



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 23, 2026 – 11:02 PM UTC

PDB ID : 8D3T / pdb_00008d3t
Title : Crystal structure of GalS1 from Populus trichocarpas
Authors : Pereira, J.H.; Prabhakar, P.K.; Urbanowicz, B.R.; Adams, P.D.
Deposited on : 2022-06-01
Resolution : 2.37 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

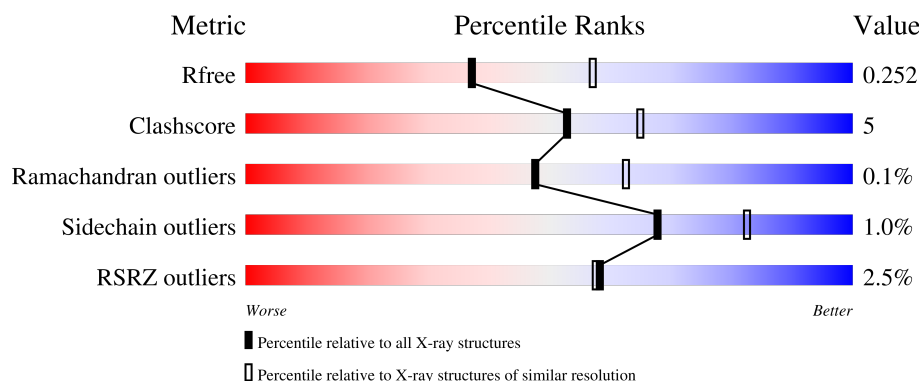
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 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.



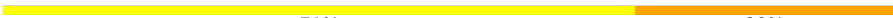
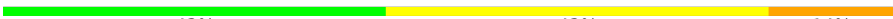
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	424	 3% 81% 12% 7%
1	B	424	 2% 84% 9% 7%
1	C	424	 .% 84% 8% 7%
1	D	424	 3% 82% 11% 8%
2	E	2	 50% 50%

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Mol	Chain	Length	Quality of chain
2	G	2	 100%
2	I	2	 100%
2	K	2	 50% 50%
3	F	7	 71% 29%
3	H	7	 43% 43% 14%
3	J	7	 43% 43% 14%
4	L	7	 14% 29% 57%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 27020 atoms, of which 12810 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 Galactan synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	393	Total	C	H	N	O	S	0	0	0
			6279	2076	3048	554	588	13			
1	B	394	Total	C	H	N	O	S	0	0	0
			6291	2080	3052	555	591	13			
1	C	394	Total	C	H	N	O	S	0	0	0
			6292	2080	3053	555	591	13			
1	D	392	Total	C	H	N	O	S	0	0	0
			6265	2071	3041	553	587	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP A0A2K2AMS7
B	1	GLY	-	expression tag	UNP A0A2K2AMS7
C	1	GLY	-	expression tag	UNP A0A2K2AMS7
D	1	GLY	-	expression tag	UNP A0A2K2AMS7

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



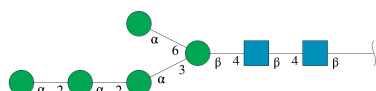
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	E	2	Total	C	H	N	O		0	0	0
			53	16	25	2	10				
2	G	2	Total	C	H	N	O		0	0	0
			53	16	25	2	10				
2	I	2	Total	C	H	N	O		0	0	0
			53	16	25	2	10				

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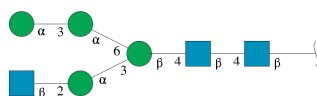
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	2	Total	C	H	N	O	0	0	0
			53	16	25	2	10			

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



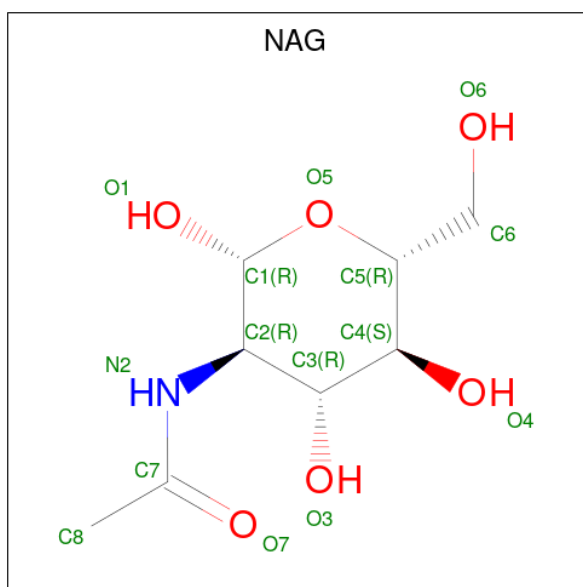
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	7	Total	C	H	N	O	0	0	0
			153	46	70	2	35			
3	H	7	Total	C	H	N	O	0	0	0
			152	46	69	2	35			
3	J	7	Total	C	H	N	O	0	0	0
			153	46	70	2	35			

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	7	Total	C	H	N	O	0	0	0
			159	48	73	3	35			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	A	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	B	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	B	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	B	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	B	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	C	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	C	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	C	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	C	1	Total	C	H	N	O	0	0
			27	8	13	1	5		

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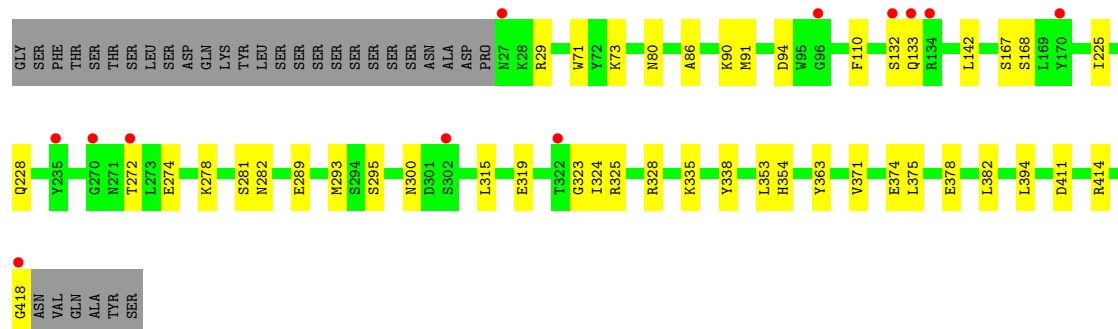
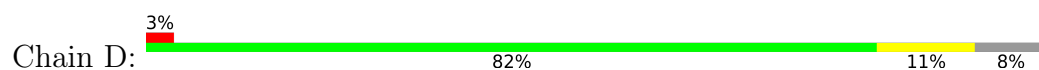
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	D	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	D	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	D	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
5	D	1	Total	C	H	N	O	0	0
			27	8	13	1	5		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	141	Total	O	0	0
			141	141		
6	B	145	Total	O	0	0
			145	145		
6	C	171	Total	O	0	0
			171	171		
6	D	121	Total	O	0	0
			121	121		

- Molecule 1: Galactan synthase





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



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



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



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



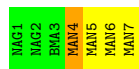
- Molecule 3: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





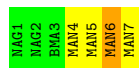
- Molecule 3: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H: 43% 43% 14%



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

Chain J: 43% 43% 14%



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

Chain L: 14% 29% 57%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	70.40Å 83.42Å 94.69Å 76.16° 69.67° 88.57°	Depositor
Resolution (Å)	80.84 – 2.37 80.84 – 2.37	Depositor EDS
% Data completeness (in resolution range)	98.2 (80.84-2.37) 98.2 (80.84-2.37)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.37Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.195 , 0.252 0.195 , 0.252	Depositor DCC
R_{free} test set	4011 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	36.6	Xtriage
Anisotropy	0.331	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.007 for h,-k,h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	27020	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 4.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, 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	0.17	0/3333	0.34	0/4525
1	B	0.12	0/3341	0.30	0/4537
1	C	0.11	0/3341	0.31	0/4537
1	D	0.10	0/3325	0.29	0/4514
All	All	0.13	0/13340	0.31	0/18113

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3231	3048	3048	34	0
1	B	3239	3052	3054	29	0
1	C	3239	3053	3052	28	0
1	D	3224	3041	3042	34	0
2	E	28	25	25	2	0
2	G	28	25	25	0	0
2	I	28	25	25	1	0
2	K	28	25	25	0	0
3	F	83	70	70	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	83	69	70	1	0
3	J	83	70	70	1	0
4	L	86	73	73	3	0
5	A	56	52	52	5	0
5	B	70	65	65	11	0
5	C	70	65	65	9	0
5	D	56	52	52	0	0
6	A	141	0	0	2	0
6	B	145	0	0	0	0
6	C	171	0	0	3	0
6	D	121	0	0	5	0
All	All	14210	12810	12813	131	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

The worst 5 of 131 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:115:ASN:O	1:A:115:ASN:OD1	1.73	1.06
1:C:171:GLY:HA2	5:C:502:NAG:H82	1.64	0.80
1:B:117:ASP:OD2	5:B:501:NAG:H83	1.83	0.79
5:A:503:NAG:HO6	3:F:4:MAN:HO4	1.26	0.75
5:C:503:NAG:H3	5:C:503:NAG:H83	1.70	0.72

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	391/424 (92%)	376 (96%)	15 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	392/424 (92%)	376 (96%)	15 (4%)	1 (0%)	36	48
1	C	392/424 (92%)	384 (98%)	7 (2%)	1 (0%)	36	48
1	D	390/424 (92%)	374 (96%)	16 (4%)	0	100	100
All	All	1565/1696 (92%)	1510 (96%)	53 (3%)	2 (0%)	48	62

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	300	ASN
1	B	282	ASN

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	337/365 (92%)	333 (99%)	4 (1%)	63	79
1	B	338/365 (93%)	336 (99%)	2 (1%)	78	88
1	C	338/365 (93%)	333 (98%)	5 (2%)	57	75
1	D	336/365 (92%)	334 (99%)	2 (1%)	78	88
All	All	1349/1460 (92%)	1336 (99%)	13 (1%)	68	82

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	172	ASN
1	C	282	ASN
1	D	319	GLU
1	C	319	GLU
1	D	300	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	159	GLN
1	D	27	ASN
1	D	271	ASN
1	D	65	HIS
1	B	133	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 ⓘ

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	$\# Z > 2$
2	NAG	E	1	1,2	14,14,15	0.44	0	17,19,21	0.75	0
2	NAG	E	2	2	14,14,15	0.24	0	17,19,21	0.47	0
3	NAG	F	1	3,1	14,14,15	0.34	0	17,19,21	0.44	0
3	NAG	F	2	3	14,14,15	0.60	1 (7%)	17,19,21	0.49	0
3	BMA	F	3	3	11,11,12	0.74	0	15,15,17	1.02	1 (6%)
3	MAN	F	4	3	11,11,12	0.67	0	15,15,17	1.11	1 (6%)
3	MAN	F	5	3	11,11,12	0.97	1 (9%)	15,15,17	1.21	1 (6%)
3	MAN	F	6	3	11,11,12	0.62	0	15,15,17	1.08	2 (13%)
3	MAN	F	7	3	11,11,12	0.65	0	15,15,17	0.92	2 (13%)
2	NAG	G	1	1,2	14,14,15	0.55	0	17,19,21	0.50	0
2	NAG	G	2	2	14,14,15	0.24	0	17,19,21	0.38	0
3	NAG	H	1	3,1	14,14,15	0.30	0	17,19,21	0.41	0
3	NAG	H	2	3	14,14,15	0.19	0	17,19,21	0.40	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BMA	H	3	3	11,11,12	0.44	0	15,15,17	0.94	0
3	MAN	H	4	3	11,11,12	0.77	0	15,15,17	1.08	1 (6%)
3	MAN	H	5	3	11,11,12	0.91	1 (9%)	15,15,17	1.12	1 (6%)
3	MAN	H	6	3	11,11,12	0.59	0	15,15,17	1.26	2 (13%)
3	MAN	H	7	3	11,11,12	0.60	0	15,15,17	1.17	2 (13%)
2	NAG	I	1	1,2	14,14,15	0.65	1 (7%)	17,19,21	0.70	0
2	NAG	I	2	2	14,14,15	0.30	0	17,19,21	0.57	0
3	NAG	J	1	3,1	14,14,15	0.47	0	17,19,21	0.63	0
3	NAG	J	2	3	14,14,15	0.36	0	17,19,21	0.46	0
3	BMA	J	3	3	11,11,12	0.49	0	15,15,17	0.90	0
3	MAN	J	4	3	11,11,12	0.59	0	15,15,17	1.41	2 (13%)
3	MAN	J	5	3	11,11,12	0.78	0	15,15,17	1.27	2 (13%)
3	MAN	J	6	3	11,11,12	0.63	0	15,15,17	1.17	2 (13%)
3	MAN	J	7	3	11,11,12	0.91	1 (9%)	15,15,17	2.02	3 (20%)
2	NAG	K	1	1,2	14,14,15	0.20	0	17,19,21	0.70	1 (5%)
2	NAG	K	2	2	14,14,15	0.43	0	17,19,21	0.49	0
4	NAG	L	1	4,1	14,14,15	0.36	0	17,19,21	0.59	0
4	NAG	L	2	4	14,14,15	0.55	0	17,19,21	0.57	0
4	BMA	L	3	4	11,11,12	0.97	0	15,15,17	1.78	1 (6%)
4	MAN	L	4	4	11,11,12	1.37	2 (18%)	15,15,17	1.67	2 (13%)
4	NAG	L	5	4	14,14,15	0.74	1 (7%)	17,19,21	0.57	0
4	MAN	L	6	4	11,11,12	0.99	0	15,15,17	1.15	2 (13%)
4	MAN	L	7	4	11,11,12	1.26	1 (9%)	15,15,17	1.63	2 (13%)

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	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
3	NAG	F	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	0/2/19/22	0/1/1/1
3	MAN	F	4	3	-	1/2/19/22	0/1/1/1
3	MAN	F	5	3	-	0/2/19/22	0/1/1/1
3	MAN	F	6	3	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MAN	F	7	3	-	0/2/19/22	0/1/1/1
2	NAG	G	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
3	NAG	H	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
3	BMA	H	3	3	-	0/2/19/22	0/1/1/1
3	MAN	H	4	3	-	2/2/19/22	0/1/1/1
3	MAN	H	5	3	-	0/2/19/22	0/1/1/1
3	MAN	H	6	3	-	0/2/19/22	0/1/1/1
3	MAN	H	7	3	-	0/2/19/22	0/1/1/1
2	NAG	I	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	2/6/23/26	0/1/1/1
3	NAG	J	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1
3	BMA	J	3	3	-	2/2/19/22	0/1/1/1
3	MAN	J	4	3	-	2/2/19/22	0/1/1/1
3	MAN	J	5	3	-	0/2/19/22	0/1/1/1
3	MAN	J	6	3	-	0/2/19/22	0/1/1/1
3	MAN	J	7	3	-	0/2/19/22	0/1/1/1
2	NAG	K	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	K	2	2	-	2/6/23/26	0/1/1/1
4	NAG	L	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	L	2	4	-	2/6/23/26	0/1/1/1
4	BMA	L	3	4	-	1/2/19/22	0/1/1/1
4	MAN	L	4	4	-	2/2/19/22	0/1/1/1
4	NAG	L	5	4	-	2/6/23/26	0/1/1/1
4	MAN	L	6	4	-	1/2/19/22	0/1/1/1
4	MAN	L	7	4	-	0/2/19/22	0/1/1/1

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	7	MAN	C1-C2	3.75	1.61	1.52
4	L	4	MAN	O5-C1	-2.98	1.38	1.43
4	L	4	MAN	C2-C3	2.74	1.56	1.52
3	F	5	MAN	O5-C1	-2.57	1.39	1.43
4	L	5	NAG	C1-C2	2.56	1.55	1.52

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	J	7	MAN	C1-O5-C5	6.35	120.70	112.19
4	L	3	BMA	C1-O5-C5	6.00	120.22	112.19
4	L	4	MAN	O2-C2-C1	4.56	119.66	109.22
4	L	7	MAN	C1-C2-C3	4.23	115.80	109.64
3	J	4	MAN	C1-O5-C5	4.20	117.82	112.19

There are no chirality outliers.

5 of 30 torsion outliers are listed below:

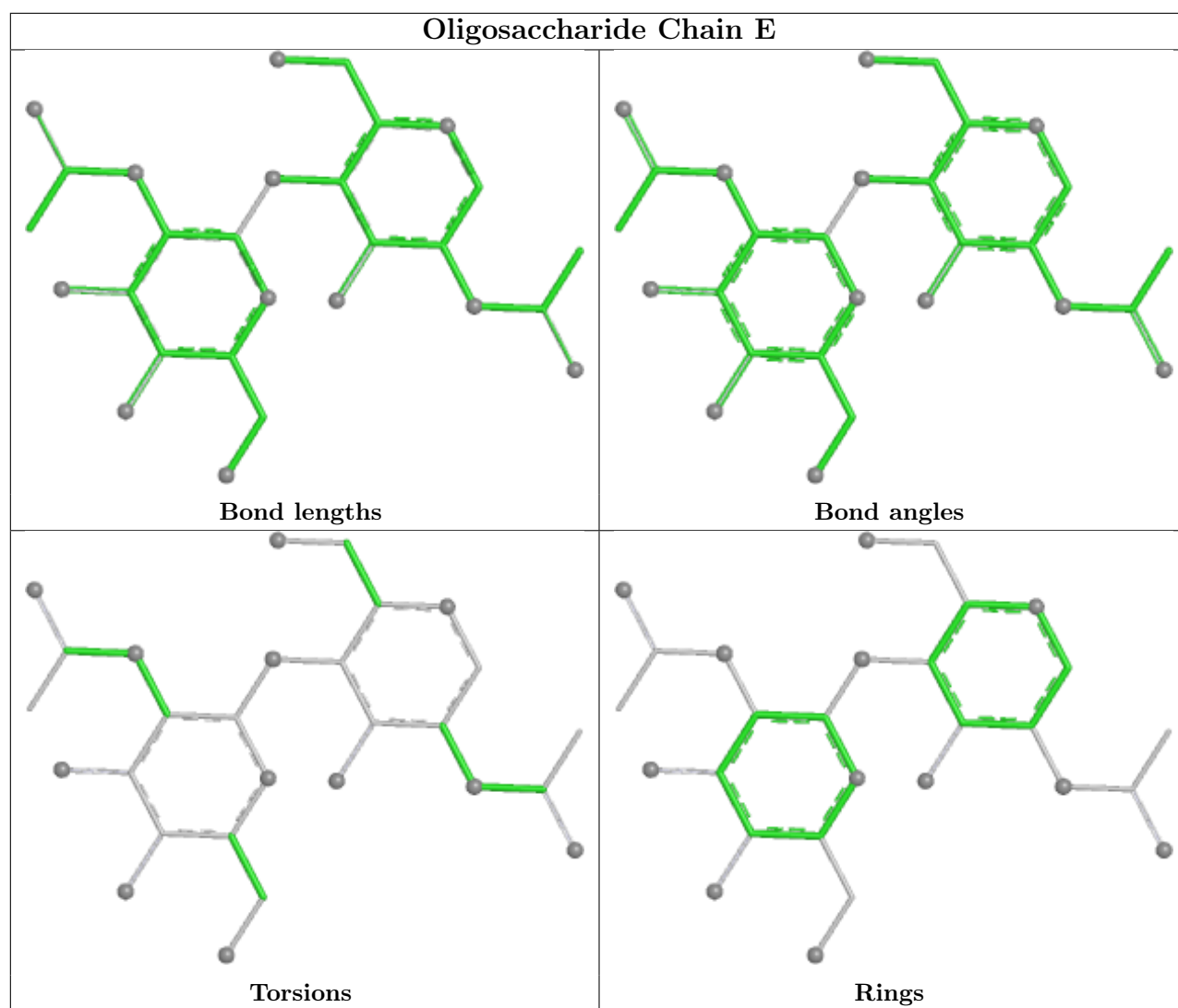
Mol	Chain	Res	Type	Atoms
4	L	5	NAG	C1-C2-N2-C7
3	H	4	MAN	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
3	J	4	MAN	C4-C5-C6-O6
4	L	4	MAN	O5-C5-C6-O6

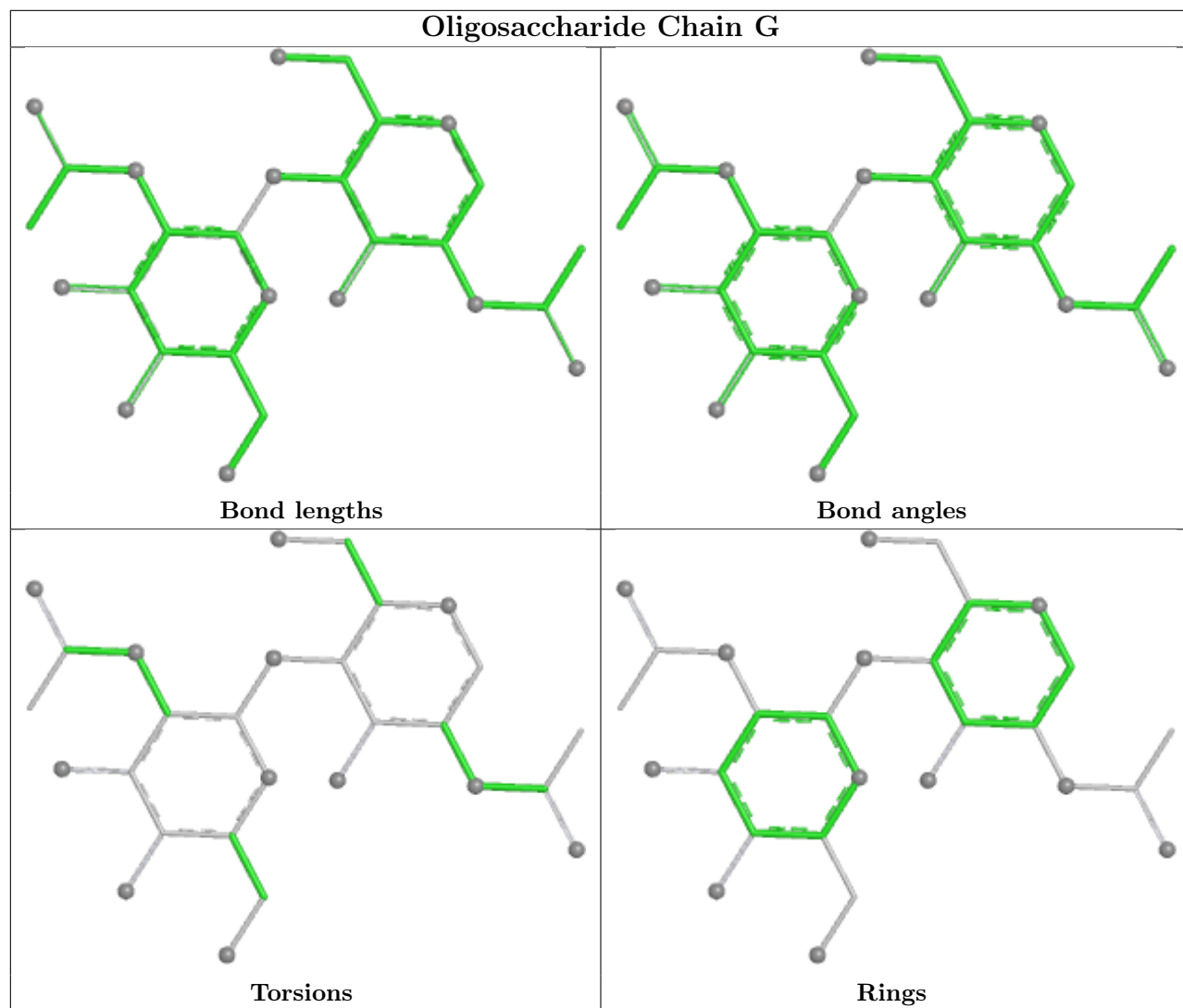
There are no ring outliers.

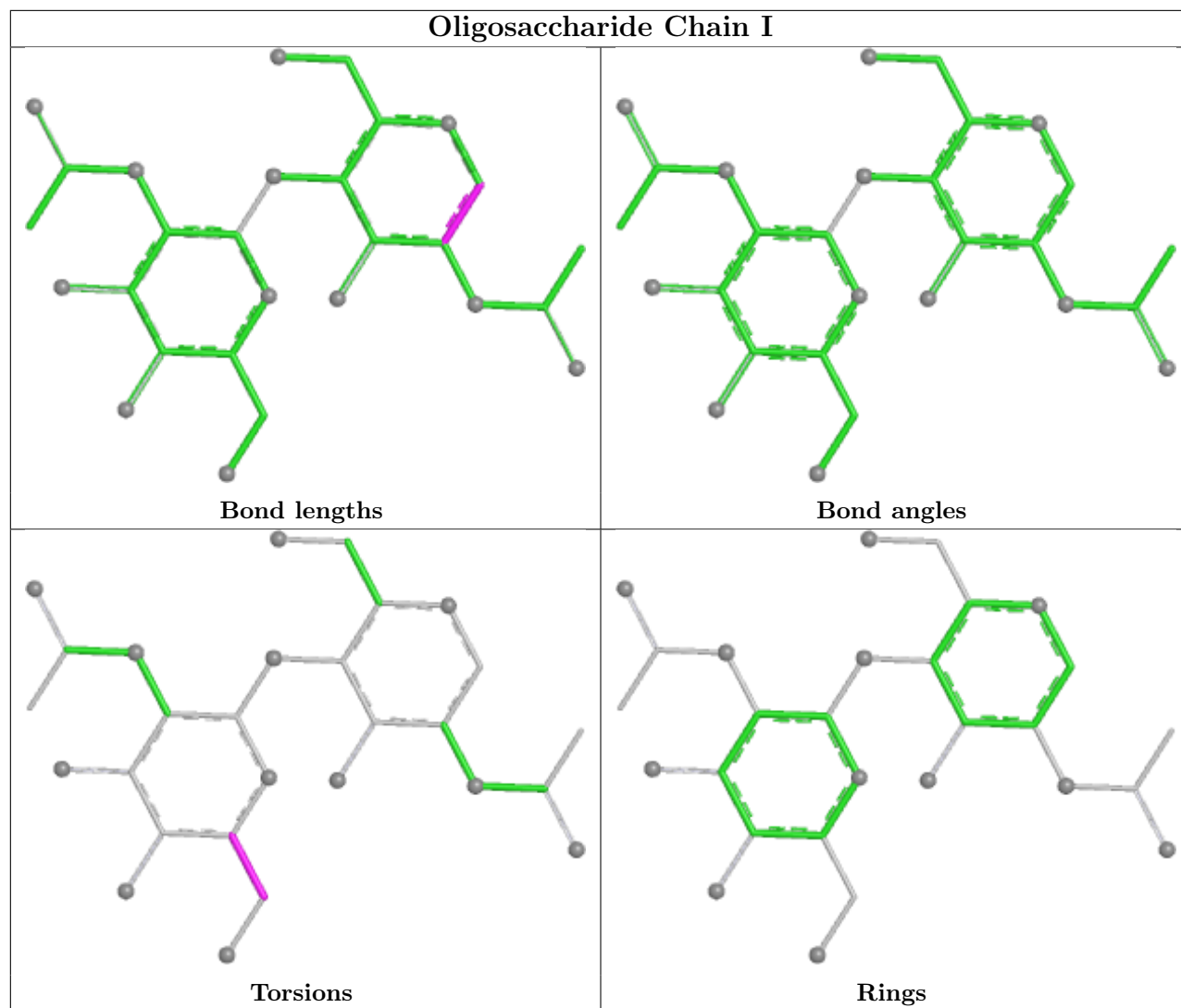
12 monomers are involved in 11 short contacts:

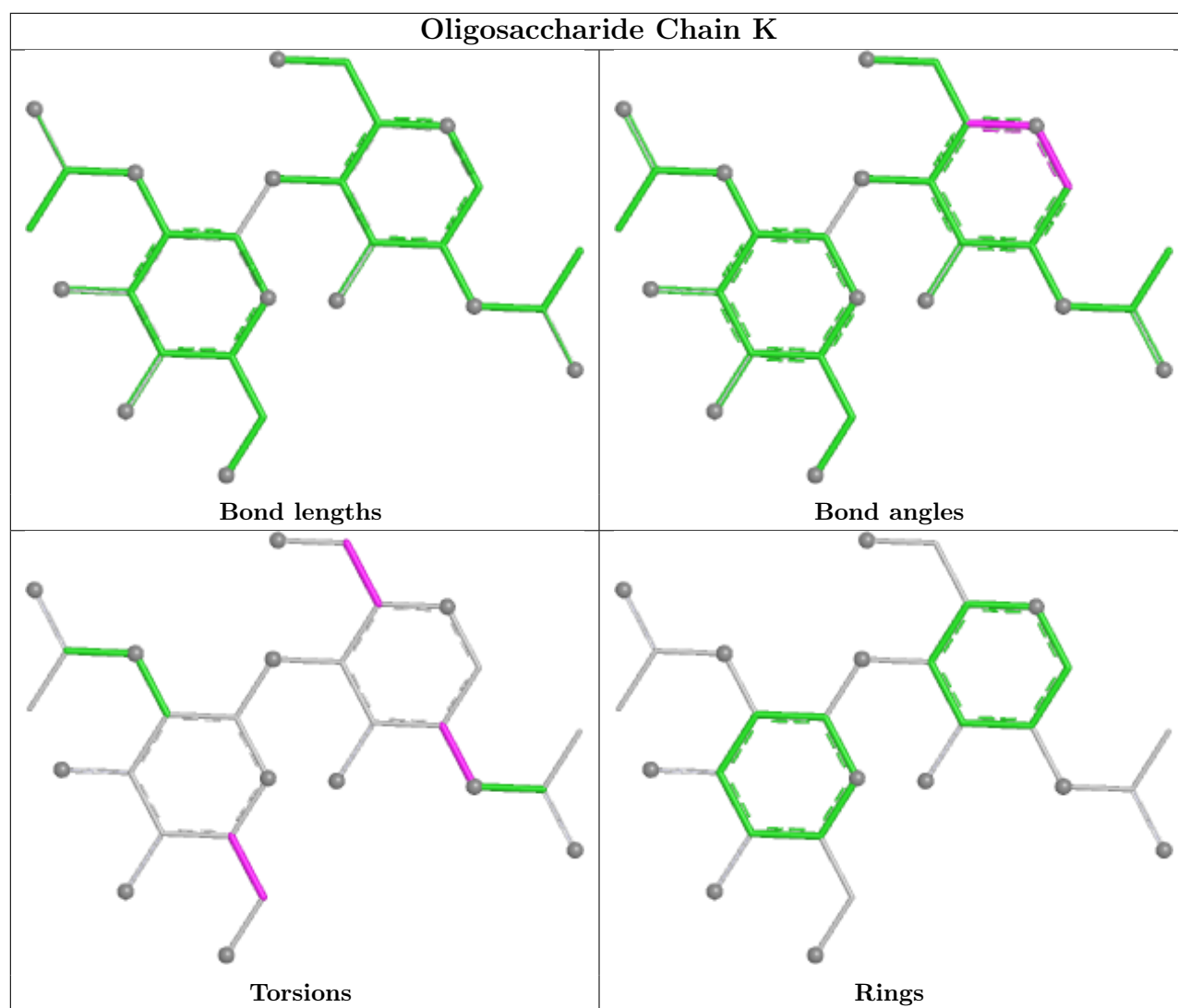
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	2	NAG	1	0
4	L	6	MAN	1	0
2	E	1	NAG	2	0
3	H	4	MAN	1	0
4	L	2	NAG	1	0
4	L	3	BMA	1	0
3	J	6	MAN	1	0
4	L	7	MAN	1	0
4	L	5	NAG	1	0
3	F	1	NAG	1	0
3	F	4	MAN	2	0
3	F	2	NAG	1	0

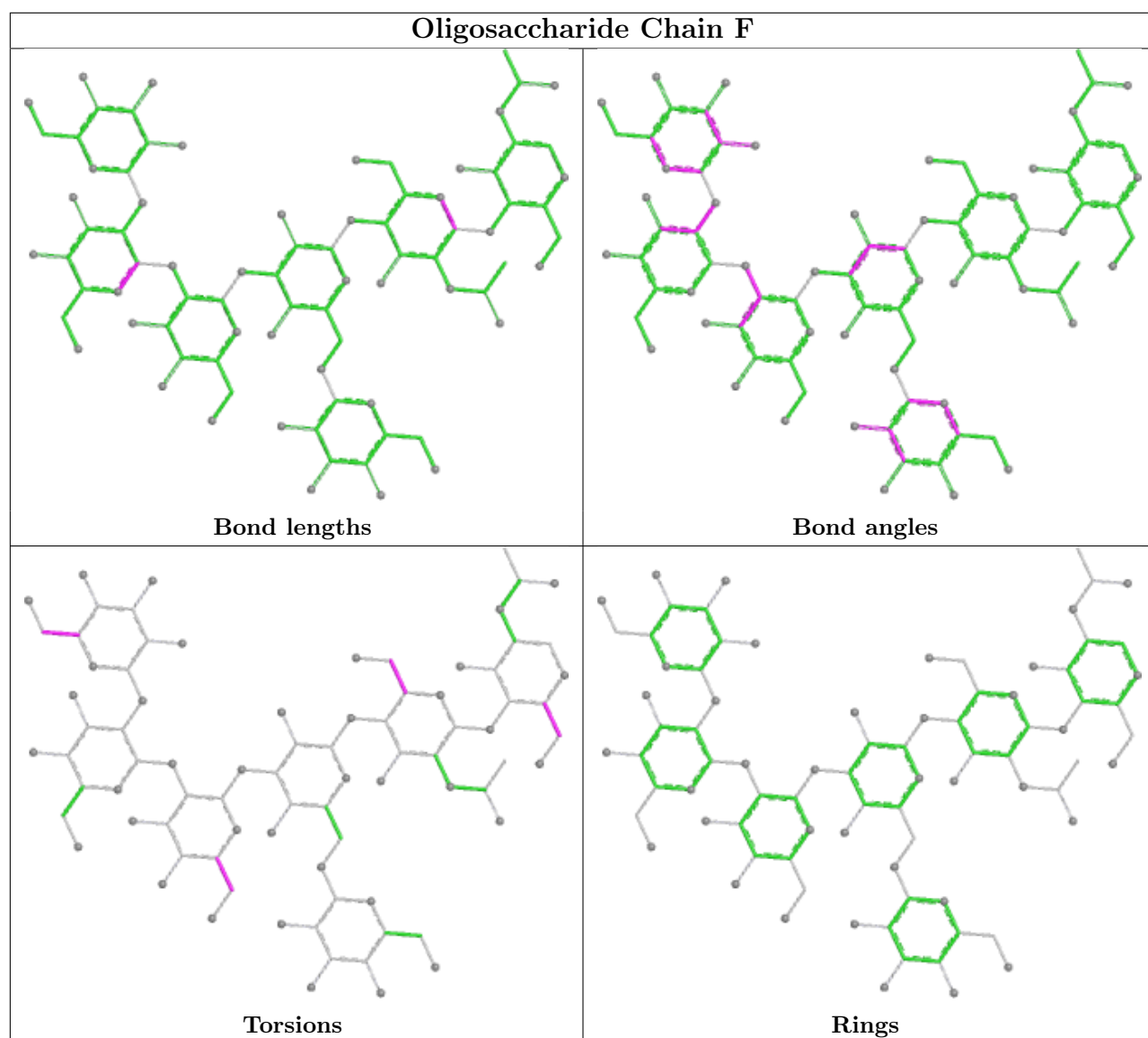
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

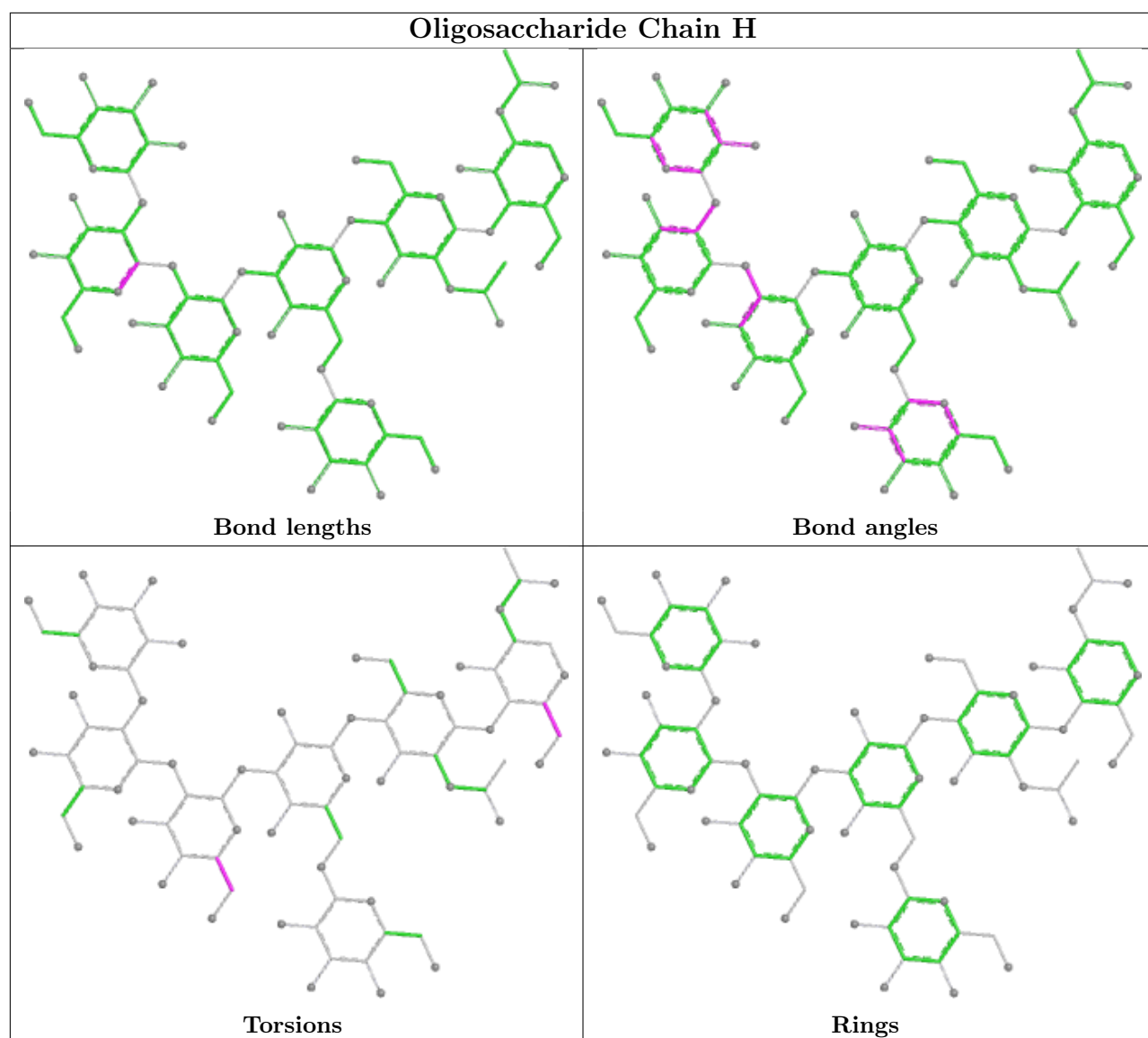


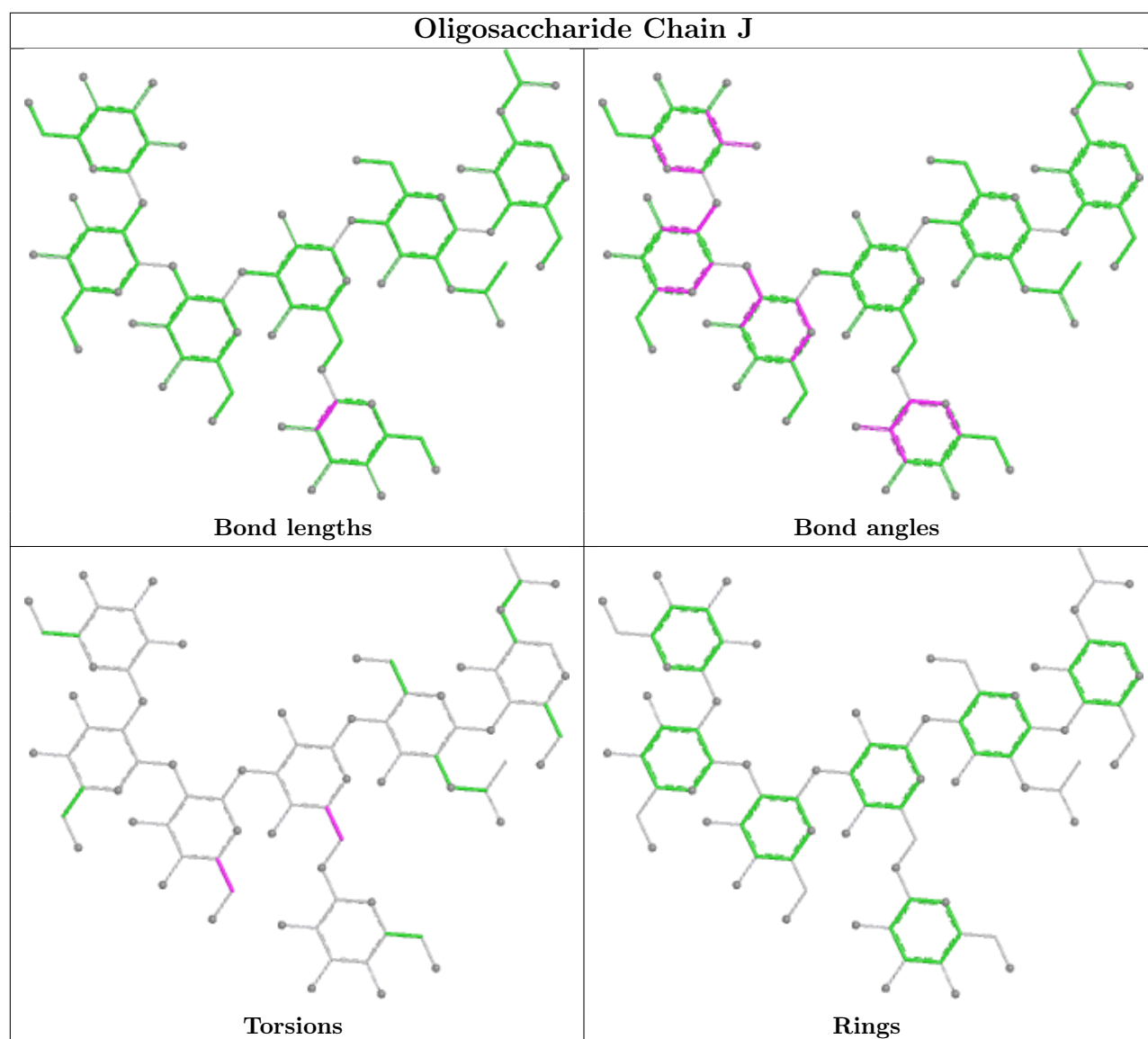


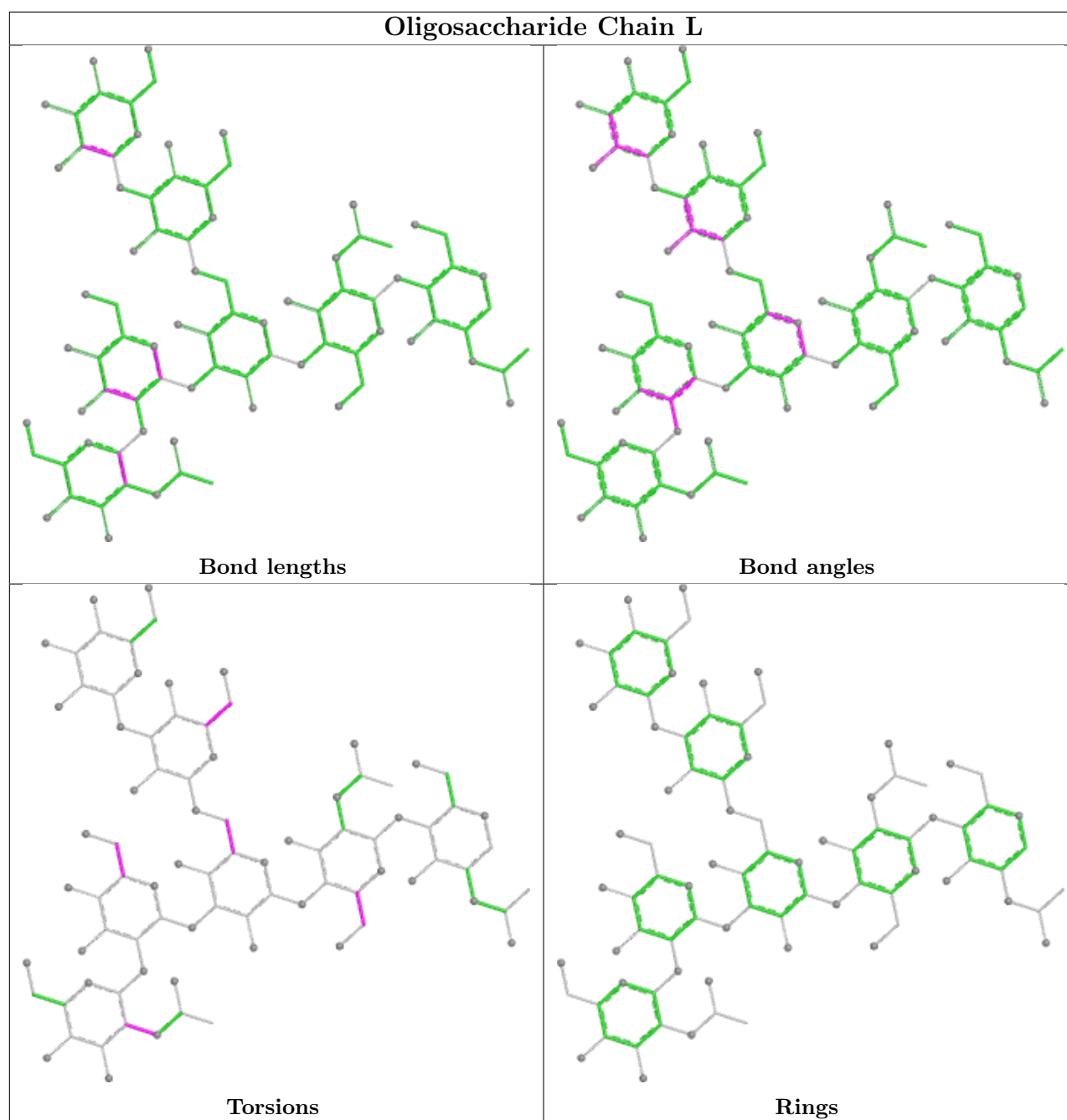












5.6 Ligand geometry [i](#)

18 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
5	NAG	B	503	1	14,14,15	0.34	0	17,19,21	0.41	0
5	NAG	D	502	1	14,14,15	0.26	0	17,19,21	0.57	0
5	NAG	A	504	1	14,14,15	0.65	0	17,19,21	0.86	1 (5%)
5	NAG	B	505	-	14,14,15	1.46	1 (7%)	17,19,21	1.03	1 (5%)
5	NAG	D	501	1	14,14,15	0.24	0	17,19,21	0.48	0
5	NAG	A	503	1	14,14,15	0.38	0	17,19,21	0.45	0
5	NAG	D	503	-	14,14,15	0.48	0	17,19,21	0.81	1 (5%)
5	NAG	D	504	1	14,14,15	0.61	0	17,19,21	1.38	2 (11%)
5	NAG	A	501	1	14,14,15	0.35	0	17,19,21	0.69	0
5	NAG	A	502	1	14,14,15	0.34	0	17,19,21	0.90	1 (5%)
5	NAG	C	502	1	14,14,15	0.46	0	17,19,21	0.84	1 (5%)
5	NAG	C	504	1	14,14,15	0.57	0	17,19,21	0.91	1 (5%)
5	NAG	C	505	-	14,14,15	0.39	0	17,19,21	0.83	1 (5%)
5	NAG	C	503	1	14,14,15	0.24	0	17,19,21	1.02	1 (5%)
5	NAG	C	501	1	14,14,15	0.89	1 (7%)	17,19,21	0.85	1 (5%)
5	NAG	B	504	-	14,14,15	0.41	0	17,19,21	0.67	0
5	NAG	B	502	1	14,14,15	0.37	0	17,19,21	0.80	1 (5%)
5	NAG	B	501	-	14,14,15	0.86	1 (7%)	17,19,21	0.94	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	B	503	1	-	2/6/23/26	0/1/1/1
5	NAG	D	502	1	-	0/6/23/26	0/1/1/1
5	NAG	A	504	1	-	2/6/23/26	0/1/1/1
5	NAG	B	505	-	-	0/6/23/26	0/1/1/1
5	NAG	D	501	1	-	2/6/23/26	0/1/1/1
5	NAG	A	503	1	-	0/6/23/26	0/1/1/1
5	NAG	D	503	-	-	4/6/23/26	0/1/1/1
5	NAG	D	504	1	-	2/6/23/26	0/1/1/1
5	NAG	A	501	1	-	0/6/23/26	0/1/1/1
5	NAG	A	502	1	-	2/6/23/26	0/1/1/1
5	NAG	C	502	1	-	1/6/23/26	0/1/1/1
5	NAG	C	504	1	-	4/6/23/26	0/1/1/1
5	NAG	C	505	-	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	C	503	1	-	6/6/23/26	0/1/1/1
5	NAG	C	501	1	-	2/6/23/26	0/1/1/1
5	NAG	B	504	-	-	2/6/23/26	0/1/1/1
5	NAG	B	502	1	-	2/6/23/26	0/1/1/1
5	NAG	B	501	-	-	3/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	505	NAG	O5-C1	-5.10	1.35	1.43
5	C	501	NAG	C1-C2	2.80	1.56	1.52
5	B	501	NAG	O5-C1	-2.76	1.39	1.43

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	504	NAG	C1-O5-C5	4.27	117.91	112.19
5	B	505	NAG	C3-C4-C5	3.51	116.60	110.23
5	C	504	NAG	C1-O5-C5	3.20	116.47	112.19
5	C	503	NAG	C2-N2-C7	3.02	126.95	122.90
5	C	501	NAG	C1-O5-C5	2.93	116.11	112.19

There are no chirality outliers.

5 of 36 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	502	NAG	O5-C5-C6-O6
5	A	502	NAG	O5-C5-C6-O6
5	D	503	NAG	C4-C5-C6-O6
5	A	502	NAG	C4-C5-C6-O6
5	C	503	NAG	O5-C5-C6-O6

There are no ring outliers.

12 monomers are involved in 25 short contacts:

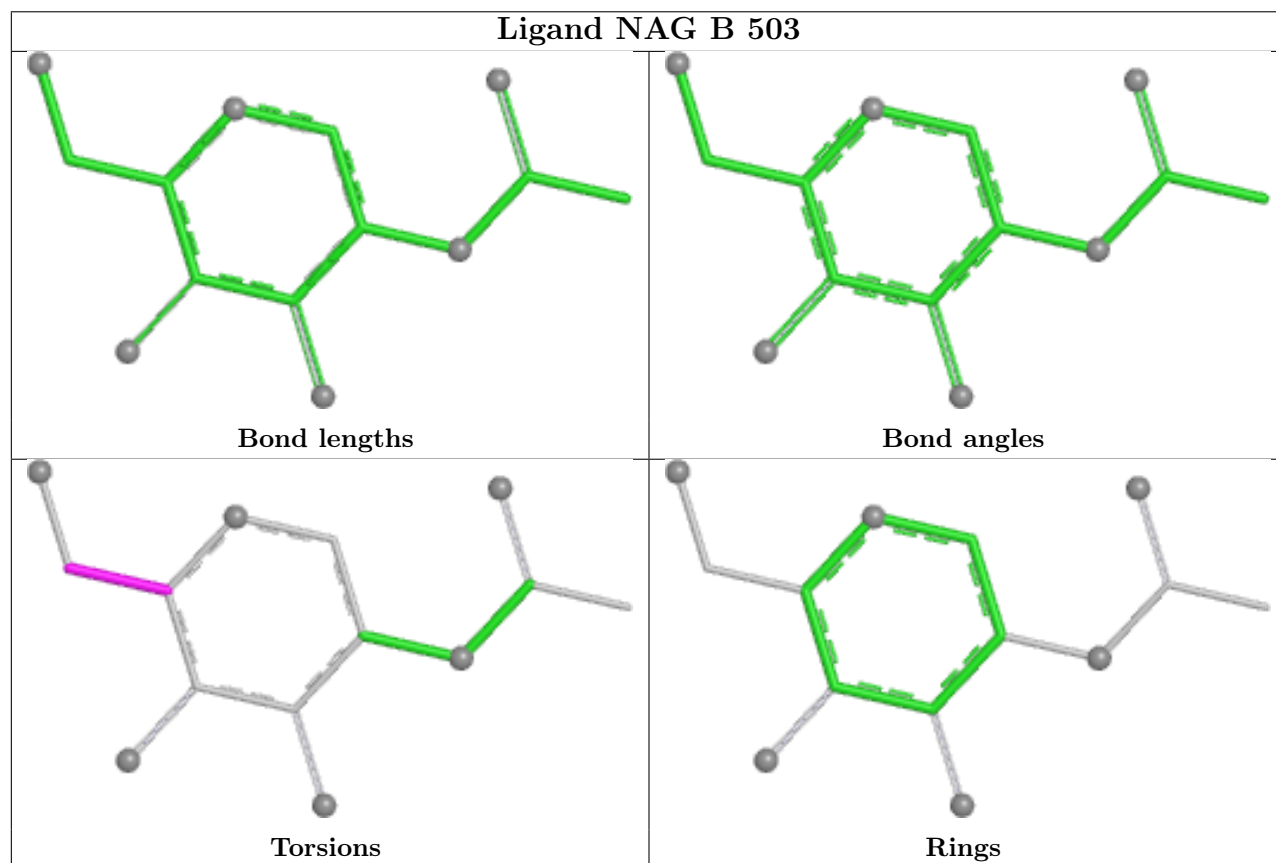
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	503	NAG	1	0
5	A	504	NAG	1	0
5	B	505	NAG	2	0
5	A	503	NAG	3	0

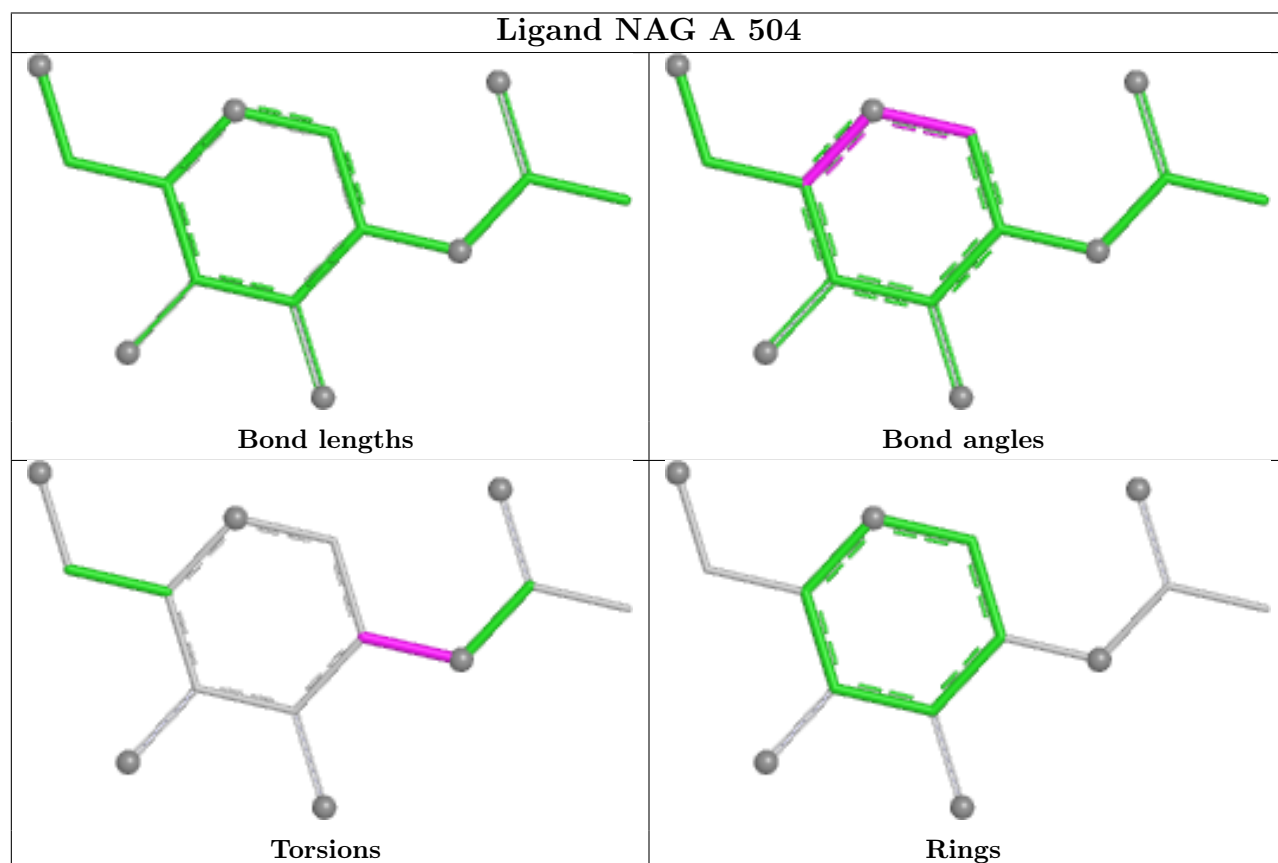
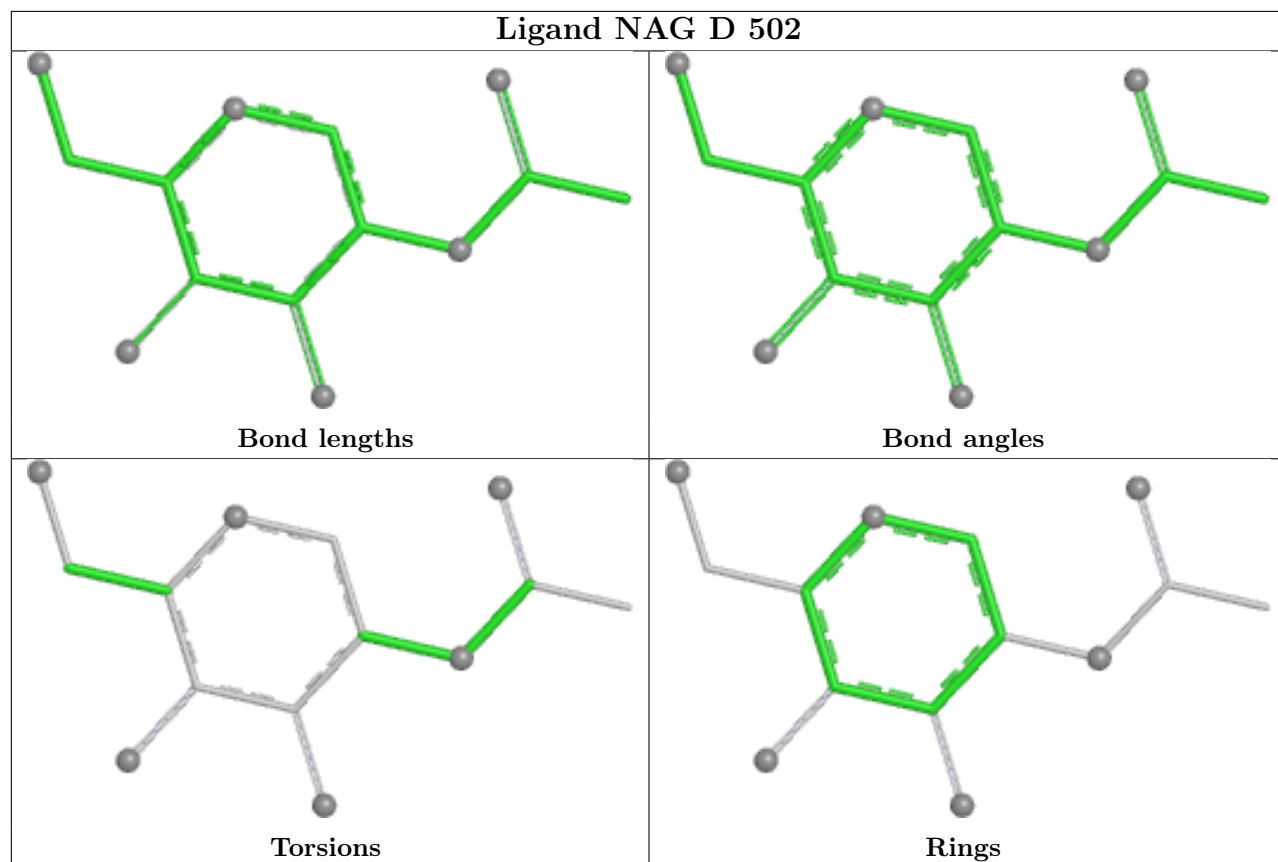
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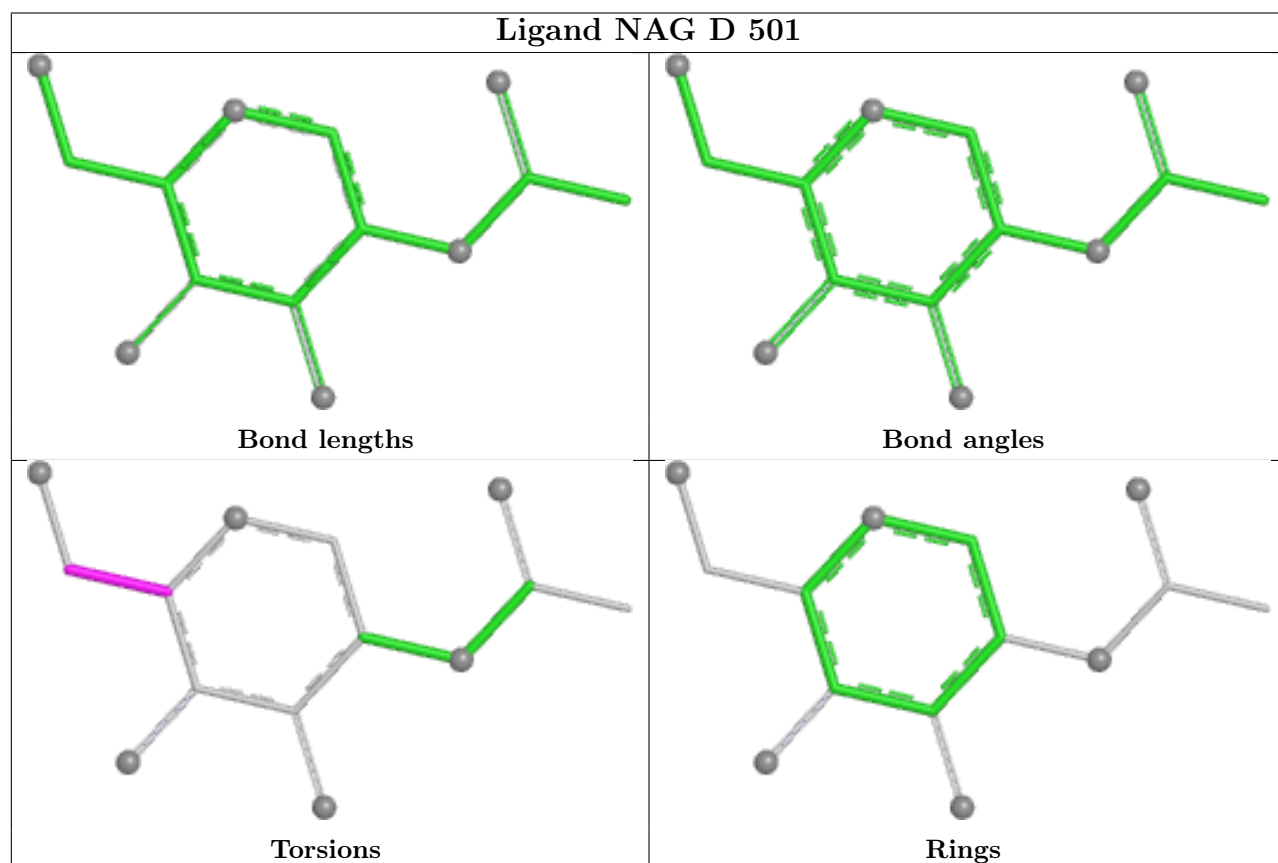
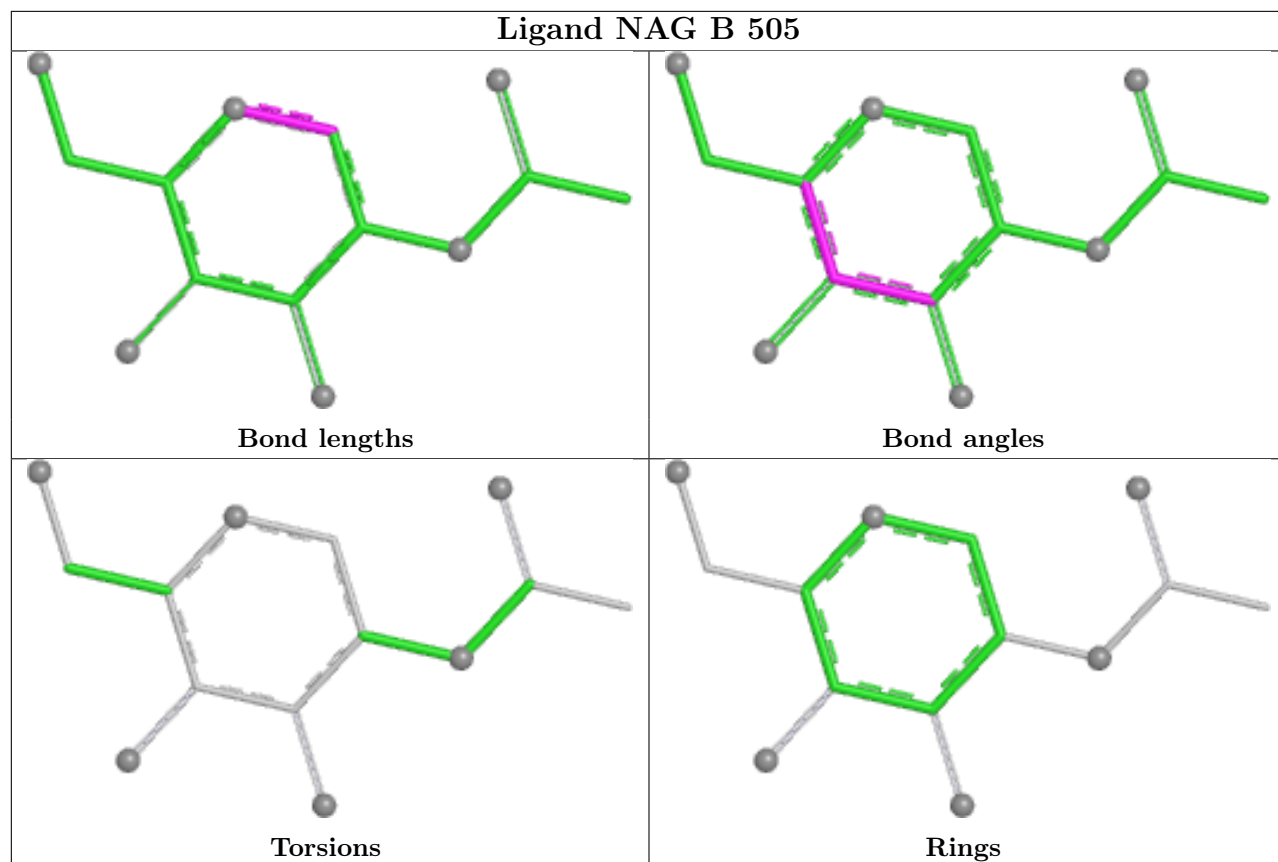
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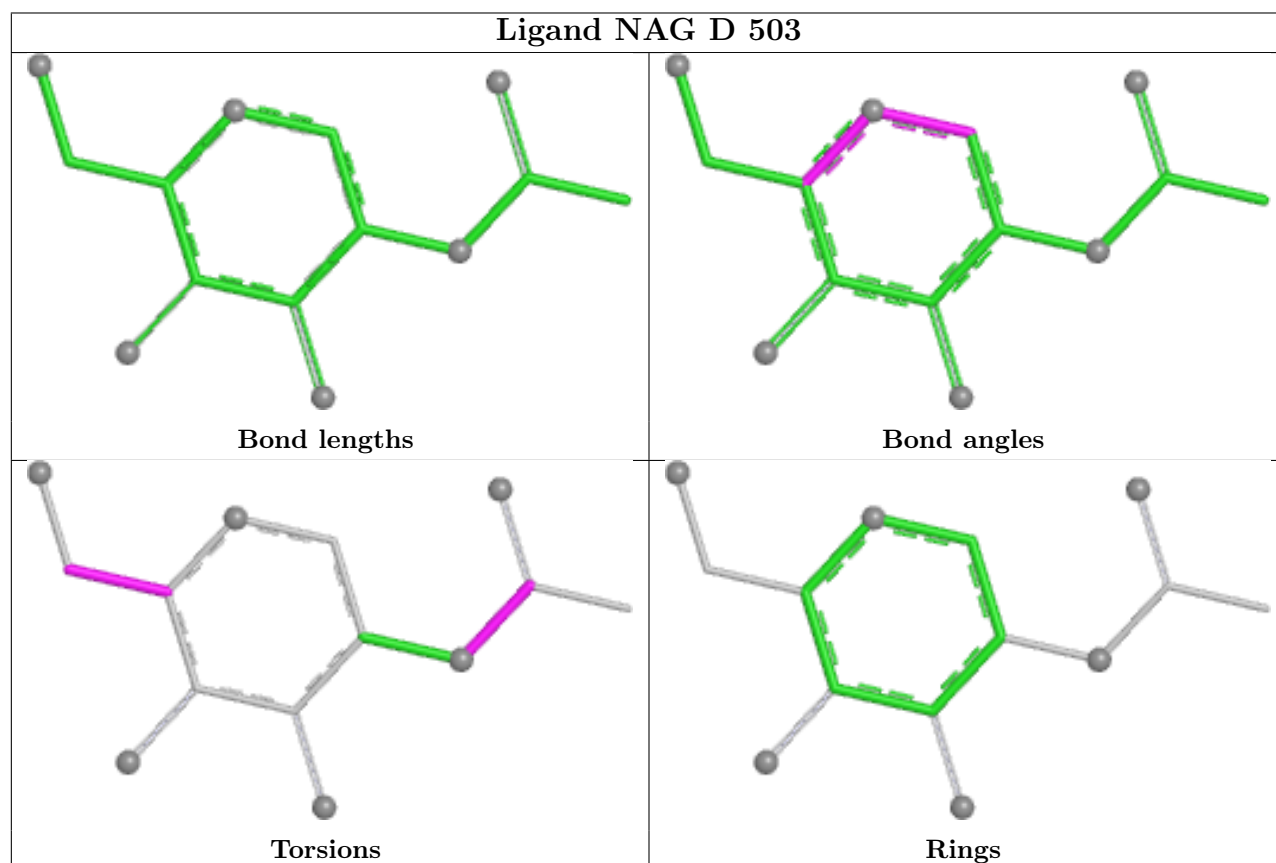
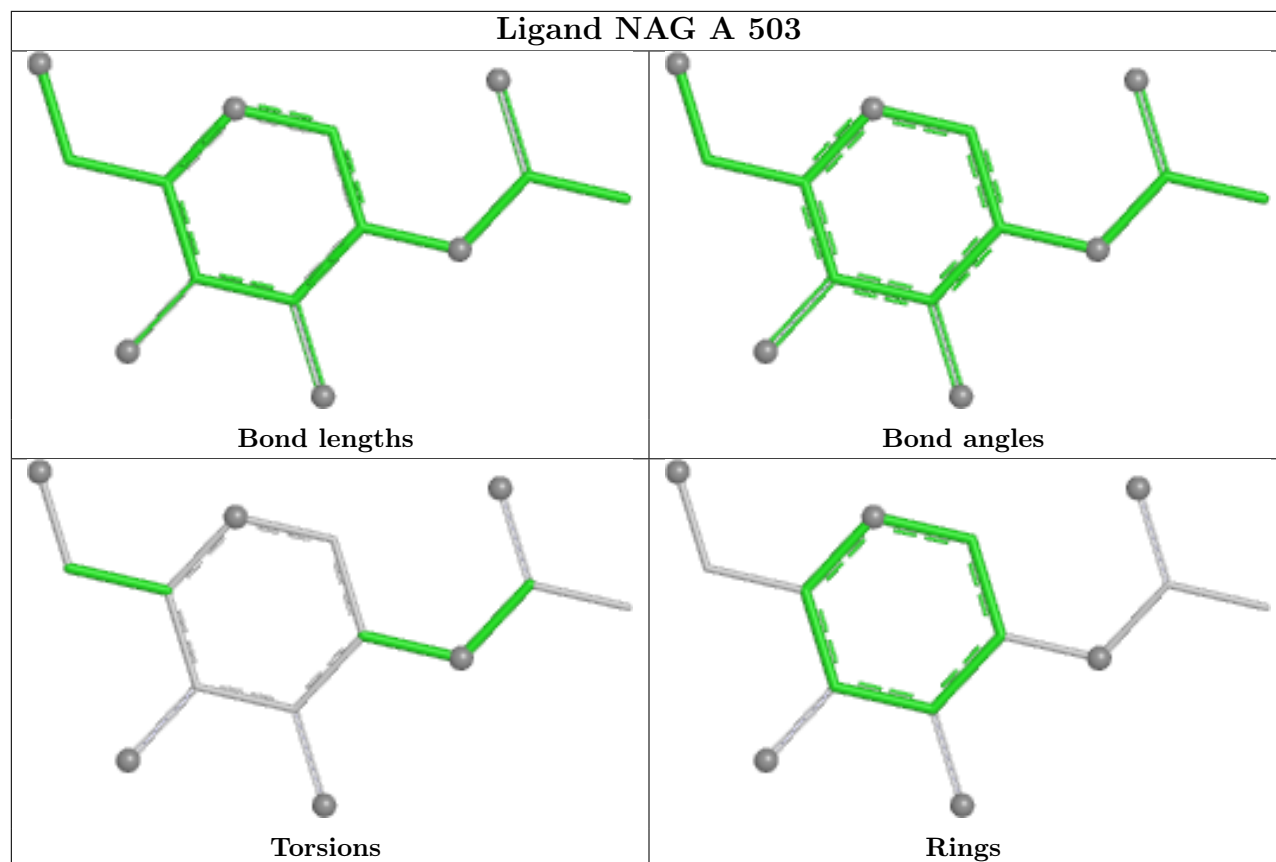
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	502	NAG	1	0
5	C	502	NAG	2	0
5	C	504	NAG	1	0
5	C	503	NAG	2	0
5	C	501	NAG	4	0
5	B	504	NAG	1	0
5	B	502	NAG	2	0
5	B	501	NAG	5	0

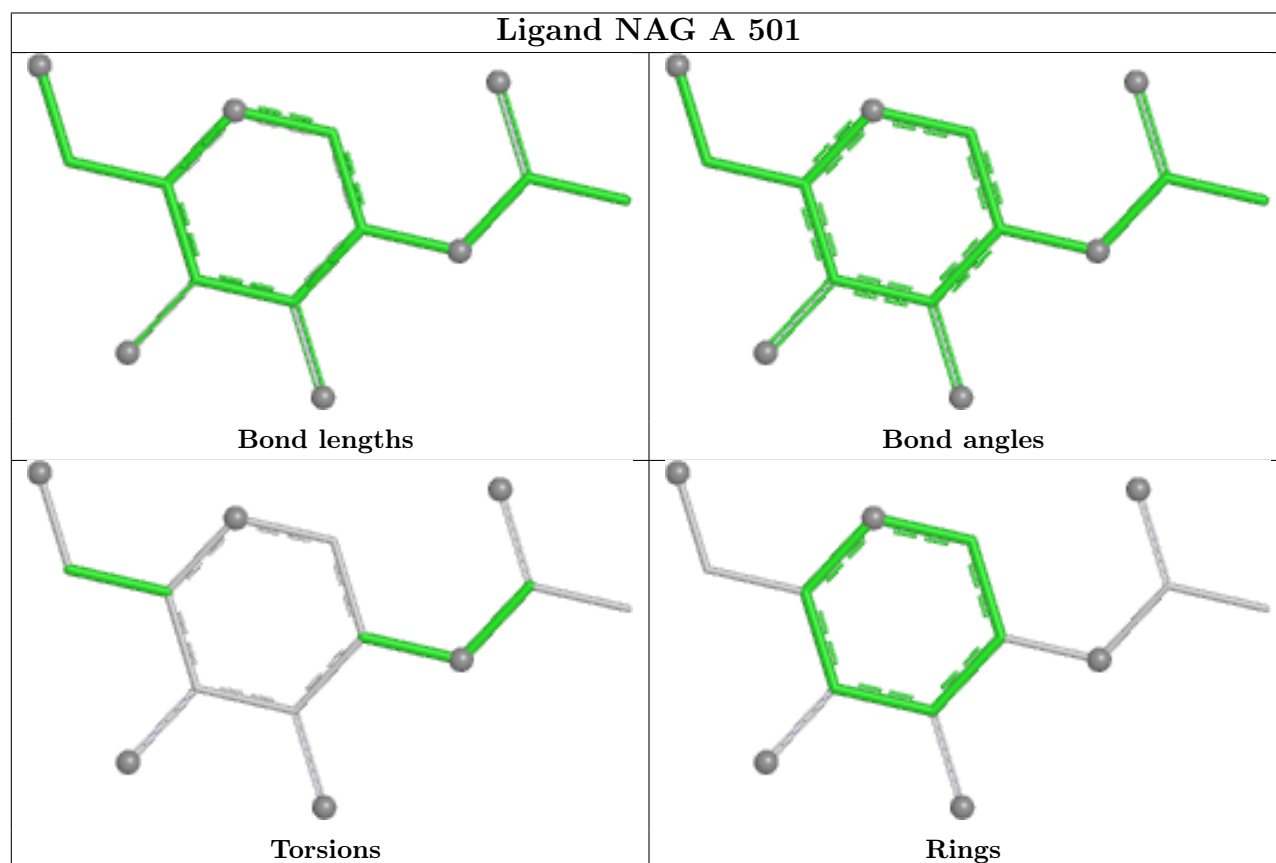
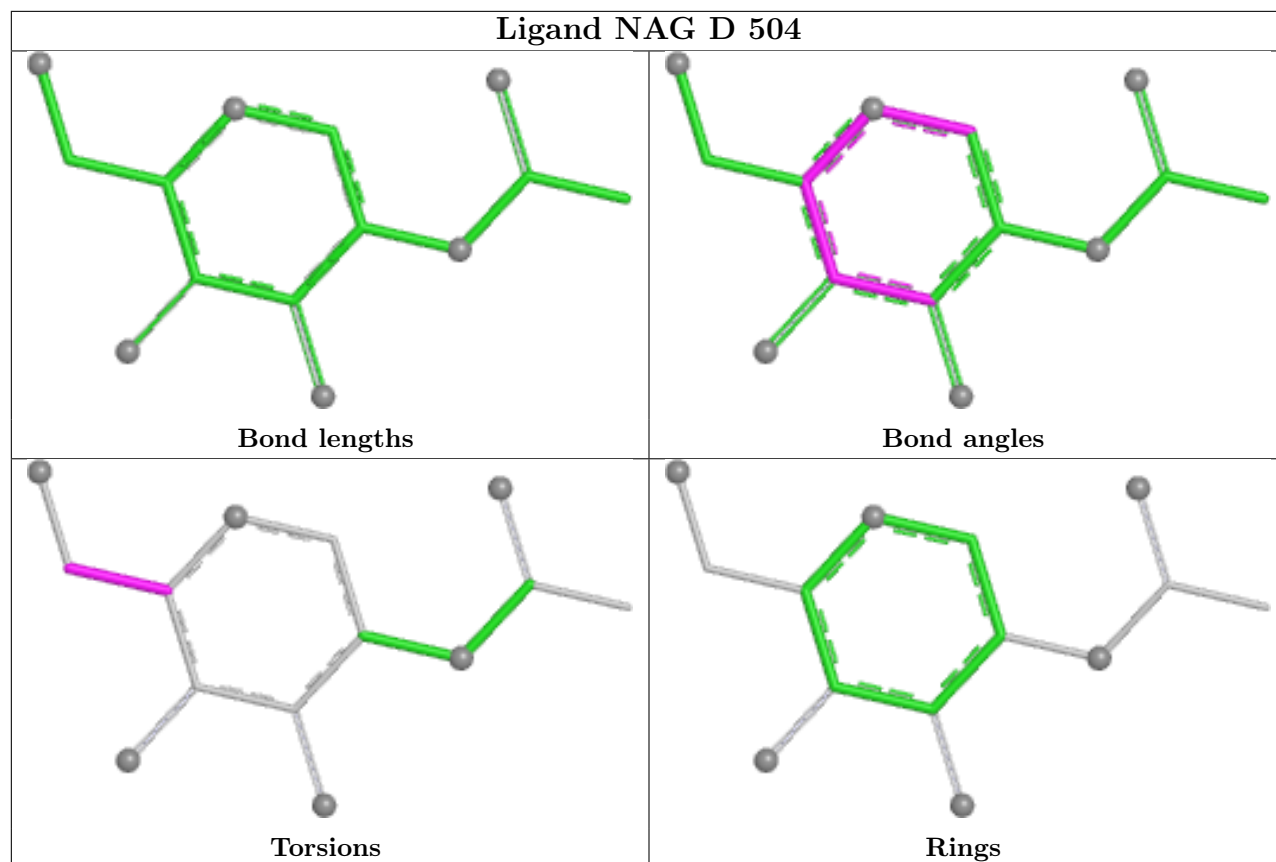
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

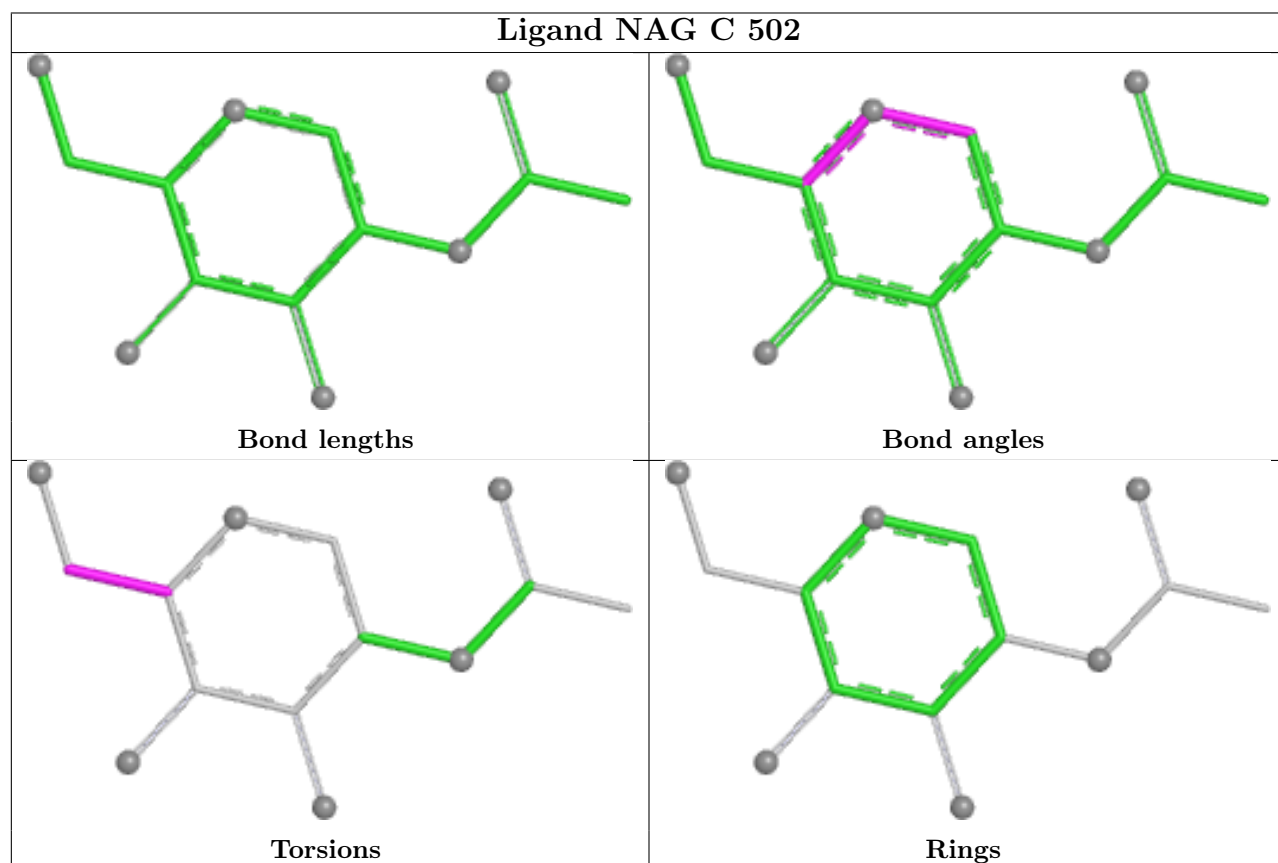
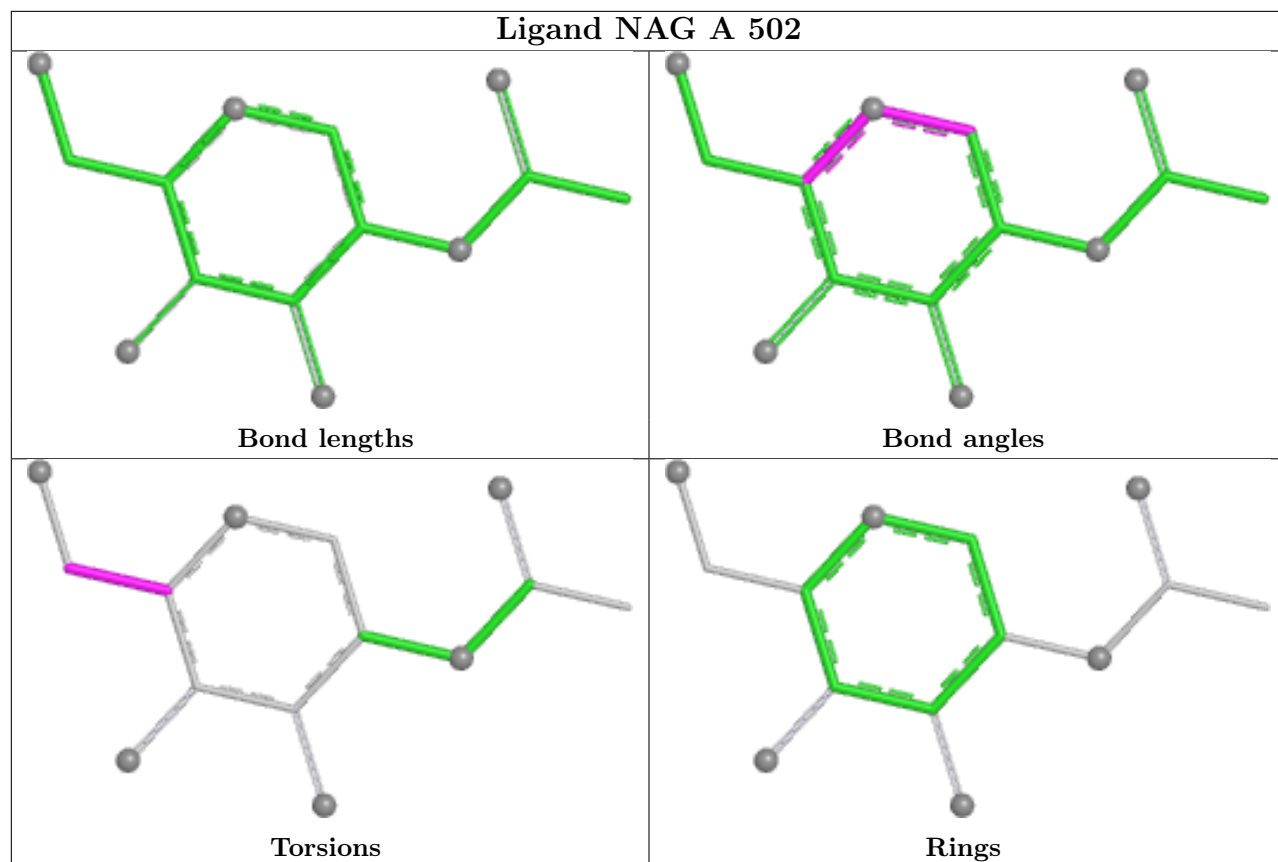


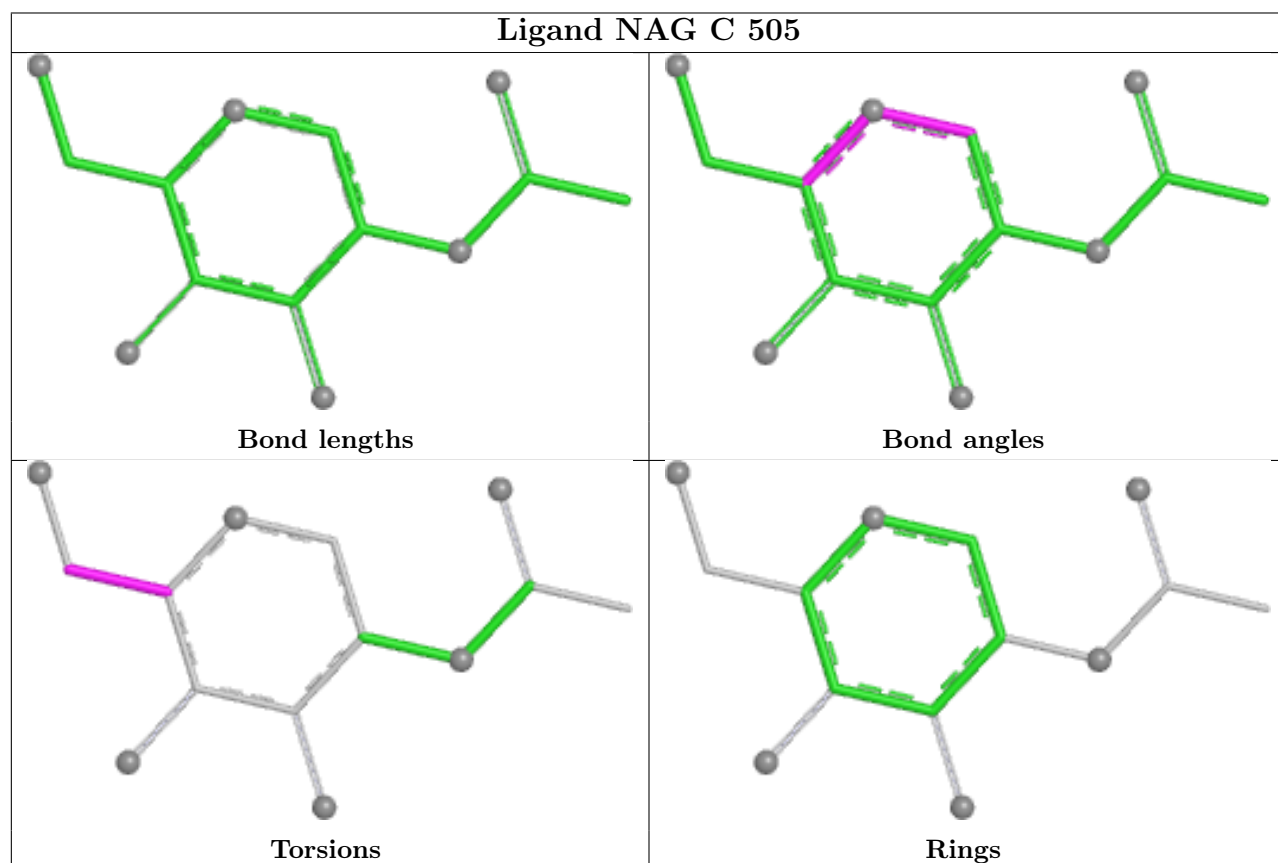
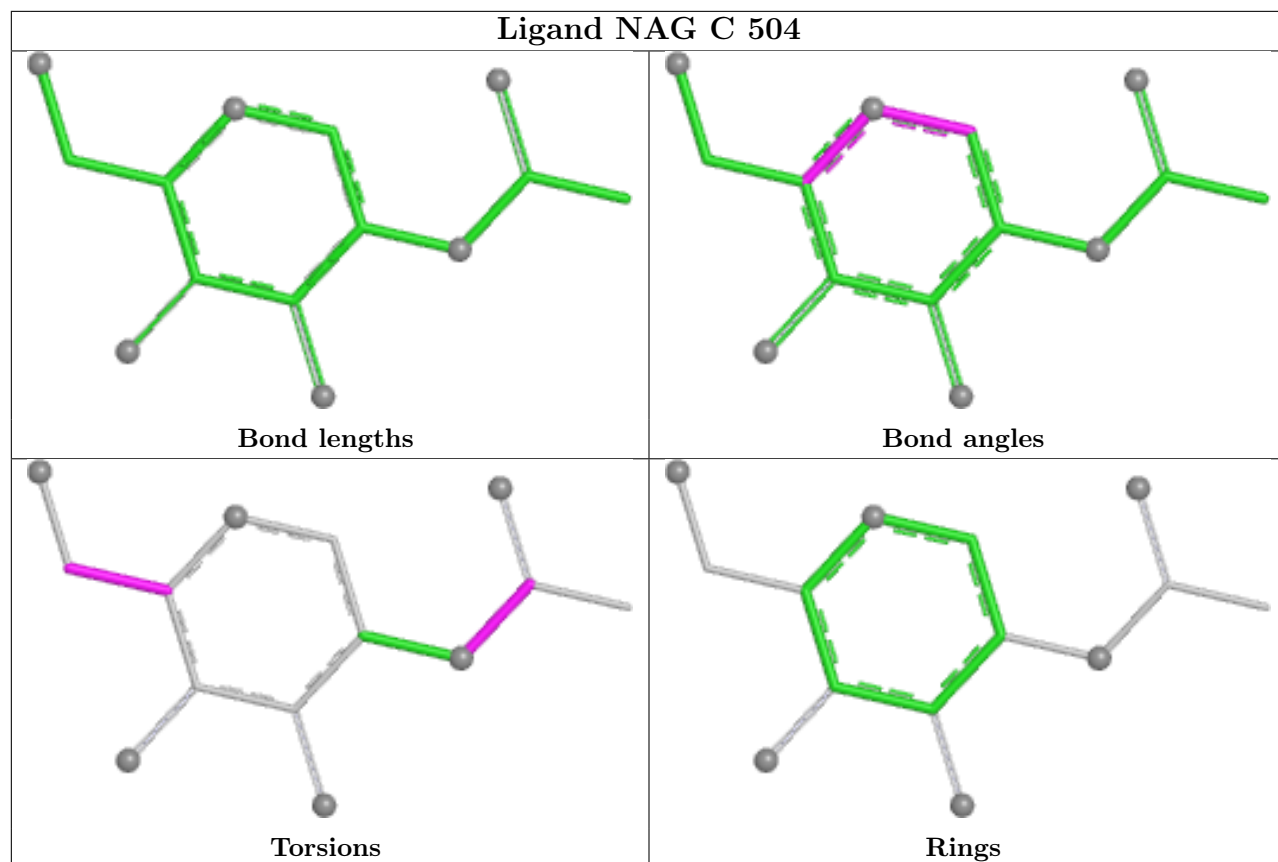


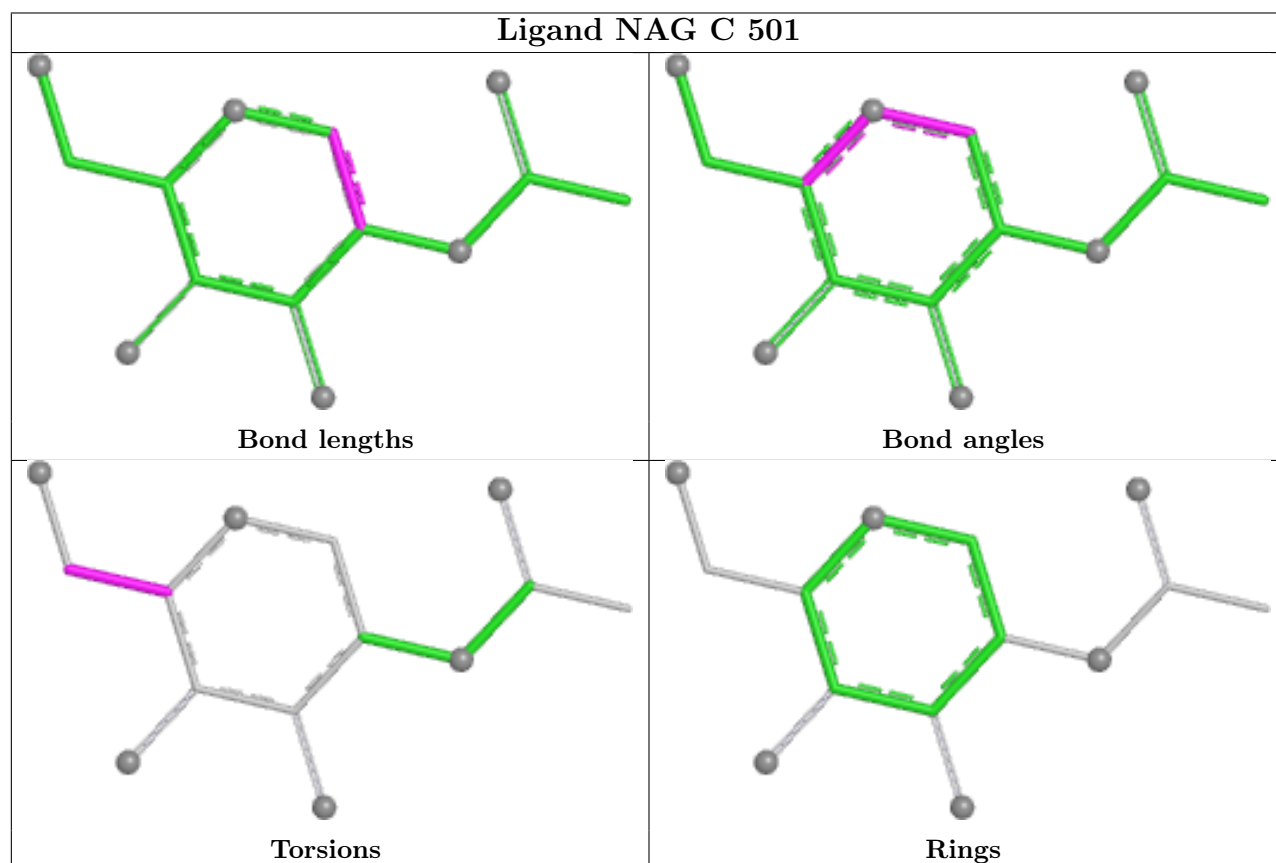
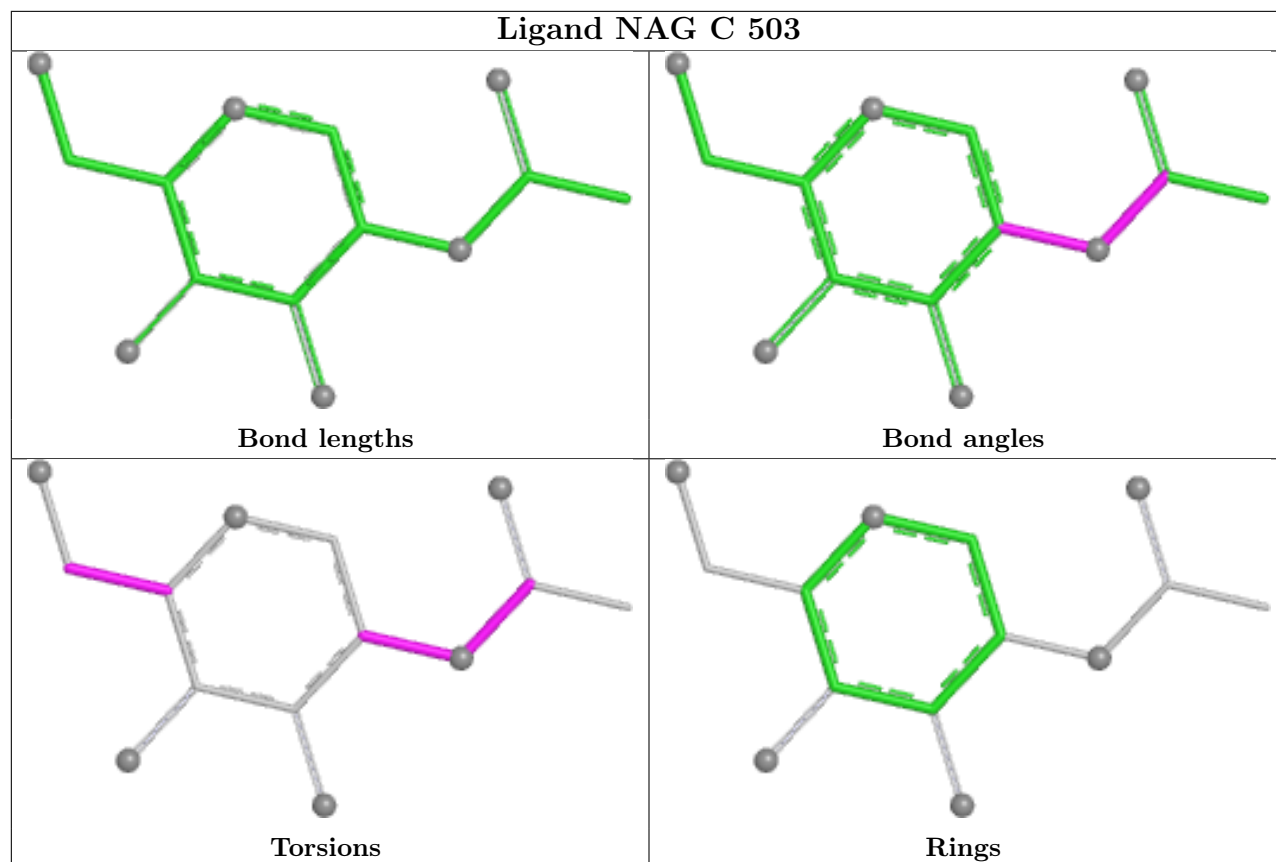


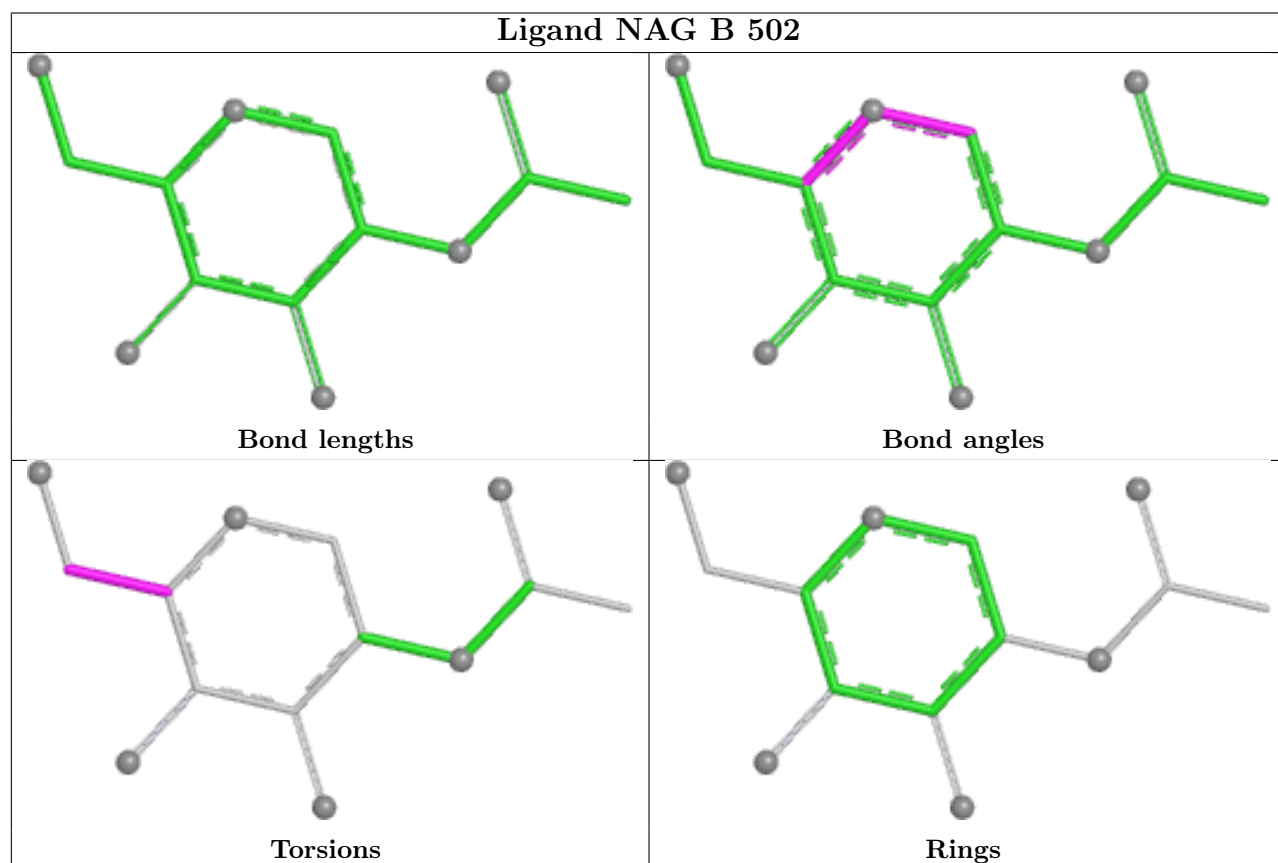
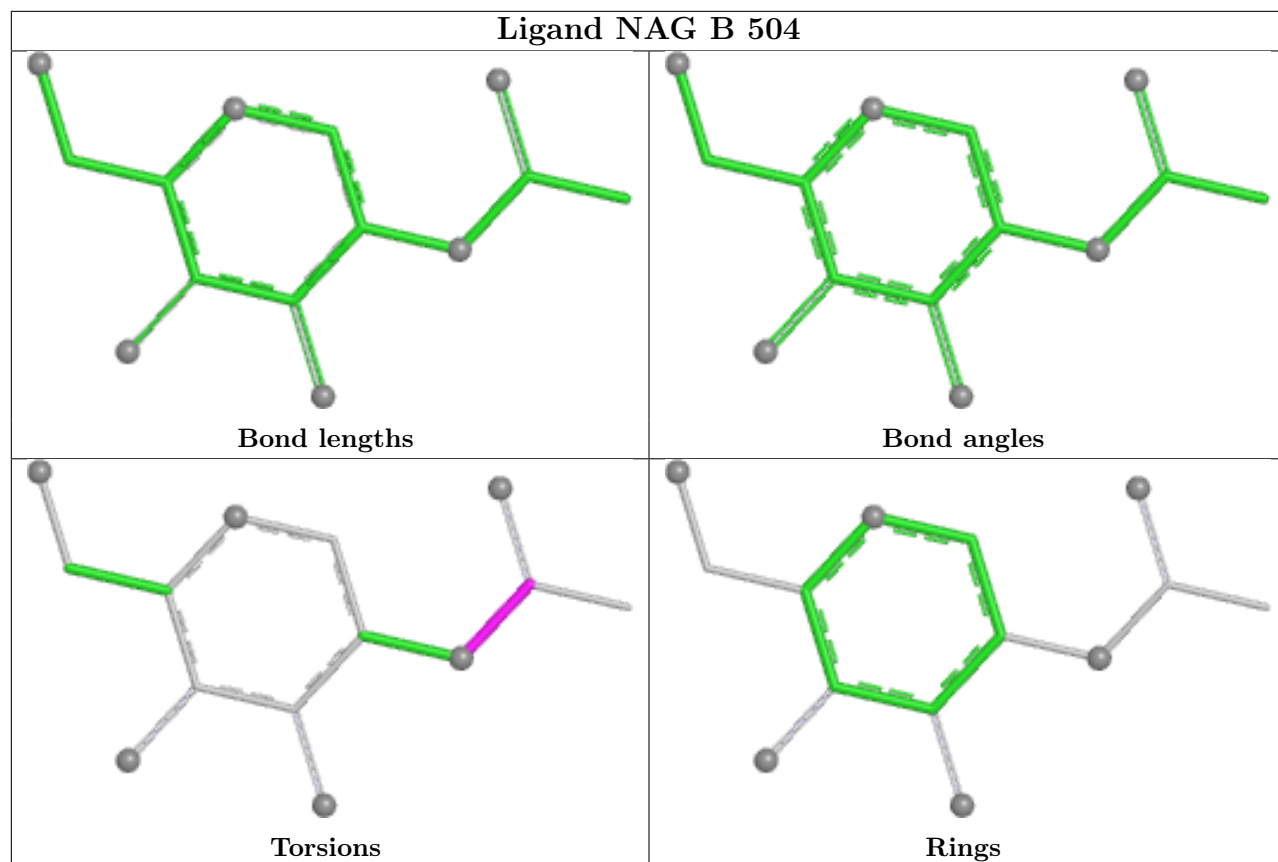


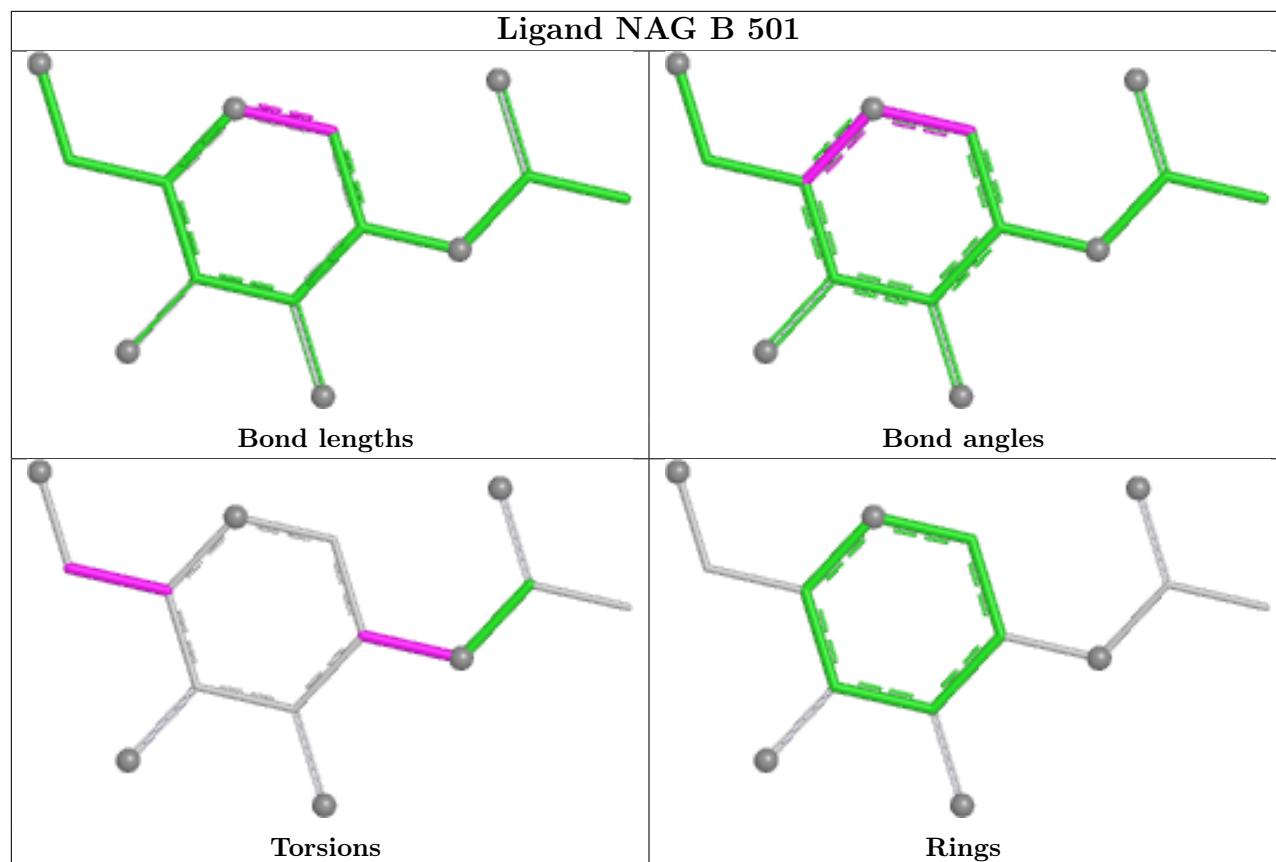












5.7 Other polymers [i](#)

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	393/424 (92%)	0.16	14 (3%) 46 46	25, 46, 83, 103	0
1	B	394/424 (92%)	0.11	7 (1%) 67 66	30, 44, 69, 99	0
1	C	394/424 (92%)	0.03	6 (1%) 72 71	27, 42, 72, 89	0
1	D	392/424 (92%)	0.24	12 (3%) 51 51	25, 47, 76, 91	0
All	All	1573/1696 (92%)	0.14	39 (2%) 58 58	25, 44, 75, 103	0

The worst 5 of 39 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	26	PRO	5.4
1	A	418	GLY	3.9
1	D	418	GLY	3.6
1	A	80	ASN	3.2
1	C	81	GLY	3.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MAN	J	7	11/12	0.20	0.27	125,139,165,167	0
3	MAN	F	7	11/12	0.39	0.22	77,92,107,113	0

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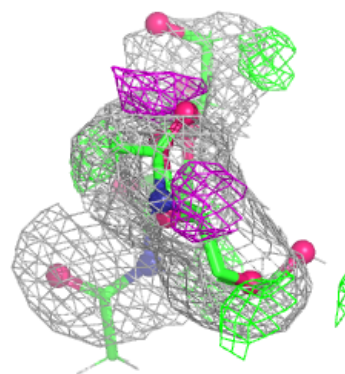
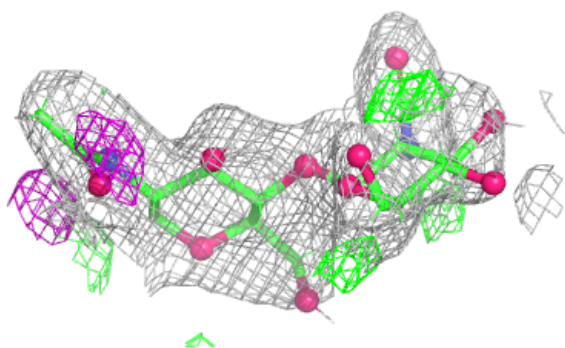
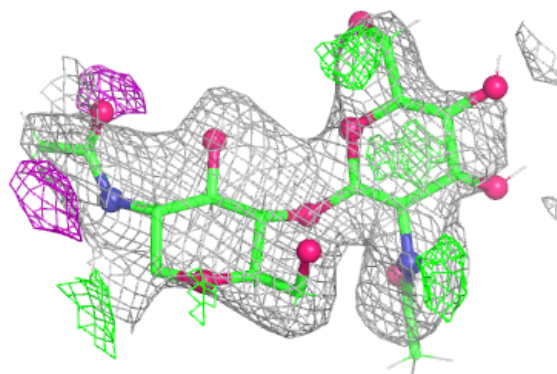
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MAN	L	7	11/12	0.54	0.17	78,93,110,116	0
2	NAG	G	2	14/15	0.55	0.21	73,89,105,112	0
4	MAN	L	6	11/12	0.61	0.16	75,91,109,112	0
3	MAN	H	7	11/12	0.61	0.17	63,79,97,105	0
4	NAG	L	5	14/15	0.62	0.21	62,78,95,95	0
2	NAG	E	2	14/15	0.71	0.16	66,81,96,103	0
2	NAG	I	2	14/15	0.74	0.15	59,81,98,109	0
3	MAN	H	6	11/12	0.76	0.15	60,72,86,88	0
2	NAG	E	1	14/15	0.77	0.16	45,68,82,94	0
4	MAN	L	4	11/12	0.79	0.17	59,71,84,85	0
3	MAN	F	6	11/12	0.79	0.13	60,74,85,94	0
3	BMA	J	3	11/12	0.80	0.16	51,72,124,124	0
2	NAG	K	2	14/15	0.80	0.14	63,77,93,96	0
2	NAG	G	1	14/15	0.81	0.15	44,58,81,89	0
2	NAG	I	1	14/15	0.83	0.14	49,60,72,81	0
2	NAG	K	1	14/15	0.84	0.13	48,62,75,85	0
3	NAG	F	2	14/15	0.84	0.13	46,57,69,86	0
3	MAN	J	5	11/12	0.85	0.15	52,63,76,91	0
3	NAG	J	2	14/15	0.85	0.13	42,54,68,76	0
3	MAN	H	5	11/12	0.86	0.13	50,63,76,80	0
4	BMA	L	3	11/12	0.86	0.14	50,61,92,92	0
3	MAN	J	6	11/12	0.86	0.13	52,63,73,85	0
3	MAN	F	4	11/12	0.89	0.10	47,56,68,73	0
3	MAN	F	5	11/12	0.89	0.10	55,66,81,84	0
3	BMA	F	3	11/12	0.89	0.11	50,64,88,88	0
3	MAN	J	4	11/12	0.89	0.11	45,57,70,73	0
4	NAG	L	2	14/15	0.90	0.10	40,51,64,66	0
3	BMA	H	3	11/12	0.90	0.11	45,52,65,68	0
3	MAN	H	4	11/12	0.91	0.09	47,56,71,82	0
3	NAG	H	2	14/15	0.92	0.09	38,44,51,57	0
3	NAG	F	1	14/15	0.92	0.09	38,46,56,64	0
3	NAG	H	1	14/15	0.94	0.08	23,36,42,50	0
3	NAG	J	1	14/15	0.94	0.08	36,47,57,72	0
4	NAG	L	1	14/15	0.94	0.08	31,38,47,50	0

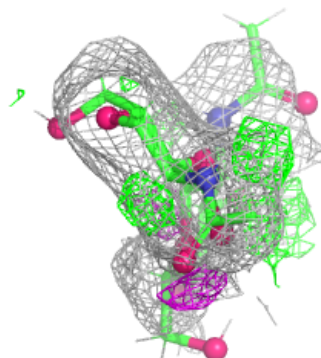
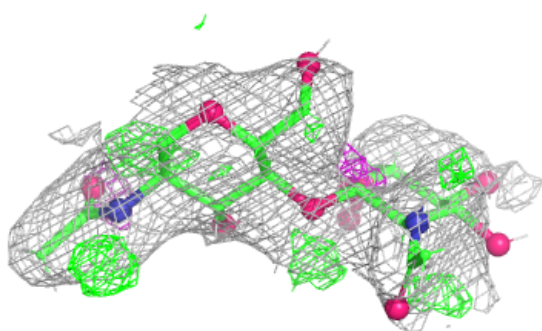
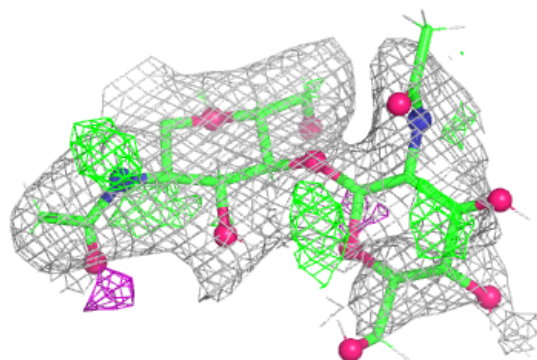
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

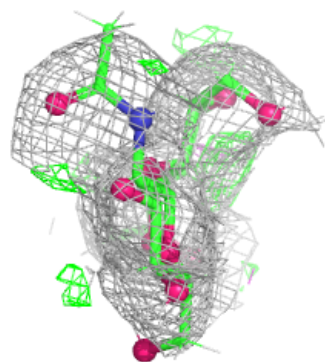
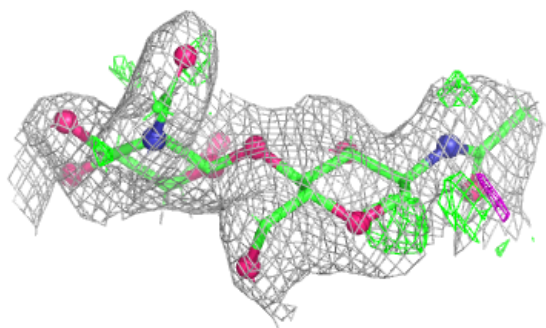
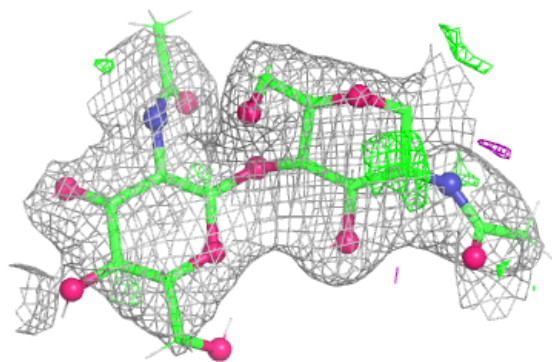
**Electron density around Chain G:**

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

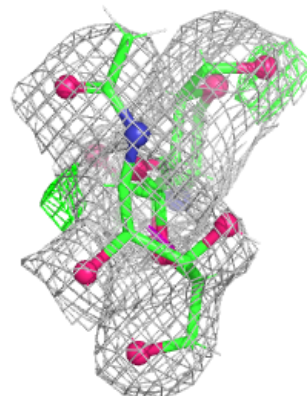
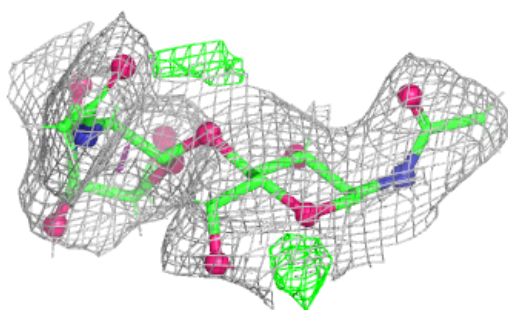
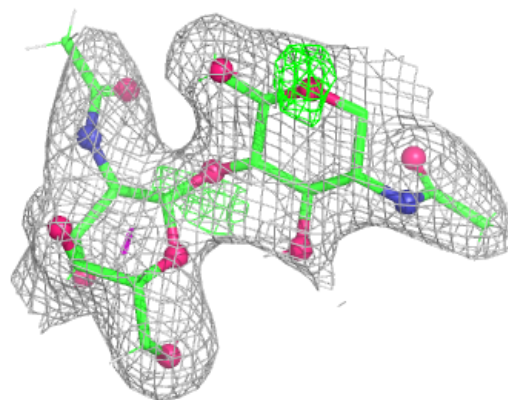


Electron density around Chain I:

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

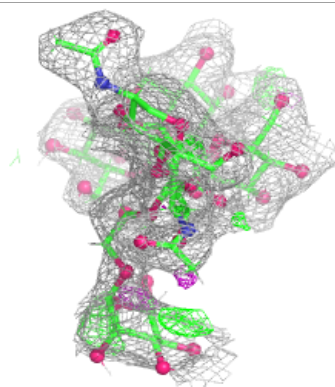
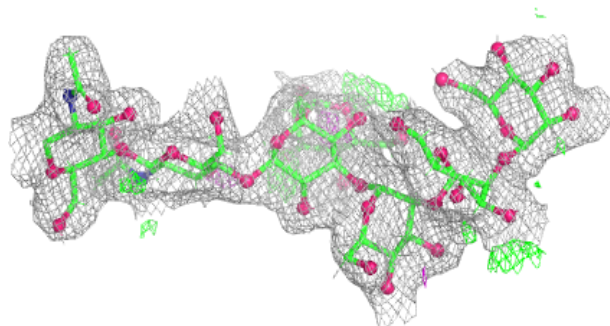
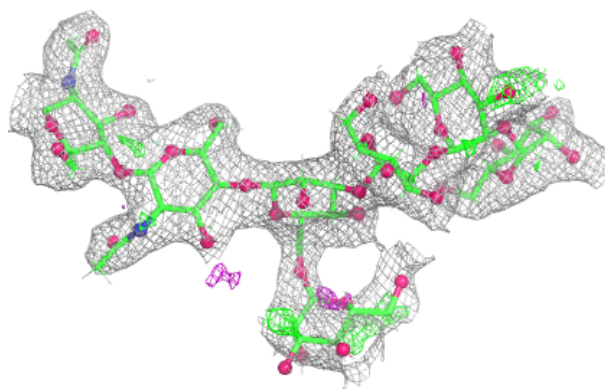
**Electron density around Chain K:**

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

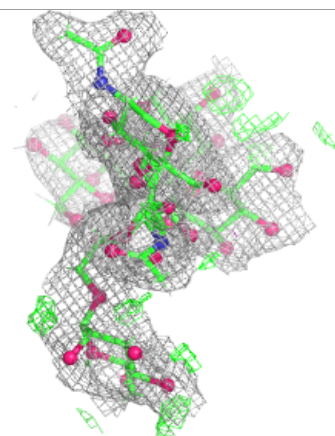
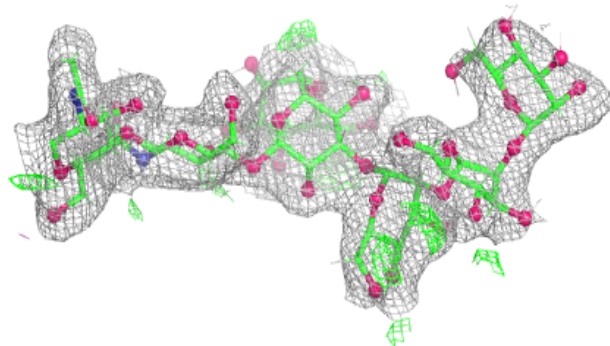
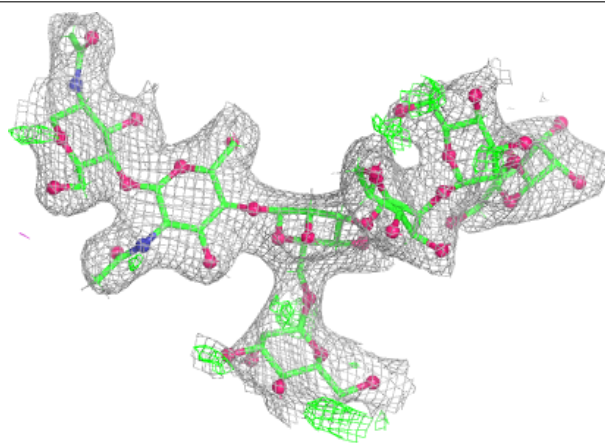


Electron density around Chain F:

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

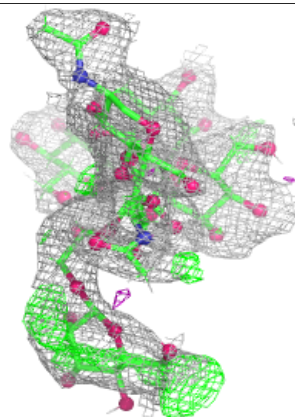
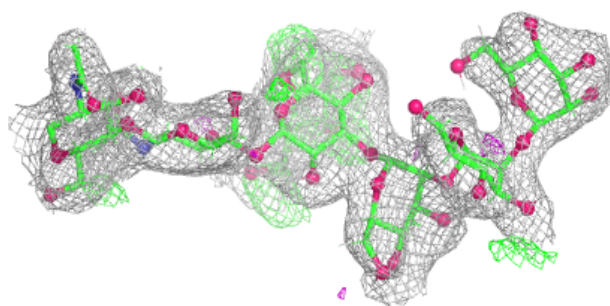
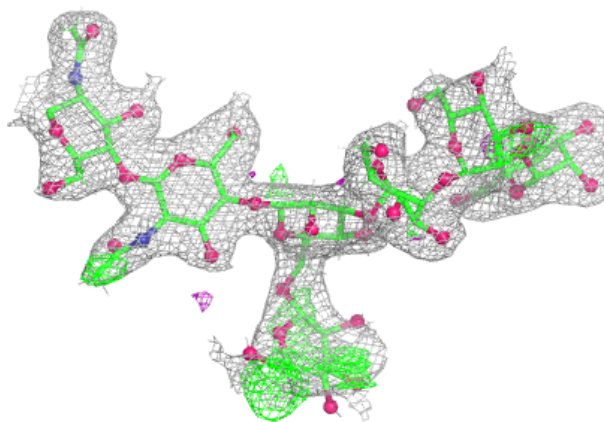
**Electron density around Chain H:**

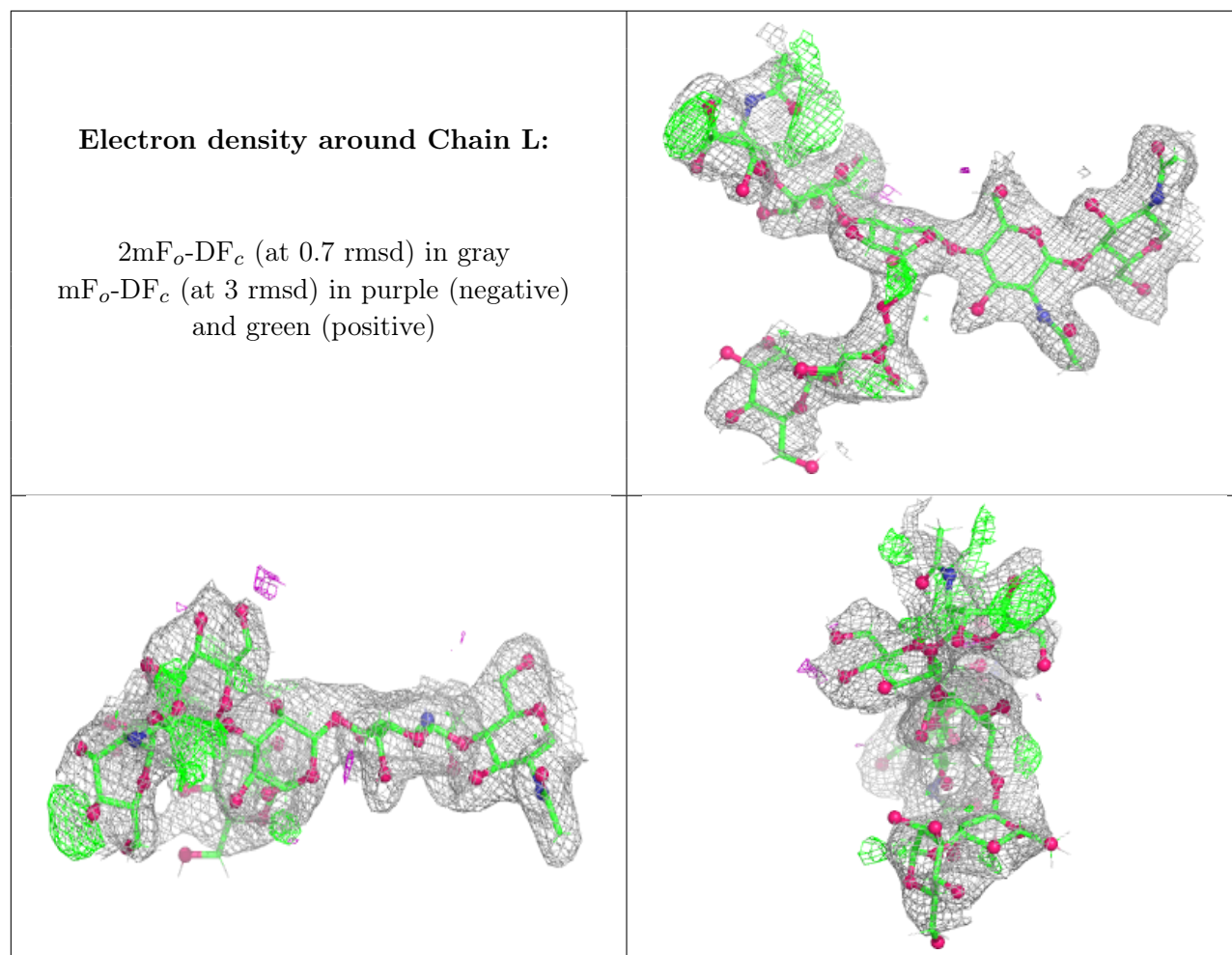
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain J:

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





6.4 Ligands [i](#)

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

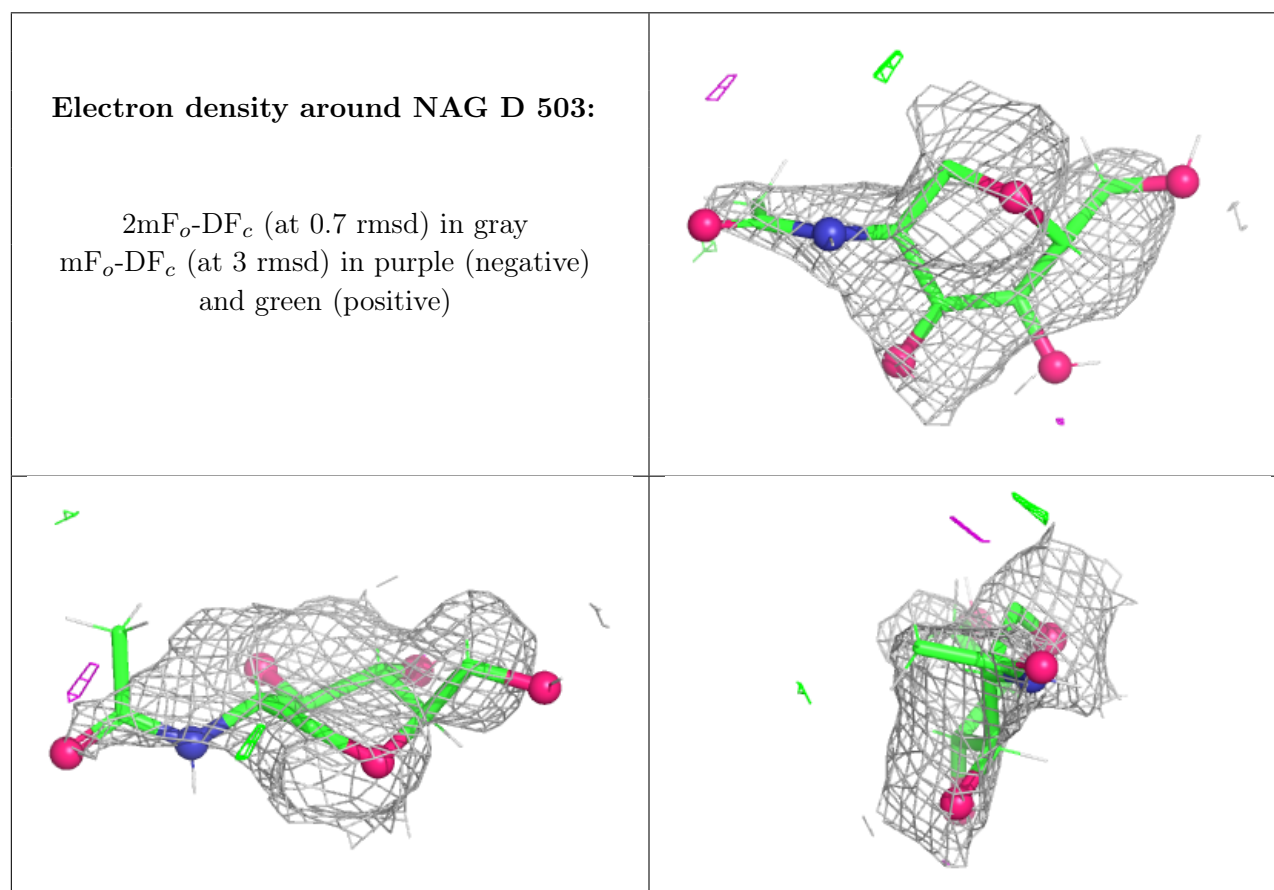
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	D	503	14/15	0.46	0.20	81,93,110,117	0
5	NAG	C	501	14/15	0.51	0.21	75,93,114,118	0
5	NAG	A	504	14/15	0.54	0.20	65,83,107,109	0
5	NAG	D	501	14/15	0.57	0.21	59,78,98,101	0
5	NAG	B	504	14/15	0.58	0.17	56,77,94,104	0
5	NAG	A	501	14/15	0.58	0.17	75,94,113,121	0
5	NAG	B	501	14/15	0.62	0.19	68,87,100,104	0
5	NAG	C	505	14/15	0.65	0.21	70,80,96,111	0
5	NAG	B	505	14/15	0.68	0.18	72,87,103,111	0
5	NAG	C	504	14/15	0.69	0.16	67,82,100,110	0

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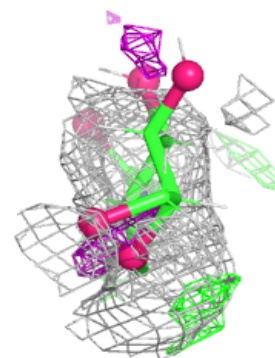
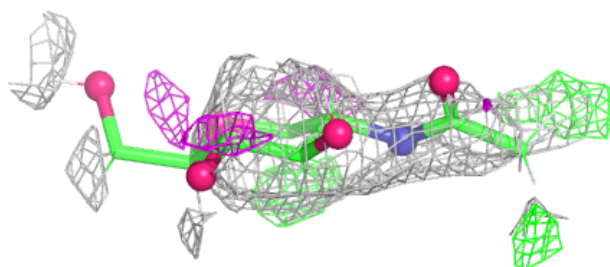
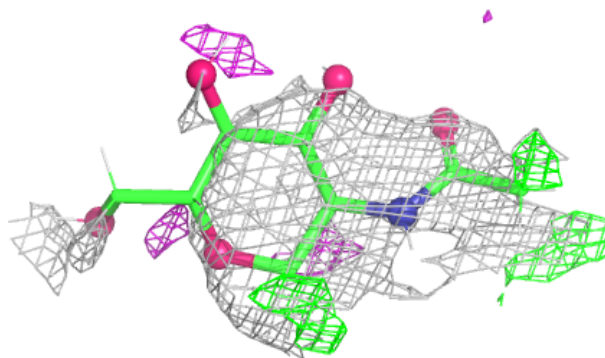
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	B	502	14/15	0.69	0.16	54,75,90,90	0
5	NAG	A	502	14/15	0.70	0.17	48,69,82,88	0
5	NAG	B	503	14/15	0.71	0.16	70,83,100,100	0
5	NAG	C	502	14/15	0.71	0.17	56,73,86,92	0
5	NAG	C	503	14/15	0.74	0.17	63,78,94,94	0
5	NAG	D	504	14/15	0.74	0.25	87,100,124,124	0
5	NAG	D	502	14/15	0.77	0.15	47,61,75,76	0
5	NAG	A	503	14/15	0.81	0.13	50,69,88,93	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.

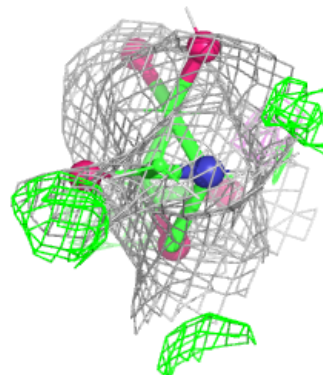
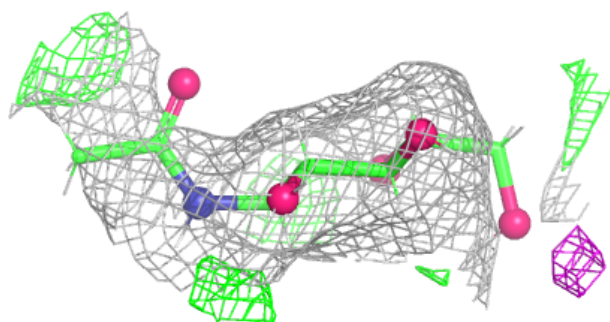
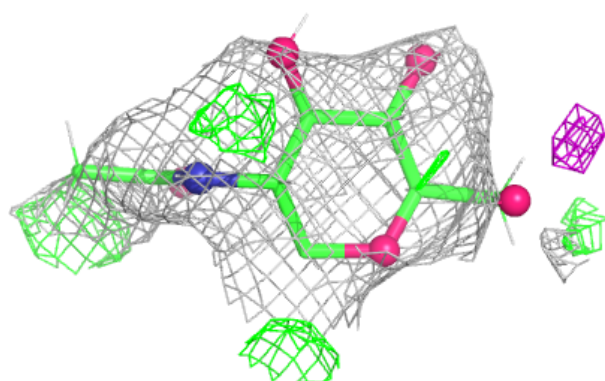


Electron density around NAG C 501:

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

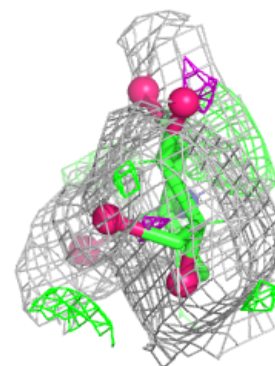
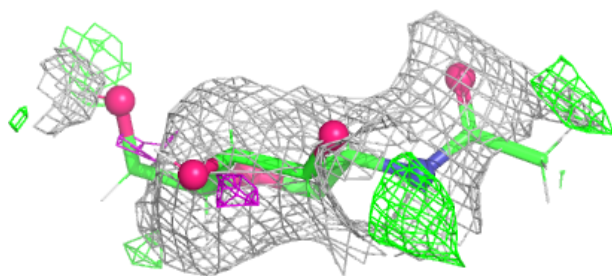
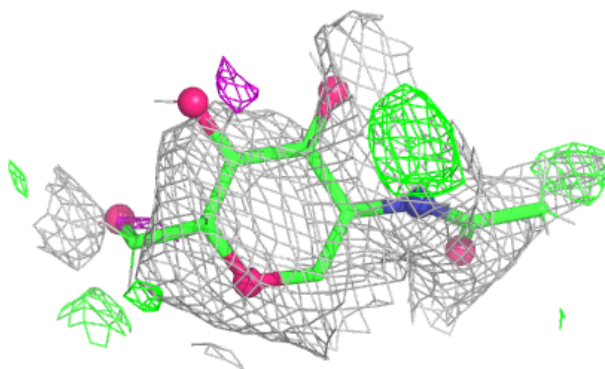
**Electron density around NAG A 504:**

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

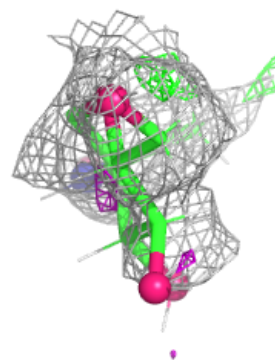
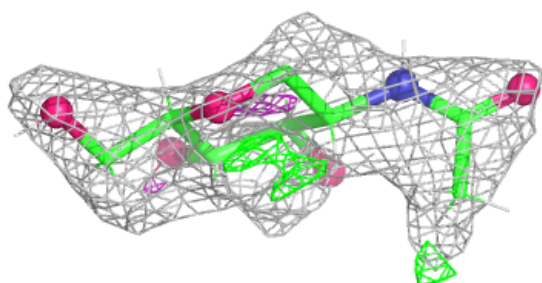
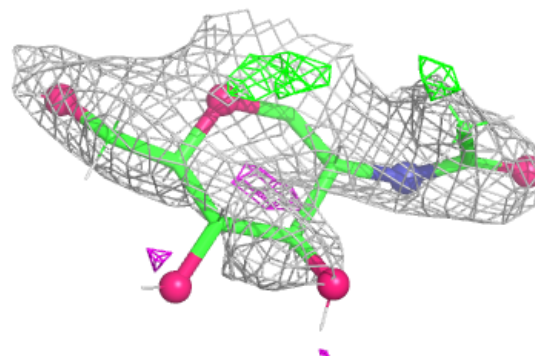


Electron density around NAG D 501:

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

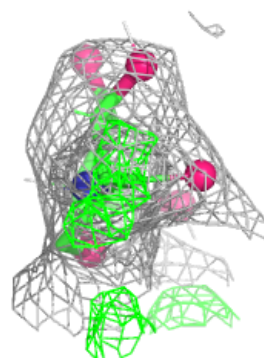
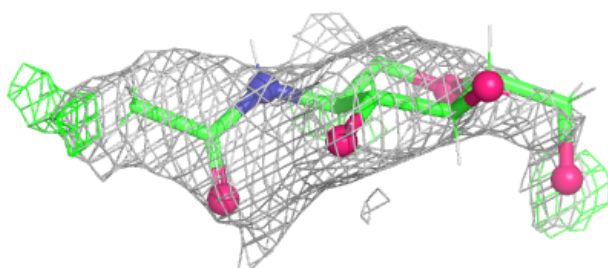
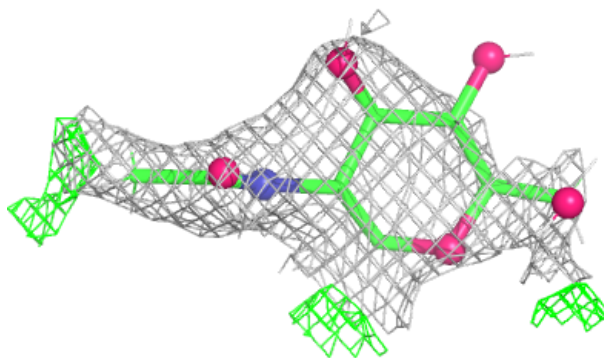
**Electron density around NAG B 504:**

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

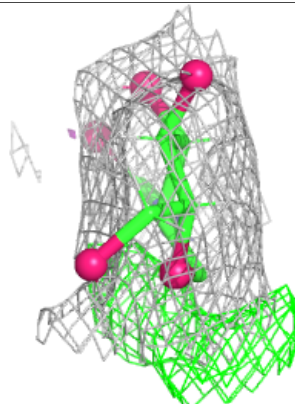
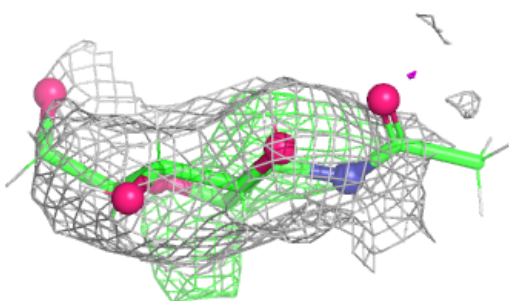
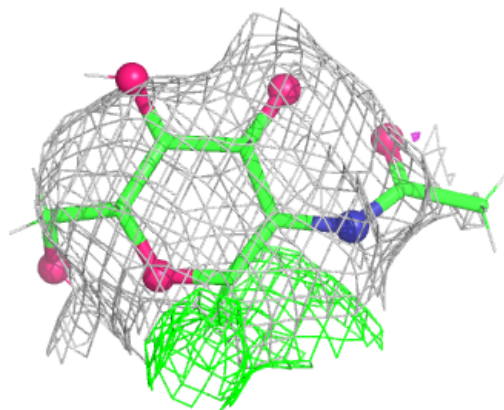


Electron density around NAG A 501:

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

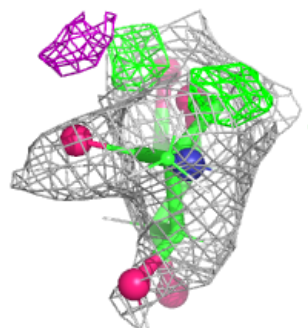
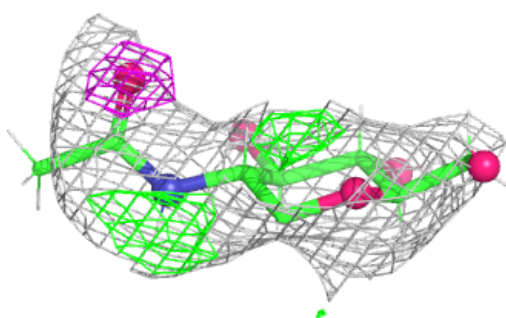
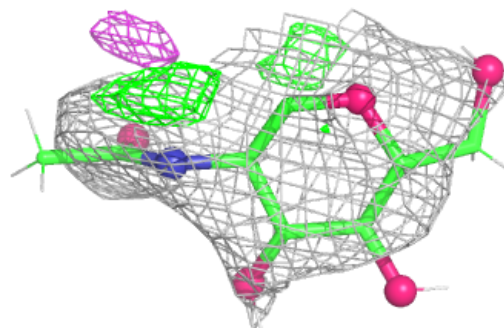
**Electron density around NAG B 501:**

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

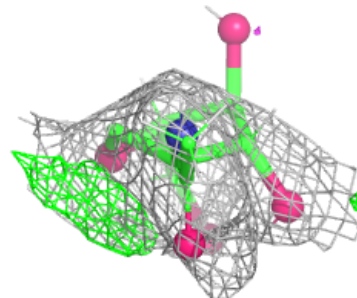
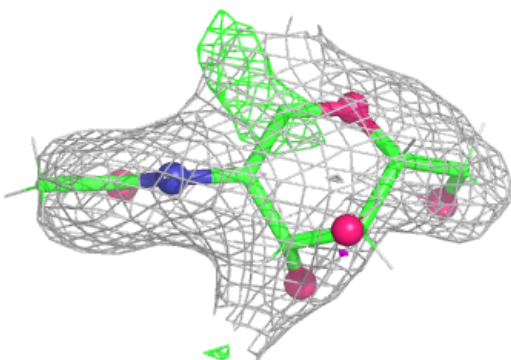
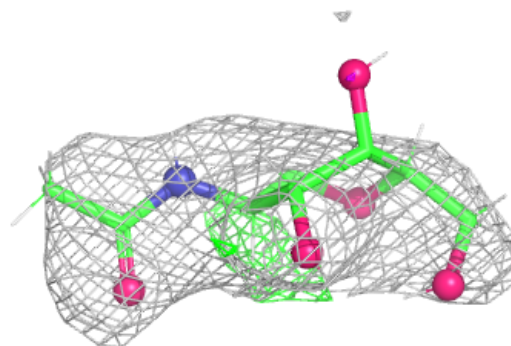


Electron density around NAG C 505:

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

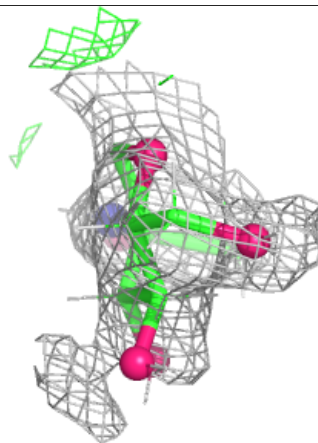
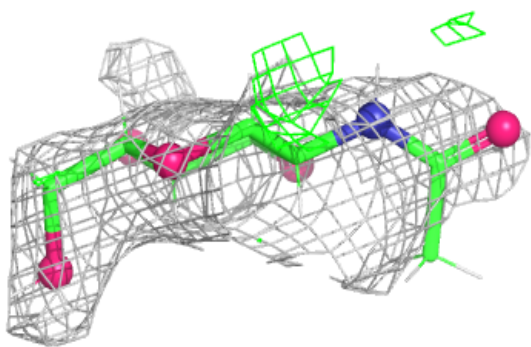
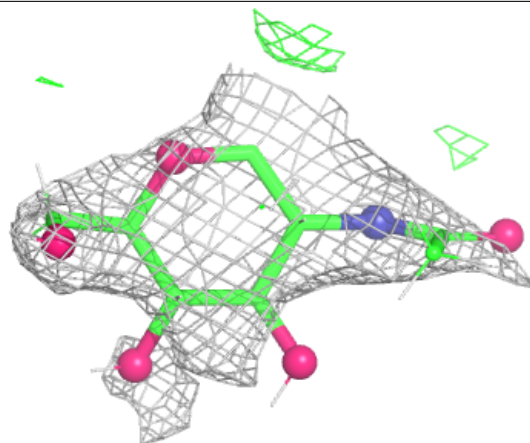
**Electron density around NAG B 505:**

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



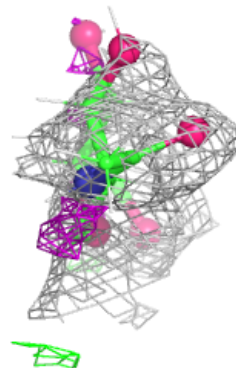
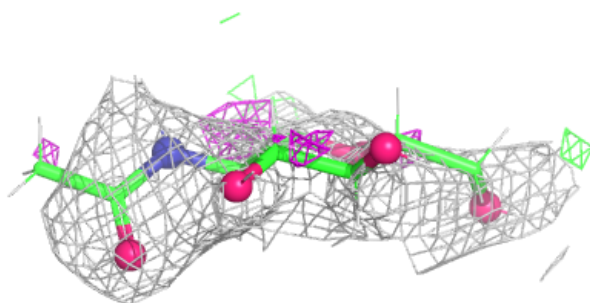
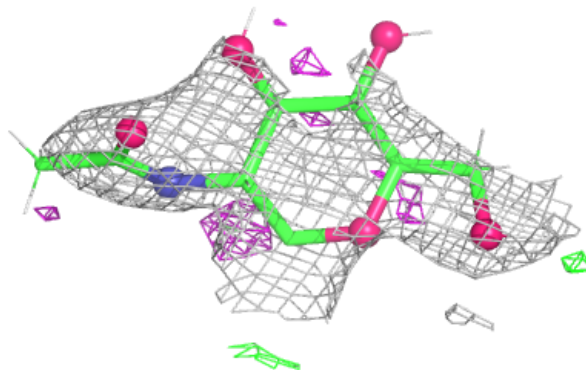
Electron density around NAG C 504:

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and green (positive)

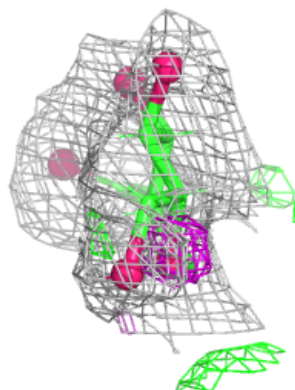
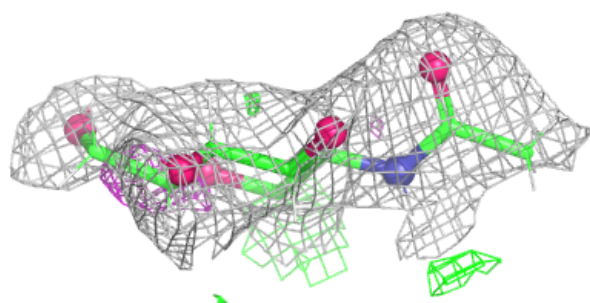
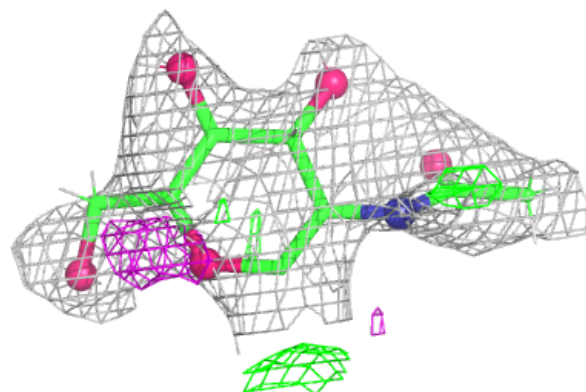


Electron density around NAG B 502:

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and green (positive)

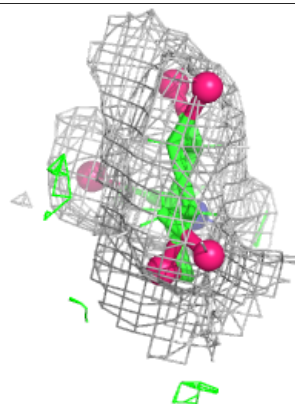
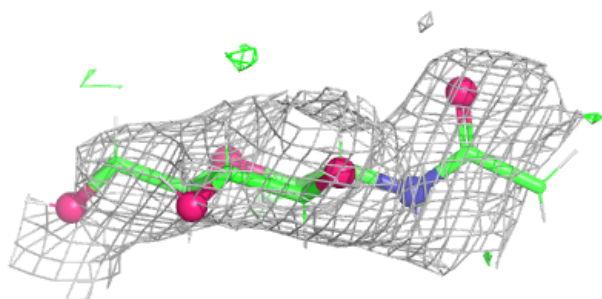
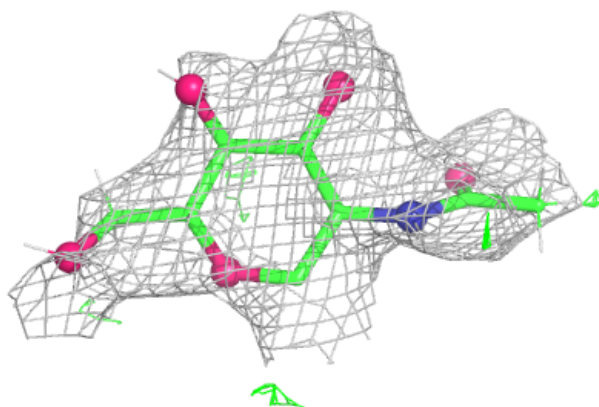
**Electron density around NAG A 502:**

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

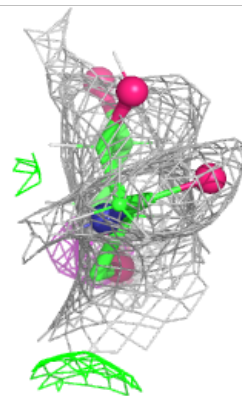
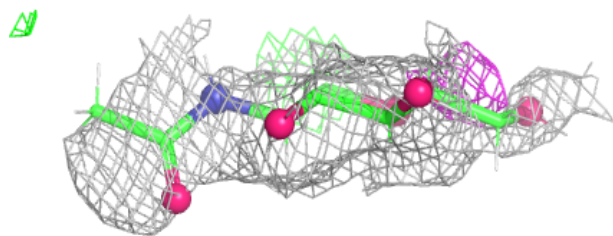
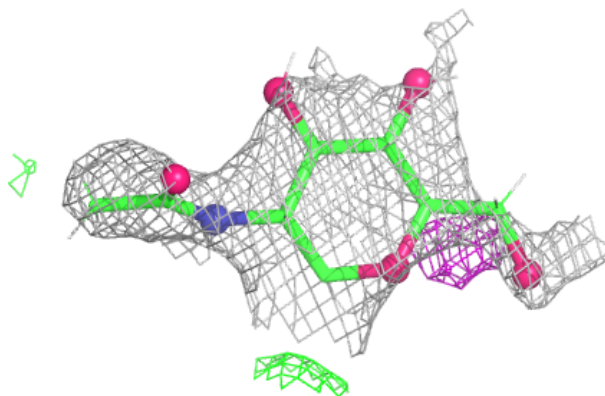


Electron density around NAG B 503:

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

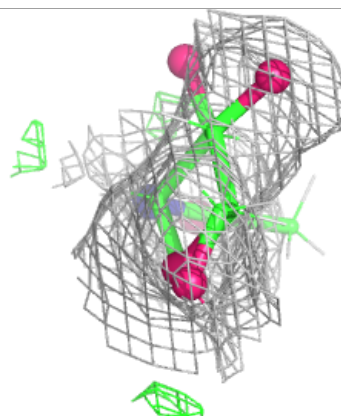
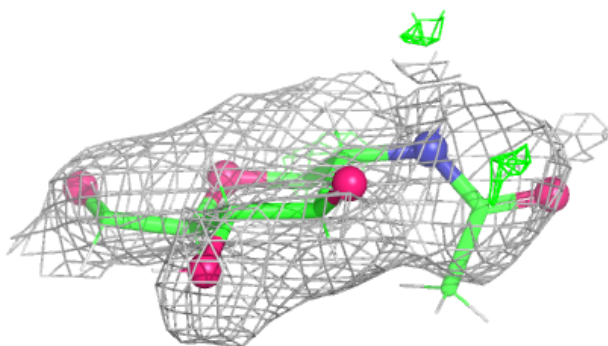
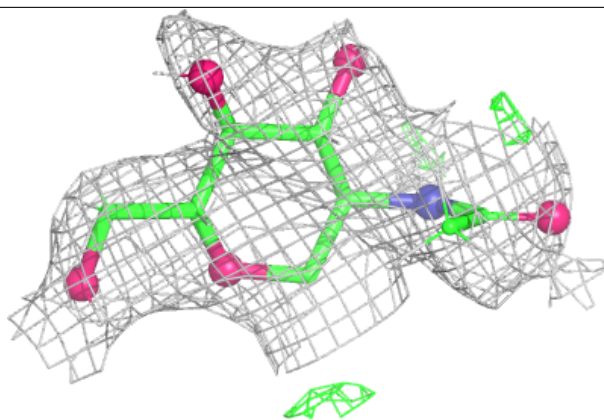
**Electron density around NAG C 502:**

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

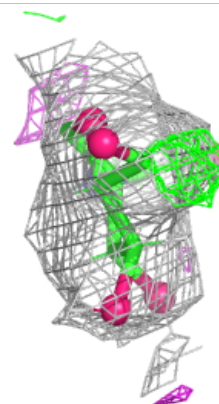
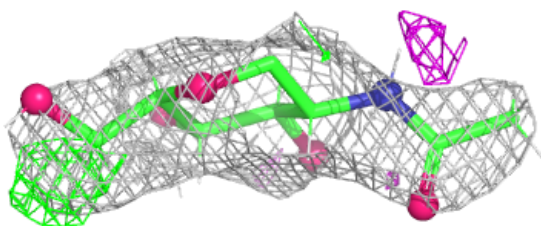
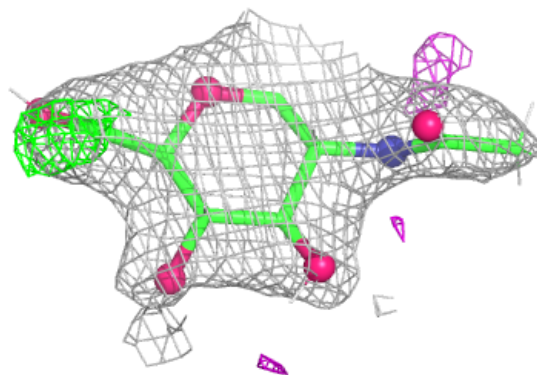


Electron density around NAG C 503:

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

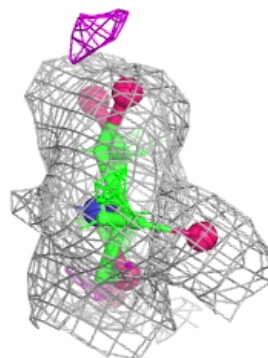
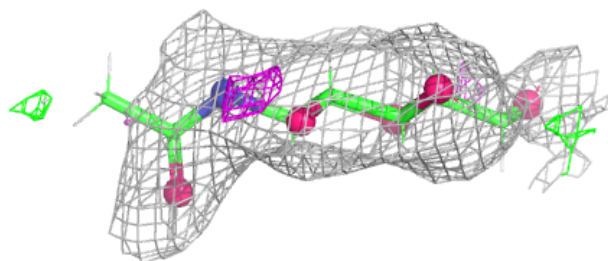
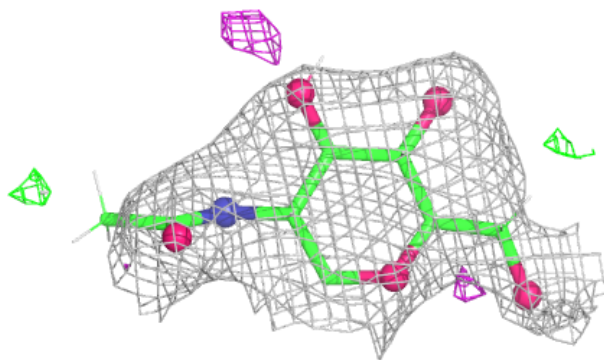
**Electron density around NAG D 504:**

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

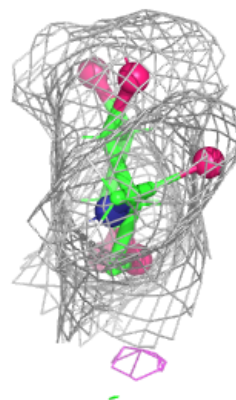
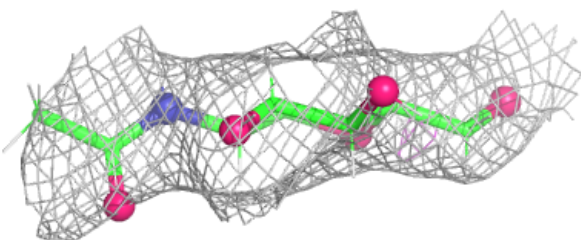
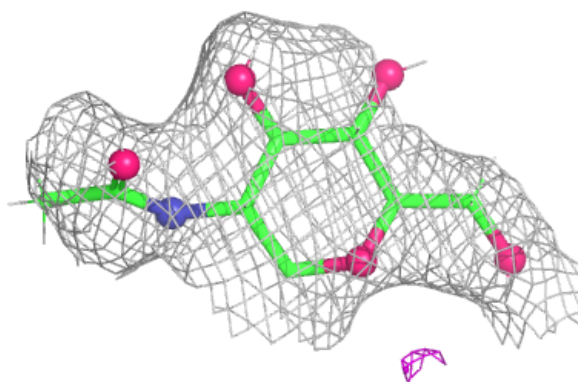


Electron density around NAG D 502:

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

**Electron density around NAG A 503:**

$2mF_o-DF_c$ (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.