM BST

PDB ID : 6TZB
Title : Crystal structure of the A/Hong Kong/1/1968 (H3N2) influenza virus hemagglutinin in complex with 6'-SLNLN
Authors : Wu, N.C.; Wilson, I.A.
Deposited on : 2019-08-12
Resolution : 2.24 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

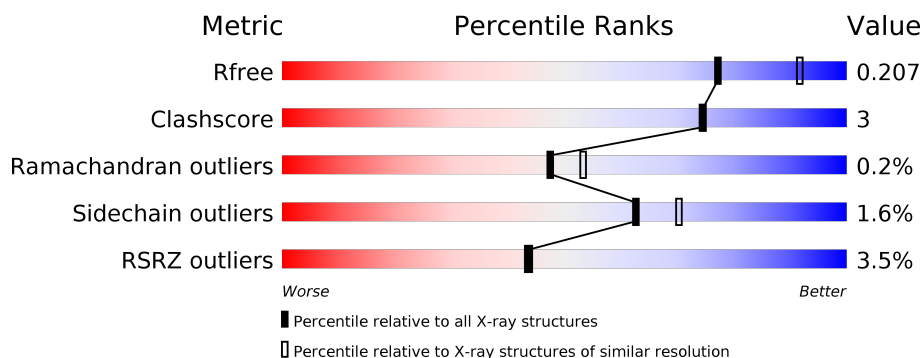
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 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}	130704	2391 (2.26-2.22)
Clashscore	141614	2539 (2.26-2.22)
Ramachandran outliers	138981	2489 (2.26-2.22)
Sidechain outliers	138945	2490 (2.26-2.22)
RSRZ outliers	127900	2353 (2.26-2.22)

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 on 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	321	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>7%</div> </div> <div>•</div> </div>
1	C	321	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>6%</div> </div> <div>••</div> </div>
1	E	321	<div> <div>0%</div> <div> <div></div> <div>89%</div> <div>9%</div> </div> <div>••</div> </div>
2	B	176	<div> <div>5%</div> <div> <div></div> <div>91%</div> <div>6%</div> </div> <div>••</div> </div>
2	D	176	<div> <div>7%</div> <div> <div></div> <div>90%</div> <div>7%</div> </div> <div>•</div> </div>
2	F	176	<div> <div>9%</div> <div> <div></div> <div>92%</div> <div>5%</div> </div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
3	G	3	 100%
3	H	3	 33% 67%
3	L	3	 100%
4	I	2	 100%
4	N	2	 50% 50%
4	P	2	 50% 50%
5	J	5	 20% 80%
5	Q	5	 20% 80%
6	K	4	 25% 75%
7	M	3	 33% 67%
8	O	4	 25% 75%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	G	2	-	-	-	X
3	NAG	H	2	-	-	-	X
3	BMA	H	3	-	-	-	X
3	BMA	L	3	-	-	-	X
4	NAG	N	2	-	-	-	X
5	NAG	J	1	-	-	-	X
5	GAL	J	2	-	-	-	X
5	NAG	Q	1	-	-	-	X
6	MAN	K	4	-	-	-	X
8	MAN	O	4	-	-	-	X
9	NAG	D	201	-	-	-	X

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 12998 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 Hemagglutinin HA1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2479	1552	438	476	13			
1	C	317	Total	C	N	O	S	0	0	0
			2443	1531	429	470	13			
1	E	317	Total	C	N	O	S	0	0	0
			2443	1531	429	470	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	PRO	-	expression tag	UNP Q91MA7
A	10	GLY	-	expression tag	UNP Q91MA7
C	9	PRO	-	expression tag	UNP Q91MA7
C	10	GLY	-	expression tag	UNP Q91MA7
E	9	PRO	-	expression tag	UNP Q91MA7
E	10	GLY	-	expression tag	UNP Q91MA7

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	172	Total	C	N	O	S	0	0	0
			1391	863	243	279	6			
2	D	171	Total	C	N	O	S	0	0	0
			1382	858	241	277	6			
2	F	171	Total	C	N	O	S	0	0	0
			1382	858	241	277	6			

There are 3 discrepancies between the modelled and reference sequences:

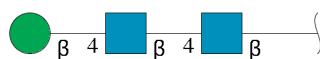
Chain	Residue	Modelled	Actual	Comment	Reference
B	123	GLY	ARG	conflict	UNP H9XC94
D	123	GLY	ARG	conflict	UNP H9XC94

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Chain	Residue	Modelled	Actual	Comment	Reference
F	123	GLY	ARG	conflict	UNP H9XC94

- 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.



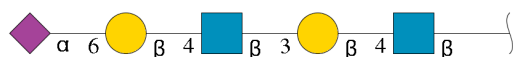
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	3	Total	C	N	O	0	0	0
			39	22	2	15			
3	H	3	Total	C	N	O	0	0	0
			39	22	2	15			
3	L	3	Total	C	N	O	0	0	0
			39	22	2	15			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	N	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	P	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



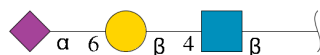
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	J	5	Total	C	N	O	0	0	0
			71	39	3	29			
5	Q	5	Total	C	N	O	0	0	0
			71	39	3	29			

- Molecule 6 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
6	K	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 7 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	M	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 8 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
8	O	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by author).



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

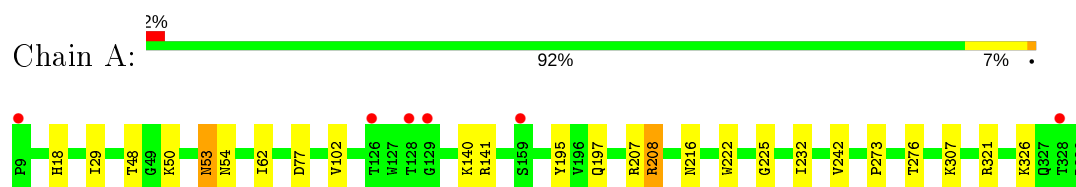
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	175	Total	O	0	0
			175	175		
10	B	97	Total	O	0	0
			97	97		
10	C	183	Total	O	0	0
			183	183		
10	D	111	Total	O	0	0
			111	111		
10	E	264	Total	O	0	0
			264	264		
10	F	89	Total	O	0	0
			89	89		

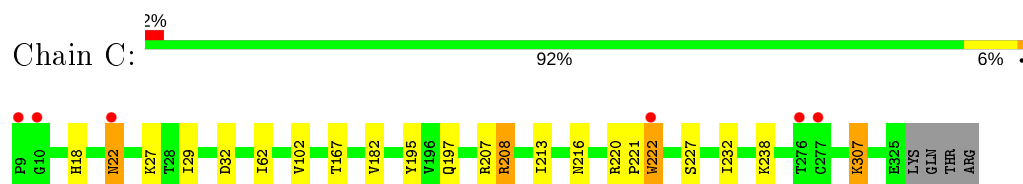
3 Residue-property plots [i](#)

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.

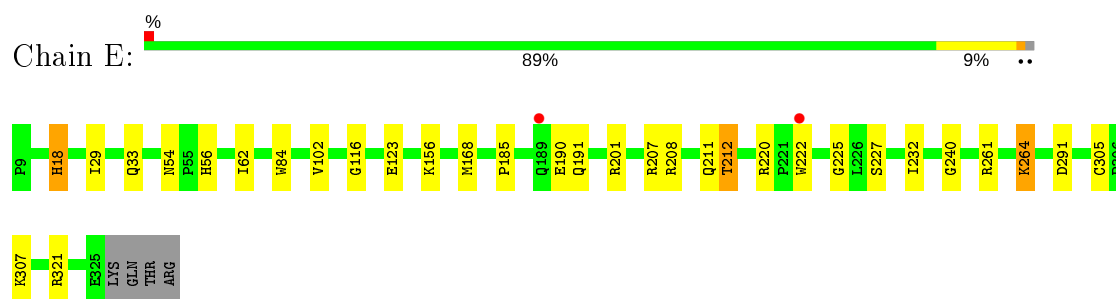
- Molecule 1: Hemagglutinin HA1 chain



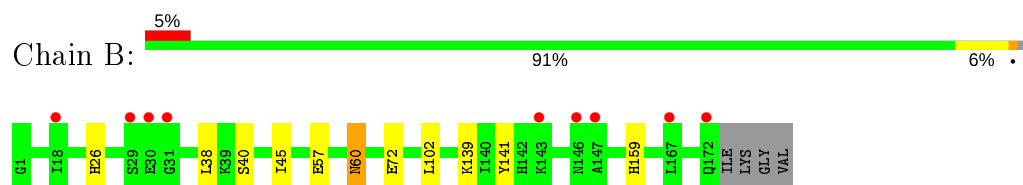
- Molecule 1: Hemagglutinin HA1 chain



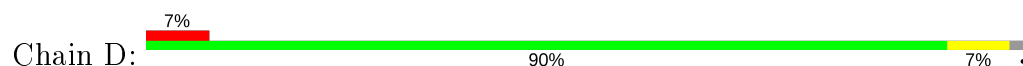
- Molecule 1: Hemagglutinin HA1 chain



- Molecule 2: Hemagglutinin HA2 chain

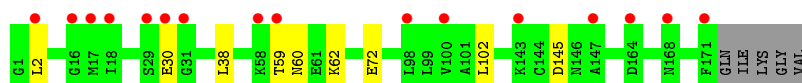
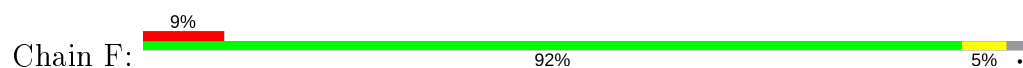


- Molecule 2: Hemagglutinin HA2 chain





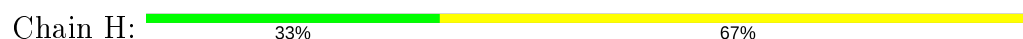
- Molecule 2: Hemagglutinin HA2 chain



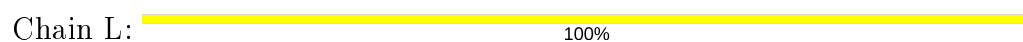
- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



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



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



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



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

Chain P:  50% 50%



- Molecule 5: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  20% 80%



- Molecule 5: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  20% 80%



- Molecule 6: 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 K:  25% 75%



- Molecule 7: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  33% 67%



- Molecule 8: 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 O:  25% 75%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	210.40 Å 132.17 Å 72.47 Å 90.00° 97.93° 90.00°	Depositor
Resolution (Å)	39.75 – 2.24 39.75 – 2.24	Depositor EDS
% Data completeness (in resolution range)	99.5 (39.75-2.24) 99.5 (39.75-2.24)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 2.24 Å)	Xtriage
Refinement program	PHENIX 1.12 _2829	Depositor
R, R_{free}	0.157 , 0.207 0.157 , 0.207	Depositor DCC
R_{free} test set	4625 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	37.2	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	12998	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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: GAL, SIA, BMA, NAG, 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	0.42	0/2536	0.60	1/3454 (0.0%)
1	C	0.40	0/2500	0.58	0/3407
1	E	0.44	0/2500	0.62	0/3407
2	B	0.40	0/1415	0.53	0/1902
2	D	0.40	0/1406	0.55	0/1890
2	F	0.42	0/1406	0.56	0/1890
All	All	0.42	0/11763	0.58	1/15950 (0.0%)

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
2	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	321	ARG	CG-CD-NE	-6.03	99.14	111.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	57	GLU	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	2479	0	2431	14	0
1	C	2443	0	2390	17	0
1	E	2443	0	2390	27	0
2	B	1391	0	1307	10	0
2	D	1382	0	1298	10	0
2	F	1382	0	1299	10	0
3	G	39	0	34	0	0
3	H	39	0	34	1	0
3	L	39	0	34	1	0
4	I	28	0	25	0	0
4	N	28	0	25	0	0
4	P	28	0	25	0	0
5	J	71	0	61	0	0
5	Q	71	0	61	1	0
6	K	50	0	43	1	0
7	M	46	0	40	0	0
8	O	50	0	43	0	0
9	A	14	0	13	0	0
9	C	28	0	26	0	0
9	D	14	0	13	0	0
9	E	14	0	13	0	0
10	A	175	0	0	1	0
10	B	97	0	0	4	0
10	C	183	0	0	3	0
10	D	111	0	0	2	0
10	E	264	0	0	10	0
10	F	89	0	0	0	0
All	All	12998	0	11605	75	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 3.

All (75) 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:E:307:LYS:NZ	10:E:601:HOH:O	1.90	1.03
2:B:60:ASN:H	2:B:60:ASN:HD22	1.31	0.79
2:D:60:ASN:HD22	2:D:62:LYS:HE2	1.50	0.75
2:F:60:ASN:HD22	2:F:62:LYS:HE2	1.55	0.71
2:D:45:ILE:HG22	10:D:301:HOH:O	1.91	0.70
1:A:77:ASP:OD2	1:A:141:ARG:NH1	2.26	0.68
1:C:307:LYS:NZ	10:C:602:HOH:O	2.25	0.68
1:E:307:LYS:NZ	2:F:60:ASN:HD21	1.92	0.68
1:E:307:LYS:HE2	1:E:307:LYS:HA	1.78	0.66
1:A:208:ARG:NH2	2:D:72:GLU:OE2	2.29	0.66
1:C:22:ASN:O	1:C:22:ASN:ND2	2.27	0.65
1:E:291:ASP:OD2	10:E:604:HOH:O	2.14	0.65
2:D:46:ASP:OD1	10:D:301:HOH:O	2.15	0.65
1:A:216:ASN:ND2	10:A:502:HOH:O	2.29	0.64
1:C:167:THR:OG1	6:K:1:NAG:H62	1.98	0.62
1:E:201:ARG:NH1	10:E:602:HOH:O	2.08	0.61
1:C:208:ARG:NH2	2:F:72:GLU:OE2	2.34	0.61
1:E:211:GLN:NE2	10:E:603:HOH:O	2.08	0.59
2:D:21:TRP:CE2	2:D:45:ILE:HD11	2.37	0.59
1:E:261:ARG:NE	10:E:609:HOH:O	2.29	0.58
2:D:54:ARG:O	2:D:57:GLU:HB2	2.04	0.57
1:E:56:HIS:ND1	10:E:611:HOH:O	2.33	0.56
1:C:207:ARG:HD2	10:C:614:HOH:O	2.07	0.54
1:E:185:PRO:HB3	1:E:190:GLU:HG3	1.88	0.54
1:A:207:ARG:HG2	1:C:221:PRO:HB2	1.90	0.53
1:C:216:ASN:HB2	1:C:220:ARG:HH22	1.74	0.53
2:B:159:HIS:ND1	10:B:204:HOH:O	2.34	0.52
2:B:60:ASN:H	2:B:60:ASN:ND2	2.03	0.52
1:C:222:TRP:CZ3	1:C:227:SER:HB3	2.45	0.51
2:B:139:LYS:HD2	2:B:141:TYR:CZ	2.46	0.51
1:E:29:ILE:HD11	2:F:102:LEU:HD23	1.93	0.51
1:A:53:ASN:HD22	1:A:53:ASN:C	2.13	0.50
2:B:26:HIS:HD2	10:B:209:HOH:O	1.95	0.49
1:A:77:ASP:CG	1:A:141:ARG:HH12	2.15	0.49
1:E:321:ARG:NH1	10:E:619:HOH:O	2.44	0.49
1:E:264:LYS:NZ	10:E:623:HOH:O	2.46	0.48
1:C:222:TRP:CE3	3:H:2:NAG:H2	2.47	0.48
2:B:72:GLU:HG3	10:B:287:HOH:O	2.12	0.48
1:E:220:ARG:HB2	1:E:227:SER:O	2.14	0.47
1:C:27:LYS:HG2	1:C:32:ASP:O	2.15	0.47
1:C:29:ILE:HD11	2:D:102:LEU:HD23	1.96	0.47
1:C:102:VAL:HG22	1:C:232:ILE:HB	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:307:LYS:HZ1	2:F:60:ASN:HD21	1.60	0.46
1:E:54:ASN:ND2	10:E:624:HOH:O	2.46	0.46
2:D:30:GLU:OE2	2:D:145:ASP:HB2	2.16	0.45
1:E:222:TRP:CE2	1:E:225:GLY:HA2	2.51	0.45
1:A:102:VAL:HG22	1:A:232:ILE:HB	1.98	0.45
1:C:182:VAL:HG21	1:C:213:ILE:HB	1.99	0.45
1:E:156:LYS:HE2	5:Q:1:NAG:H62	1.98	0.45
1:A:53:ASN:HD22	1:A:54:ASN:N	2.15	0.44
2:F:30:GLU:OE2	2:F:145:ASP:HB2	2.16	0.44
1:C:238:LYS:HD2	2:F:72:GLU:OE2	2.17	0.44
1:A:222:TRP:CZ2	1:A:225:GLY:HA2	2.53	0.44
1:E:18:HIS:HE1	10:E:689:HOH:O	2.01	0.44
1:E:185:PRO:HG2	1:E:191:GLN:OE1	2.18	0.44
1:E:207:ARG:NH1	1:E:240:GLY:O	2.51	0.43
1:E:84:TRP:CE2	1:E:116:GLY:HA2	2.53	0.43
1:A:307:LYS:HE3	2:B:60:ASN:OD1	2.18	0.43
1:E:102:VAL:HG22	1:E:232:ILE:HB	1.98	0.43
2:B:45:ILE:HD12	10:B:232:HOH:O	2.19	0.43
3:L:2:NAG:H4	3:L:3:BMA:H2	1.71	0.43
1:C:216:ASN:HB2	1:C:220:ARG:NH2	2.32	0.43
1:E:222:TRP:CZ2	1:E:225:GLY:HA2	2.54	0.43
1:C:207:ARG:NH2	10:C:610:HOH:O	2.43	0.42
2:D:3:PHE:CZ	2:F:2:LEU:HB3	2.55	0.42
1:A:195:TYR:O	1:A:197:GLN:N	2.51	0.42
1:A:216:ASN:OD1	1:E:212:THR:HG21	2.19	0.41
2:B:38:LEU:HD23	2:B:38:LEU:HA	1.93	0.41
1:E:212:THR:HG22	1:E:212:THR:O	2.20	0.41
1:A:29:ILE:HD11	2:B:102:LEU:HD23	2.02	0.41
2:D:3:PHE:HZ	2:F:2:LEU:HB3	1.86	0.41
1:A:50:LYS:HG2	1:A:273:PRO:HG2	2.03	0.40
1:E:305:CYS:O	2:F:59:THR:HB	2.21	0.40
1:C:195:TYR:O	1:C:197:GLN:N	2.54	0.40
1:E:123:GLU:OE1	1:E:168:MET:HG2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	319/321 (99%)	311 (98%)	7 (2%)	1 (0%)	41	44
1	C	315/321 (98%)	306 (97%)	8 (2%)	1 (0%)	41	44
1	E	315/321 (98%)	306 (97%)	8 (2%)	1 (0%)	41	44
2	B	170/176 (97%)	163 (96%)	7 (4%)	0	100	100
2	D	169/176 (96%)	160 (95%)	9 (5%)	0	100	100
2	F	169/176 (96%)	160 (95%)	9 (5%)	0	100	100
All	All	1457/1491 (98%)	1406 (96%)	48 (3%)	3 (0%)	47	53

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	62	ILE
1	E	62	ILE
1	A	62	ILE

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	282/282 (100%)	274 (97%)	8 (3%)	43	49
1	C	278/282 (99%)	273 (98%)	5 (2%)	59	66
1	E	278/282 (99%)	273 (98%)	5 (2%)	59	66
2	B	146/149 (98%)	144 (99%)	2 (1%)	67	74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	145/149 (97%)	145 (100%)	0	100	100
2	F	145/149 (97%)	144 (99%)	1 (1%)	84	88
All	All	1274/1293 (98%)	1253 (98%)	21 (2%)	62	70

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	48	THR
1	A	53	ASN
1	A	140	LYS
1	A	208	ARG
1	A	242	VAL
1	A	276	THR
1	A	326	LYS
2	B	40	SER
2	B	60	ASN
1	C	18	HIS
1	C	22	ASN
1	C	208	ARG
1	C	222	TRP
1	C	307	LYS
1	E	18	HIS
1	E	33	GLN
1	E	208	ARG
1	E	212	THR
1	E	264	LYS
2	F	38	LEU

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

Mol	Chain	Res	Type
1	A	53	ASN
2	B	60	ASN
2	F	60	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

36 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	G	1	1,3	14,14,15	0.28	0	17,19,21	0.37	0
3	NAG	G	2	3	14,14,15	0.28	0	17,19,21	0.57	0
3	BMA	G	3	3	11,11,12	0.82	0	15,15,17	0.78	0
3	NAG	H	1	1,3	14,14,15	0.47	0	17,19,21	0.64	0
3	NAG	H	2	3	14,14,15	0.28	0	17,19,21	0.50	0
3	BMA	H	3	3	11,11,12	1.04	1 (9%)	15,15,17	1.68	3 (20%)
4	NAG	I	1	1,4	14,14,15	0.34	0	17,19,21	0.61	0
4	NAG	I	2	4	14,14,15	0.49	0	17,19,21	0.37	0
5	NAG	J	1	5	15,15,15	0.13	0	21,21,21	0.33	0
5	GAL	J	2	5	11,11,12	1.11	2 (18%)	15,15,17	1.62	2 (13%)
5	NAG	J	3	5	14,14,15	0.55	0	17,19,21	0.83	1 (5%)
5	GAL	J	4	5	11,11,12	0.87	0	15,15,17	1.05	1 (6%)
5	SIA	J	5	5	17,20,21	1.01	1 (5%)	21,28,31	1.15	2 (9%)
6	NAG	K	1	1,6	14,14,15	0.48	0	17,19,21	0.58	0
6	NAG	K	2	6	14,14,15	0.28	0	17,19,21	0.60	0
6	BMA	K	3	6	11,11,12	1.41	2 (18%)	15,15,17	1.79	4 (26%)
6	MAN	K	4	6	11,11,12	1.41	3 (27%)	15,15,17	1.54	2 (13%)
3	NAG	L	1	1,3	14,14,15	0.75	1 (7%)	17,19,21	0.36	0
3	NAG	L	2	3	14,14,15	0.20	0	17,19,21	0.54	0
3	BMA	L	3	3	11,11,12	0.87	0	15,15,17	0.73	0
7	NAG	M	1	7	15,15,15	0.45	0	21,21,21	0.51	0
7	GAL	M	2	7	11,11,12	0.64	0	15,15,17	1.06	2 (13%)
7	SIA	M	3	7	17,20,21	1.05	0	21,28,31	1.12	2 (9%)
4	NAG	N	1	1,4	14,14,15	0.34	0	17,19,21	0.84	1 (5%)
4	NAG	N	2	4	14,14,15	0.33	0	17,19,21	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	O	1	1,8	14,14,15	0.31	0	17,19,21	0.73	1 (5%)
8	NAG	O	2	8	14,14,15	0.64	0	17,19,21	0.56	0
8	BMA	O	3	8	11,11,12	0.89	1 (9%)	15,15,17	1.57	3 (20%)
8	MAN	O	4	8	11,11,12	0.97	0	15,15,17	1.19	1 (6%)
4	NAG	P	1	1,4	14,14,15	0.62	1 (7%)	17,19,21	0.62	0
4	NAG	P	2	4	14,14,15	0.42	0	17,19,21	0.41	0
5	NAG	Q	1	5	15,15,15	0.19	0	21,21,21	0.18	0
5	GAL	Q	2	5	11,11,12	0.59	0	15,15,17	0.93	0
5	NAG	Q	3	5	14,14,15	0.74	1 (7%)	17,19,21	0.59	0
5	GAL	Q	4	5	11,11,12	0.71	0	15,15,17	1.52	3 (20%)
5	SIA	Q	5	5	17,20,21	0.94	1 (5%)	21,28,31	1.08	1 (4%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	G	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	BMA	G	3	3	-	1/2/19/22	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1
3	BMA	H	3	3	-	2/2/19/22	0/1/1/1
4	NAG	I	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	I	2	4	-	1/6/23/26	0/1/1/1
5	NAG	J	1	5	-	0/6/26/26	0/1/1/1
5	GAL	J	2	5	-	2/2/19/22	0/1/1/1
5	NAG	J	3	5	-	0/6/23/26	0/1/1/1
5	GAL	J	4	5	-	2/2/19/22	0/1/1/1
5	SIA	J	5	5	-	4/14/34/38	0/1/1/1
6	NAG	K	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	K	2	6	-	2/6/23/26	0/1/1/1
6	BMA	K	3	6	-	0/2/19/22	0/1/1/1
6	MAN	K	4	6	-	2/2/19/22	0/1/1/1
3	NAG	L	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	L	2	3	-	2/6/23/26	0/1/1/1
3	BMA	L	3	3	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	M	1	7	-	2/6/26/26	0/1/1/1
7	GAL	M	2	7	-	1/2/19/22	0/1/1/1
7	SIA	M	3	7	-	0/14/34/38	0/1/1/1
4	NAG	N	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	N	2	4	-	0/6/23/26	0/1/1/1
8	NAG	O	1	1,8	-	0/6/23/26	0/1/1/1
8	NAG	O	2	8	-	1/6/23/26	0/1/1/1
8	BMA	O	3	8	-	2/2/19/22	0/1/1/1
8	MAN	O	4	8	-	2/2/19/22	0/1/1/1
4	NAG	P	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	P	2	4	-	0/6/23/26	0/1/1/1
5	NAG	Q	1	5	-	0/6/26/26	0/1/1/1
5	GAL	Q	2	5	-	0/2/19/22	0/1/1/1
5	NAG	Q	3	5	-	0/6/23/26	0/1/1/1
5	GAL	Q	4	5	-	0/2/19/22	0/1/1/1
5	SIA	Q	5	5	-	0/14/34/38	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	K	3	BMA	C1-C2	3.84	1.61	1.52
6	K	4	MAN	C1-C2	2.93	1.58	1.52
3	H	3	BMA	C1-C2	2.80	1.58	1.52
5	Q	3	NAG	O5-C1	-2.53	1.39	1.43
6	K	4	MAN	O5-C5	2.35	1.48	1.43
5	Q	5	SIA	C7-C6	2.30	1.55	1.53
4	P	1	NAG	C1-C2	2.27	1.55	1.52
5	J	2	GAL	C1-C2	2.27	1.57	1.52
8	O	3	BMA	C2-C3	2.11	1.55	1.52
6	K	3	BMA	C2-C3	2.11	1.55	1.52
5	J	2	GAL	C2-C3	2.09	1.55	1.52
6	K	4	MAN	C2-C3	2.09	1.55	1.52
5	J	5	SIA	C6-C5	2.06	1.56	1.53
3	L	1	NAG	O5-C1	2.02	1.46	1.43

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	J	2	GAL	C1-C2-C3	4.95	115.75	109.67
6	K	4	MAN	C1-O5-C5	4.22	117.92	112.19
6	K	3	BMA	C1-C2-C3	4.06	114.65	109.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	Q	4	GAL	C1-O5-C5	3.71	117.22	112.19
8	O	4	MAN	C1-O5-C5	3.60	117.07	112.19
8	O	3	BMA	C1-C2-C3	3.59	114.08	109.67
3	H	3	BMA	C1-C2-C3	3.52	113.99	109.67
3	H	3	BMA	C1-O5-C5	3.50	116.94	112.19
6	K	3	BMA	O5-C1-C2	3.41	116.04	110.77
3	H	3	BMA	O5-C1-C2	2.93	115.30	110.77
4	N	1	NAG	C1-O5-C5	2.85	116.06	112.19
7	M	3	SIA	C4-C3-C2	2.64	114.54	109.81
8	O	3	BMA	C2-C3-C4	2.63	115.45	110.89
7	M	2	GAL	C1-O5-C5	2.63	115.75	112.19
7	M	3	SIA	C6-O6-C2	2.61	116.92	111.34
5	Q	4	GAL	C1-C2-C3	2.58	112.84	109.67
5	Q	4	GAL	O5-C5-C6	-2.55	103.20	107.20
6	K	3	BMA	C1-O5-C5	2.50	115.58	112.19
5	J	4	GAL	C1-C2-C3	2.40	112.61	109.67
8	O	3	BMA	C1-O5-C5	2.30	115.31	112.19
6	K	3	BMA	O5-C5-C4	-2.27	105.29	110.83
5	J	3	NAG	C3-C4-C5	2.18	114.12	110.24
5	Q	5	SIA	C3-C2-C1	2.17	116.67	111.93
5	J	5	SIA	O4-C4-C5	-2.16	104.80	109.77
5	J	2	GAL	C1-O5-C5	2.15	115.11	112.19
8	O	1	NAG	C1-O5-C5	2.11	115.05	112.19
6	K	4	MAN	C1-C2-C3	2.09	112.23	109.67
7	M	2	GAL	O2-C2-C1	2.02	113.28	109.15
5	J	5	SIA	C3-C4-C5	-2.02	109.02	111.46

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	K	2	NAG	O5-C5-C6-O6
3	L	2	NAG	O5-C5-C6-O6
5	J	4	GAL	O5-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
6	K	2	NAG	C4-C5-C6-O6
3	L	2	NAG	C4-C5-C6-O6
6	K	4	MAN	C4-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
5	J	4	GAL	C4-C5-C6-O6
3	L	1	NAG	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6

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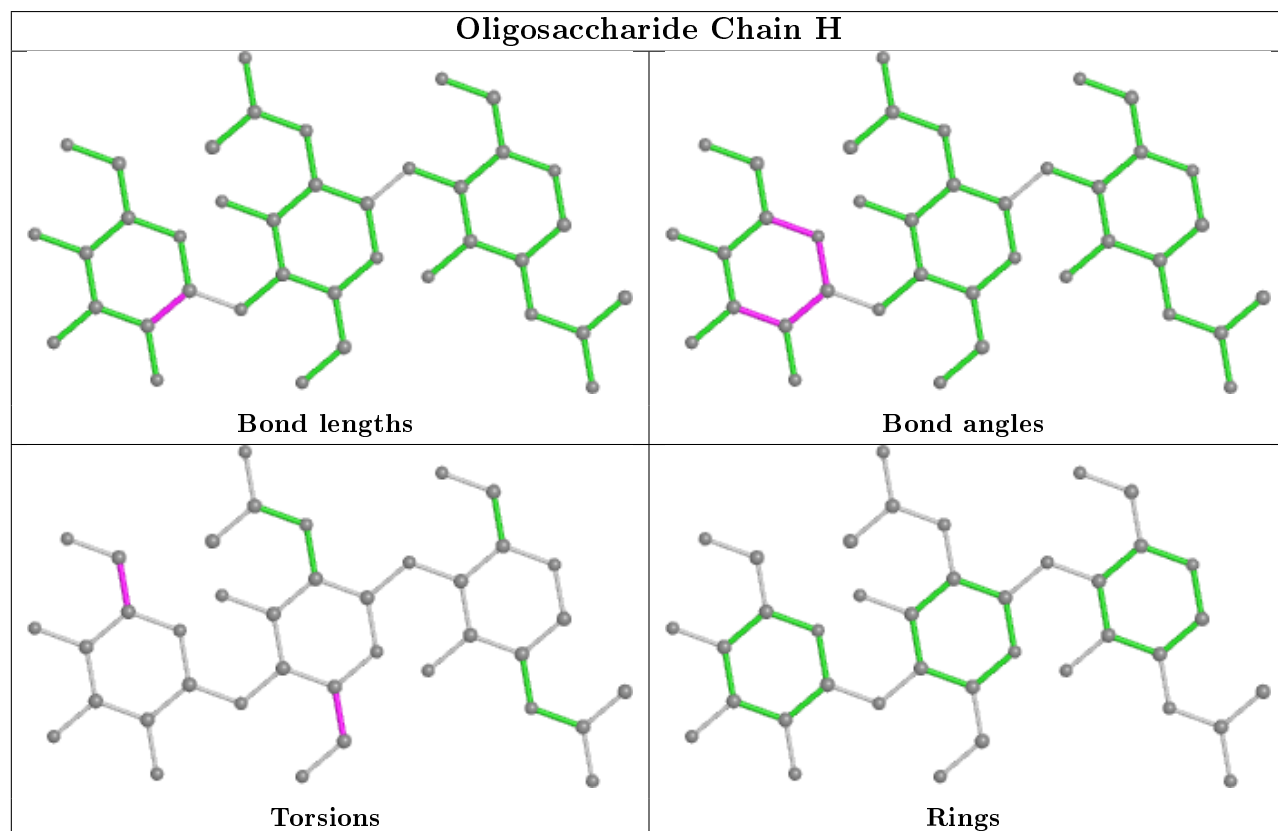
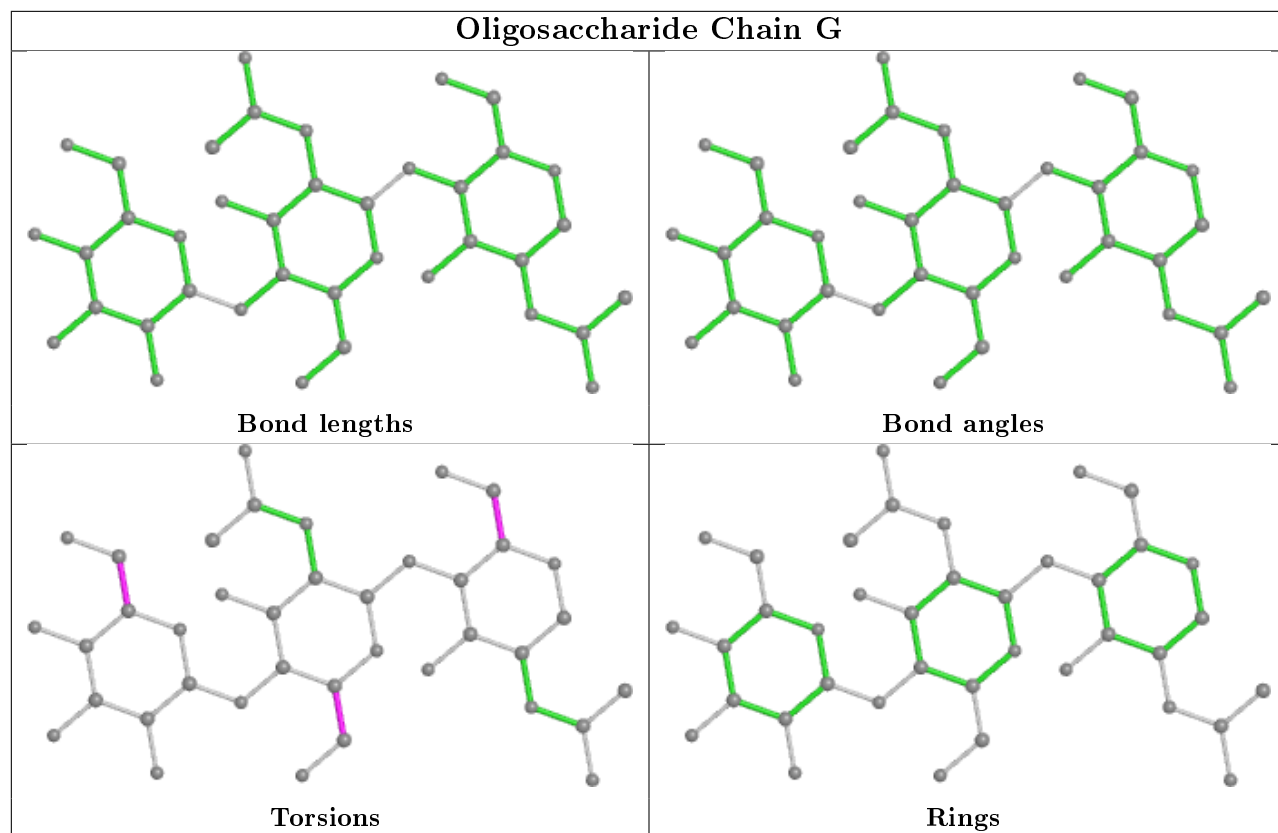
Mol	Chain	Res	Type	Atoms
6	K	4	MAN	O5-C5-C6-O6
3	L	1	NAG	C4-C5-C6-O6
8	O	4	MAN	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
5	J	5	SIA	C6-C7-C8-O8
5	J	2	GAL	O5-C5-C6-O6
5	J	5	SIA	O7-C7-C8-O8
8	O	3	BMA	C4-C5-C6-O6
8	O	3	BMA	O5-C5-C6-O6
3	H	3	BMA	O5-C5-C6-O6
4	I	2	NAG	O5-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
3	G	3	BMA	O5-C5-C6-O6
7	M	1	NAG	C4-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6
8	O	4	MAN	C4-C5-C6-O6
8	O	2	NAG	C4-C5-C6-O6
3	H	3	BMA	C4-C5-C6-O6
5	J	5	SIA	O7-C7-C8-C9
5	J	5	SIA	C6-C7-C8-C9
7	M	1	NAG	O5-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
5	J	2	GAL	C4-C5-C6-O6
7	M	2	GAL	O5-C5-C6-O6

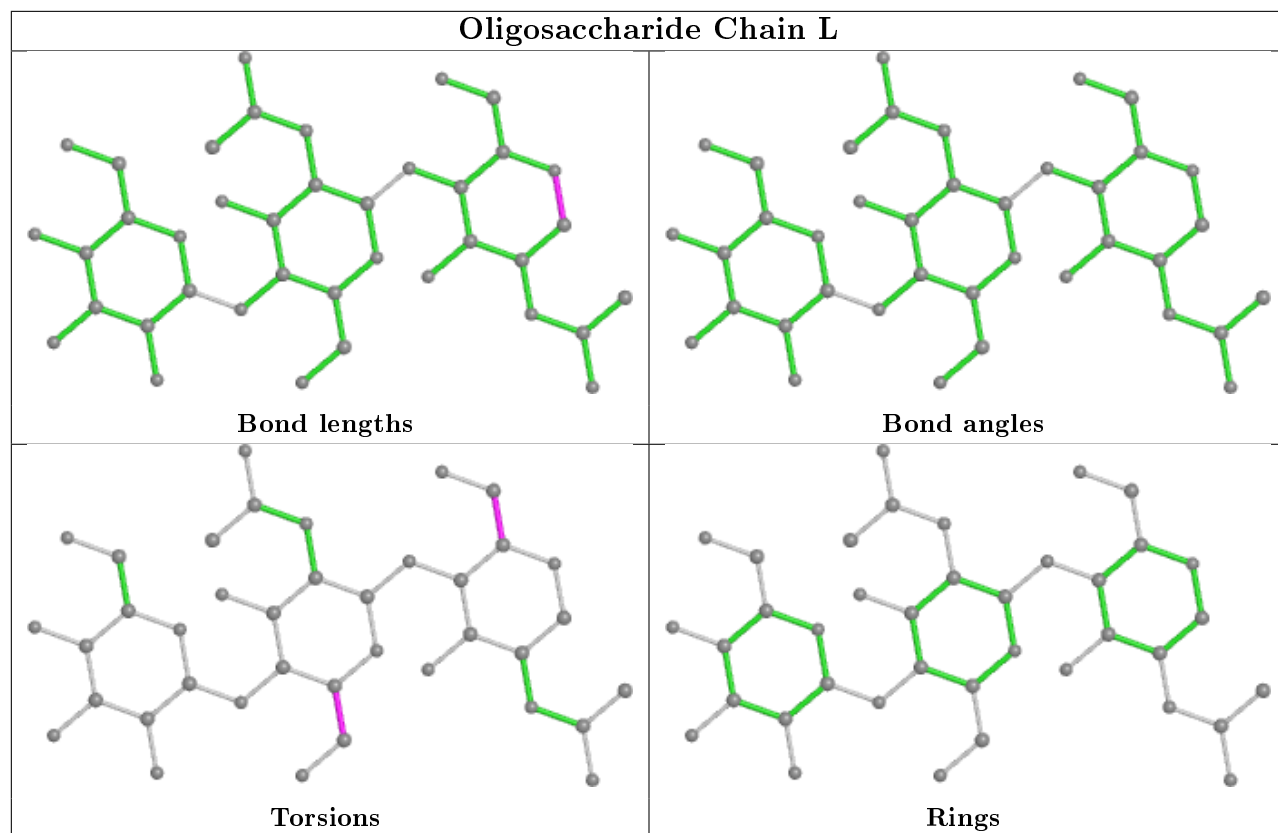
There are no ring outliers.

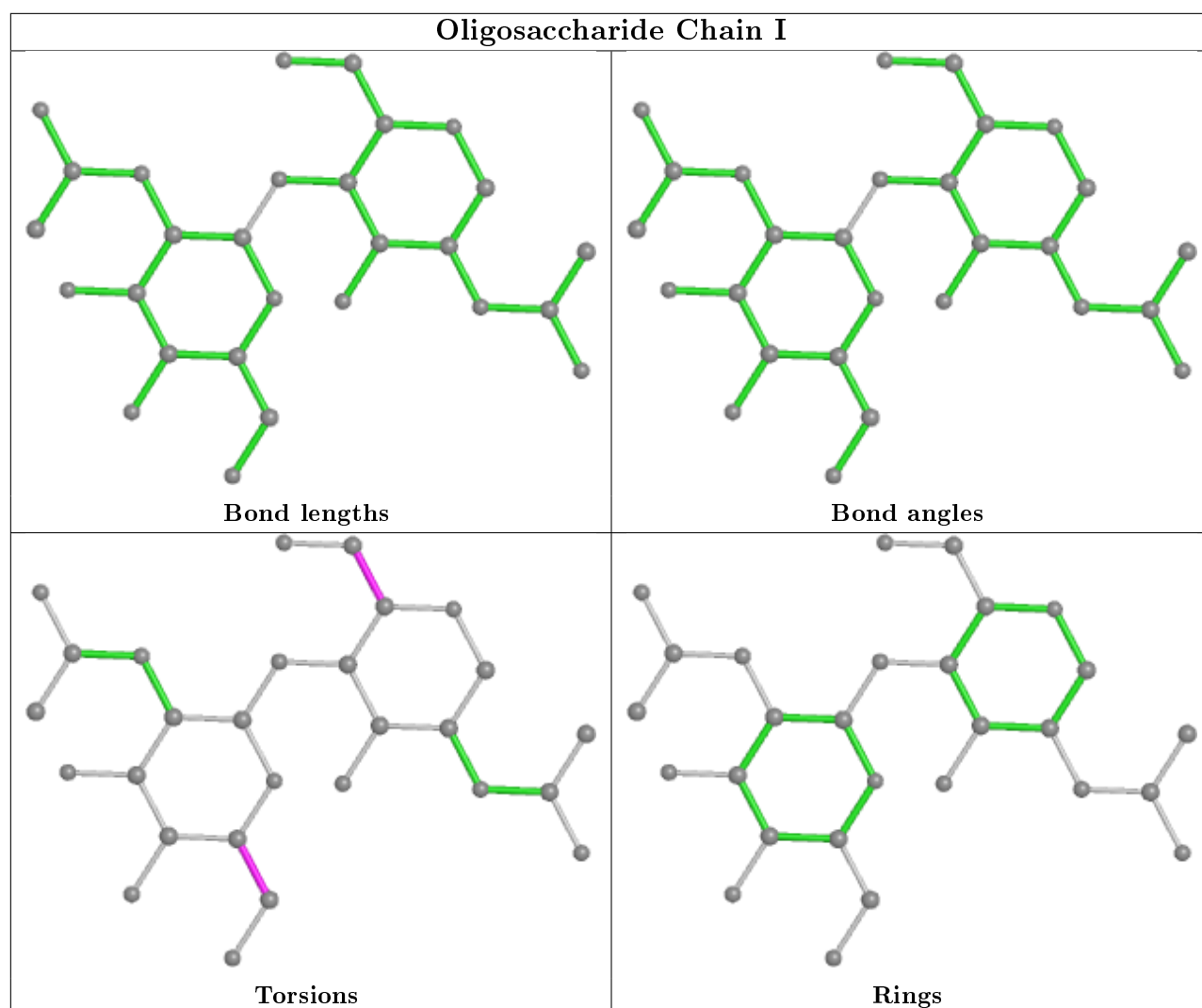
5 monomers are involved in 4 short contacts:

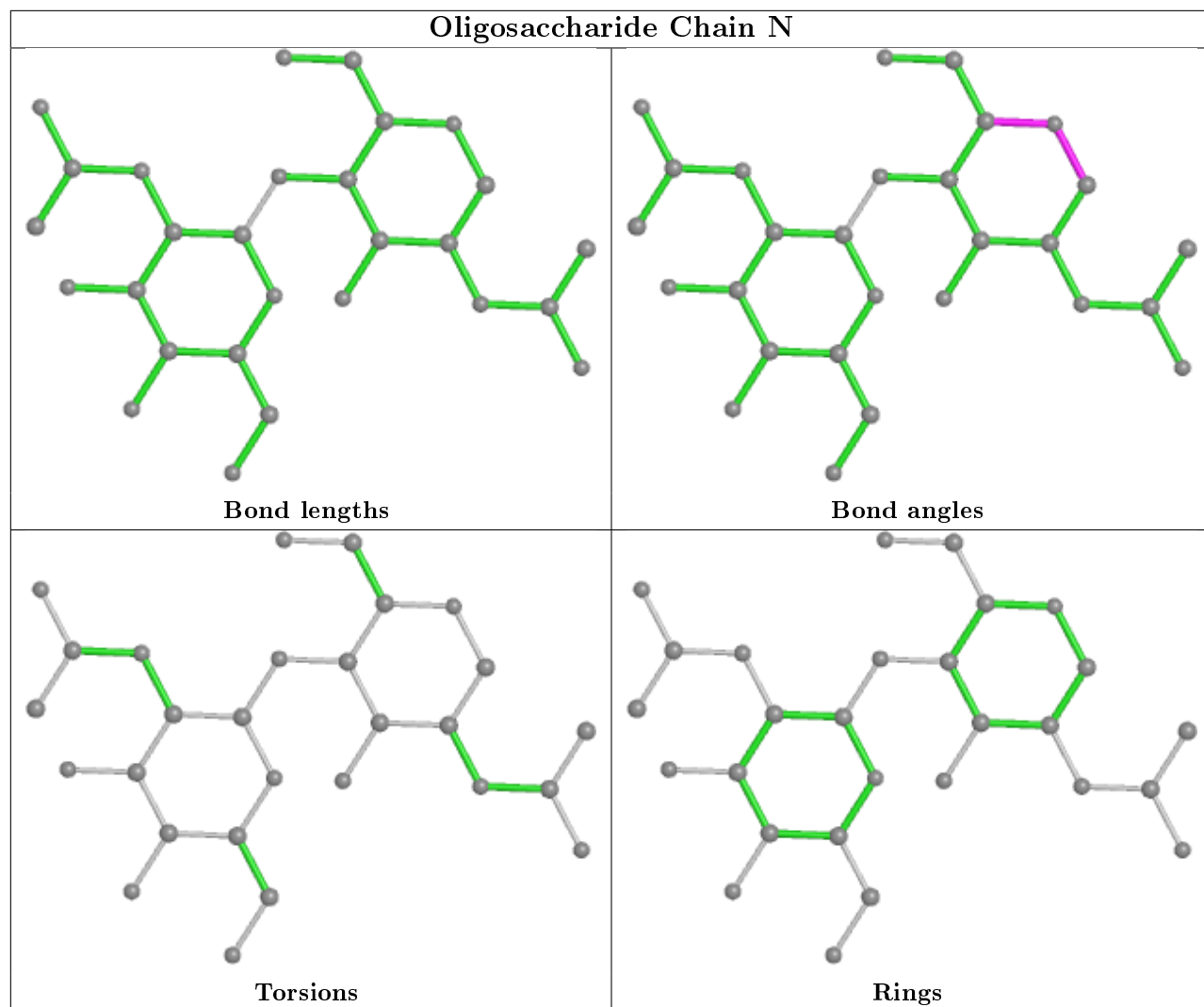
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	2	NAG	1	0
6	K	1	NAG	1	0
3	L	3	BMA	1	0
5	Q	1	NAG	1	0
3	L	2	NAG	1	0

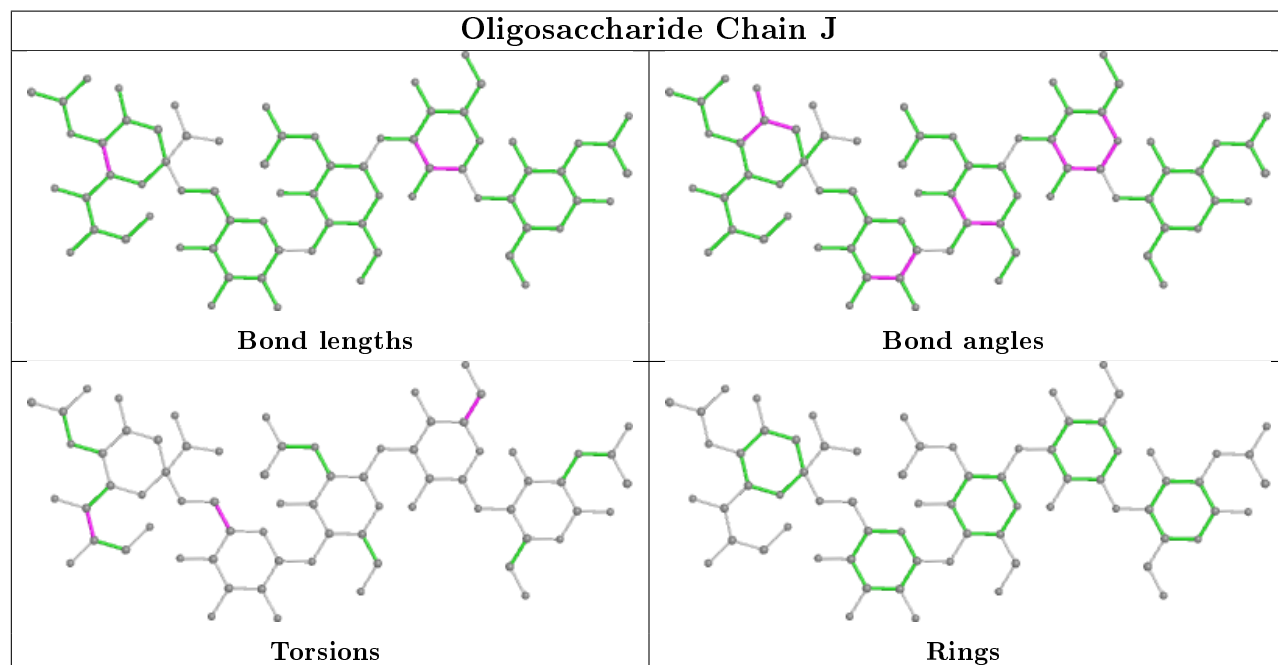
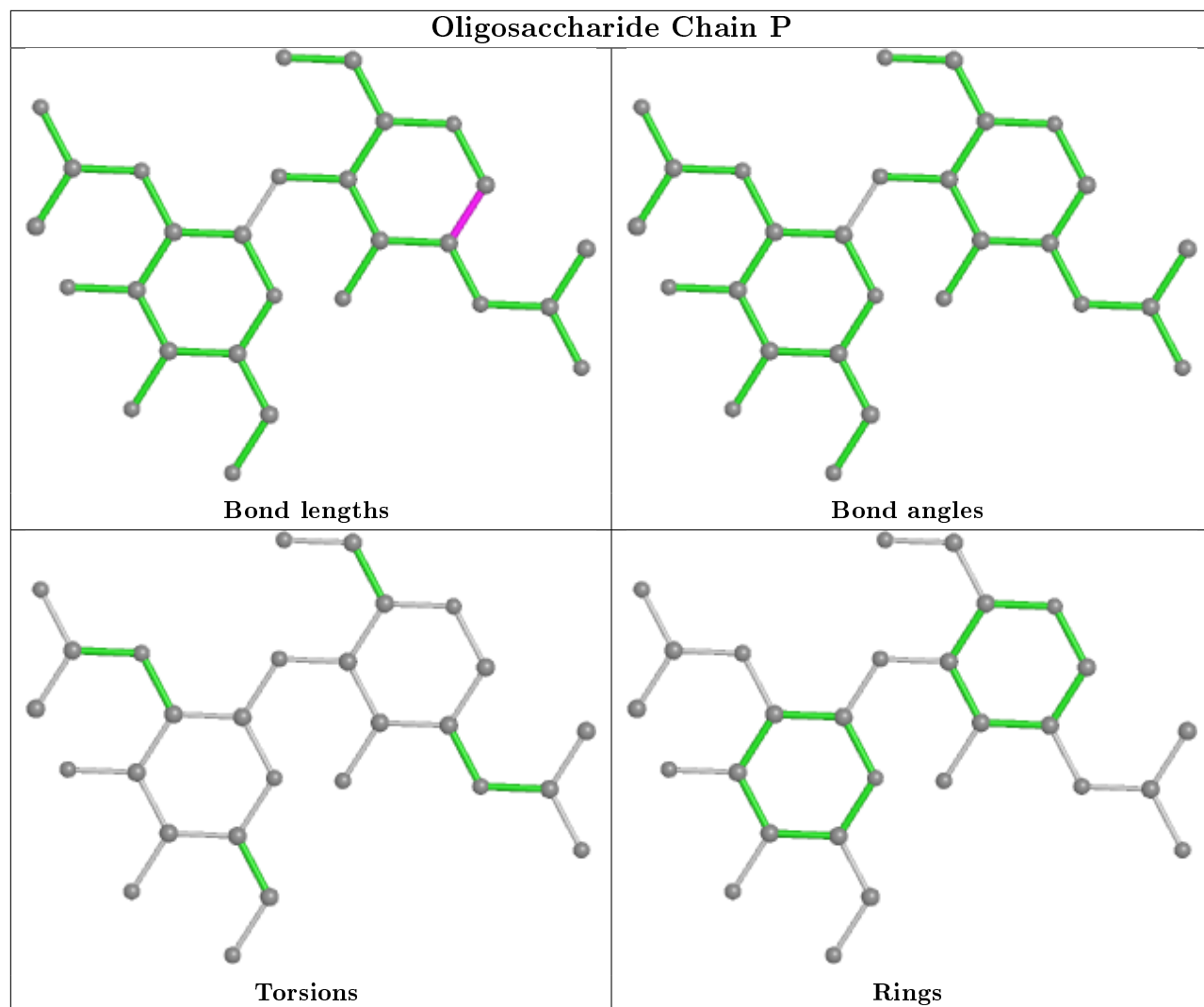
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



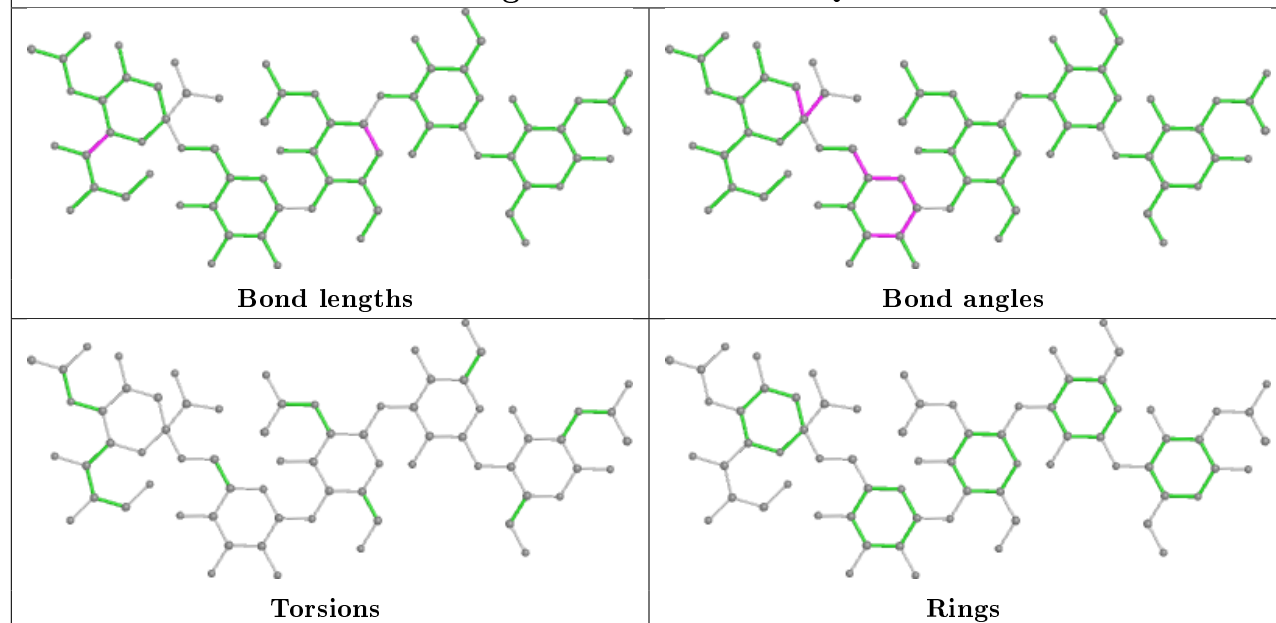




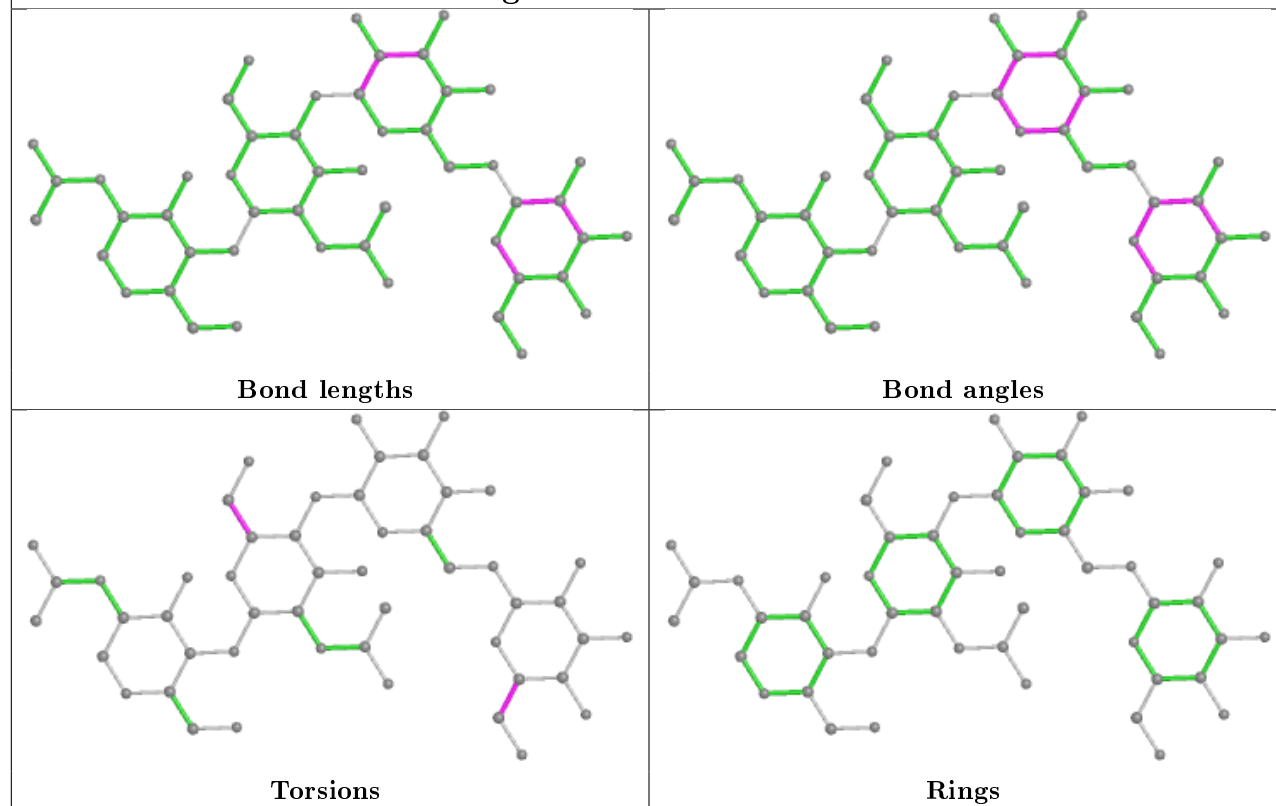


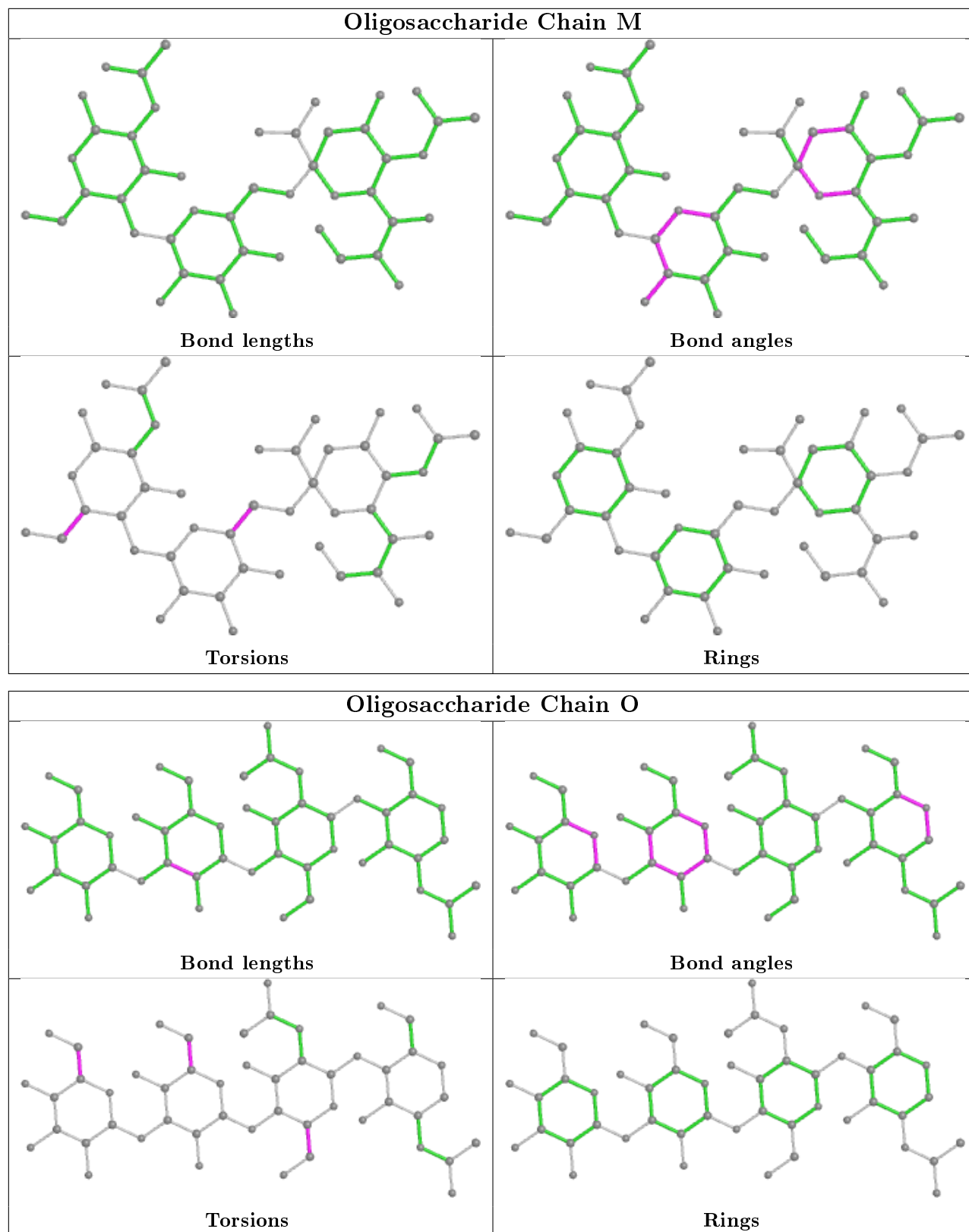


Oligosaccharide Chain Q



Oligosaccharide Chain K





5.6 Ligand geometry [i](#)

5 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	D	201	2	14,14,15	0.94	2 (14%)	17,19,21	0.61	0
9	NAG	C	501	1	14,14,15	0.86	1 (7%)	17,19,21	1.34	1 (5%)
9	NAG	C	502	1	14,14,15	1.07	1 (7%)	17,19,21	1.66	1 (5%)
9	NAG	A	404	1	14,14,15	0.74	1 (7%)	17,19,21	0.74	1 (5%)
9	NAG	E	509	1	14,14,15	0.96	2 (14%)	17,19,21	0.72	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	D	201	2	-	3/6/23/26	0/1/1/1
9	NAG	C	501	1	-	2/6/23/26	0/1/1/1
9	NAG	C	502	1	-	0/6/23/26	0/1/1/1
9	NAG	A	404	1	-	2/6/23/26	0/1/1/1
9	NAG	E	509	1	-	2/6/23/26	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	C	502	NAG	O5-C1	3.86	1.49	1.43
9	C	501	NAG	O5-C1	3.03	1.48	1.43
9	E	509	NAG	O5-C1	2.87	1.48	1.43
9	D	201	NAG	C1-C2	2.63	1.56	1.52
9	A	404	NAG	O5-C1	2.57	1.47	1.43
9	D	201	NAG	O5-C1	2.16	1.47	1.43
9	E	509	NAG	C1-C2	2.03	1.55	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	C	502	NAG	C1-O5-C5	6.33	120.77	112.19
9	C	501	NAG	C1-O5-C5	5.16	119.19	112.19
9	A	404	NAG	C1-O5-C5	2.40	115.44	112.19

There are no chirality outliers.

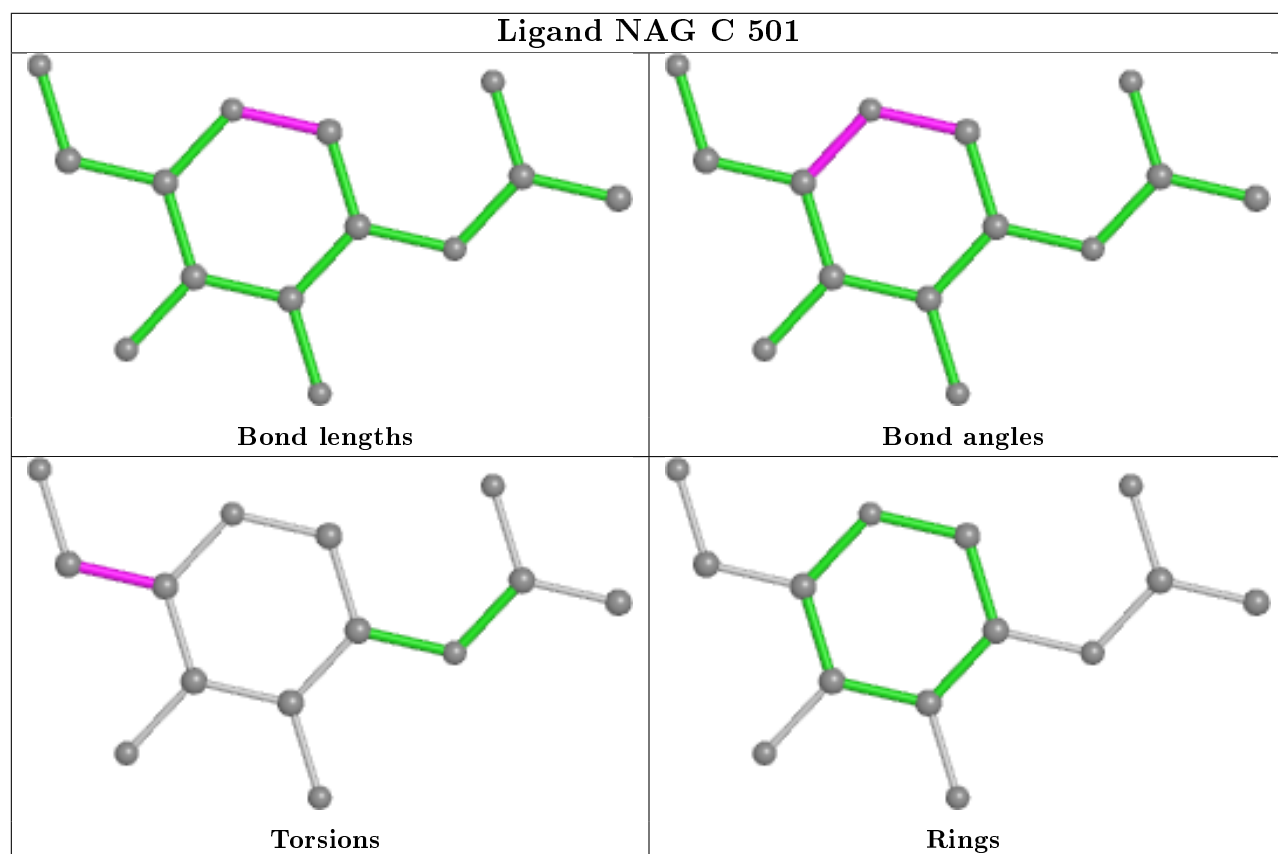
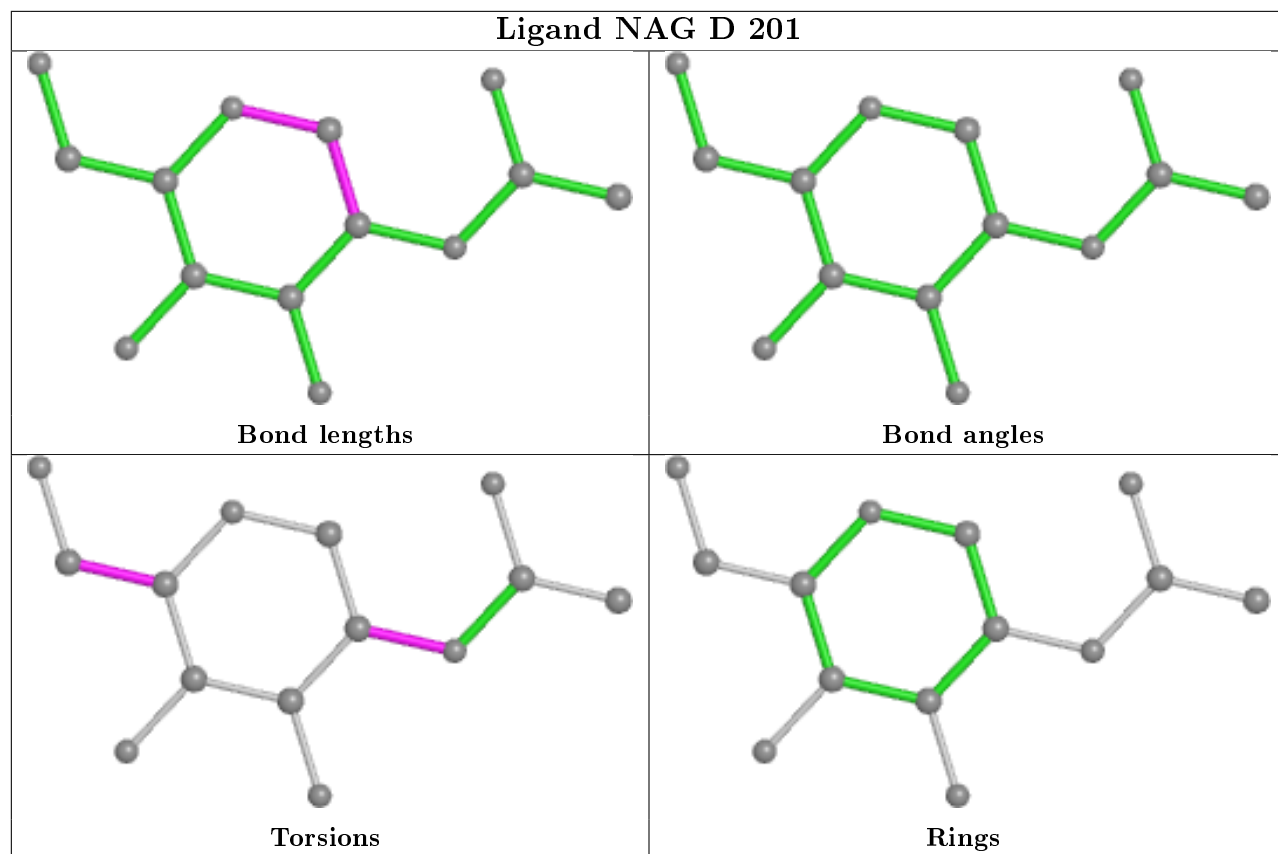
All (9) torsion outliers are listed below:

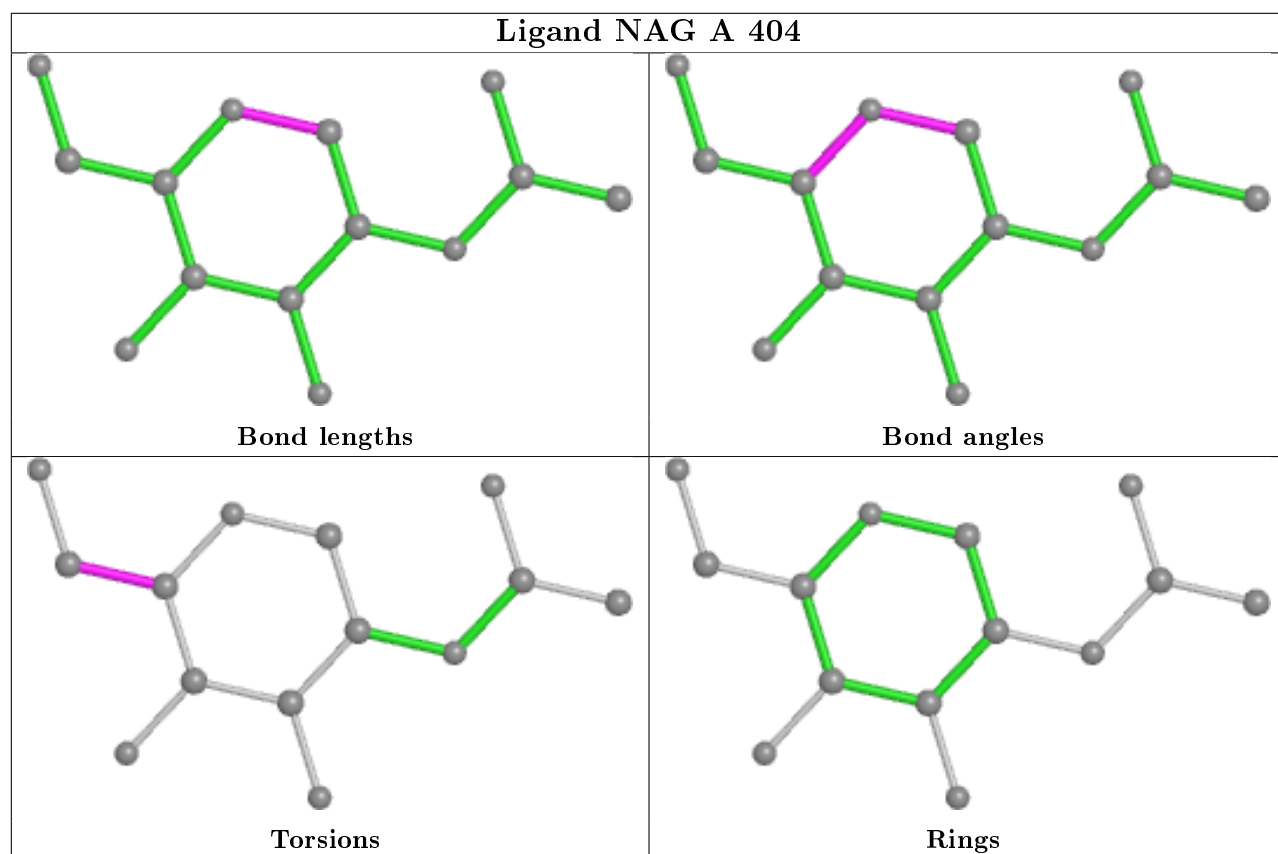
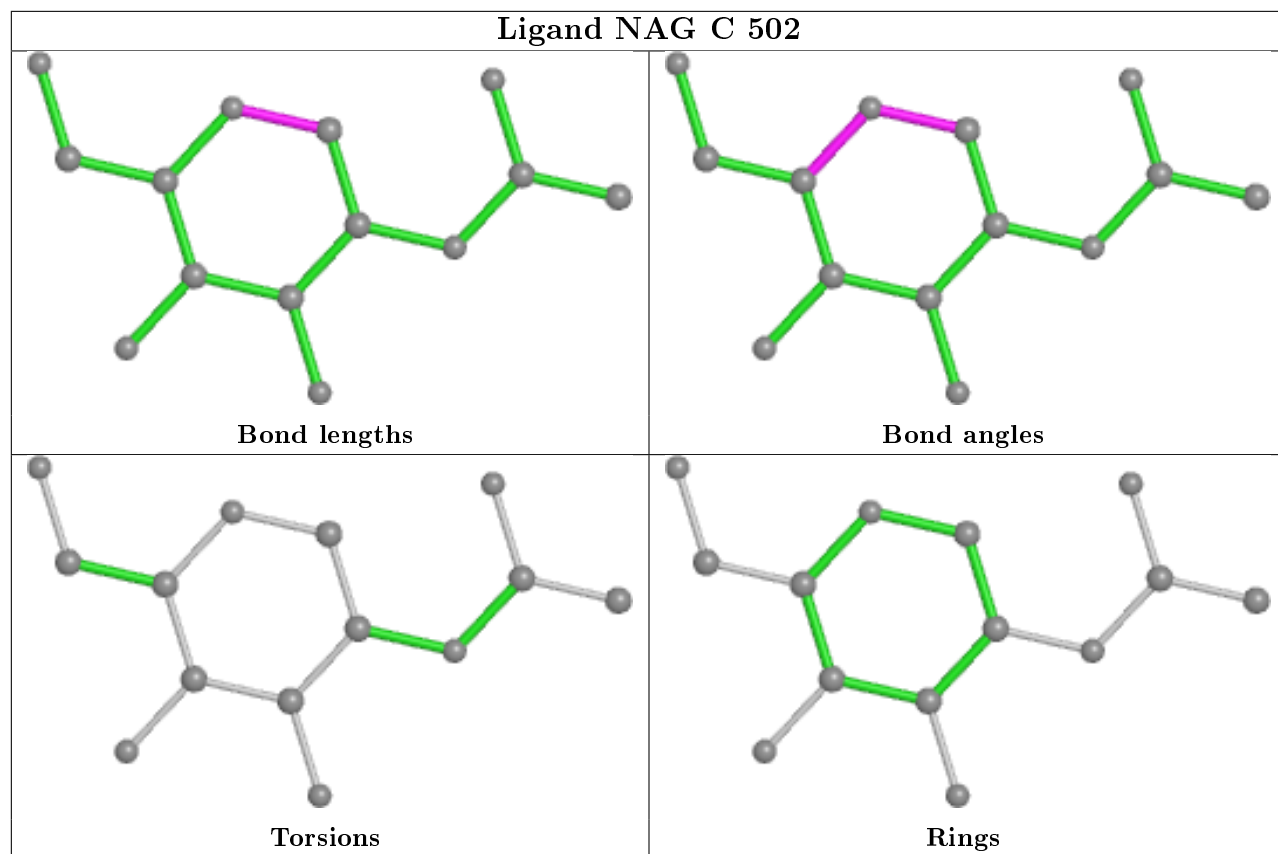
Mol	Chain	Res	Type	Atoms
9	E	509	NAG	O5-C5-C6-O6
9	D	201	NAG	O5-C5-C6-O6
9	D	201	NAG	C4-C5-C6-O6
9	E	509	NAG	C4-C5-C6-O6
9	A	404	NAG	O5-C5-C6-O6
9	A	404	NAG	C4-C5-C6-O6
9	C	501	NAG	C4-C5-C6-O6
9	D	201	NAG	C3-C2-N2-C7
9	C	501	NAG	O5-C5-C6-O6

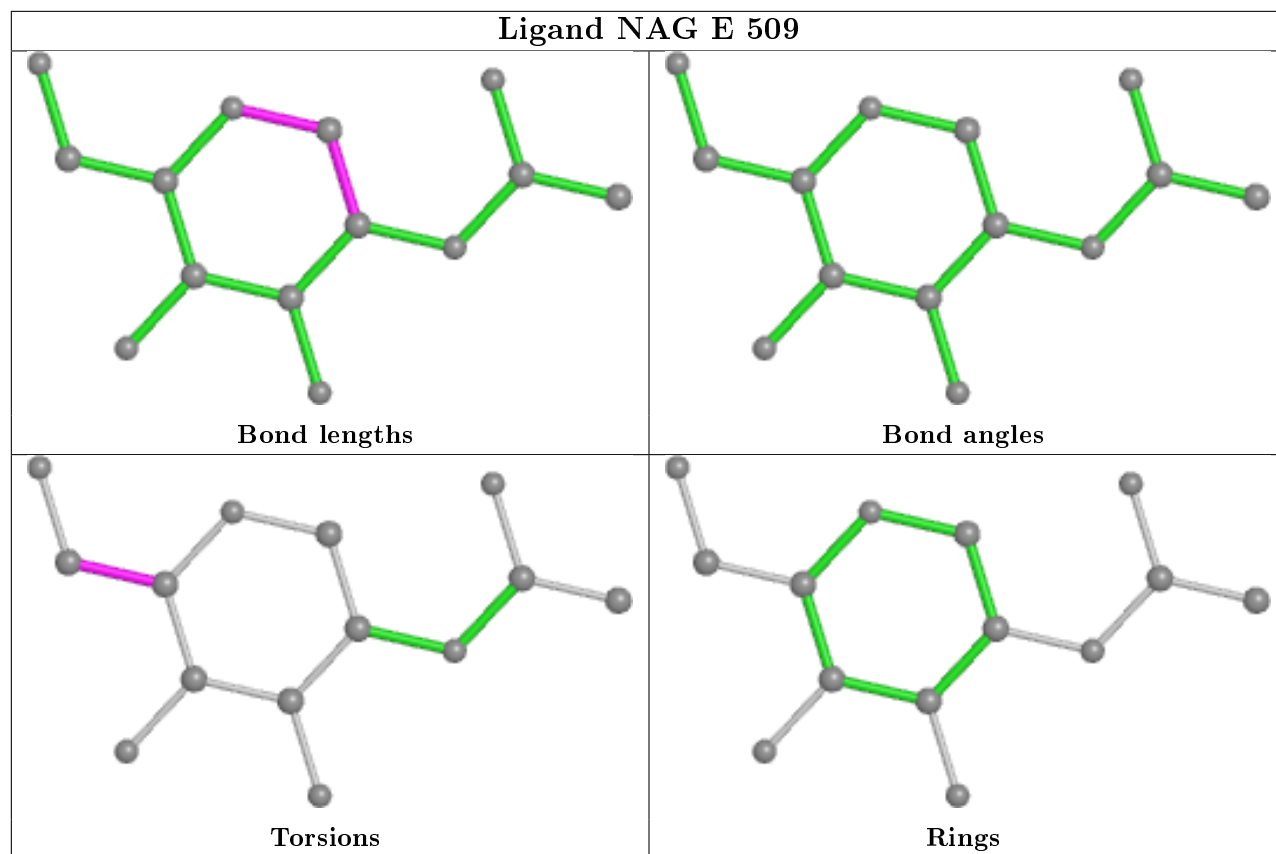
There are no ring outliers.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	321/321 (100%)	-0.11	6 (1%) 66 68	23, 46, 82, 115	0
1	C	317/321 (98%)	-0.20	6 (1%) 66 68	26, 41, 62, 105	0
1	E	317/321 (98%)	-0.22	2 (0%) 89 89	22, 34, 57, 107	0
2	B	172/176 (97%)	0.18	9 (5%) 27 26	22, 55, 88, 126	0
2	D	171/176 (97%)	0.40	12 (7%) 16 15	24, 51, 84, 100	0
2	F	171/176 (97%)	0.38	16 (9%) 8 8	22, 53, 92, 103	0
All	All	1469/1491 (98%)	-0.00	51 (3%) 44 43	22, 42, 82, 126	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	147	ALA	5.2
1	C	9	PRO	4.7
1	C	22	ASN	4.1
2	F	16	GLY	4.0
1	A	159	SER	3.8
2	F	18	ILE	3.6
2	F	171	PHE	3.5
2	B	18	ILE	3.5
2	D	58	LYS	3.4
1	A	328	THR	3.4
2	B	29	SER	3.3
2	F	147	ALA	3.1
2	B	31	GLY	3.1
2	B	172	GLN	3.1
2	F	2	LEU	3.1
1	C	276	THR	3.0
2	D	157	TYR	2.9
2	F	31	GLY	2.9
2	D	156	THR	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	222	TRP	2.9
1	A	9	PRO	2.8
1	C	277	CYS	2.7
2	F	168	ASN	2.7
2	B	147	ALA	2.6
2	F	59	THR	2.6
2	B	143	LYS	2.6
2	D	143	LYS	2.6
2	F	164	ASP	2.4
2	F	30	GLU	2.4
2	D	99	LEU	2.3
2	F	17	MET	2.3
1	A	126	THR	2.3
1	E	222	TRP	2.3
2	B	30	GLU	2.3
2	D	33	GLY	2.2
2	D	162	TYR	2.2
2	F	58	LYS	2.2
1	E	189	GLN	2.2
1	A	128	THR	2.2
2	F	143	LYS	2.2
2	F	98	LEU	2.2
2	D	57	GLU	2.2
2	B	146	ASN	2.1
1	C	10	GLY	2.1
2	D	161	VAL	2.1
2	B	167	LEU	2.1
1	A	129	GLY	2.1
2	F	29	SER	2.1
2	F	100	VAL	2.1
2	D	150	GLU	2.0
2	D	106	HIS	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,

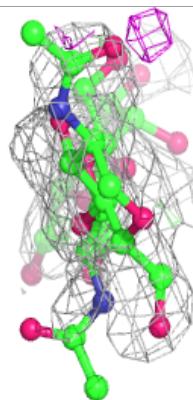
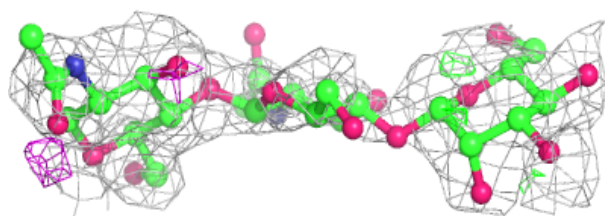
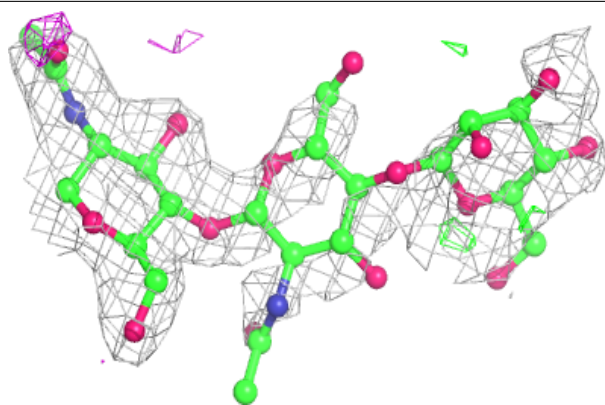
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	MAN	K	4	11/12	0.43	0.61	130,140,145,148	0
3	BMA	H	3	11/12	0.62	0.49	119,125,132,132	0
3	NAG	H	2	14/15	0.64	0.46	118,126,130,133	0
4	NAG	N	2	14/15	0.67	0.51	114,129,134,136	0
3	BMA	G	3	11/12	0.67	0.29	117,124,130,132	0
5	NAG	J	1	15/15	0.69	0.44	115,132,136,137	0
3	BMA	L	3	11/12	0.70	0.50	120,130,134,136	0
5	NAG	Q	1	15/15	0.70	0.58	135,140,142,143	0
6	BMA	K	3	11/12	0.74	0.21	103,110,117,124	0
5	GAL	J	2	11/12	0.75	0.42	122,131,136,138	0
7	NAG	M	1	15/15	0.75	0.30	68,87,97,99	0
5	GAL	Q	2	11/12	0.76	0.36	105,120,125,128	0
4	NAG	I	2	14/15	0.76	0.36	71,100,116,122	0
8	MAN	O	4	11/12	0.76	0.52	127,130,133,133	0
8	NAG	O	2	14/15	0.76	0.21	65,84,95,98	0
3	NAG	G	2	14/15	0.79	0.55	118,124,131,133	0
3	NAG	L	2	14/15	0.79	0.30	87,99,112,124	0
5	NAG	J	3	14/15	0.81	0.27	83,95,110,111	0
4	NAG	N	1	14/15	0.82	0.34	66,79,99,110	0
4	NAG	P	2	14/15	0.82	0.34	82,98,109,114	0
8	BMA	O	3	11/12	0.84	0.20	91,96,108,120	0
3	NAG	H	1	14/15	0.86	0.26	70,91,100,112	0
5	GAL	J	4	11/12	0.88	0.22	56,73,89,89	0
8	NAG	O	1	14/15	0.89	0.22	52,69,78,80	0
5	NAG	Q	3	14/15	0.89	0.18	67,79,89,92	0
3	NAG	G	1	14/15	0.90	0.29	60,74,91,103	0
4	NAG	P	1	14/15	0.92	0.20	40,49,63,73	0
3	NAG	L	1	14/15	0.93	0.18	36,56,69,83	0
4	NAG	I	1	14/15	0.93	0.22	39,50,69,76	0
7	GAL	M	2	11/12	0.94	0.18	46,53,66,68	0
5	SIA	J	5	20/21	0.94	0.15	47,53,59,63	0
6	NAG	K	2	14/15	0.95	0.19	39,62,88,91	0
6	NAG	K	1	14/15	0.95	0.15	41,56,67,74	0
7	SIA	M	3	20/21	0.96	0.11	30,39,45,53	0
5	GAL	Q	4	11/12	0.96	0.18	37,57,65,73	0
5	SIA	Q	5	20/21	0.97	0.10	32,40,48,49	0

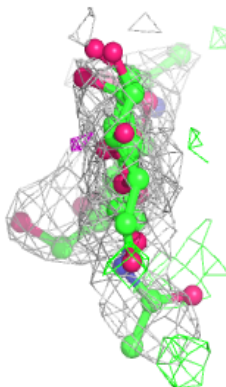
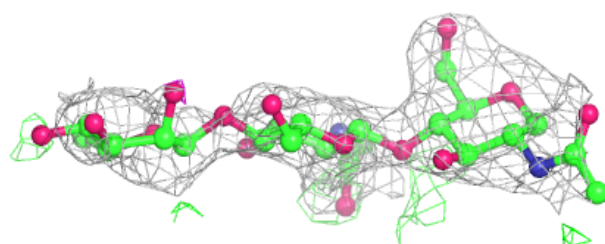
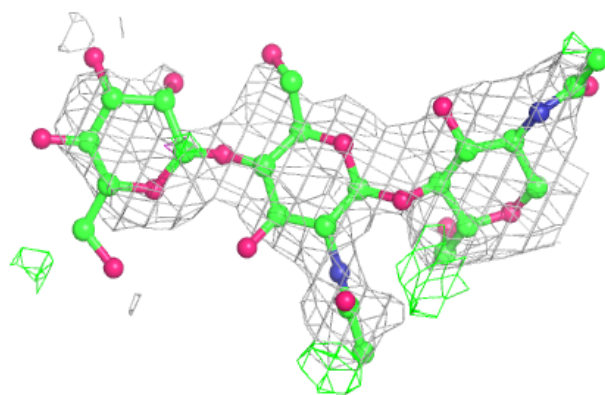
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain G:

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

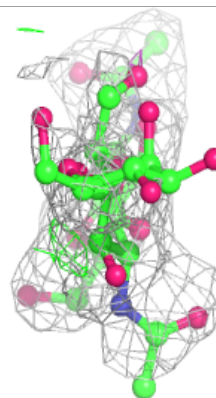
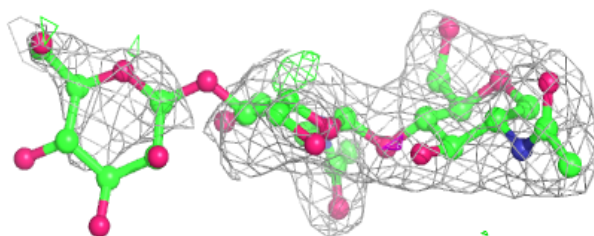
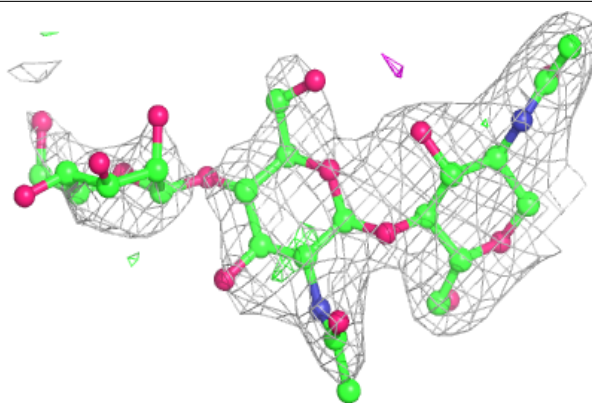
**Electron density around Chain H:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

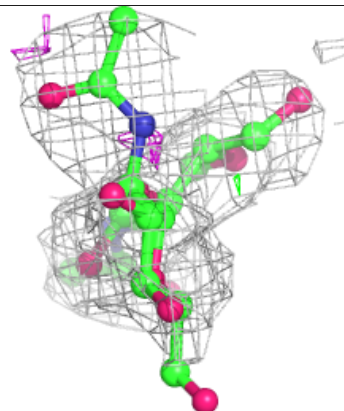
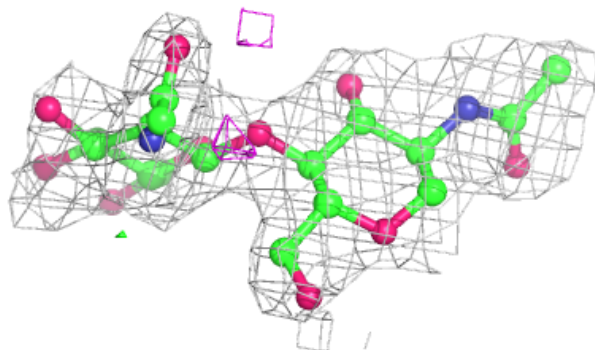
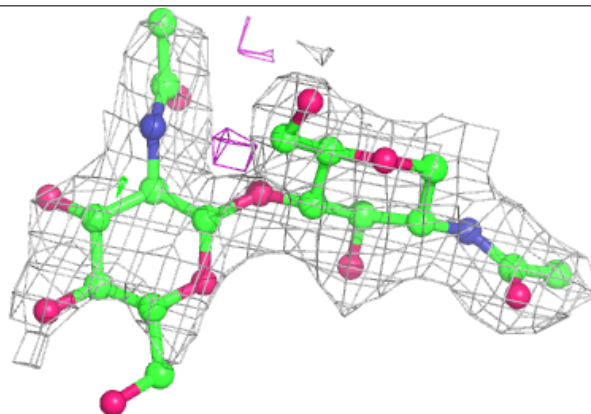


Electron density around Chain L:

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

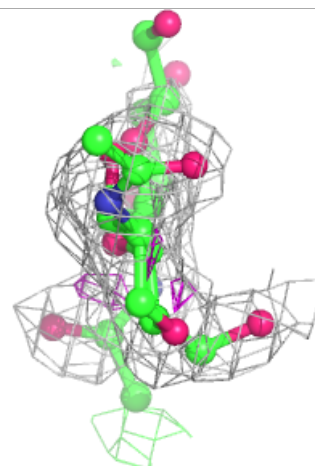
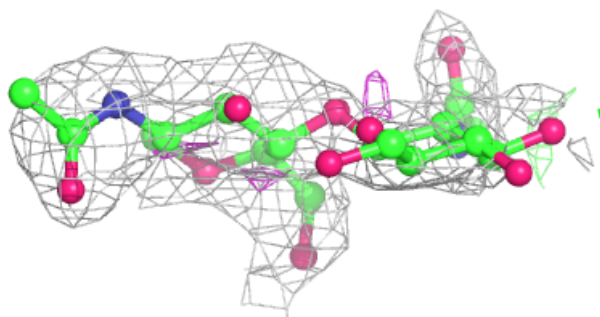
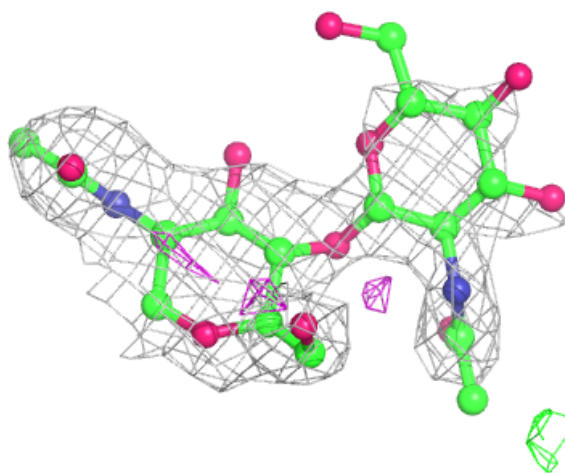
**Electron density around Chain I:**

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



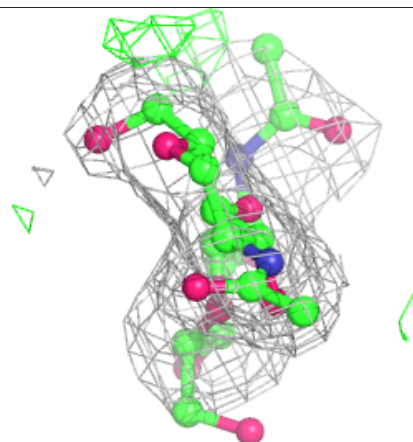
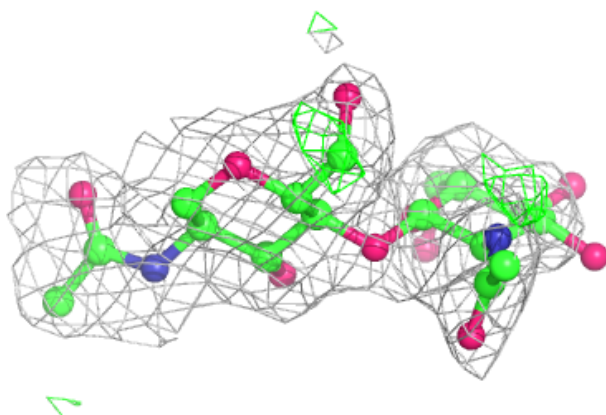
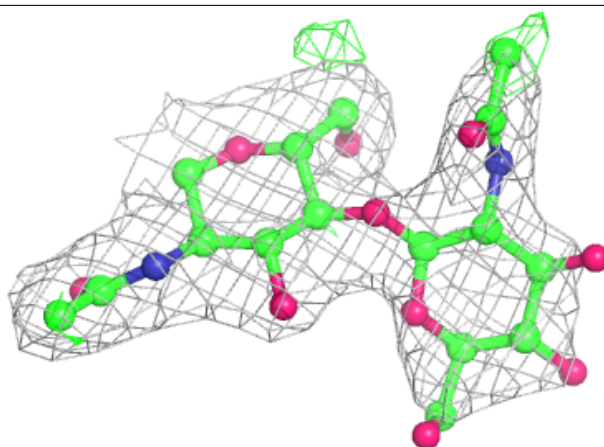
Electron density around Chain N:

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and green (positive)

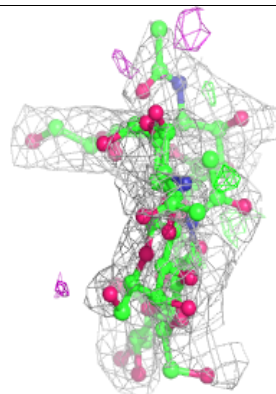
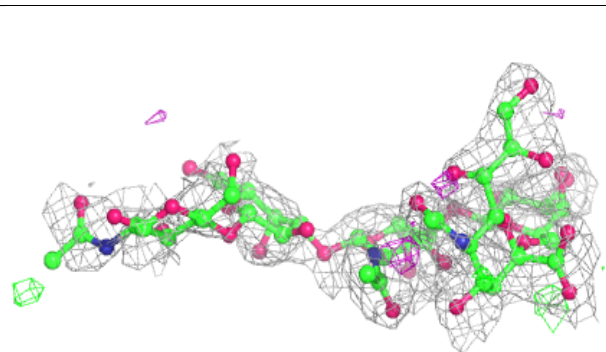
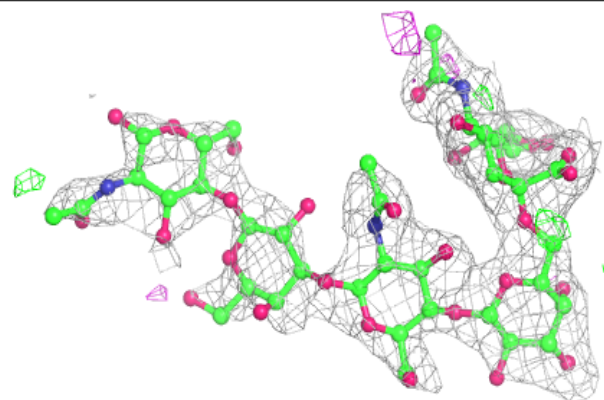


Electron density around Chain P:

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

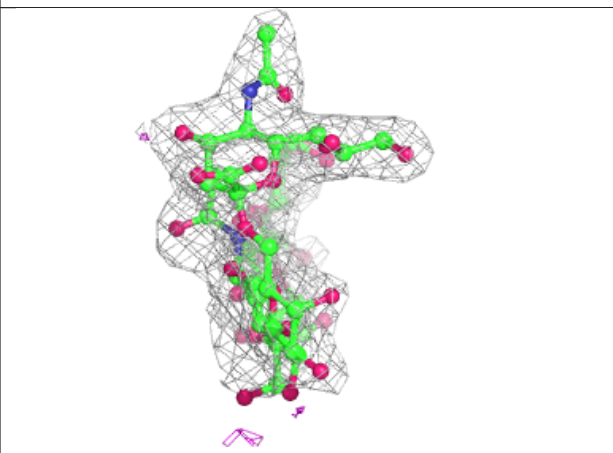
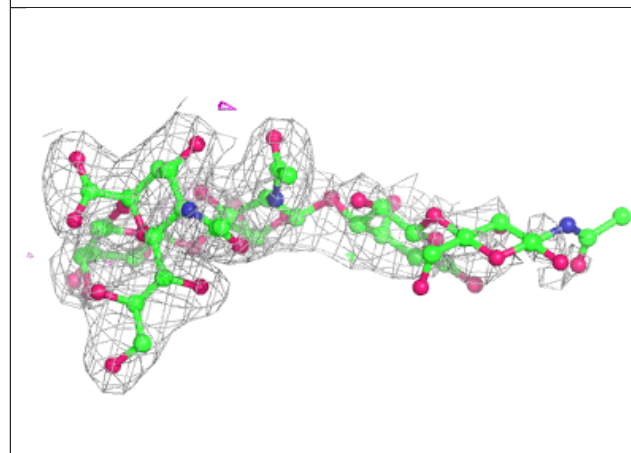
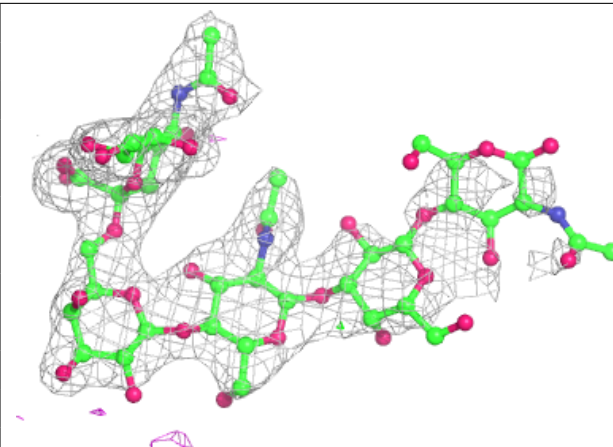
**Electron density around Chain J:**

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

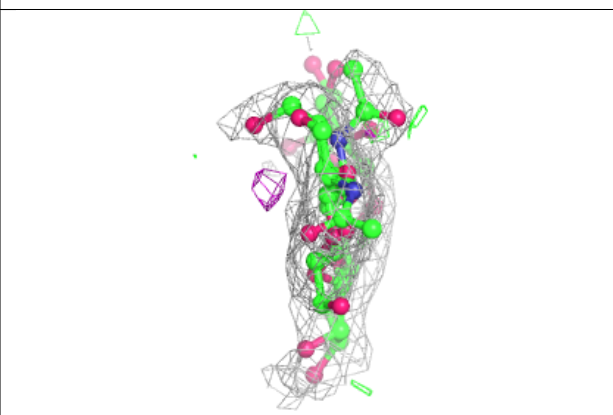
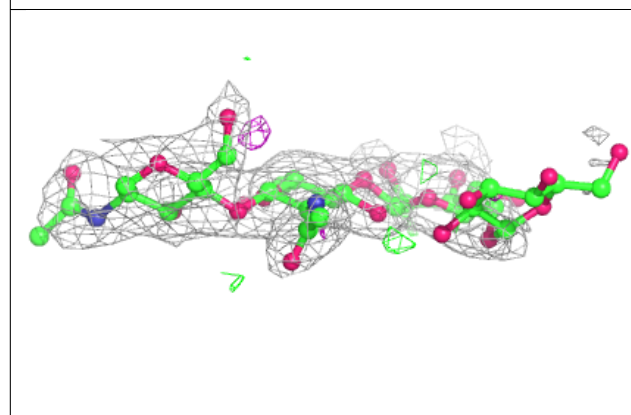
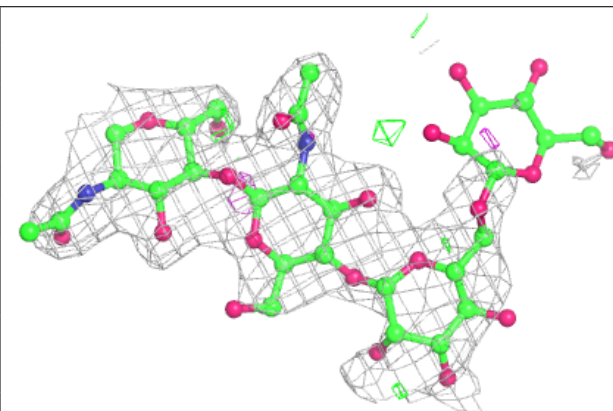


Electron density around Chain Q:

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

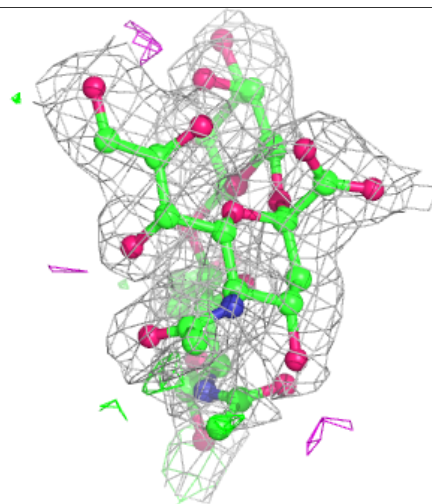
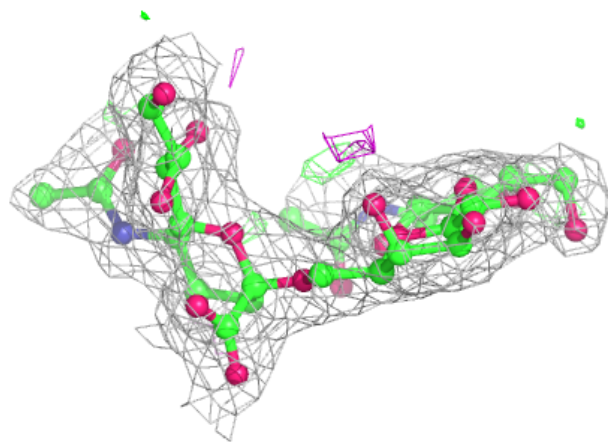
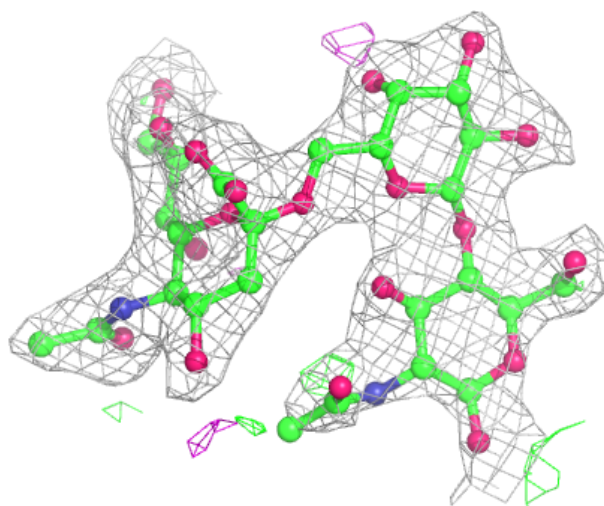
**Electron density around Chain K:**

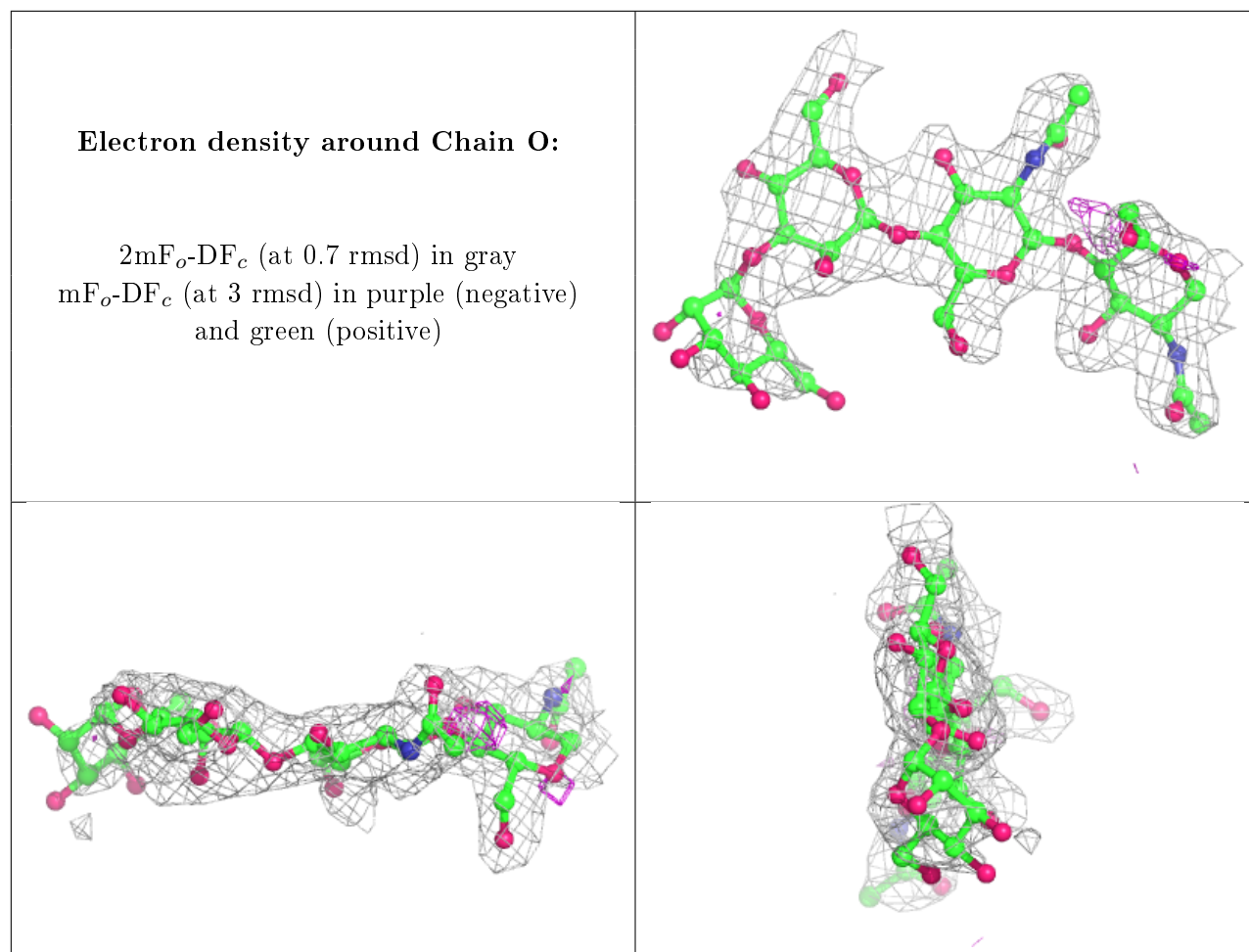
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain M:

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





6.4 Ligands ⓘ

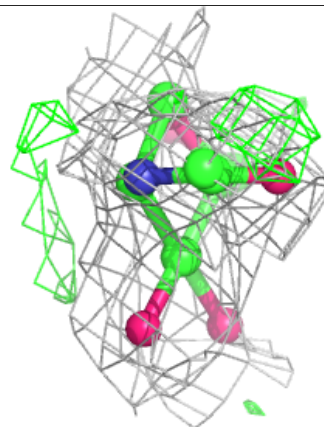
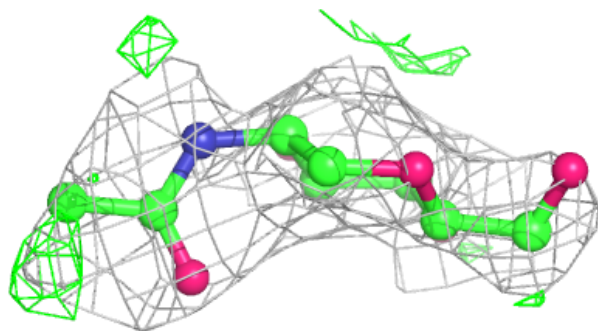
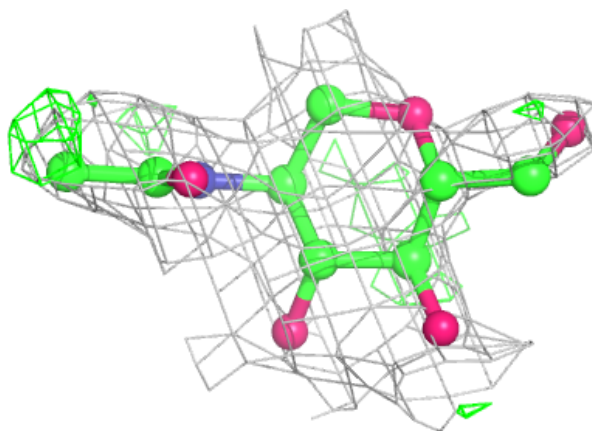
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
9	NAG	D	201	14/15	0.63	0.40	78,101,107,109	0
9	NAG	C	501	14/15	0.78	0.29	72,87,92,94	0
9	NAG	C	502	14/15	0.79	0.26	88,101,106,107	0
9	NAG	E	509	14/15	0.82	0.18	62,72,78,79	0
9	NAG	A	404	14/15	0.84	0.15	61,74,81,82	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.

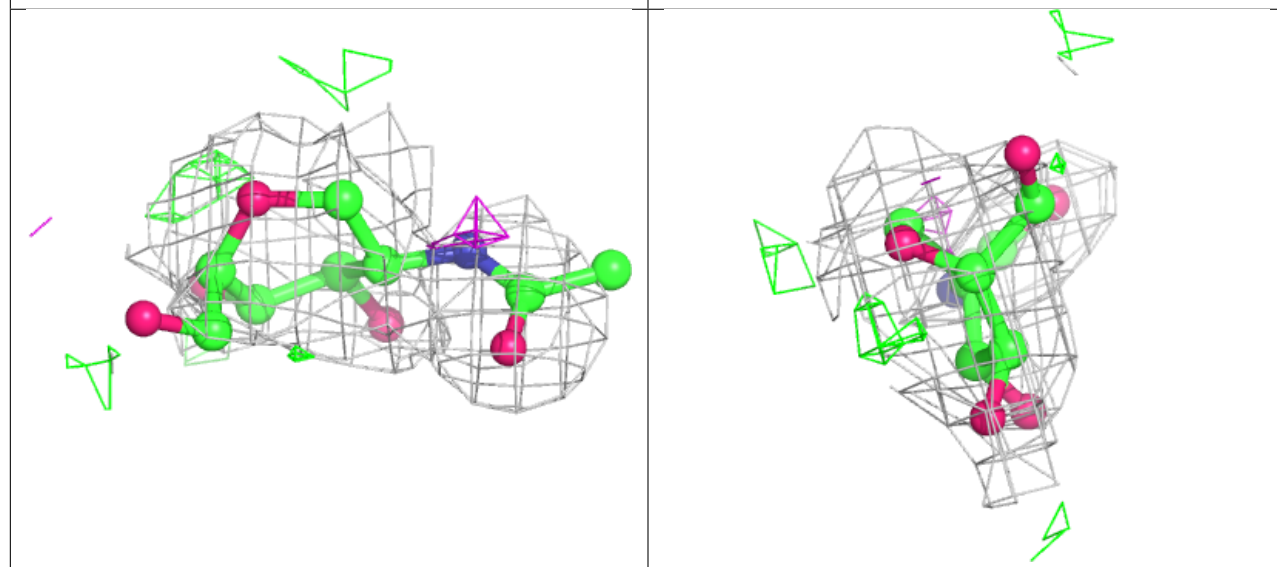
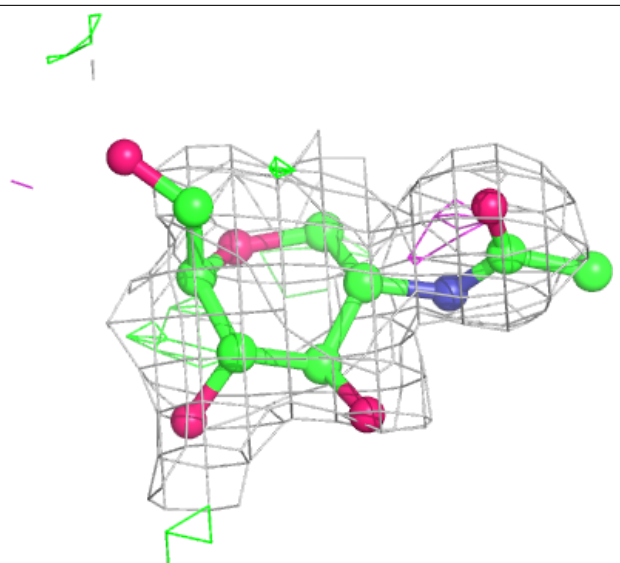
Electron density around NAG D 201:

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



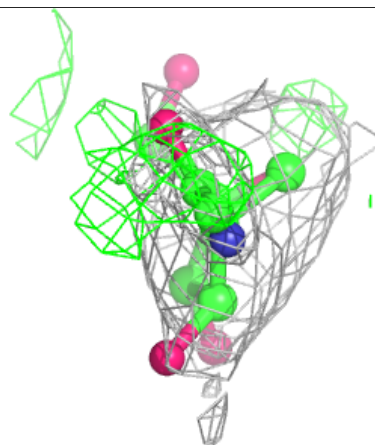
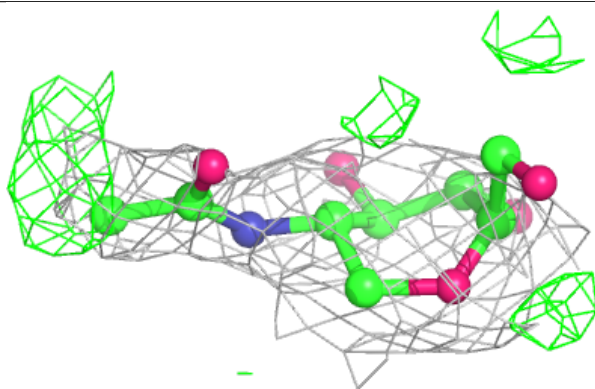
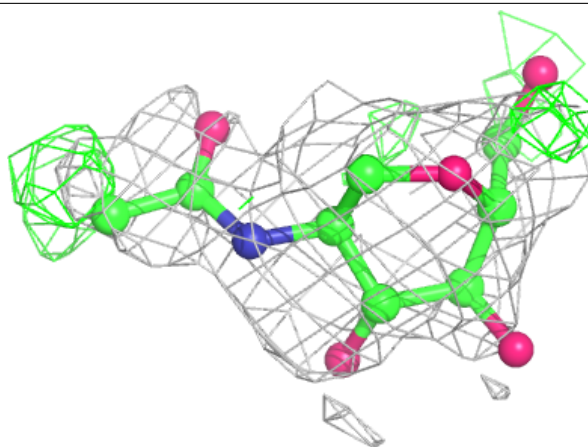
Electron density around NAG C 501:

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

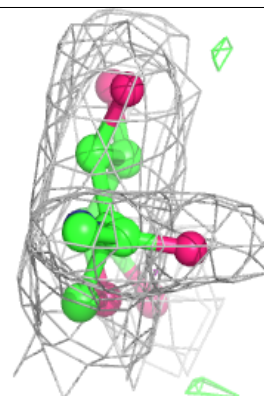
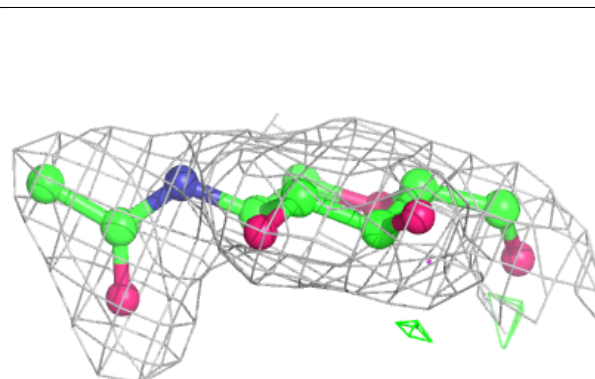
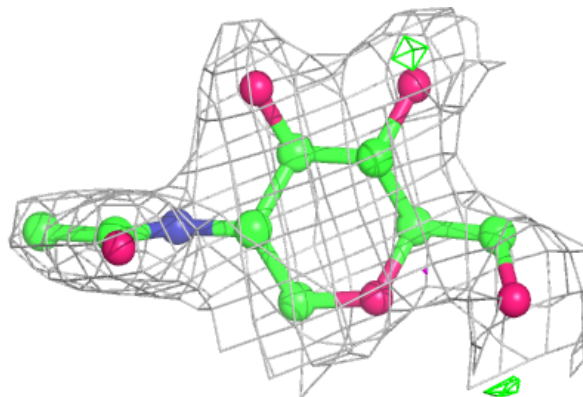


Electron density around NAG C 502:

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

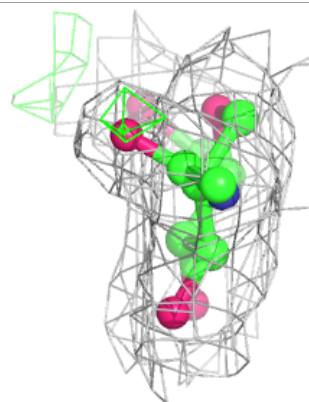
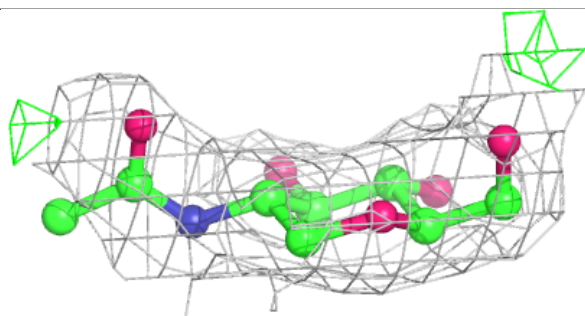
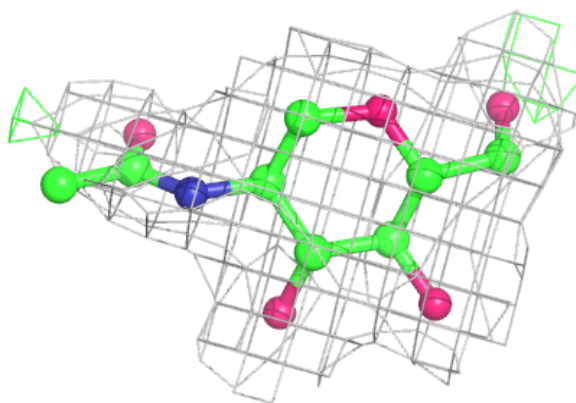
**Electron density around NAG E 509:**

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



Electron density around NAG A 404:

$2mF_o-DF_c$ (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.