



Full wwPDB EM Validation Report ⓘ

Mar 29, 2026 – 11:29 PM UTC

PDB ID : 8ZY0 / pdb_00008zy0
EMDB ID : EMD-60551
Title : Sarbecovirus BtKY72 Spike Trimer in a Locked Conformation
Authors : Wang, J.; Xiong, X.
Deposited on : 2024-06-16
Resolution : 3.00 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 : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

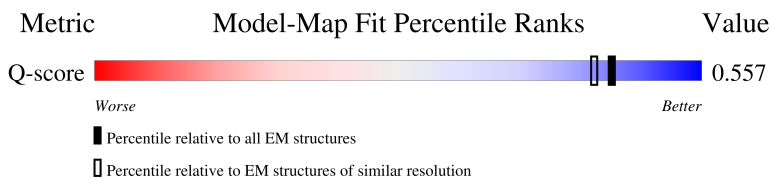
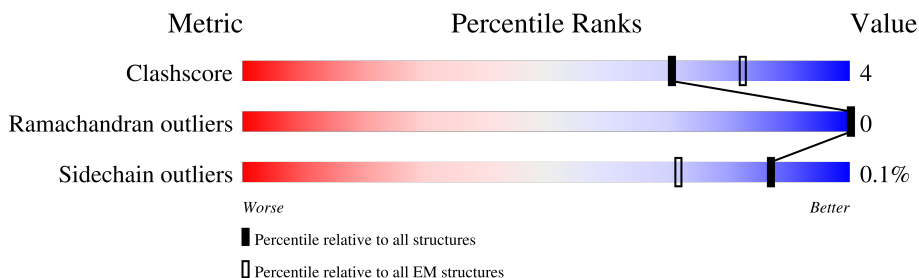
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.00 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14081 (2.50 - 3.50)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1193	
1	B	1193	
1	C	1193	
2	D	2	

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Mol	Chain	Length	Quality of chain
2	F	2	 100%
2	H	2	 100%
3	E	5	 40%40%60%
4	G	4	 75%25%
4	I	4	 75%25%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 26611 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.

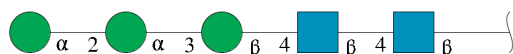
Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1092	Total	C	N	O	S	0	0
			8509	5419	1413	1633	44		
1	B	1092	Total	C	N	O	S	0	0
			8509	5419	1413	1633	44		
1	C	1092	Total	C	N	O	S	0	0
			8505	5416	1412	1633	44		

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



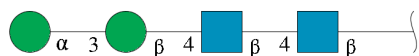
Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	2	Total	C	N	O	0	0
			28	16	2	10		
2	F	2	Total	C	N	O	0	0
			28	16	2	10		
2	H	2	Total	C	N	O	0	0
			28	16	2	10		

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



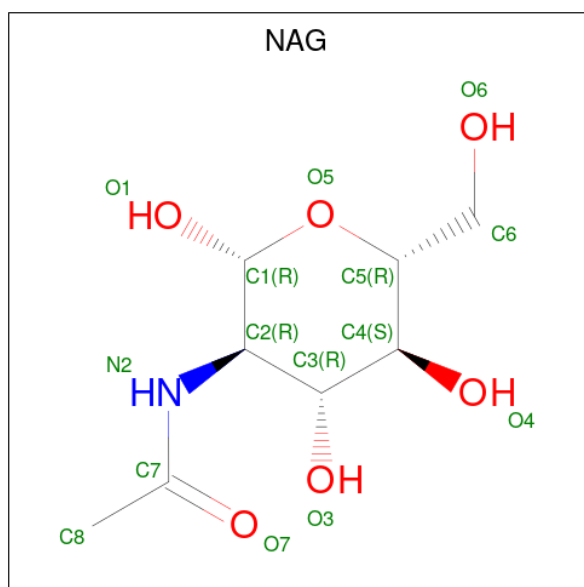
Mol	Chain	Residues	Atoms				AltConf	Trace
3	E	5	Total	C	N	O	0	0
			61	34	2	25		

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



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

- 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				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	

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

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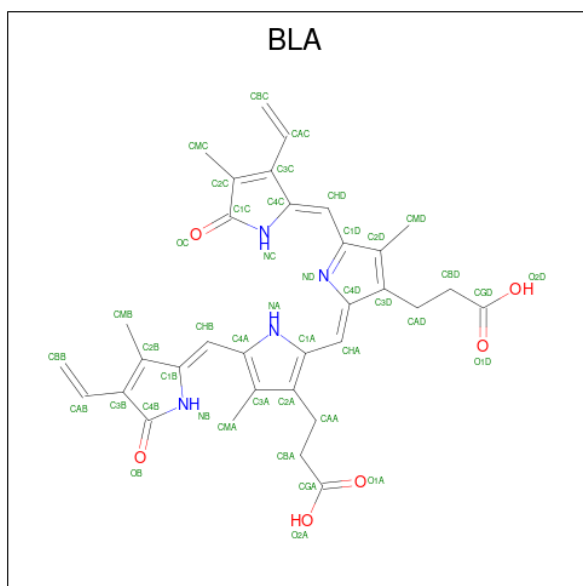
Mol	Chain	Residues	Atoms				AltConf
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5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

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

- Molecule 6 is BILIVERDINE IX ALPHA (CCD ID: BLA) (formula: $C_{33}H_{34}N_4O_6$) (labeled as "Ligand of Interest" by depositor).

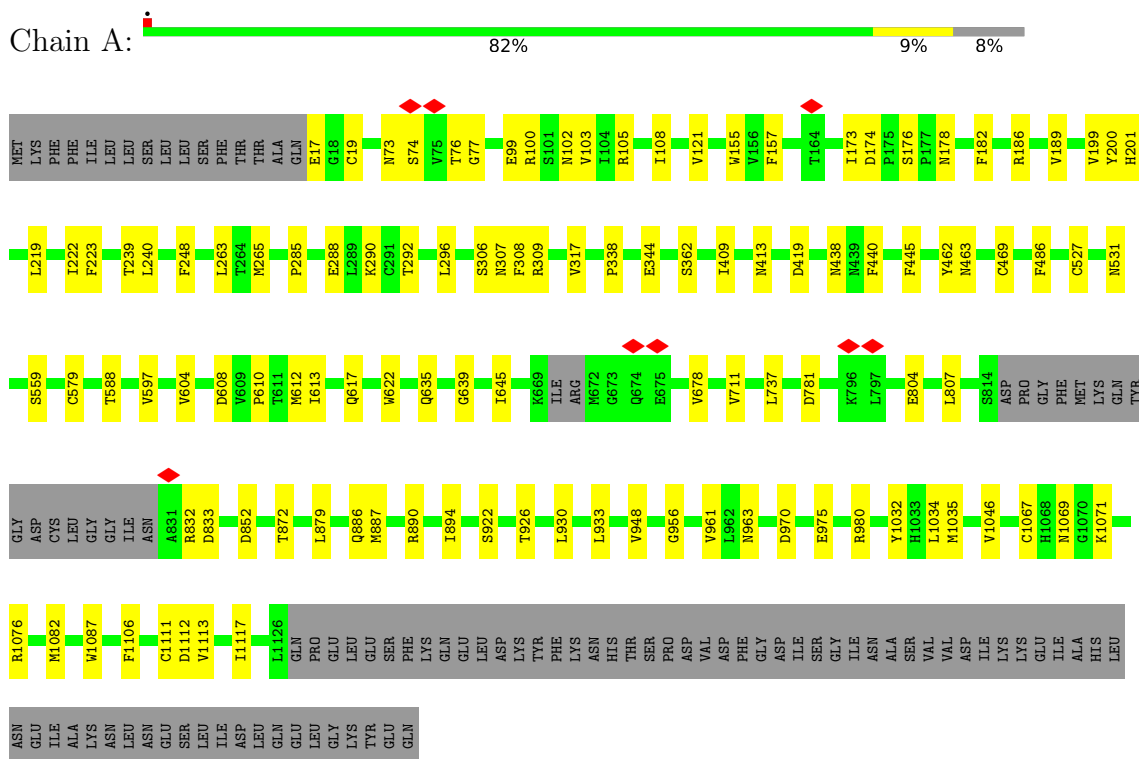


Mol	Chain	Residues	Atoms				AltConf
6	A	1	Total	C	N	O	0
			43	33	4	6	
6	B	1	Total	C	N	O	0
			43	33	4	6	
6	C	1	Total	C	N	O	0
			43	33	4	6	

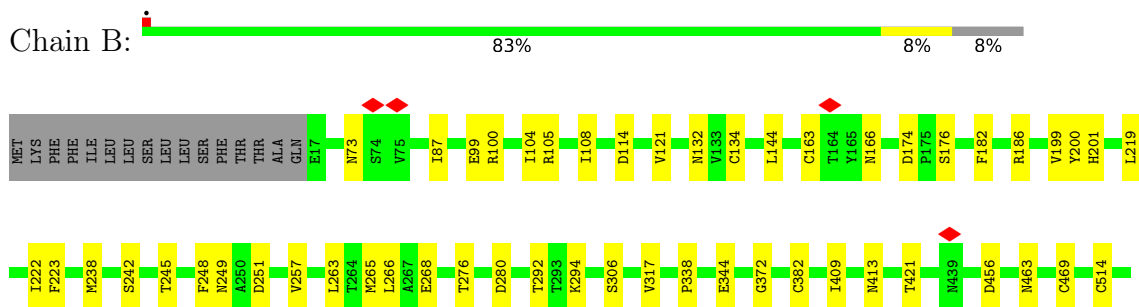
3 Residue-property plots [i](#)

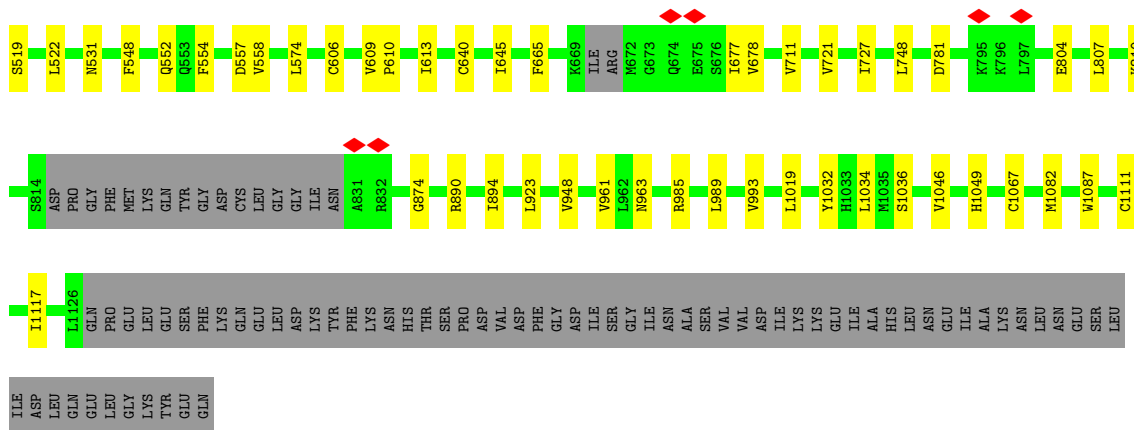
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Spike glycoprotein

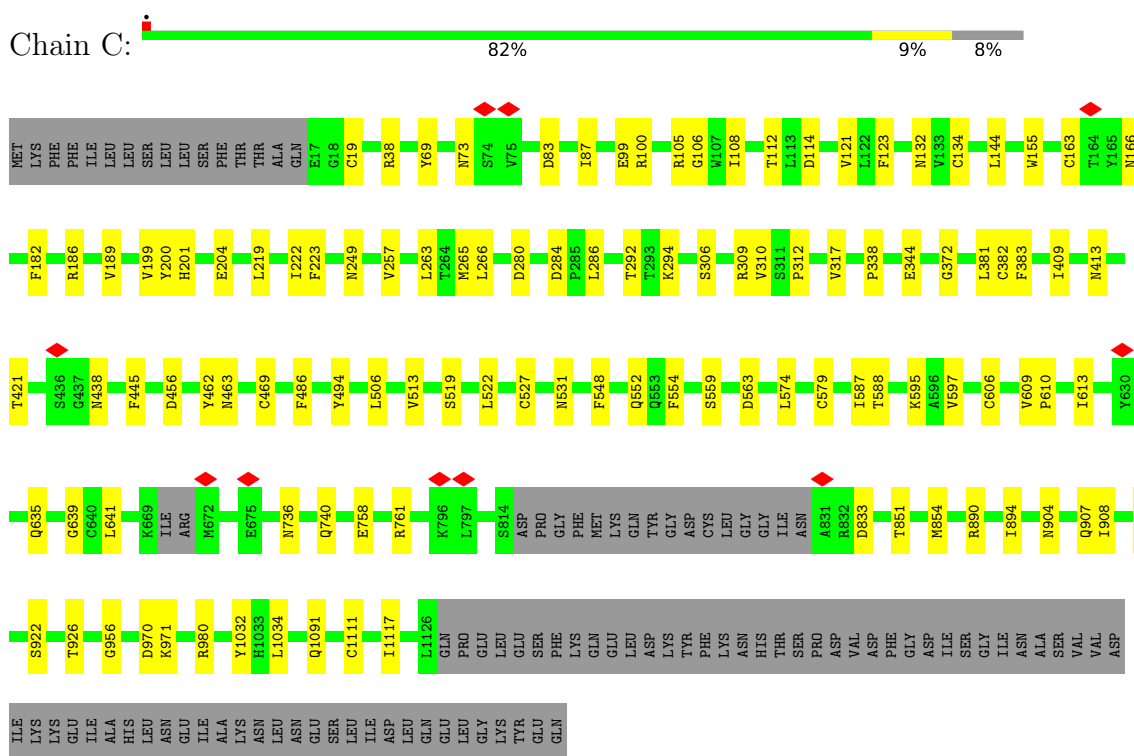


• Molecule 1: Spike glycoprotein

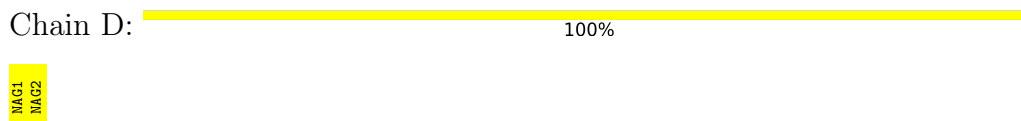




- Molecule 1: Spike glycoprotein



- 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



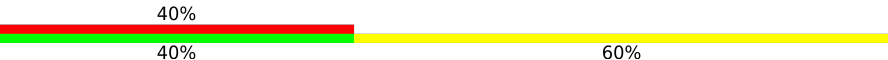
MAG1
MAG2

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

Chain H:  100%

MAG1
MAG2

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

Chain E:  40% 40% 60%

MAG1
MAG2
BMA3
MAN4
MAN5

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

Chain G:  75% 25%

MAG1
MAG2
BMA3
MAN4

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

Chain I:  75% 25%

MAG1
MAG2
BMA3
MAN4

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	98797	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.267	Depositor
Minimum map value	-0.156	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.024	Depositor
Map size (Å)	280.32, 280.32, 280.32	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.095, 1.095, 1.095	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, MAN, BLA, BMA

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	0.15	0/8706	0.39	0/11843
1	B	0.20	0/8706	0.40	0/11843
1	C	0.15	0/8702	0.37	0/11839
All	All	0.17	0/26114	0.39	0/35525

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	8509	0	8238	67	0
1	B	8509	0	8238	55	0
1	C	8505	0	8227	62	0
2	D	28	0	25	0	0
2	F	28	0	25	0	0
2	H	28	0	25	0	0
3	E	61	0	52	0	0
4	G	50	0	43	0	0
4	I	50	0	43	0	0
5	A	238	0	221	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	238	0	221	0	0
5	C	238	0	221	0	0
6	A	43	0	32	4	0
6	B	43	0	32	0	0
6	C	43	0	32	3	0
All	All	26611	0	25675	185	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 4.

All (185) 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:382:CYS:HA	1:B:514:CYS:HB3	1.48	0.95
6:A:1218:BLA:HHA	6:A:1218:BLA:HBD2	1.70	0.73
6:A:1218:BLA:HHD	6:A:1218:BLA:HBC1	1.71	0.73
1:C:382:CYS:SG	1:C:513:VAL:O	2.50	0.70
1:B:104:ILE:HG22	1:B:238:MET:HG2	1.72	0.70
1:A:887:MET:HE1	1:A:1035:MET:HE1	1.74	0.69
1:A:240:LEU:HD13	1:A:248:PHE:HB3	1.74	0.69
1:A:317:VAL:HG12	1:A:531:ASN:HB3	1.76	0.68
1:B:317:VAL:HG12	1:B:531:ASN:HB3	1.78	0.65
1:A:463:ASN:HB2	1:A:469:CYS:HB3	1.77	0.64
1:B:874:GLY:HA3	1:B:1019:LEU:HD12	1.78	0.63
1:B:557:ASP:OD1	1:B:558:VAL:N	2.32	0.63
1:B:748:LEU:HG	1:B:993:VAL:HG21	1.84	0.60
1:A:173:ILE:HG22	6:A:1218:BLA:HMD2	1.84	0.60
1:B:265:MET:HG2	1:B:280:ASP:HA	1.84	0.59
1:C:265:MET:HG2	1:C:280:ASP:HA	1.84	0.59
1:C:317:VAL:HG12	1:C:531:ASN:HB3	1.84	0.58
1:A:99:GLU:OE2	1:A:186:ARG:NH1	2.35	0.58
1:A:438:ASN:HD21	1:A:486:PHE:HB2	1.69	0.58
1:B:317:VAL:HG23	1:B:519:SER:HA	1.84	0.57
1:B:552:GLN:NE2	1:B:554:PHE:O	2.36	0.57
1:A:199:VAL:HB	1:A:223:PHE:HB2	1.86	0.57
6:C:1217:BLA:HBD1	6:C:1217:BLA:HHA	1.86	0.56
1:B:372:GLY:HA3	1:B:421:THR:HG23	1.87	0.56
1:C:372:GLY:HA3	1:C:421:THR:HG23	1.87	0.56
1:C:552:GLN:NE2	1:C:554:PHE:O	2.37	0.56
1:C:99:GLU:OE2	1:C:186:ARG:NH1	2.39	0.56
1:C:890:ARG:NH1	1:C:1034:LEU:O	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:894:ILE:HD12	1:C:1032:TYR:HB3	1.89	0.55
1:C:199:VAL:HB	1:C:223:PHE:HB2	1.87	0.55
1:C:610:PRO:HA	1:C:613:ILE:HG22	1.88	0.55
1:C:438:ASN:ND2	1:C:486:PHE:O	2.40	0.55
1:C:907:GLN:NE2	1:C:911:GLN:OE1	2.40	0.55
1:C:736:ASN:O	1:C:740:GLN:NE2	2.40	0.55
1:B:99:GLU:OE2	1:B:186:ARG:NH1	2.40	0.54
1:A:419:ASP:OD2	1:C:971:LYS:NZ	2.40	0.54
1:B:890:ARG:NH1	1:B:1034:LEU:O	2.40	0.54
1:C:338:PRO:HG3	1:C:344:GLU:HG3	1.89	0.54
1:C:310:VAL:HG12	1:C:312:PRO:HD3	1.90	0.54
1:C:606:CYS:HA	1:C:609:VAL:HG23	1.89	0.54
1:B:711:VAL:HG12	1:B:1046:VAL:HG22	1.90	0.53
1:A:100:ARG:NH1	1:A:178:ASN:O	2.42	0.53
1:A:338:PRO:HG3	1:A:344:GLU:HG3	1.91	0.53
1:A:610:PRO:HA	1:A:613:ILE:HG22	1.91	0.53
1:B:199:VAL:HB	1:B:223:PHE:HB2	1.90	0.53
1:C:292:THR:HG21	1:C:306:SER:H	1.74	0.52
1:C:463:ASN:HB2	1:C:469:CYS:HB3	1.91	0.52
1:B:548:PHE:HB3	1:B:552:GLN:HB3	1.90	0.52
1:B:132:ASN:OD1	1:B:166:ASN:ND2	2.40	0.52
1:B:727:ILE:O	1:B:985:ARG:NH1	2.42	0.52
1:C:105:ARG:HD2	1:C:144:LEU:HD21	1.91	0.52
1:C:548:PHE:HB3	1:C:552:GLN:HB3	1.92	0.52
1:B:268:GLU:OE2	1:B:276:THR:OG1	2.24	0.51
1:A:100:ARG:HB3	1:A:182:PHE:HA	1.92	0.51
1:A:1069:ASN:OD1	1:A:1071:LYS:NZ	2.40	0.51
1:B:73:ASN:ND2	1:B:249:ASN:OD1	2.44	0.51
1:B:894:ILE:HD12	1:B:1032:TYR:HB3	1.93	0.51
1:C:317:VAL:HG23	1:C:519:SER:HA	1.93	0.51
1:C:73:ASN:ND2	1:C:249:ASN:OD1	2.44	0.51
1:A:200:TYR:HB3	1:A:219:LEU:HB3	1.92	0.50
1:B:804:GLU:HA	1:B:807:LEU:HG	1.94	0.50
1:A:409:ILE:HD13	1:A:413:ASN:HD22	1.76	0.50
1:C:132:ASN:OD1	1:C:166:ASN:ND2	2.40	0.50
1:A:608:ASP:N	1:A:608:ASP:OD1	2.44	0.50
1:A:362:SER:O	1:C:494:TYR:OH	2.29	0.50
1:C:266:LEU:HD11	1:C:294:LYS:HA	1.93	0.49
1:A:922:SER:O	1:A:926:THR:OG1	2.31	0.49
1:B:1111:CYS:HB2	1:B:1117:ILE:HD13	1.95	0.49
1:C:106:GLY:HA3	1:C:123:PHE:HA	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:C:1217:BLA:HHD	6:C:1217:BLA:HBC1	1.94	0.49
1:A:103:VAL:HG23	1:A:239:THR:HB	1.94	0.49
1:B:266:LEU:HD11	1:B:294:LYS:HA	1.94	0.49
1:B:610:PRO:HA	1:B:613:ILE:HG22	1.93	0.49
1:C:1111:CYS:HB2	1:C:1117:ILE:HD13	1.94	0.49
1:C:851:THR:HG22	1:C:854:MET:HE2	1.94	0.48
1:C:922:SER:O	1:C:926:THR:OG1	2.31	0.48
1:A:19:CYS:HA	1:A:155:TRP:HB3	1.94	0.48
1:A:102:ASN:O	1:A:105:ARG:NH2	2.46	0.48
1:A:17:GLU:N	1:A:157:PHE:O	2.47	0.48
1:C:833:ASP:N	1:C:833:ASP:OD1	2.45	0.48
1:A:737:LEU:HD11	1:A:975:GLU:HG3	1.96	0.48
1:B:134:CYS:HA	1:B:163:CYS:HB3	1.95	0.47
1:A:956:GLY:O	1:A:980:ARG:NH1	2.47	0.47
1:C:456:ASP:OD2	1:C:456:ASP:N	2.45	0.47
1:A:108:ILE:HG12	1:A:121:VAL:HG12	1.96	0.47
1:B:263:LEU:HD11	1:B:265:MET:HE3	1.97	0.47
1:B:338:PRO:HG3	1:B:344:GLU:HG3	1.95	0.47
1:C:19:CYS:HA	1:C:155:TRP:HB3	1.97	0.47
1:A:970:ASP:OD1	1:A:970:ASP:N	2.47	0.47
1:B:201:HIS:HB2	1:B:222:ILE:HD13	1.97	0.46
1:B:242:SER:OG	1:B:245:THR:OG1	2.30	0.46
1:B:456:ASP:OD2	1:B:456:ASP:N	2.49	0.46
1:A:201:HIS:HB2	1:A:222:ILE:HD13	1.98	0.46
1:B:1067:CYS:HB2	1:B:1111:CYS:HB2	1.83	0.46
1:C:100:ARG:HB3	1:C:182:PHE:HD1	1.80	0.46
1:A:804:GLU:HA	1:A:807:LEU:HG	1.98	0.46
1:C:382:CYS:SG	1:C:383:PHE:N	2.89	0.45
1:B:242:SER:HB2	1:B:248:PHE:H	1.81	0.45
1:C:758:GLU:HG3	1:C:761:ARG:HH21	1.80	0.45
1:C:956:GLY:O	1:C:980:ARG:NH1	2.49	0.45
1:A:711:VAL:HG12	1:A:1046:VAL:HG22	1.99	0.45
1:C:263:LEU:HD11	1:C:265:MET:HE3	1.98	0.45
1:A:894:ILE:HD12	1:A:1032:TYR:HB3	1.99	0.45
1:A:886:GLN:HE21	1:A:890:ARG:HE	1.63	0.45
1:B:174:ASP:OD2	1:B:176:SER:OG	2.32	0.45
1:B:105:ARG:HD2	1:B:144:LEU:HD21	1.99	0.45
1:A:781:ASP:OD1	1:A:781:ASP:N	2.50	0.44
1:C:87:ILE:HG21	1:C:257:VAL:HG11	1.99	0.44
6:C:1217:BLA:HBD1	6:C:1217:BLA:CHA	2.47	0.44
1:B:292:THR:HG21	1:B:306:SER:H	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:527:CYS:HB2	1:C:579:CYS:HB2	1.79	0.44
1:A:1076:ARG:HH21	1:A:1106:PHE:HB3	1.83	0.44
1:B:606:CYS:HA	1:B:609:VAL:HG23	1.99	0.44
1:B:1036:SER:OG	1:B:1049:HIS:ND1	2.43	0.44
1:A:645:ILE:HG13	1:A:678:VAL:HG12	2.00	0.44
1:B:961:VAL:HG12	1:B:963:ASN:H	1.83	0.44
1:A:635:GLN:NE2	1:A:639:GLY:O	2.50	0.43
1:A:833:ASP:OD1	1:A:833:ASP:N	2.49	0.43
1:A:930:LEU:HD23	1:A:933:LEU:HD23	2.00	0.43
1:C:134:CYS:HA	1:C:163:CYS:HB3	2.00	0.43
1:C:588:THR:HB	1:C:597:VAL:HG12	2.00	0.43
1:A:832:ARG:NH2	1:C:563:ASP:OD2	2.40	0.43
1:C:200:TYR:HB3	1:C:219:LEU:HB3	1.99	0.43
1:C:587:ILE:HD12	1:C:641:LEU:HD21	2.00	0.43
1:A:189:VAL:HG22	1:A:200:TYR:HB2	2.00	0.43
1:C:970:ASP:OD1	1:C:970:ASP:N	2.47	0.43
1:A:438:ASN:HD22	1:A:440:PHE:HD2	1.66	0.43
1:A:174:ASP:OD2	1:A:176:SER:OG	2.31	0.43
6:A:1218:BLA:HA	6:A:1218:BLA:HB	1.66	0.43
1:C:38:ARG:NH1	1:C:204:GLU:OE2	2.50	0.43
1:A:1067:CYS:HB2	1:A:1111:CYS:HB2	1.83	0.43
1:B:87:ILE:HG21	1:B:257:VAL:HG11	2.00	0.43
1:B:200:TYR:HB3	1:B:219:LEU:HB3	2.00	0.43
1:C:445:PHE:HB3	1:C:462:TYR:CD2	2.54	0.43
1:C:635:GLN:NE2	1:C:639:GLY:O	2.52	0.43
1:A:292:THR:HG21	1:A:306:SER:H	1.82	0.42
1:C:189:VAL:HG22	1:C:200:TYR:HB2	2.01	0.42
1:A:527:CYS:HB2	1:A:579:CYS:HB2	1.96	0.42
1:A:527:CYS:H	1:A:617:GLN:HE22	1.67	0.42
1:B:463:ASN:HB2	1:B:469:CYS:HB2	2.00	0.42
1:A:73:ASN:OD1	1:A:74:SER:N	2.52	0.42
1:C:286:LEU:HD21	1:C:595:LYS:HD2	2.00	0.42
1:B:108:ILE:HG12	1:B:121:VAL:HG12	2.02	0.42
1:B:409:ILE:HD13	1:B:413:ASN:HD22	1.85	0.42
1:A:445:PHE:HB3	1:A:462:TYR:CD2	2.55	0.42
1:B:748:LEU:HD21	1:B:989:LEU:HB3	2.02	0.42
1:B:810:LYS:NZ	1:B:923:LEU:O	2.35	0.42
1:C:69:TYR:HB3	1:C:83:ASP:HB2	2.01	0.42
1:A:307:ASN:HD22	1:A:308:PHE:N	2.18	0.42
1:B:522:LEU:HD21	1:B:574:LEU:HD11	2.02	0.42
1:C:201:HIS:HB2	1:C:222:ILE:HD13	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:904:ASN:O	1:C:908:ILE:HG12	2.20	0.42
1:A:263:LEU:HD11	1:A:265:MET:HE3	2.01	0.41
1:A:240:LEU:HD23	1:A:240:LEU:HA	1.81	0.41
1:A:1112:ASP:OD1	1:A:1113:VAL:N	2.53	0.41
1:B:721:VAL:HG21	1:B:989:LEU:HD21	2.01	0.41
1:B:1082:MET:HB3	1:B:1087:TRP:CD2	2.55	0.41
1:C:284:ASP:OD1	1:C:284:ASP:N	2.52	0.41
1:C:309:ARG:H	1:C:309:ARG:HG2	1.65	0.41
1:C:409:ILE:HA	1:C:413:ASN:HB2	2.02	0.41
1:C:112:THR:OG1	1:C:114:ASP:OD1	2.37	0.41
1:C:381:LEU:HD21	1:C:506:LEU:HD11	2.02	0.41
1:A:559:SER:HB3	1:B:948:VAL:HG22	2.03	0.41
1:A:588:THR:HB	1:A:597:VAL:HG12	2.03	0.41
1:A:948:VAL:HG22	1:C:559:SER:HB3	2.02	0.41
1:B:251:ASP:N	1:B:251:ASP:OD1	2.53	0.41
1:B:645:ILE:HG13	1:B:678:VAL:HG12	2.03	0.41
1:A:604:VAL:HG11	1:A:612:MET:HE3	2.02	0.41
1:A:872:THR:HG21	1:A:879:LEU:HB2	2.02	0.41
1:B:100:ARG:HB3	1:B:182:PHE:HD1	1.85	0.41
1:B:114:ASP:N	1:B:114:ASP:OD1	2.52	0.41
1:A:285:PRO:HD3	1:A:622:TRP:CE2	2.56	0.40
1:A:290:LYS:HD2	1:A:296:LEU:HA	2.02	0.40
1:A:852:ASP:N	1:A:852:ASP:OD1	2.53	0.40
1:A:288:GLU:HG3	1:A:308:PHE:HD2	1.86	0.40
1:A:309:ARG:H	1:A:309:ARG:HG2	1.63	0.40
1:B:665:PHE:HD1	1:B:677:ILE:HG22	1.86	0.40
1:C:108:ILE:HG12	1:C:121:VAL:HG12	2.02	0.40
1:A:1111:CYS:HB2	1:A:1117:ILE:HD13	2.04	0.40
1:A:961:VAL:HG12	1:A:963:ASN:H	1.87	0.40
1:B:781:ASP:OD1	1:B:781:ASP:N	2.49	0.40
1:C:522:LEU:HD21	1:C:574:LEU:HD11	2.04	0.40
1:A:76:THR:OG1	1:A:77:GLY:N	2.54	0.40
1:A:890:ARG:NH1	1:A:1034:LEU:O	2.51	0.40
1:A:1082:MET:HB3	1:A:1087:TRP:CD2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	1086/1193 (91%)	1035 (95%)	51 (5%)	0	100	100
1	B	1086/1193 (91%)	1038 (96%)	48 (4%)	0	100	100
1	C	1086/1193 (91%)	1035 (95%)	51 (5%)	0	100	100
All	All	3258/3579 (91%)	3108 (95%)	150 (5%)	0	100	100

There are no Ramachandran outliers to report.

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 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	952/1042 (91%)	952 (100%)	0	100	100
1	B	952/1042 (91%)	951 (100%)	1 (0%)	88	93
1	C	951/1042 (91%)	950 (100%)	1 (0%)	88	93
All	All	2855/3126 (91%)	2853 (100%)	2 (0%)	87	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	640	CYS
1	C	1091	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	A	428	ASN
1	A	525	ASN
1	A	553	GLN
1	A	617	GLN
1	A	635	GLN
1	A	736	ASN
1	A	769	GLN
1	A	841	ASN
1	A	886	GLN
1	A	920	GLN
1	A	977	GLN
1	A	1033	HIS
1	A	1039	GLN
1	A	1098	GLN
1	B	188	HIS
1	B	379	ASN
1	B	428	ASN
1	B	525	ASN
1	B	531	ASN
1	B	553	GLN
1	B	841	ASN
1	B	920	GLN
1	B	950	GLN
1	B	977	GLN
1	B	990	GLN
1	B	1033	HIS
1	B	1039	GLN
1	B	1098	GLN
1	C	63	ASN
1	C	379	ASN
1	C	428	ASN
1	C	463	ASN
1	C	525	ASN
1	C	553	GLN
1	C	602	GLN
1	C	617	GLN
1	C	635	GLN
1	C	747	GLN
1	C	841	ASN
1	C	886	GLN
1	C	898	GLN
1	C	907	GLN

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Mol	Chain	Res	Type
1	C	950	GLN
1	C	1033	HIS
1	C	1039	GLN
1	C	1091	GLN
1	C	1098	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 ⓘ

19 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	2,1	14,14,15	0.67	1 (7%)	17,19,21	2.09	2 (11%)
2	NAG	D	2	2	14,14,15	0.40	0	17,19,21	0.86	1 (5%)
3	NAG	E	1	3,1	14,14,15	0.27	0	17,19,21	0.55	0
3	NAG	E	2	3	14,14,15	0.33	0	17,19,21	0.58	0
3	BMA	E	3	3	11,11,12	1.20	2 (18%)	15,15,17	1.19	2 (13%)
3	MAN	E	4	3	11,11,12	1.68	3 (27%)	15,15,17	2.25	2 (13%)
3	MAN	E	5	3	11,11,12	0.70	0	15,15,17	1.03	2 (13%)
2	NAG	F	1	2,1	14,14,15	0.66	0	17,19,21	2.10	2 (11%)
2	NAG	F	2	2	14,14,15	0.38	0	17,19,21	0.87	1 (5%)
4	NAG	G	1	4,1	14,14,15	0.26	0	17,19,21	0.54	0
4	NAG	G	2	4	14,14,15	0.23	0	17,19,21	0.50	0
4	BMA	G	3	4	11,11,12	0.70	0	15,15,17	0.80	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	G	4	4	11,11,12	0.97	0	15,15,17	1.06	2 (13%)
2	NAG	H	1	2,1	14,14,15	0.69	1 (7%)	17,19,21	2.09	2 (11%)
2	NAG	H	2	2	14,14,15	0.39	0	17,19,21	0.86	1 (5%)
4	NAG	I	1	4,1	14,14,15	0.24	0	17,19,21	0.55	0
4	NAG	I	2	4	14,14,15	0.24	0	17,19,21	0.50	0
4	BMA	I	3	4	11,11,12	0.69	0	15,15,17	0.82	0
4	MAN	I	4	4	11,11,12	0.96	0	15,15,17	1.10	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	D	1	2,1	-	6/6/23/26	0/1/1/1
2	NAG	D	2	2	-	4/6/23/26	0/1/1/1
3	NAG	E	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	BMA	E	3	3	-	2/2/19/22	0/1/1/1
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
3	MAN	E	5	3	-	2/2/19/22	0/1/1/1
2	NAG	F	1	2,1	-	6/6/23/26	0/1/1/1
2	NAG	F	2	2	-	4/6/23/26	0/1/1/1
4	NAG	G	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
4	BMA	G	3	4	-	0/2/19/22	0/1/1/1
4	MAN	G	4	4	-	0/2/19/22	1/1/1/1
2	NAG	H	1	2,1	-	6/6/23/26	0/1/1/1
2	NAG	H	2	2	-	4/6/23/26	0/1/1/1
4	NAG	I	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	I	2	4	-	2/6/23/26	0/1/1/1
4	BMA	I	3	4	-	0/2/19/22	0/1/1/1
4	MAN	I	4	4	-	0/2/19/22	1/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	4	MAN	O5-C1	3.43	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	4	MAN	O5-C5	3.23	1.49	1.43
3	E	4	MAN	C1-C2	2.59	1.58	1.52
3	E	3	BMA	C2-C3	2.50	1.56	1.52
2	H	1	NAG	C1-C2	2.09	1.55	1.52
3	E	3	BMA	C1-C2	2.07	1.57	1.52
2	D	1	NAG	C1-C2	2.01	1.55	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	4	MAN	C1-O5-C5	7.79	122.63	112.19
2	F	1	NAG	C2-N2-C7	7.37	132.78	122.90
2	D	1	NAG	C2-N2-C7	7.33	132.73	122.90
2	H	1	NAG	C2-N2-C7	7.31	132.70	122.90
2	F	1	NAG	C1-C2-N2	3.57	116.05	110.43
2	D	1	NAG	C1-C2-N2	3.56	116.04	110.43
2	H	1	NAG	C1-C2-N2	3.49	115.93	110.43
4	I	4	MAN	C1-O5-C5	3.19	116.47	112.19
4	G	4	MAN	C1-O5-C5	3.02	116.23	112.19
3	E	5	MAN	C1-O5-C5	2.86	116.02	112.19
3	E	4	MAN	O2-C2-C3	-2.74	104.47	110.15
2	H	2	NAG	C2-N2-C7	2.60	126.38	122.90
2	F	2	NAG	C2-N2-C7	2.58	126.36	122.90
2	D	2	NAG	C2-N2-C7	2.57	126.34	122.90
3	E	3	BMA	C1-C2-C3	2.55	113.36	109.64
3	E	5	MAN	O2-C2-C3	-2.20	105.58	110.15
4	I	4	MAN	O2-C2-C3	-2.20	105.59	110.15
4	G	4	MAN	O2-C2-C3	-2.18	105.63	110.15
3	E	3	BMA	C1-O5-C5	2.04	114.92	112.19

There are no chirality outliers.

All (46) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1	NAG	O5-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
2	H	1	NAG	O5-C5-C6-O6
3	E	2	NAG	O5-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6
4	I	2	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	E	1	NAG	C4-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
4	I	1	NAG	C4-C5-C6-O6
4	I	2	NAG	C4-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
4	G	2	NAG	C4-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
2	H	1	NAG	C4-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	D	1	NAG	C8-C7-N2-C2
2	D	1	NAG	O7-C7-N2-C2
2	F	1	NAG	C8-C7-N2-C2
2	F	1	NAG	O7-C7-N2-C2
2	H	1	NAG	C8-C7-N2-C2
2	H	1	NAG	O7-C7-N2-C2
3	E	5	MAN	O5-C5-C6-O6
2	H	2	NAG	C1-C2-N2-C7
2	F	1	NAG	C3-C2-N2-C7
2	H	1	NAG	C3-C2-N2-C7
3	E	5	MAN	C4-C5-C6-O6
3	E	3	BMA	C4-C5-C6-O6
2	D	1	NAG	C1-C2-N2-C7
2	D	2	NAG	C1-C2-N2-C7
2	F	1	NAG	C1-C2-N2-C7
2	F	2	NAG	C1-C2-N2-C7
2	H	1	NAG	C1-C2-N2-C7
2	D	1	NAG	C3-C2-N2-C7
2	D	2	NAG	C3-C2-N2-C7
2	F	2	NAG	C3-C2-N2-C7
2	H	2	NAG	C3-C2-N2-C7
3	E	3	BMA	O5-C5-C6-O6

All (2) ring outliers are listed below:

Continued on next page...

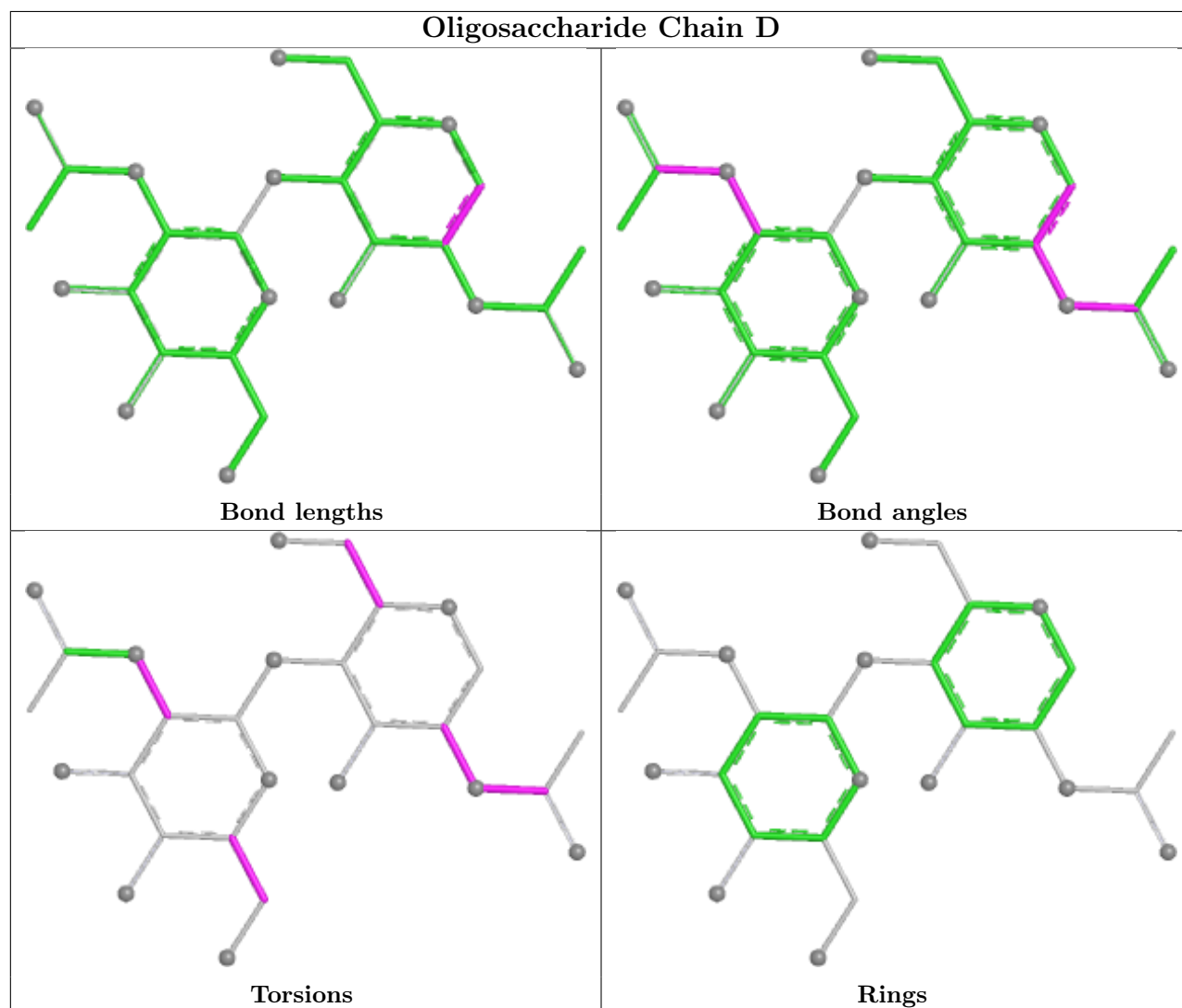
Continued from previous page...

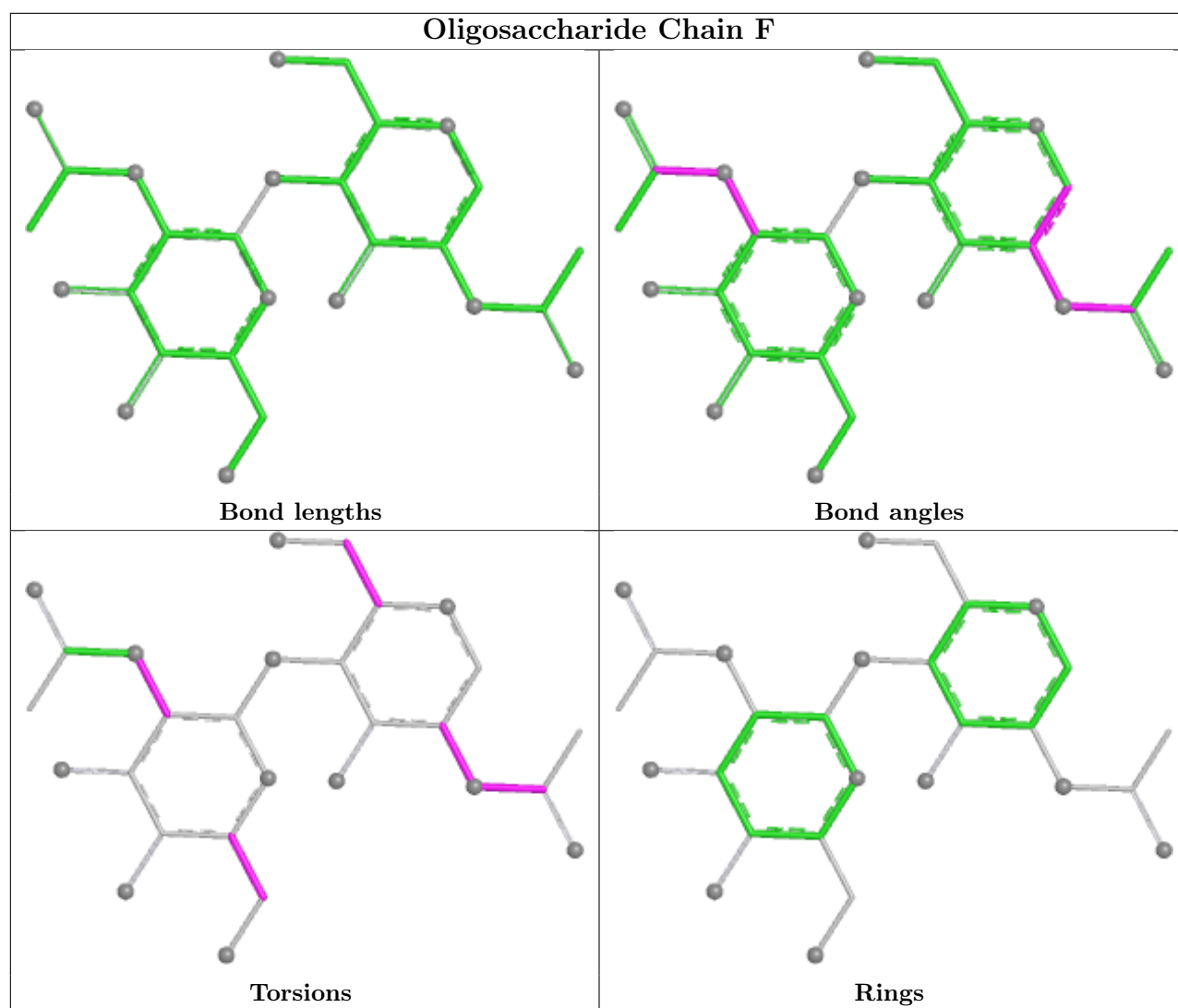
Mol	Chain	Res	Type	Atoms
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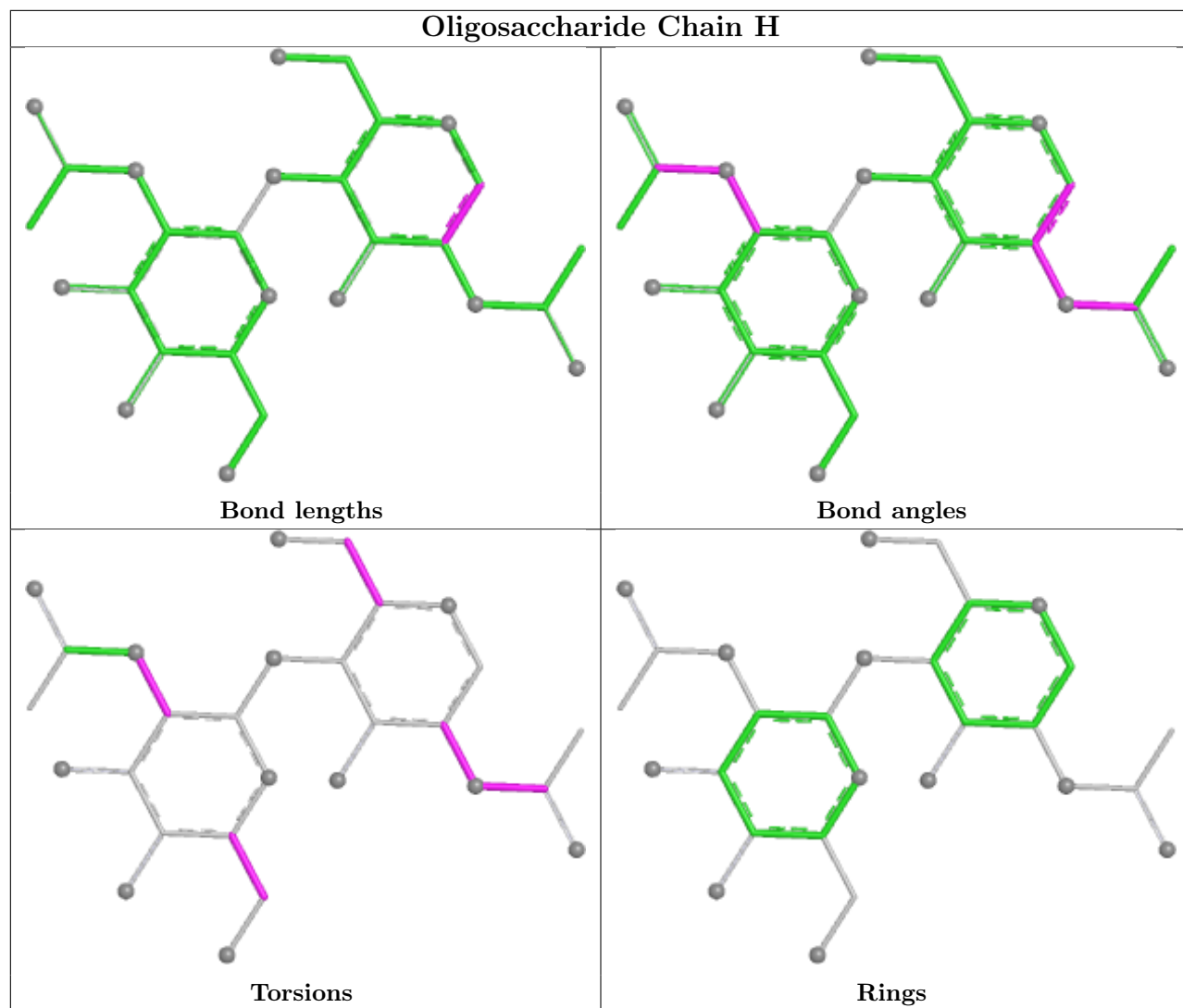
Mol	Chain	Res	Type	Atoms
4	G	4	MAN	C1-C2-C3-C4-C5-O5
4	I	4	MAN	C1-C2-C3-C4-C5-O5

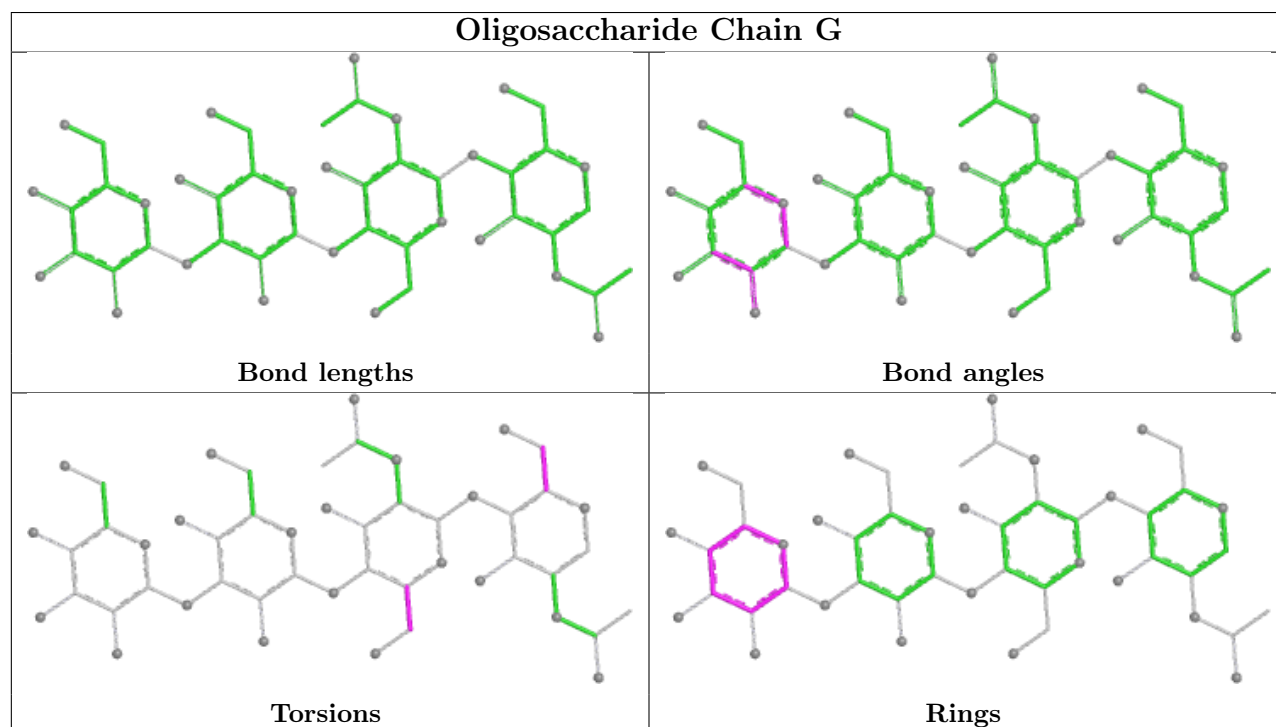
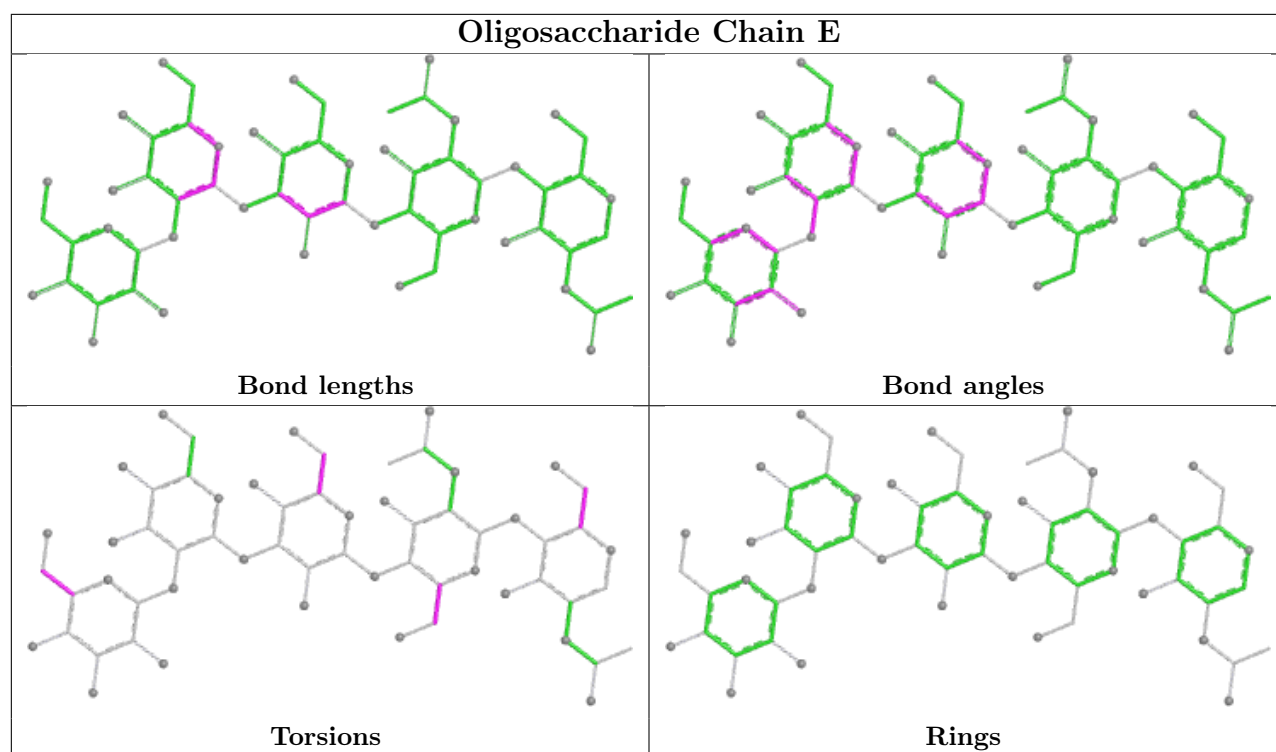
No monomer is involved in short contacts.

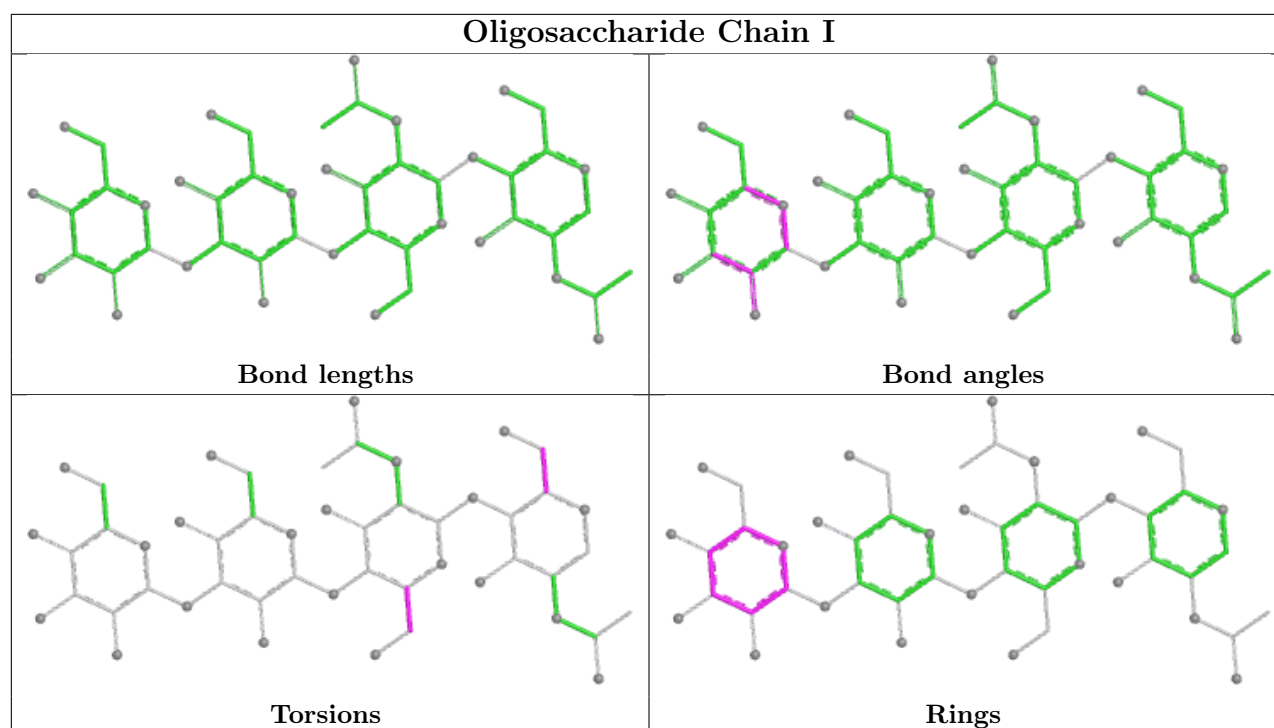
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

54 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	$\# Z > 2$
6	BLA	B	1217	-	46,46,46	1.06	3 (6%)	66,67,67	1.01	4 (6%)
5	NAG	C	1218	1	14,14,15	0.40	0	17,19,21	0.88	1 (5%)
5	NAG	C	1204	1	14,14,15	0.29	0	17,19,21	0.48	0
5	NAG	A	1204	1	14,14,15	0.24	0	17,19,21	0.43	0
5	NAG	C	1207	1	14,14,15	0.43	0	17,19,21	0.54	0
5	NAG	C	1211	1	14,14,15	0.38	0	17,19,21	0.87	1 (5%)
5	NAG	A	1205	1	14,14,15	0.28	0	17,19,21	0.49	0
5	NAG	B	1201	1	14,14,15	0.27	0	17,19,21	0.50	0
5	NAG	C	1201	1	14,14,15	0.28	0	17,19,21	0.50	0
5	NAG	C	1214	1	14,14,15	0.28	0	17,19,21	0.49	0
5	NAG	B	1202	1	14,14,15	0.35	0	17,19,21	0.47	0
5	NAG	A	1213	1	14,14,15	0.27	0	17,19,21	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	C	1209	1	14,14,15	0.34	0	17,19,21	0.50	0
5	NAG	B	1213	1	14,14,15	0.44	0	17,19,21	0.90	1 (5%)
5	NAG	B	1203	1	14,14,15	0.76	1 (7%)	17,19,21	2.10	2 (11%)
5	NAG	B	1206	1	14,14,15	0.38	0	17,19,21	0.88	1 (5%)
6	BLA	A	1218	-	46,46,46	1.06	3 (6%)	66,67,67	1.02	3 (4%)
5	NAG	A	1215	1	14,14,15	0.29	0	17,19,21	0.49	0
5	NAG	B	1212	1	14,14,15	0.28	0	17,19,21	0.49	0
5	NAG	B	1214	1	14,14,15	0.28	0	17,19,21	0.49	0
6	BLA	C	1217	-	46,46,46	1.07	3 (6%)	66,67,67	1.01	3 (4%)
5	NAG	A	1208	1	14,14,15	0.44	0	17,19,21	0.55	0
5	NAG	A	1212	1	14,14,15	0.38	0	17,19,21	0.86	1 (5%)
5	NAG	B	1215	1	14,14,15	0.29	0	17,19,21	0.47	0
5	NAG	B	1207	1	14,14,15	0.44	0	17,19,21	0.55	0
5	NAG	C	1216	1	14,14,15	0.27	0	17,19,21	0.49	0
5	NAG	A	1203	1	14,14,15	0.76	1 (7%)	17,19,21	2.11	2 (11%)
5	NAG	B	1205	1	14,14,15	0.28	0	17,19,21	0.49	0
5	NAG	A	1207	1	14,14,15	0.38	0	17,19,21	0.86	1 (5%)
5	NAG	A	1201	1	14,14,15	0.26	0	17,19,21	0.50	0
5	NAG	C	1208	1	14,14,15	0.33	0	17,19,21	0.85	1 (5%)
5	NAG	B	1210	1	14,14,15	0.25	0	17,19,21	0.46	0
5	NAG	B	1208	1	14,14,15	0.34	0	17,19,21	0.85	1 (5%)
5	NAG	C	1203	1	14,14,15	0.26	0	17,19,21	0.45	0
5	NAG	A	1206	1	14,14,15	0.36	0	17,19,21	0.47	0
5	NAG	A	1209	1	14,14,15	0.34	0	17,19,21	0.85	1 (5%)
5	NAG	B	1216	1	14,14,15	0.27	0	17,19,21	0.48	0
5	NAG	A	1210	1	14,14,15	0.34	0	17,19,21	0.52	0
5	NAG	B	1218	1	14,14,15	0.27	0	17,19,21	0.39	0
5	NAG	C	1212	1	14,14,15	0.28	0	17,19,21	0.48	0
5	NAG	A	1202	1	14,14,15	0.33	0	17,19,21	0.47	0
5	NAG	B	1204	1	14,14,15	0.25	0	17,19,21	0.44	0
5	NAG	B	1209	1	14,14,15	0.32	0	17,19,21	0.49	0
5	NAG	C	1205	1	14,14,15	0.33	0	17,19,21	0.45	0
5	NAG	A	1217	1	14,14,15	0.28	0	17,19,21	0.49	0
5	NAG	B	1211	1	14,14,15	0.38	0	17,19,21	0.87	1 (5%)
5	NAG	C	1215	1	14,14,15	0.27	0	17,19,21	0.49	0
5	NAG	C	1213	1	14,14,15	0.43	0	17,19,21	0.90	1 (5%)
5	NAG	C	1202	1	14,14,15	0.35	0	17,19,21	0.46	0
5	NAG	A	1216	1	14,14,15	0.28	0	17,19,21	0.49	0
5	NAG	A	1214	1	14,14,15	0.44	0	17,19,21	0.90	1 (5%)
5	NAG	C	1210	1	14,14,15	0.26	0	17,19,21	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	C	1206	1	14,14,15	0.37	0	17,19,21	0.89	1 (5%)
5	NAG	A	1211	1	14,14,15	0.25	0	17,19,21	0.45	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
6	BLA	B	1217	-	-	6/26/74/74	0/4/4/4
5	NAG	C	1218	1	-	3/6/23/26	0/1/1/1
5	NAG	C	1204	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1204	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1207	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1211	1	-	3/6/23/26	0/1/1/1
5	NAG	A	1205	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1201	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1201	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1214	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1202	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1213	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1209	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1213	1	-	4/6/23/26	0/1/1/1
5	NAG	B	1203	1	-	4/6/23/26	0/1/1/1
5	NAG	B	1206	1	-	2/6/23/26	0/1/1/1
6	BLA	A	1218	-	-	7/26/74/74	0/4/4/4
5	NAG	A	1215	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1212	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1214	1	-	2/6/23/26	0/1/1/1
6	BLA	C	1217	-	-	10/26/74/74	0/4/4/4
5	NAG	A	1208	1	-	4/6/23/26	0/1/1/1
5	NAG	A	1212	1	-	3/6/23/26	0/1/1/1
5	NAG	B	1215	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1207	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1216	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1203	1	-	4/6/23/26	0/1/1/1
5	NAG	B	1205	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1207	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1201	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1208	1	-	3/6/23/26	0/1/1/1
5	NAG	B	1210	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1208	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1203	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1206	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1209	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1216	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1210	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1218	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1212	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1202	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1204	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1209	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1205	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1217	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1211	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1215	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1213	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1202	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1216	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1214	1	-	4/6/23/26	0/1/1/1
5	NAG	C	1210	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1206	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1211	1	-	2/6/23/26	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1218	BLA	CAB-C3B	-2.76	1.40	1.47
6	C	1217	BLA	CAB-C3B	-2.76	1.40	1.47
6	B	1217	BLA	CAB-C3B	-2.72	1.40	1.47
6	C	1217	BLA	CAC-C3C	-2.70	1.40	1.47
6	B	1217	BLA	CAC-C3C	-2.69	1.40	1.47
6	A	1218	BLA	CAC-C3C	-2.65	1.40	1.47
6	A	1218	BLA	C3B-C2B	2.59	1.42	1.37
6	B	1217	BLA	C3B-C2B	2.57	1.42	1.37
6	C	1217	BLA	C3B-C2B	2.56	1.42	1.37
5	B	1203	NAG	C1-C2	2.38	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1203	NAG	C1-C2	2.33	1.55	1.52

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1203	NAG	C2-N2-C7	7.34	132.73	122.90
5	B	1203	NAG	C2-N2-C7	7.29	132.67	122.90
6	A	1218	BLA	CMC-C2C-C1C	3.69	128.93	121.21
6	C	1217	BLA	CMC-C2C-C1C	3.64	128.81	121.21
6	B	1217	BLA	CMC-C2C-C1C	3.63	128.80	121.21
5	A	1203	NAG	C1-C2-N2	3.52	115.98	110.43
5	B	1203	NAG	C1-C2-N2	3.51	115.97	110.43
6	B	1217	BLA	CMB-C2B-C1B	2.68	127.42	124.16
6	A	1218	BLA	CMB-C2B-C1B	2.67	127.41	124.16
6	C	1217	BLA	CMB-C2B-C1B	2.65	127.38	124.16
5	B	1206	NAG	C2-N2-C7	2.63	126.43	122.90
6	C	1217	BLA	C4A-CHB-C1B	-2.63	123.25	130.82
5	C	1211	NAG	C2-N2-C7	2.63	126.42	122.90
5	C	1206	NAG	C2-N2-C7	2.61	126.40	122.90
5	B	1208	NAG	C2-N2-C7	2.61	126.39	122.90
5	B	1211	NAG	C2-N2-C7	2.60	126.39	122.90
5	C	1208	NAG	C2-N2-C7	2.59	126.38	122.90
5	A	1207	NAG	C2-N2-C7	2.58	126.36	122.90
5	A	1209	NAG	C2-N2-C7	2.58	126.36	122.90
5	A	1212	NAG	C2-N2-C7	2.58	126.36	122.90
5	A	1214	NAG	C2-N2-C7	2.57	126.34	122.90
5	B	1213	NAG	C2-N2-C7	2.57	126.34	122.90
5	C	1213	NAG	C2-N2-C7	2.54	126.31	122.90
5	C	1218	NAG	C2-N2-C7	2.51	126.26	122.90
6	B	1217	BLA	C4A-CHB-C1B	-2.48	123.70	130.82
6	A	1218	BLA	C4A-CHB-C1B	-2.34	124.07	130.82
6	B	1217	BLA	C4C-CHD-C1D	-2.11	122.87	128.06

There are no chirality outliers.

All (115) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1213	NAG	C4-C5-C6-O6
5	C	1213	NAG	C4-C5-C6-O6
6	A	1218	BLA	C4D-C3D-CAD-CBD
6	C	1217	BLA	C4D-C3D-CAD-CBD
5	A	1217	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	B	1213	NAG	O5-C5-C6-O6
5	C	1216	NAG	O5-C5-C6-O6
5	A	1214	NAG	O5-C5-C6-O6
5	A	1201	NAG	C4-C5-C6-O6
5	B	1216	NAG	O5-C5-C6-O6
5	B	1201	NAG	C4-C5-C6-O6
5	B	1214	NAG	O5-C5-C6-O6
5	C	1210	NAG	O5-C5-C6-O6
5	C	1214	NAG	O5-C5-C6-O6
5	C	1213	NAG	O5-C5-C6-O6
5	C	1201	NAG	C4-C5-C6-O6
5	C	1216	NAG	C4-C5-C6-O6
5	A	1217	NAG	C4-C5-C6-O6
5	C	1210	NAG	C4-C5-C6-O6
5	A	1201	NAG	O5-C5-C6-O6
5	A	1215	NAG	O5-C5-C6-O6
5	A	1214	NAG	C4-C5-C6-O6
5	A	1201	NAG	C8-C7-N2-C2
5	A	1201	NAG	O7-C7-N2-C2
5	A	1203	NAG	C8-C7-N2-C2
5	A	1203	NAG	O7-C7-N2-C2
5	A	1208	NAG	C8-C7-N2-C2
5	A	1208	NAG	O7-C7-N2-C2
5	B	1201	NAG	C8-C7-N2-C2
5	B	1201	NAG	O7-C7-N2-C2
5	B	1203	NAG	C8-C7-N2-C2
5	B	1203	NAG	O7-C7-N2-C2
5	B	1207	NAG	C8-C7-N2-C2
5	B	1207	NAG	O7-C7-N2-C2
5	C	1201	NAG	C8-C7-N2-C2
5	C	1201	NAG	O7-C7-N2-C2
5	C	1207	NAG	C8-C7-N2-C2
5	C	1207	NAG	O7-C7-N2-C2
5	A	1208	NAG	O5-C5-C6-O6
5	B	1214	NAG	C4-C5-C6-O6
5	B	1216	NAG	C4-C5-C6-O6
5	B	1201	NAG	O5-C5-C6-O6
5	C	1214	NAG	C4-C5-C6-O6
5	C	1201	NAG	O5-C5-C6-O6
6	A	1218	BLA	C2D-C3D-CAD-CBD
6	C	1217	BLA	C2D-C3D-CAD-CBD
5	C	1207	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	B	1207	NAG	O5-C5-C6-O6
5	A	1202	NAG	C4-C5-C6-O6
5	A	1202	NAG	O5-C5-C6-O6
5	B	1211	NAG	C4-C5-C6-O6
5	B	1211	NAG	O5-C5-C6-O6
5	A	1215	NAG	C4-C5-C6-O6
6	C	1217	BLA	C2D-C1D-CHD-C4C
5	A	1208	NAG	C4-C5-C6-O6
5	C	1218	NAG	O5-C5-C6-O6
6	B	1217	BLA	C2D-C1D-CHD-C4C
6	C	1217	BLA	ND-C1D-CHD-C4C
5	C	1205	NAG	O5-C5-C6-O6
6	B	1217	BLA	NC-C4C-CHD-C1D
5	A	1211	NAG	C4-C5-C6-O6
6	B	1217	BLA	ND-C1D-CHD-C4C
5	A	1209	NAG	C1-C2-N2-C7
5	B	1208	NAG	C1-C2-N2-C7
5	C	1208	NAG	C1-C2-N2-C7
6	B	1217	BLA	C3C-C4C-CHD-C1D
5	A	1211	NAG	O5-C5-C6-O6
6	C	1217	BLA	NC-C4C-CHD-C1D
6	A	1218	BLA	C4C-C3C-CAC-CBC
5	C	1207	NAG	C4-C5-C6-O6
5	A	1203	NAG	C3-C2-N2-C7
5	B	1203	NAG	C3-C2-N2-C7
5	B	1213	NAG	C3-C2-N2-C7
5	B	1205	NAG	C4-C5-C6-O6
5	B	1207	NAG	C4-C5-C6-O6
5	B	1210	NAG	C4-C5-C6-O6
6	B	1217	BLA	CAA-CBA-CGA-O2A
6	B	1217	BLA	CAA-CBA-CGA-O1A
6	C	1217	BLA	C3C-C4C-CHD-C1D
6	A	1218	BLA	CAA-CBA-CGA-O1A
5	A	1203	NAG	C1-C2-N2-C7
5	A	1207	NAG	C1-C2-N2-C7
5	A	1212	NAG	C1-C2-N2-C7
5	A	1214	NAG	C1-C2-N2-C7
5	B	1203	NAG	C1-C2-N2-C7
5	B	1206	NAG	C1-C2-N2-C7
5	B	1211	NAG	C1-C2-N2-C7
5	B	1213	NAG	C1-C2-N2-C7
5	C	1206	NAG	C1-C2-N2-C7

Continued on next page...

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Mol	Chain	Res	Type	Atoms
5	C	1211	NAG	C1-C2-N2-C7
5	C	1213	NAG	C1-C2-N2-C7
5	C	1218	NAG	C1-C2-N2-C7
5	B	1210	NAG	O5-C5-C6-O6
6	A	1218	BLA	CAA-CBA-CGA-O2A
6	C	1217	BLA	CAA-CBA-CGA-O2A
5	A	1207	NAG	C3-C2-N2-C7
5	A	1209	NAG	C3-C2-N2-C7
5	A	1212	NAG	C3-C2-N2-C7
5	A	1214	NAG	C3-C2-N2-C7
5	B	1206	NAG	C3-C2-N2-C7
5	B	1208	NAG	C3-C2-N2-C7
5	B	1211	NAG	C3-C2-N2-C7
5	C	1206	NAG	C3-C2-N2-C7
5	C	1208	NAG	C3-C2-N2-C7
5	C	1211	NAG	C3-C2-N2-C7
5	C	1213	NAG	C3-C2-N2-C7
5	C	1218	NAG	C3-C2-N2-C7
6	C	1217	BLA	CAD-CBD-CGD-O2D
6	A	1218	BLA	CAD-CBD-CGD-O2D
6	A	1218	BLA	CAD-CBD-CGD-O1D
6	C	1217	BLA	CAA-CBA-CGA-O1A
6	C	1217	BLA	CAD-CBD-CGD-O1D
5	C	1208	NAG	O5-C5-C6-O6
5	A	1212	NAG	C4-C5-C6-O6
5	C	1211	NAG	C4-C5-C6-O6

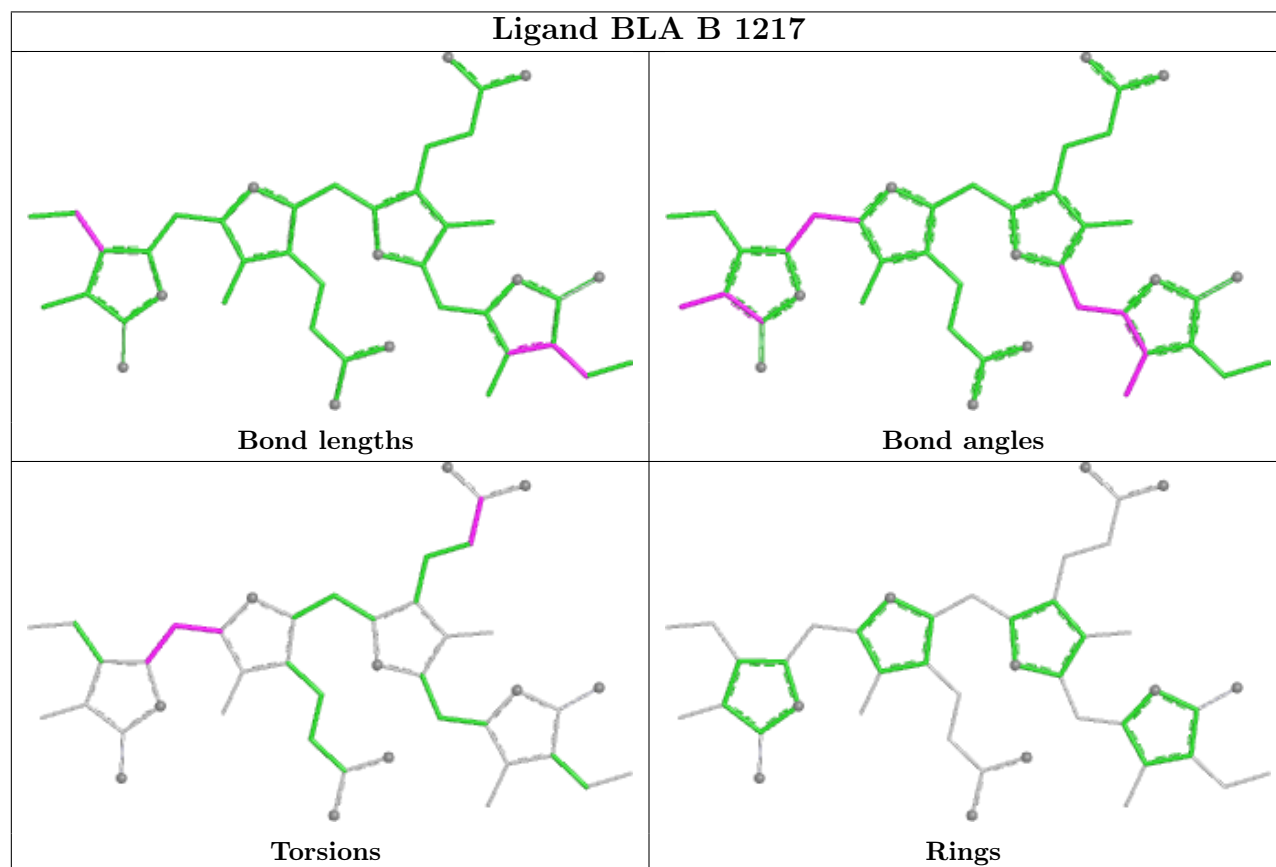
There are no ring outliers.

2 monomers are involved in 7 short contacts:

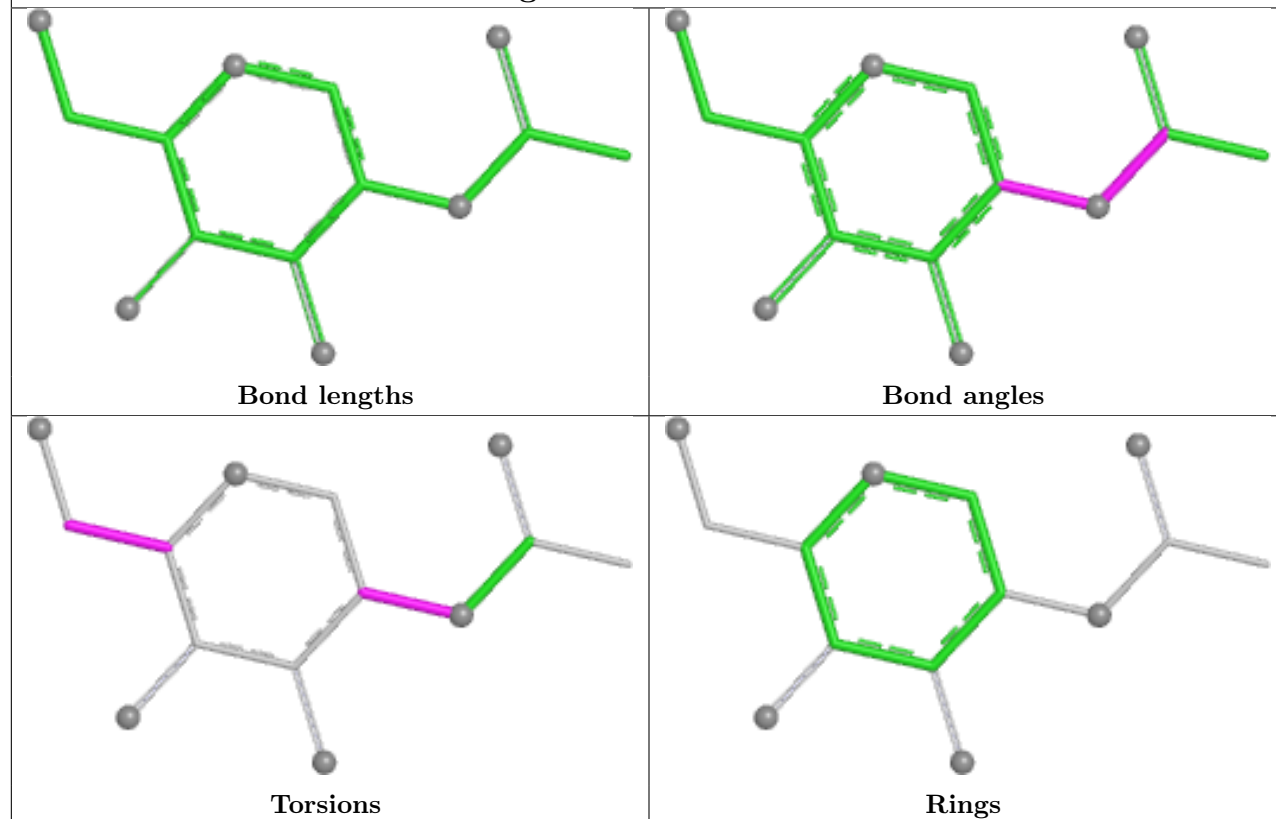
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1218	BLA	4	0
6	C	1217	BLA	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

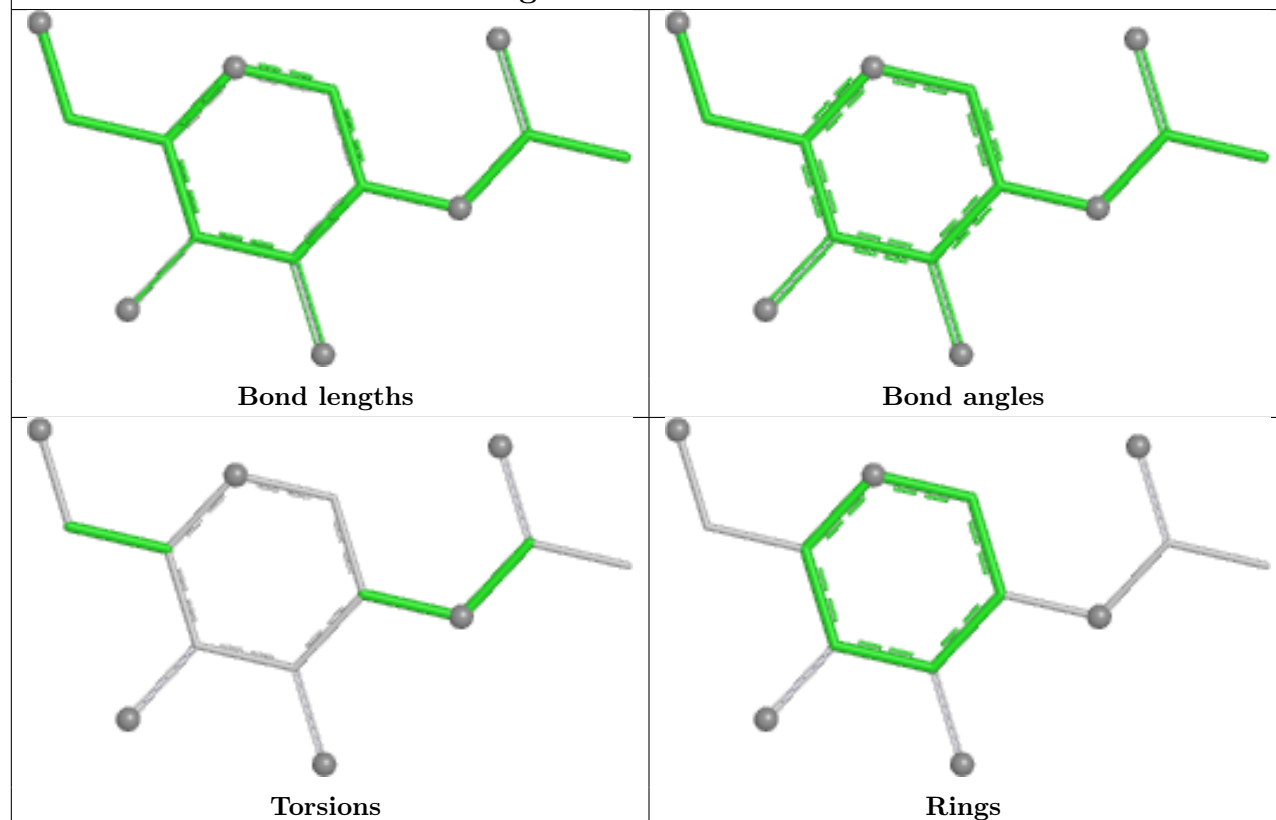
average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



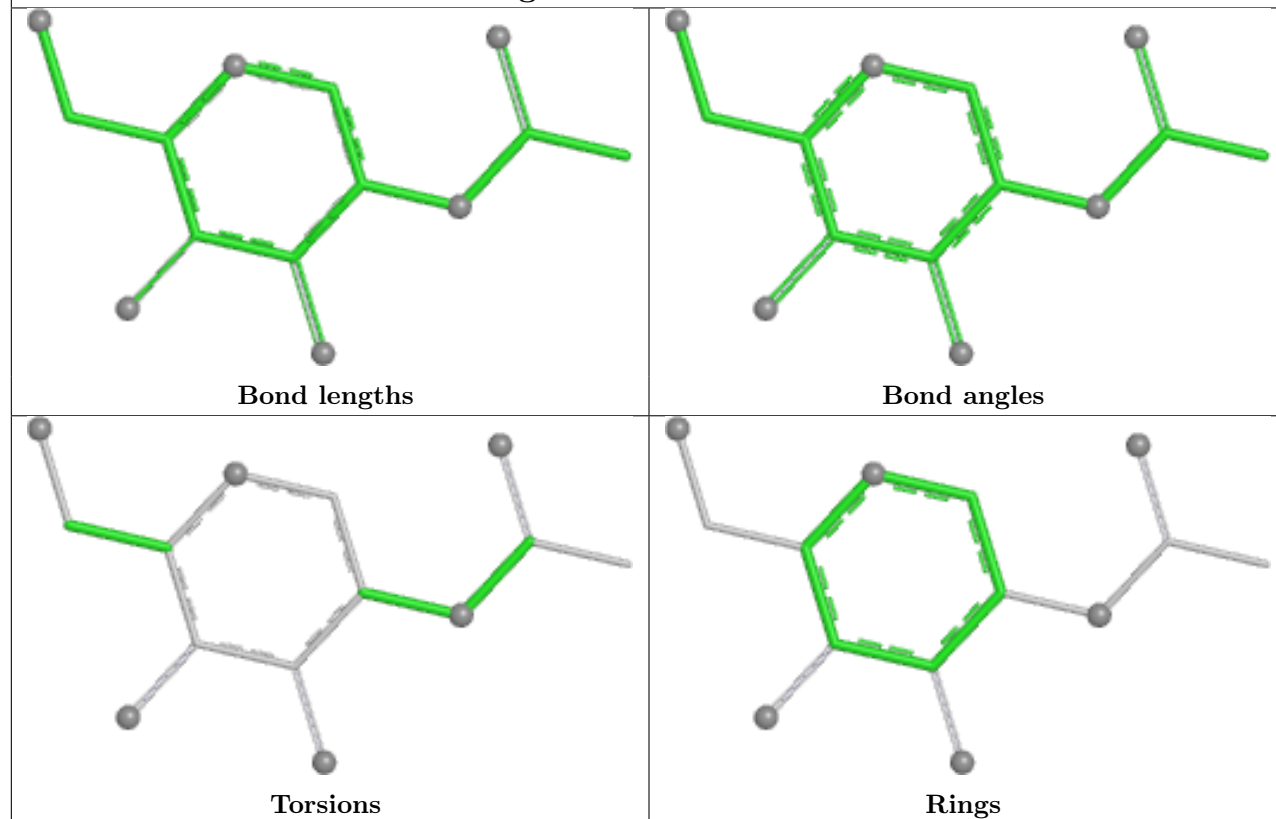
Ligand NAG C 1218



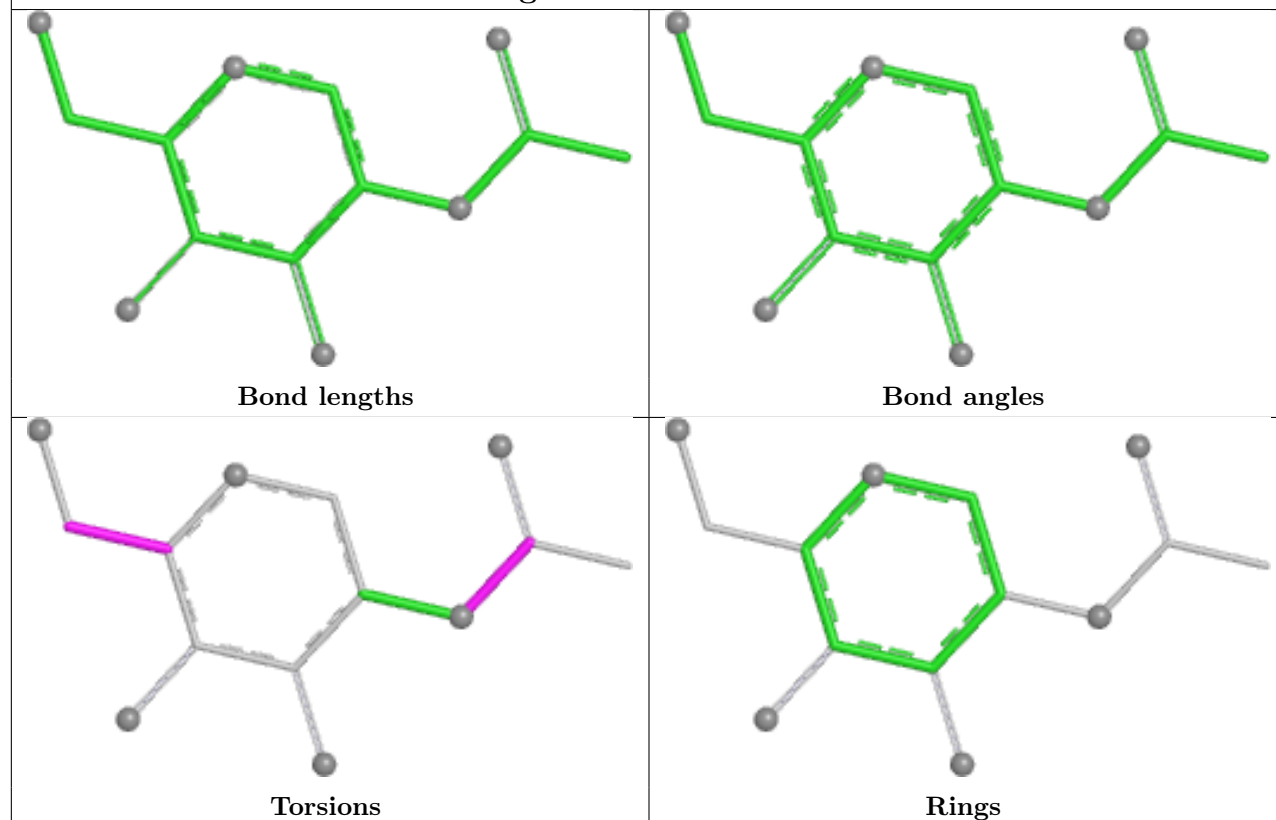
Ligand NAG C 1204



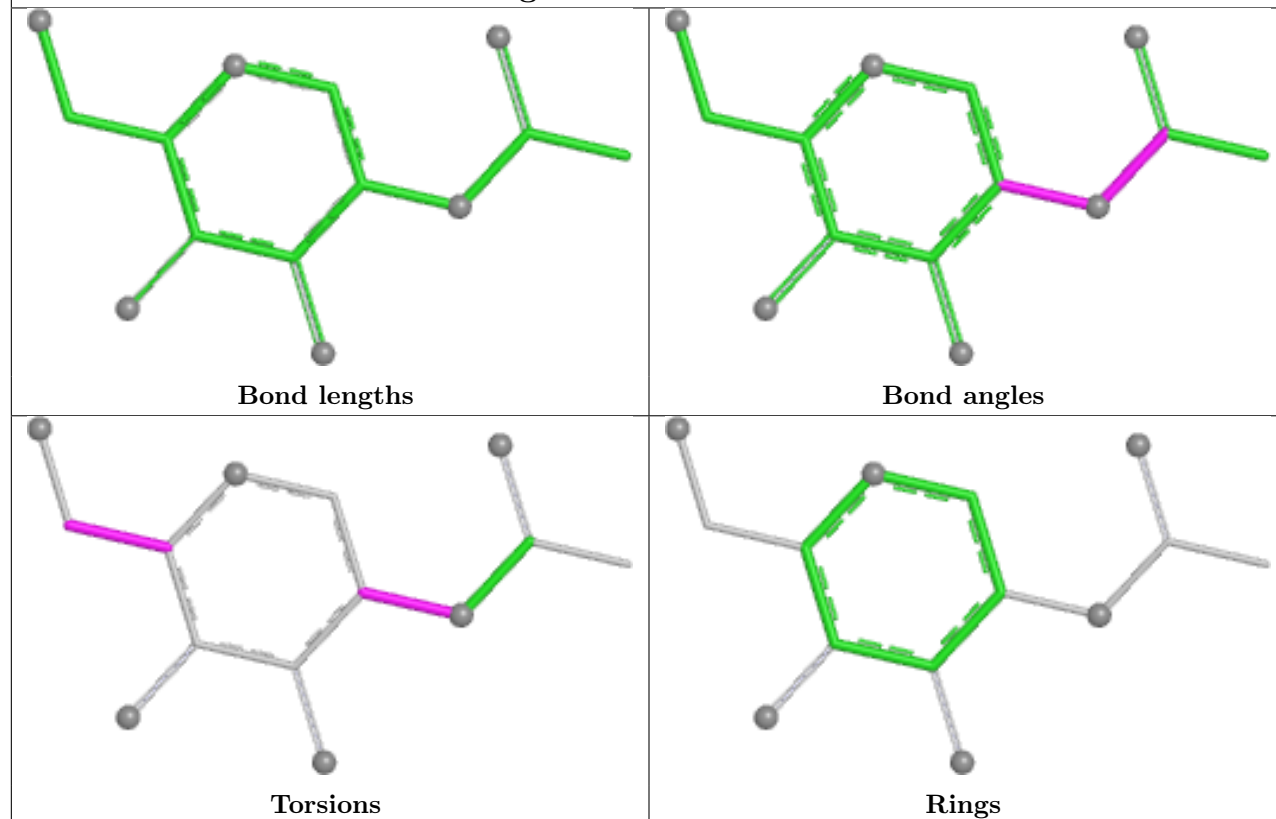
Ligand NAG A 1204



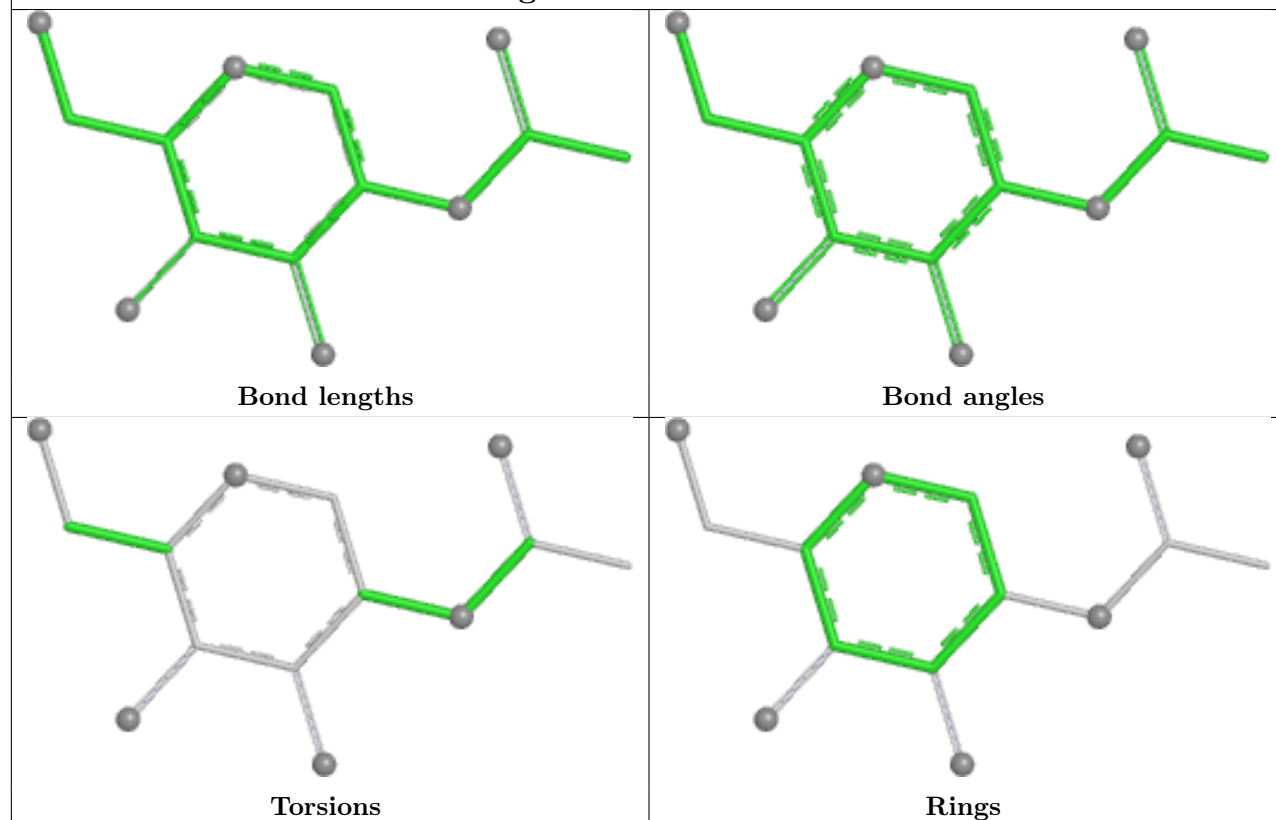
Ligand NAG C 1207

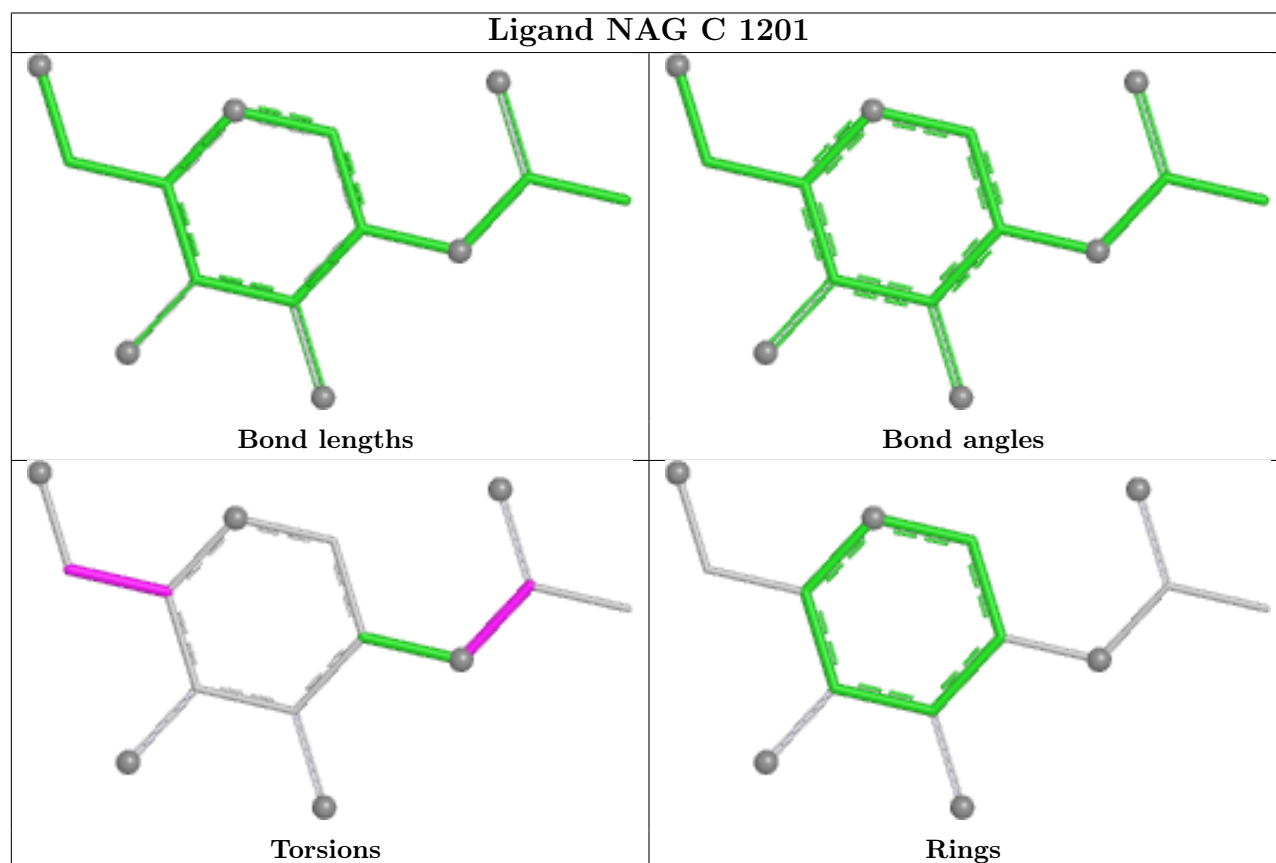
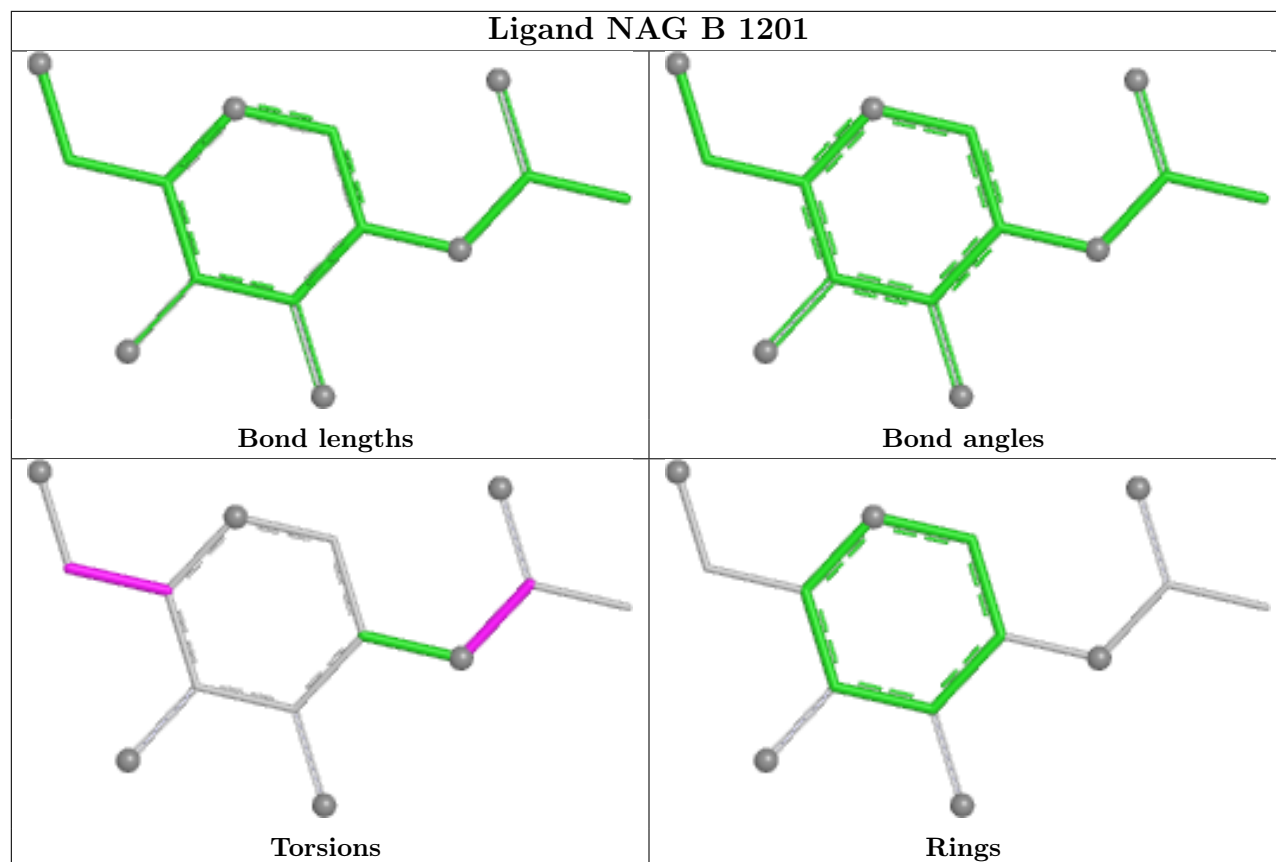


Ligand NAG C 1211

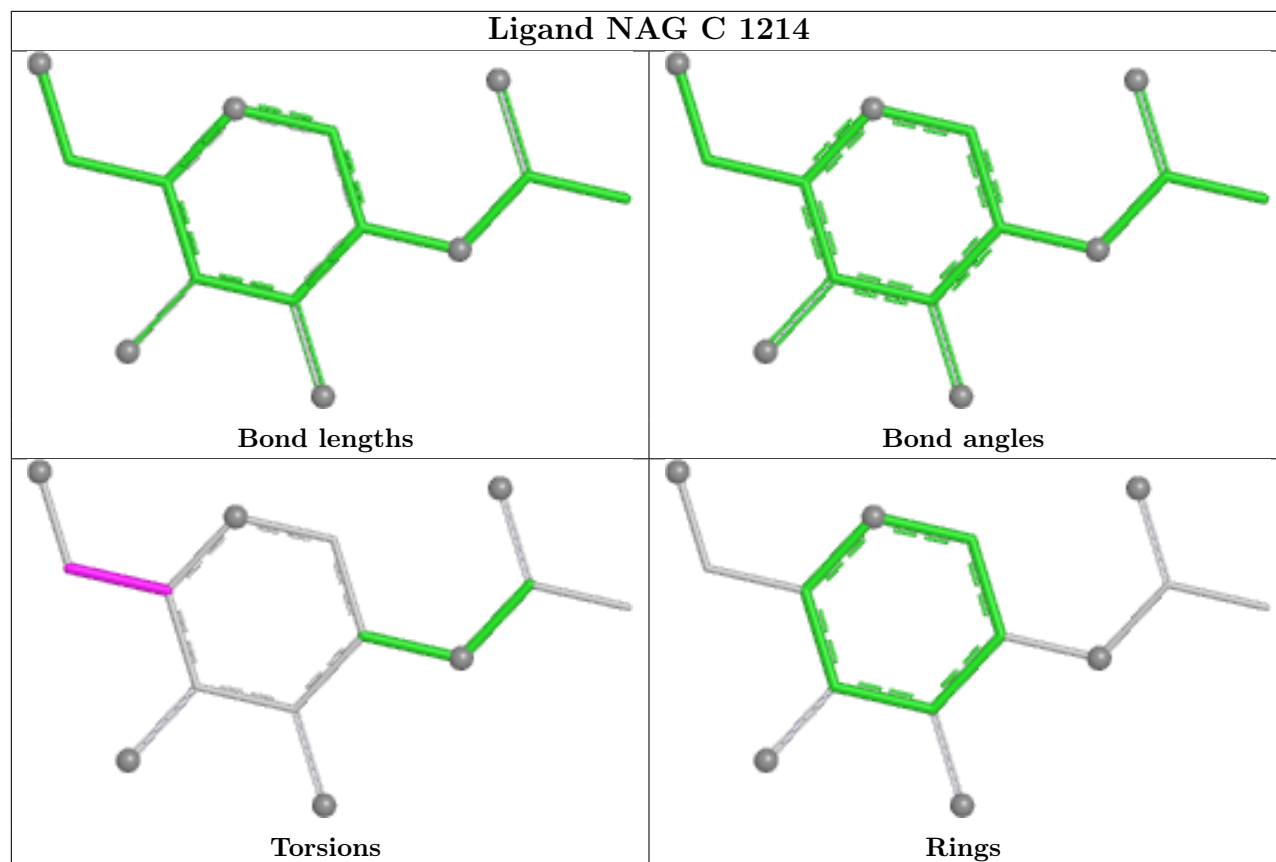


Ligand NAG A 1205

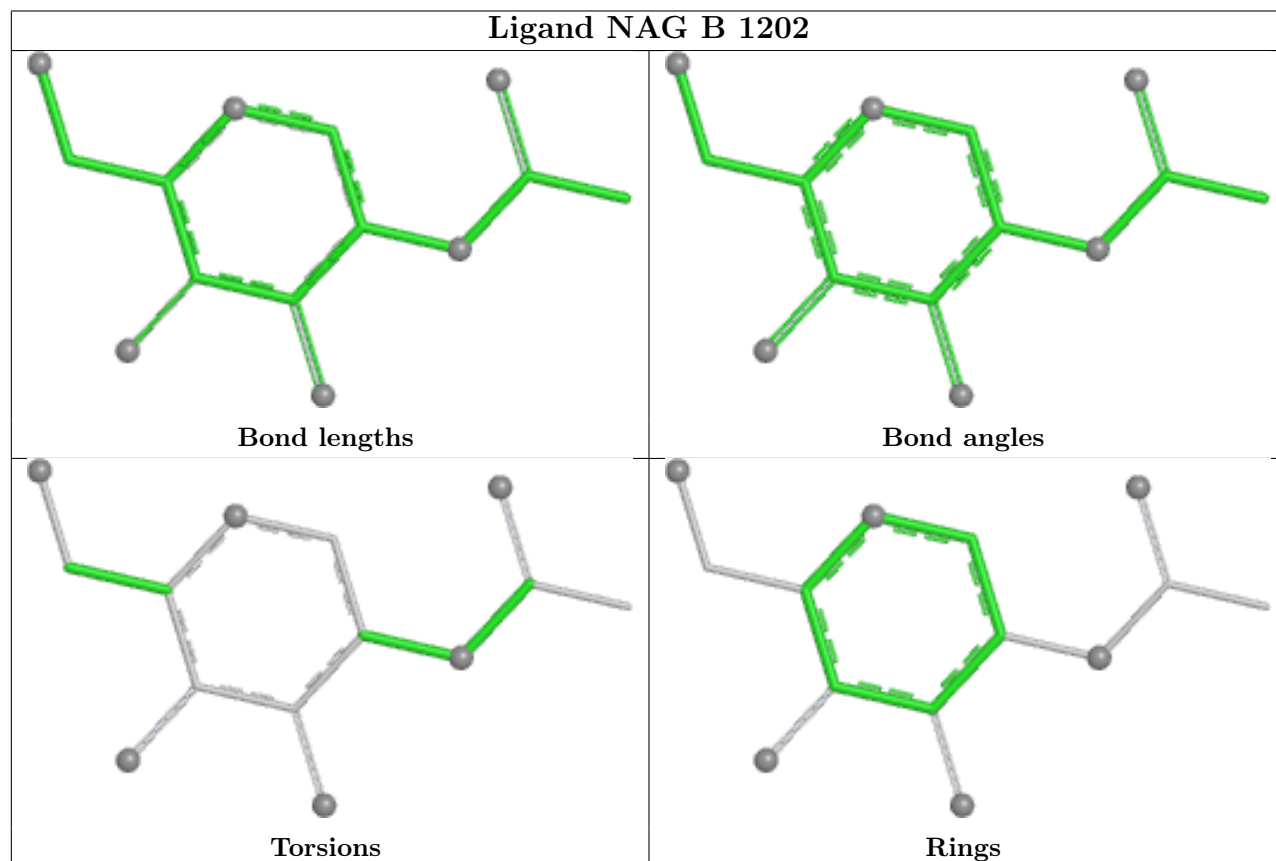




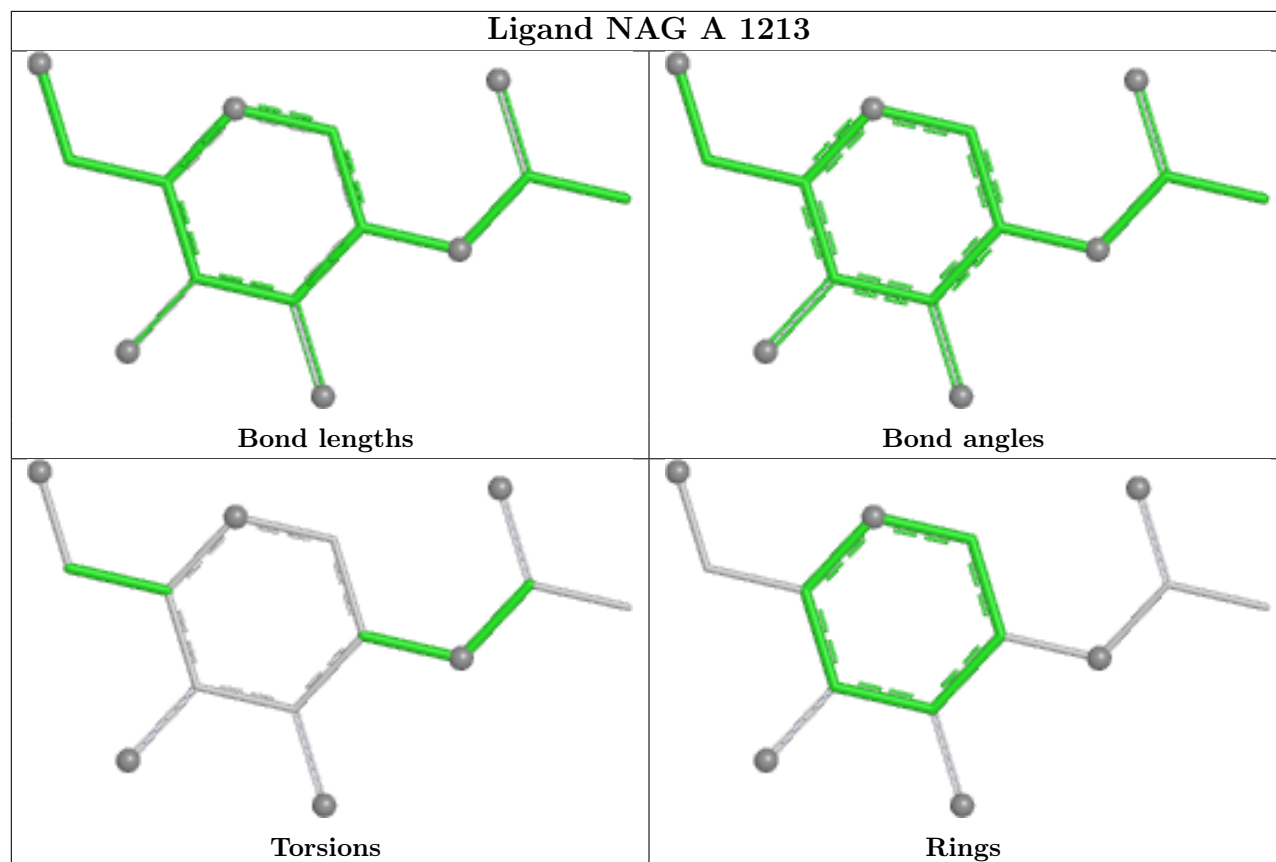
Ligand NAG C 1214



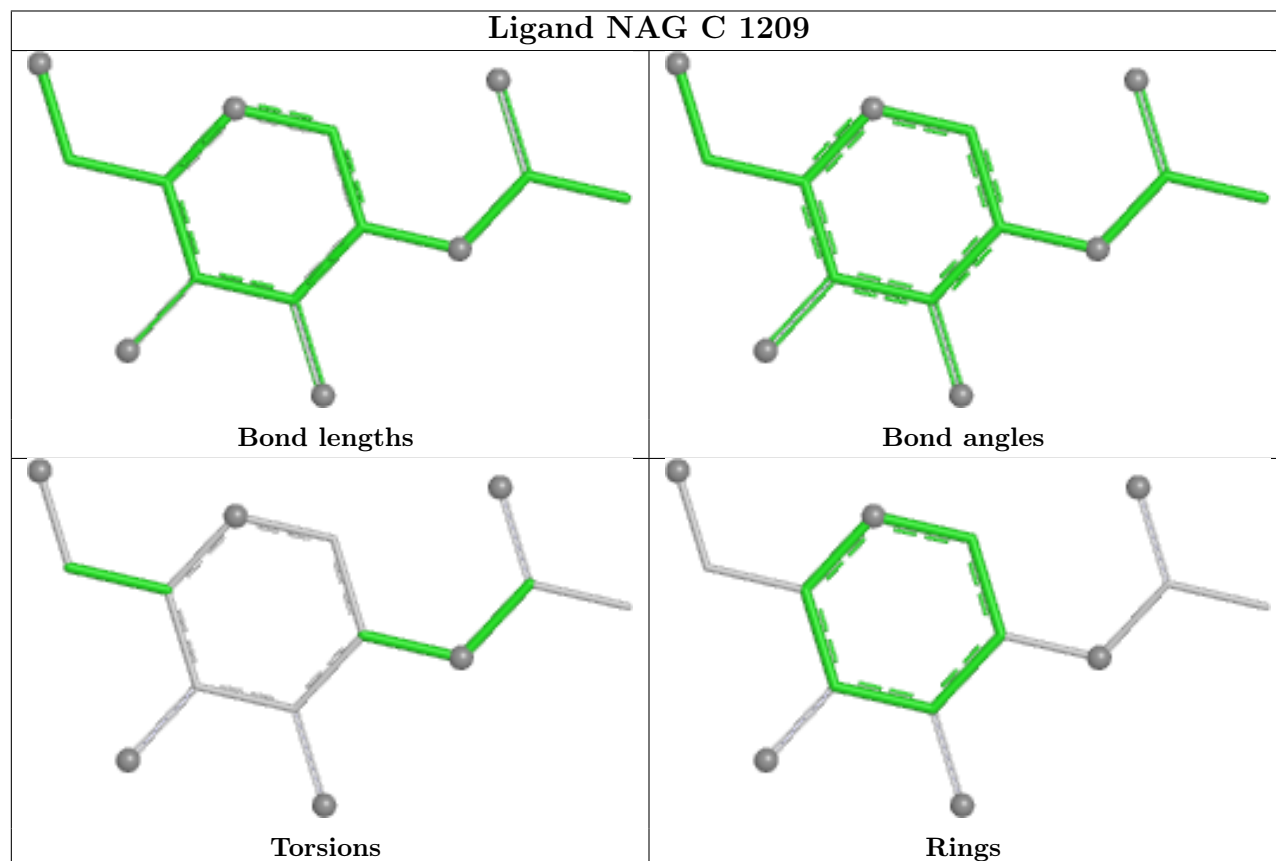
Ligand NAG B 1202

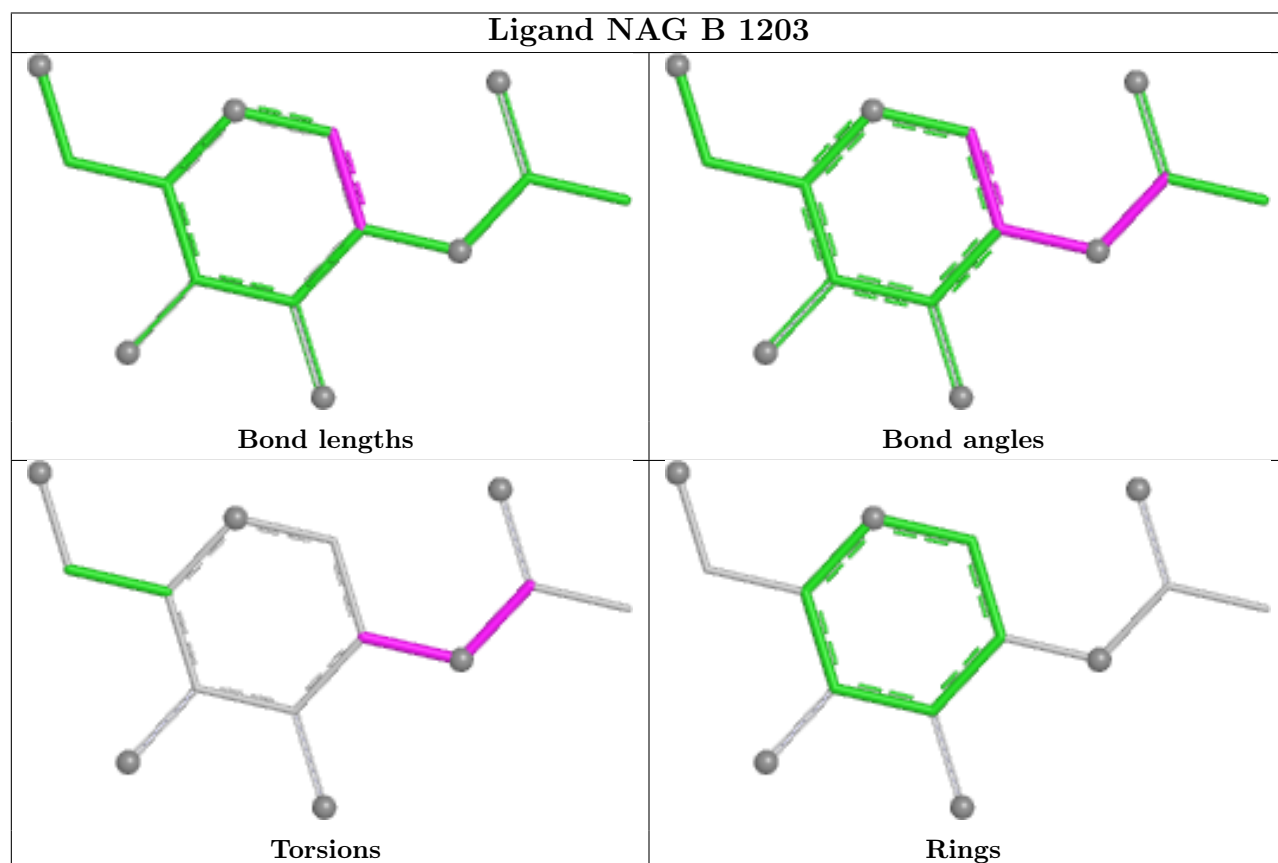
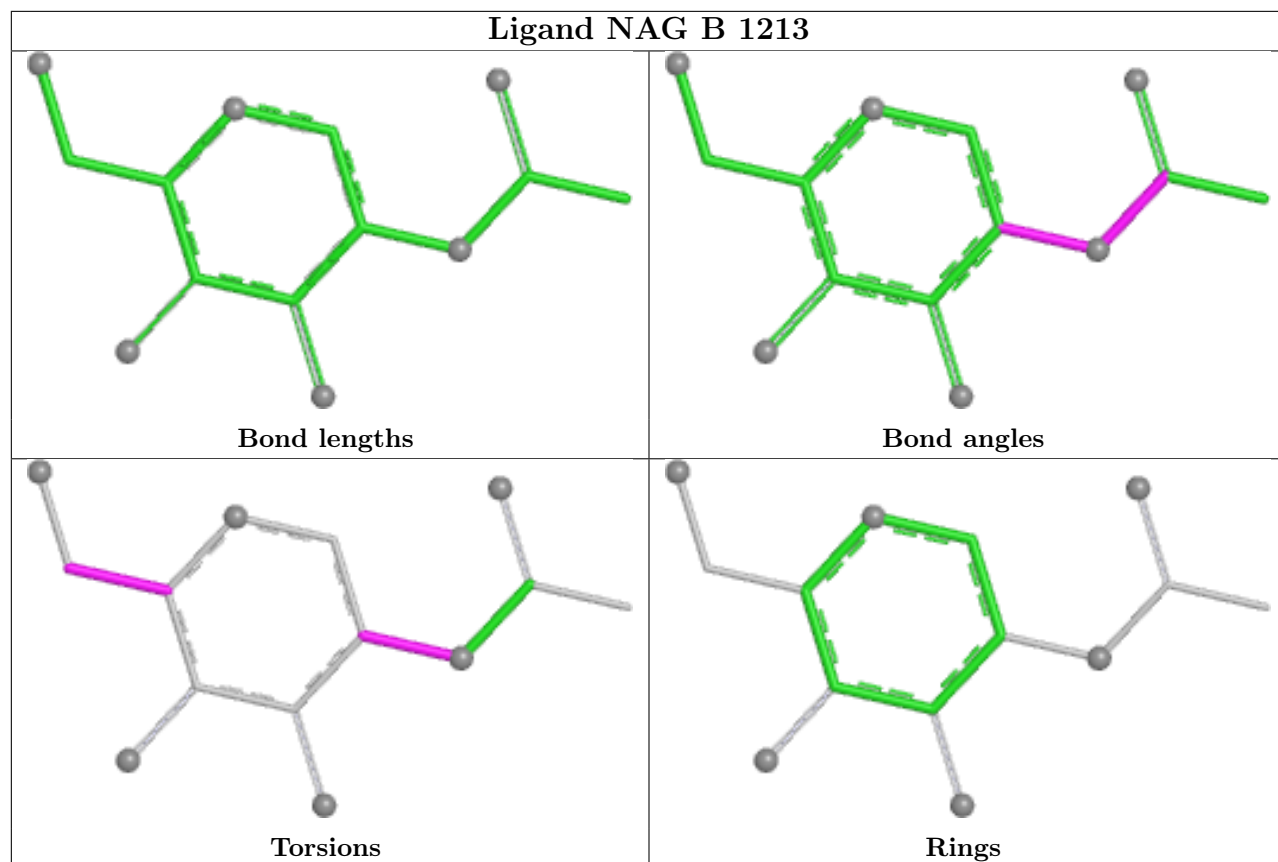


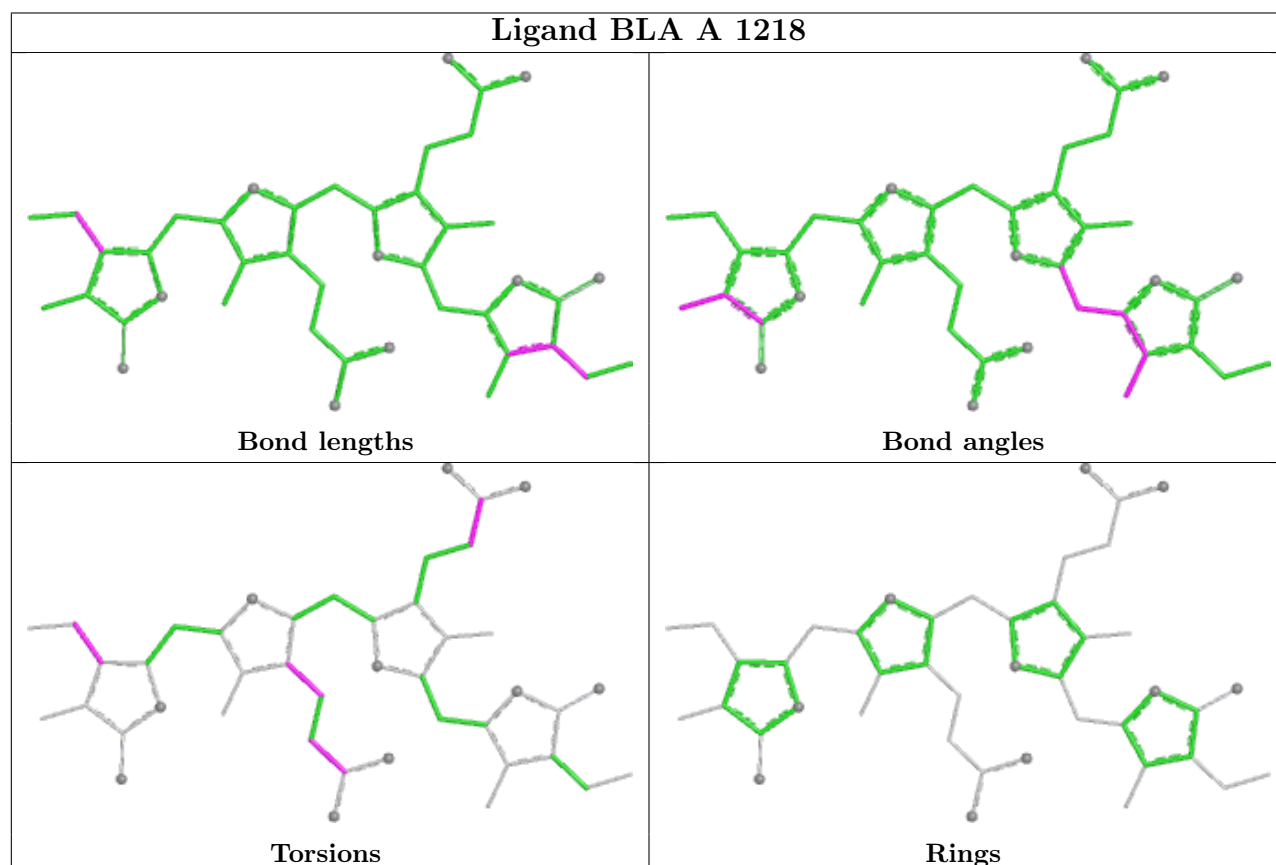
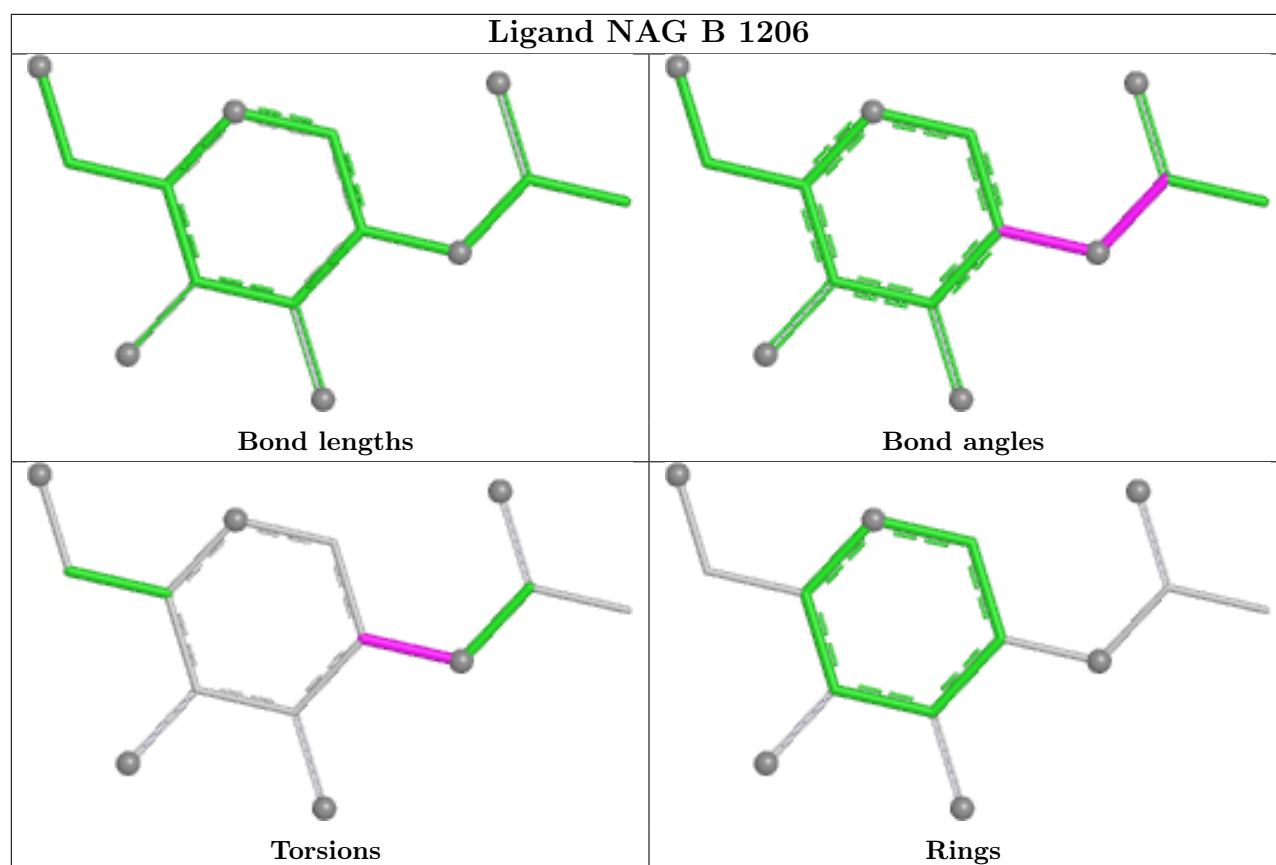
Ligand NAG A 1213



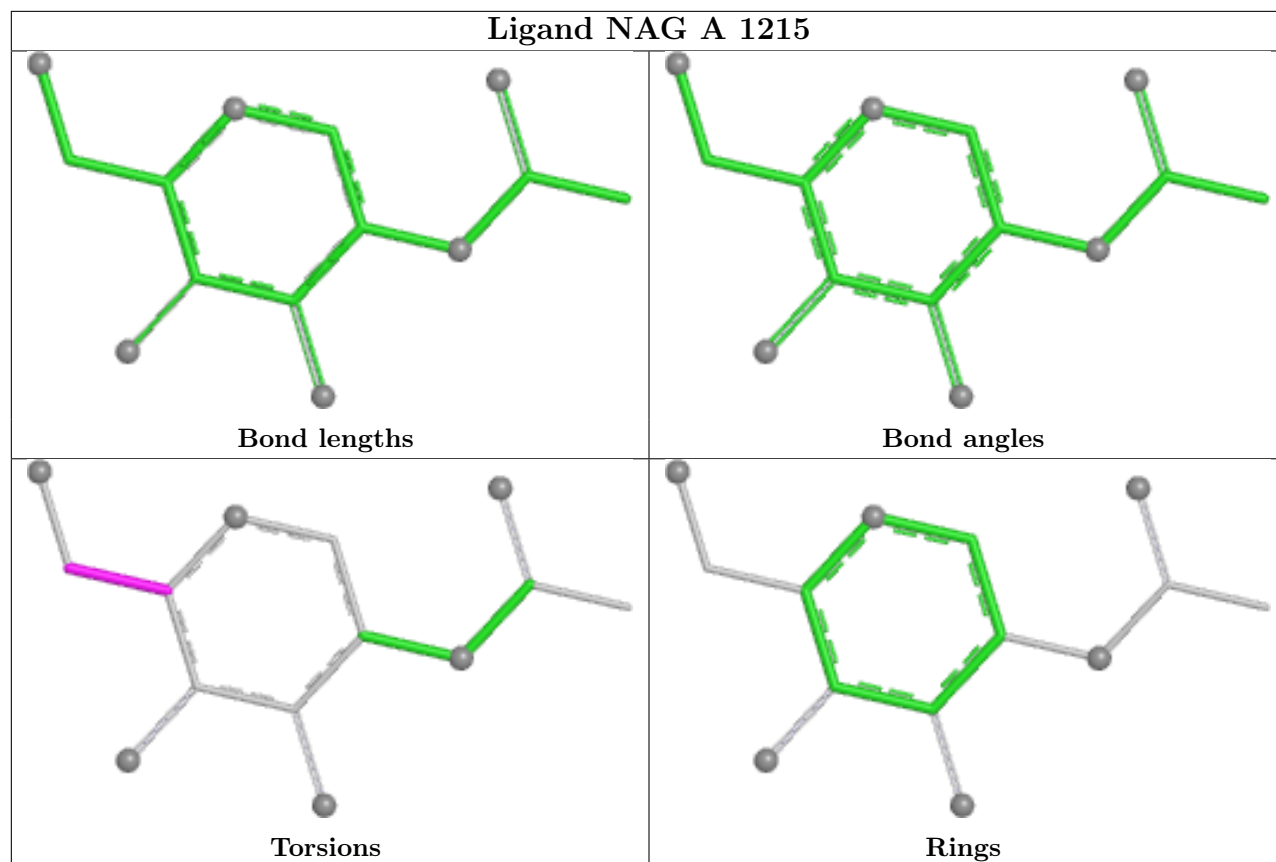
Ligand NAG C 1209



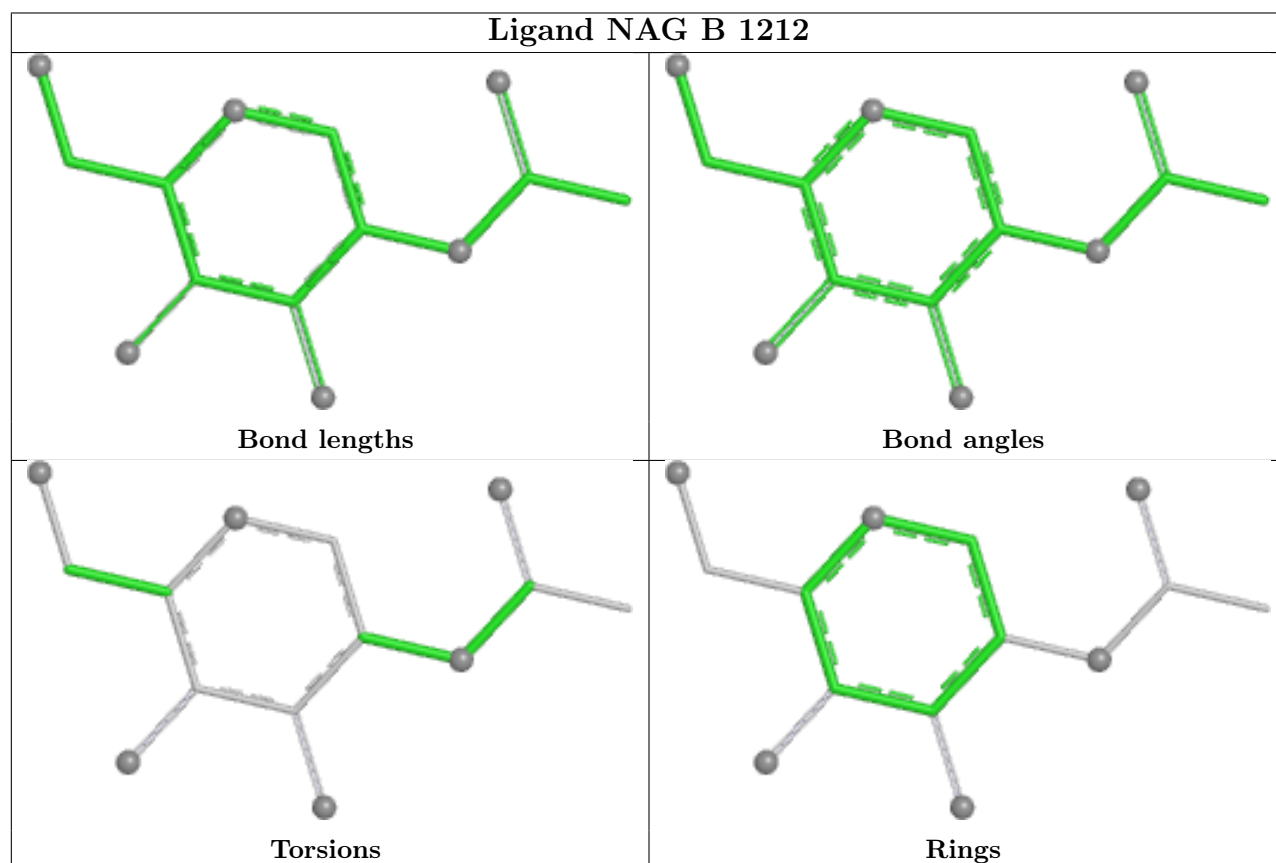


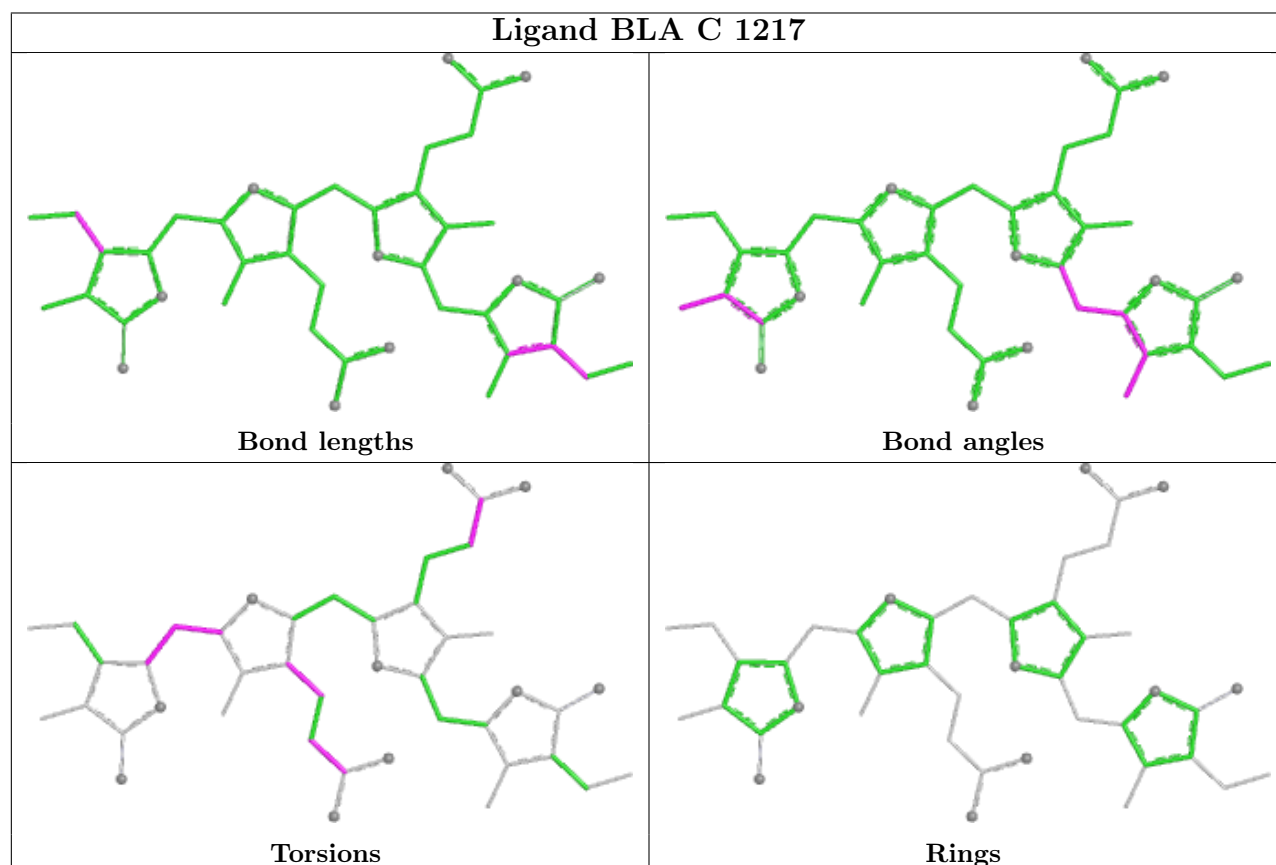
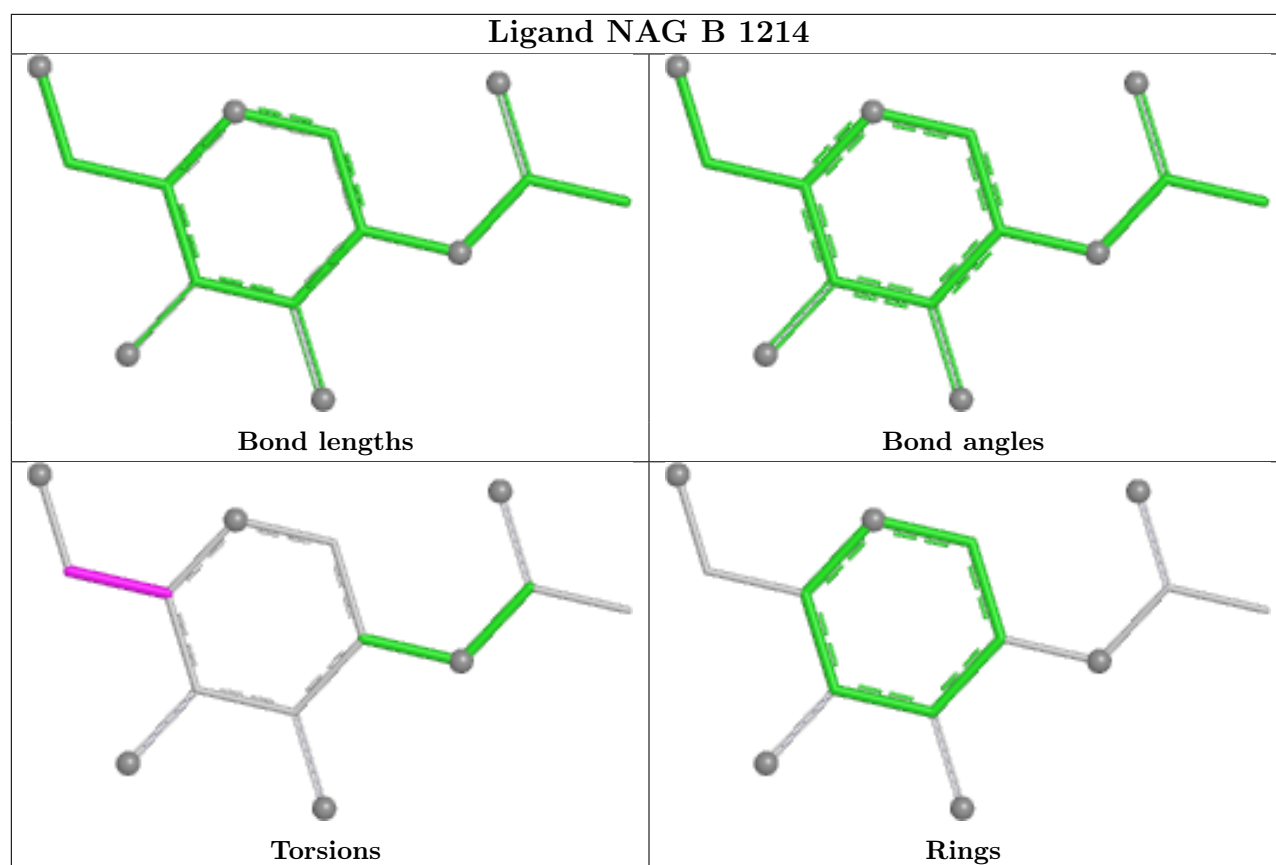


Ligand NAG A 1215

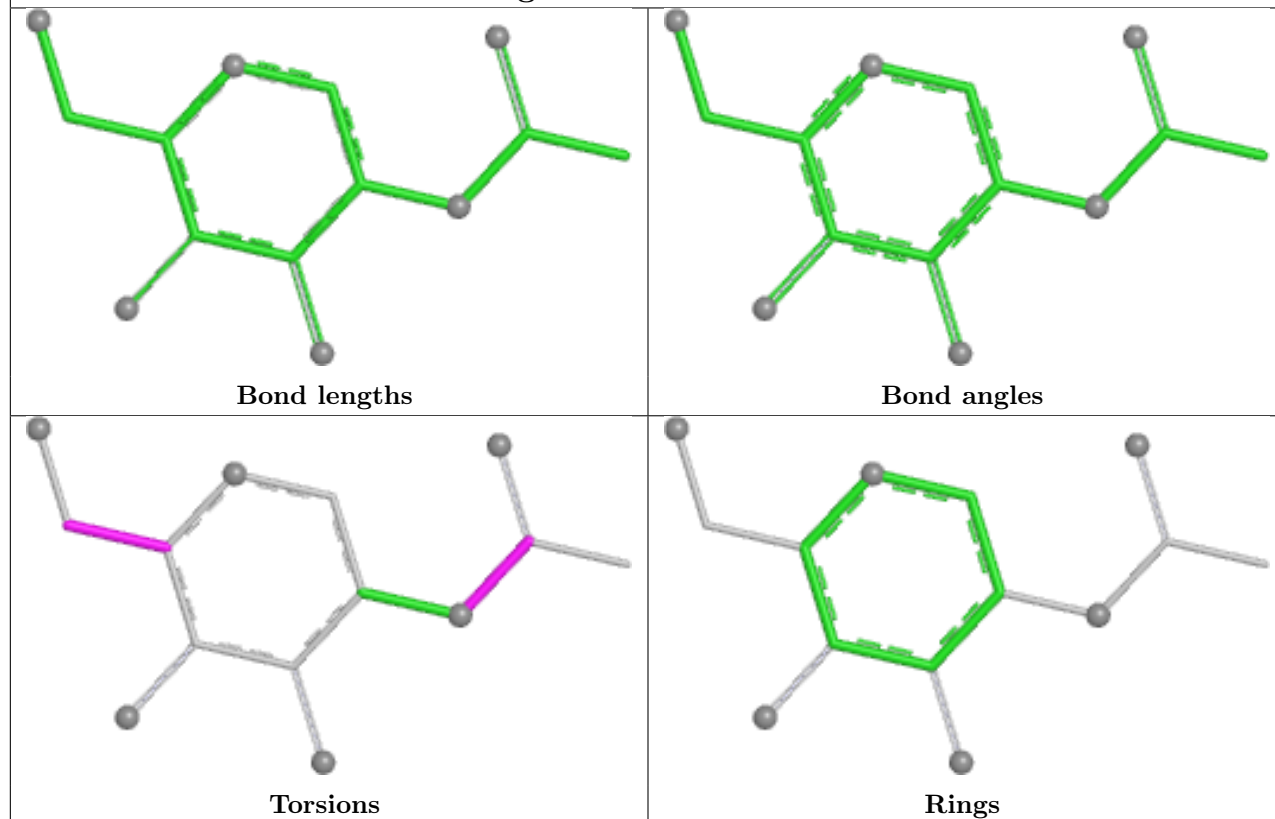


Ligand NAG B 1212

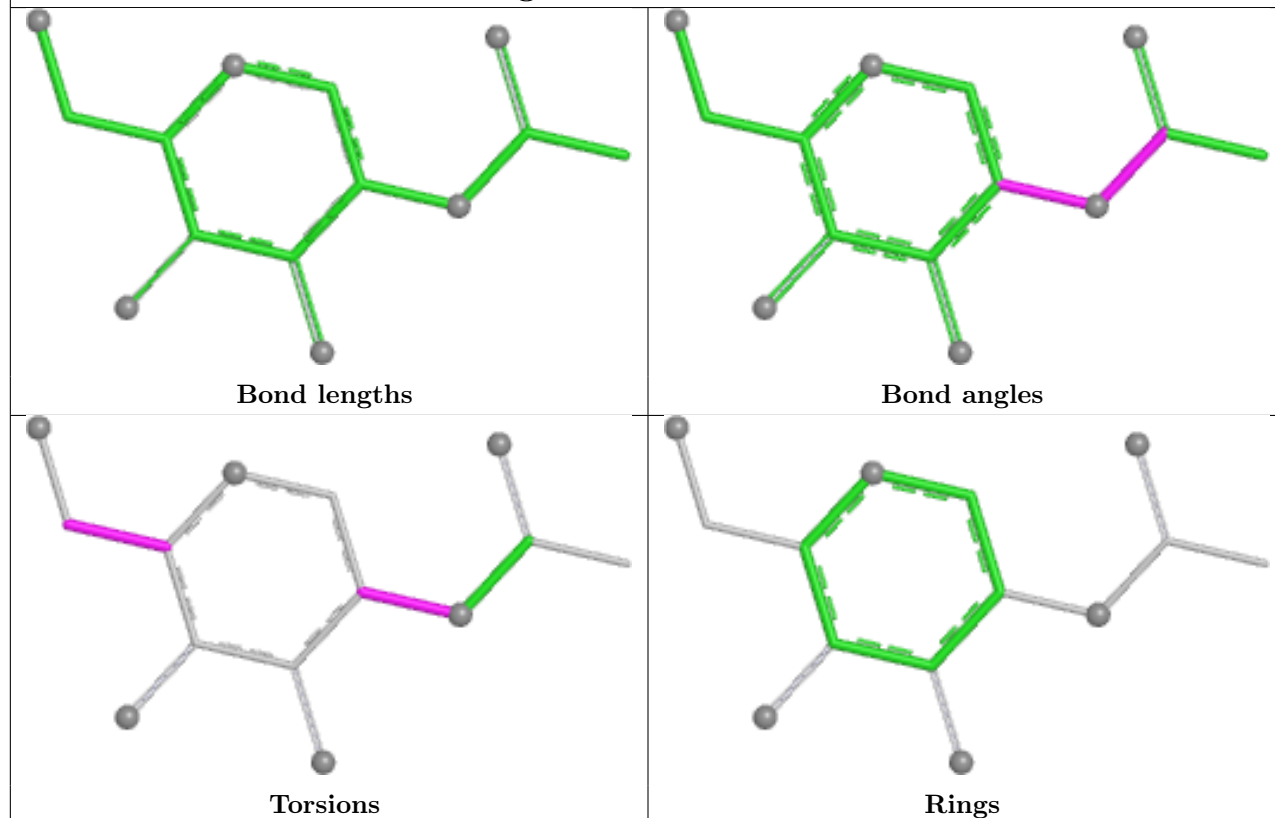


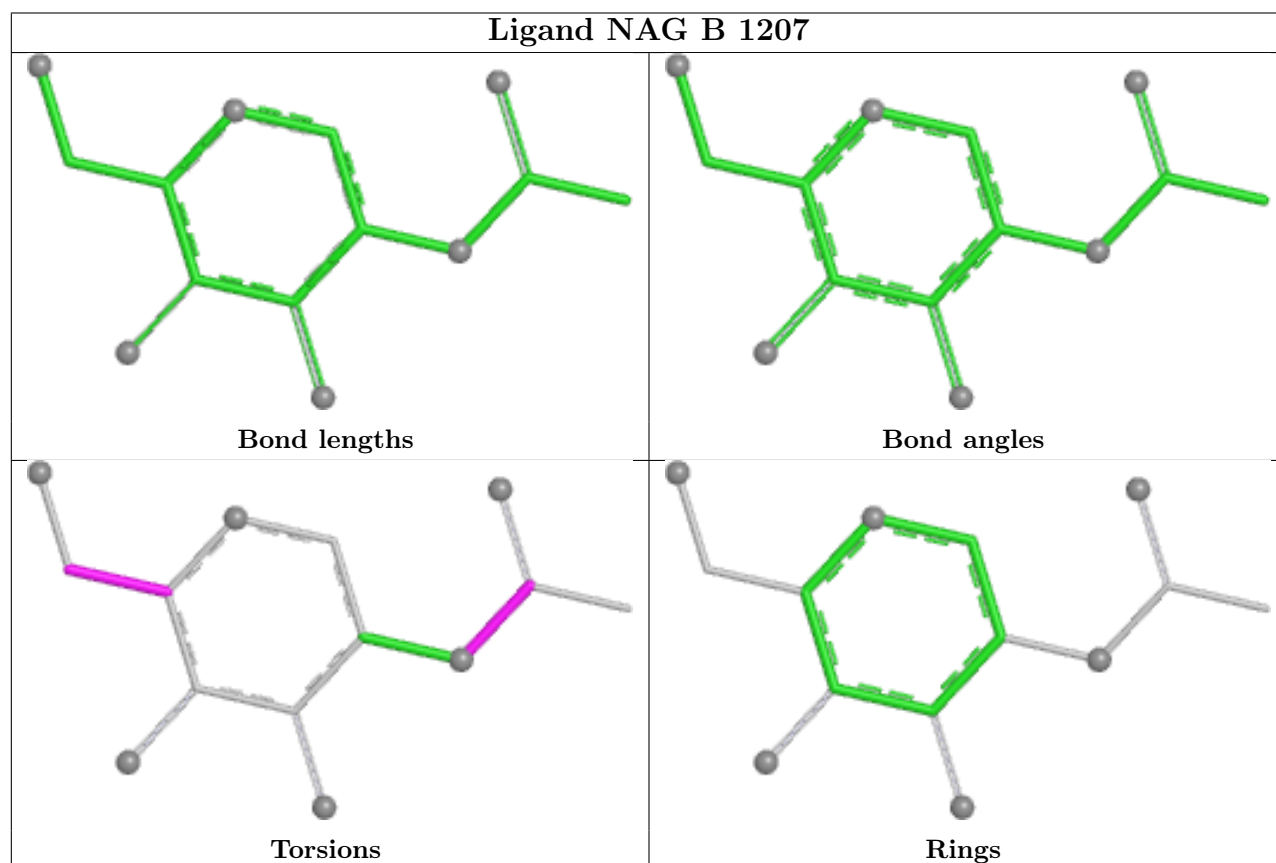
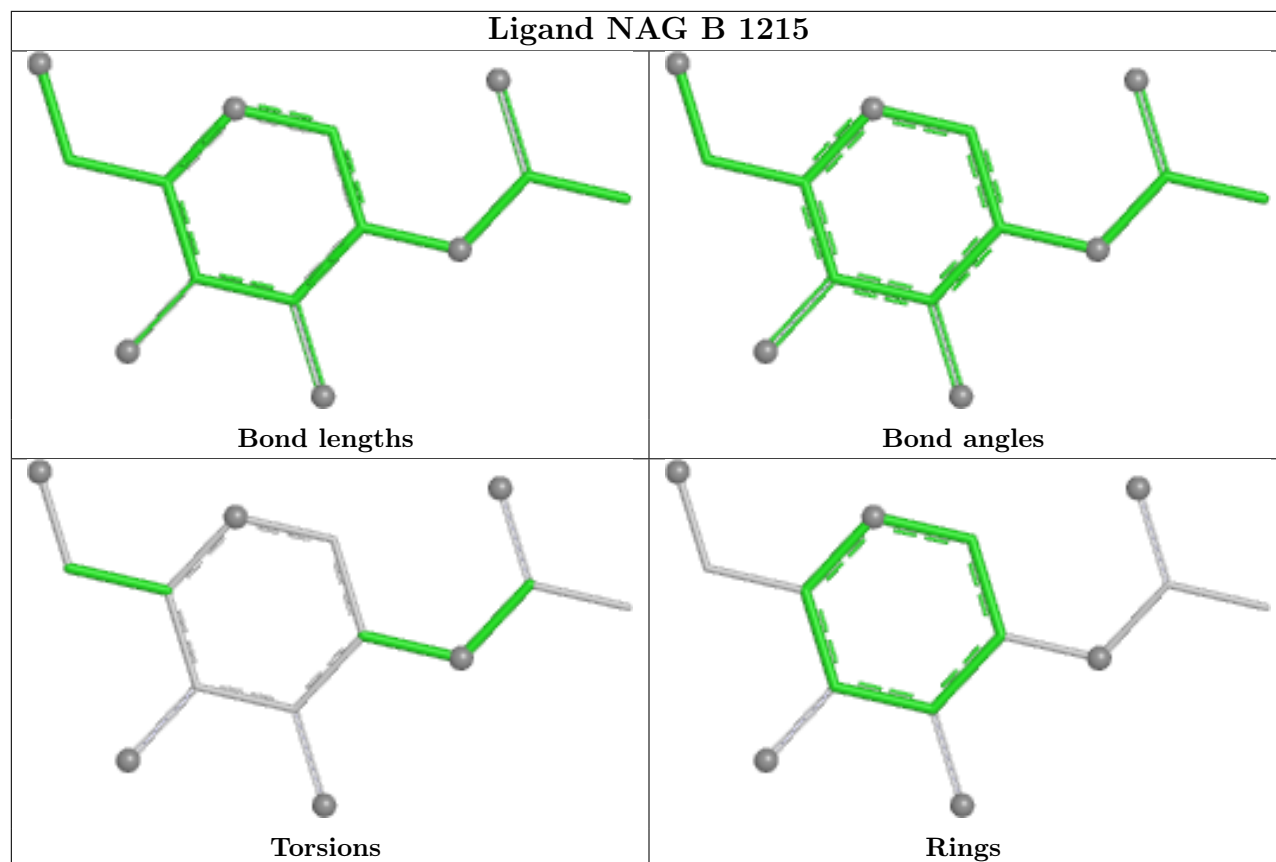


Ligand NAG A 1208

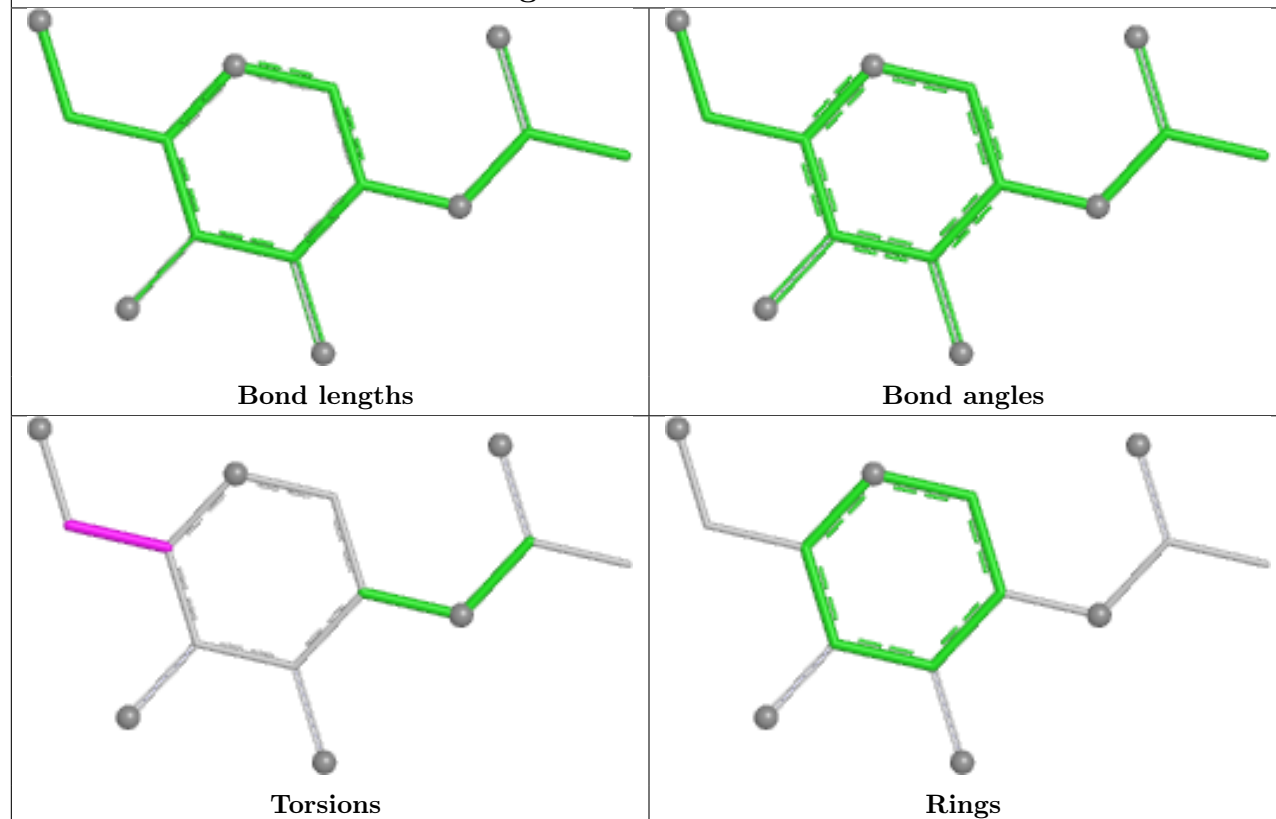


Ligand NAG A 1212

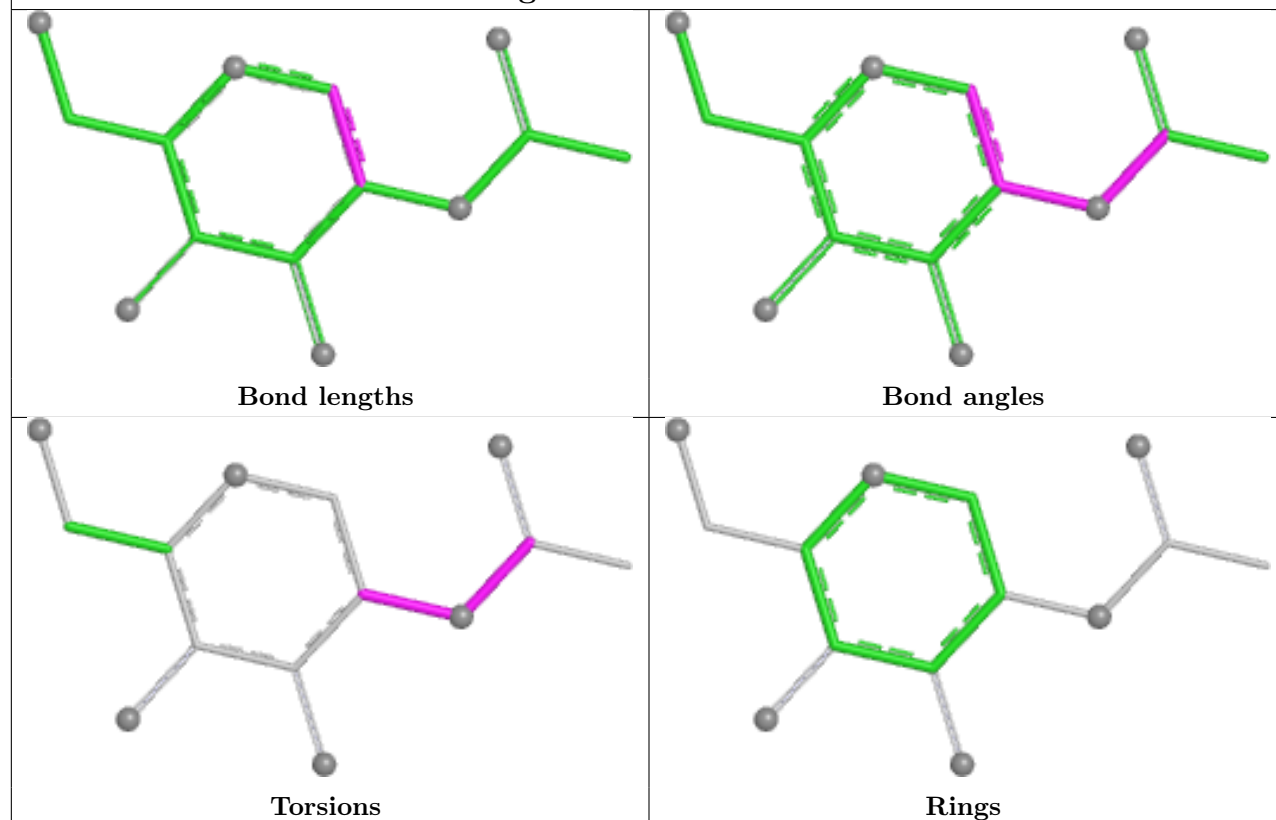




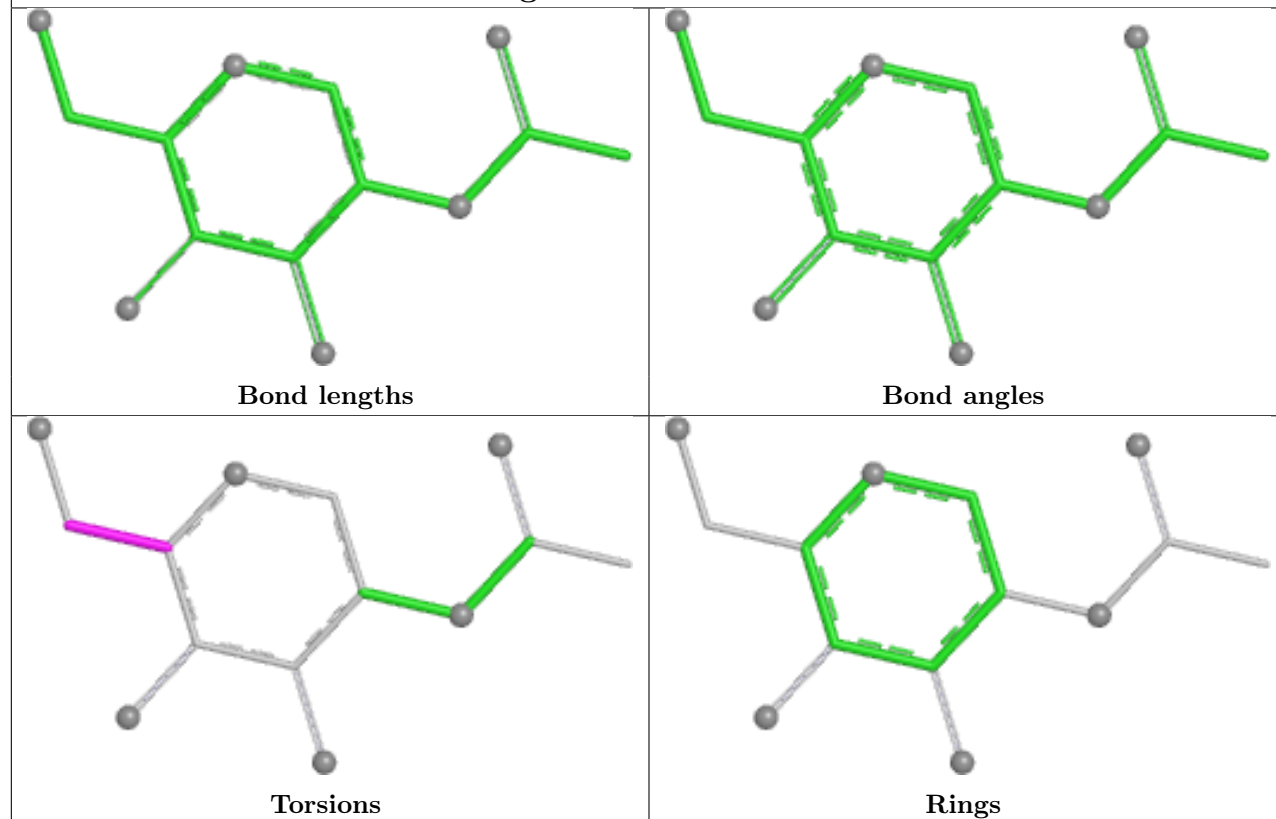
Ligand NAG C 1216



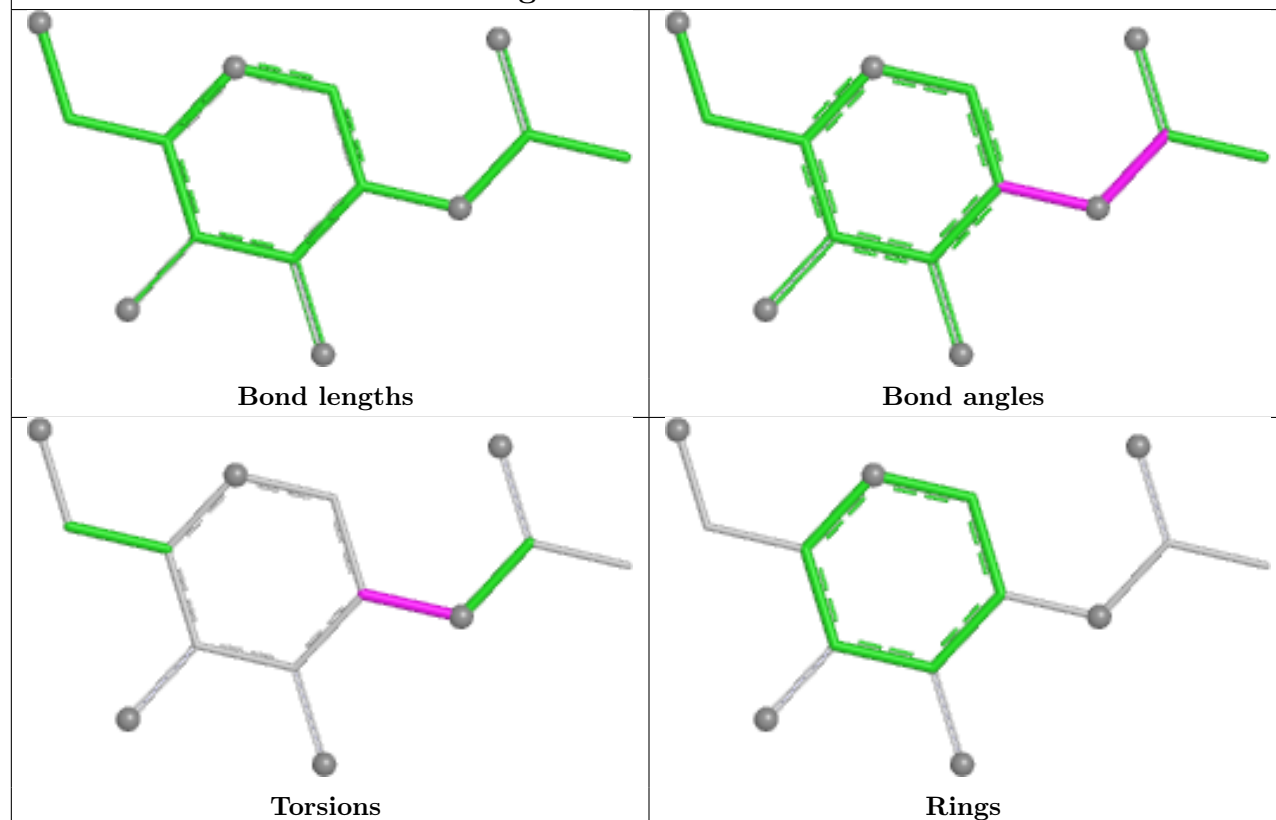
Ligand NAG A 1203



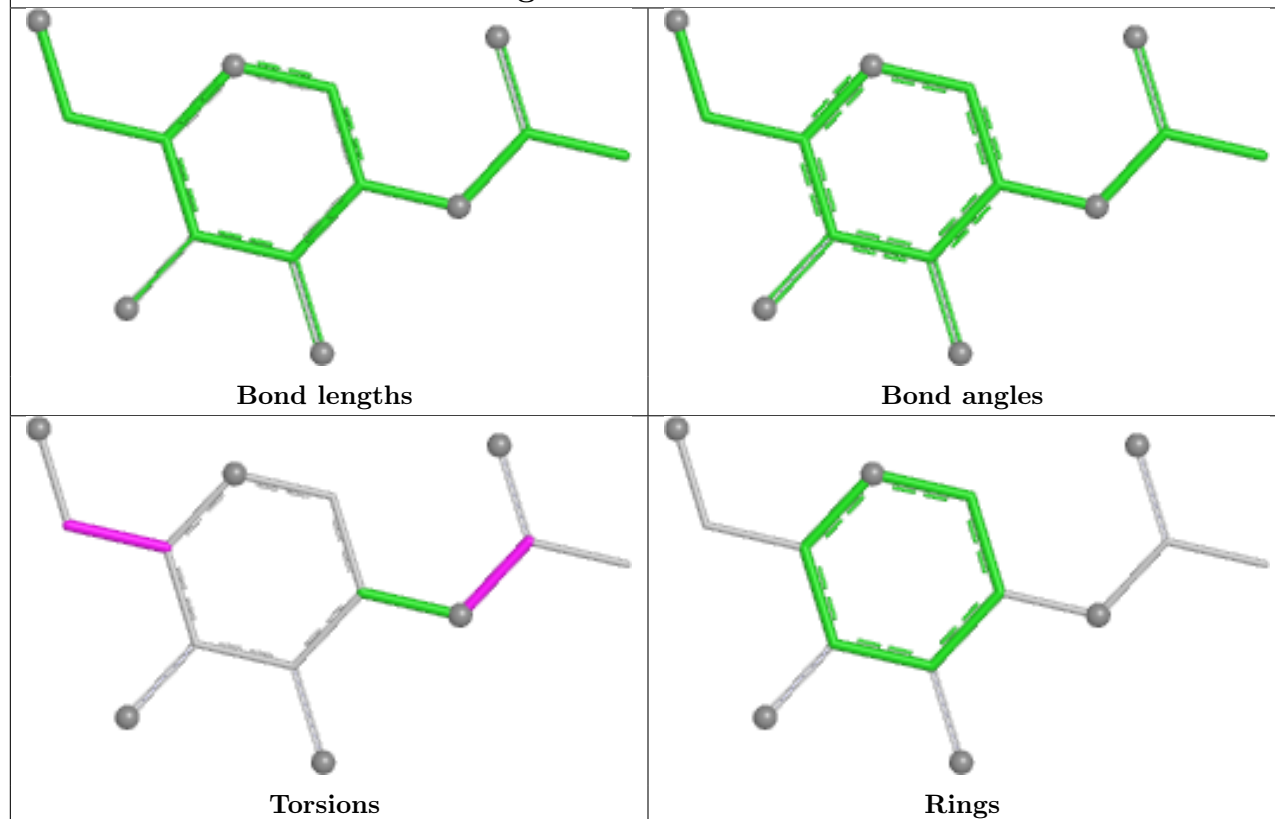
Ligand NAG B 1205



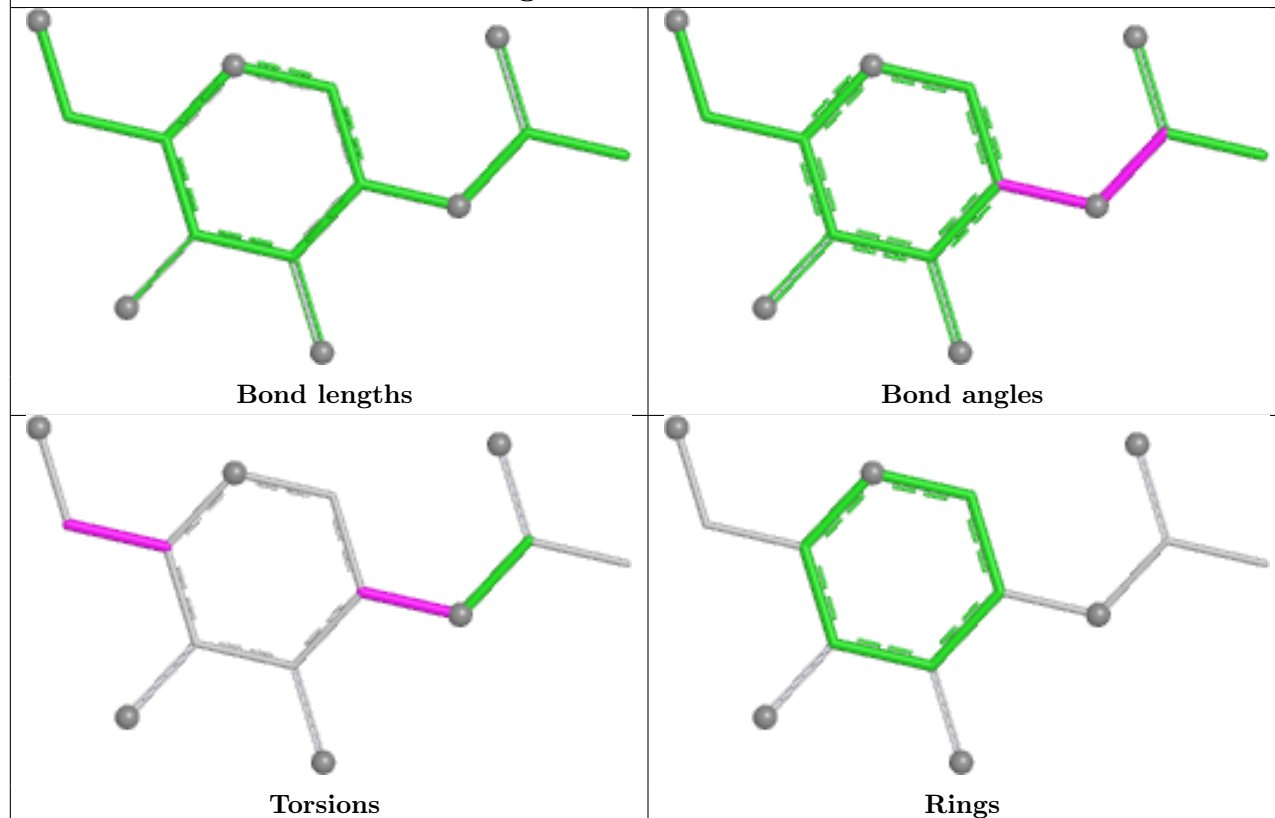
Ligand NAG A 1207



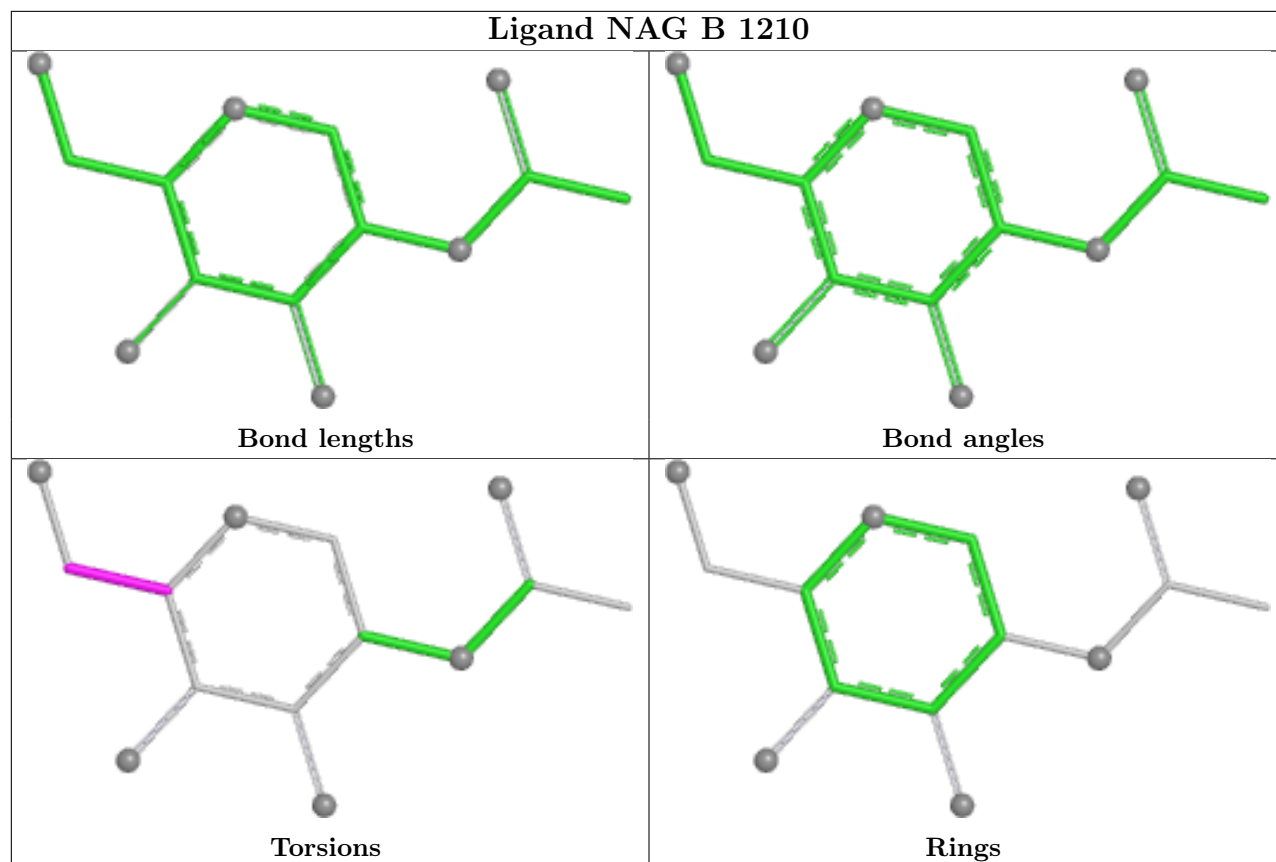
Ligand NAG A 1201



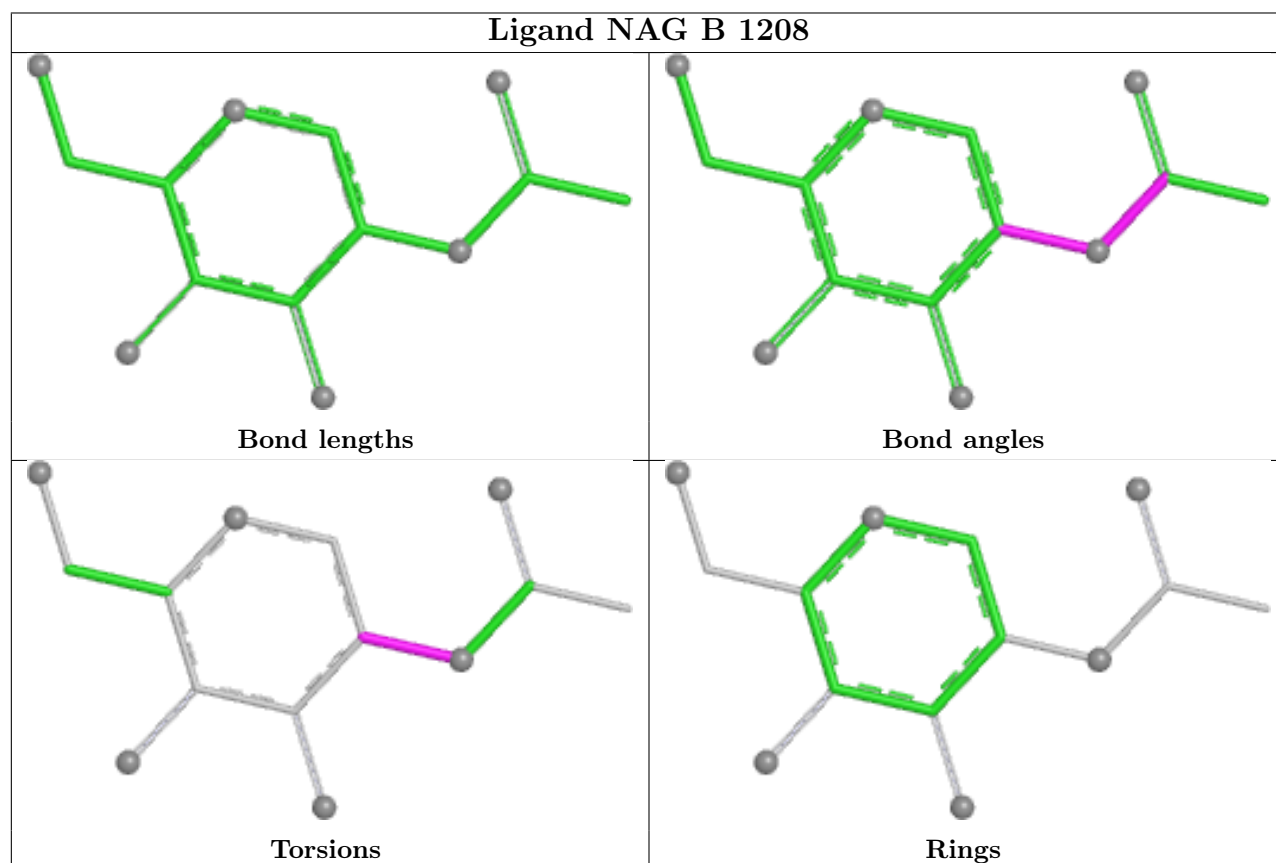
Ligand NAG C 1208

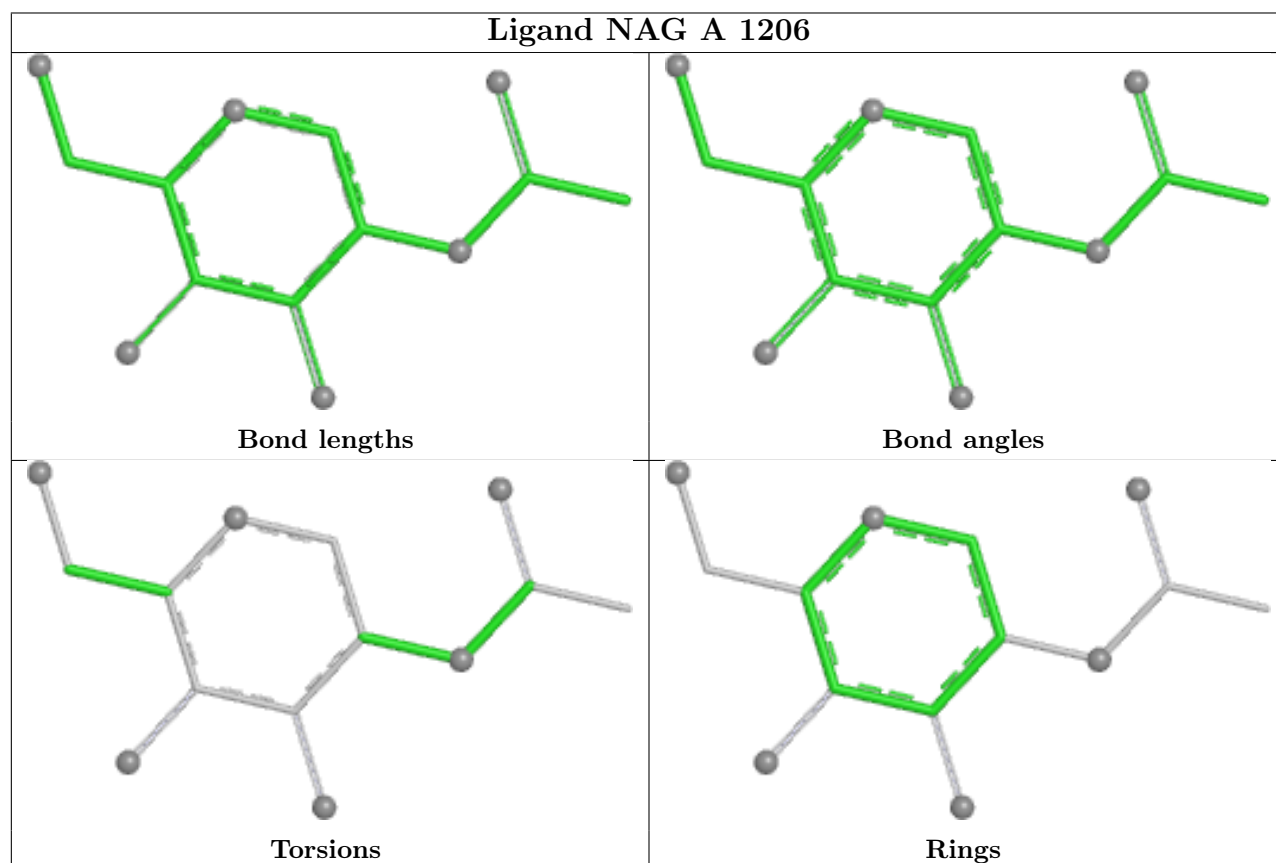
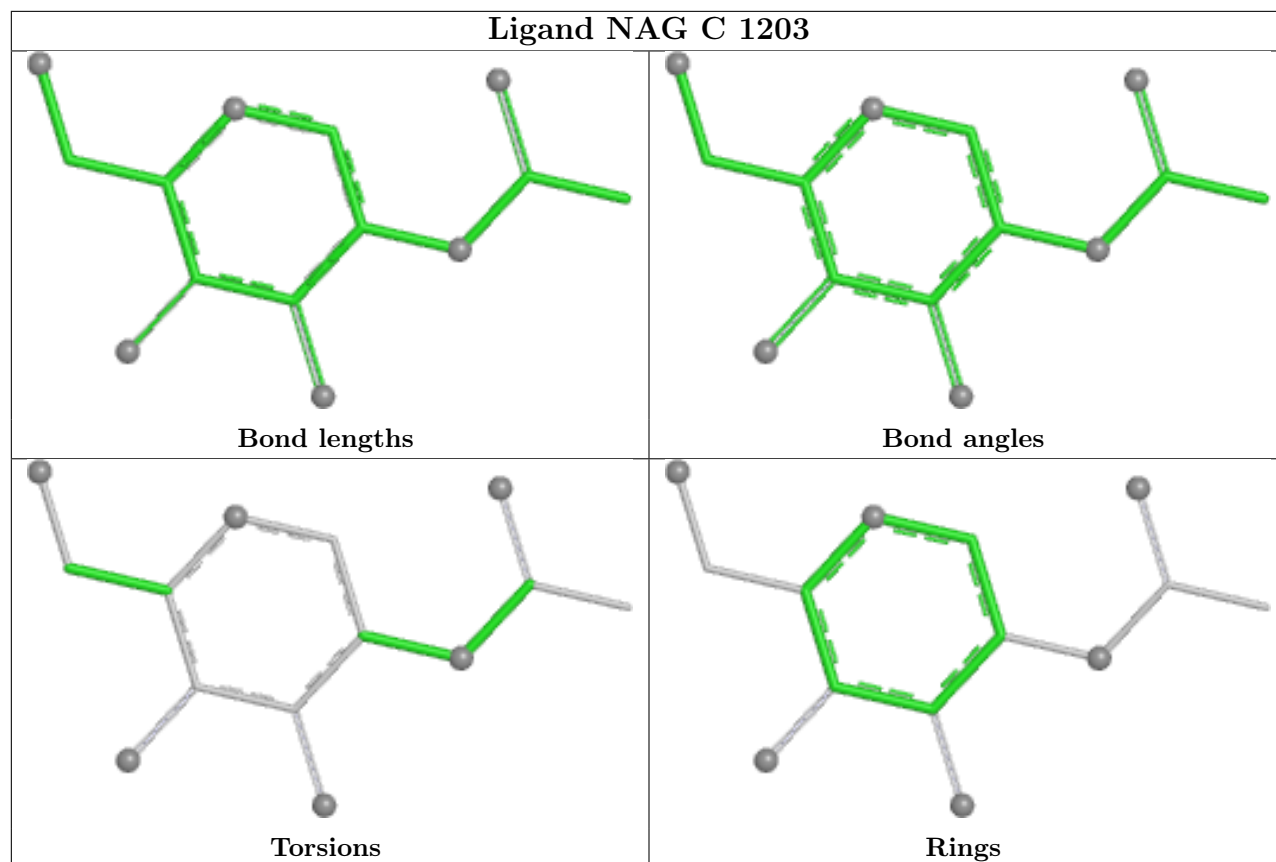


Ligand NAG B 1210

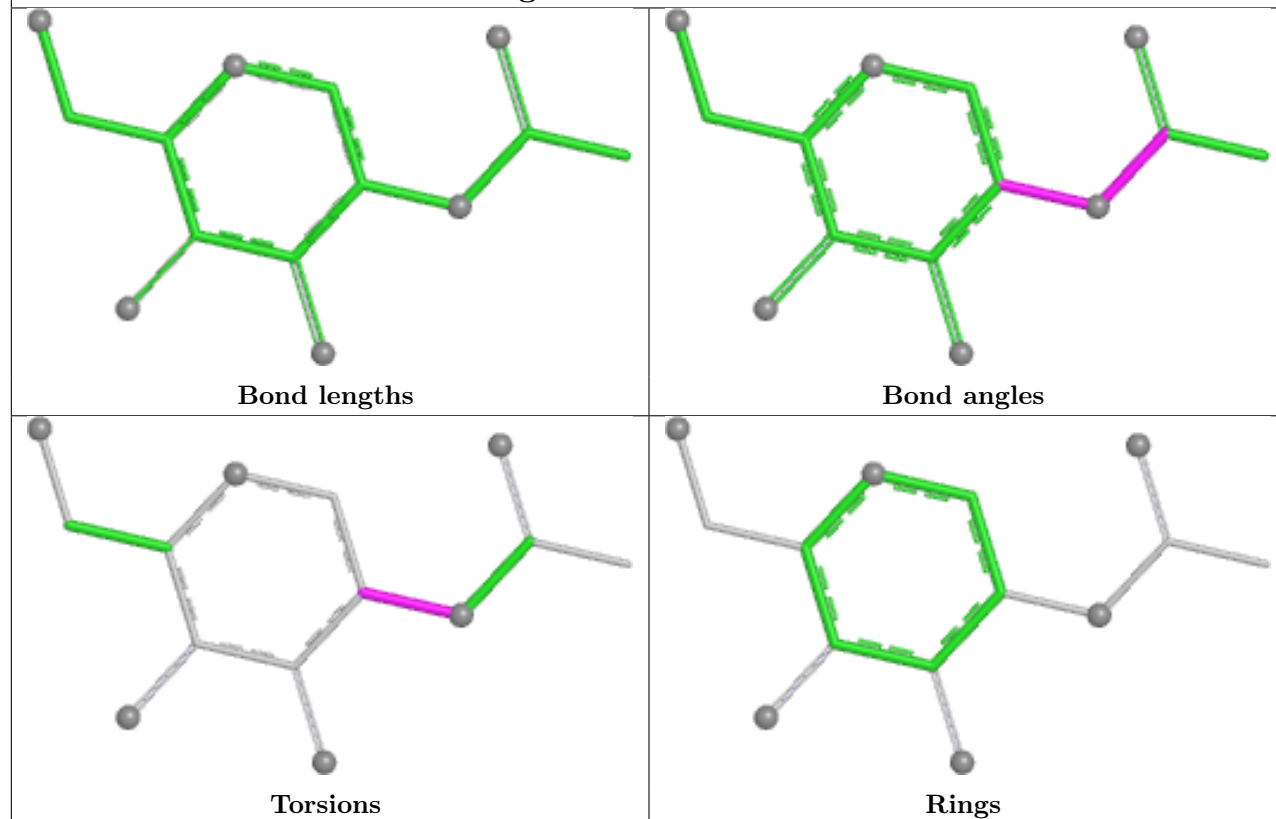


Ligand NAG B 1208

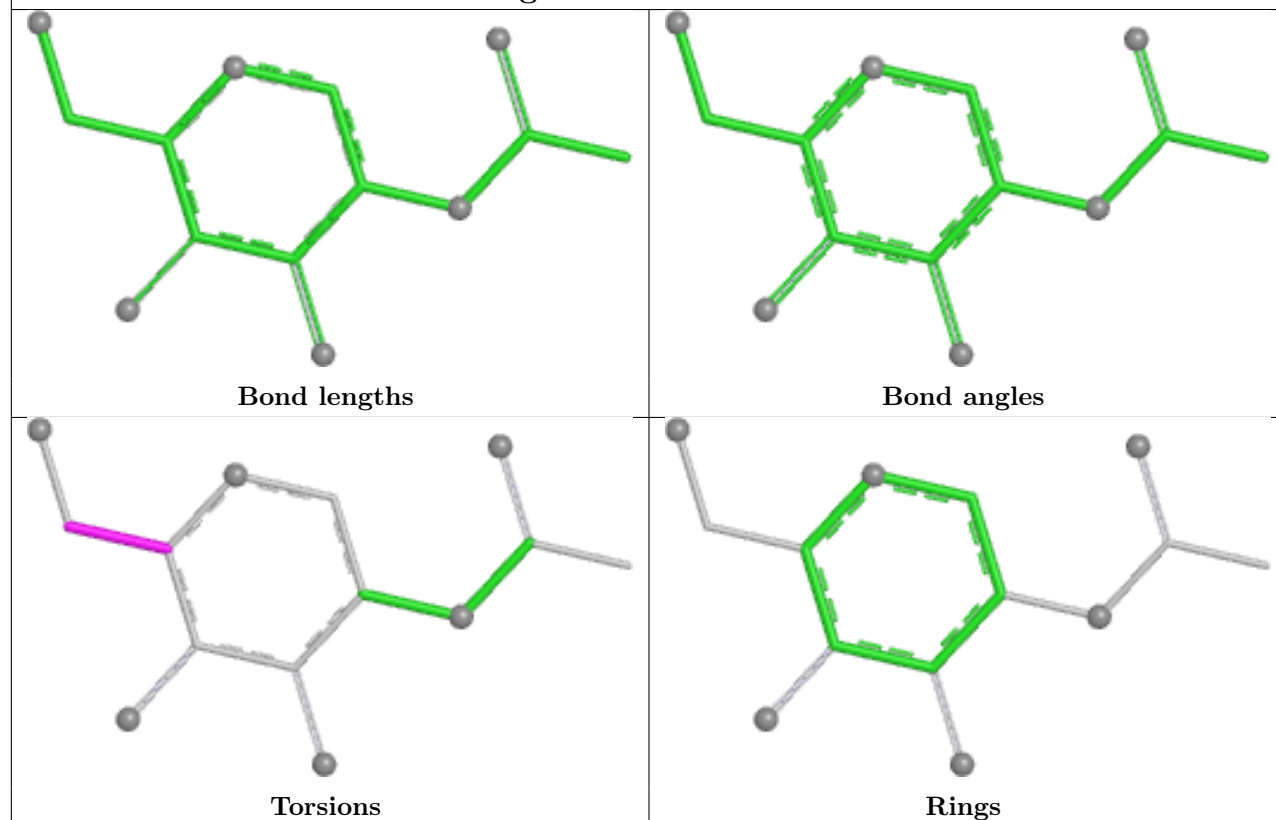




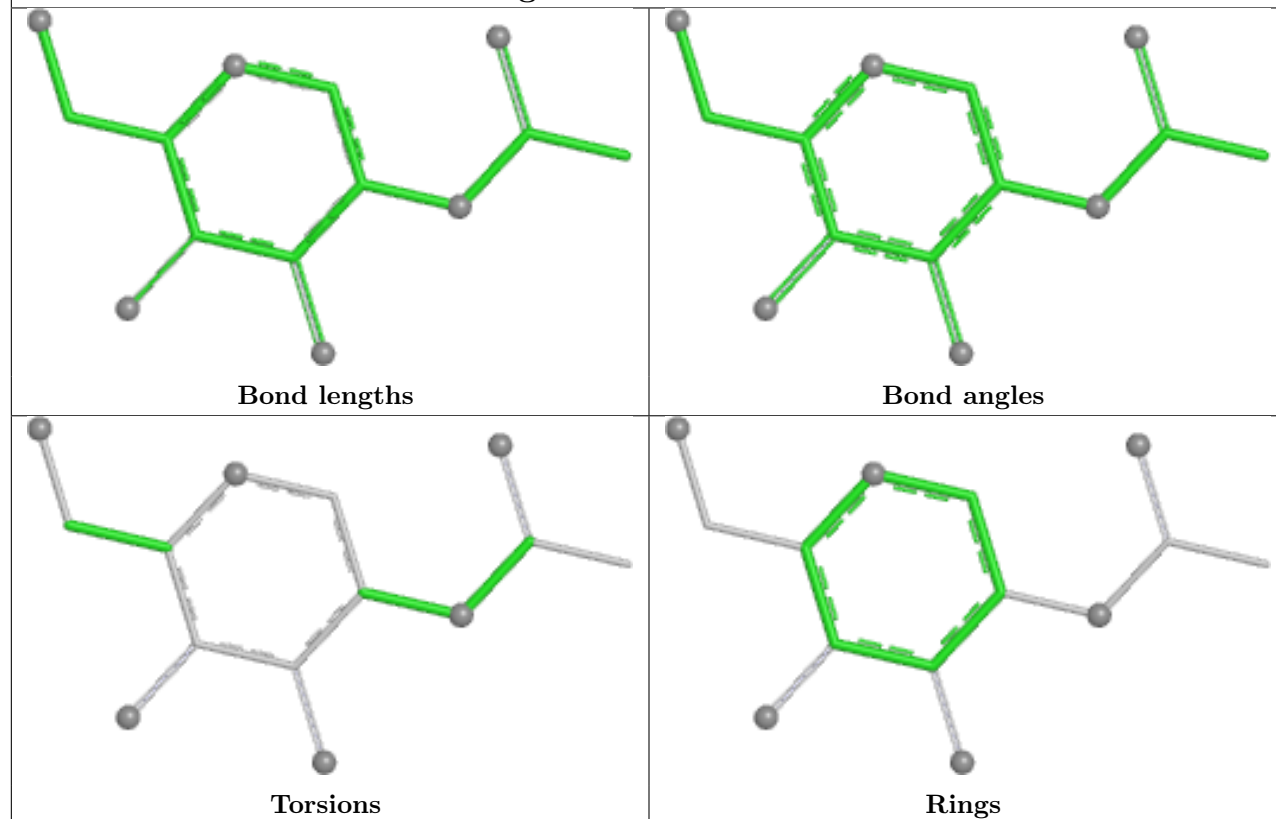
Ligand NAG A 1209



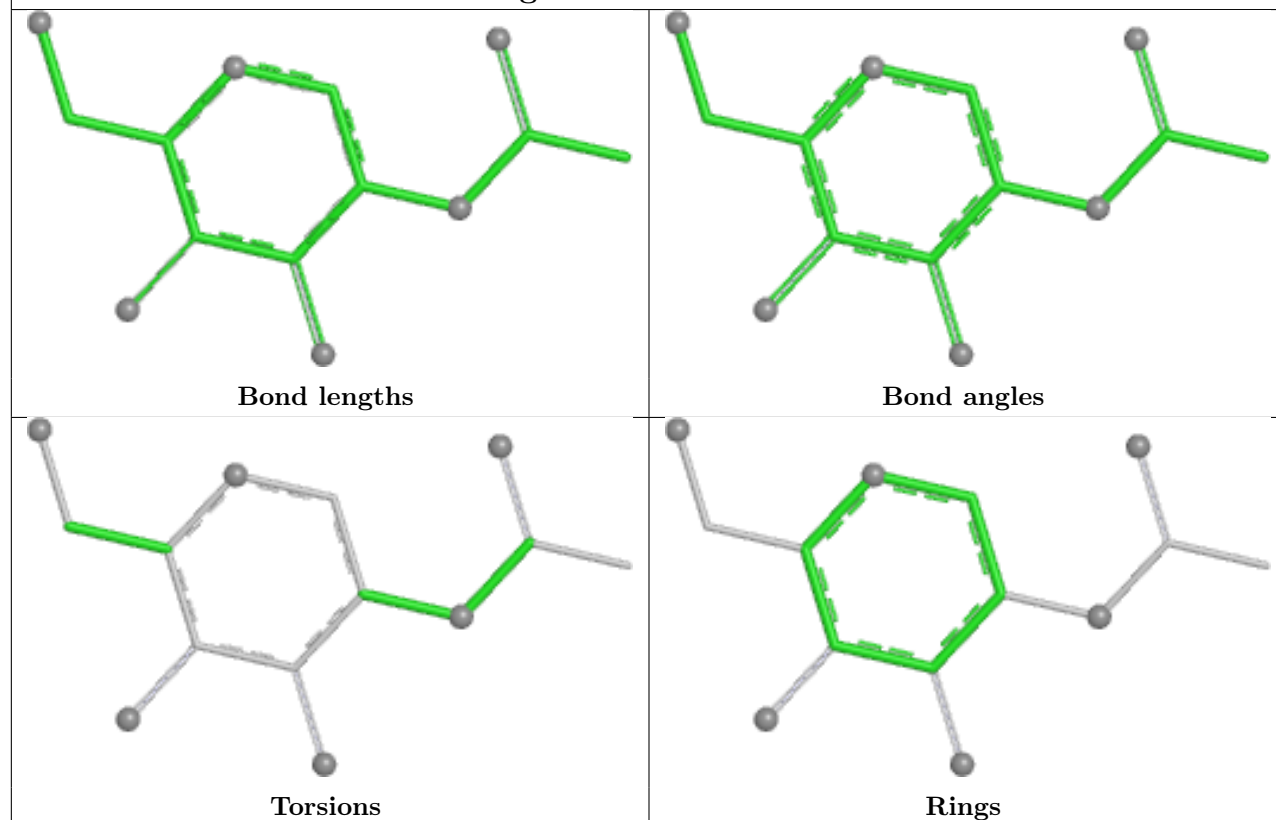
Ligand NAG B 1216

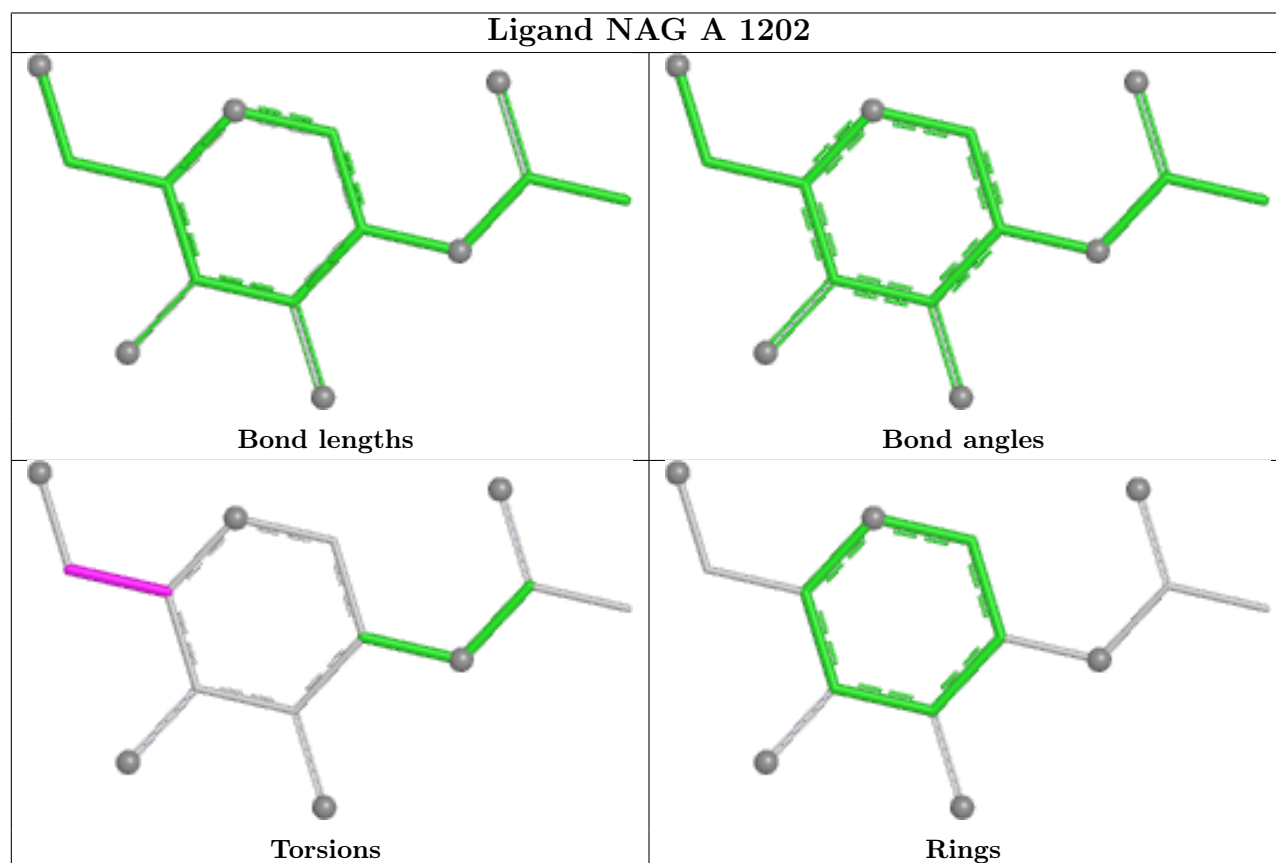
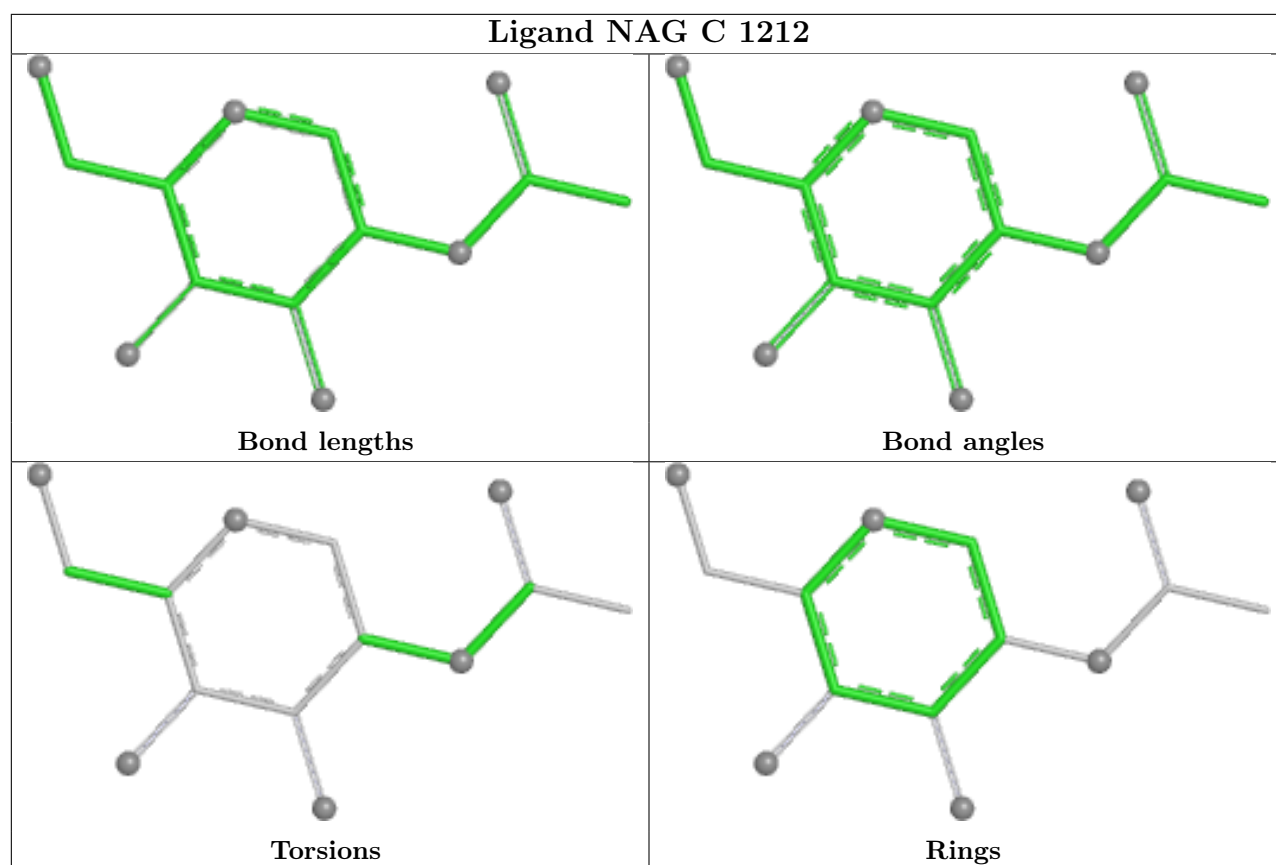


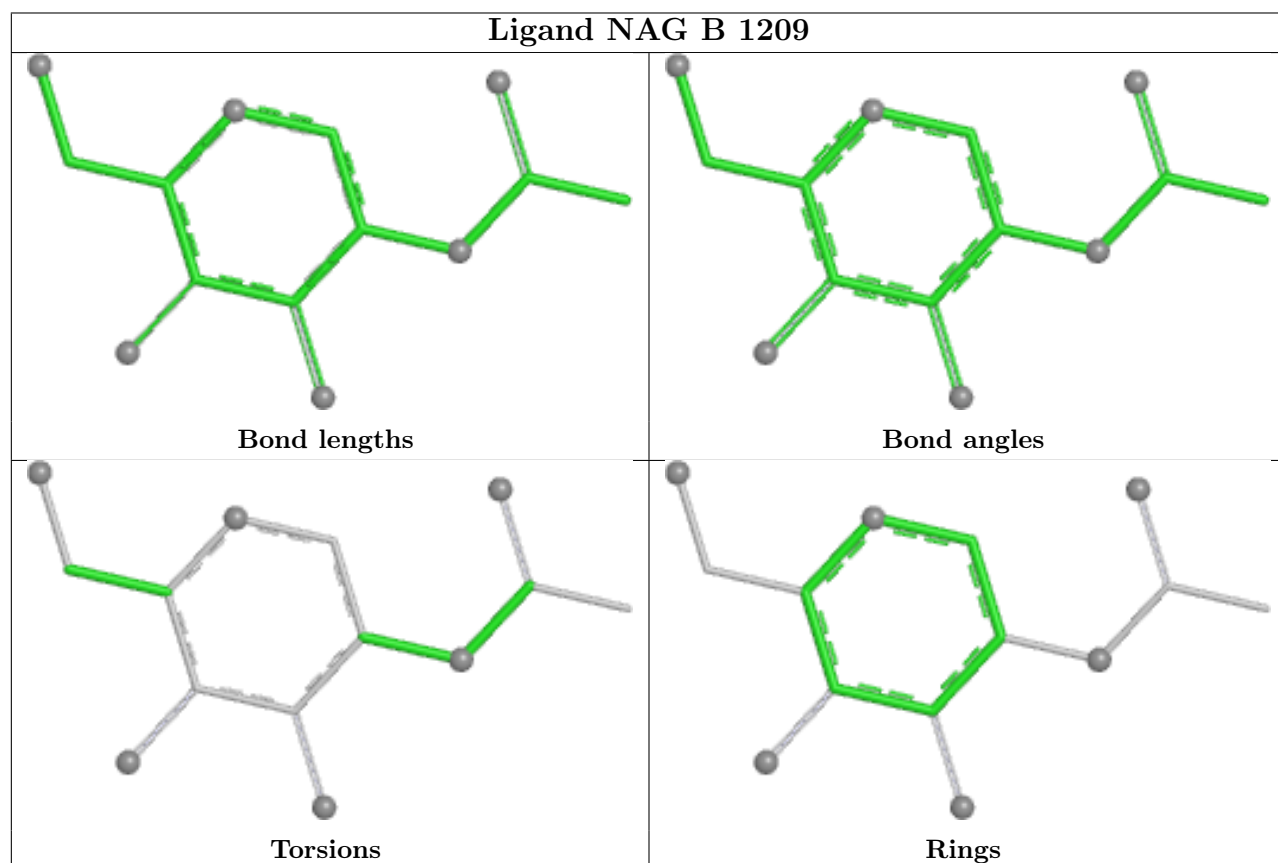
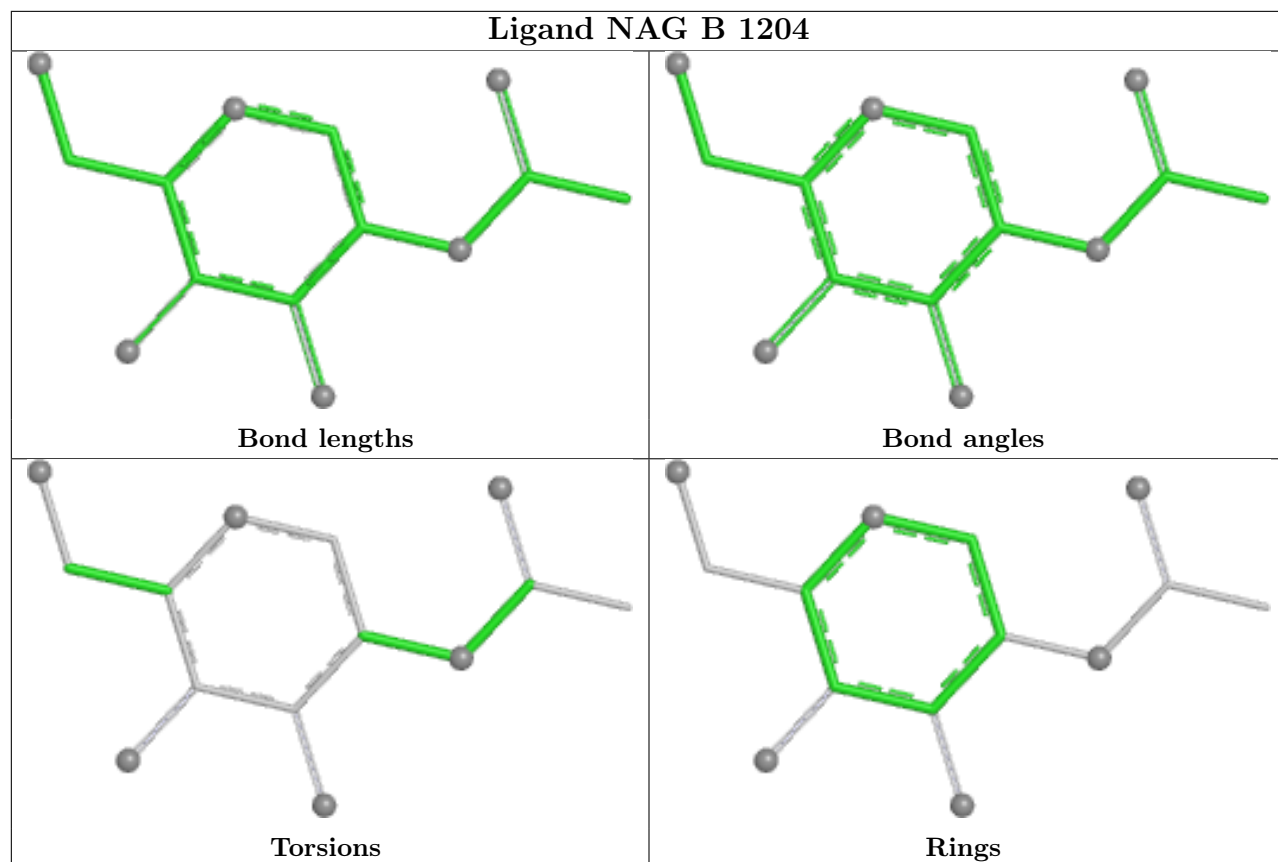
Ligand NAG A 1210



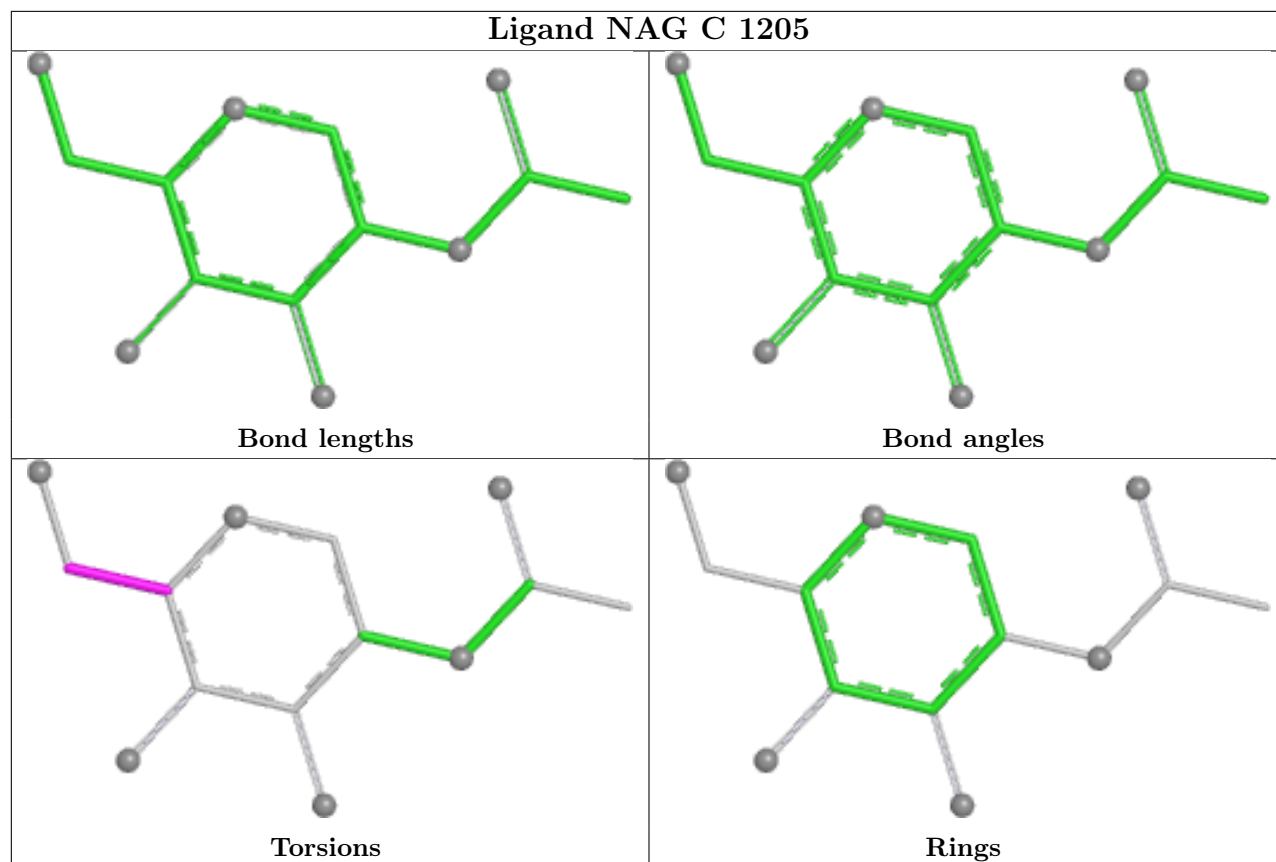
Ligand NAG B 1218



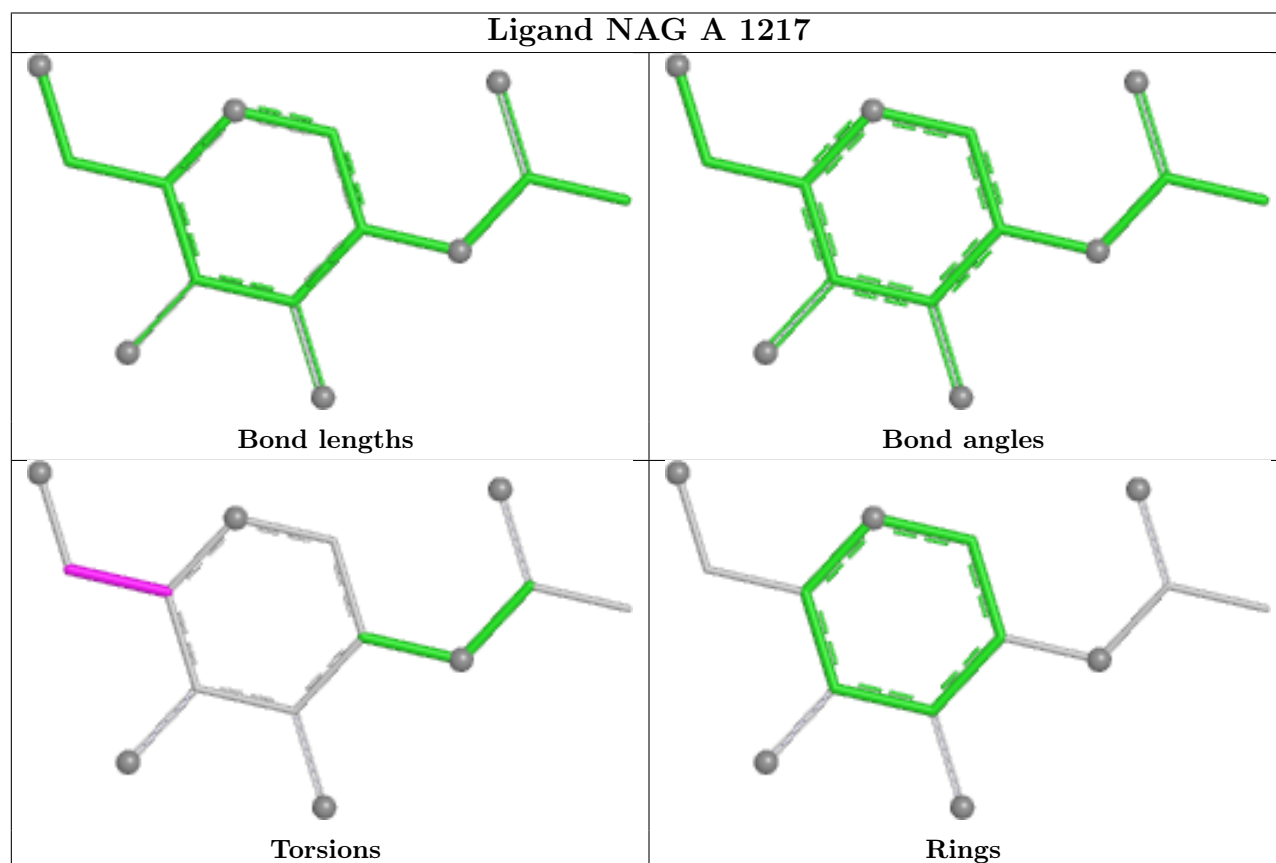




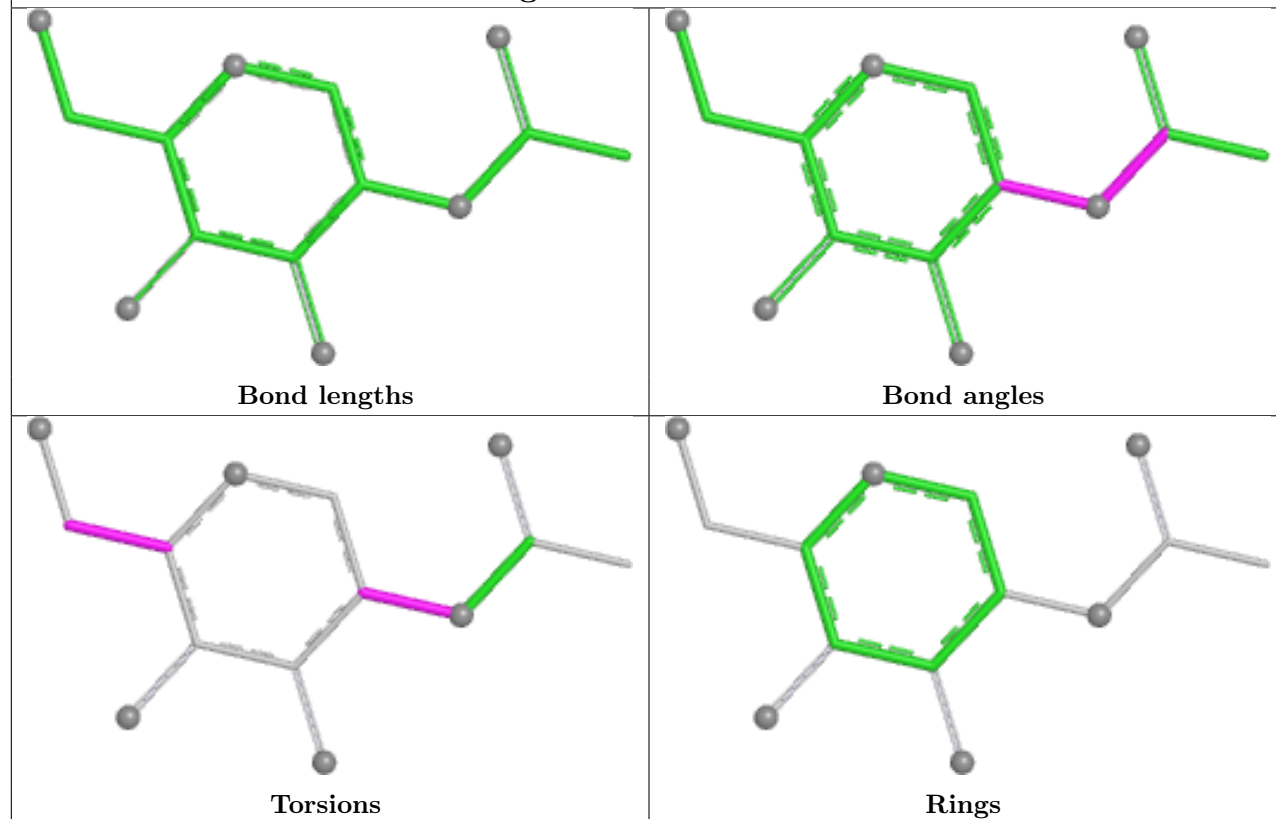
Ligand NAG C 1205



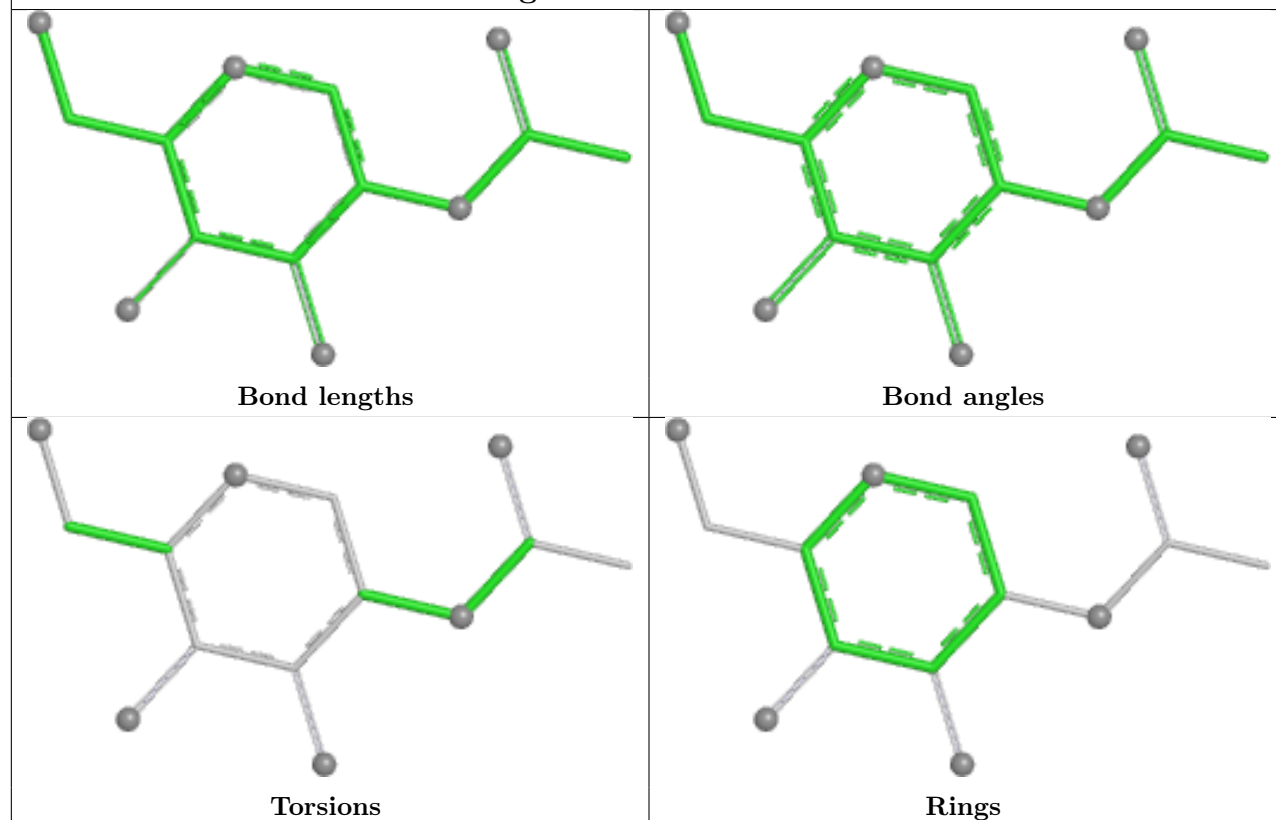
Ligand NAG A 1217

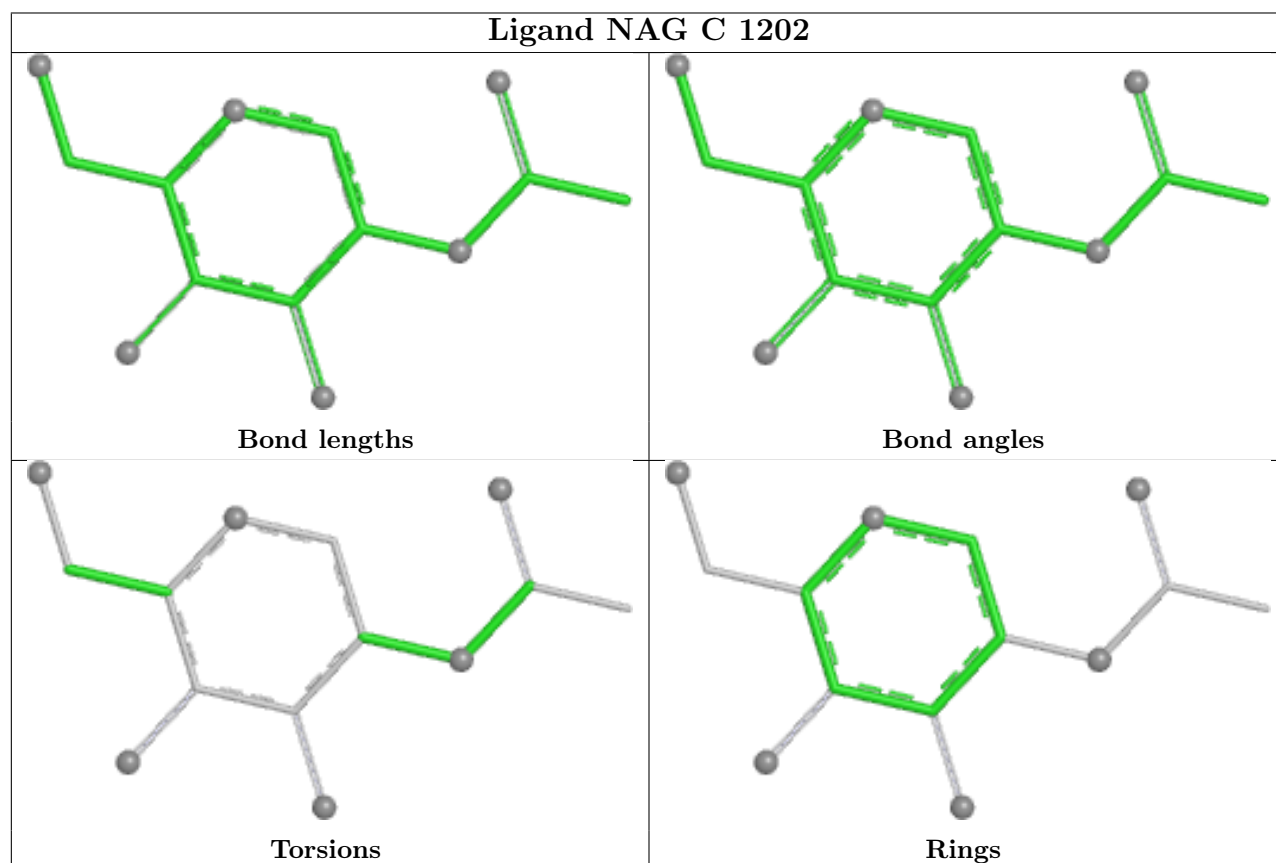
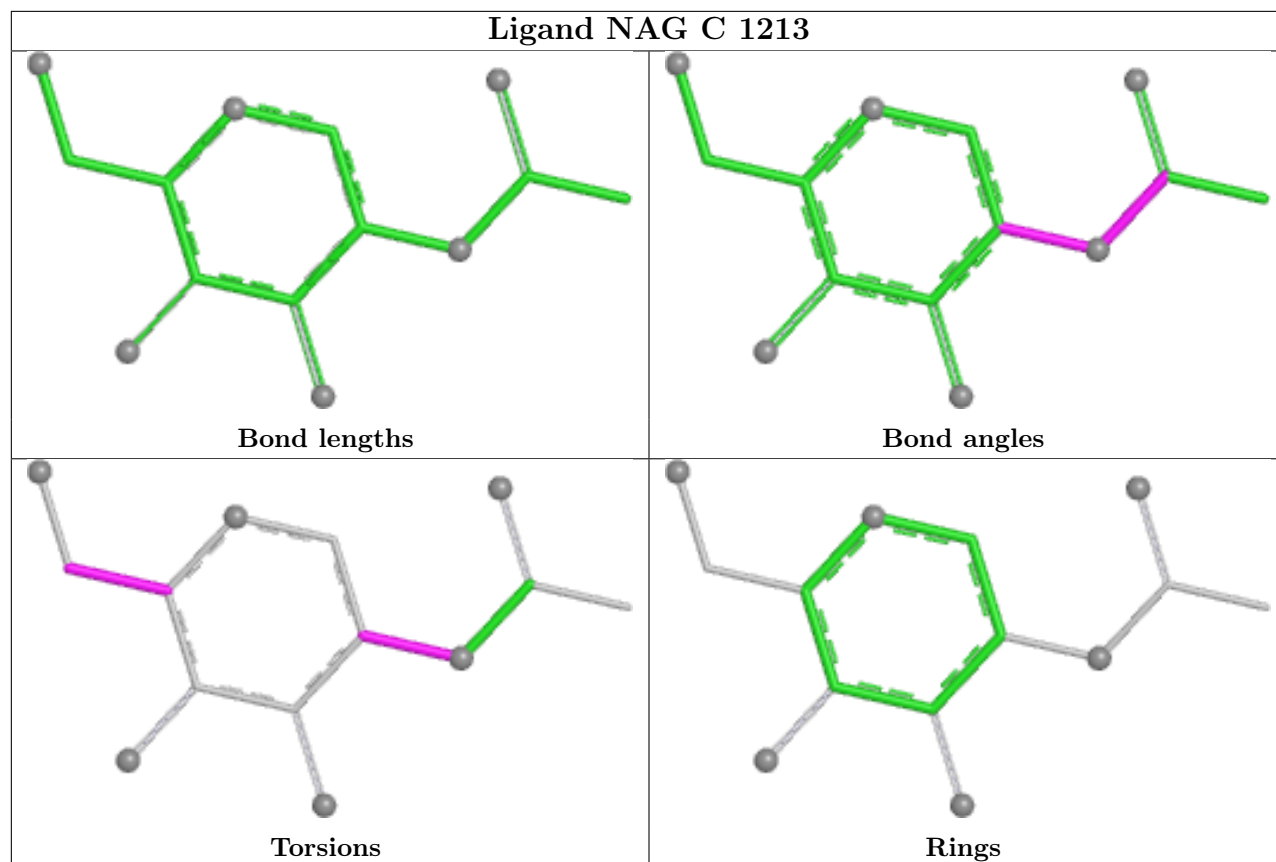


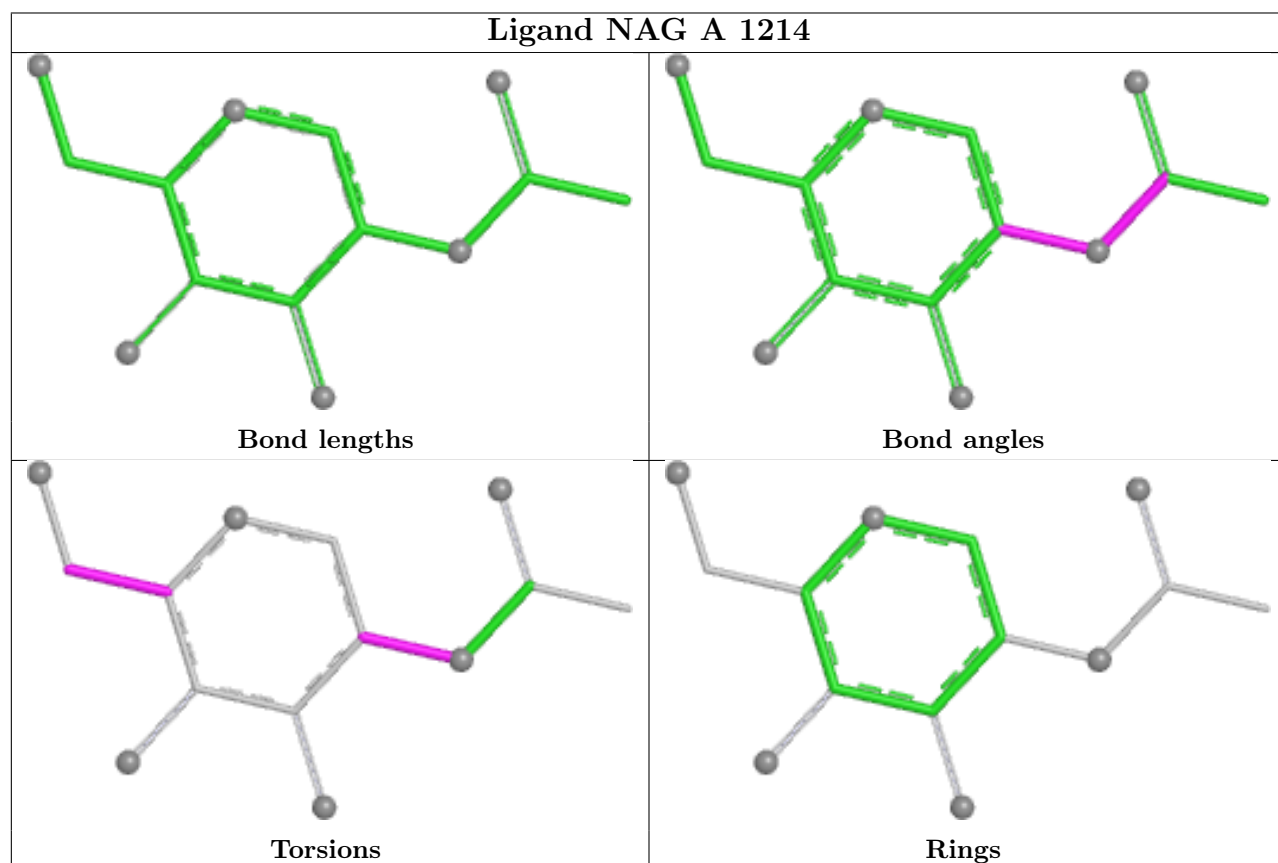
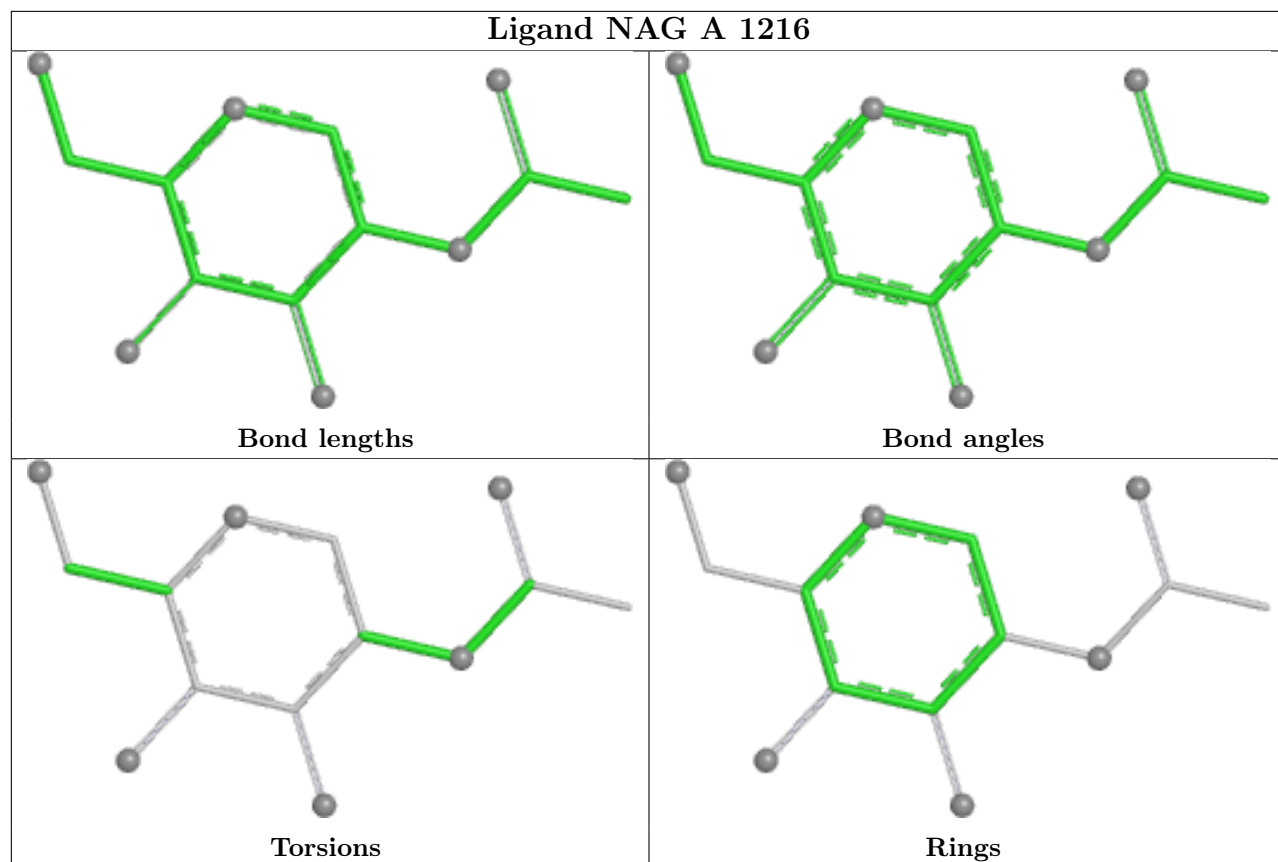
Ligand NAG B 1211



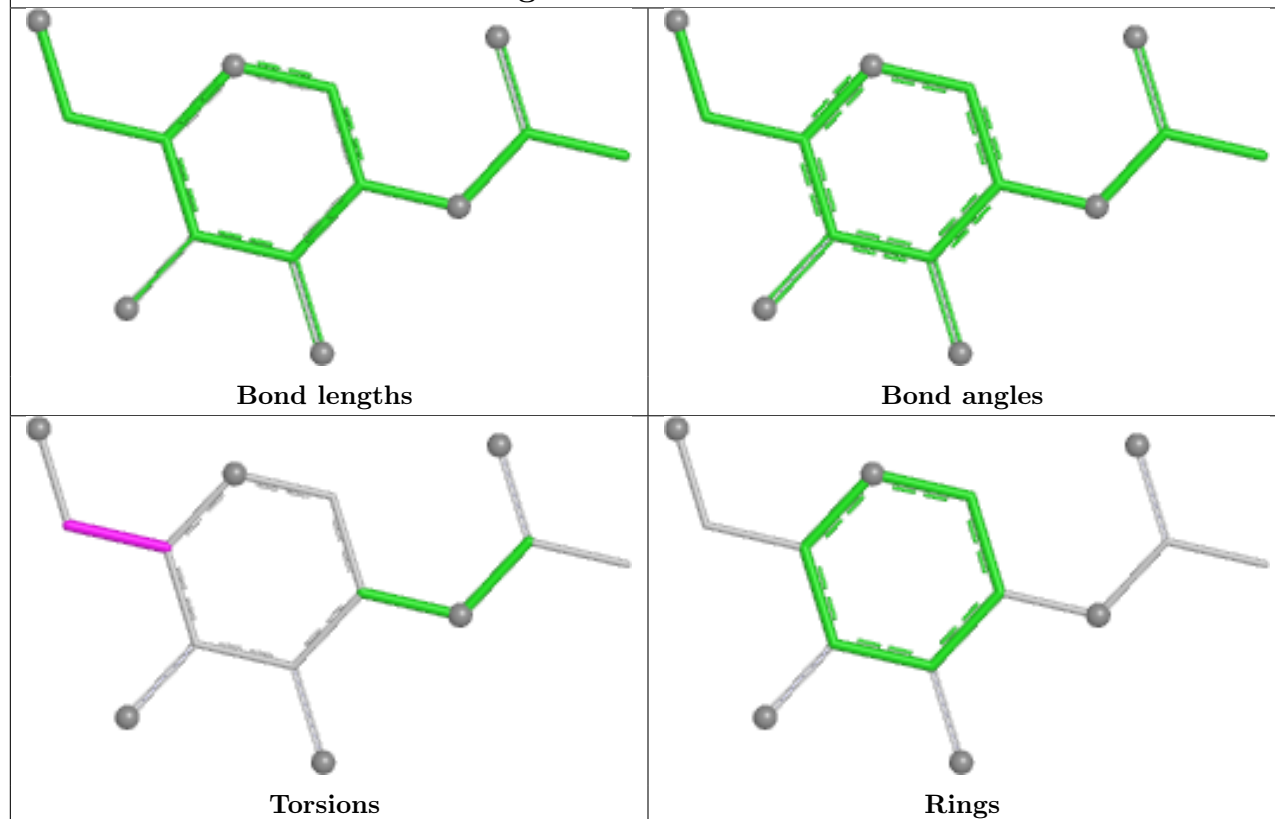
Ligand NAG C 1215



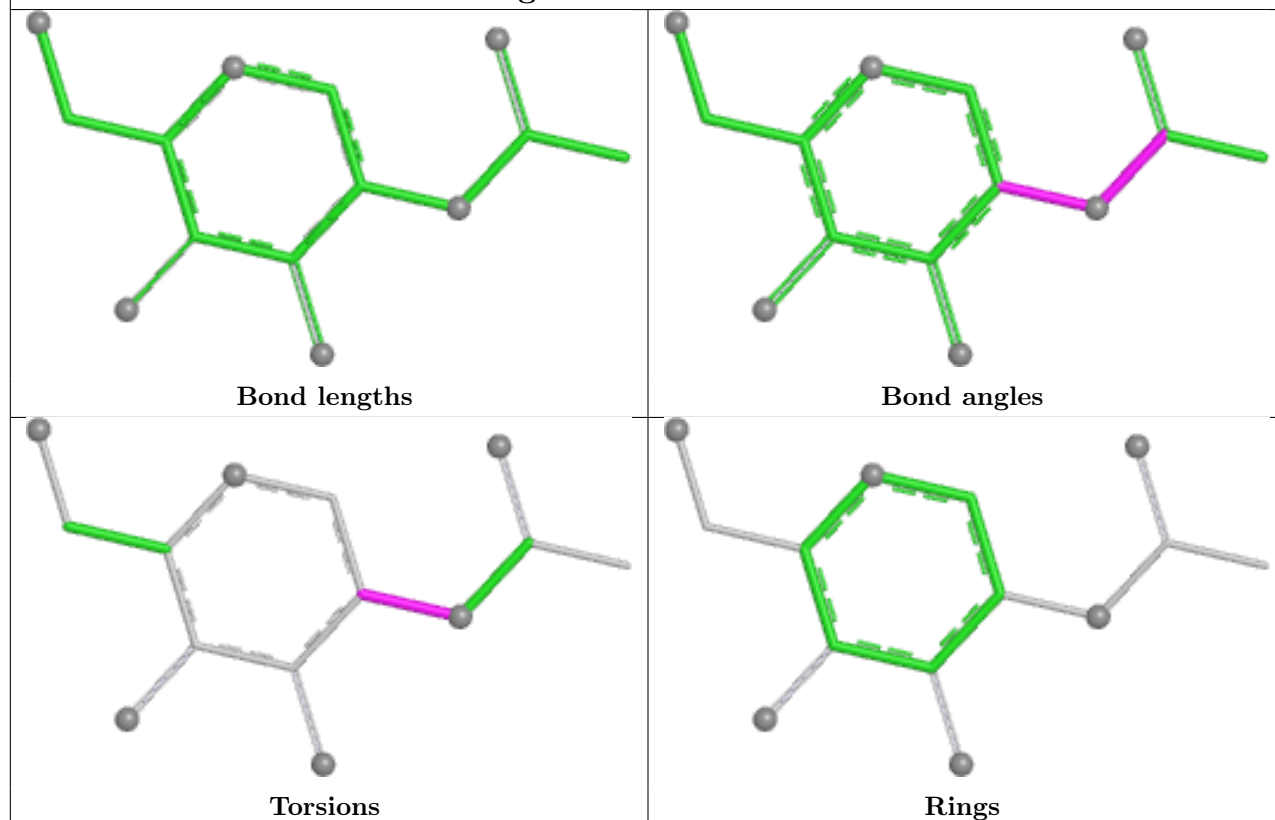


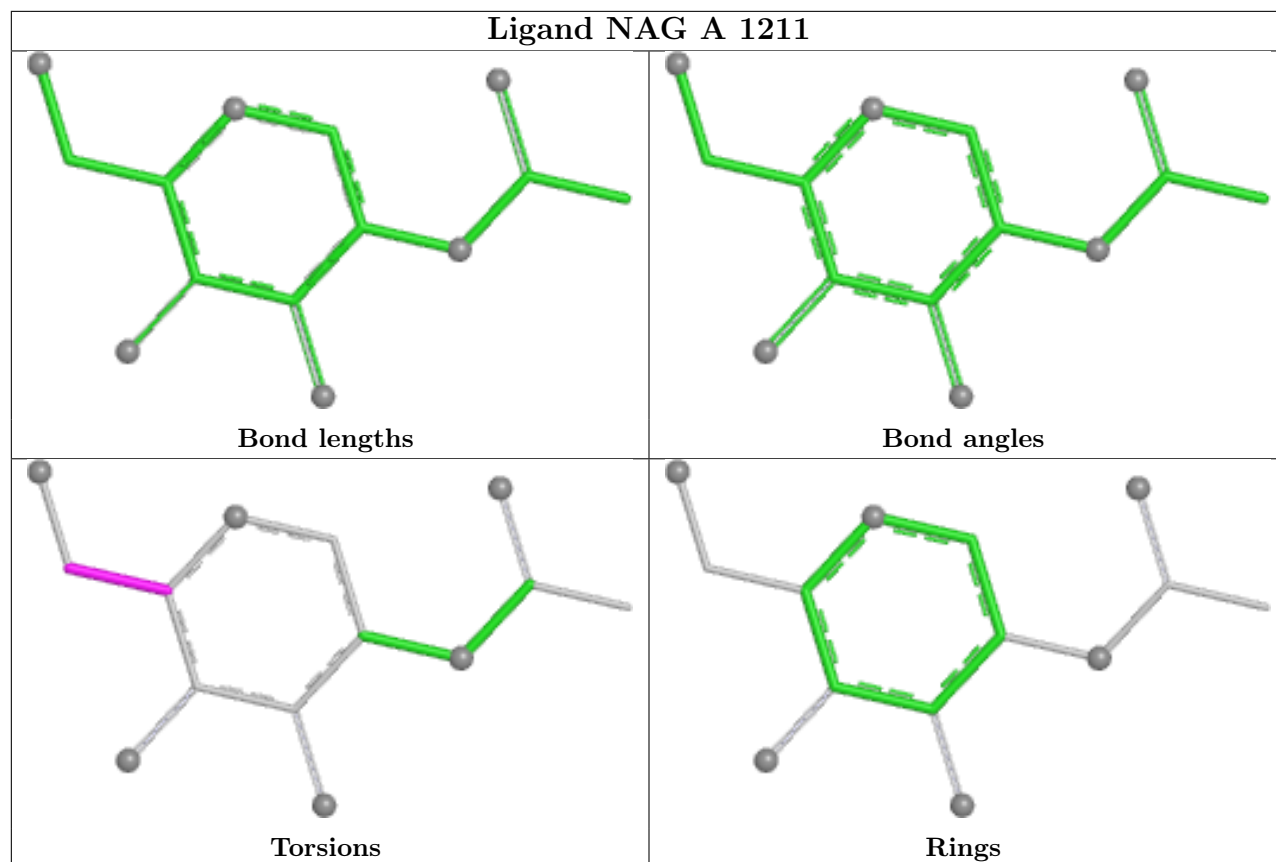


Ligand NAG C 1210



Ligand NAG C 1206





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.

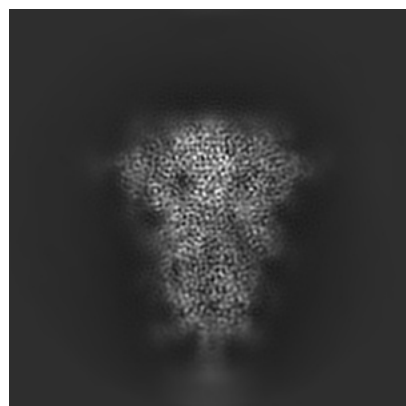
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-60551. These allow visual inspection of the internal detail of the map and identification of artifacts.

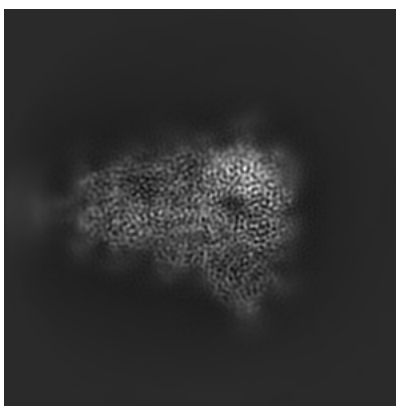
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

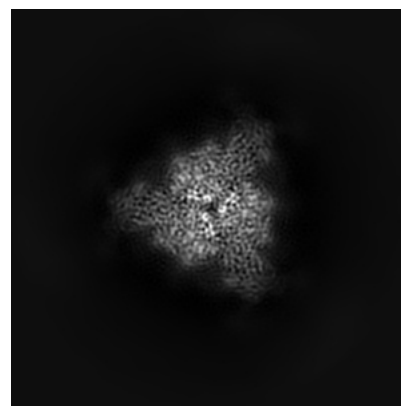
6.1.1 Primary map



X

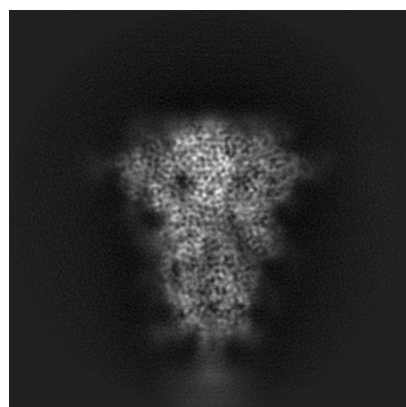


Y

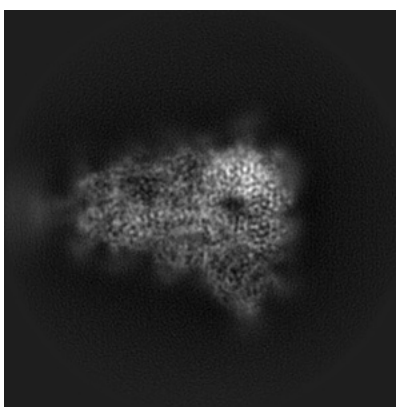


Z

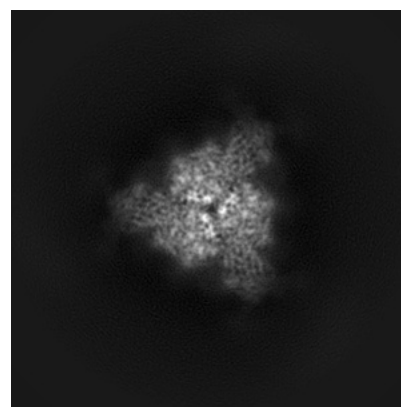
6.1.2 Raw map



X



Y

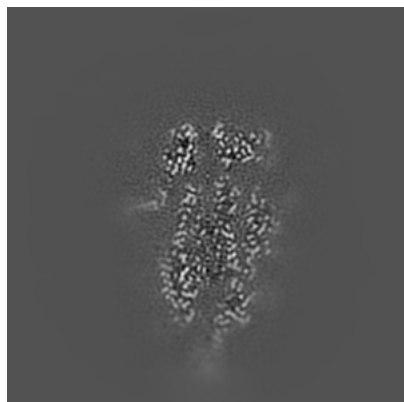


Z

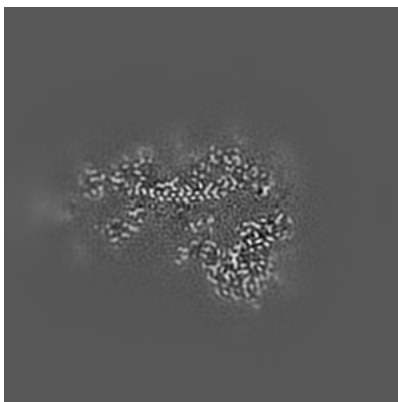
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

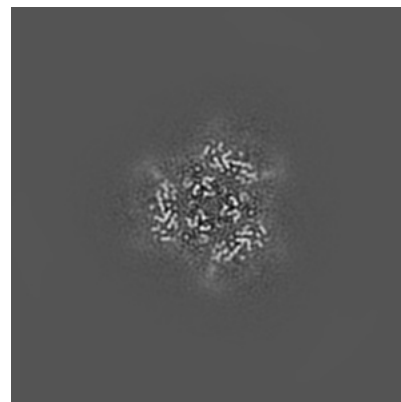
6.2.1 Primary map



X Index: 128

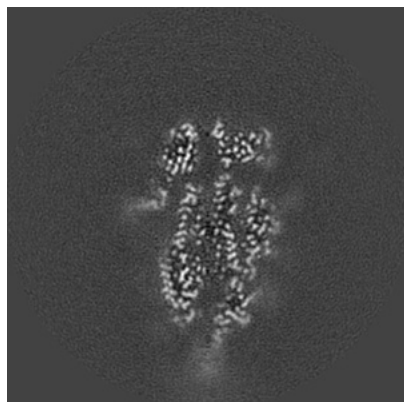


Y Index: 128

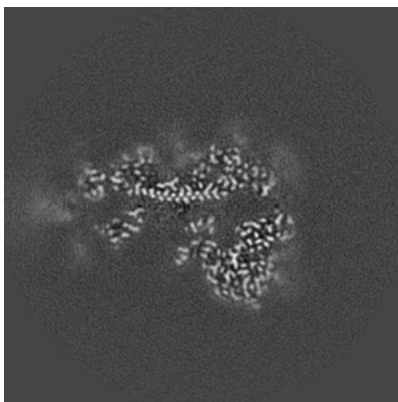


Z Index: 128

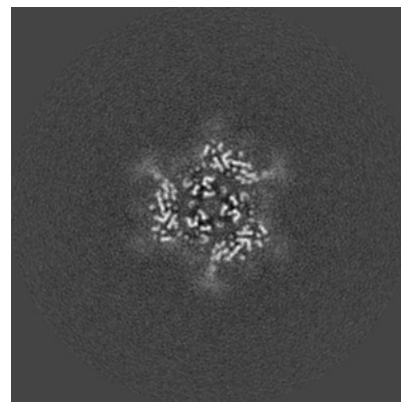
6.2.2 Raw map



X Index: 128



Y Index: 128

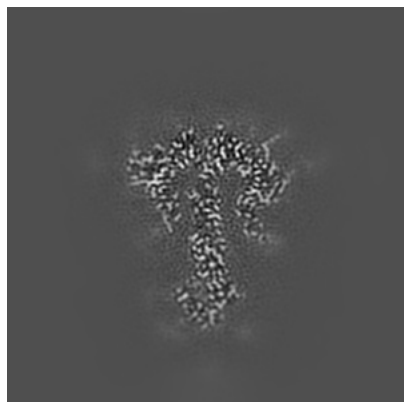


Z Index: 128

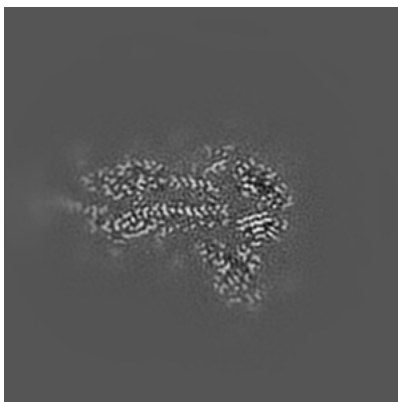
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

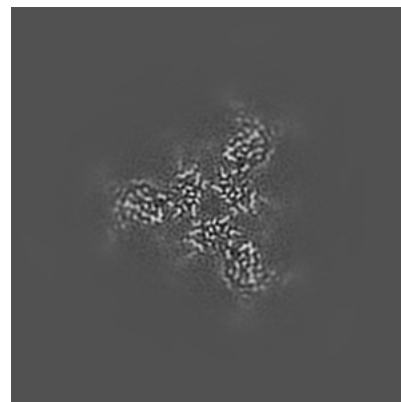
6.3.1 Primary map



X Index: 141

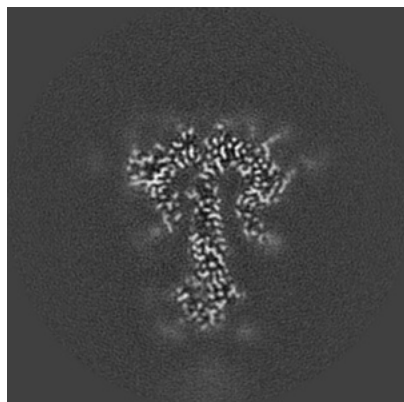


Y Index: 135

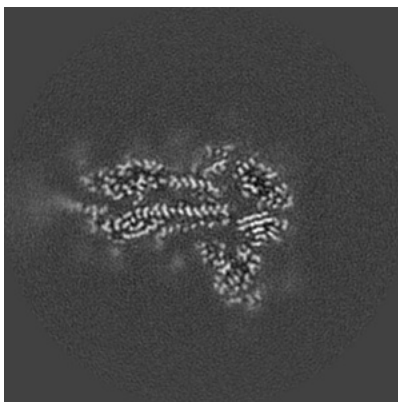


Z Index: 154

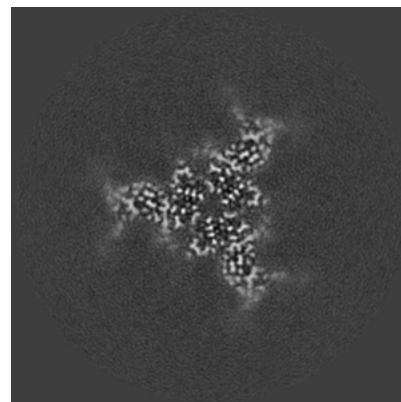
6.3.2 Raw map



X Index: 141



Y Index: 135

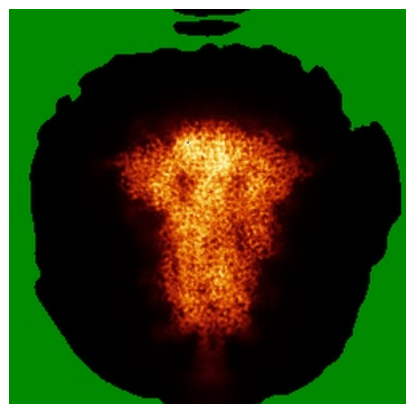


Z Index: 158

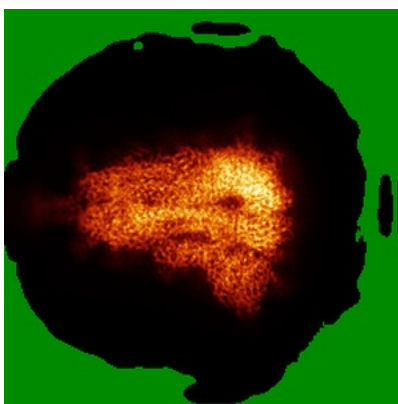
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

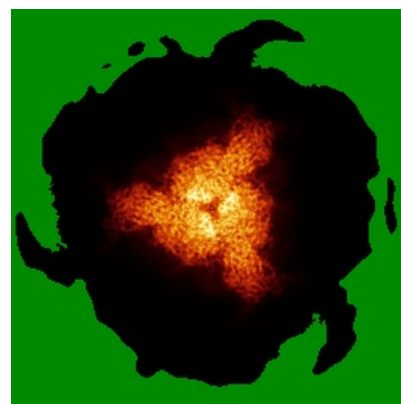
6.4.1 Primary map



X

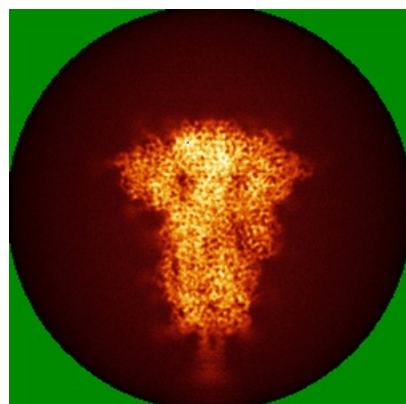


Y

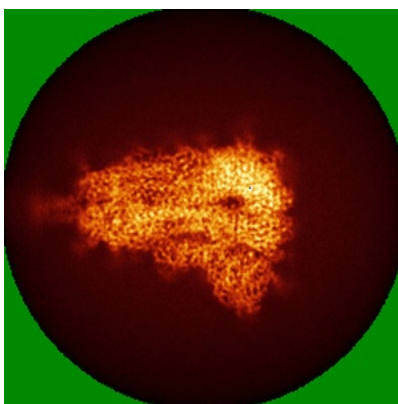


Z

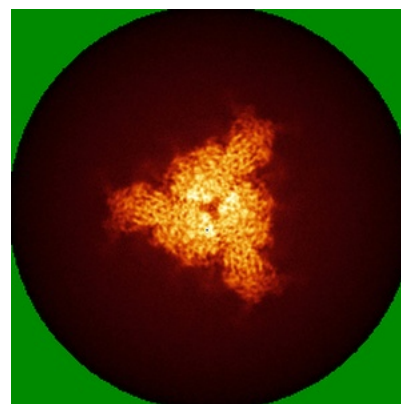
6.4.2 Raw map



X



Y

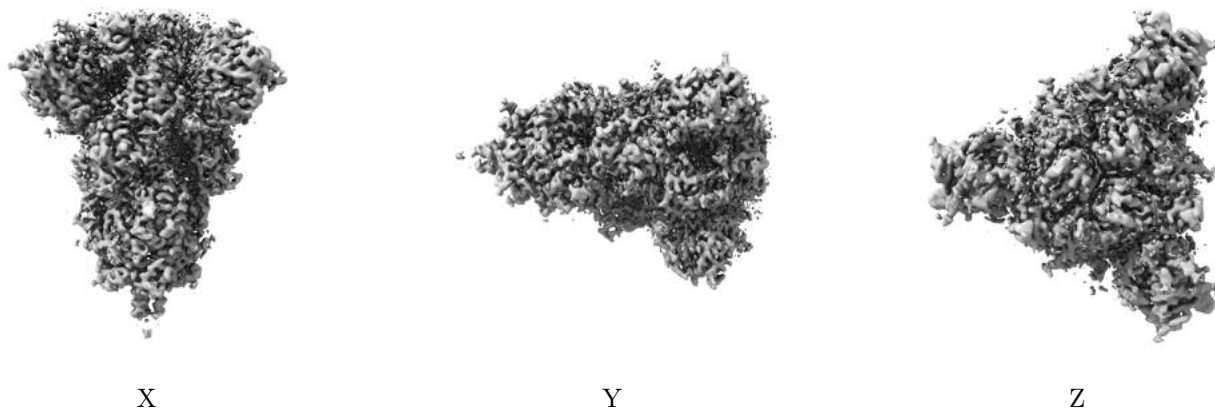


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

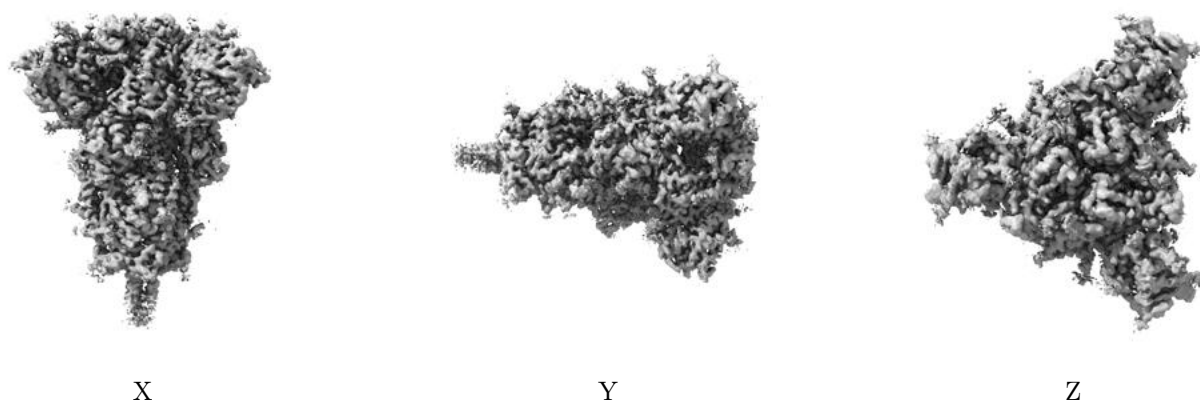
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.024. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

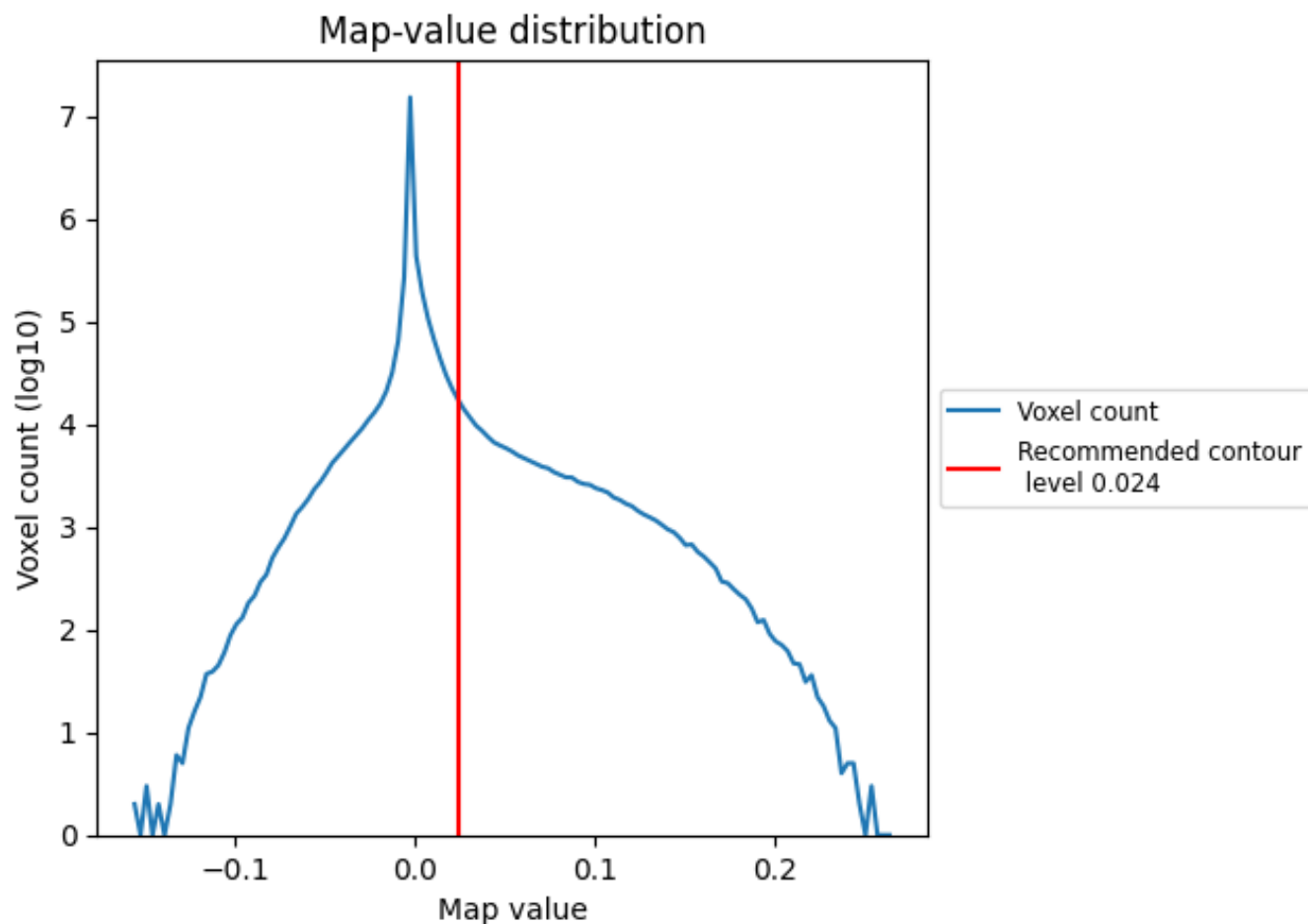
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

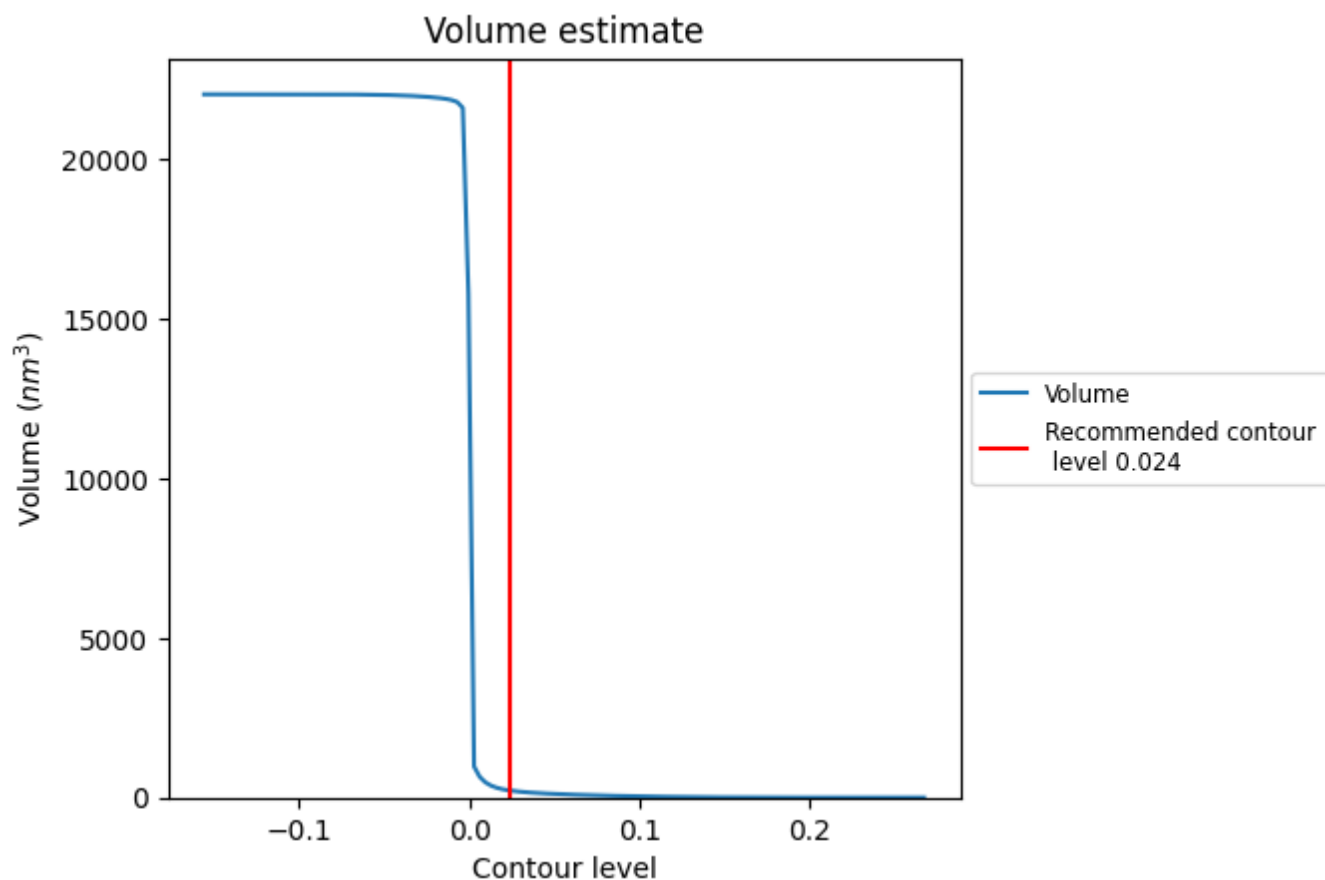
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

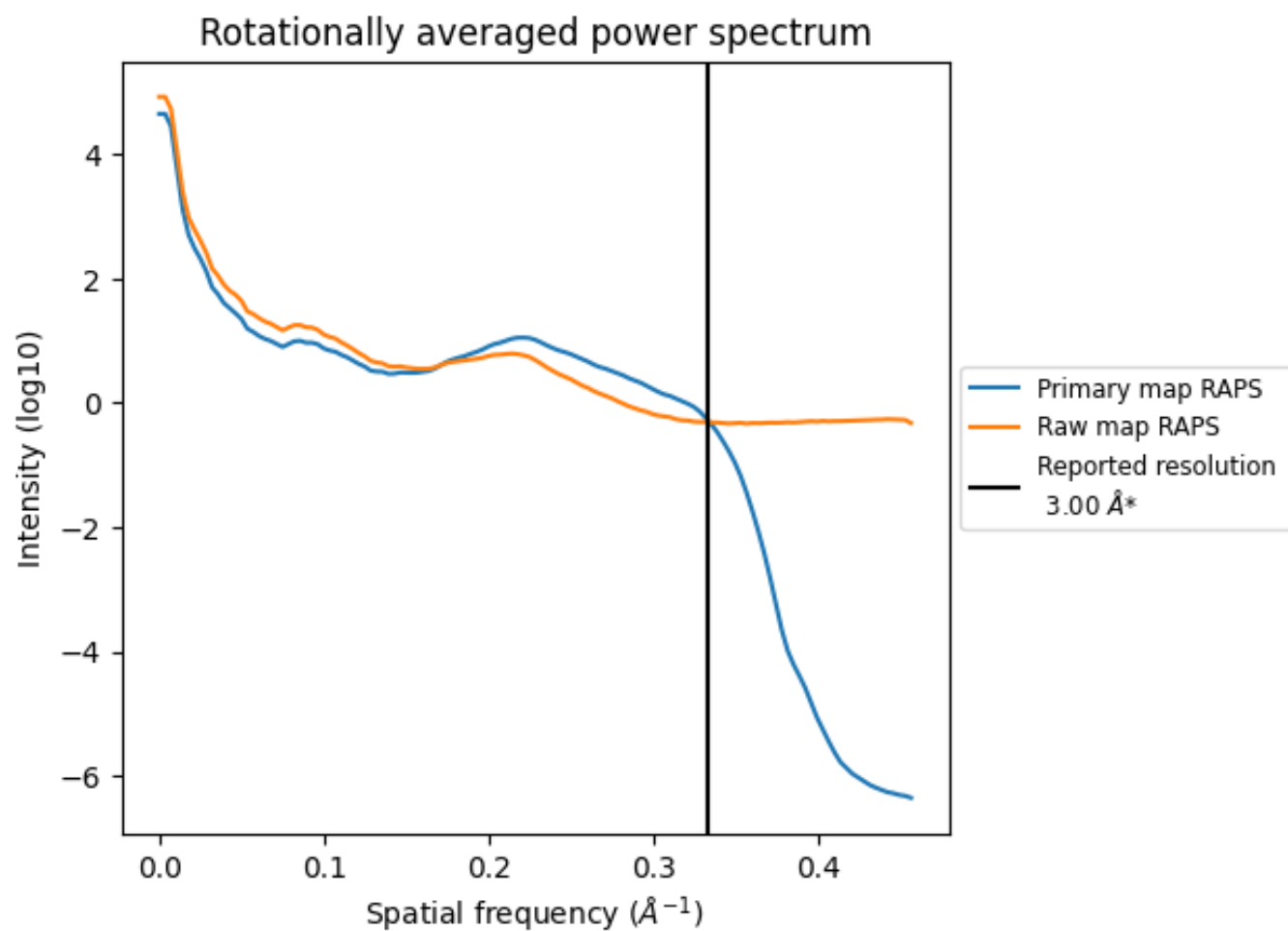
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 221 nm³; this corresponds to an approximate mass of 200 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

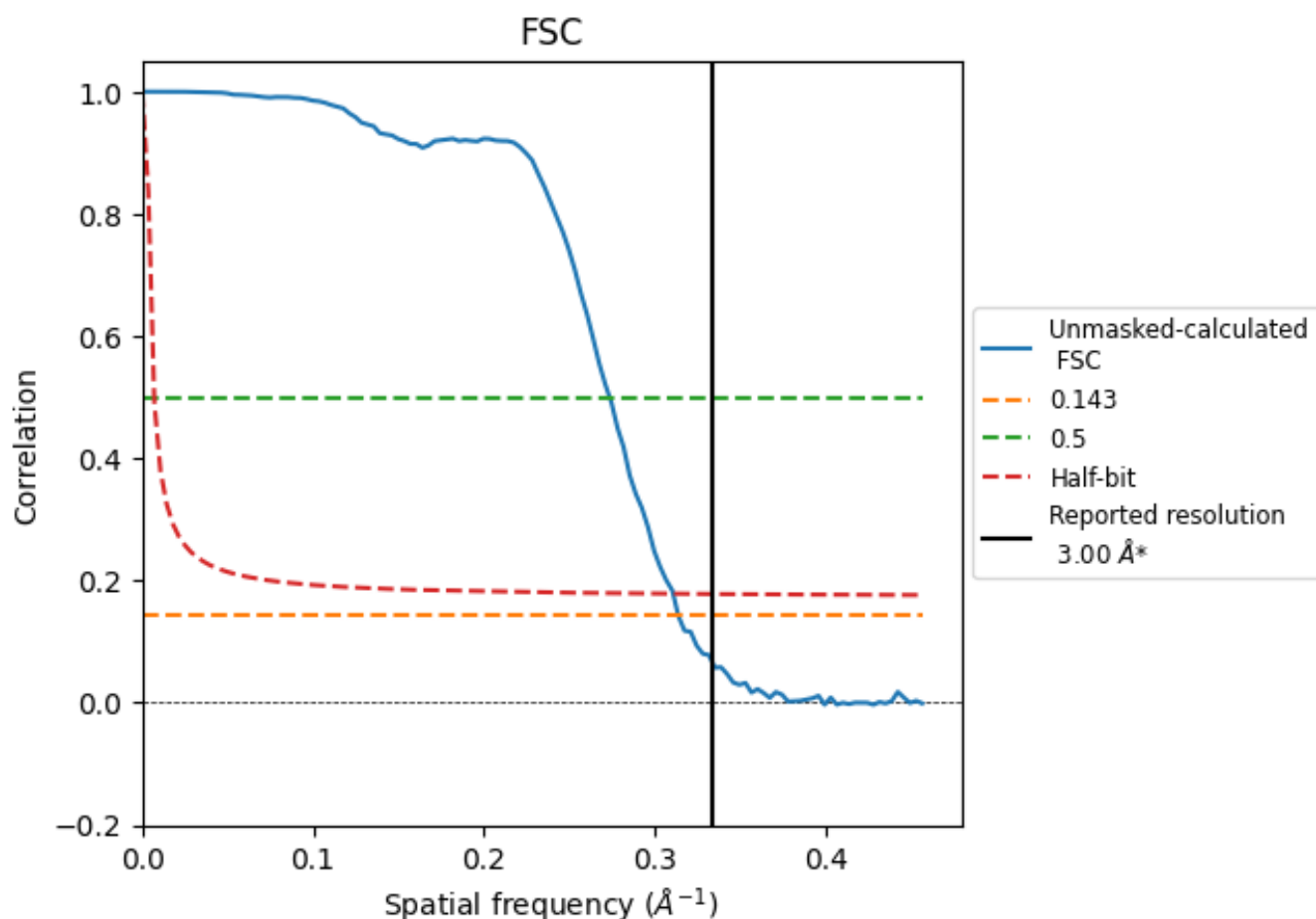


*Reported resolution corresponds to spatial frequency of $0.333 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.333 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

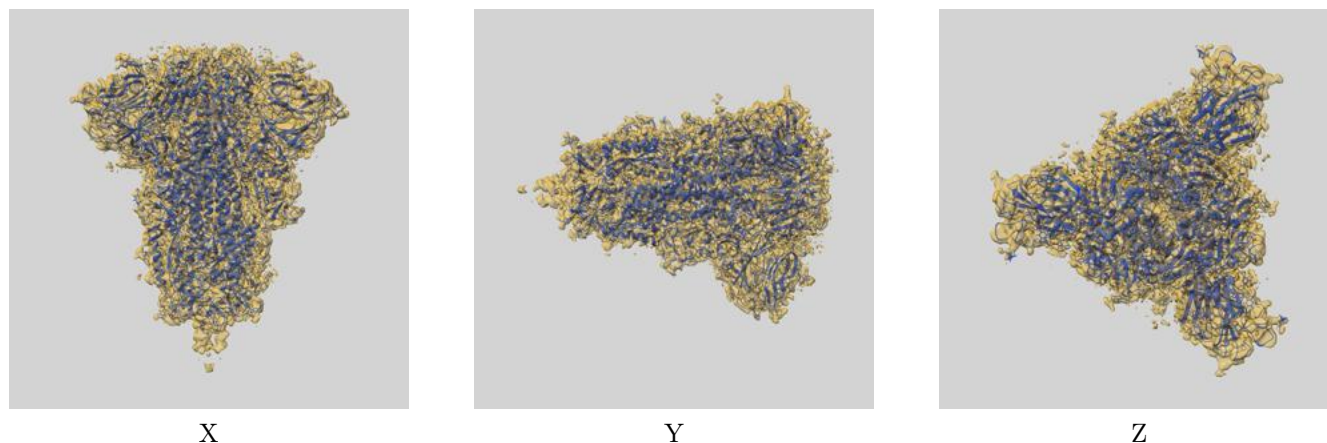
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.19	3.65	3.22

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

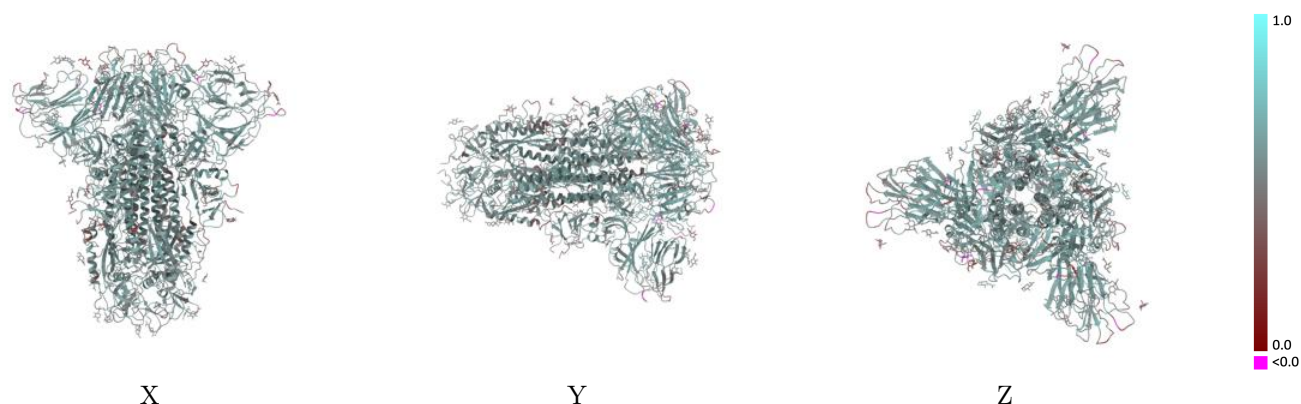
This section contains information regarding the fit between EMDB map EMD-60551 and PDB model 8ZY0. Per-residue inclusion information can be found in section [3](#) on page [9](#).

9.1 Map-model overlay [i](#)



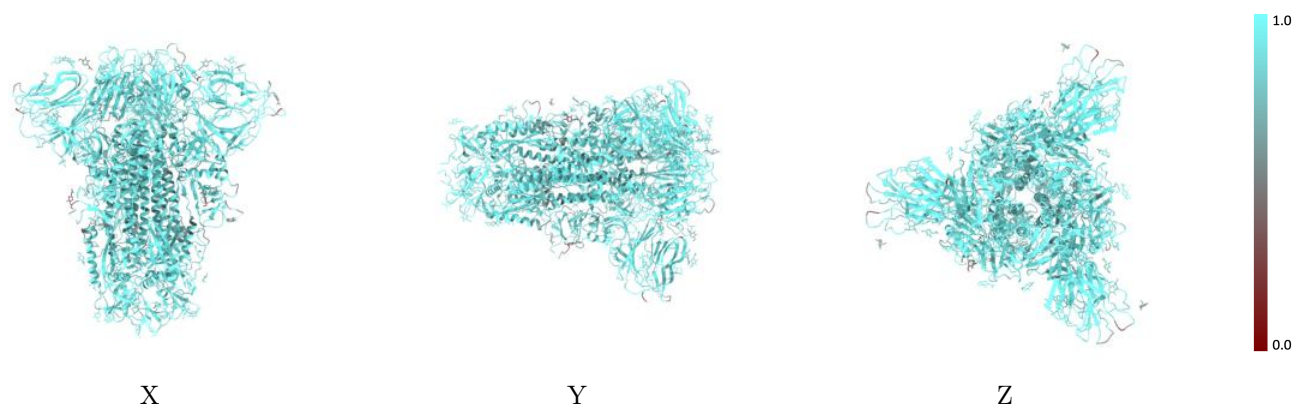
The images above show the 3D surface view of the map at the recommended contour level 0.024 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



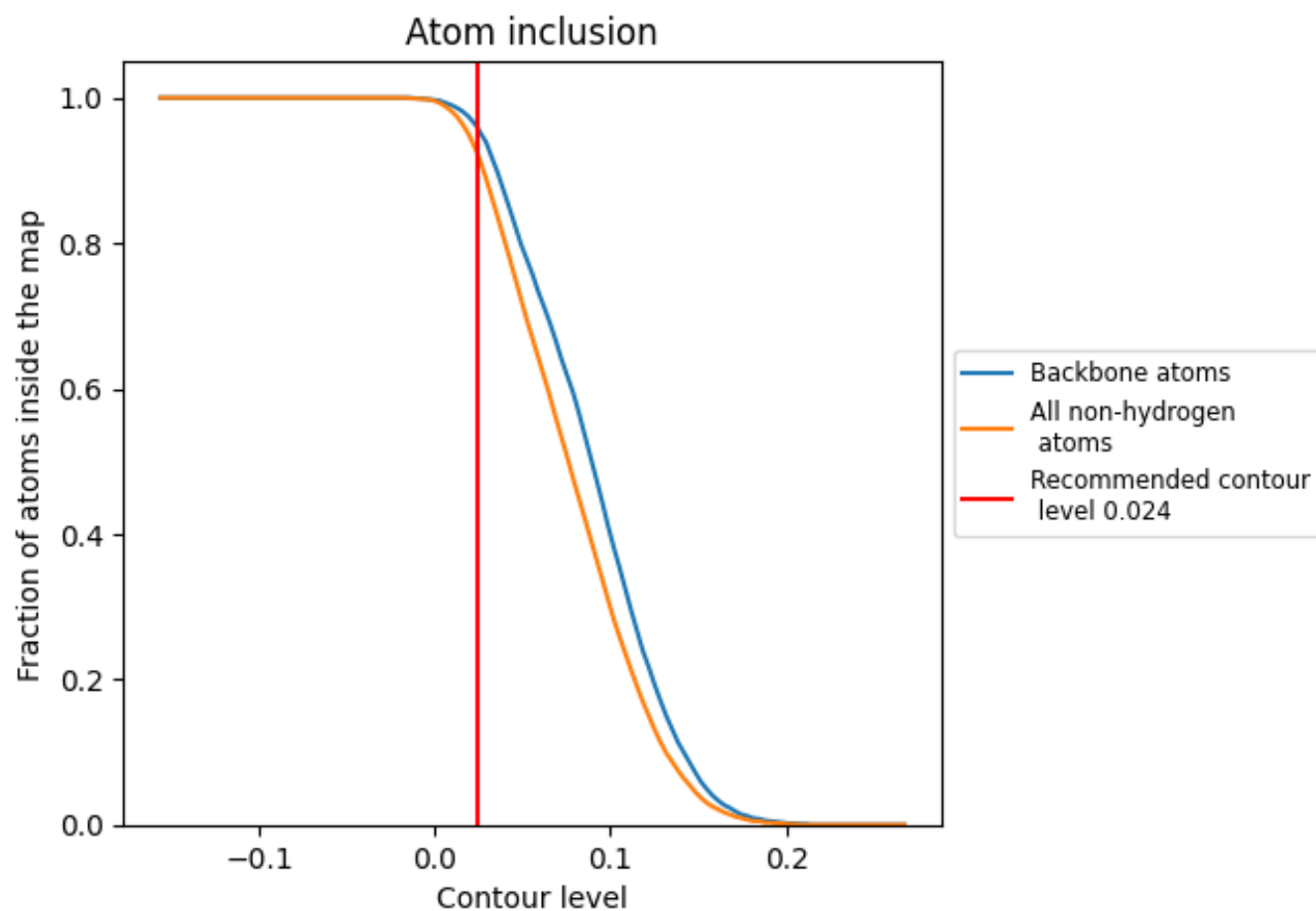
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.024).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 93% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.024) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9260	0.5570
A	0.9280	0.5580
B	0.9260	0.5580
C	0.9270	0.5580
D	0.8210	0.5220
E	0.7210	0.3560
F	0.8210	0.5130
G	0.9400	0.4960
H	0.8210	0.5210
I	0.9400	0.4880

1.0

0.0

<0.0