



Full wwPDB EM Validation Report ⓘ

Nov 18, 2024 – 01:02 PM JST

PDB ID : 8Y32
EMDB ID : EMD-38866
Title : BA.2.86-V483 RBD in complex with bACE2
Authors : Yue, C.; Liu, P.
Deposited on : 2024-01-28
Resolution : 4.13 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) : 2.39

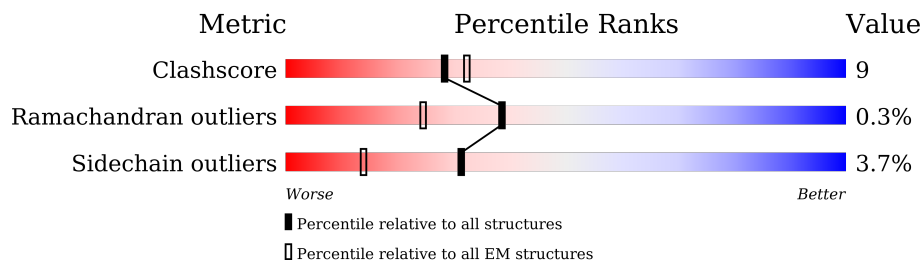
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.13 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	786	
2	B	211	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6571 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 Angiotensin-converting enzyme.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	596	4906	3136	815	925	30	0	0

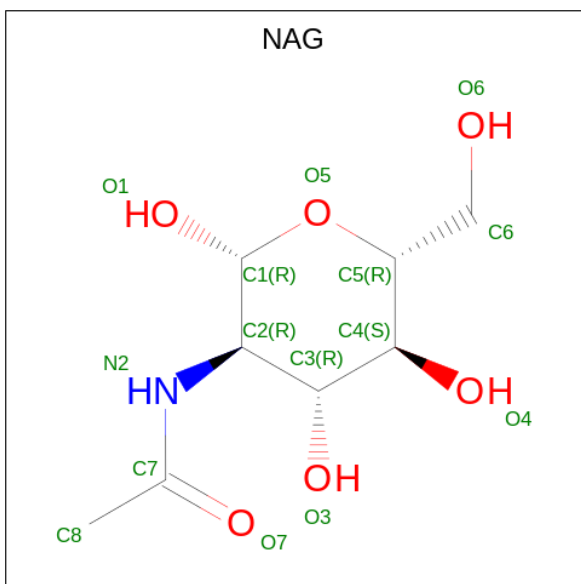
- Molecule 2 is a protein called Spike protein S1 RBD.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	195	1566	1015	262	281	8	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		AltConf
			Total	Cl	
3	A	1	1	1	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



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

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	597508	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2000	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, CL

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	0.28	0/5043	0.50	0/6838
2	B	0.31	0/1618	0.49	0/2203
All	All	0.29	0/6661	0.50	0/9041

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	4906	0	4691	92	0
2	B	1566	0	1486	27	0
3	A	1	0	0	0	0
4	A	70	0	65	0	0
4	B	28	0	26	0	0
All	All	6571	0	6268	119	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 9.

All (119) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:TRP:HE1	1:A:305:ARG:HB3	1.45	0.81
2:B:379:CYS:HB2	2:B:432:CYS:HA	1.68	0.75
2:B:350:VAL:HG21	2:B:418:ILE:HD11	1.73	0.71
2:B:422:ASN:HD21	2:B:454:ARG:H	1.37	0.71
1:A:283:PRO:HG3	1:A:439:LEU:HD23	1.73	0.70
1:A:238:GLN:OE1	1:A:598:ASN:ND2	2.25	0.69
2:B:339:HIS:ND1	2:B:340:GLU:OE1	2.26	0.66
1:A:151:MET:HB2	1:A:164:TRP:HE1	1.61	0.65
2:B:474:GLN:HE22	2:B:480:CYS:H	1.46	0.64
2:B:442:ASP:OD1	2:B:509:ARG:NH2	2.32	0.63
1:A:76:GLN:NE2	1:A:79:MET:SD	2.73	0.61
1:A:370:THR:O	1:A:374:GLU:HG2	2.00	0.60
1:A:470:GLN:N	1:A:470:GLN:OE1	2.34	0.60
1:A:173:LYS:HA	1:A:176:ARG:HE	1.67	0.60
1:A:216:TYR:CZ	1:A:566:THR:HG21	2.39	0.57
2:B:454:ARG:NH1	2:B:467:ASP:OD2	2.37	0.56
1:A:296:MET:HG3	1:A:301:TRP:CE3	2.40	0.56
1:A:477:TRP:CD2	1:A:488:GLU:HG2	2.40	0.56
1:A:296:MET:HG3	1:A:301:TRP:HE3	1.71	0.56
1:A:572:ILE:HG23	1:A:573:VAL:HG23	1.88	0.55
1:A:232:ILE:HG22	1:A:580:VAL:HG11	1.90	0.53
1:A:405:GLU:OE1	1:A:517:ARG:NH1	2.40	0.53
1:A:490:LEU:H	1:A:612:TYR:HB3	1.73	0.53
2:B:344:ALA:HB3	2:B:347:PHE:HE1	1.73	0.53
1:A:156:ASP:O	1:A:160:ARG:HD2	2.09	0.53
1:A:292:VAL:H	1:A:365:MET:HE3	1.75	0.52
2:B:497:PHE:CZ	2:B:507:PRO:HB3	2.45	0.51
1:A:244:ARG:NH2	1:A:604:GLY:O	2.45	0.50
1:A:292:VAL:H	1:A:365:MET:CE	2.24	0.50
2:B:422:ASN:HD21	2:B:454:ARG:N	2.08	0.50
1:A:596:GLU:O	1:A:599:ARG:NH2	2.45	0.49
1:A:202:TRP:CZ3	1:A:510:SER:HA	2.47	0.49
2:B:342:PHE:CE1	2:B:511:VAL:HG11	2.48	0.49
1:A:227:ARG:NH2	1:A:578:MET:O	2.45	0.49
1:A:542:ASP:OD2	1:A:544:SER:OG	2.22	0.49
1:A:147:LEU:HA	1:A:150:ILE:HG22	1.94	0.48
1:A:134:ASP:OD2	1:A:163:ALA:HA	2.13	0.48
1:A:350:LEU:HB2	1:A:354:ASP:HB3	1.96	0.47
1:A:307:PHE:HZ	1:A:332:LEU:HD13	1.78	0.47
1:A:22:GLU:HG2	1:A:88:ILE:HD12	1.96	0.47
1:A:202:TRP:CH2	1:A:510:SER:HA	2.50	0.47
1:A:401:GLU:HB2	1:A:517:ARG:HD3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:422:ASN:ND2	2:B:454:ARG:H	2.11	0.47
1:A:122:LYS:HA	1:A:122:LYS:HD2	1.74	0.46
1:A:335:PRO:HB3	1:A:339:ARG:CG	2.46	0.46
1:A:397:GLU:OE2	1:A:397:GLU:N	2.47	0.46
1:A:37:GLU:OE2	1:A:392:ARG:NH1	2.49	0.46
1:A:340:LYS:N	1:A:340:LYS:HD3	2.31	0.46
2:B:391:CYS:HB2	2:B:525:CYS:HB3	1.59	0.46
1:A:301:TRP:HZ2	1:A:309:GLU:HG3	1.81	0.45
2:B:439:ASN:OD1	2:B:506:GLN:NE2	2.48	0.45
1:A:403:VAL:HA	1:A:406:ILE:HD12	1.98	0.45
1:A:310:ALA:O	1:A:314:PHE:HD2	1.98	0.45
1:A:563:GLU:OE1	1:A:567:LEU:HB3	2.15	0.45
1:A:534:HIS:CE1	1:A:541:CYS:HG	2.34	0.45
1:A:83:TYR:O	1:A:101:GLN:NE2	2.50	0.45
1:A:23:GLU:O	1:A:27:THR:HG23	2.16	0.45
1:A:366:ASP:O	1:A:370:THR:OG1	2.28	0.45
1:A:375:MET:HA	1:A:378:ILE:HG22	1.98	0.45
2:B:402:ILE:HG23	2:B:403:LYS:O	2.17	0.45
2:B:503:VAL:HA	2:B:506:GLN:HG3	1.97	0.45
1:A:442:ALA:O	1:A:446:VAL:HG22	2.18	0.44
1:A:534:HIS:CG	1:A:541:CYS:HG	2.33	0.44
1:A:50:TYR:CE1	1:A:54:ILE:HG12	2.52	0.44
1:A:292:VAL:HG11	1:A:422:LEU:HB3	1.99	0.44
2:B:461:LEU:HD23	2:B:461:LEU:HA	1.82	0.44
1:A:164:TRP:CE3	1:A:167:TRP:HB2	2.52	0.44
1:A:333:THR:OG1	1:A:334:GLU:N	2.50	0.44
1:A:481:ARG:HH11	1:A:487:VAL:HG23	1.81	0.44
1:A:483:ILE:HD12	1:A:484:VAL:HG23	1.99	0.44
2:B:409:GLN:HG2	2:B:419:ALA:HB2	1.99	0.44
1:A:510:SER:HB2	1:A:513:ARG:HE	1.83	0.44
1:A:523:GLN:HG2	1:A:582:PRO:HG2	2.00	0.43
1:A:108:LEU:HD12	1:A:108:LEU:H	1.83	0.43
1:A:223:LYS:HA	1:A:226:GLU:HG2	2.00	0.43
1:A:108:LEU:HD11	1:A:189:MET:HG2	2.00	0.43
2:B:379:CYS:HB2	2:B:432:CYS:CA	2.35	0.43
1:A:570:GLU:HB2	1:A:576:LYS:NZ	2.34	0.43
1:A:216:TYR:CE1	1:A:566:THR:HG21	2.52	0.43
2:B:474:GLN:NE2	2:B:480:CYS:H	2.14	0.43
1:A:246:LYS:HD3	1:A:246:LYS:HA	1.73	0.43
1:A:534:HIS:ND1	1:A:541:CYS:SG	2.85	0.43
1:A:161:LEU:HD23	1:A:161:LEU:HA	1.77	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:594:LEU:HD12	1:A:594:LEU:HA	1.84	0.43
1:A:354:ASP:OD2	1:A:356:ARG:NE	2.52	0.42
1:A:350:LEU:HD23	1:A:356:ARG:HD2	2.01	0.42
1:A:434:GLU:HB2	1:A:539:PHE:CE2	2.55	0.42
1:A:568:ALA:O	1:A:572:ILE:HG22	2.19	0.42
1:A:431:ASN:O	1:A:434:GLU:HG2	2.19	0.42
1:A:151:MET:HB2	1:A:164:TRP:NE1	2.31	0.42
1:A:554:LEU:HD12	1:A:554:LEU:O	2.19	0.42
2:B:360:ASN:OD1	2:B:523:THR:HB	2.20	0.42
1:A:344:HIS:O	1:A:346:THR:HG23	2.20	0.42
1:A:434:GLU:HB2	1:A:539:PHE:HE2	1.85	0.42
1:A:108:LEU:HB3	1:A:112:LYS:HB2	2.02	0.41
2:B:474:GLN:OE1	2:B:480:CYS:N	2.53	0.41
2:B:342:PHE:HB3	2:B:371:PHE:CE2	2.56	0.41
1:A:171:VAL:O	1:A:175:LEU:HD13	2.20	0.41
1:A:329:ASN:OD1	1:A:356:ARG:HG2	2.20	0.41
1:A:374:GLU:HA	1:A:377:HIS:ND1	2.36	0.41
1:A:548:GLU:HG2	1:A:549:ALA:N	2.35	0.41
2:B:477:ASN:OD1	2:B:477:ASN:N	2.52	0.41
1:A:147:LEU:HG	1:A:164:TRP:CZ2	2.55	0.41
1:A:263:ALA:N	1:A:487:VAL:O	2.53	0.41
1:A:489:PRO:HB2	1:A:490:LEU:HD12	2.02	0.41
1:A:233:LYS:HB3	1:A:233:LYS:HE3	1.85	0.41
1:A:380:TYR:CD1	1:A:557:LEU:HD22	2.56	0.41
1:A:406:ILE:HG13	1:A:521:GLN:HB2	2.02	0.41
2:B:427:ASP:OD1	2:B:427:ASP:N	2.53	0.41
1:A:301:TRP:NE1	1:A:305:ARG:HB3	2.24	0.41
1:A:456:GLU:HG2	1:A:511:PHE:HB3	2.02	0.41
2:B:403:LYS:HB3	2:B:403:LYS:HE2	1.78	0.41
1:A:141:LEU:HD12	1:A:141:LEU:HA	1.89	0.41
1:A:314:PHE:O	1:A:317:ILE:HG22	2.22	0.40
1:A:420:LEU:HD12	1:A:420:LEU:HA	1.94	0.40
1:A:530:LYS:HE3	1:A:530:LYS:HB3	1.90	0.40
2:B:492:LEU:H	2:B:492:LEU:HG	1.74	0.40
1:A:291:ASP:OD1	1:A:291:ASP:N	2.54	0.40
1:A:317:ILE:HD12	1:A:317:ILE:HA	1.81	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	594/786 (76%)	558 (94%)	34 (6%)	2 (0%)	37	71
2	B	193/211 (92%)	176 (91%)	17 (9%)	0	100	100
All	All	787/997 (79%)	734 (93%)	51 (6%)	2 (0%)	38	71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	255	ILE
1	A	490	LEU

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	525/697 (75%)	507 (97%)	18 (3%)	32	54
2	B	169/185 (91%)	161 (95%)	8 (5%)	22	46
All	All	694/882 (79%)	668 (96%)	26 (4%)	31	52

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	LEU
1	A	167	TRP
1	A	176	ARG
1	A	197	ASP

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Mol	Chain	Res	Type
1	A	248	MET
1	A	270	TRP
1	A	307	PHE
1	A	399	PHE
1	A	400	HIS
1	A	441	GLN
1	A	472	TRP
1	A	479	MET
1	A	490	LEU
1	A	511	PHE
1	A	529	CYS
1	A	563	GLU
1	A	599	ARG
1	A	605	TRP
2	B	379	CYS
2	B	380	TYR
2	B	384	PRO
2	B	427	ASP
2	B	429	PHE
2	B	438	SER
2	B	495	TYR
2	B	514	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	422	ASN

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](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 for Mogul analysis.

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
4	NAG	A	902	1	14,14,15	0.41	0	17,19,21	0.42	0
4	NAG	B	601	2	14,14,15	0.58	0	17,19,21	1.06	1 (5%)
4	NAG	A	906	1	14,14,15	0.25	0	17,19,21	0.57	0
4	NAG	A	903	1	14,14,15	0.32	0	17,19,21	0.48	0
4	NAG	A	904	-	14,14,15	0.37	0	17,19,21	0.71	1 (5%)
4	NAG	A	905	1	14,14,15	0.16	0	17,19,21	0.48	0
4	NAG	B	602	2	14,14,15	0.63	0	17,19,21	1.69	2 (11%)

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
4	NAG	A	902	1	-	2/6/23/26	0/1/1/1
4	NAG	B	601	2	-	4/6/23/26	0/1/1/1
4	NAG	A	906	1	-	3/6/23/26	0/1/1/1
4	NAG	A	903	1	-	0/6/23/26	0/1/1/1
4	NAG	A	904	-	-	2/6/23/26	0/1/1/1
4	NAG	A	905	1	-	3/6/23/26	0/1/1/1
4	NAG	B	602	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	602	NAG	C4-C3-C2	-4.20	104.86	111.02
4	B	602	NAG	O5-C1-C2	-3.95	105.05	111.29
4	B	601	NAG	O5-C1-C2	2.25	114.84	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	904	NAG	C1-O5-C5	2.10	115.04	112.19

There are no chirality outliers.

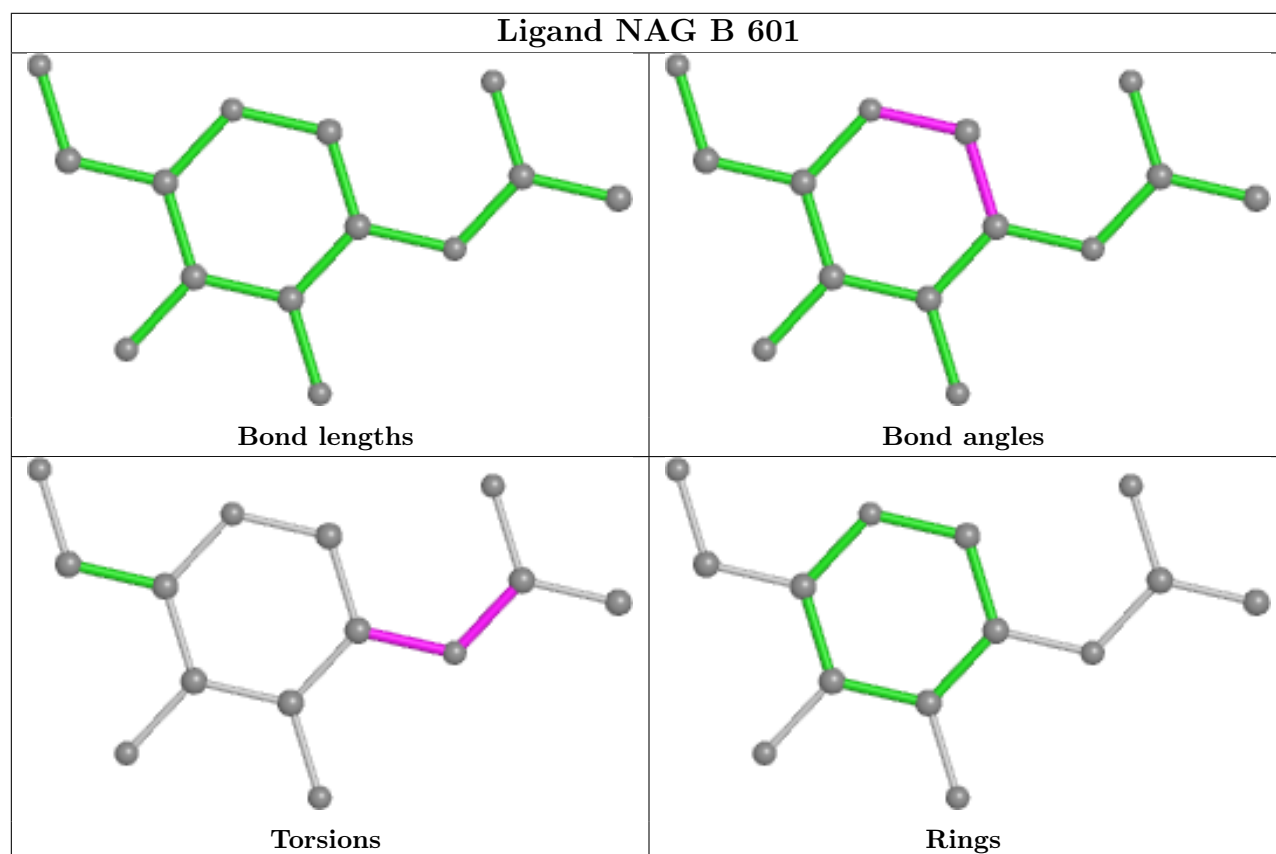
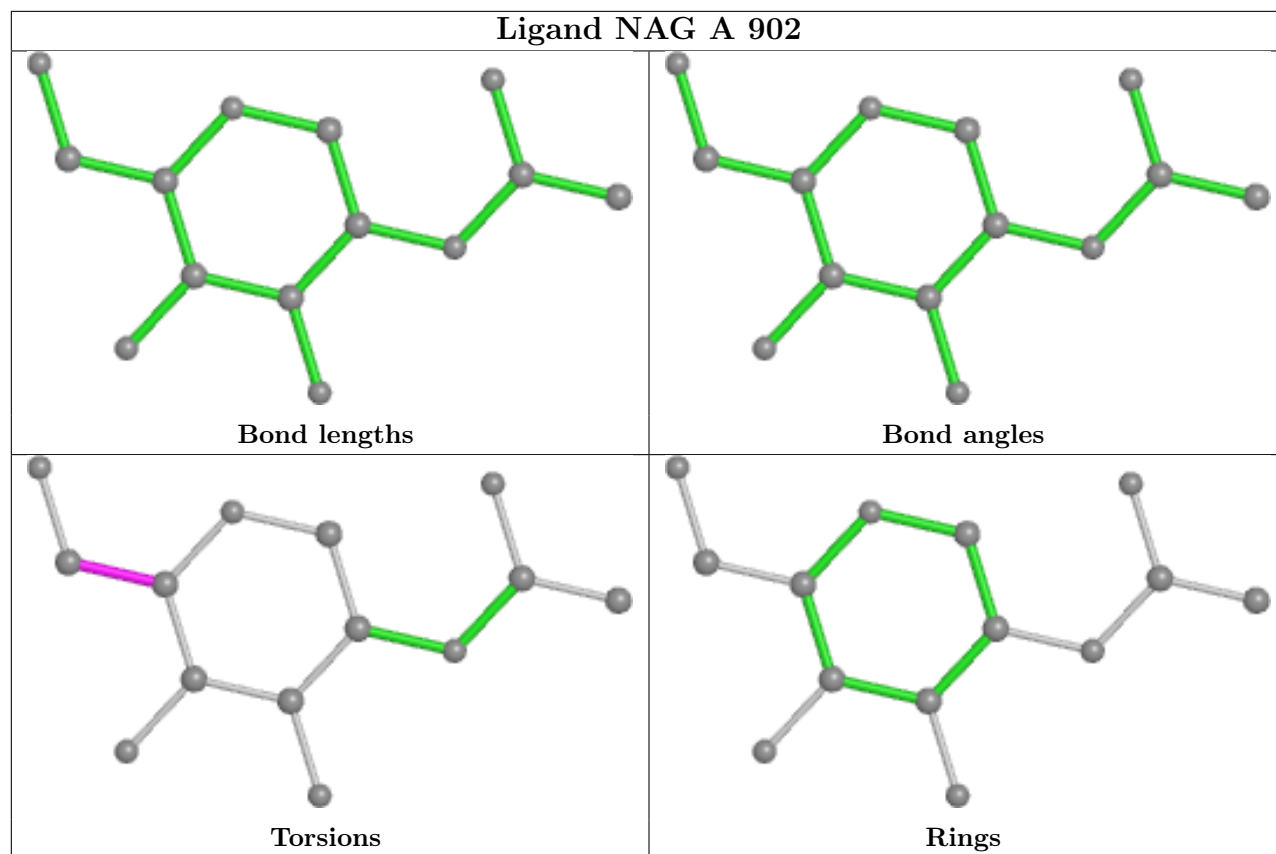
All (16) torsion outliers are listed below:

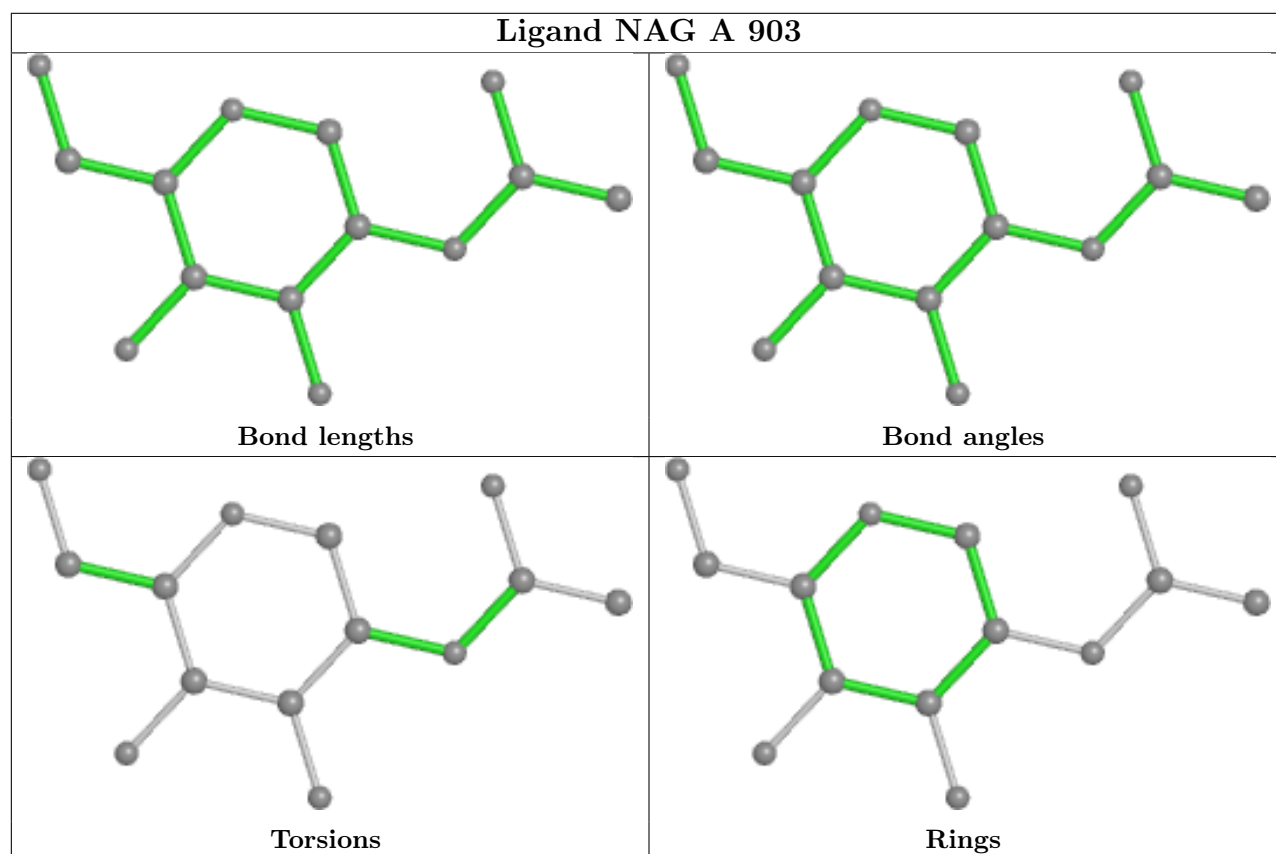
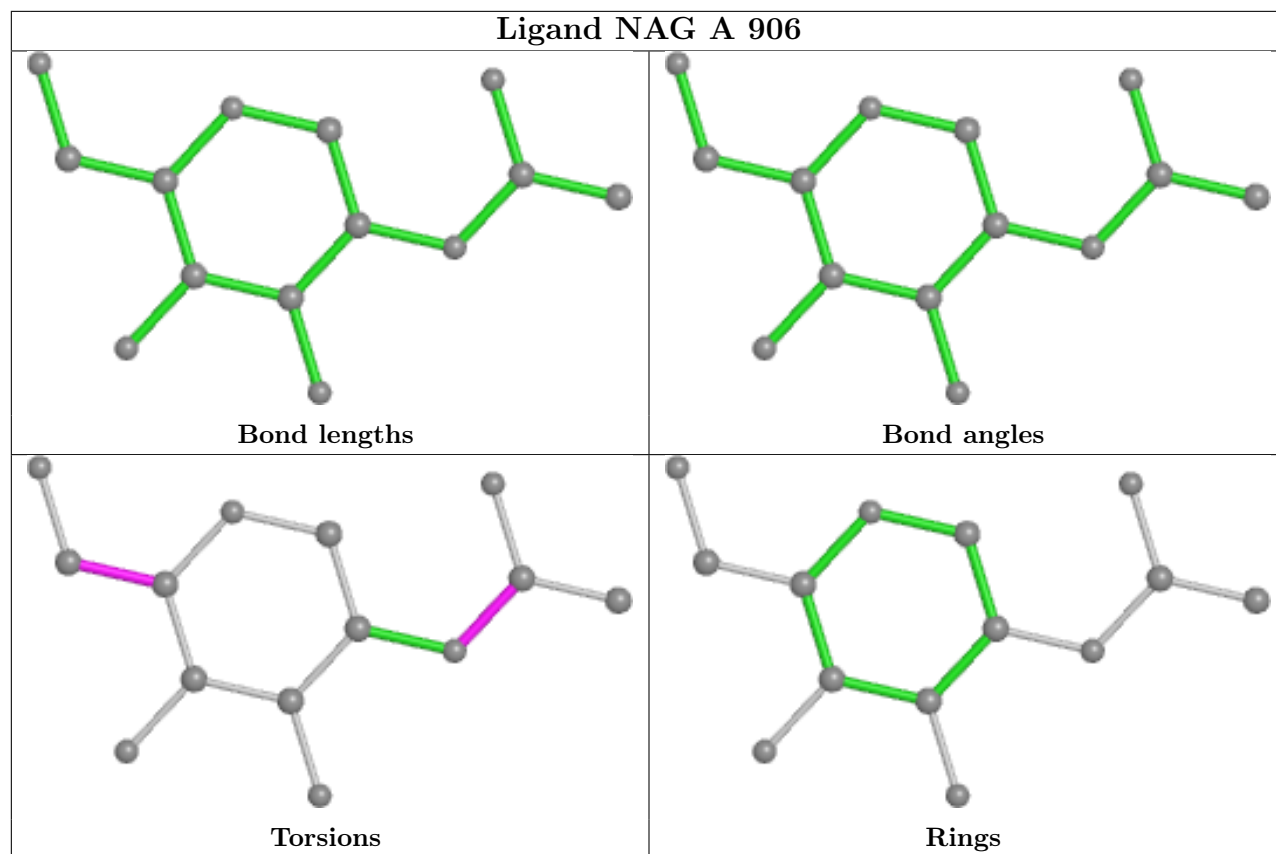
Mol	Chain	Res	Type	Atoms
4	B	601	NAG	C8-C7-N2-C2
4	B	601	NAG	O7-C7-N2-C2
4	A	902	NAG	O5-C5-C6-O6
4	A	904	NAG	C8-C7-N2-C2
4	A	904	NAG	O7-C7-N2-C2
4	A	905	NAG	C8-C7-N2-C2
4	A	905	NAG	O7-C7-N2-C2
4	A	906	NAG	C8-C7-N2-C2
4	A	906	NAG	O7-C7-N2-C2
4	A	905	NAG	O5-C5-C6-O6
4	B	602	NAG	C8-C7-N2-C2
4	A	902	NAG	C4-C5-C6-O6
4	A	906	NAG	O5-C5-C6-O6
4	B	601	NAG	C1-C2-N2-C7
4	B	602	NAG	O7-C7-N2-C2
4	B	601	NAG	C3-C2-N2-C7

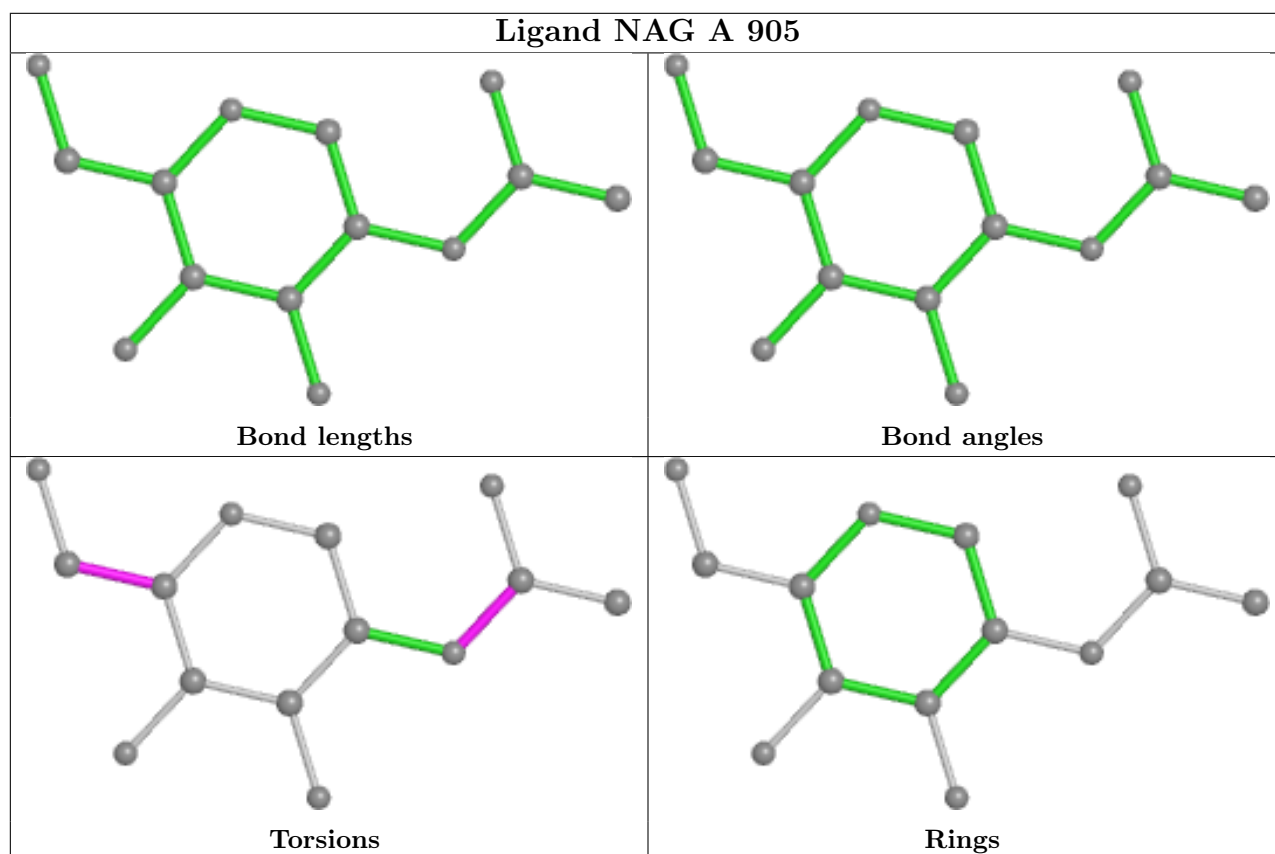
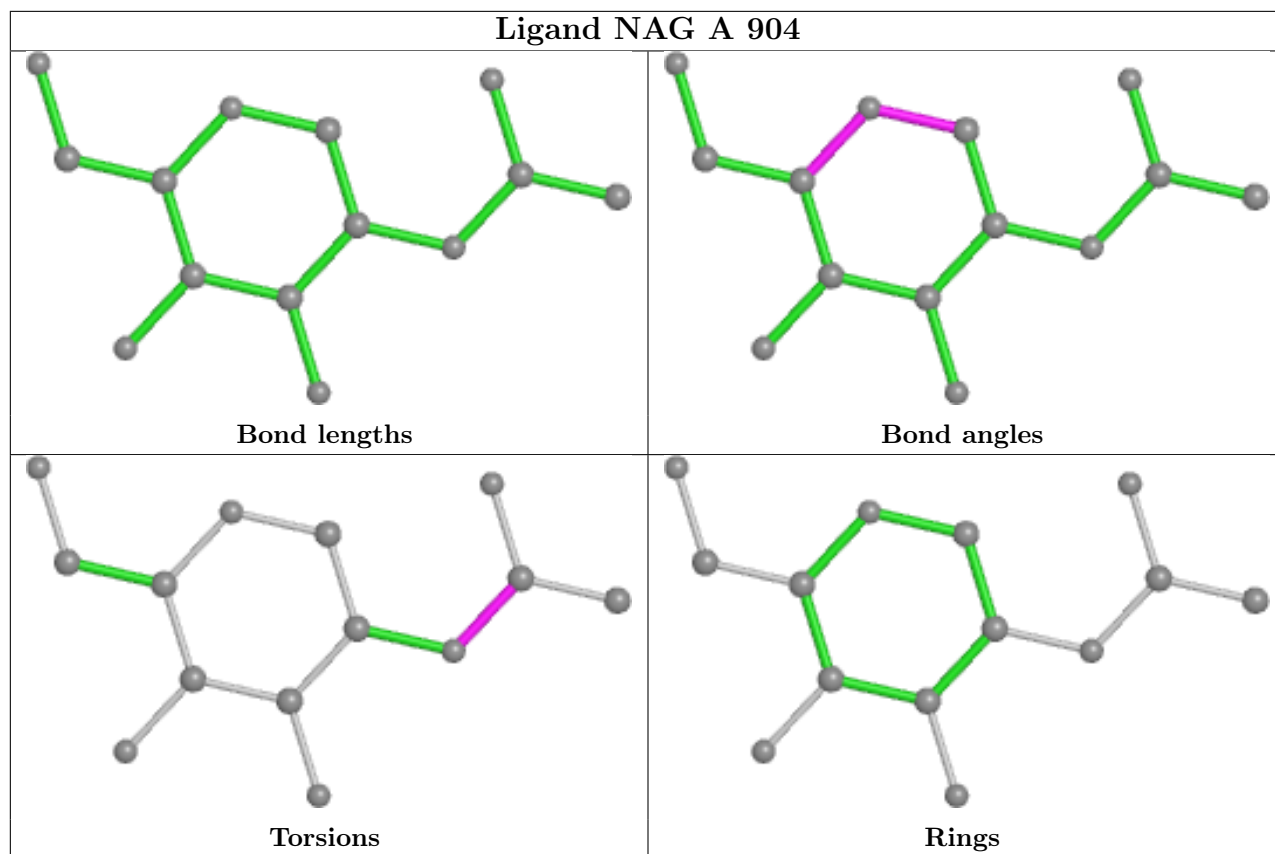
There are no ring outliers.

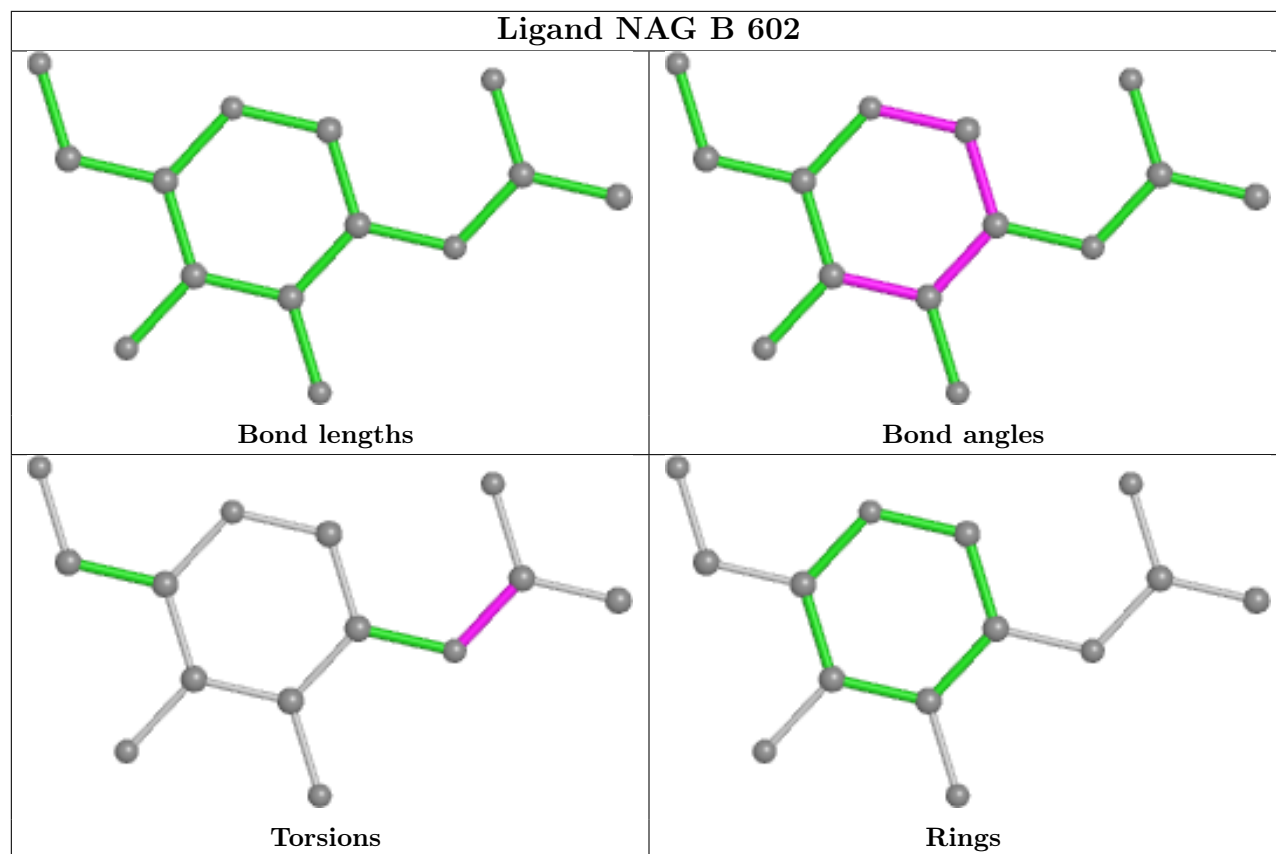
No monomer is involved in short contacts.

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.