



Full wwPDB EM Validation Report ⓘ

Feb 19, 2025 – 08:10 PM JST

PDB ID : 8ZY5
EMDB ID : EMD-60556
Title : Sarbecovirus RmYN02 Spike Trimer in a Locked Conformation
Authors : Wang, J.; Xiong, X.
Deposited on : 2024-06-16
Resolution : 3.80 Å (reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>

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:

EMDB validation analysis : **FAILED**
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.41.2

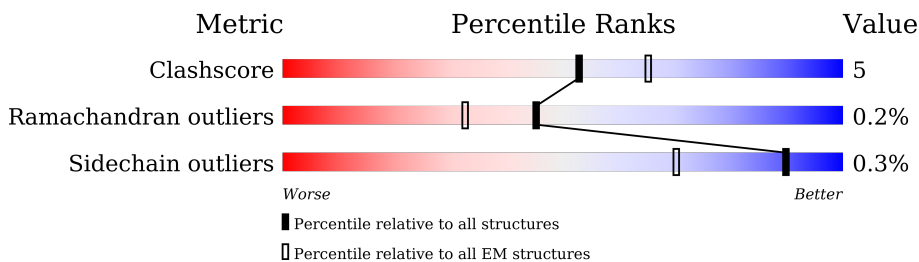
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.80 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)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	1241	
1	B	1241	
1	C	1241	
2	D	2	
2	E	2	
2	F	2	
2	G	2	
2	H	2	
2	I	2	

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Mol	Chain	Length	Quality of chain
2	J	2	 50% 50%
2	K	2	 50% 50%
2	L	2	 50% 50%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 24450 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Spike glycoprotein.

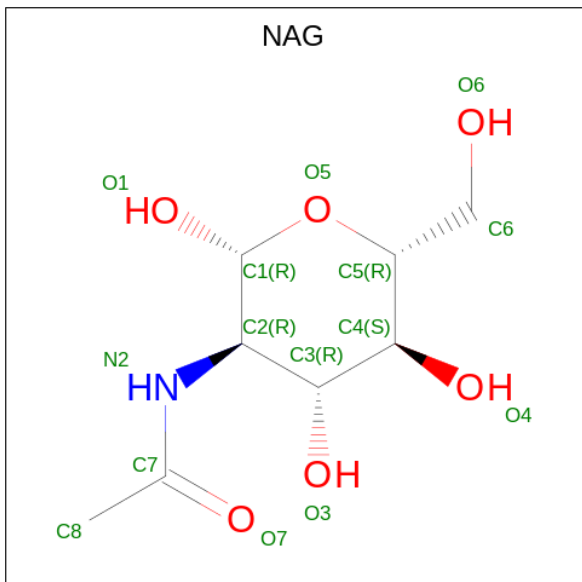
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1010	7856	5018	1306	1497	35	0	0
1	B	1010	7856	5018	1306	1497	35	0	0
1	C	1010	7856	5018	1306	1497	35	0	0

- 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				AltConf	Trace
			Total	C	N	O		
2	D	2	28	16	2	10	0	0
2	E	2	28	16	2	10	0	0
2	F	2	28	16	2	10	0	0
2	G	2	28	16	2	10	0	0
2	H	2	28	16	2	10	0	0
2	I	2	28	16	2	10	0	0
2	J	2	28	16	2	10	0	0
2	K	2	28	16	2	10	0	0
2	L	2	28	16	2	10	0	0

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).



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

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

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Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0
3	C	1	Total 14	8	1	5	0

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

Chain D:  100%

MAG1
MAG2

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

Chain E:  50% 50%

MAG1
MAG2

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

Chain F:  50% 50%

MAG1
MAG2

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

Chain G:  100%

MAG1
MAG2

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

Chain H:  50% 50%

MAG1
MAG2

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

Chain I:  50% 50%

MAG1
MAG2

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

Chain J:  50% 50%

A legend for the validation report showing two colored boxes: a green box labeled 'MAG1' and a yellow box labeled 'MAG2'.

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

Chain K: A horizontal bar representing the validation for Chain K. The bar is split into two equal halves, each labeled '50%'. The left half is green (MAG1) and the right half is yellow (MAG2).

A legend for the validation report showing two colored boxes: a green box labeled 'MAG1' and a yellow box labeled 'MAG2'.

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

Chain L: A horizontal bar representing the validation for Chain L. The bar is split into two equal halves, each labeled '50%'. The left half is green (MAG1) and the right half is yellow (MAG2).

A legend for the validation report showing two colored boxes: a green box labeled 'MAG1' and a yellow box labeled 'MAG2'.

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	27757	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor

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: NAG

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	0.29	0/8027	0.58	0/10926
1	B	0.29	0/8027	0.58	0/10926
1	C	0.30	0/8027	0.58	0/10926
All	All	0.30	0/24081	0.58	0/32778

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	7856	0	7702	83	0
1	B	7856	0	7702	92	0
1	C	7856	0	7702	85	0
2	D	28	0	25	0	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
2	I	28	0	25	0	0
2	J	28	0	25	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	K	28	0	25	0	0
2	L	28	0	25	0	0
3	A	210	0	195	1	0
3	B	210	0	195	1	0
3	C	210	0	195	1	0
All	All	24450	0	23916	248	0

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

All (248) 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:B:572:TYR:O	1:B:608:GLY:HA3	1.77	0.84
1:C:572:TYR:O	1:C:608:GLY:HA3	1.77	0.84
1:A:572:TYR:O	1:A:608:GLY:HA3	1.80	0.82
1:B:1047:GLY:HA3	1:B:1059:THR:O	1.86	0.76
1:A:1047:GLY:HA3	1:A:1059:THR:O	1.86	0.74
1:B:707:LEU:HD13	1:B:951:ILE:HD11	1.76	0.68
1:C:125:GLN:HG3	3:C:1312:NAG:HN2	1.60	0.67
1:A:125:GLN:HG3	3:A:1312:NAG:HN2	1.61	0.66
1:B:707:LEU:HD12	1:B:710:TYR:HB2	1.78	0.65
1:C:1047:GLY:HA3	1:C:1059:THR:O	2.00	0.62
1:A:696:ILE:HG21	1:A:951:ILE:HG13	1.82	0.61
1:B:81:ILE:HD11	1:B:246:VAL:HG13	1.82	0.61
1:B:125:GLN:HG3	3:B:1312:NAG:HN2	1.66	0.61
1:B:696:ILE:HG21	1:B:951:ILE:HG13	1.83	0.60
1:C:696:ILE:HG21	1:C:951:ILE:HG13	1.84	0.60
1:A:422:ILE:HD12	1:A:459:PRO:HD3	1.83	0.60
1:A:807:GLN:HG3	1:A:812:LEU:O	2.02	0.60
1:B:815:LEU:HD12	1:B:816:PRO:HD2	1.82	0.59
1:B:485:CYS:SG	1:B:486:GLY:N	2.77	0.58
1:C:389:ILE:HD13	1:C:470:VAL:HG11	1.86	0.58
1:A:758:GLN:NE2	1:A:889:GLN:OE1	2.37	0.57
1:C:485:CYS:SG	1:C:486:GLY:N	2.77	0.57
1:C:397:ILE:O	1:C:398:ALA:HB3	2.05	0.57
1:A:769:ARG:NH1	1:A:774:ASP:OD1	2.38	0.57
1:A:485:CYS:SG	1:A:486:GLY:N	2.77	0.57
1:A:808:LYS:O	1:A:809:PHE:HB2	2.05	0.57
1:A:278:LEU:HD21	1:A:287:VAL:HG11	1.88	0.56
1:B:1000:GLY:HA3	1:B:1020:THR:HB	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:SER:HB3	1:C:245:ALA:HB3	1.88	0.56
1:C:769:ARG:NH1	1:C:774:ASP:OD1	2.39	0.56
1:C:278:LEU:HD21	1:C:287:VAL:HG11	1.87	0.55
1:A:748:ILE:HG22	1:A:750:ASP:H	1.71	0.55
1:C:172:LYS:HD3	1:C:197:ILE:HG13	1.89	0.55
1:B:204:SER:OG	1:B:205:LEU:N	2.40	0.55
1:C:38:ASP:OD2	1:C:42:ARG:NH1	2.40	0.55
1:C:1000:GLY:HA3	1:C:1020:THR:HB	1.88	0.55
1:B:374:VAL:HG22	1:B:475:PHE:HB3	1.89	0.55
1:A:389:ILE:HD13	1:A:470:VAL:HG11	1.89	0.55
1:C:779:LYS:NZ	1:C:895:THR:O	2.40	0.55
1:B:769:ARG:NH1	1:B:774:ASP:OD1	2.40	0.54
1:A:431:ARG:NH1	1:A:433:VAL:O	2.40	0.54
1:B:38:ASP:OD2	1:B:42:ARG:NH1	2.40	0.54
1:B:779:LYS:NZ	1:B:895:THR:O	2.39	0.54
1:B:748:ILE:HG22	1:B:750:ASP:H	1.73	0.54
1:A:1000:GLY:HA3	1:A:1020:THR:HB	1.89	0.54
1:B:389:ILE:HD13	1:B:470:VAL:HG11	1.89	0.54
1:B:431:ARG:NH1	1:B:433:VAL:O	2.41	0.54
1:C:748:ILE:HG22	1:C:750:ASP:H	1.72	0.54
1:A:38:ASP:OD2	1:A:42:ARG:NH1	2.40	0.54
1:C:939:ASP:OD1	1:C:939:ASP:N	2.41	0.54
1:C:60:LEU:HD23	1:C:245:ALA:HB1	1.90	0.53
1:C:125:GLN:HB3	1:C:147:SER:HB3	1.90	0.53
1:C:431:ARG:NH1	1:C:433:VAL:O	2.41	0.53
1:A:747:VAL:HG13	1:A:748:ILE:HG12	1.90	0.53
1:B:1045:ARG:NH1	1:B:1072:ASP:O	2.42	0.53
1:A:27:VAL:HG13	1:A:62:TRP:HE1	1.74	0.53
1:A:939:ASP:OD1	1:A:939:ASP:N	2.40	0.53
1:B:60:LEU:HD23	1:B:245:ALA:HB1	1.91	0.53
1:A:497:ARG:NH1	1:A:587:ASP:OD2	2.40	0.53
1:B:84:SER:HB3	1:B:245:ALA:HB3	1.89	0.53
1:B:537:ARG:HD3	1:B:542:LEU:HD23	1.91	0.53
1:A:707:LEU:HD13	1:A:951:ILE:HD11	1.90	0.53
1:A:837:THR:HG23	1:C:661:TYR:HB2	1.91	0.53
1:B:298:ARG:NH2	1:B:552:PHE:O	2.41	0.52
1:C:497:ARG:NH1	1:C:587:ASP:OD2	2.42	0.52
1:C:747:VAL:HG13	1:C:748:ILE:HG12	1.91	0.52
1:B:296:ASN:HD21	1:B:551:SER:HB3	1.73	0.52
1:A:374:VAL:HG22	1:A:475:PHE:HB3	1.90	0.52
1:B:747:VAL:HG13	1:B:748:ILE:HG12	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:851:PRO:HA	1:C:661:TYR:HE1	1.74	0.52
1:C:298:ARG:NH2	1:C:552:PHE:O	2.43	0.52
1:A:296:ASN:HD21	1:A:551:SER:HB3	1.74	0.52
1:A:537:ARG:HD3	1:A:542:LEU:HD23	1.92	0.52
1:B:687:LYS:HE2	1:B:815:LEU:HD23	1.92	0.51
1:C:49:PHE:O	1:C:253:THR:HA	2.10	0.51
1:C:537:ARG:HD3	1:C:542:LEU:HD23	1.91	0.51
1:B:369:LEU:HD11	1:C:937:ARG:HA	1.92	0.51
1:C:807:GLN:O	1:C:810:ASN:HB2	2.10	0.51
1:B:211:PRO:HB3	1:B:264:ILE:HB	1.93	0.51
1:B:497:ARG:NH1	1:B:587:ASP:OD2	2.43	0.51
1:B:661:TYR:HE1	1:C:851:PRO:HA	1.75	0.51
1:C:81:ILE:HD11	1:C:246:VAL:HG13	1.91	0.51
1:C:710:TYR:O	1:C:712:SER:N	2.44	0.51
1:B:661:TYR:HB2	1:C:837:THR:HG23	1.93	0.51
1:C:444:SER:HA	1:C:452:LEU:HD23	1.93	0.51
1:C:517:LYS:HB2	1:C:544:ILE:HG21	1.93	0.51
1:B:517:LYS:HB2	1:B:544:ILE:HG21	1.91	0.51
1:C:716:GLN:OE1	1:C:719:ARG:NH1	2.44	0.51
1:A:661:TYR:HB2	1:B:837:THR:HG23	1.93	0.51
1:C:1045:ARG:NH1	1:C:1072:ASP:O	2.43	0.51
1:A:517:LYS:HB2	1:A:544:ILE:HG21	1.93	0.51
1:B:27:VAL:HG13	1:B:62:TRP:HE1	1.76	0.51
1:A:298:ARG:NH2	1:A:552:PHE:O	2.43	0.51
1:A:586:LEU:O	1:A:594:ARG:NH1	2.44	0.51
1:B:516:ASP:OD1	1:B:516:ASP:N	2.43	0.51
1:C:551:SER:OG	1:C:552:PHE:N	2.44	0.50
1:C:211:PRO:HB3	1:C:264:ILE:HB	1.93	0.50
1:A:41:TYR:HB3	1:C:479:ASN:HD22	1.76	0.50
1:C:296:ASN:HD21	1:C:551:SER:HB3	1.75	0.50
1:A:130:PRO:HB2	1:A:230:VAL:HG13	1.93	0.50
1:A:937:ARG:HA	1:C:369:LEU:HD11	1.93	0.50
1:B:444:SER:HA	1:B:452:LEU:HD23	1.94	0.50
1:C:895:THR:HG22	1:C:897:SER:H	1.76	0.50
2:J:1:NAG:H61	2:J:2:NAG:H61	1.93	0.50
1:B:982:LYS:NZ	1:B:996:PHE:O	2.45	0.50
1:B:586:LEU:O	1:B:594:ARG:NH1	2.44	0.50
1:A:125:GLN:HB3	1:A:147:SER:HB3	1.94	0.50
1:B:526:GLY:N	1:B:535:SER:O	2.45	0.50
1:A:982:LYS:NZ	1:A:996:PHE:O	2.45	0.49
1:B:278:LEU:HD21	1:B:287:VAL:HG11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:PHE:O	1:A:253:THR:HA	2.12	0.49
1:A:1045:ARG:NH1	1:A:1072:ASP:O	2.45	0.49
1:B:680:VAL:HG22	1:B:1015:VAL:HG22	1.94	0.49
1:C:306:VAL:HG22	1:C:502:ASN:HB3	1.94	0.49
1:A:551:SER:OG	1:A:552:PHE:N	2.44	0.49
1:A:526:GLY:N	1:A:535:SER:O	2.44	0.49
1:A:60:LEU:HD23	1:A:245:ALA:HB1	1.95	0.49
1:A:444:SER:HA	1:A:452:LEU:HD23	1.94	0.49
1:B:125:GLN:HB3	1:B:147:SER:HB3	1.95	0.49
1:B:551:SER:OG	1:B:552:PHE:N	2.45	0.49
1:C:347:LEU:O	1:C:351:THR:OG1	2.31	0.49
1:A:211:PRO:HB3	1:A:264:ILE:HB	1.95	0.49
1:A:895:THR:HG22	1:A:897:SER:H	1.78	0.49
1:A:347:LEU:O	1:A:351:THR:OG1	2.31	0.48
1:C:815:LEU:HD12	1:C:816:PRO:HD2	1.95	0.48
1:C:130:PRO:HB2	1:C:230:VAL:HG13	1.95	0.48
1:A:680:VAL:HG22	1:A:1015:VAL:HG22	1.96	0.48
1:B:172:LYS:HD3	1:B:197:ILE:HG13	1.94	0.48
1:A:661:TYR:HE1	1:B:851:PRO:HA	1.78	0.48
1:B:895:THR:HG22	1:B:897:SER:H	1.77	0.48
1:A:708:LEU:HD11	1:C:50:THR:HG23	1.95	0.48
1:B:604:GLN:NE2	1:B:608:GLY:O	2.46	0.48
1:B:347:LEU:O	1:B:351:THR:OG1	2.31	0.47
1:C:526:GLY:N	1:C:535:SER:O	2.44	0.47
1:A:370:CYS:HB3	1:A:485:CYS:HB2	1.79	0.47
1:C:604:GLN:NE2	1:C:608:GLY:O	2.47	0.47
1:B:939:ASP:N	1:B:939:ASP:OD1	2.40	0.47
1:C:123:TYR:O	1:C:148:SER:HA	2.15	0.47
1:A:123:TYR:O	1:A:148:SER:HA	2.14	0.47
1:B:130:PRO:HB2	1:B:230:VAL:HG13	1.96	0.47
1:C:577:CYS:HB2	1:C:609:CYS:HB3	1.72	0.47
1:B:234:ARG:HA	1:B:234:ARG:HD3	1.77	0.47
1:B:254:ILE:HG22	1:B:269:ASP:HA	1.95	0.47
1:B:577:CYS:HB2	1:B:609:CYS:HB3	1.73	0.47
1:A:172:LYS:HD3	1:A:197:ILE:HG13	1.97	0.47
1:B:1036:CYS:HB3	1:B:1088:ASN:HB3	1.97	0.47
1:B:538:ASP:OD1	1:B:538:ASP:N	2.49	0.46
1:B:815:LEU:CD1	1:B:816:PRO:HD2	2.45	0.46
1:C:680:VAL:HG22	1:C:1015:VAL:HG22	1.96	0.46
1:A:604:GLN:NE2	1:A:608:GLY:O	2.47	0.46
1:B:38:ASP:OD1	1:B:38:ASP:N	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:ILE:HD12	1:B:459:PRO:HD3	1.96	0.46
1:C:370:CYS:HB3	1:C:485:CYS:HB2	1.79	0.46
1:C:982:LYS:NZ	1:C:996:PHE:O	2.49	0.46
1:A:538:ASP:N	1:A:538:ASP:OD1	2.47	0.46
1:C:422:ILE:HD12	1:C:459:PRO:HD3	1.97	0.46
1:A:352:SER:HG	1:C:384:SER:HG	1.57	0.46
1:A:866:THR:H	1:A:1060:GLN:HE22	1.64	0.46
1:C:586:LEU:O	1:C:594:ARG:NH1	2.49	0.46
1:B:123:TYR:O	1:B:148:SER:HA	2.16	0.45
1:A:369:LEU:HD11	1:B:937:ARG:HA	1.97	0.45
1:A:762:ASP:OD1	1:A:762:ASP:N	2.49	0.45
1:C:841:THR:OG1	1:C:844:ILE:O	2.35	0.45
1:A:254:ILE:HG22	1:A:269:ASP:HA	1.97	0.45
1:A:781:THR:O	1:A:903:GLN:NE2	2.50	0.45
1:B:781:THR:O	1:B:903:GLN:NE2	2.49	0.45
1:C:234:ARG:HA	1:C:234:ARG:HD3	1.77	0.45
2:J:1:NAG:H4	2:J:2:NAG:H2	1.61	0.45
1:B:873:ASN:HB2	1:B:876:LEU:HB2	1.99	0.45
1:C:1036:CYS:HB3	1:C:1088:ASN:HB3	1.98	0.45
1:A:144:VAL:HG13	1:A:145:TYR:HD1	1.82	0.44
1:A:1036:CYS:HB3	1:A:1088:ASN:HB3	1.99	0.44
1:B:700:SER:O	1:B:704:ASN:ND2	2.50	0.44
1:B:866:THR:H	1:B:1060:GLN:HE22	1.65	0.44
1:C:254:ILE:HG22	1:C:269:ASP:HA	1.99	0.44
1:A:1047:GLY:CA	1:A:1059:THR:O	2.63	0.44
1:B:367:ILE:HG23	1:B:487:PRO:HD2	1.99	0.44
1:B:676:VAL:HA	1:B:1018:HIS:O	2.18	0.44
1:A:306:VAL:HG22	1:A:502:ASN:HB3	2.00	0.44
1:A:49:PHE:HB3	1:A:51:VAL:HG13	2.00	0.44
1:C:95:TRP:HB3	1:C:227:PHE:HE1	1.82	0.43
1:C:676:VAL:HA	1:C:1018:HIS:O	2.17	0.43
1:B:1081:ASP:OD1	1:B:1081:ASP:N	2.51	0.43
1:C:894:SER:OG	1:C:895:THR:N	2.52	0.43
1:B:373:SER:OG	1:B:476:GLU:OE2	2.37	0.43
1:A:348:TYR:HB3	1:C:396:VAL:HG13	2.00	0.43
1:A:427:TYR:HB3	1:A:455:TYR:HE1	1.84	0.43
1:B:894:SER:OG	1:B:895:THR:N	2.51	0.43
1:B:1047:GLY:CA	1:B:1059:THR:O	2.63	0.43
1:C:49:PHE:HB3	1:C:51:VAL:HG13	2.00	0.43
1:A:717:LEU:HD22	1:A:962:VAL:HG21	2.00	0.43
1:B:306:VAL:HG22	1:B:502:ASN:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:765:LYS:HD2	1:A:765:LYS:HA	1.85	0.43
1:B:510:GLY:HA2	1:B:550:CYS:HB3	2.01	0.43
1:B:717:LEU:HD22	1:B:962:VAL:HG21	2.01	0.42
1:C:579:ASP:OD1	1:C:579:ASP:N	2.49	0.42
1:B:463:LEU:HD23	1:B:463:LEU:HA	1.86	0.42
1:C:370:CYS:HA	1:C:485:CYS:HA	2.00	0.42
1:A:873:ASN:HB2	1:A:876:LEU:HB2	2.01	0.42
1:B:24:LYS:HG2	1:B:63:TYR:HD1	1.85	0.42
1:C:866:THR:H	1:C:1060:GLN:HE22	1.67	0.42
1:A:894:SER:OG	1:A:895:THR:N	2.52	0.42
1:B:370:CYS:HB3	1:B:485:CYS:HB2	1.80	0.42
1:B:427:TYR:HB3	1:B:455:TYR:HE1	1.84	0.42
1:B:707:LEU:HB2	1:B:951:ILE:HD11	2.01	0.42
1:A:80:GLY:HA2	1:A:179:PHE:O	2.20	0.42
1:B:29:SER:HB2	1:B:58:SER:H	1.85	0.42
1:B:707:LEU:CD1	1:B:951:ILE:HD11	2.48	0.42
1:C:396:VAL:HG11	1:C:427:TYR:HE2	1.85	0.42
1:C:692:CYS:SG	1:C:718:ASN:ND2	2.92	0.42
1:A:38:ASP:OD1	1:A:38:ASP:N	2.45	0.42
1:A:173:ARG:HH21	1:A:192:TYR:HB3	1.84	0.42
1:A:1081:ASP:OD1	1:A:1081:ASP:N	2.51	0.42
1:B:49:PHE:O	1:B:253:THR:HA	2.20	0.42
1:A:577:CYS:HB2	1:A:609:CYS:HB3	1.88	0.42
1:B:174:LEU:HB2	1:B:195:HIS:HB2	2.01	0.42
1:B:690:VAL:HA	1:B:811:GLY:O	2.19	0.42
1:C:738:GLN:NE2	1:C:984:SER:OG	2.44	0.42
1:A:937:ARG:HB3	1:C:361:VAL:HG12	2.00	0.42
1:B:66:TRP:H	1:B:69:ALA:HB2	1.85	0.41
1:B:370:CYS:HA	1:B:485:CYS:HA	2.02	0.41
1:B:538:ASP:HA	1:B:539:PRO:HD3	1.95	0.41
1:B:49:PHE:HB3	1:B:51:VAL:HG13	2.01	0.41
1:C:427:TYR:HB3	1:C:455:TYR:HE1	1.85	0.41
1:C:700:SER:O	1:C:704:ASN:ND2	2.49	0.41
1:A:688:THR:HG21	1:A:913:LEU:HD21	2.01	0.41
1:C:66:TRP:H	1:C:69:ALA:HB2	1.86	0.41
1:C:321:PHE:HE1	1:C:471:VAL:HG11	1.85	0.41
1:A:29:SER:HB2	1:A:58:SER:H	1.86	0.41
1:A:373:SER:OG	1:A:476:GLU:OE2	2.39	0.41
1:A:676:VAL:HA	1:A:1018:HIS:O	2.21	0.41
1:B:694:MET:HG2	1:B:809:PHE:O	2.20	0.41
1:C:781:THR:O	1:C:903:GLN:NE2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:CYS:HA	1:A:485:CYS:HA	2.03	0.41
1:A:174:LEU:HB2	1:A:195:HIS:HB2	2.02	0.41
1:B:97:PHE:HB3	1:B:224:ILE:HD12	2.03	0.41
1:C:669:PRO:HA	1:C:1025:GLN:O	2.21	0.41
1:C:873:ASN:HB2	1:C:876:LEU:HB2	2.02	0.41
1:A:498:CYS:HB2	1:A:550:CYS:HB2	1.94	0.41
1:C:55:ARG:O	1:C:58:SER:OG	2.29	0.41
1:C:398:ALA:HA	1:C:402:TYR:O	2.21	0.41
1:C:538:ASP:OD1	1:C:538:ASP:N	2.48	0.40
1:B:589:ILE:HD12	1:B:593:TRP:HB3	2.03	0.40
1:B:765:LYS:HA	1:B:765:LYS:HD2	1.86	0.40
1:A:566:SER:O	1:A:566:SER:OG	2.32	0.40
1:B:70:TYR:HE1	1:B:231:LEU:HG	1.86	0.40
1:C:707:LEU:HD13	1:C:951:ILE:HD11	2.03	0.40
1:C:831:LEU:HD23	1:C:842:PHE:HE2	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	996/1241 (80%)	929 (93%)	64 (6%)	3 (0%)	37 69
1	B	996/1241 (80%)	928 (93%)	68 (7%)	0	100 100
1	C	996/1241 (80%)	927 (93%)	67 (7%)	2 (0%)	44 74
All	All	2988/3723 (80%)	2784 (93%)	199 (7%)	5 (0%)	45 74

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	810	ASN
1	C	711	GLY

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Mol	Chain	Res	Type
1	A	809	PHE
1	C	396	VAL
1	A	811	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	881/1066 (83%)	879 (100%)	2 (0%)	92	94
1	B	881/1066 (83%)	879 (100%)	2 (0%)	92	94
1	C	881/1066 (83%)	878 (100%)	3 (0%)	91	92
All	All	2643/3198 (83%)	2636 (100%)	7 (0%)	90	92

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	807	GLN
1	A	809	PHE
1	B	807	GLN
1	B	809	PHE
1	C	75	MET
1	C	396	VAL
1	C	397	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates i

18 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	D	1	1,2	14,14,15	0.35	0	17,19,21	0.86	1 (5%)
2	NAG	D	2	2	14,14,15	0.81	1 (7%)	17,19,21	0.68	1 (5%)
2	NAG	E	1	1,2	14,14,15	0.25	0	17,19,21	0.49	0
2	NAG	E	2	2	14,14,15	0.50	0	17,19,21	0.99	1 (5%)
2	NAG	F	1	1,2	14,14,15	0.22	0	17,19,21	0.58	0
2	NAG	F	2	2	14,14,15	0.28	0	17,19,21	0.71	1 (5%)
2	NAG	G	1	1,2	14,14,15	0.38	0	17,19,21	0.88	1 (5%)
2	NAG	G	2	2	14,14,15	0.81	1 (7%)	17,19,21	0.70	1 (5%)
2	NAG	H	1	1,2	14,14,15	0.26	0	17,19,21	0.49	0
2	NAG	H	2	2	14,14,15	0.49	0	17,19,21	0.99	1 (5%)
2	NAG	I	1	1,2	14,14,15	0.22	0	17,19,21	0.64	0
2	NAG	I	2	2	14,14,15	0.27	0	17,19,21	0.71	1 (5%)
2	NAG	J	1	1,2	14,14,15	0.28	0	17,19,21	0.66	0
2	NAG	J	2	2	14,14,15	1.60	2 (14%)	17,19,21	1.41	1 (5%)
2	NAG	K	1	1,2	14,14,15	0.26	0	17,19,21	0.51	0
2	NAG	K	2	2	14,14,15	0.51	0	17,19,21	1.00	1 (5%)
2	NAG	L	1	1,2	14,14,15	0.26	0	17,19,21	0.65	0
2	NAG	L	2	2	14,14,15	0.27	0	17,19,21	0.71	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
2	NAG	D	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	2	2	-	1/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	4/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
2	NAG	H	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	H	2	2	-	1/6/23/26	0/1/1/1
2	NAG	I	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	I	2	2	-	4/6/23/26	0/1/1/1
2	NAG	J	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	NAG	K	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	K	2	2	-	1/6/23/26	0/1/1/1
2	NAG	L	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	L	2	2	-	4/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	2	NAG	O5-C1	5.37	1.52	1.43
2	J	2	NAG	C1-C2	2.56	1.56	1.52
2	G	2	NAG	C1-C2	2.44	1.56	1.52
2	D	2	NAG	C1-C2	2.40	1.55	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	2	NAG	C1-O5-C5	5.59	119.77	112.19
2	K	2	NAG	C2-N2-C7	3.09	127.30	122.90
2	E	2	NAG	C2-N2-C7	3.06	127.26	122.90
2	H	2	NAG	C2-N2-C7	3.05	127.24	122.90
2	G	1	NAG	C1-O5-C5	2.76	115.93	112.19
2	D	1	NAG	C1-O5-C5	2.74	115.91	112.19
2	L	2	NAG	C1-O5-C5	2.39	115.43	112.19
2	F	2	NAG	C1-O5-C5	2.37	115.40	112.19
2	G	2	NAG	C1-O5-C5	2.36	115.40	112.19
2	I	2	NAG	C1-O5-C5	2.36	115.39	112.19
2	D	2	NAG	C1-O5-C5	2.27	115.26	112.19

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	2	NAG	O5-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
2	I	1	NAG	O5-C5-C6-O6
2	L	2	NAG	O5-C5-C6-O6
2	K	1	NAG	O5-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	I	2	NAG	O5-C5-C6-O6
2	I	1	NAG	C4-C5-C6-O6
2	H	1	NAG	O5-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
2	L	2	NAG	C4-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	L	1	NAG	O5-C5-C6-O6
2	K	1	NAG	C4-C5-C6-O6
2	L	1	NAG	C4-C5-C6-O6
2	F	2	NAG	C8-C7-N2-C2
2	F	2	NAG	O7-C7-N2-C2
2	I	2	NAG	C8-C7-N2-C2
2	I	2	NAG	O7-C7-N2-C2
2	L	2	NAG	C8-C7-N2-C2
2	L	2	NAG	O7-C7-N2-C2
2	F	1	NAG	O5-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	H	1	NAG	C4-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	I	2	NAG	C4-C5-C6-O6
2	J	2	NAG	C4-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
2	G	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C3-C2-N2-C7
2	H	2	NAG	C3-C2-N2-C7
2	K	2	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

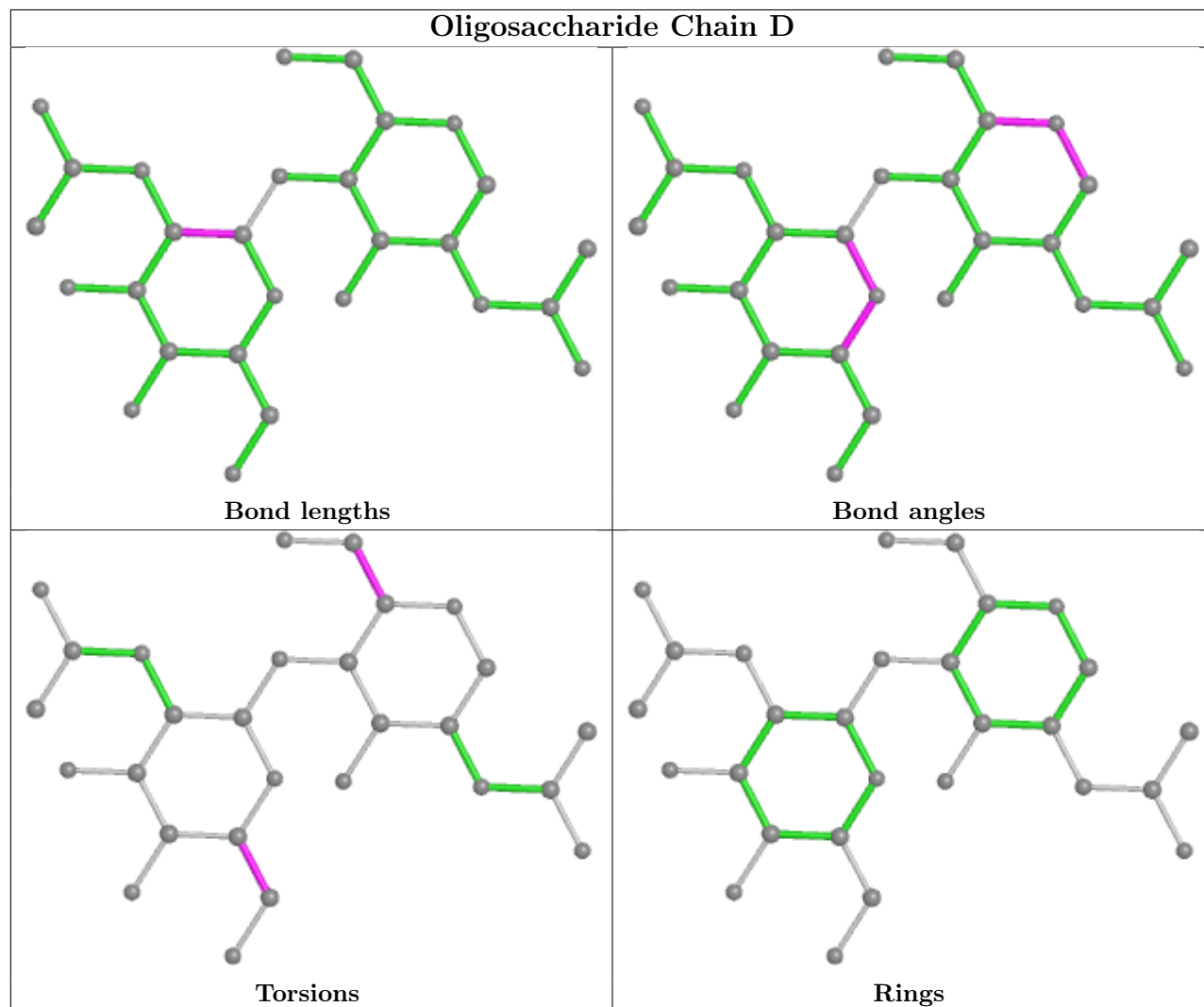
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	1	NAG	2	0

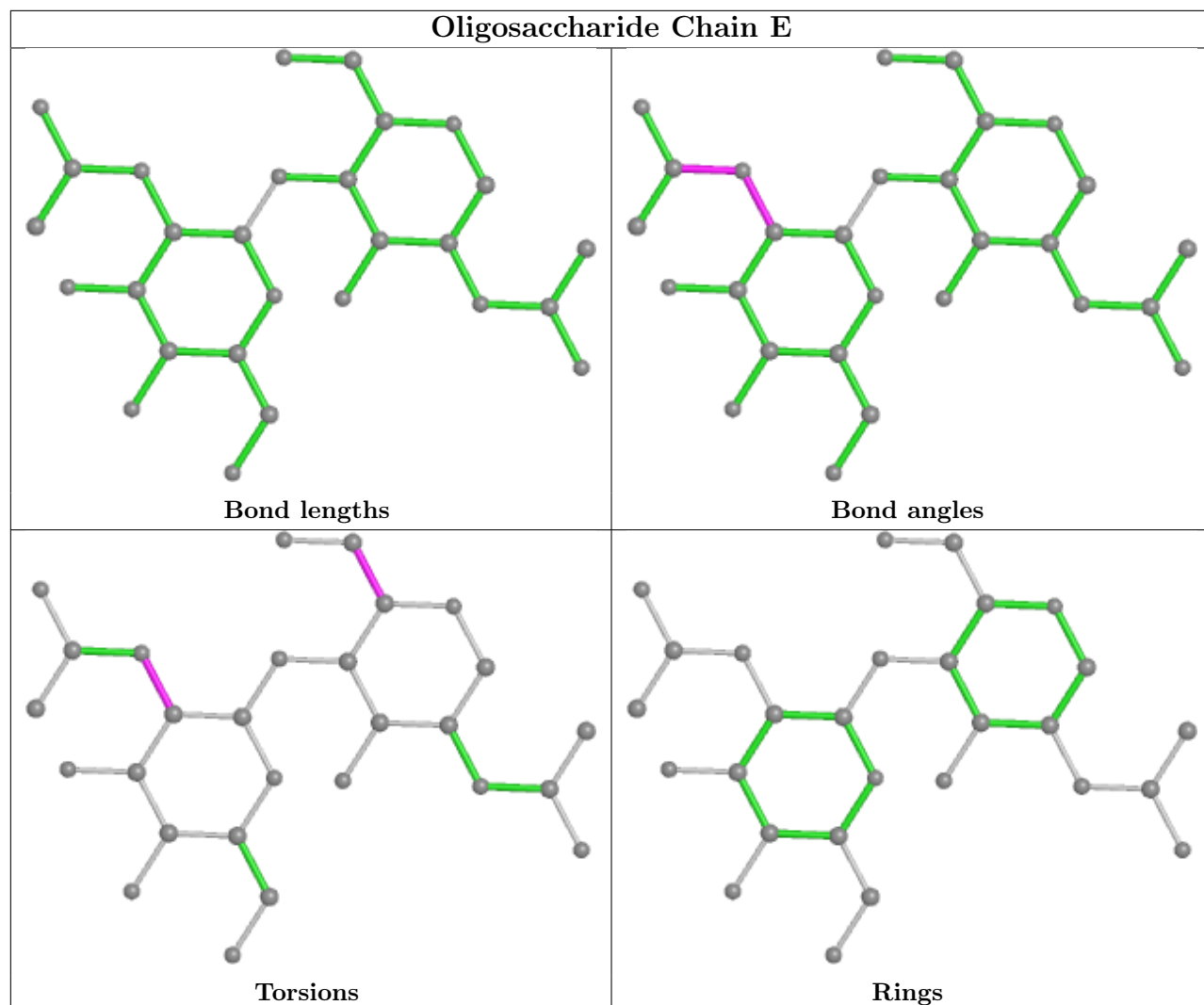
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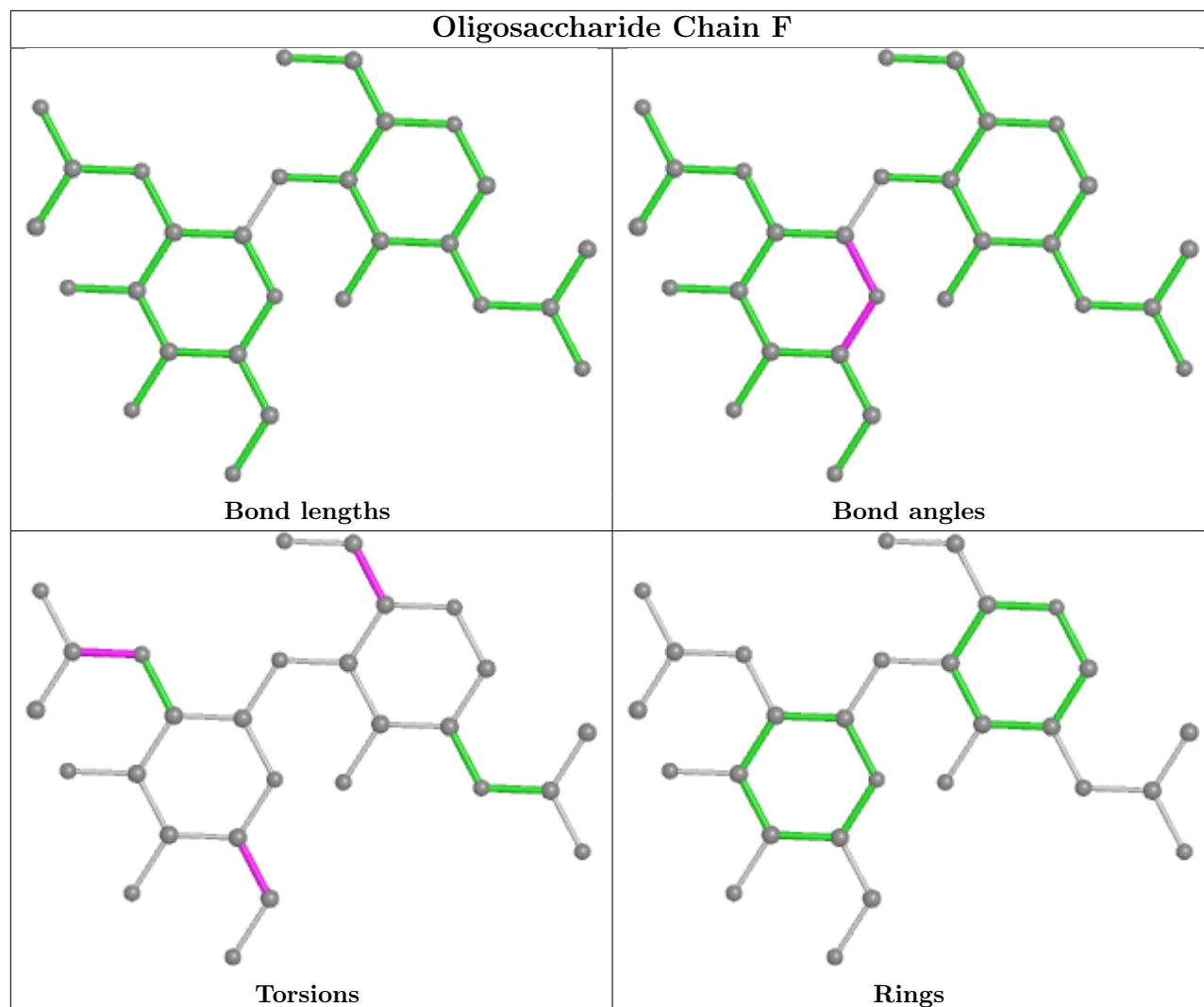
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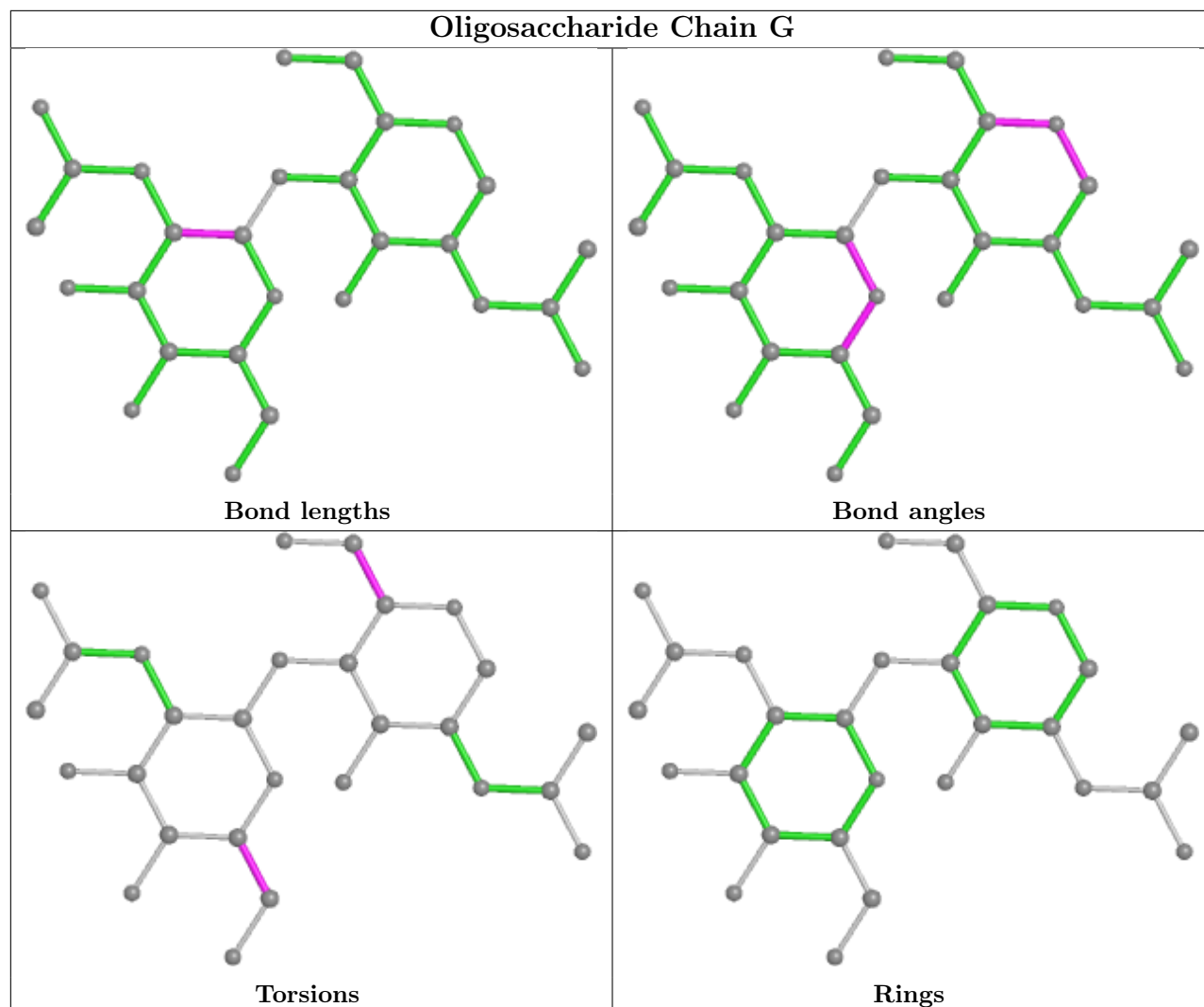
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	2	NAG	2	0

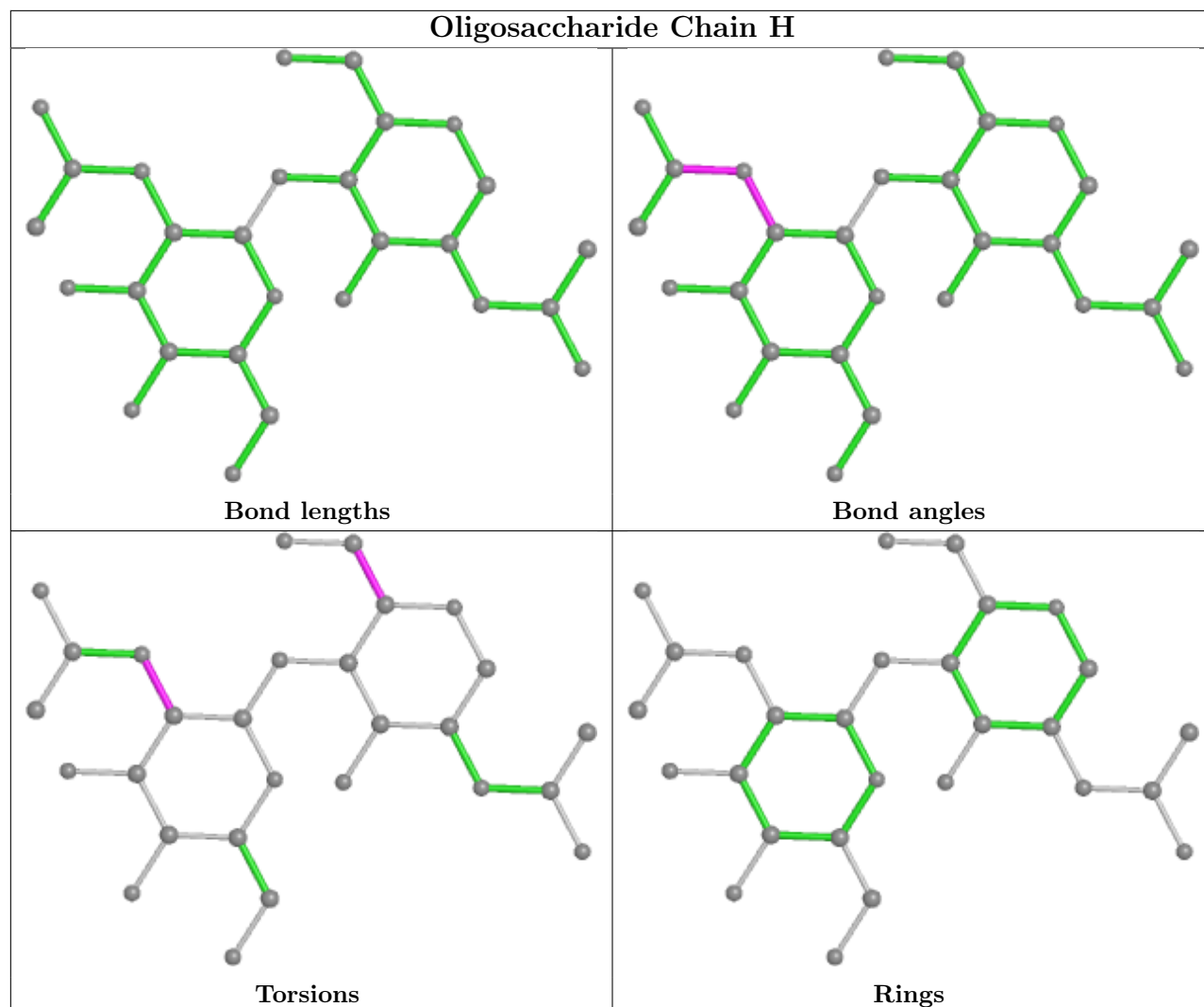
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

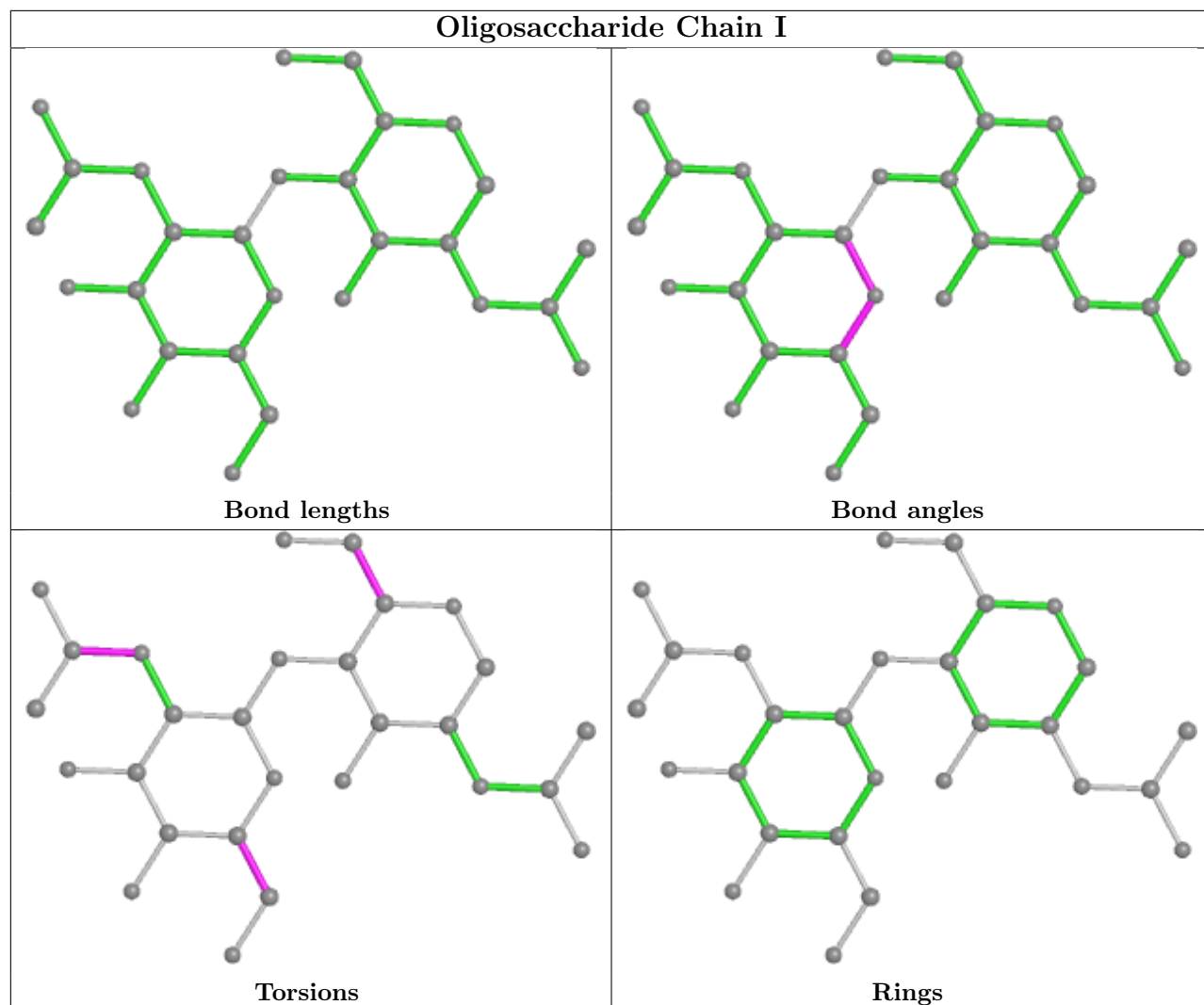


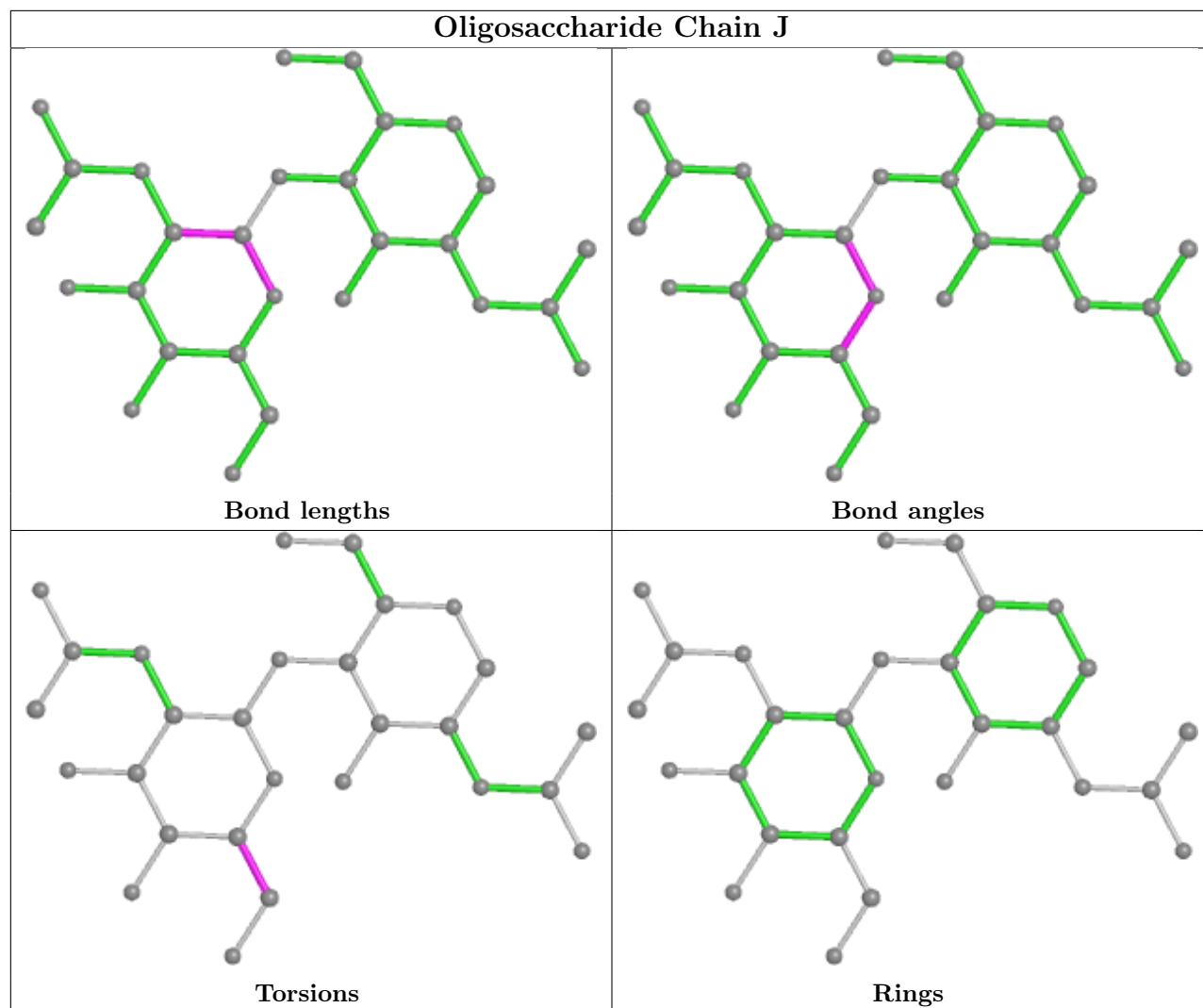


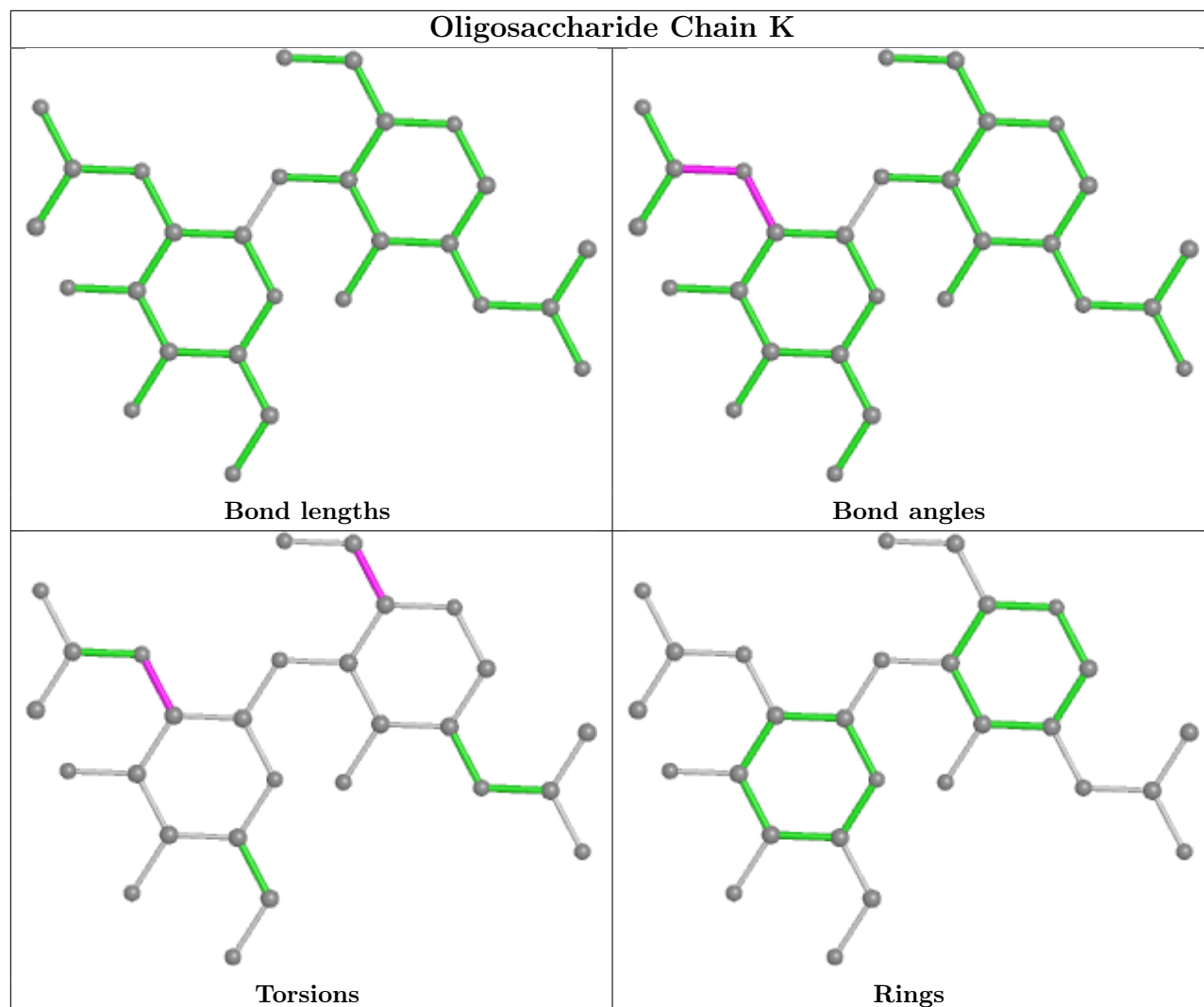


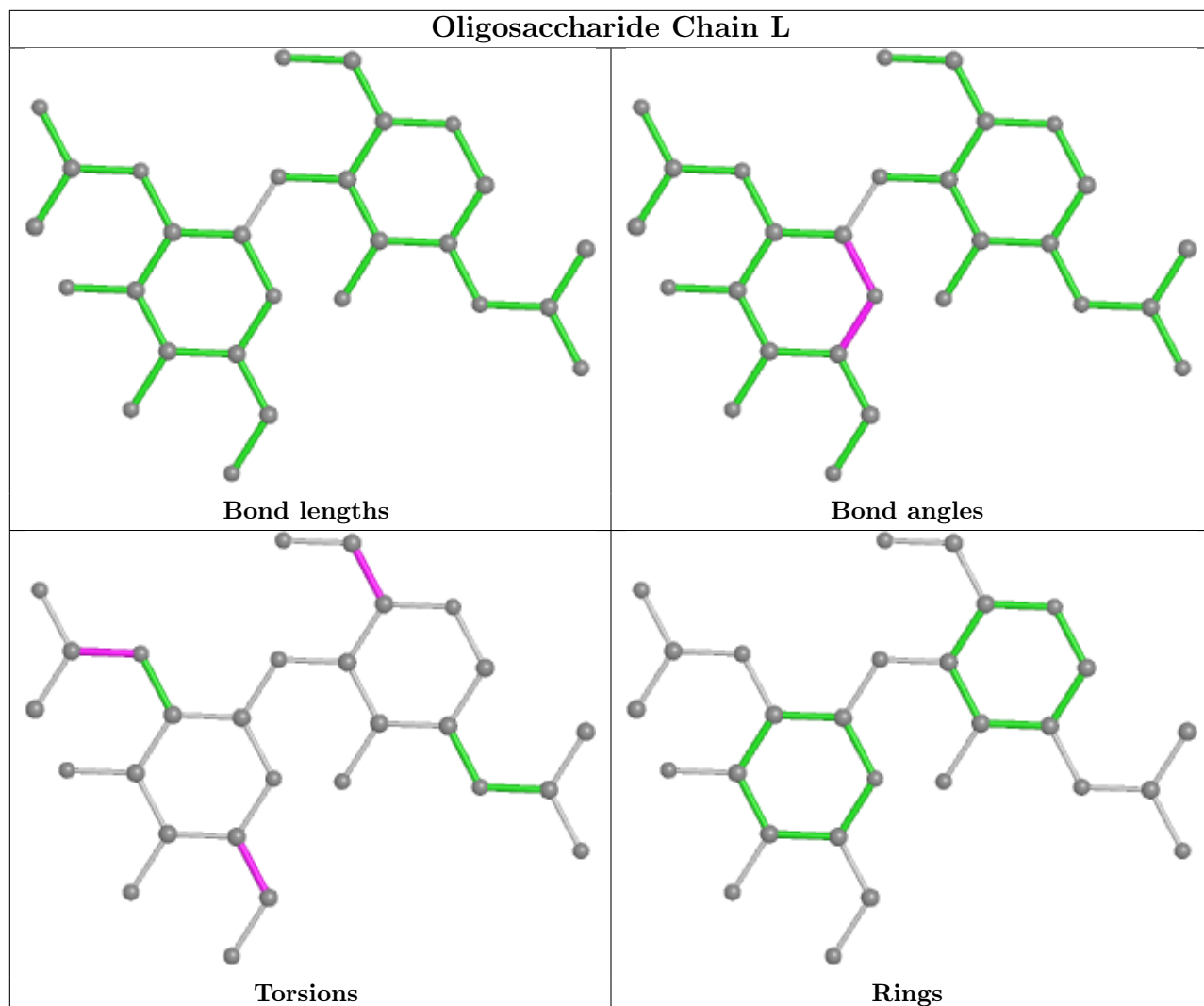












5.6 Ligand geometry [i](#)

45 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	$\# Z > 2$
3	NAG	B	1301	1	14,14,15	0.32	0	17,19,21	0.47	0
3	NAG	C	1305	1	14,14,15	0.37	0	17,19,21	0.51	0
3	NAG	A	1302	1	14,14,15	0.86	1 (7%)	17,19,21	0.66	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	C	1308	1	14,14,15	0.34	0	17,19,21	0.43	0
3	NAG	C	1315	1	14,14,15	0.33	0	17,19,21	0.55	0
3	NAG	B	1309	1	14,14,15	0.38	0	17,19,21	0.97	1 (5%)
3	NAG	B	1315	1	14,14,15	0.31	0	17,19,21	0.55	0
3	NAG	A	1301	1	14,14,15	0.32	0	17,19,21	0.48	0
3	NAG	A	1309	1	14,14,15	0.38	0	17,19,21	0.97	1 (5%)
3	NAG	A	1312	1	14,14,15	0.44	0	17,19,21	0.55	0
3	NAG	C	1309	1	14,14,15	0.38	0	17,19,21	0.97	1 (5%)
3	NAG	A	1303	-	14,14,15	0.33	0	17,19,21	0.47	0
3	NAG	C	1311	1	14,14,15	0.78	1 (7%)	17,19,21	1.15	2 (11%)
3	NAG	A	1308	1	14,14,15	0.32	0	17,19,21	0.40	0
3	NAG	A	1315	1	14,14,15	0.33	0	17,19,21	0.56	0
3	NAG	B	1311	1	14,14,15	0.51	0	17,19,21	1.01	1 (5%)
3	NAG	C	1310	1	14,14,15	0.53	0	17,19,21	0.41	0
3	NAG	B	1308	1	14,14,15	0.31	0	17,19,21	0.41	0
3	NAG	C	1301	1	14,14,15	0.33	0	17,19,21	0.48	0
3	NAG	C	1307	1	14,14,15	0.34	0	17,19,21	0.50	0
3	NAG	C	1304	1	14,14,15	0.40	0	17,19,21	0.52	0
3	NAG	B	1314	1	14,14,15	0.46	0	17,19,21	0.52	0
3	NAG	C	1306	1	14,14,15	0.48	0	17,19,21	0.57	0
3	NAG	B	1305	1	14,14,15	0.38	0	17,19,21	0.53	0
3	NAG	A	1314	1	14,14,15	0.46	0	17,19,21	0.48	0
3	NAG	C	1313	1	14,14,15	0.31	0	17,19,21	0.50	0
3	NAG	B	1306	1	14,14,15	0.46	0	17,19,21	0.57	0
3	NAG	A	1305	1	14,14,15	0.37	0	17,19,21	0.52	0
3	NAG	C	1302	1	14,14,15	0.86	1 (7%)	17,19,21	0.65	1 (5%)
3	NAG	C	1312	1	14,14,15	0.44	0	17,19,21	0.55	0
3	NAG	B	1307	1	14,14,15	0.34	0	17,19,21	0.49	0
3	NAG	B	1313	1	14,14,15	0.29	0	17,19,21	0.46	0
3	NAG	B	1303	-	14,14,15	0.32	0	17,19,21	0.48	0
3	NAG	A	1311	1	14,14,15	0.47	0	17,19,21	1.00	1 (5%)
3	NAG	B	1310	1	14,14,15	0.49	0	17,19,21	0.58	0
3	NAG	A	1307	1	14,14,15	0.33	0	17,19,21	0.48	0
3	NAG	C	1303	-	14,14,15	0.32	0	17,19,21	0.48	0
3	NAG	A	1310	1	14,14,15	0.47	0	17,19,21	0.58	0
3	NAG	A	1304	1	14,14,15	0.42	0	17,19,21	0.54	0
3	NAG	B	1302	1	14,14,15	0.86	1 (7%)	17,19,21	0.65	1 (5%)
3	NAG	A	1306	1	14,14,15	0.46	0	17,19,21	0.57	0
3	NAG	B	1312	1	14,14,15	0.52	0	17,19,21	0.57	0
3	NAG	B	1304	1	14,14,15	0.44	0	17,19,21	0.54	0
3	NAG	A	1313	1	14,14,15	0.27	0	17,19,21	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	C	1314	1	14,14,15	0.46	0	17,19,21	0.51	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	1301	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1305	1	-	1/6/23/26	0/1/1/1
3	NAG	A	1302	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1308	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1315	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1309	1	-	3/6/23/26	0/1/1/1
3	NAG	B	1315	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1301	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1309	1	-	3/6/23/26	0/1/1/1
3	NAG	A	1312	1	-	4/6/23/26	0/1/1/1
3	NAG	C	1309	1	-	3/6/23/26	0/1/1/1
3	NAG	A	1303	-	-	0/6/23/26	0/1/1/1
3	NAG	C	1311	1	-	1/6/23/26	0/1/1/1
3	NAG	A	1308	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1315	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1311	1	-	3/6/23/26	0/1/1/1
3	NAG	C	1310	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1308	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1301	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1307	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1304	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1314	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1306	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1305	1	-	1/6/23/26	0/1/1/1
3	NAG	A	1314	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1313	1	-	4/6/23/26	0/1/1/1
3	NAG	B	1306	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1305	1	-	1/6/23/26	0/1/1/1
3	NAG	C	1302	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1312	1	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	1307	1	-	1/6/23/26	0/1/1/1
3	NAG	B	1313	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1303	-	-	0/6/23/26	0/1/1/1
3	NAG	A	1311	1	-	3/6/23/26	0/1/1/1
3	NAG	B	1310	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1307	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1303	-	-	0/6/23/26	0/1/1/1
3	NAG	A	1310	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1304	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1302	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1306	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1312	1	-	3/6/23/26	0/1/1/1
3	NAG	B	1304	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1313	1	-	4/6/23/26	0/1/1/1
3	NAG	C	1314	1	-	0/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1302	NAG	O5-C1	2.85	1.48	1.43
3	B	1302	NAG	O5-C1	2.83	1.48	1.43
3	A	1302	NAG	O5-C1	2.82	1.48	1.43
3	C	1311	NAG	C1-C2	2.41	1.55	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1309	NAG	C2-N2-C7	3.16	127.41	122.90
3	A	1309	NAG	C2-N2-C7	3.15	127.39	122.90
3	C	1309	NAG	C2-N2-C7	3.13	127.35	122.90
3	B	1311	NAG	C2-N2-C7	3.05	127.25	122.90
3	A	1311	NAG	C2-N2-C7	3.05	127.24	122.90
3	C	1311	NAG	C2-N2-C7	3.03	127.21	122.90
3	C	1311	NAG	C1-O5-C5	2.94	116.17	112.19
3	A	1302	NAG	C1-O5-C5	2.42	115.47	112.19
3	B	1302	NAG	C1-O5-C5	2.41	115.46	112.19
3	C	1302	NAG	C1-O5-C5	2.38	115.42	112.19

There are no chirality outliers.

All (85) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1315	NAG	O5-C5-C6-O6
3	B	1310	NAG	O5-C5-C6-O6
3	C	1309	NAG	O5-C5-C6-O6
3	A	1311	NAG	O5-C5-C6-O6
3	B	1311	NAG	O5-C5-C6-O6
3	A	1309	NAG	O5-C5-C6-O6
3	A	1310	NAG	O5-C5-C6-O6
3	B	1309	NAG	O5-C5-C6-O6
3	B	1315	NAG	O5-C5-C6-O6
3	C	1315	NAG	O5-C5-C6-O6
3	A	1311	NAG	C4-C5-C6-O6
3	C	1309	NAG	C4-C5-C6-O6
3	C	1312	NAG	O5-C5-C6-O6
3	A	1315	NAG	C4-C5-C6-O6
3	B	1310	NAG	C4-C5-C6-O6
3	A	1312	NAG	O5-C5-C6-O6
3	A	1309	NAG	C4-C5-C6-O6
3	C	1302	NAG	C4-C5-C6-O6
3	B	1309	NAG	C4-C5-C6-O6
3	A	1302	NAG	C4-C5-C6-O6
3	B	1301	NAG	C4-C5-C6-O6
3	A	1310	NAG	C4-C5-C6-O6
3	A	1307	NAG	C4-C5-C6-O6
3	B	1311	NAG	C4-C5-C6-O6
3	C	1315	NAG	C4-C5-C6-O6
3	A	1312	NAG	C8-C7-N2-C2
3	A	1312	NAG	O7-C7-N2-C2
3	A	1313	NAG	C8-C7-N2-C2
3	A	1313	NAG	O7-C7-N2-C2
3	B	1312	NAG	C8-C7-N2-C2
3	B	1312	NAG	O7-C7-N2-C2
3	B	1313	NAG	C8-C7-N2-C2
3	B	1313	NAG	O7-C7-N2-C2
3	C	1312	NAG	C8-C7-N2-C2
3	C	1312	NAG	O7-C7-N2-C2
3	C	1313	NAG	C8-C7-N2-C2
3	C	1313	NAG	O7-C7-N2-C2
3	A	1301	NAG	C4-C5-C6-O6
3	B	1302	NAG	C4-C5-C6-O6
3	B	1315	NAG	C4-C5-C6-O6
3	C	1301	NAG	C4-C5-C6-O6
3	A	1304	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	B	1304	NAG	C4-C5-C6-O6
3	A	1307	NAG	O5-C5-C6-O6
3	C	1302	NAG	O5-C5-C6-O6
3	A	1312	NAG	C4-C5-C6-O6
3	C	1306	NAG	C4-C5-C6-O6
3	C	1313	NAG	O5-C5-C6-O6
3	A	1302	NAG	O5-C5-C6-O6
3	B	1301	NAG	O5-C5-C6-O6
3	C	1313	NAG	C4-C5-C6-O6
3	A	1301	NAG	O5-C5-C6-O6
3	C	1301	NAG	O5-C5-C6-O6
3	A	1306	NAG	C4-C5-C6-O6
3	C	1312	NAG	C4-C5-C6-O6
3	B	1302	NAG	O5-C5-C6-O6
3	A	1314	NAG	C4-C5-C6-O6
3	A	1304	NAG	O5-C5-C6-O6
3	B	1304	NAG	O5-C5-C6-O6
3	C	1305	NAG	O5-C5-C6-O6
3	B	1306	NAG	C4-C5-C6-O6
3	C	1304	NAG	C4-C5-C6-O6
3	B	1312	NAG	O5-C5-C6-O6
3	A	1305	NAG	O5-C5-C6-O6
3	B	1305	NAG	O5-C5-C6-O6
3	B	1314	NAG	C4-C5-C6-O6
3	C	1306	NAG	O5-C5-C6-O6
3	A	1308	NAG	C4-C5-C6-O6
3	A	1306	NAG	O5-C5-C6-O6
3	A	1314	NAG	O5-C5-C6-O6
3	B	1306	NAG	O5-C5-C6-O6
3	C	1304	NAG	O5-C5-C6-O6
3	A	1308	NAG	O5-C5-C6-O6
3	B	1314	NAG	O5-C5-C6-O6
3	C	1311	NAG	C3-C2-N2-C7
3	A	1313	NAG	C4-C5-C6-O6
3	B	1308	NAG	C4-C5-C6-O6
3	B	1307	NAG	C4-C5-C6-O6
3	A	1313	NAG	O5-C5-C6-O6
3	A	1309	NAG	C3-C2-N2-C7
3	A	1311	NAG	C3-C2-N2-C7
3	B	1309	NAG	C3-C2-N2-C7
3	B	1311	NAG	C3-C2-N2-C7
3	C	1309	NAG	C3-C2-N2-C7

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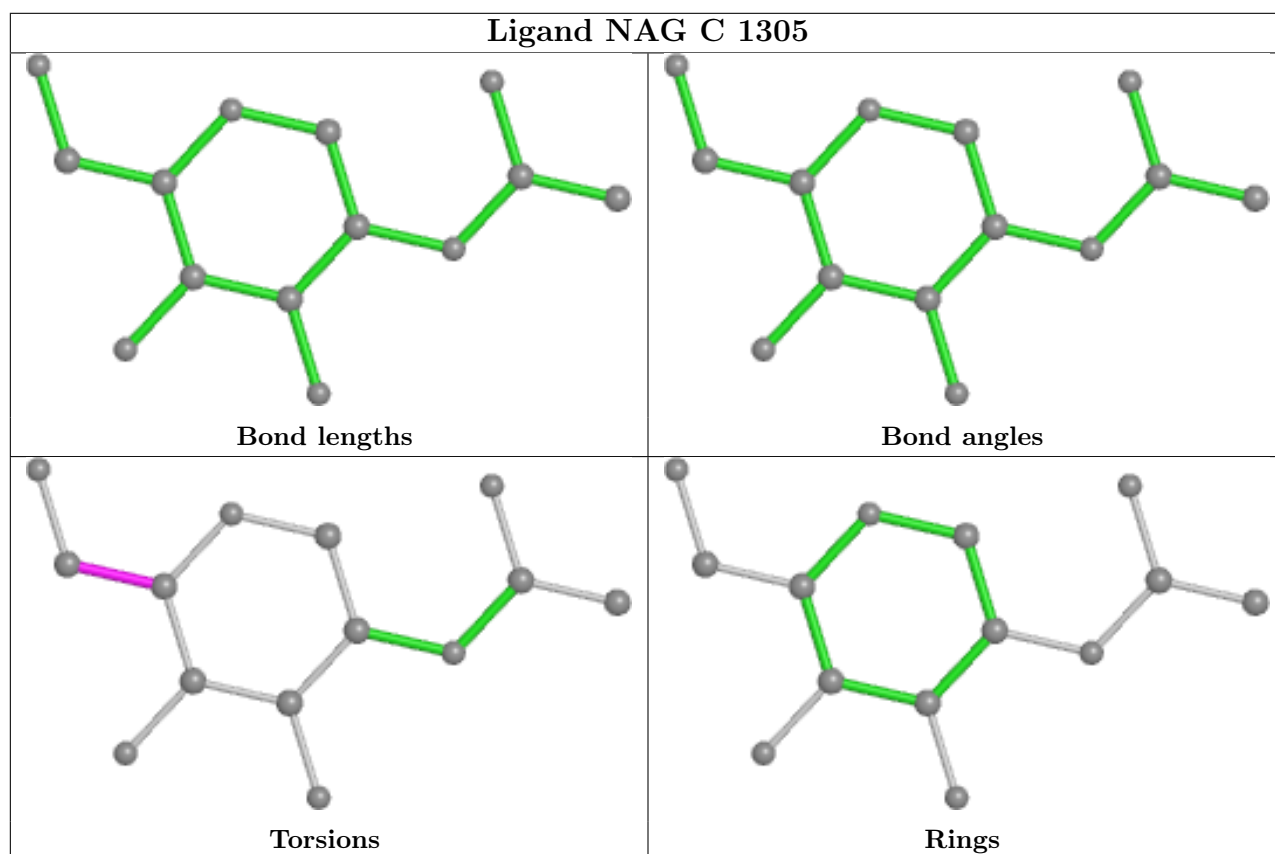
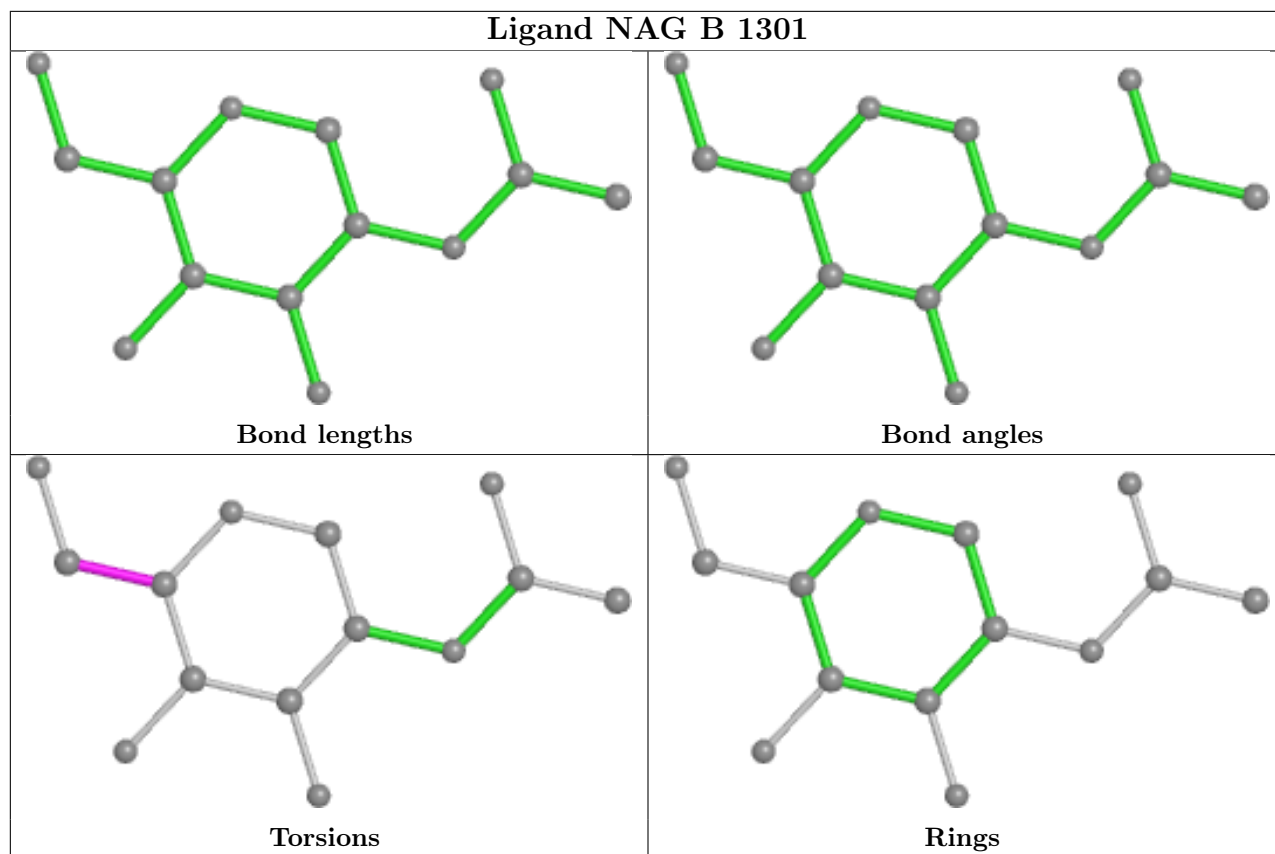
Mol	Chain	Res	Type	Atoms
3	B	1308	NAG	O5-C5-C6-O6

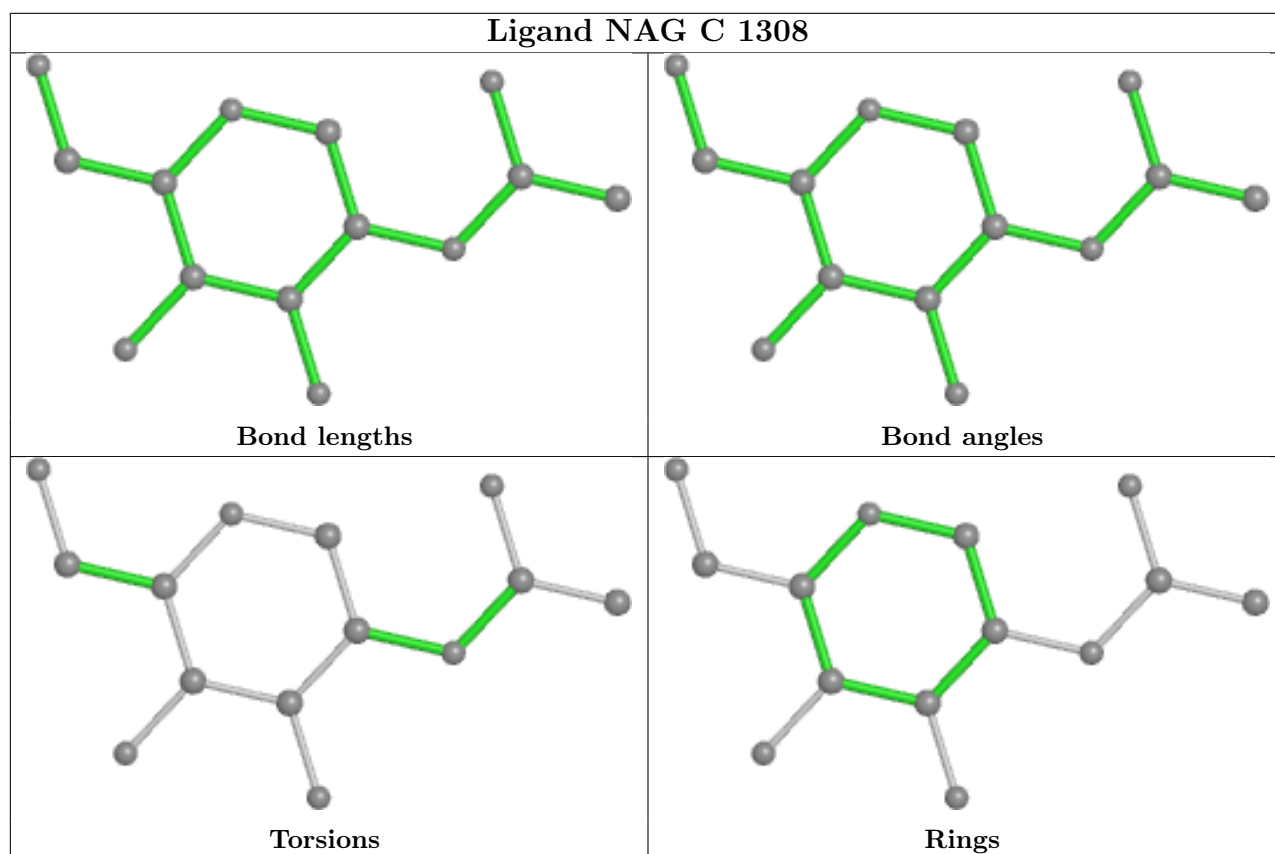
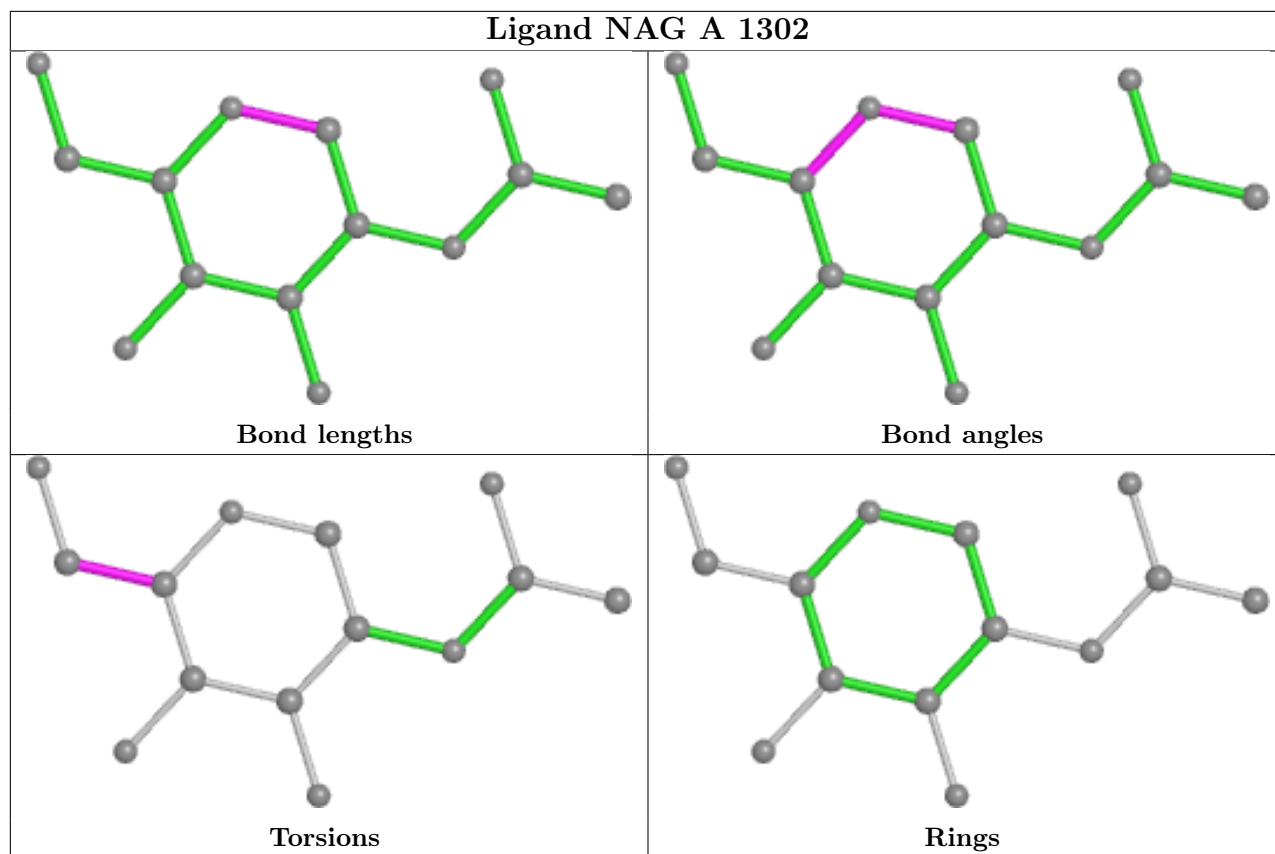
There are no ring outliers.

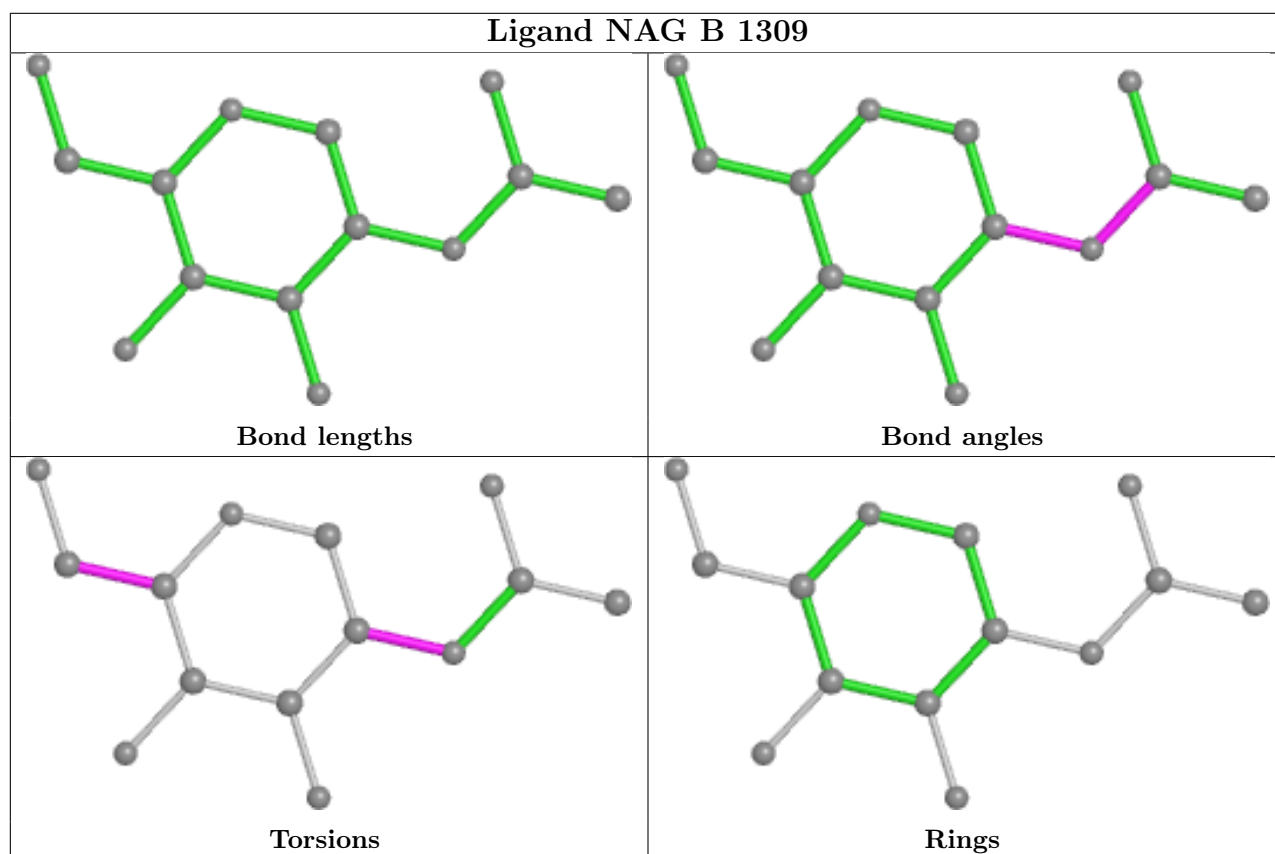
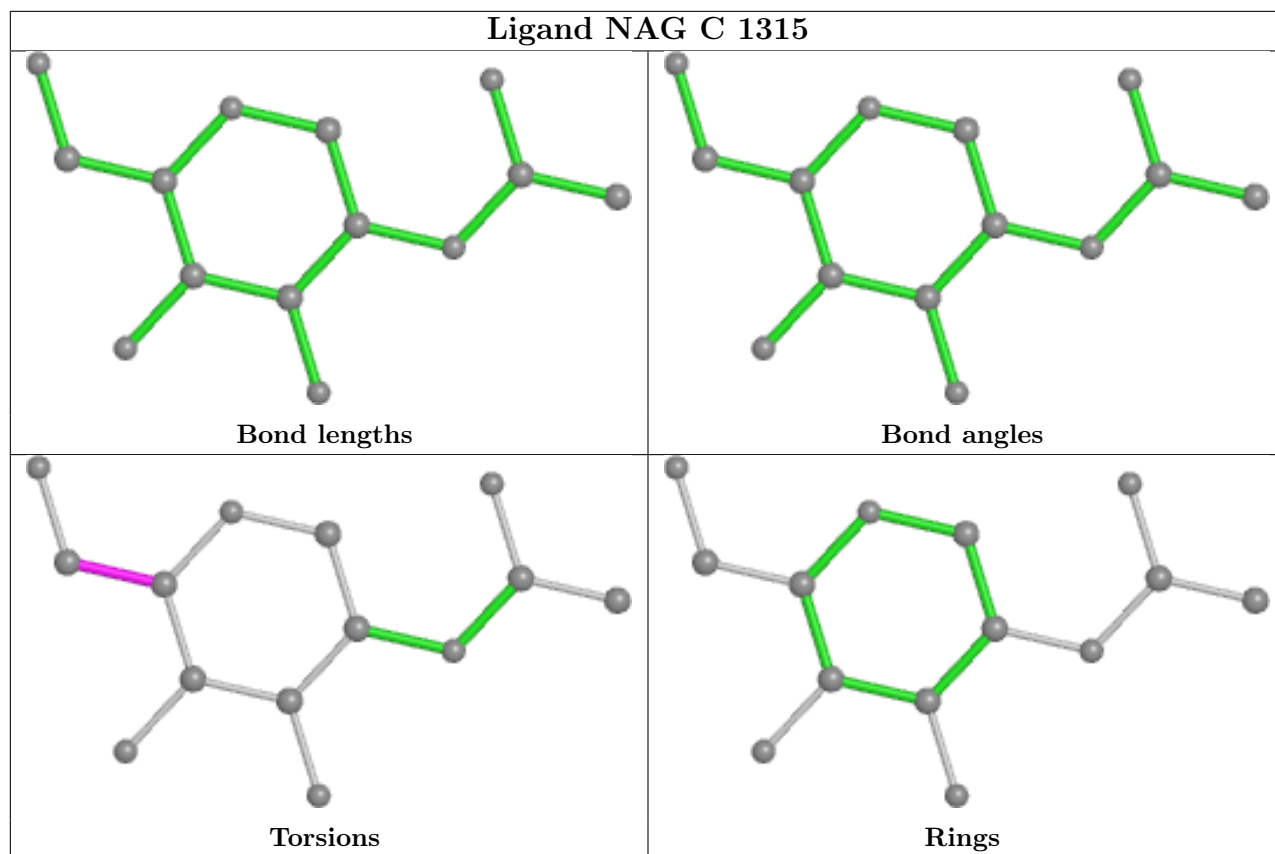
3 monomers are involved in 3 short contacts:

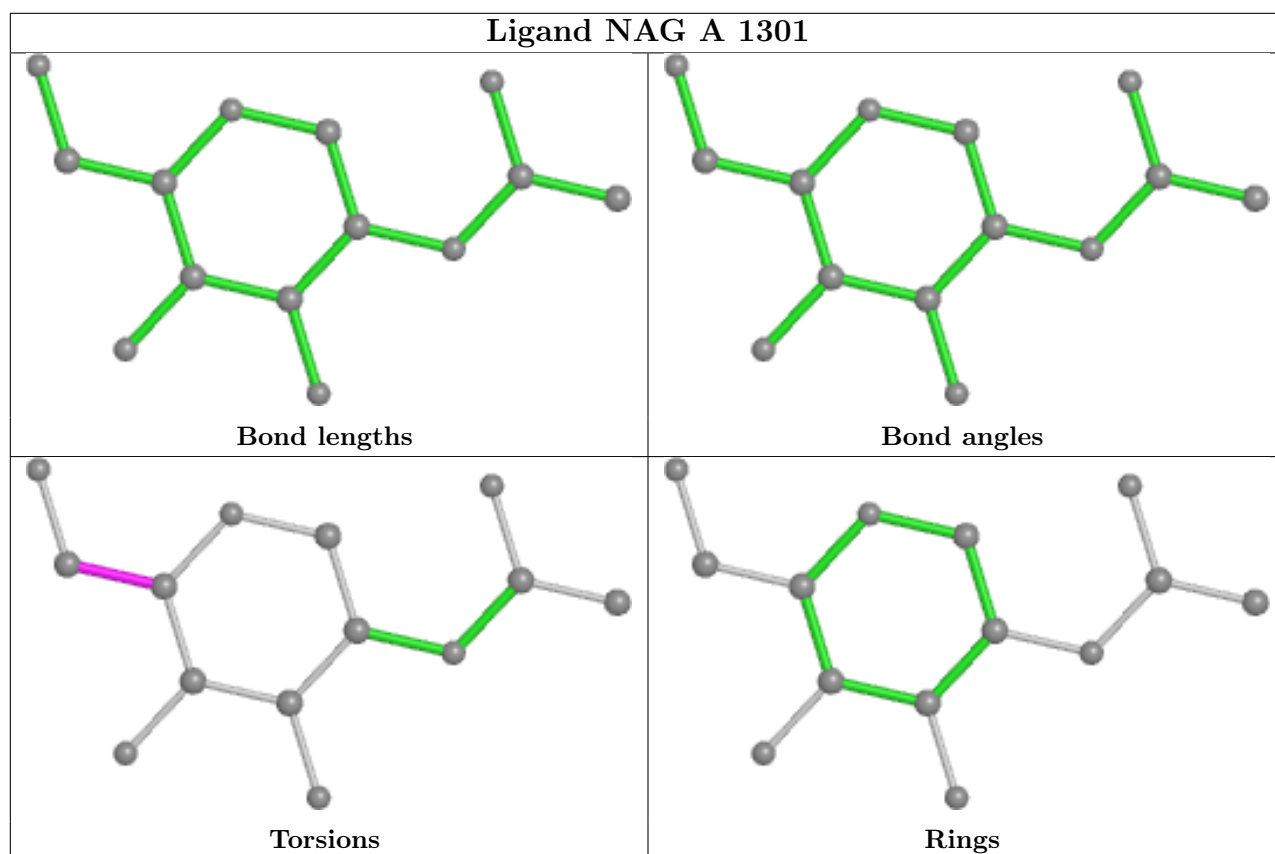
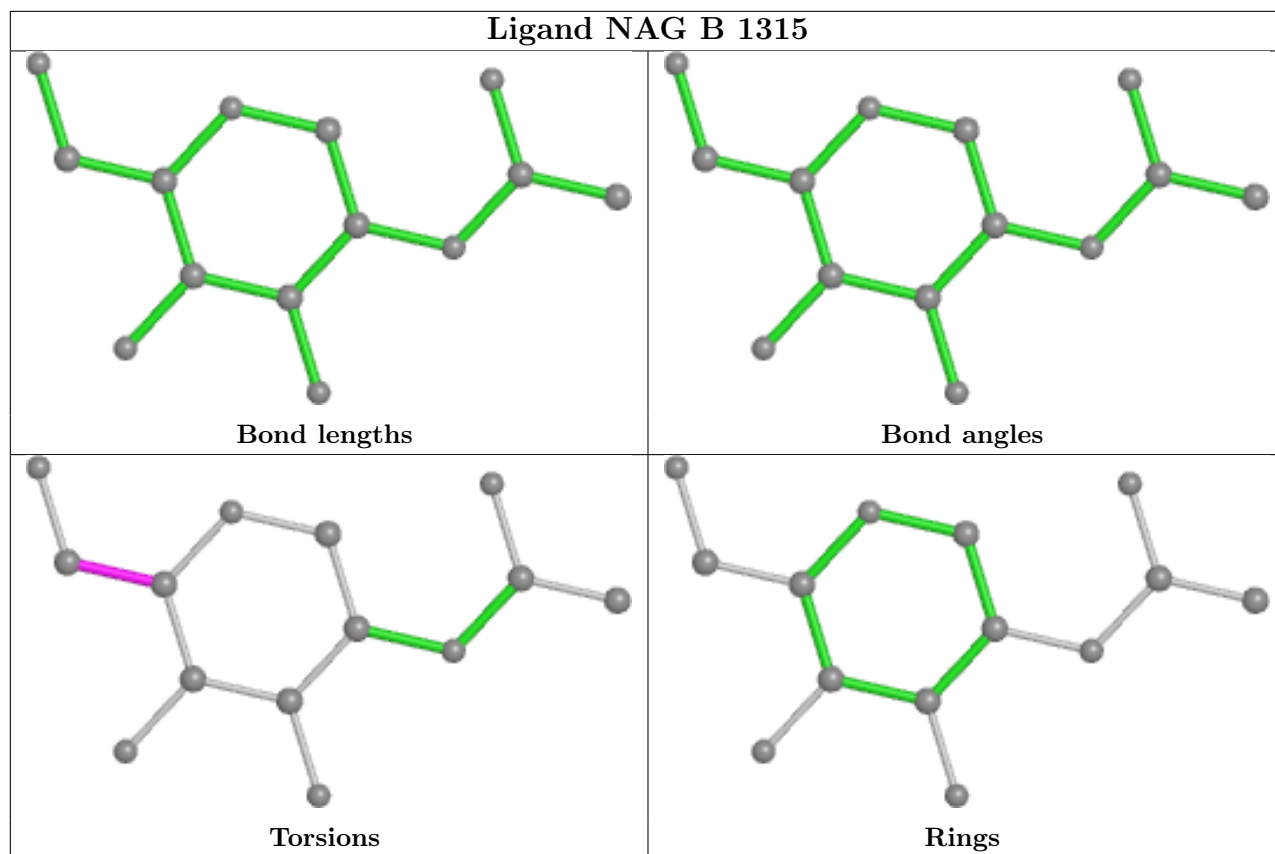
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1312	NAG	1	0
3	C	1312	NAG	1	0
3	B	1312	NAG	1	0

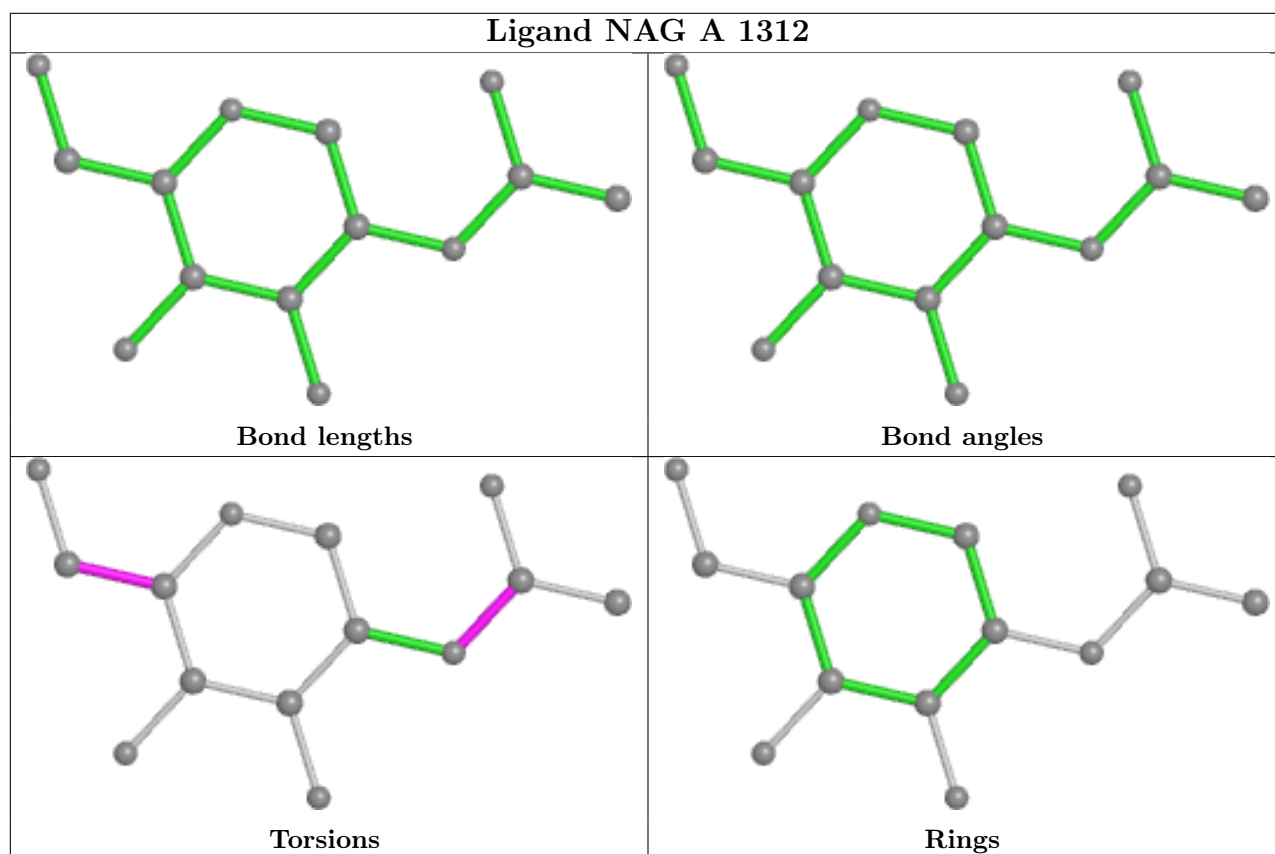
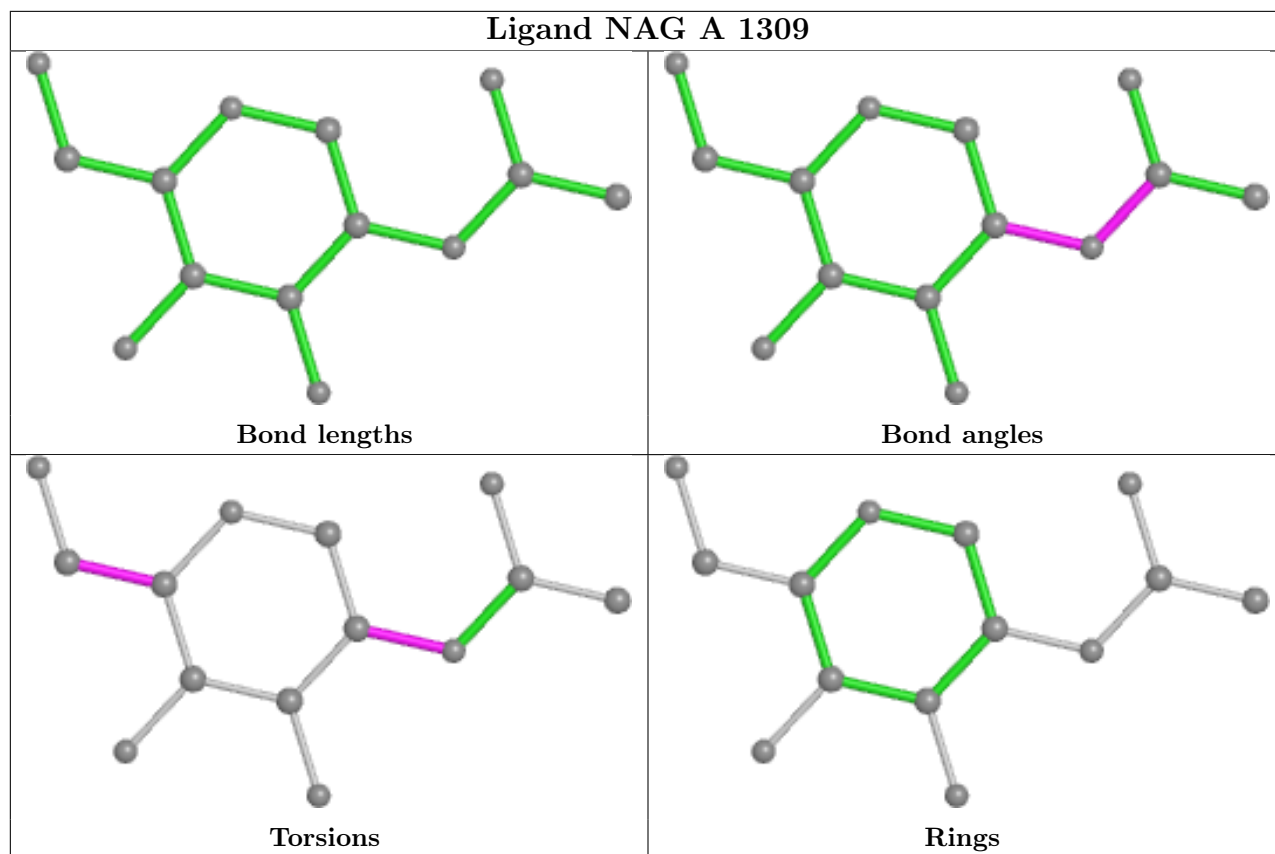
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

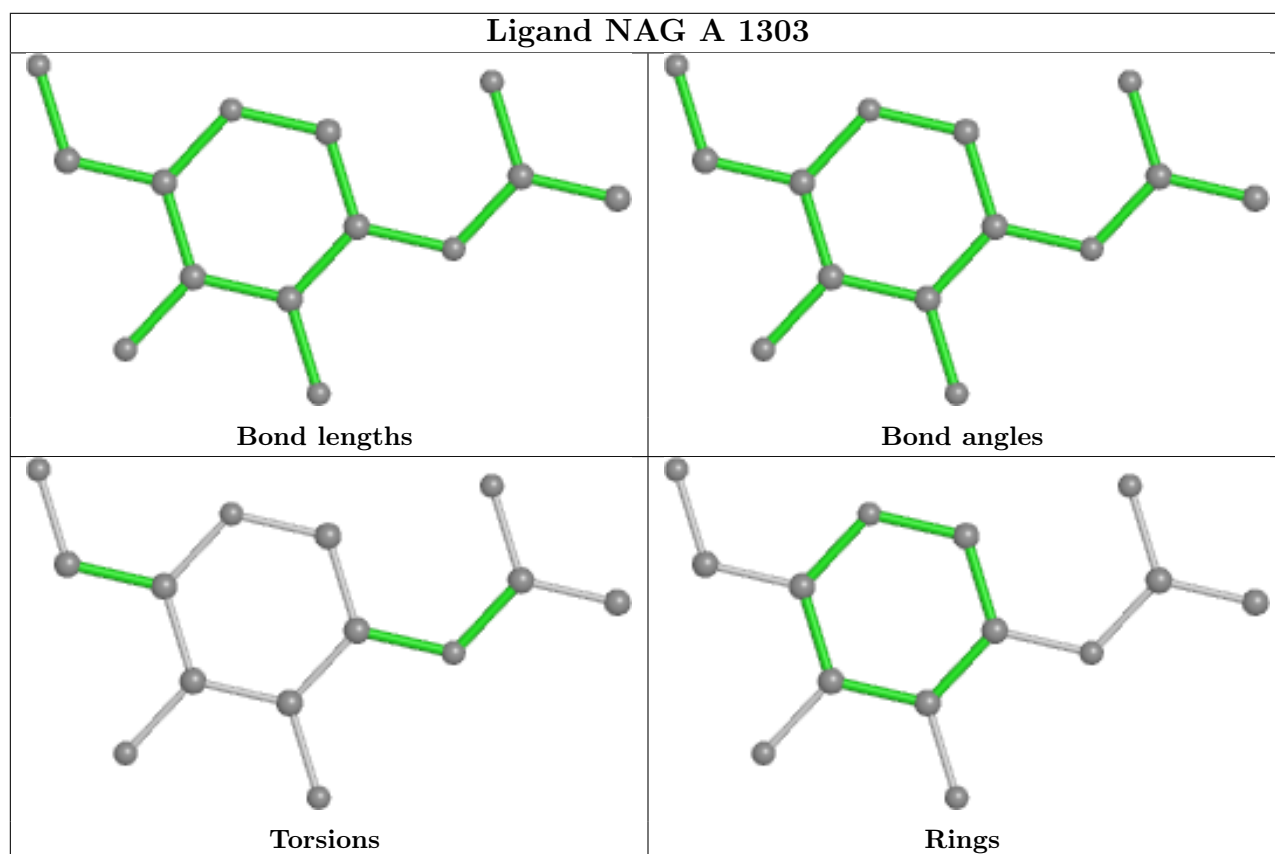
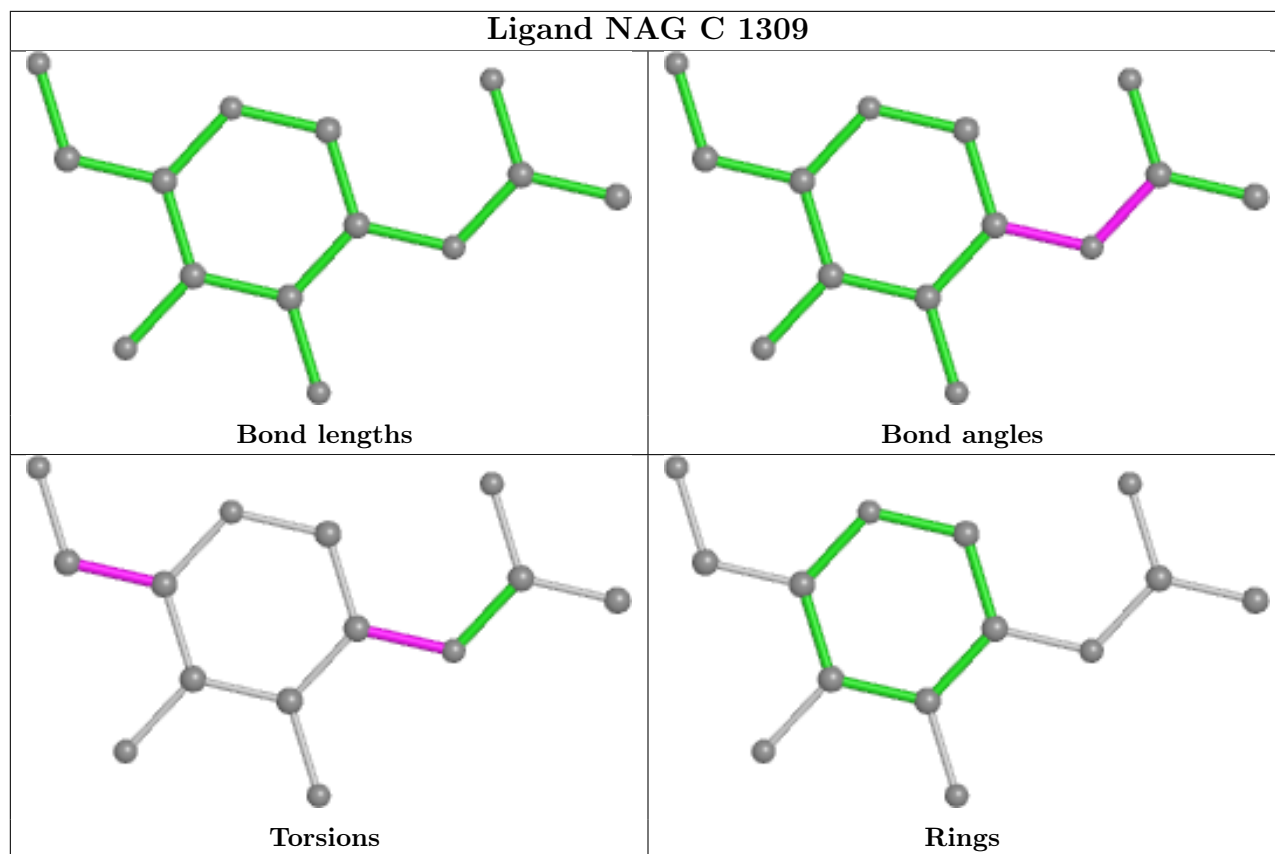


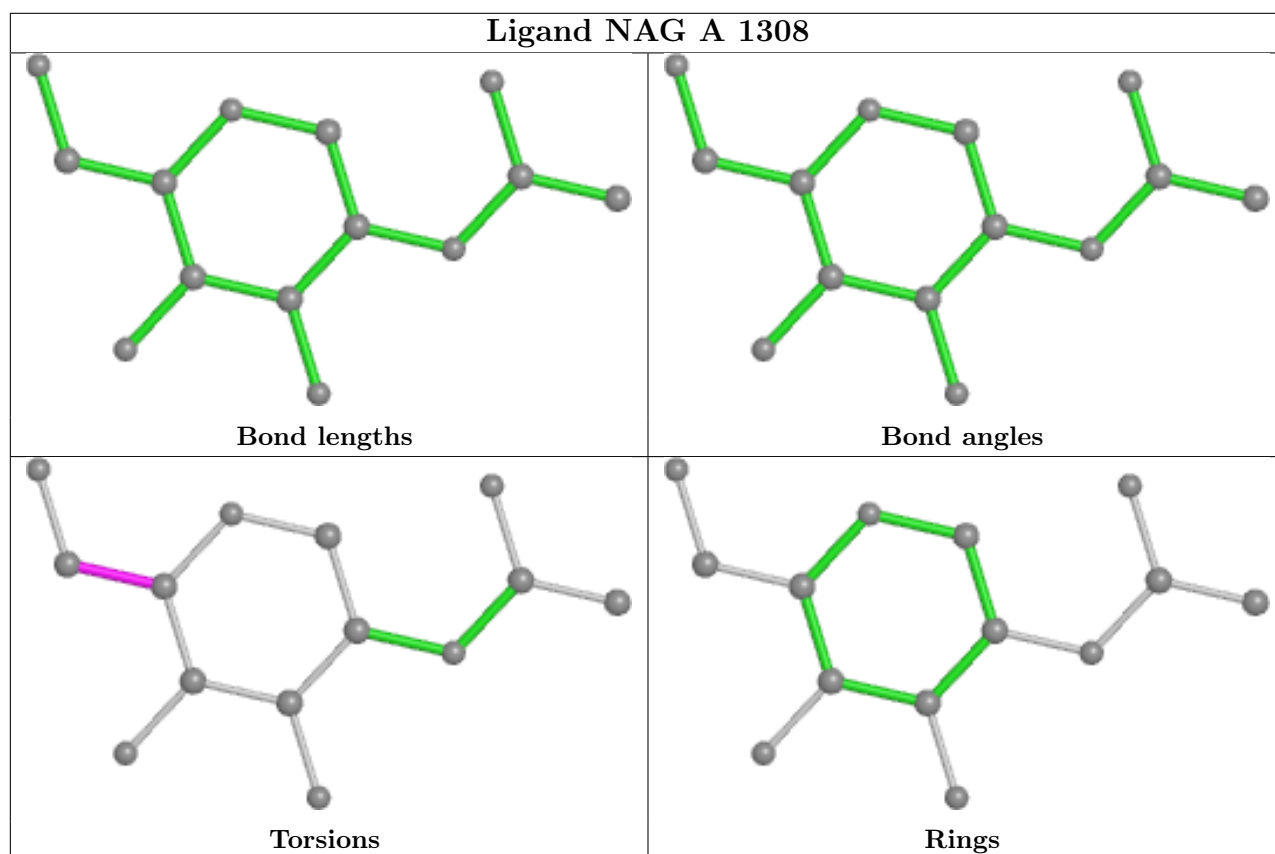
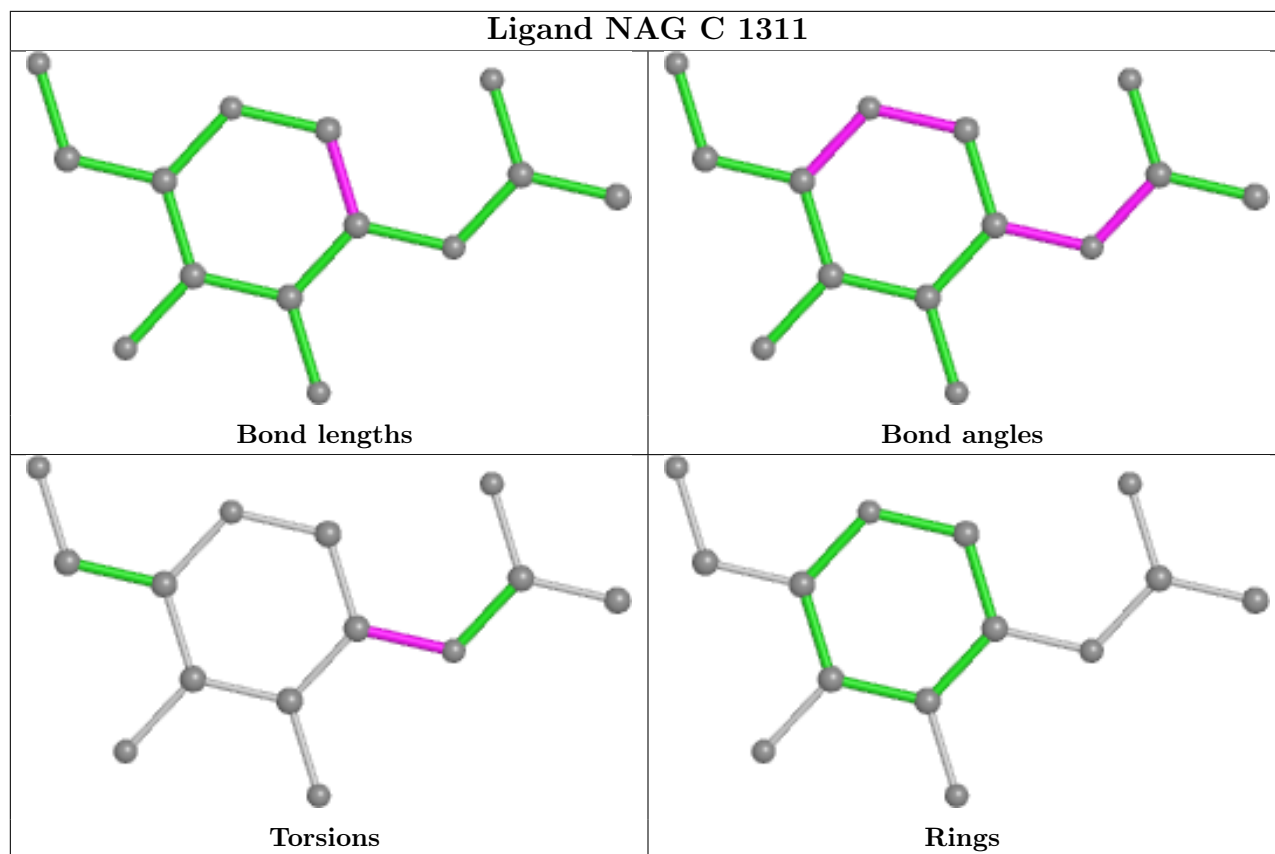


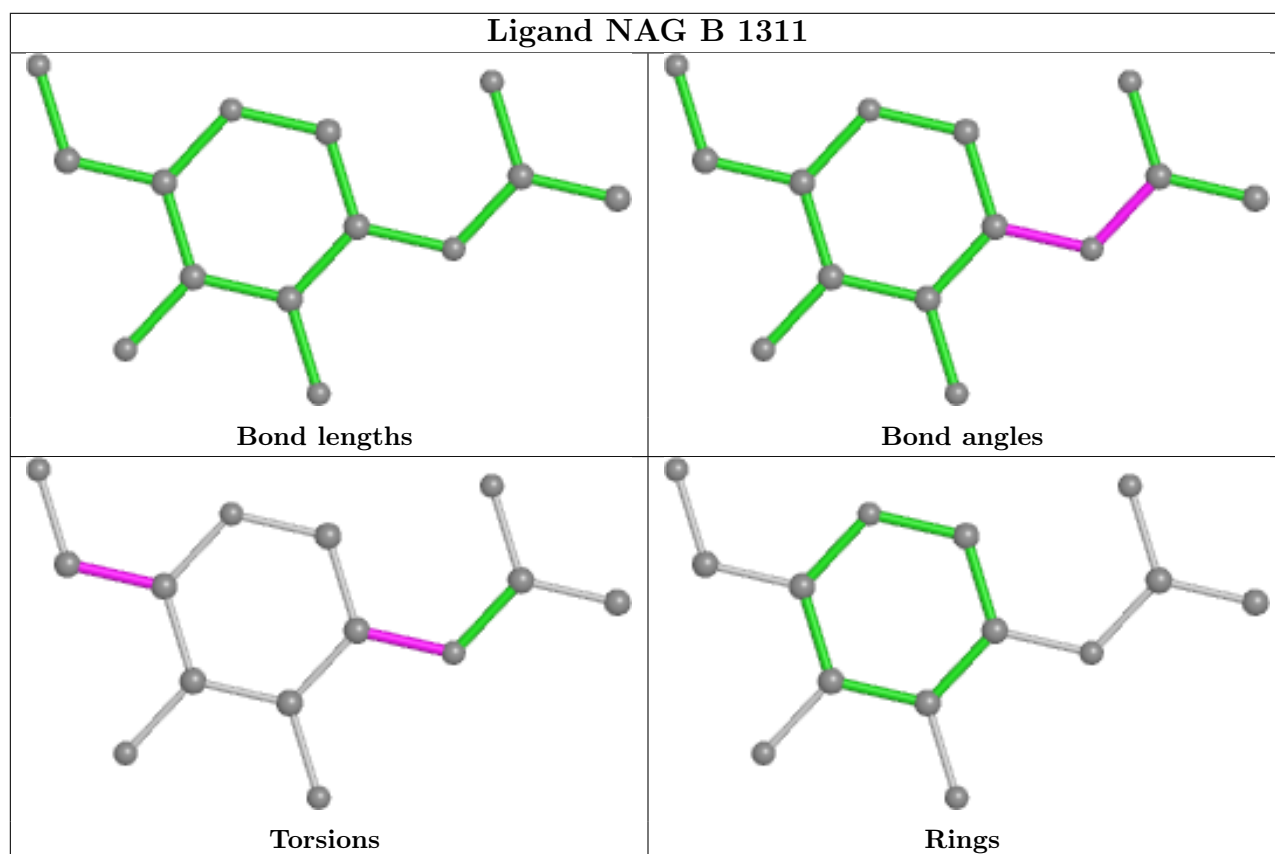
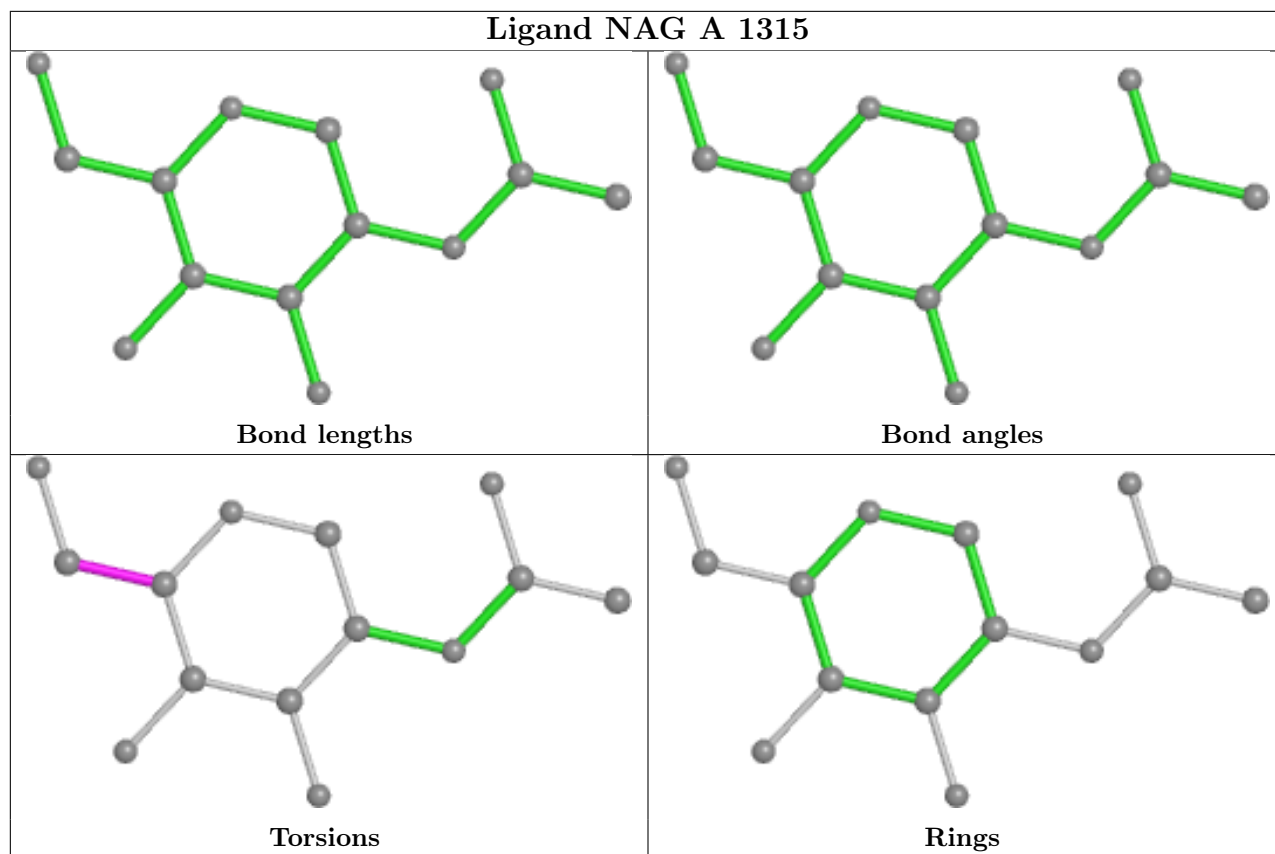


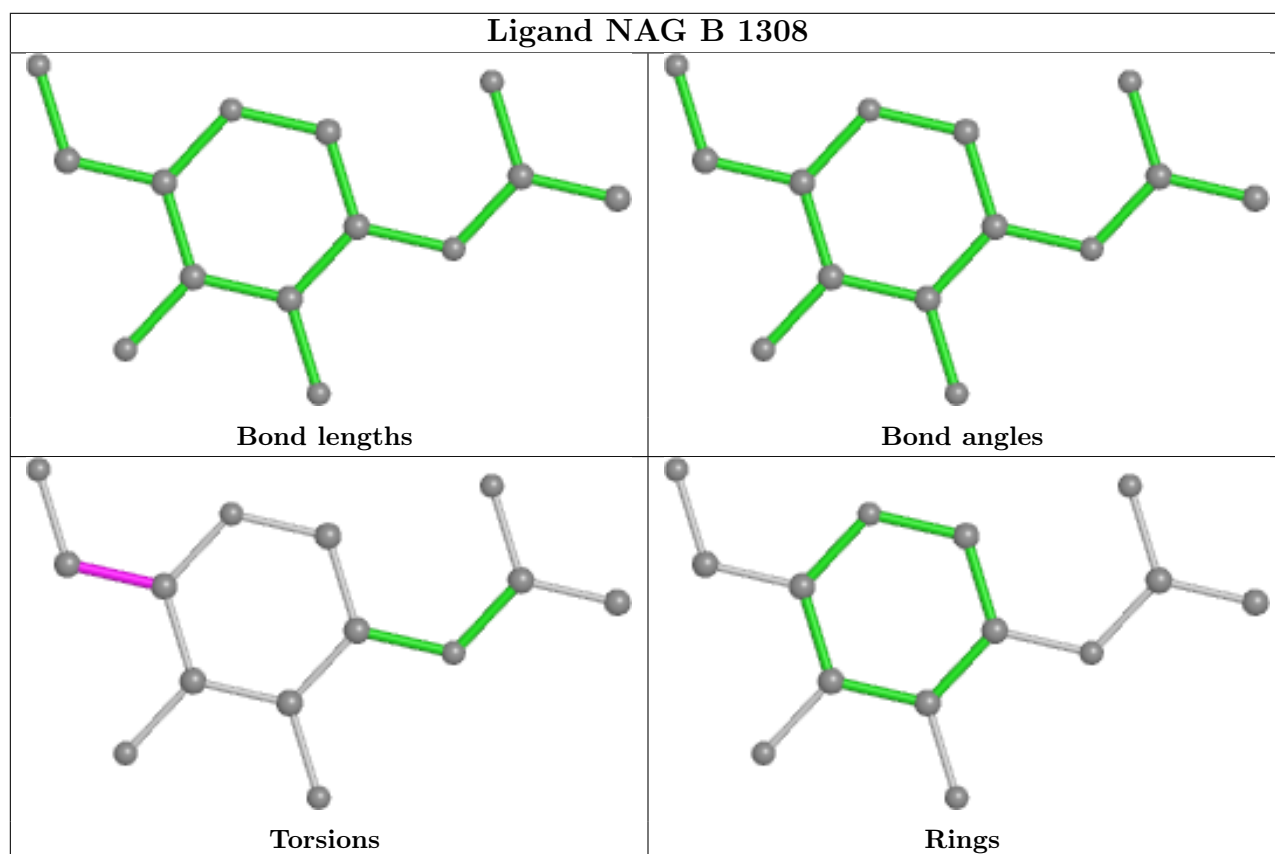
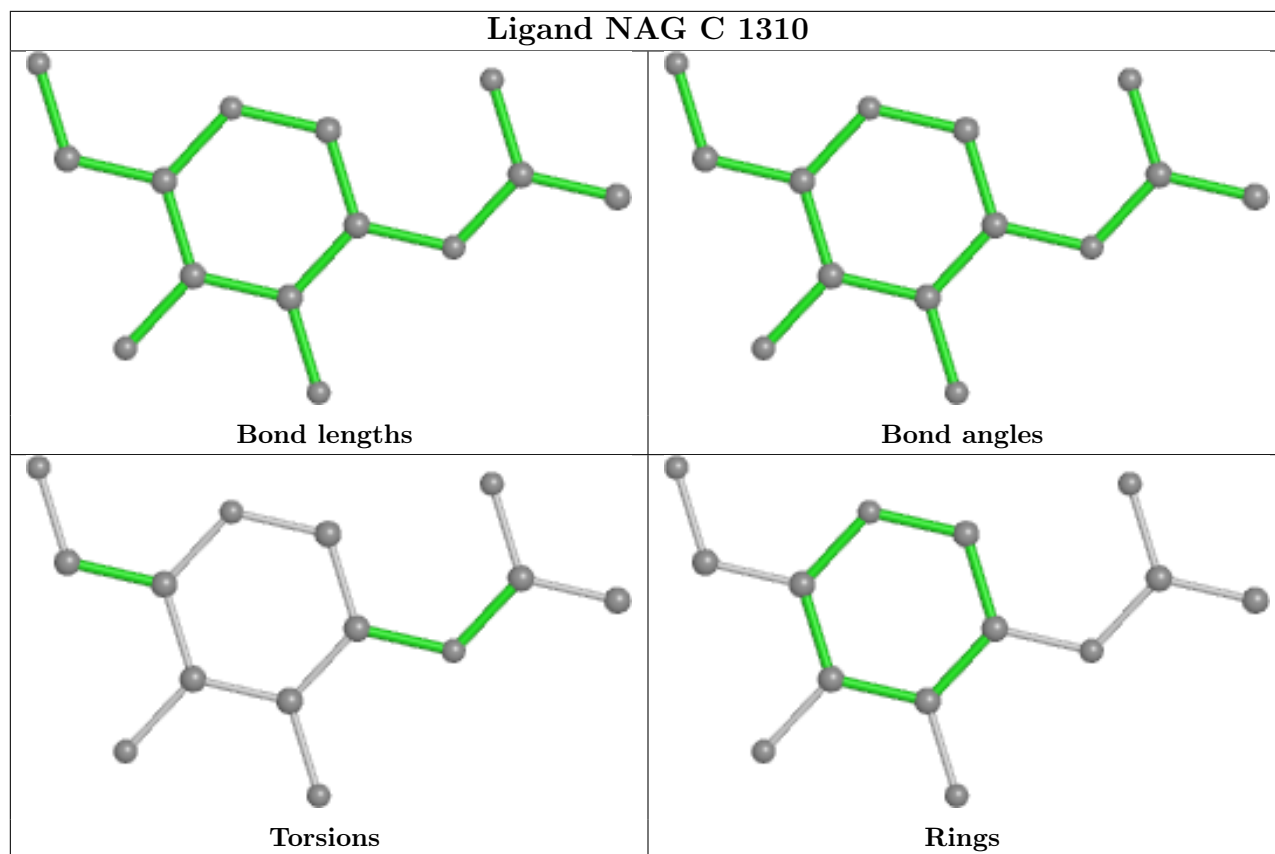


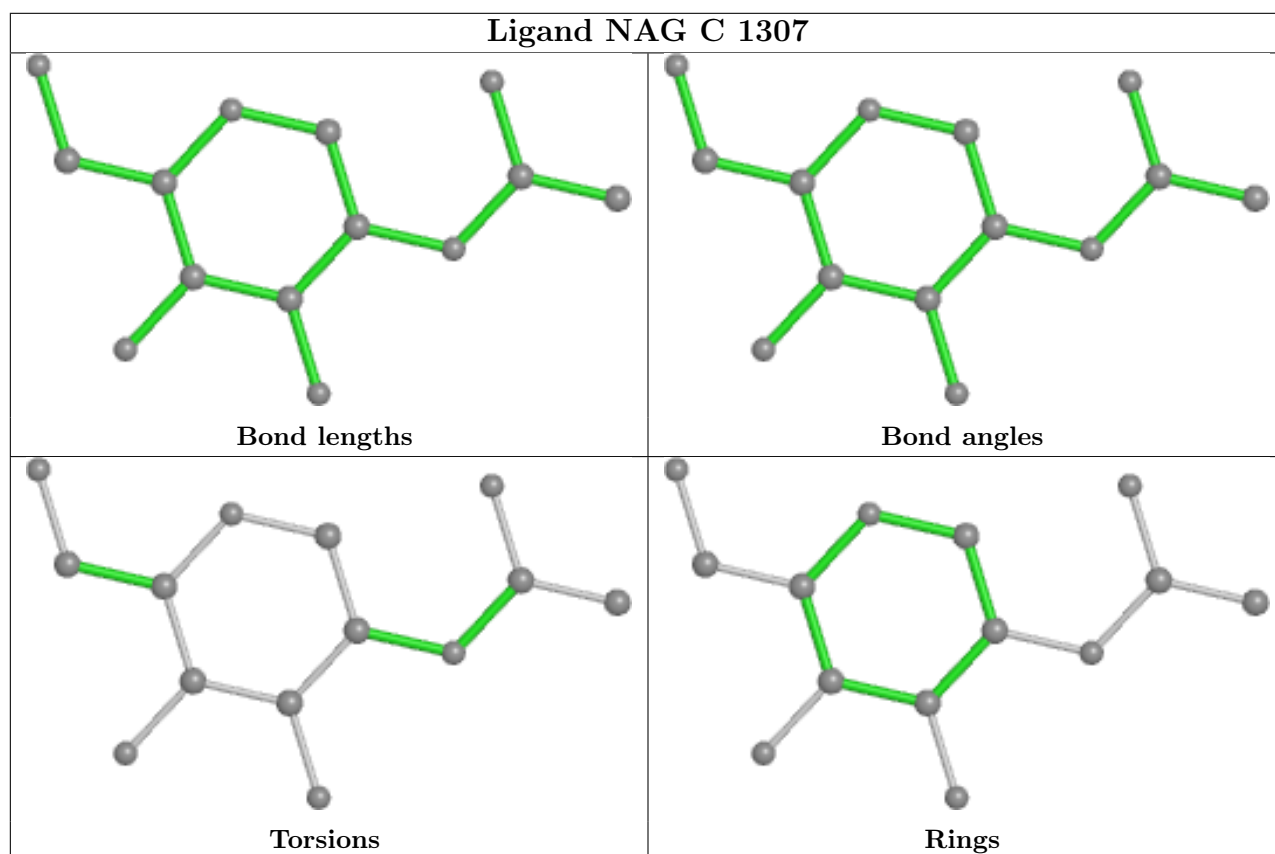
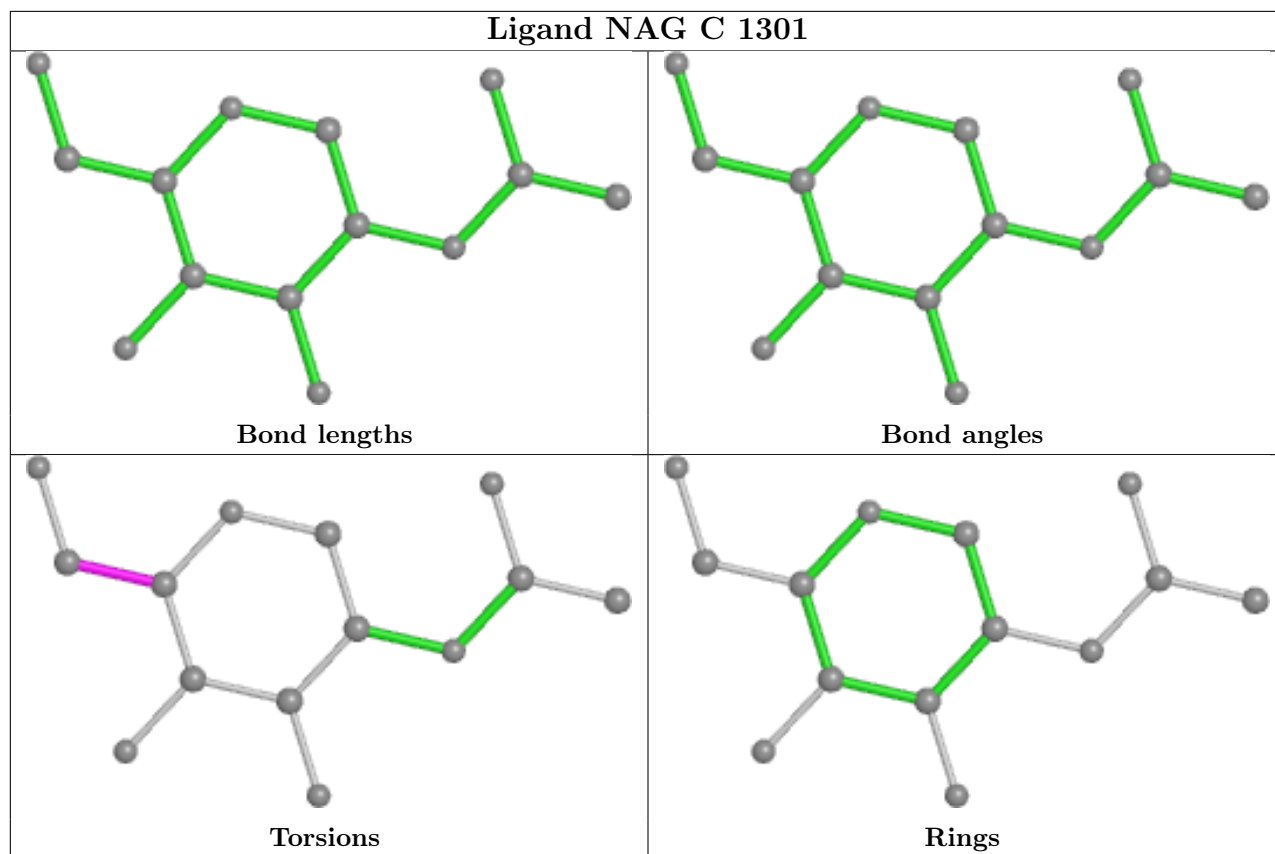


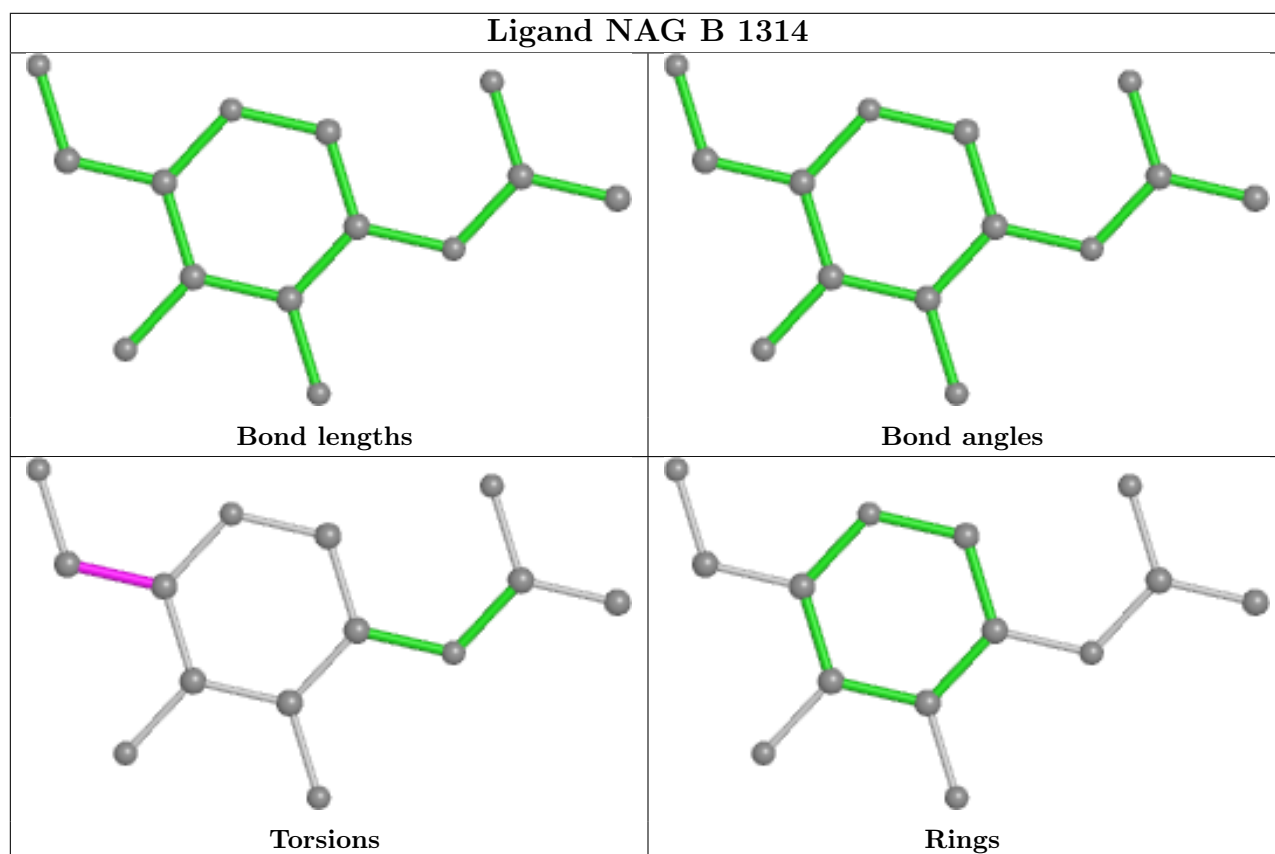
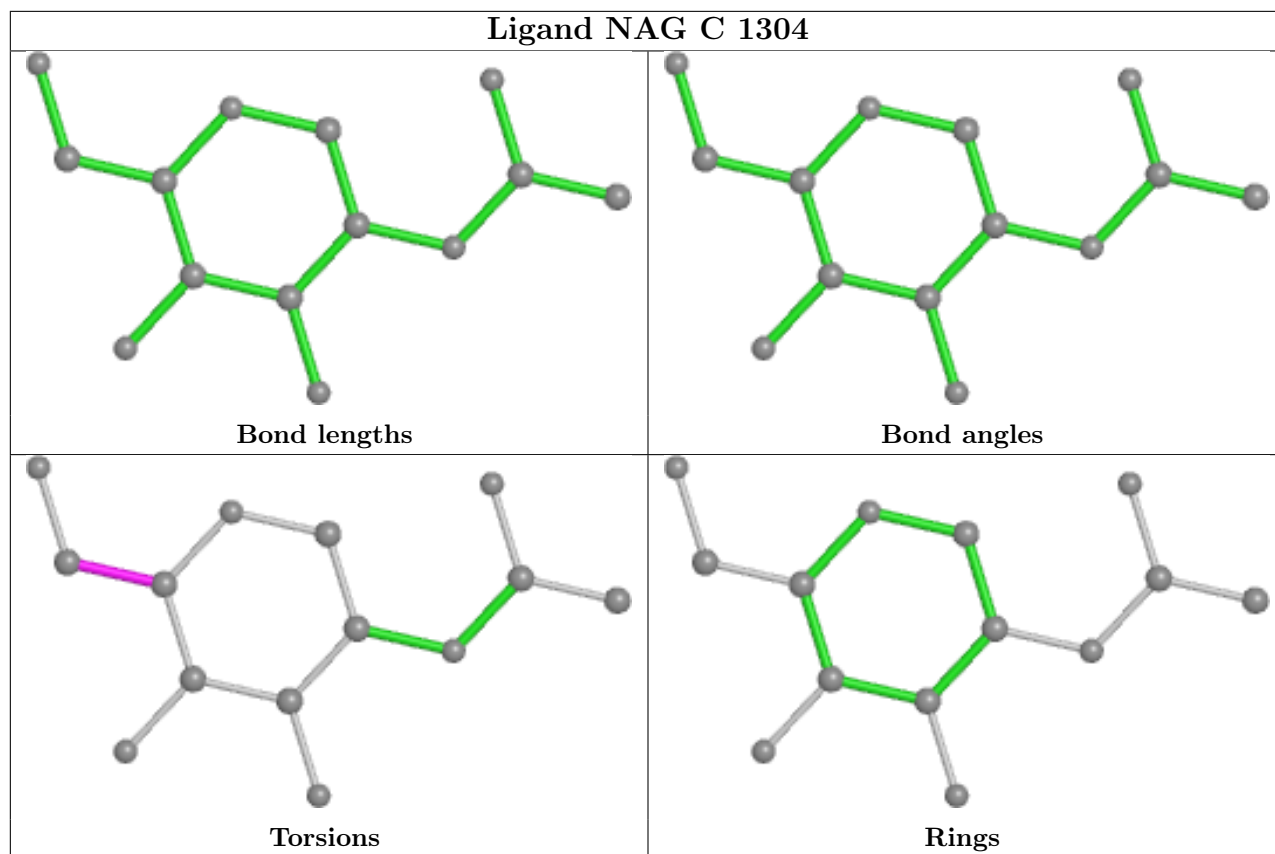


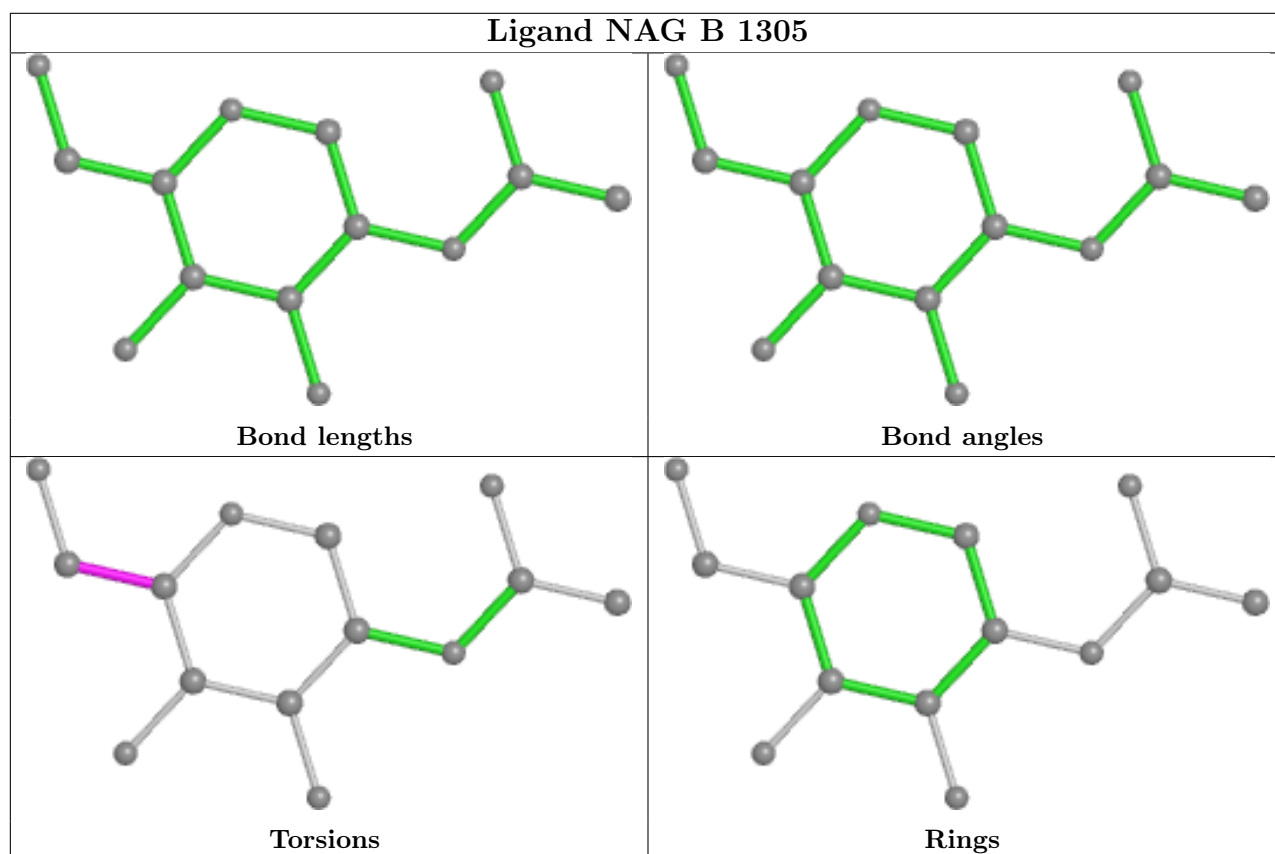
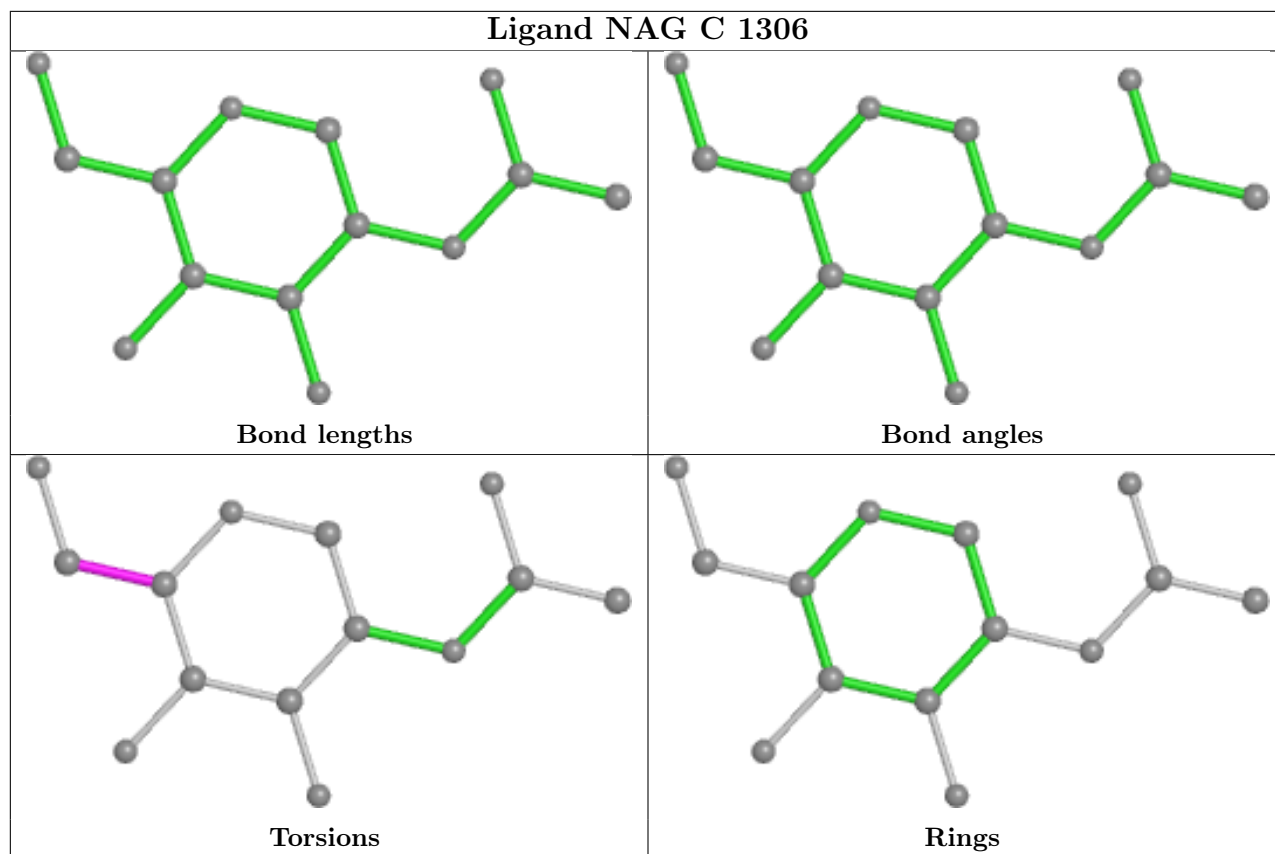


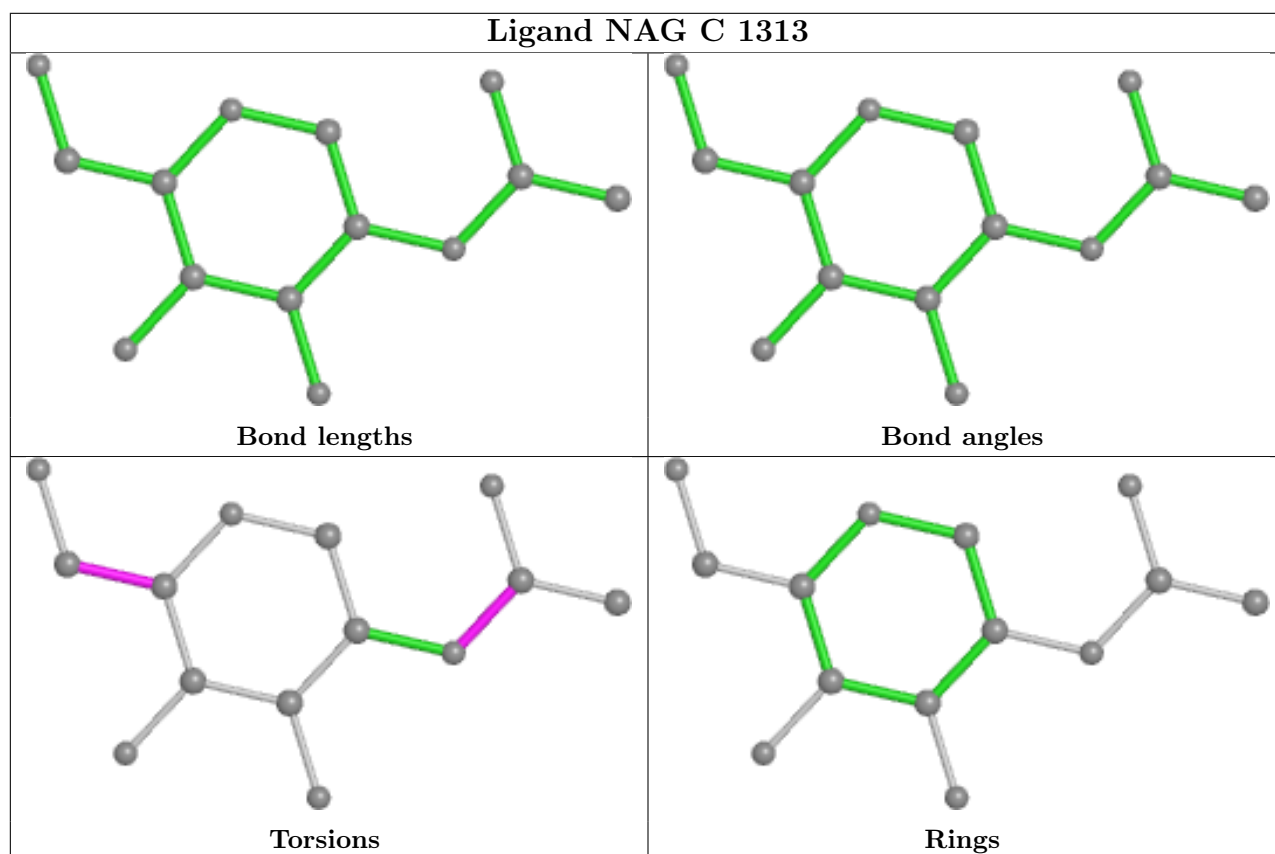
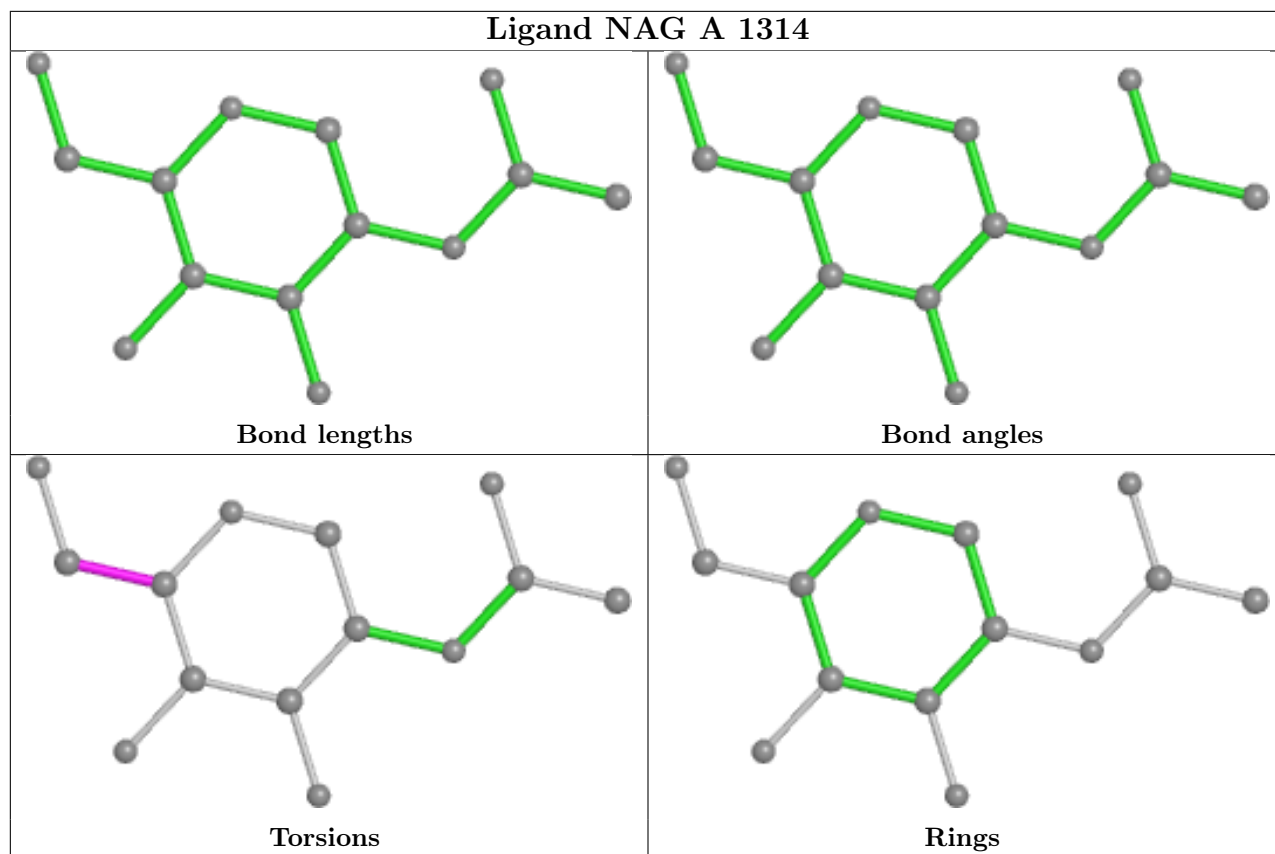


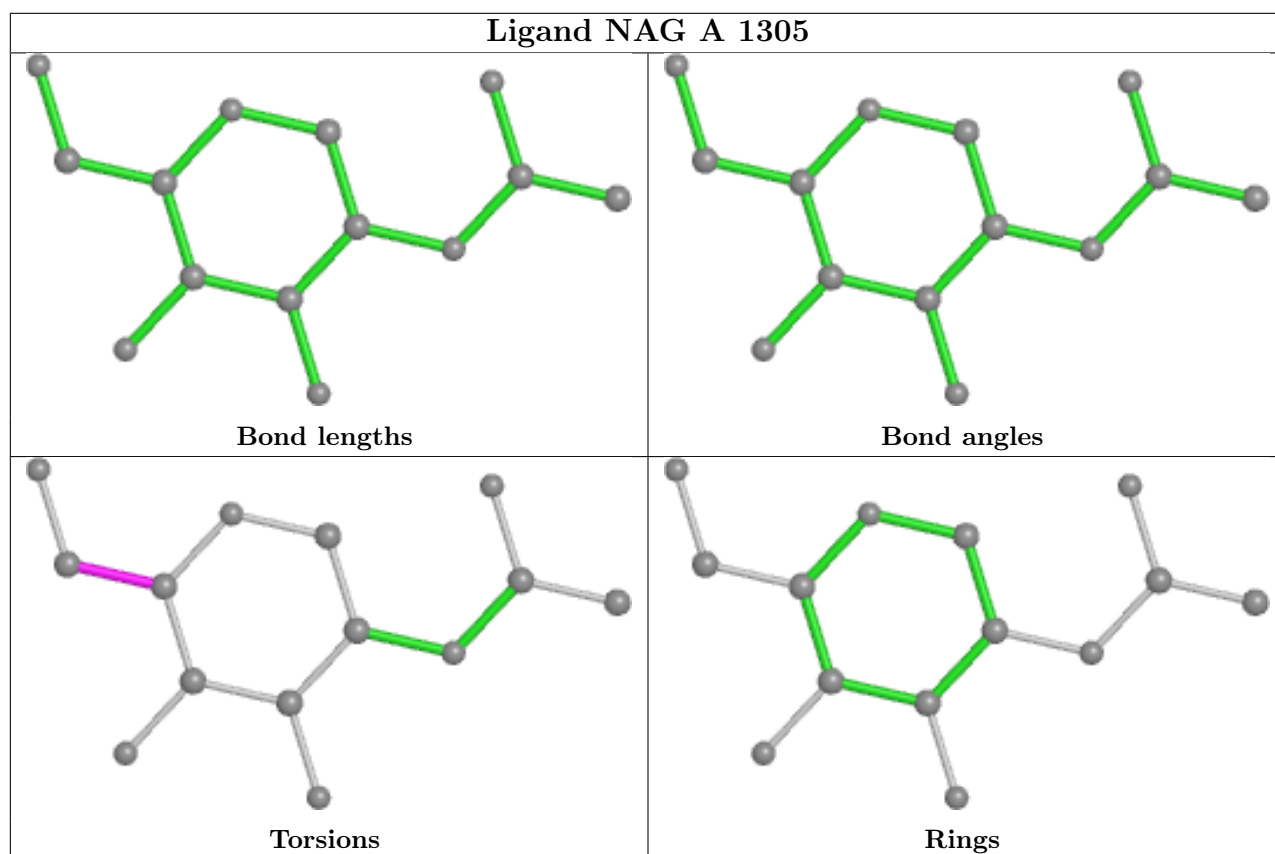
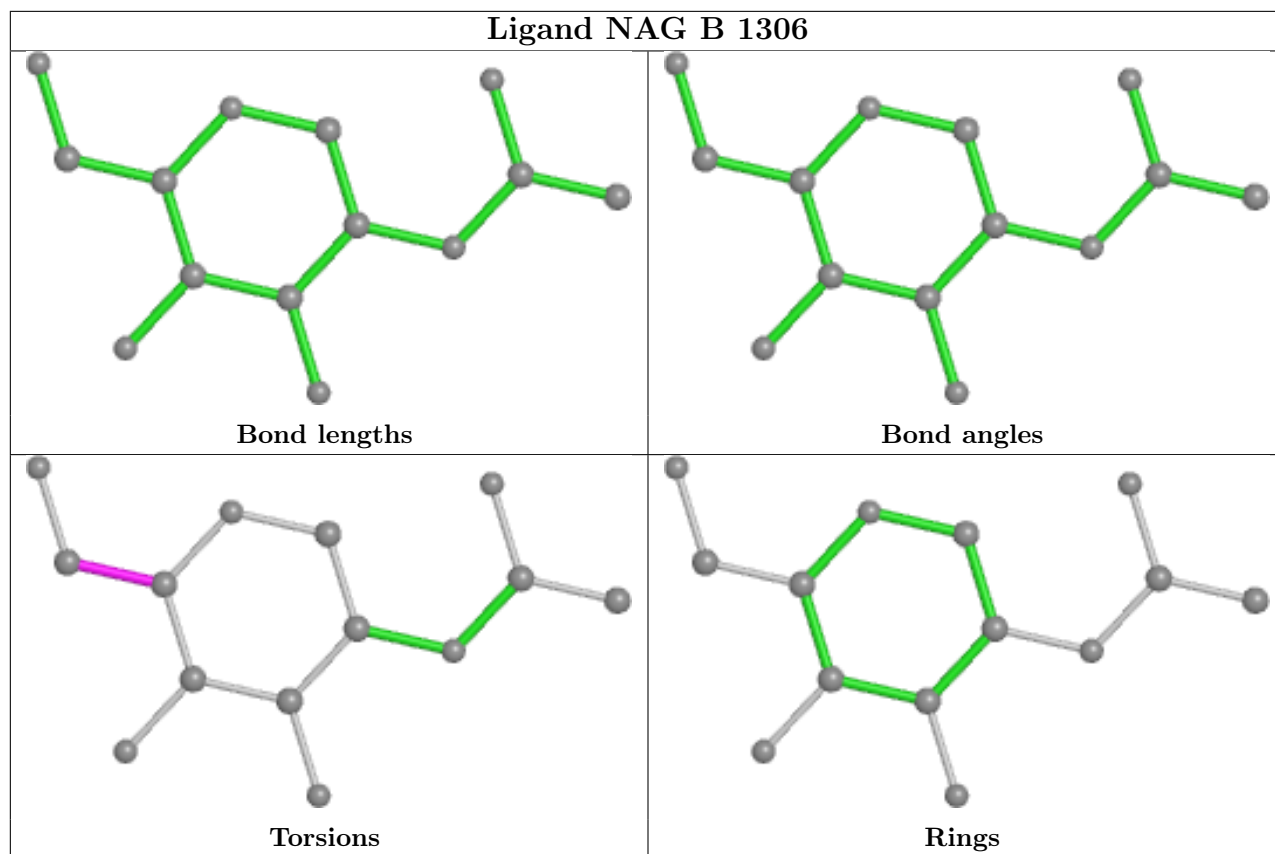


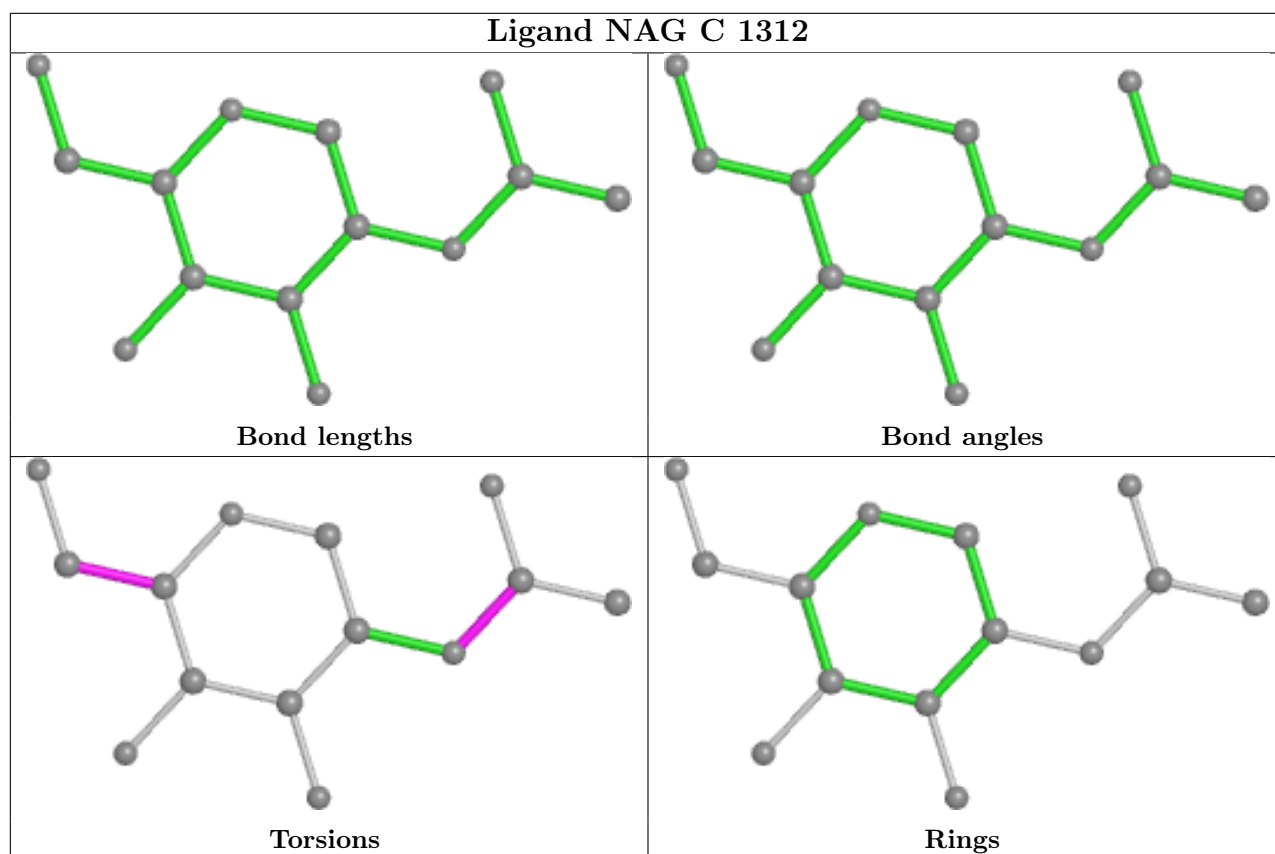
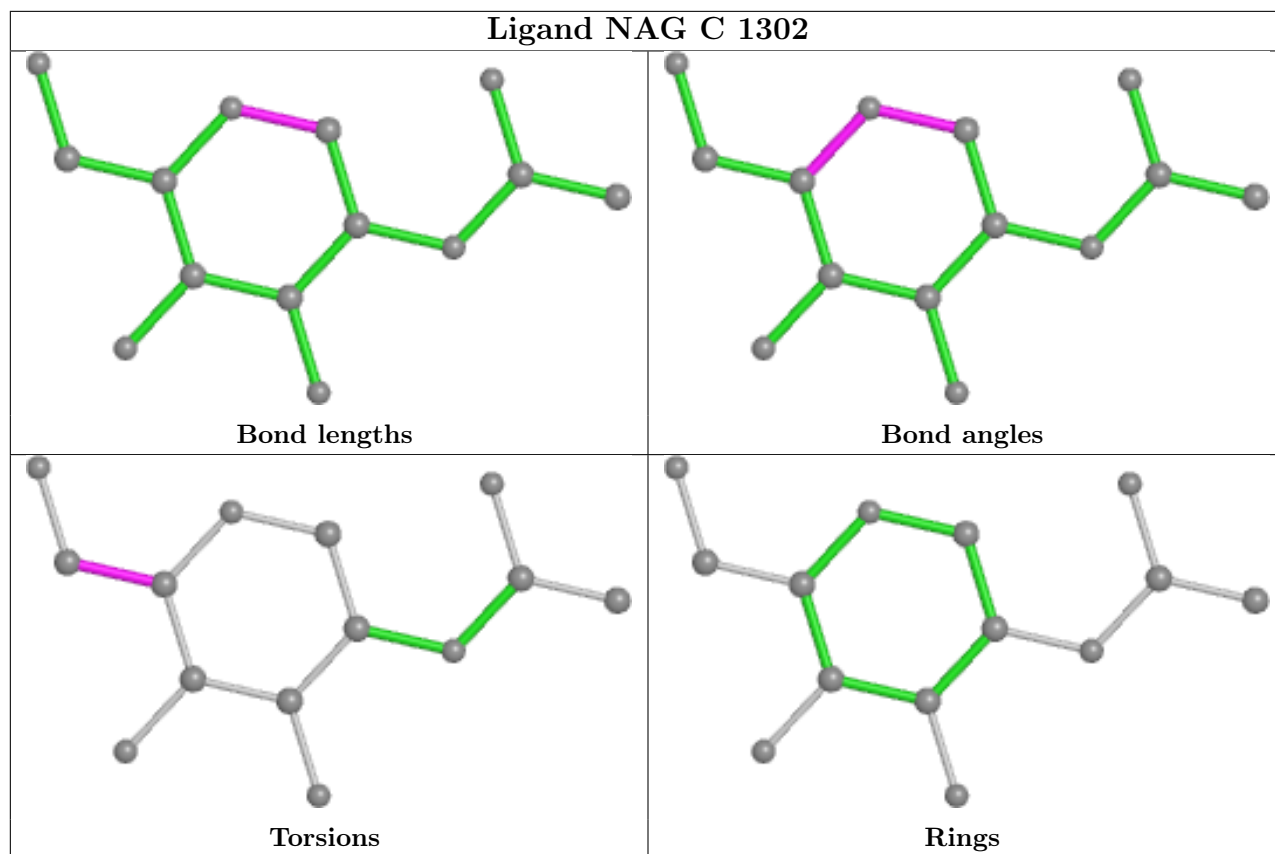


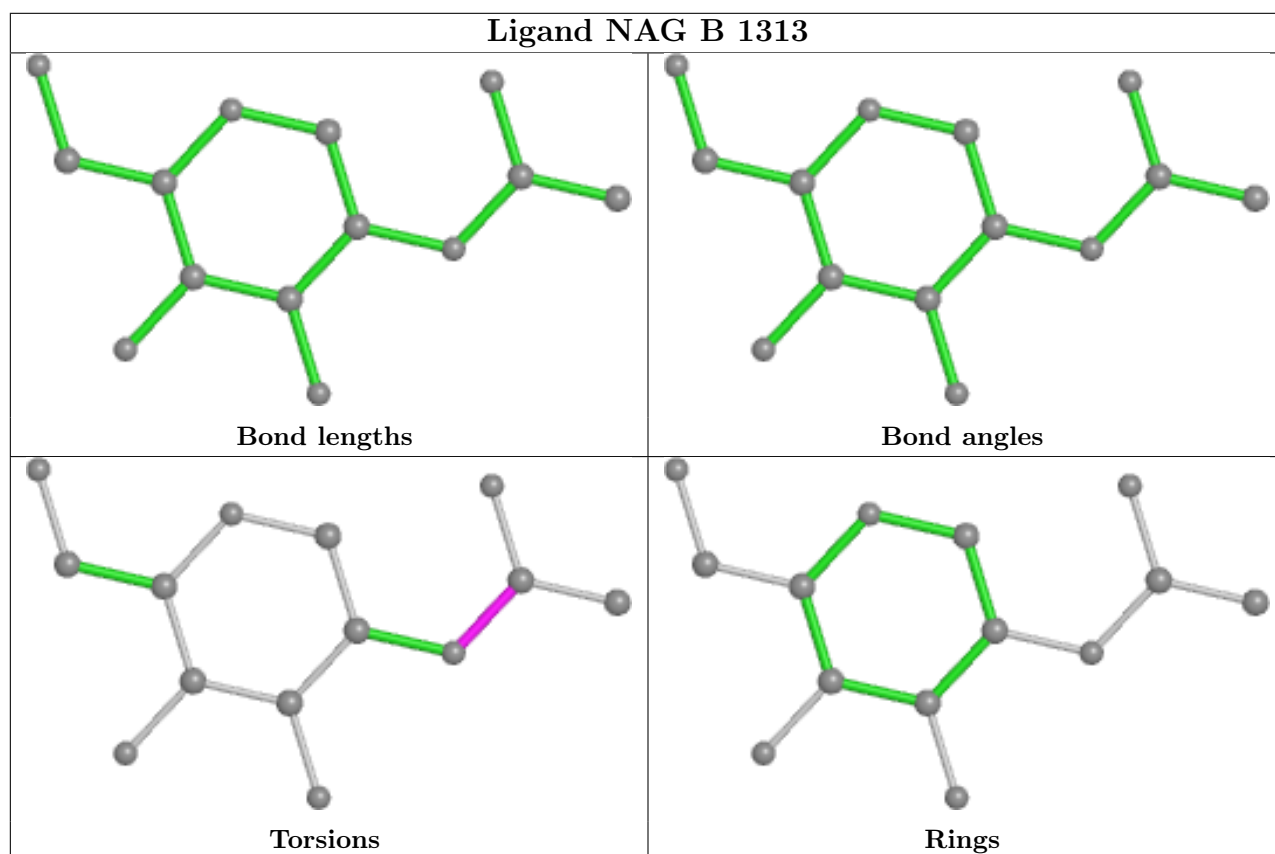
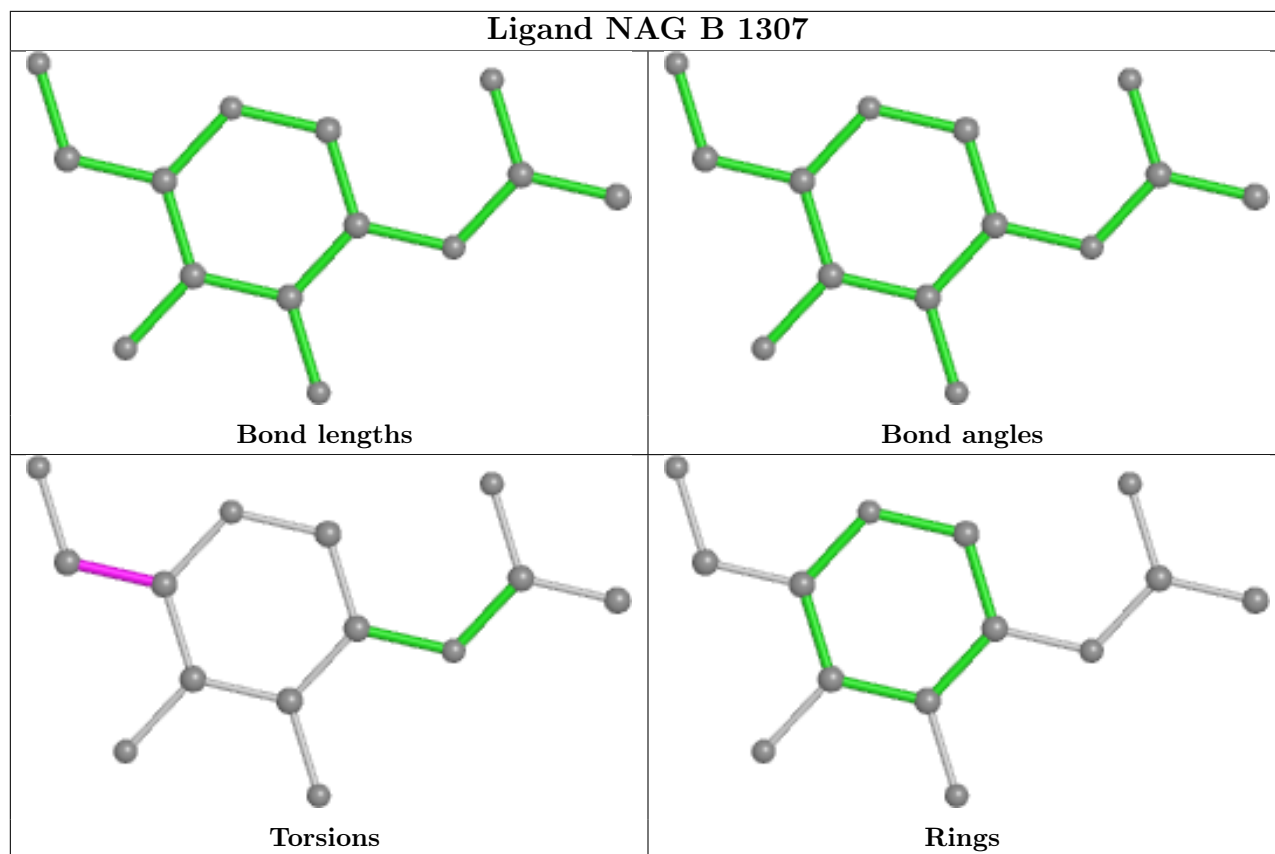


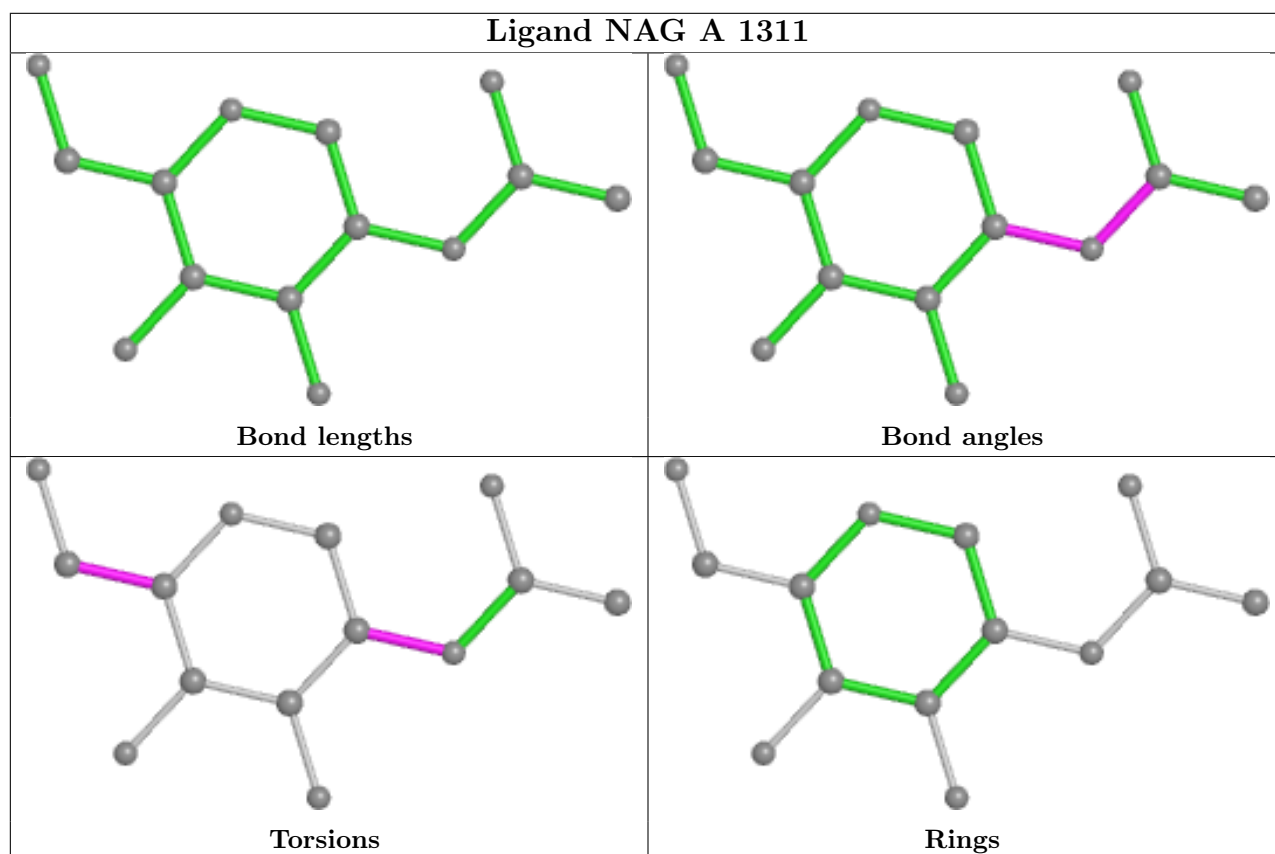
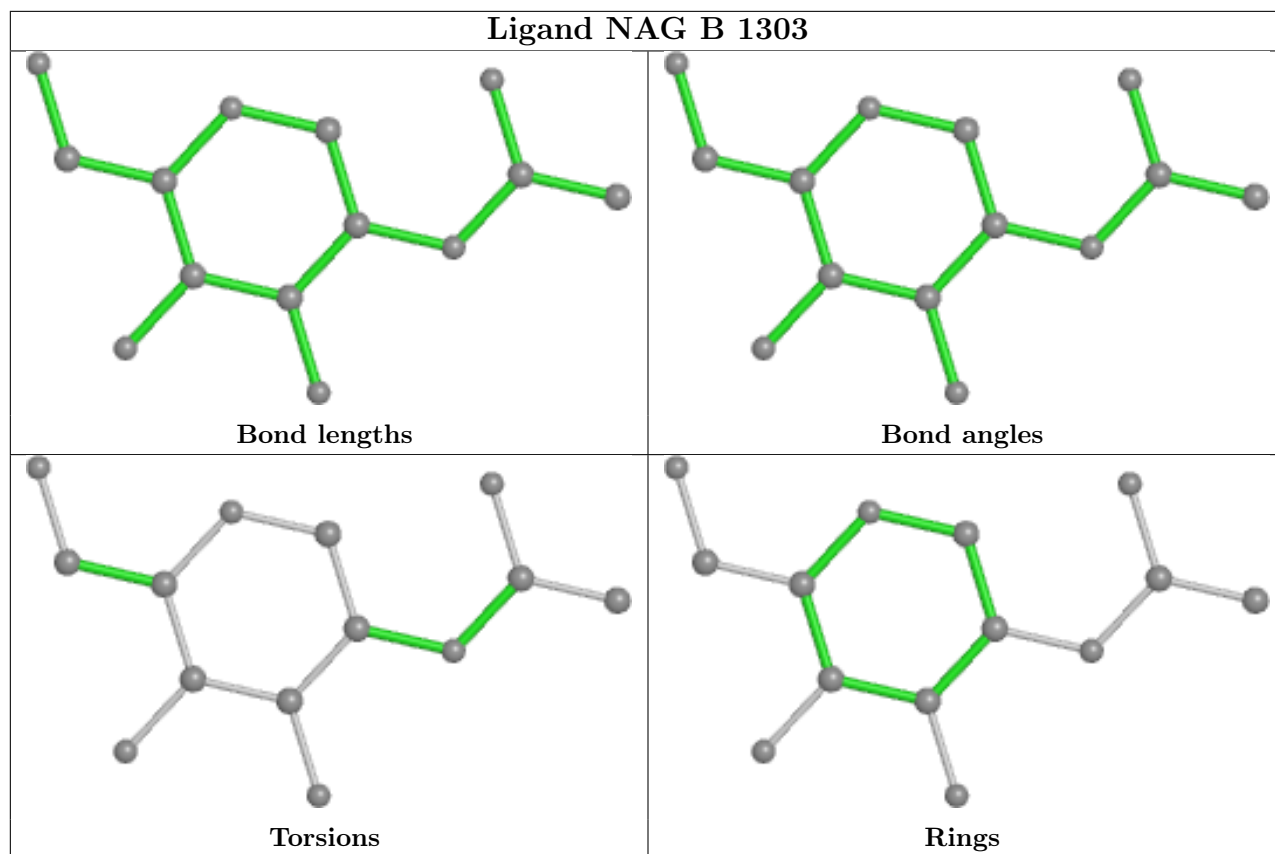


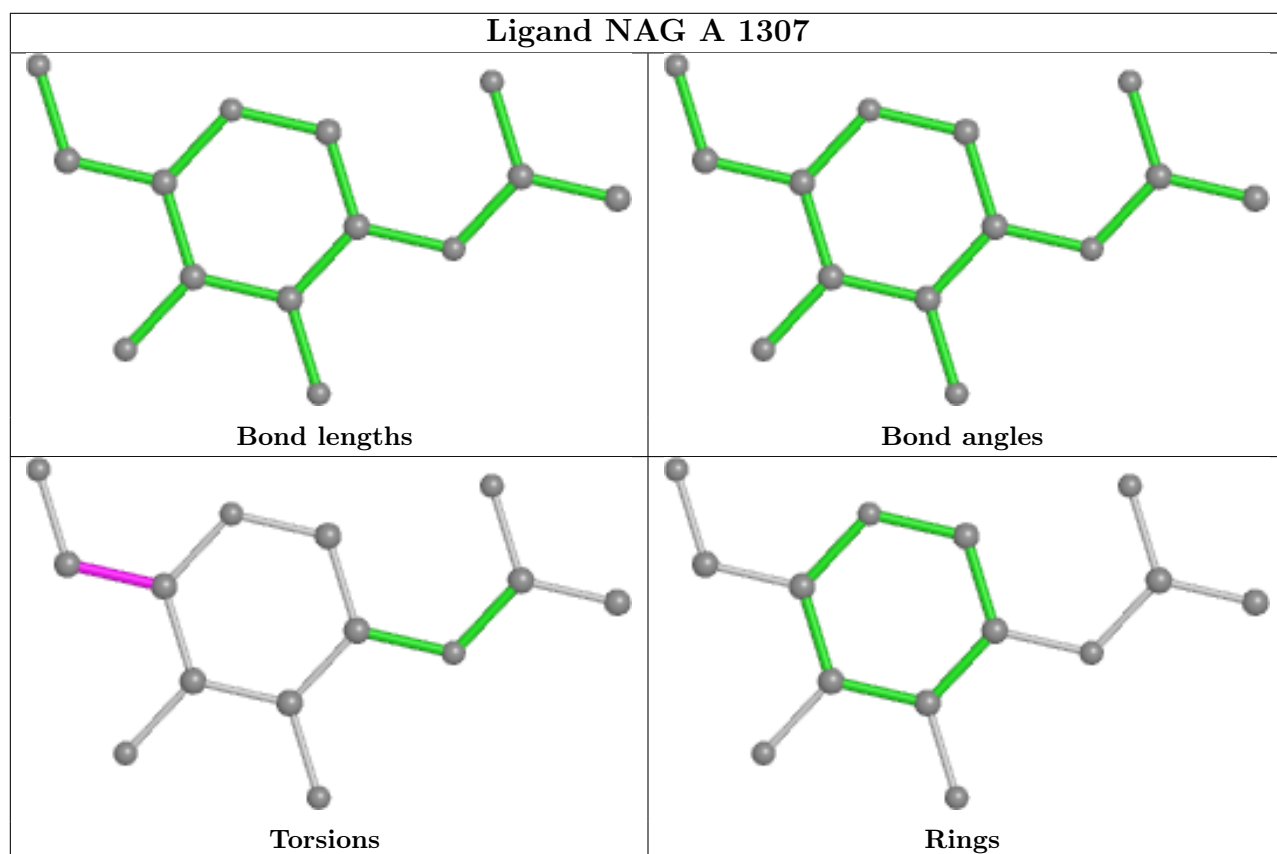
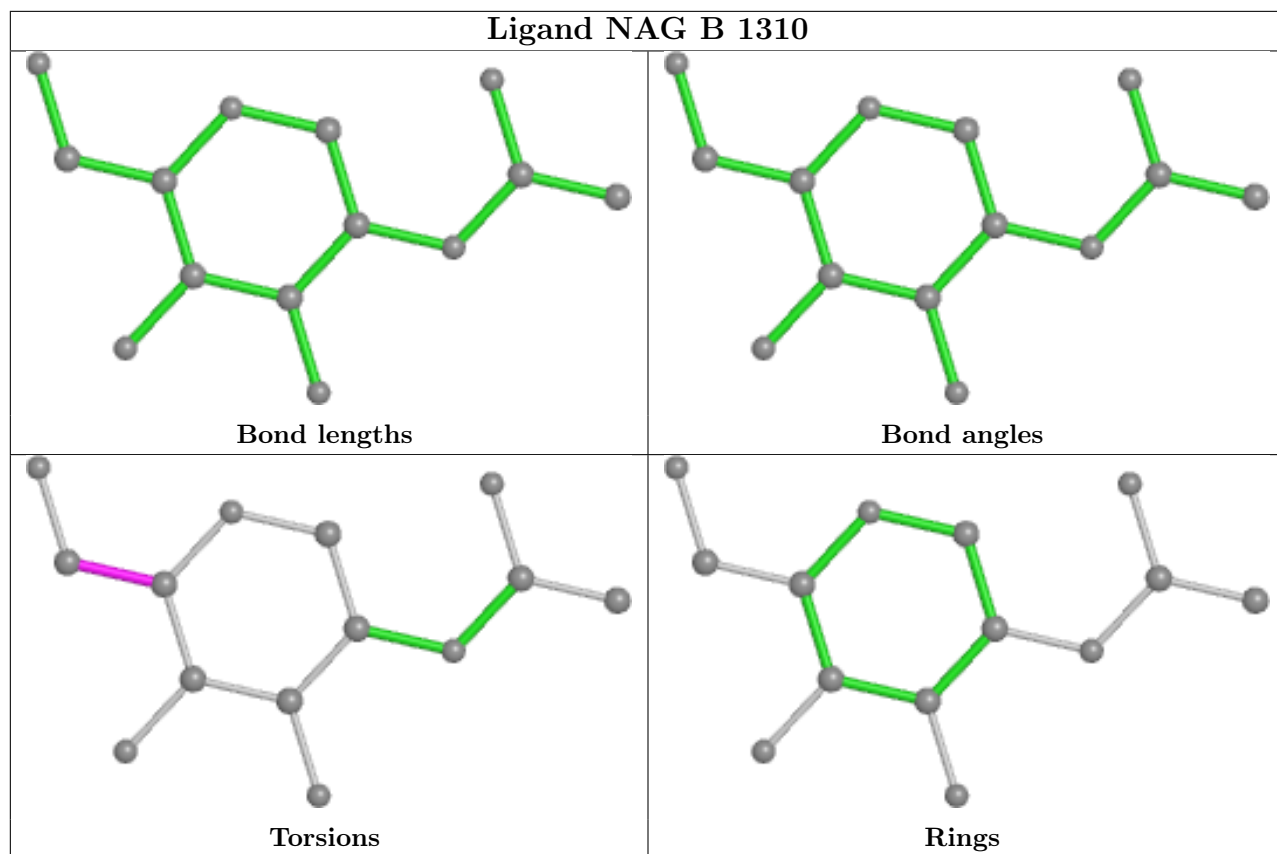


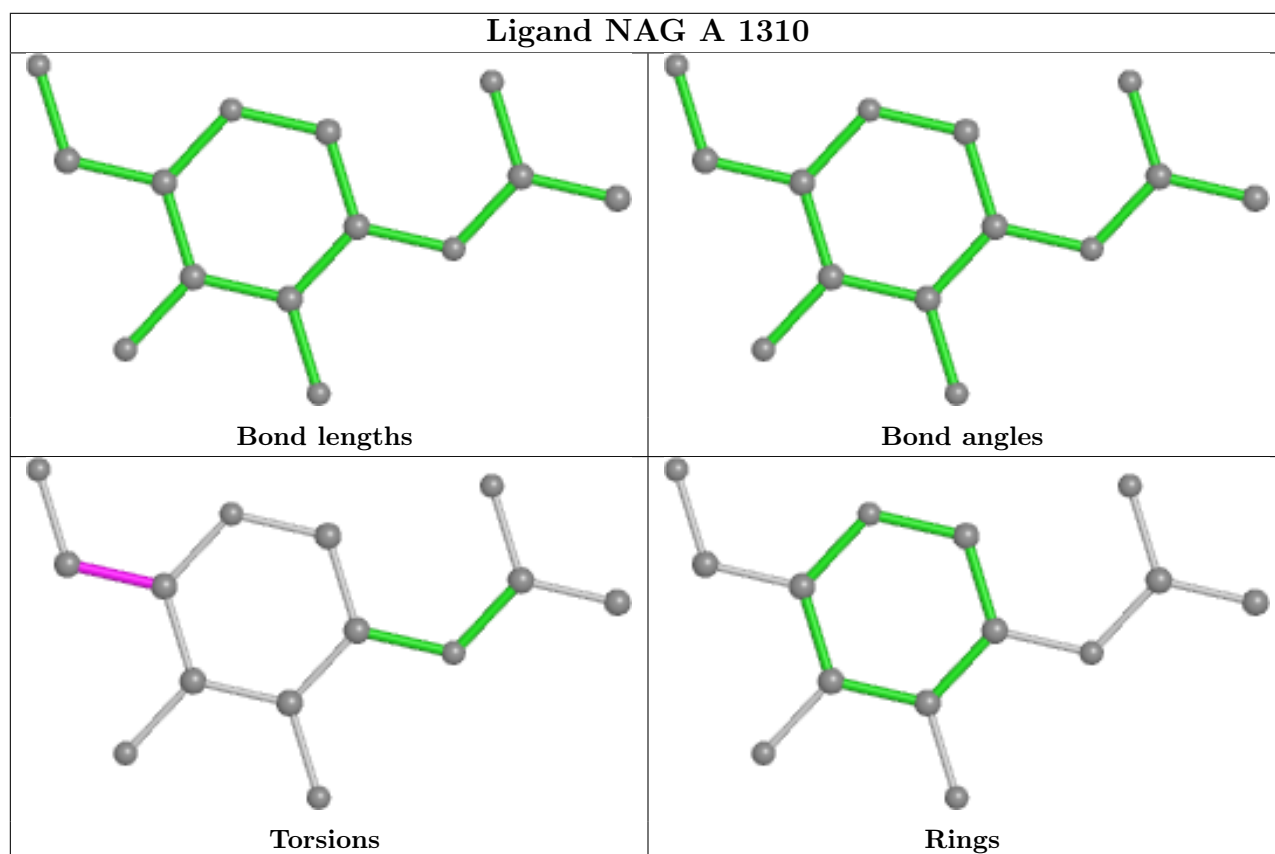
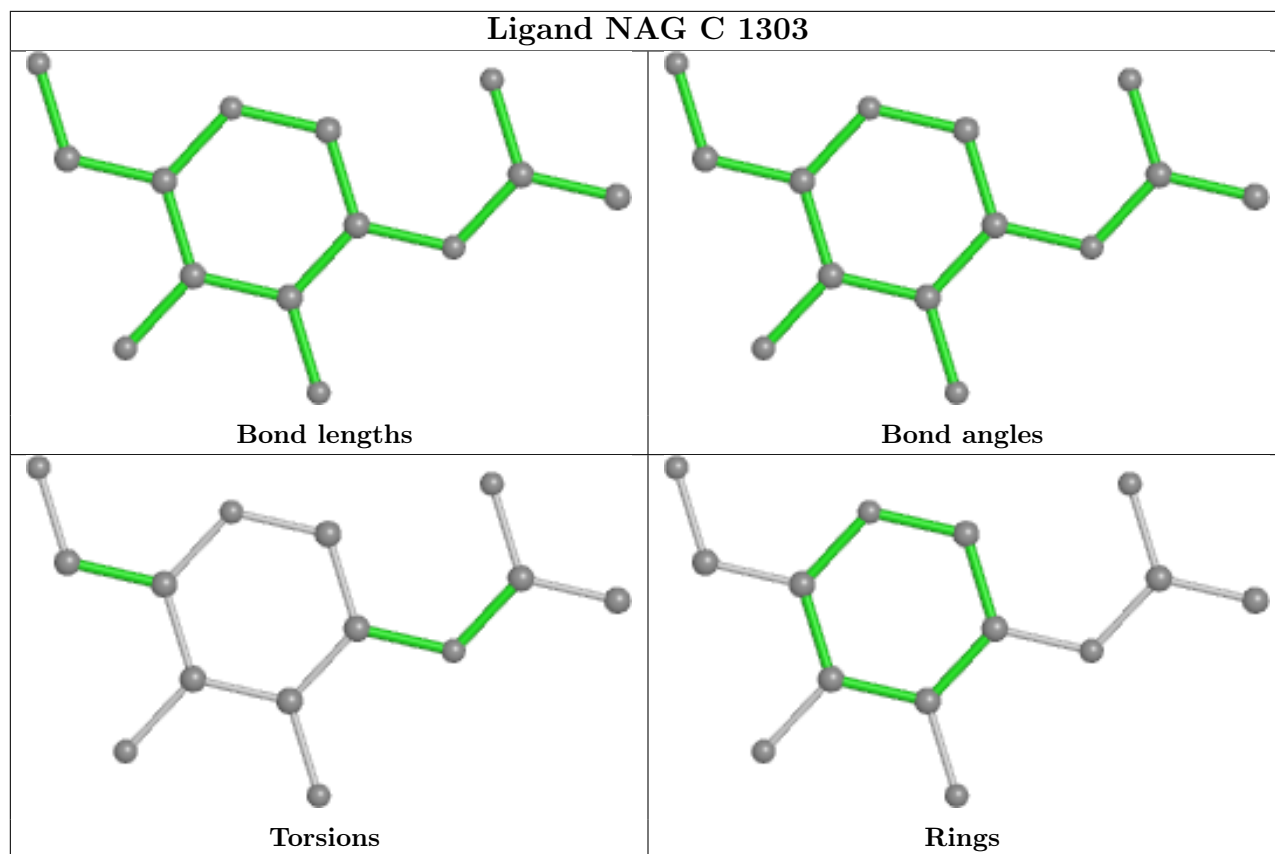


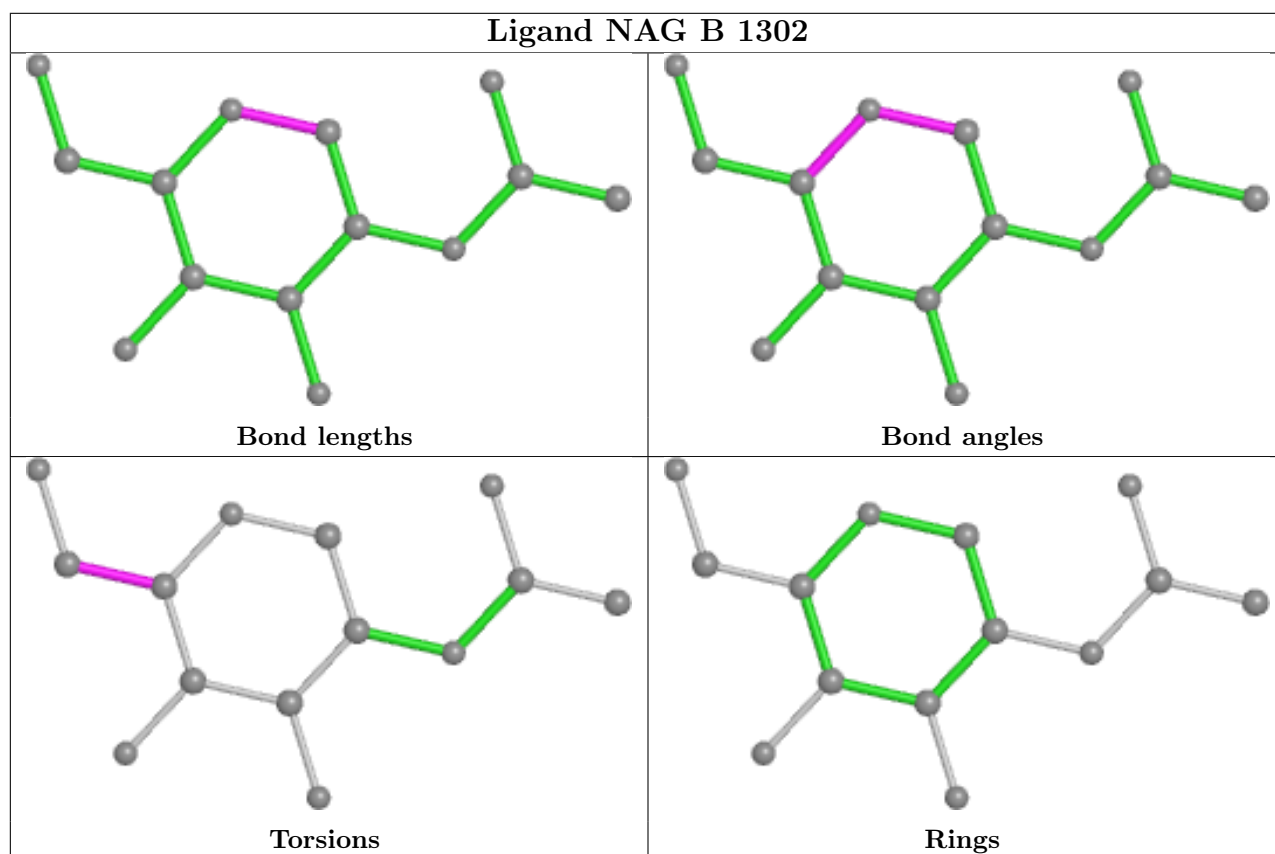
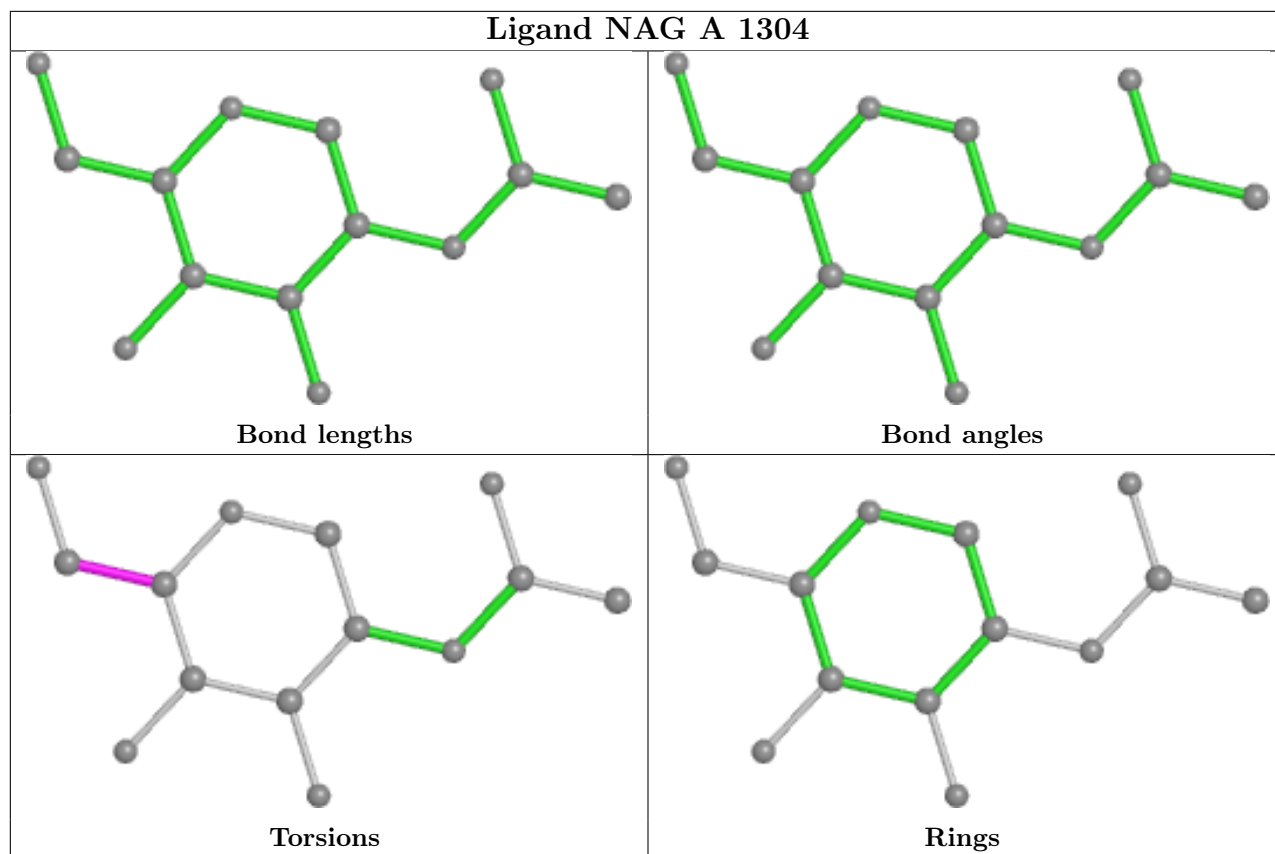


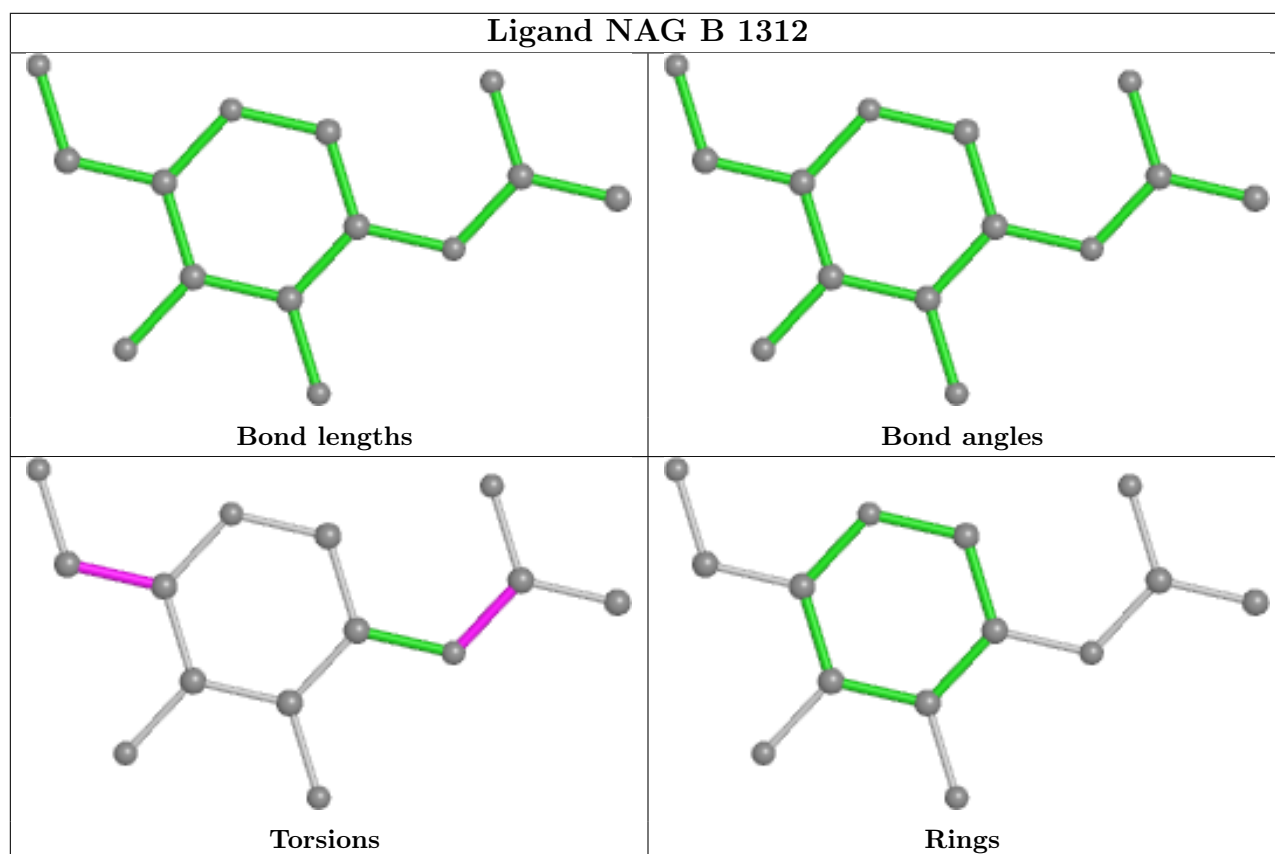
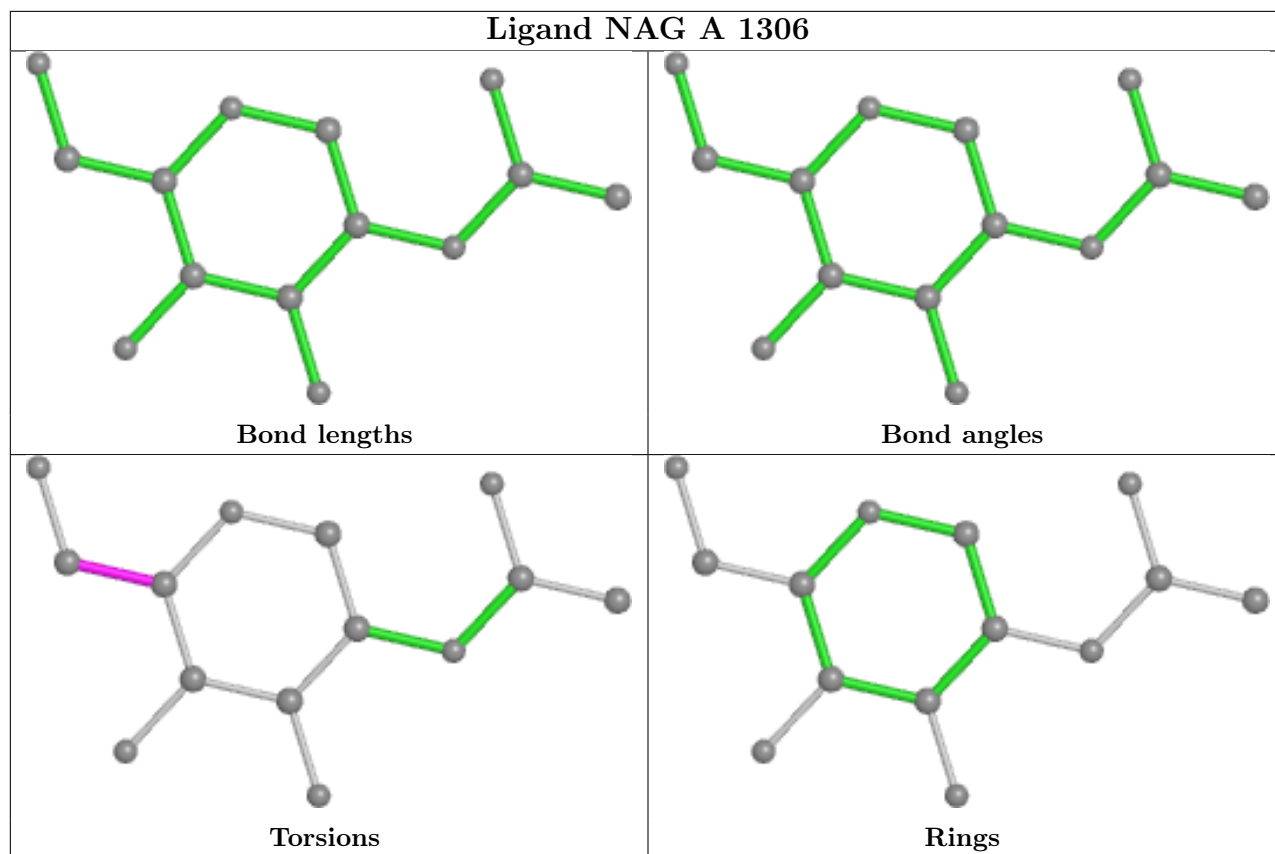


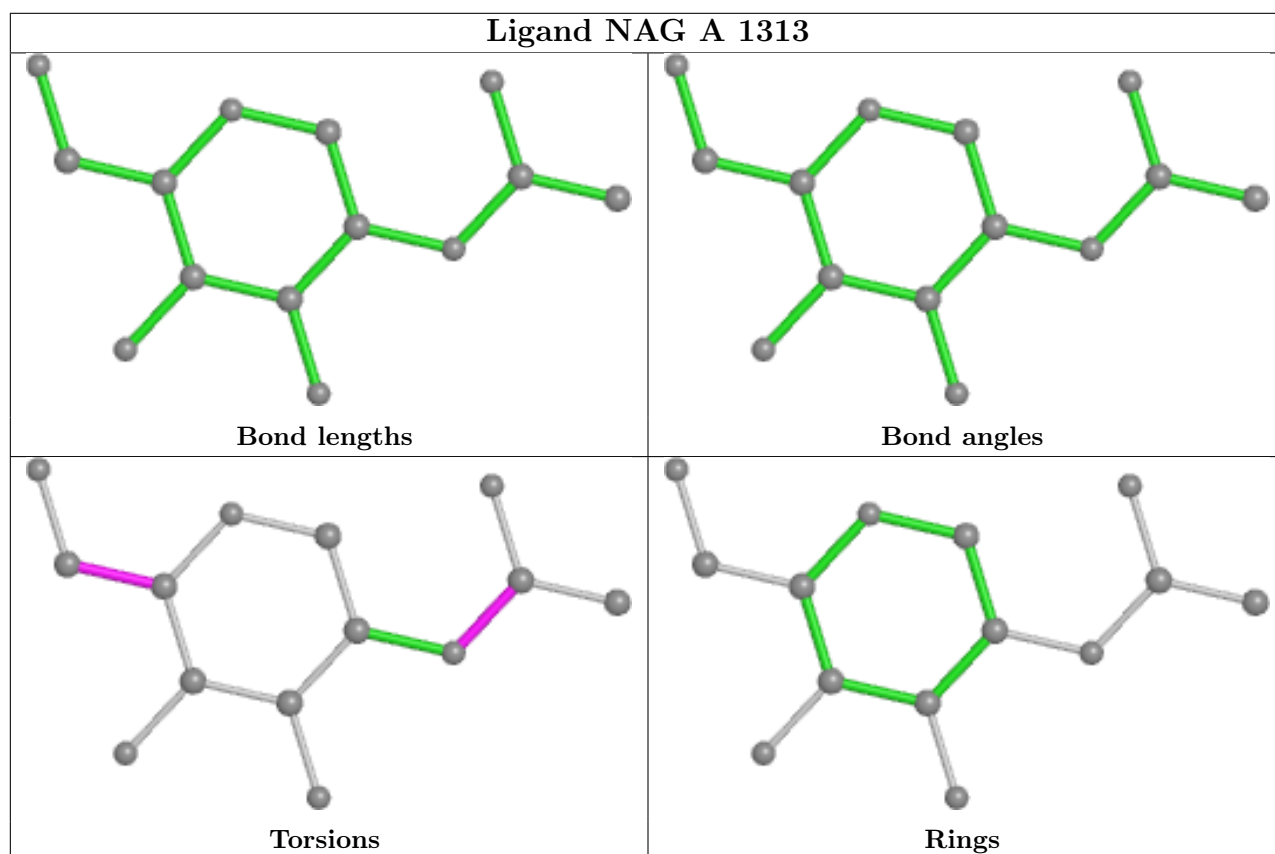
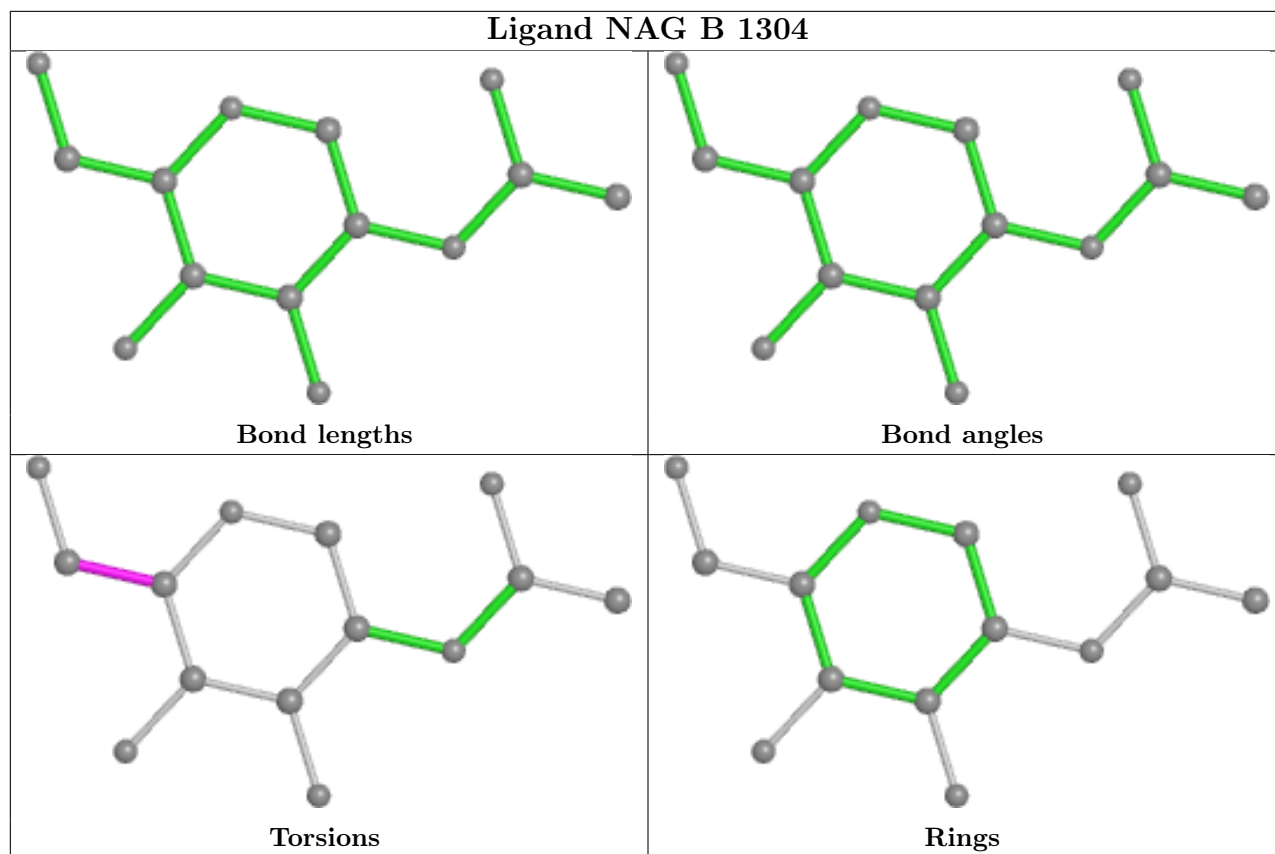


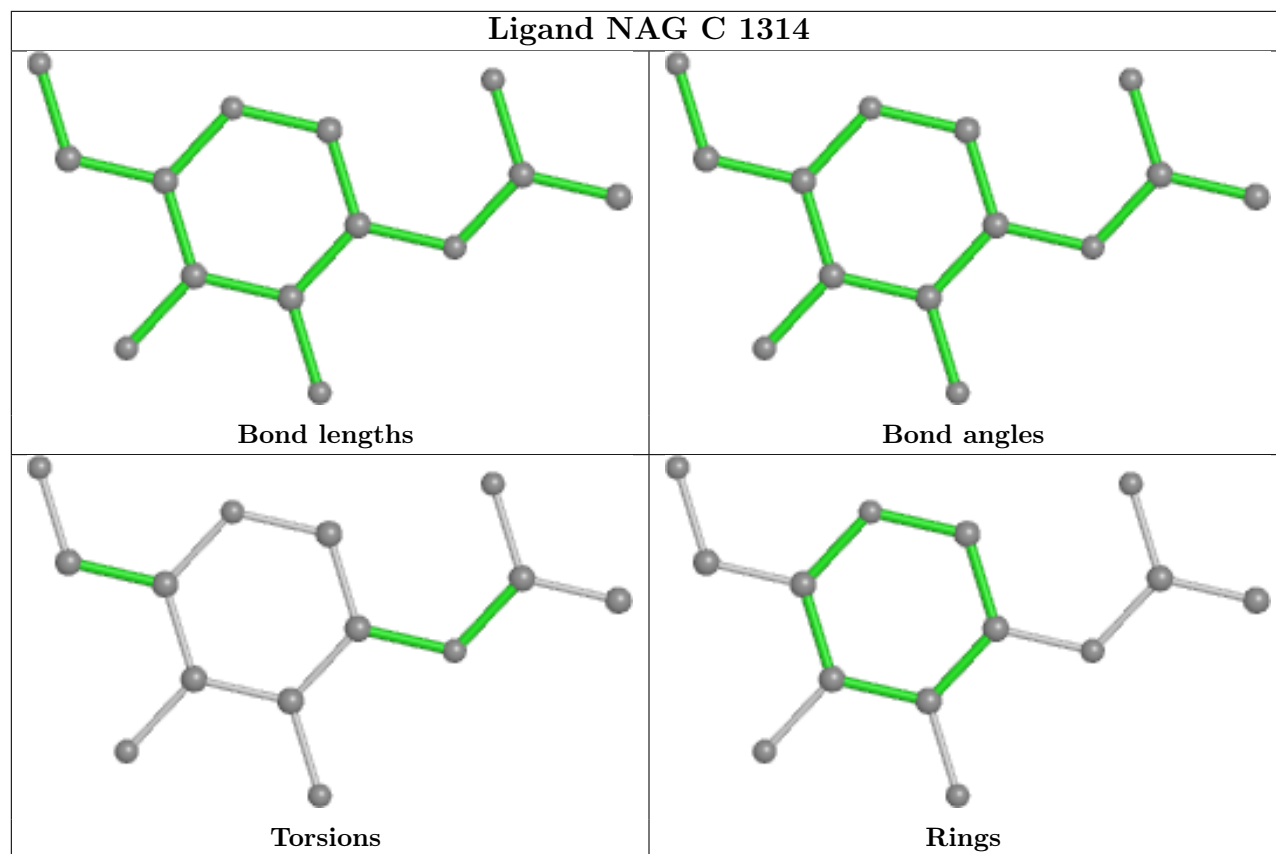












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.