



wwPDB EM Validation Summary Report ⓘ

Mar 5, 2026 – 04:52 PM UTC

PDB ID : 8SQW / pdb_00008sqw
EMDB ID : EMD-40714
Title : particulate methane monooxygenase crosslinked with 2,2,2-trifluoroethanol bound
Authors : Tucci, F.J.; Rosenzweig, A.C.
Deposited on : 2023-05-04
Resolution : 2.16 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at
<https://www.wwpdb.org/validation/2017/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

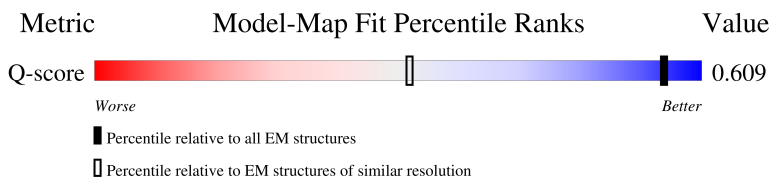
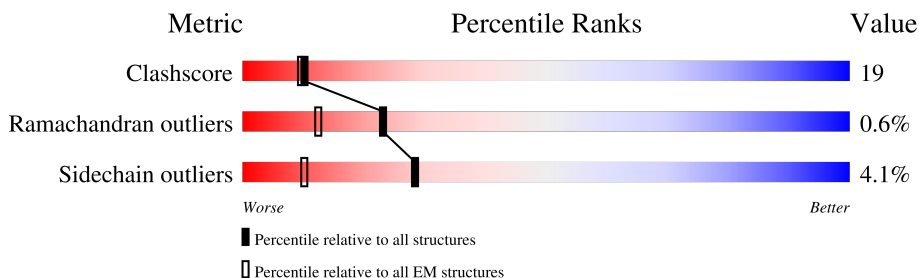
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.16 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.



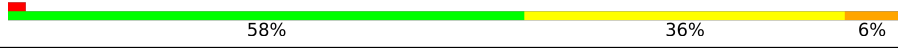
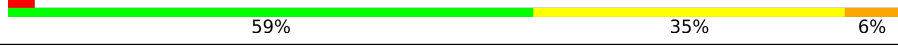
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2609 (1.67 - 2.66)

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	382	 82% 17% .
1	E	382	 80% 19% .
1	I	382	 81% 18% .
2	B	241	 69% 29% .

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Mol	Chain	Length	Quality of chain
2	F	241	
2	J	241	
3	C	236	
3	G	236	
3	K	236	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	P1O	B	307	-	-	X	-
7	P1O	F	308	-	-	X	-
7	P1O	J	301	-	-	X	-
9	ETF	C	310	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 25319 atoms, of which 2469 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 Particulate methane monooxygenase alpha subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	382	Total 3017	C 1938	N 513	O 551	S 15	0	0
1	E	382	Total 3017	C 1938	N 513	O 551	S 15	0	0
1	I	382	Total 3017	C 1938	N 513	O 551	S 15	0	0

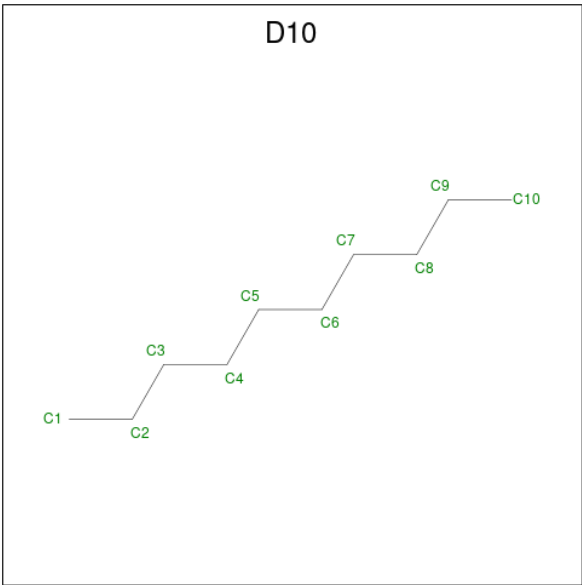
- Molecule 2 is a protein called Particulate methane monooxygenase beta subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	241	Total 1977	C 1329	N 315	O 322	S 11	0	0
2	F	241	Total 1977	C 1329	N 315	O 322	S 11	0	0
2	J	241	Total 1977	C 1329	N 315	O 322	S 11	0	0

- Molecule 3 is a protein called Ammonia monooxygenase/methane monooxygenase, subunit C family protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	236	Total 1972	C 1339	N 299	O 329	S 5	0	0
3	G	236	Total 1972	C 1339	N 299	O 329	S 5	0	0
3	K	236	Total 1972	C 1339	N 299	O 329	S 5	0	0

- Molecule 4 is DECANE (CCD ID: D10) (formula: C₁₀H₂₂).



Mol	Chain	Residues	Atoms			AltConf
4	A	1	Total	C	H	0
			32	10	22	
4	B	1	Total	C	H	0
			32	10	22	
4	B	1	Total	C	H	0
			32	10	22	
4	B	1	Total	C	H	0
			32	10	22	
4	B	1	Total	C	H	0
			32	10	22	
4	C	1	Total	C	H	0
			32	10	22	
4	E	1	Total	C	H	0
			32	10	22	
4	I	1	Total	C	H	0
			32	10	22	
4	F	1	Total	C	H	0
			32	10	22	
4	F	1	Total	C	H	0
			32	10	22	
4	F	1	Total	C	H	0
			32	10	22	
4	F	1	Total	C	H	0
			32	10	22	
4	J	1	Total	C	H	0
			32	10	22	
4	J	1	Total	C	H	0
			32	10	22	

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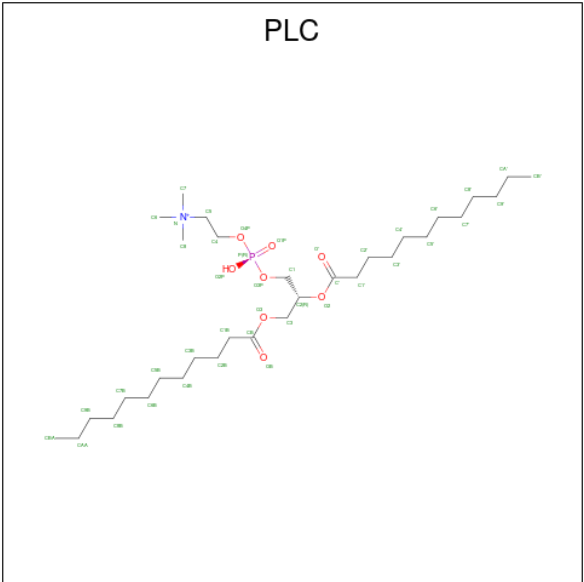
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Mol	Chain	Residues	Atoms			AltConf
4	J	1	Total	C	H	0
			32	10	22	
4	J	1	Total	C	H	0
			32	10	22	
4	G	1	Total	C	H	0
			32	10	22	
4	K	1	Total	C	H	0
			32	10	22	

- Molecule 5 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

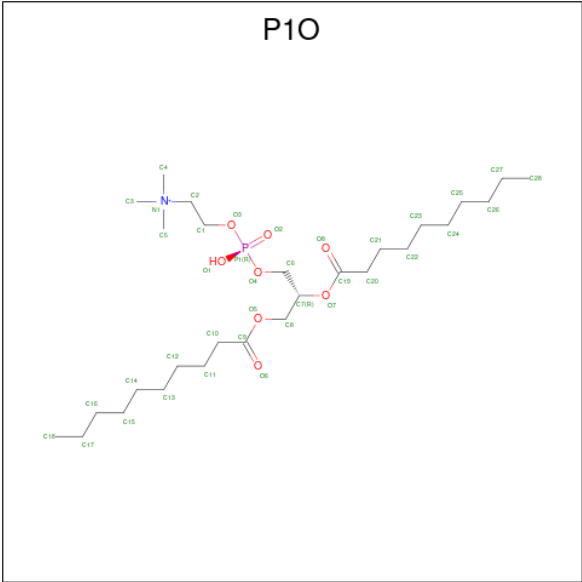
Mol	Chain	Residues	Atoms		AltConf
5	A	2	Total	Cu	0
			2	2	
5	C	1	Total	Cu	0
			1	1	
5	E	2	Total	Cu	0
			2	2	
5	I	2	Total	Cu	0
			2	2	
5	G	1	Total	Cu	0
			1	1	
5	K	1	Total	Cu	0
			1	1	

- Molecule 6 is DIUNDECYL PHOSPHATIDYL CHOLINE (CCD ID: PLC) (formula: C₃₂H₆₅NO₈P).



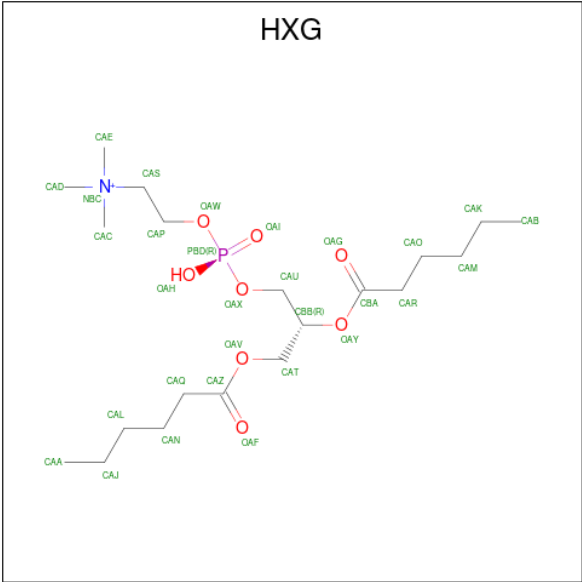
Mol	Chain	Residues	Atoms						AltConf
6	B	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	B	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	C	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	C	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	C	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	C	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	F	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	F	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	J	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	J	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	G	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	G	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	G	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	G	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	K	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	K	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	K	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	K	1	Total 106	C 32	H 64	N 1	O 8	P 1	0

- Molecule 7 is 1,2-DIDECANOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: P1O) (formula: C₂₈H₅₇NO₈P).



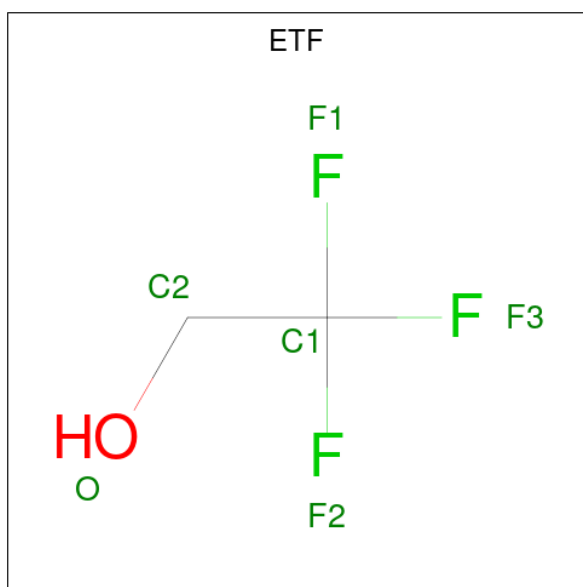
Mol	Chain	Residues	Atoms						AltConf
7	B	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	B	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	C	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	C	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	F	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	F	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	J	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	J	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	G	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	G	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	K	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	K	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	

- Molecule 8 is 1,2-dihexanoyl-sn-glycero-3-phosphocholine (CCD ID: HXG) (formula: C₂₀H₄₁NO₈P).



Mol	Chain	Residues	Atoms						AltConf
8	C	1	Total	C	H	N	O	P	0
			70	20	40	1	8	1	
8	C	1	Total	C	H	N	O	P	0
			70	20	40	1	8	1	
8	G	1	Total	C	H	N	O	P	0
			70	20	40	1	8	1	
8	G	1	Total	C	H	N	O	P	0
			70	20	40	1	8	1	
8	K	1	Total	C	H	N	O	P	0
			70	20	40	1	8	1	
8	K	1	Total	C	H	N	O	P	0
			70	20	40	1	8	1	

- Molecule 9 is TRIFLUOROETHANOL (CCD ID: ETF) (formula: C₂H₃F₃O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
9	C	1	Total	C	F	H	O	0
			9	2	3	3	1	
9	G	1	Total	C	F	H	O	0
			9	2	3	3	1	
9	K	1	Total	C	F	H	O	0
			9	2	3	3	1	

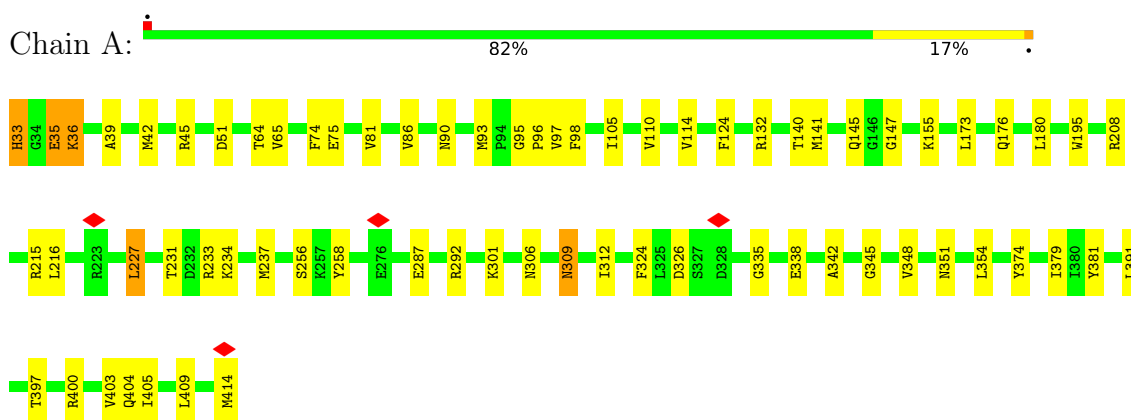
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		AltConf
10	A	70	Total	O	0
			70	70	
10	B	35	Total	O	0
			35	35	
10	C	14	Total	O	0
			14	14	
10	E	70	Total	O	0
			70	70	
10	I	69	Total	O	0
			69	69	
10	F	34	Total	O	0
			34	34	
10	J	36	Total	O	0
			36	36	
10	G	13	Total	O	0
			13	13	
10	K	12	Total	O	0
			12	12	

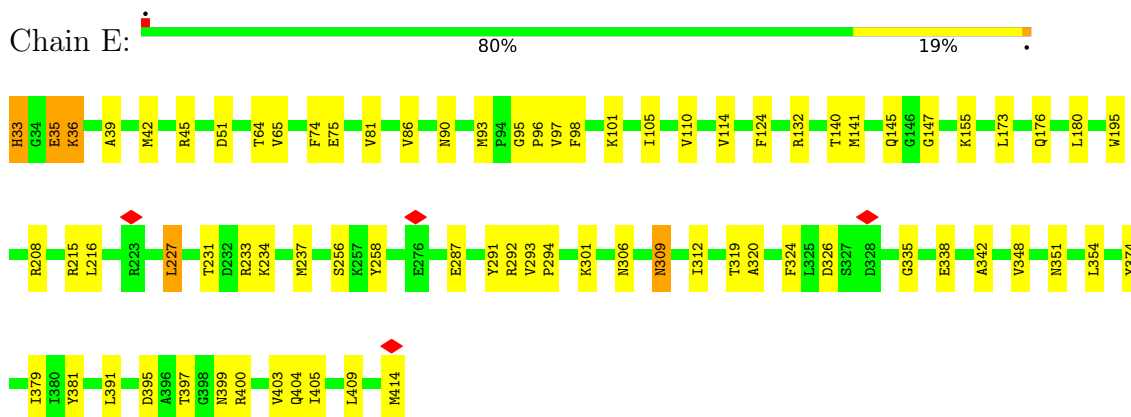
3 Residue-property plots

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.

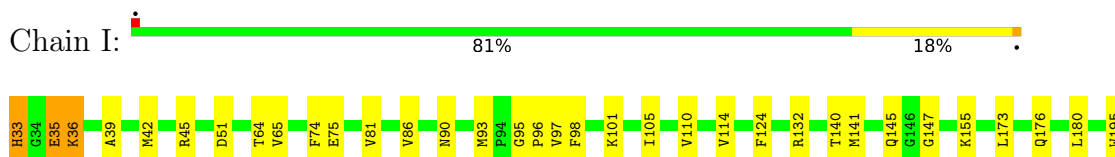
- Molecule 1: Particulate methane monooxygenase alpha subunit

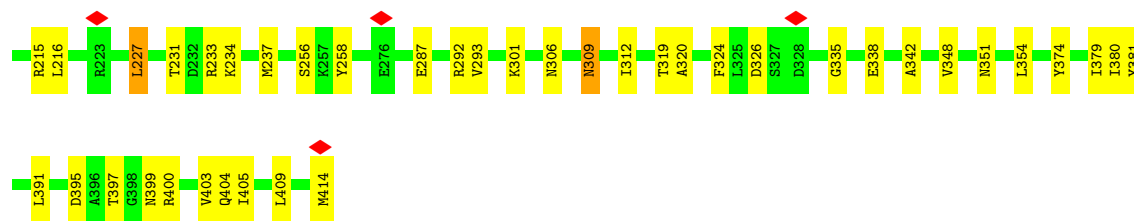


- Molecule 1: Particulate methane monooxygenase alpha subunit

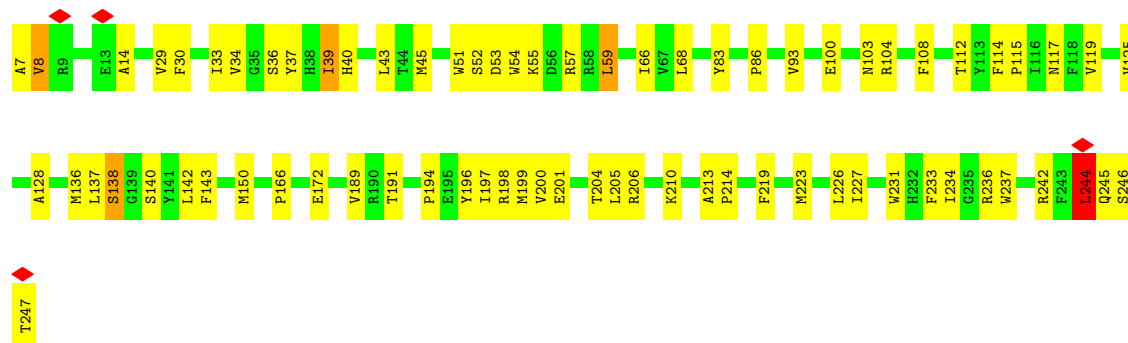


- Molecule 1: Particulate methane monooxygenase alpha subunit

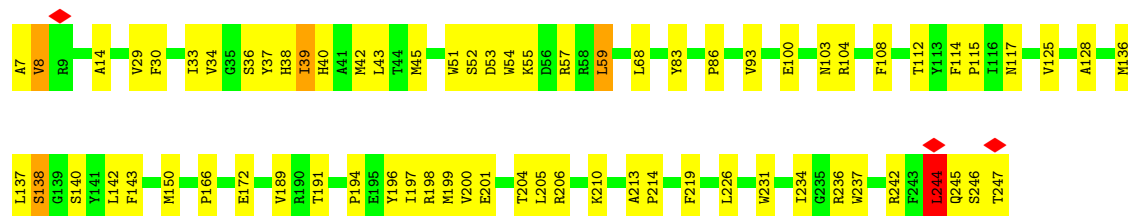




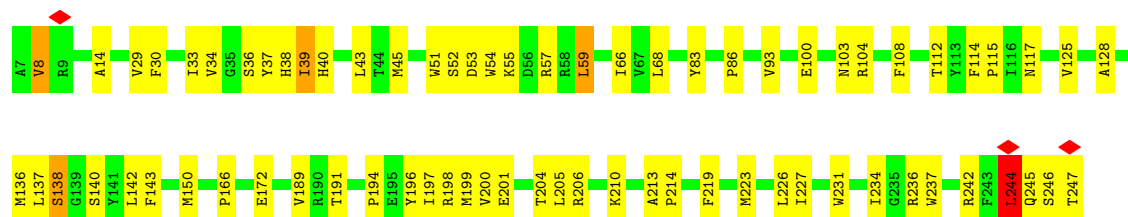
- Molecule 2: Particulate methane monooxygenase beta subunit



- Molecule 2: Particulate methane monooxygenase beta subunit

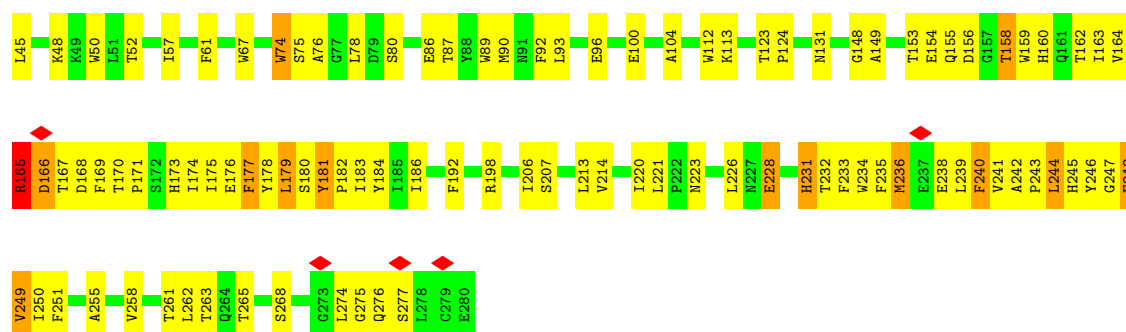


- Molecule 2: Particulate methane monooxygenase beta subunit



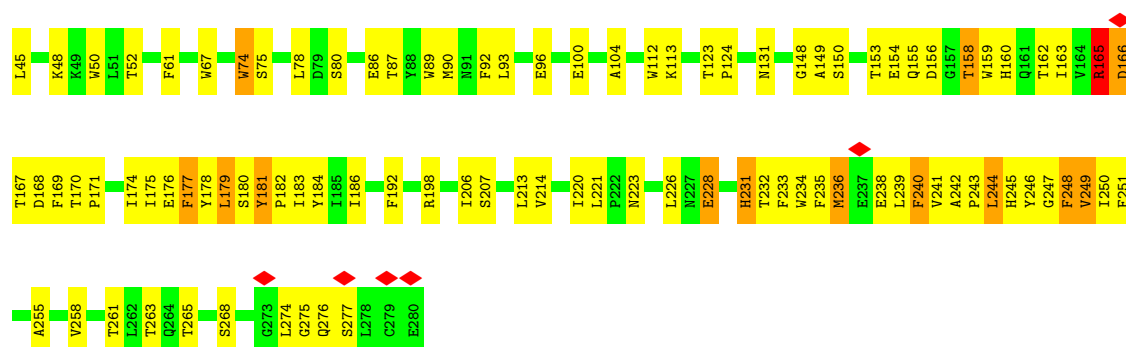
- Molecule 3: Ammonia monooxygenase/methane monooxygenase, subunit C family protein





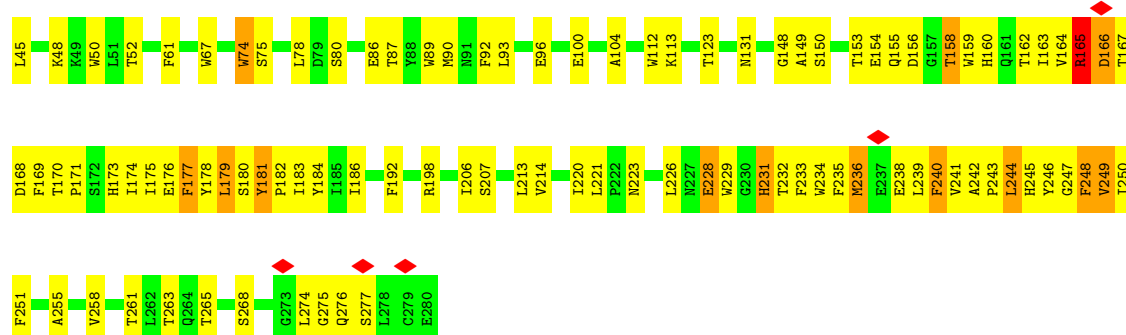
- Molecule 3: Ammonia monooxygenase/methane monooxygenase, subunit C family protein

Chain G: 59% 35% 6%



- Molecule 3: Ammonia monooxygenase/methane monooxygenase, subunit C family protein

Chain K: 58% 36% 6%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	500000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.226	Depositor
Minimum map value	-0.110	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.0342	Depositor
Map size (Å)	270.8992, 270.8992, 270.8992	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.5291, 0.5291, 0.5291	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: P1O, HXG, CU, ETF, PLC, D10

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	0.22	0/3099	0.41	0/4215
1	E	0.22	0/3099	0.41	0/4215
1	I	0.22	0/3099	0.41	0/4215
2	B	0.25	0/2053	0.43	0/2810
2	F	0.25	0/2053	0.43	0/2810
2	J	0.25	0/2053	0.43	0/2810
3	C	0.45	0/2051	0.77	6/2810 (0.2%)
3	G	0.45	0/2051	0.77	6/2810 (0.2%)
3	K	0.45	0/2051	0.77	6/2810 (0.2%)
All	All	0.31	0/21609	0.54	18/29505 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	C	0	1
3	G	0	1
3	K	0	1
All	All	0	3

There are no bond length outliers.

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	248	PHE	CA-C-N	-8.51	109.66	120.56
3	G	248	PHE	C-N-CA	-8.51	109.66	120.56
3	K	248	PHE	CA-C-N	-8.51	109.67	120.56
3	K	248	PHE	C-N-CA	-8.51	109.67	120.56

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	C	248	PHE	CA-C-N	-8.50	109.68	120.56

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	165	ARG	Sidechain
3	G	165	ARG	Sidechain
3	K	165	ARG	Sidechain

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	3017	0	2980	50	0
1	E	3017	0	2980	55	0
1	I	3017	0	2980	52	0
2	B	1977	0	1936	111	0
2	F	1977	0	1936	106	0
2	J	1977	0	1936	109	0
3	C	1972	0	1904	140	0
3	G	1972	0	1904	139	0
3	K	1972	0	1904	138	0
4	A	10	22	22	0	0
4	B	40	88	88	9	0
4	C	10	22	22	3	0
4	E	10	22	22	0	0
4	F	40	88	88	7	0
4	G	10	22	22	3	0
4	I	10	22	22	0	0
4	J	40	88	88	8	0
4	K	10	22	22	3	0
5	A	2	0	0	0	0
5	C	1	0	0	0	0
5	E	2	0	0	0	0
5	G	1	0	0	0	0
5	I	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	K	1	0	0	0	0
6	B	84	128	128	3	0
6	C	168	256	256	19	0
6	F	84	128	128	3	0
6	G	168	256	256	19	0
6	J	84	128	128	3	0
6	K	168	256	256	21	0
7	B	76	112	112	35	0
7	C	76	112	112	21	0
7	F	76	112	112	37	0
7	G	76	112	112	23	0
7	J	76	112	112	34	0
7	K	76	112	112	22	0
8	C	60	80	80	23	0
8	G	60	80	80	23	0
8	K	60	80	80	24	0
9	C	6	3	2	4	0
9	G	6	3	2	1	0
9	K	6	3	2	1	0
10	A	70	0	0	4	0
10	B	35	0	0	2	0
10	C	14	0	0	7	0
10	E	70	0	0	4	0
10	F	34	0	0	5	0
10	G	13	0	0	4	0
10	I	69	0	0	3	0
10	J	36	0	0	2	0
10	K	12	0	0	4	0
All	All	22850	2469	22926	871	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 19.

The worst 5 of 871 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)
3:C:67:TRP:CD1	8:C:306:HXG:H41	1.81	1.16
3:G:67:TRP:CD1	8:G:307:HXG:H41	1.81	1.16
3:K:67:TRP:CD1	8:K:307:HXG:H41	1.81	1.15
6:K:303:PLC:H73	8:K:304:HXG:H36	1.32	1.10
6:G:303:PLC:H73	8:G:304:HXG:H36	1.32	1.10

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	380/382 (100%)	364 (96%)	16 (4%)	0	100	100
1	E	380/382 (100%)	364 (96%)	16 (4%)	0	100	100
1	I	380/382 (100%)	365 (96%)	15 (4%)	0	100	100
2	B	239/241 (99%)	227 (95%)	11 (5%)	1 (0%)	30	26
2	F	239/241 (99%)	227 (95%)	11 (5%)	1 (0%)	30	26
2	J	239/241 (99%)	228 (95%)	10 (4%)	1 (0%)	30	26
3	C	234/236 (99%)	204 (87%)	26 (11%)	4 (2%)	7	2
3	G	234/236 (99%)	204 (87%)	26 (11%)	4 (2%)	7	2
3	K	234/236 (99%)	205 (88%)	25 (11%)	4 (2%)	7	2
All	All	2559/2577 (99%)	2388 (93%)	156 (6%)	15 (1%)	23	15

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	277	SER
3	G	277	SER
3	K	277	SER
2	B	244	LEU
3	C	233	PHE

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	323/323 (100%)	313 (97%)	10 (3%)	35	37
1	E	323/323 (100%)	313 (97%)	10 (3%)	35	37
1	I	323/323 (100%)	313 (97%)	10 (3%)	35	37
2	B	206/206 (100%)	201 (98%)	5 (2%)	43	47
2	F	206/206 (100%)	201 (98%)	5 (2%)	43	47
2	J	206/206 (100%)	201 (98%)	5 (2%)	43	47
3	C	200/200 (100%)	185 (92%)	15 (8%)	12	8
3	G	200/200 (100%)	185 (92%)	15 (8%)	12	8
3	K	200/200 (100%)	185 (92%)	15 (8%)	12	8
All	All	2187/2187 (100%)	2097 (96%)	90 (4%)	28	26

5 of 90 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	J	59	LEU
3	G	236	MET
2	J	244	LEU
3	G	158	THR
3	K	45	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
2	J	174	ASN
2	J	187	ASN
1	E	168	ASN
1	I	168	ASN
2	F	174	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 66 ligands modelled in this entry, 9 are monoatomic - leaving 57 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	$\# Z > 2$
6	PLC	C	305	-	41,41,41	0.90	0	47,49,49	0.68	1 (2%)
6	PLC	K	306	-	41,41,41	0.90	0	47,49,49	0.68	1 (2%)
8	HXG	K	304	-	29,29,29	0.34	0	35,37,37	0.36	0
6	PLC	G	301	-	41,41,41	1.05	2 (4%)	47,49,49	1.06	4 (8%)
6	PLC	J	308	-	41,41,41	0.84	0	47,49,49	0.75	1 (2%)
7	P1O	B	302	-	37,37,37	1.11	2 (5%)	43,45,45	1.12	3 (6%)
8	HXG	K	307	-	29,29,29	0.35	0	35,37,37	0.37	0
6	PLC	K	301	-	41,41,41	1.05	2 (4%)	47,49,49	1.05	3 (6%)
7	P1O	J	303	-	37,37,37	1.11	2 (5%)	43,45,45	1.11	3 (6%)
4	D10	B	303	-	9,9,9	0.20	0	8,8,8	0.56	0
8	HXG	C	306	-	29,29,29	0.36	0	35,37,37	0.37	0
4	D10	I	501	-	9,9,9	0.20	0	8,8,8	0.56	0
7	P1O	C	309	-	37,37,37	1.13	2 (5%)	43,45,45	1.14	3 (6%)
4	D10	E	501	-	9,9,9	0.20	0	8,8,8	0.55	0
6	PLC	C	311	-	41,41,41	1.05	2 (4%)	47,49,49	1.06	4 (8%)
7	P1O	F	308	-	37,37,37	1.12	2 (5%)	43,45,45	1.10	3 (6%)
7	P1O	B	307	-	37,37,37	1.12	2 (5%)	43,45,45	1.10	3 (6%)
6	PLC	F	301	-	41,41,41	0.84	0	47,49,49	0.75	1 (2%)
7	P1O	C	308	-	37,37,37	1.11	2 (5%)	43,45,45	1.13	4 (9%)
4	D10	A	501	-	9,9,9	0.20	0	8,8,8	0.55	0
6	PLC	B	301	-	41,41,41	0.91	0	47,49,49	0.84	1 (2%)
6	PLC	J	302	-	41,41,41	0.91	0	47,49,49	0.84	1 (2%)
7	P1O	J	301	-	37,37,37	1.12	2 (5%)	43,45,45	1.10	3 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	P1O	K	309	-	37,37,37	1.10	2 (5%)	43,45,45	1.13	4 (9%)
4	D10	C	304	-	9,9,9	0.21	0	8,8,8	0.57	0
4	D10	G	305	-	9,9,9	0.21	0	8,8,8	0.57	0
6	PLC	K	303	-	41,41,41	1.06	2 (4%)	47,49,49	1.11	4 (8%)
8	HXG	G	304	-	29,29,29	0.34	0	35,37,37	0.37	0
9	ETF	G	311	5	5,5,5	0.27	0	7,7,7	0.10	0
4	D10	K	305	-	9,9,9	0.21	0	8,8,8	0.57	0
9	ETF	K	311	5	5,5,5	0.27	0	7,7,7	0.10	0
6	PLC	C	302	-	41,41,41	1.06	2 (4%)	47,49,49	1.11	3 (6%)
6	PLC	B	308	-	41,41,41	0.84	0	47,49,49	0.76	1 (2%)
8	HXG	C	303	-	29,29,29	0.34	0	35,37,37	0.36	0
4	D10	F	306	-	9,9,9	0.20	0	8,8,8	0.55	0
4	D10	B	305	-	9,9,9	0.21	0	8,8,8	0.55	0
4	D10	J	307	-	9,9,9	0.20	0	8,8,8	0.56	0
4	D10	J	305	-	9,9,9	0.21	0	8,8,8	0.56	0
4	D10	F	307	-	9,9,9	0.19	0	8,8,8	0.55	0
6	PLC	G	303	-	41,41,41	1.06	2 (4%)	47,49,49	1.10	4 (8%)
4	D10	F	305	-	9,9,9	0.21	0	8,8,8	0.56	0
4	D10	B	304	-	9,9,9	0.21	0	8,8,8	0.56	0
4	D10	J	306	-	9,9,9	0.21	0	8,8,8	0.55	0
6	PLC	G	306	-	41,41,41	0.90	0	47,49,49	0.68	1 (2%)
4	D10	J	304	-	9,9,9	0.20	0	8,8,8	0.56	0
7	P1O	G	310	-	37,37,37	1.13	2 (5%)	43,45,45	1.13	3 (6%)
7	P1O	F	303	-	37,37,37	1.11	2 (5%)	43,45,45	1.11	3 (6%)
7	P1O	G	309	-	37,37,37	1.11	2 (5%)	43,45,45	1.13	4 (9%)
4	D10	F	304	-	9,9,9	0.20	0	8,8,8	0.56	0
4	D10	B	306	-	9,9,9	0.19	0	8,8,8	0.55	0
6	PLC	G	308	-	41,41,41	0.89	0	47,49,49	0.69	2 (4%)
7	P1O	K	310	-	37,37,37	1.12	2 (5%)	43,45,45	1.14	3 (6%)
8	HXG	G	307	-	29,29,29	0.36	0	35,37,37	0.37	0
6	PLC	K	308	-	41,41,41	0.89	0	47,49,49	0.69	2 (4%)
6	PLC	C	307	-	41,41,41	0.89	0	47,49,49	0.69	2 (4%)
6	PLC	F	302	-	41,41,41	0.91	0	47,49,49	0.85	1 (2%)
9	ETF	C	310	5	5,5,5	0.27	0	7,7,7	0.10	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	PLC	C	305	-	-	14/45/45/45	-
6	PLC	K	306	-	-	14/45/45/45	-
8	HXG	K	304	-	-	9/33/33/33	-
6	PLC	G	301	-	-	25/45/45/45	-
6	PLC	J	308	-	-	15/45/45/45	-
7	P1O	B	302	-	-	26/41/41/41	-
8	HXG	K	307	-	-	7/33/33/33	-
6	PLC	K	301	-	-	25/45/45/45	-
7	P1O	J	303	-	-	26/41/41/41	-
4	D10	B	303	-	-	0/7/7/7	-
8	HXG	C	306	-	-	7/33/33/33	-
4	D10	I	501	-	-	0/7/7/7	-
7	P1O	C	309	-	-	26/41/41/41	-
4	D10	E	501	-	-	0/7/7/7	-
6	PLC	C	311	-	-	25/45/45/45	-
7	P1O	F	308	-	-	20/41/41/41	-
7	P1O	B	307	-	-	20/41/41/41	-
6	PLC	F	301	-	-	15/45/45/45	-
7	P1O	C	308	-	-	19/41/41/41	-
4	D10	A	501	-	-	0/7/7/7	-
6	PLC	B	301	-	-	15/45/45/45	-
6	PLC	J	302	-	-	15/45/45/45	-
7	P1O	J	301	-	-	20/41/41/41	-
7	P1O	K	309	-	-	19/41/41/41	-
4	D10	C	304	-	-	0/7/7/7	-
4	D10	G	305	-	-	0/7/7/7	-
6	PLC	K	303	-	-	22/45/45/45	-
8	HXG	G	304	-	-	9/33/33/33	-
9	ETF	G	311	5	-	0/3/3/3	-
4	D10	K	305	-	-	0/7/7/7	-
9	ETF	K	311	5	-	0/3/3/3	-
6	PLC	C	302	-	-	22/45/45/45	-
6	PLC	B	308	-	-	15/45/45/45	-
8	HXG	C	303	-	-	9/33/33/33	-
4	D10	F	306	-	-	6/7/7/7	-
4	D10	B	305	-	-	6/7/7/7	-
4	D10	J	307	-	-	5/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	D10	J	305	-	-	2/7/7/7	-
4	D10	F	307	-	-	5/7/7/7	-
6	PLC	G	303	-	-	22/45/45/45	-
4	D10	F	305	-	-	2/7/7/7	-
4	D10	B	304	-	-	2/7/7/7	-
4	D10	J	306	-	-	6/7/7/7	-
6	PLC	G	306	-	-	14/45/45/45	-
4	D10	J	304	-	-	0/7/7/7	-
7	P1O	G	310	-	-	26/41/41/41	-
7	P1O	F	303	-	-	26/41/41/41	-
7	P1O	G	309	-	-	19/41/41/41	-
4	D10	F	304	-	-	0/7/7/7	-
4	D10	B	306	-	-	5/7/7/7	-
6	PLC	G	308	-	-	16/45/45/45	-
7	P1O	K	310	-	-	26/41/41/41	-
8	HXG	G	307	-	-	7/33/33/33	-
6	PLC	K	308	-	-	16/45/45/45	-
6	PLC	C	307	-	-	16/45/45/45	-
6	PLC	F	302	-	-	15/45/45/45	-
9	ETF	C	310	5	-	0/3/3/3	-

The worst 5 of 36 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	F	308	P1O	O5-C9	4.34	1.46	1.33
7	J	301	P1O	O5-C9	4.32	1.45	1.33
7	B	307	P1O	O5-C9	4.31	1.45	1.33
7	F	303	P1O	O5-C9	4.31	1.45	1.33
7	J	303	P1O	O5-C9	4.29	1.45	1.33

The worst 5 of 76 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	307	P1O	O7-C19-C20	4.24	120.66	111.48
7	F	308	P1O	O7-C19-C20	4.22	120.61	111.48
7	J	301	P1O	O7-C19-C20	4.21	120.59	111.48
7	K	309	P1O	O7-C19-C20	4.18	120.52	111.48
7	C	308	P1O	O7-C19-C20	4.16	120.48	111.48

There are no chirality outliers.

5 of 681 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	301	PLC	C1'-C'-O2-C2
6	B	301	PLC	C1-O3P-P-O2P
6	B	301	PLC	C1-O3P-P-O4P
6	B	308	PLC	C1'-C'-O2-C2
6	B	308	PLC	C1-O3P-P-O1P

There are no ring outliers.

48 monomers are involved in 333 short contacts:

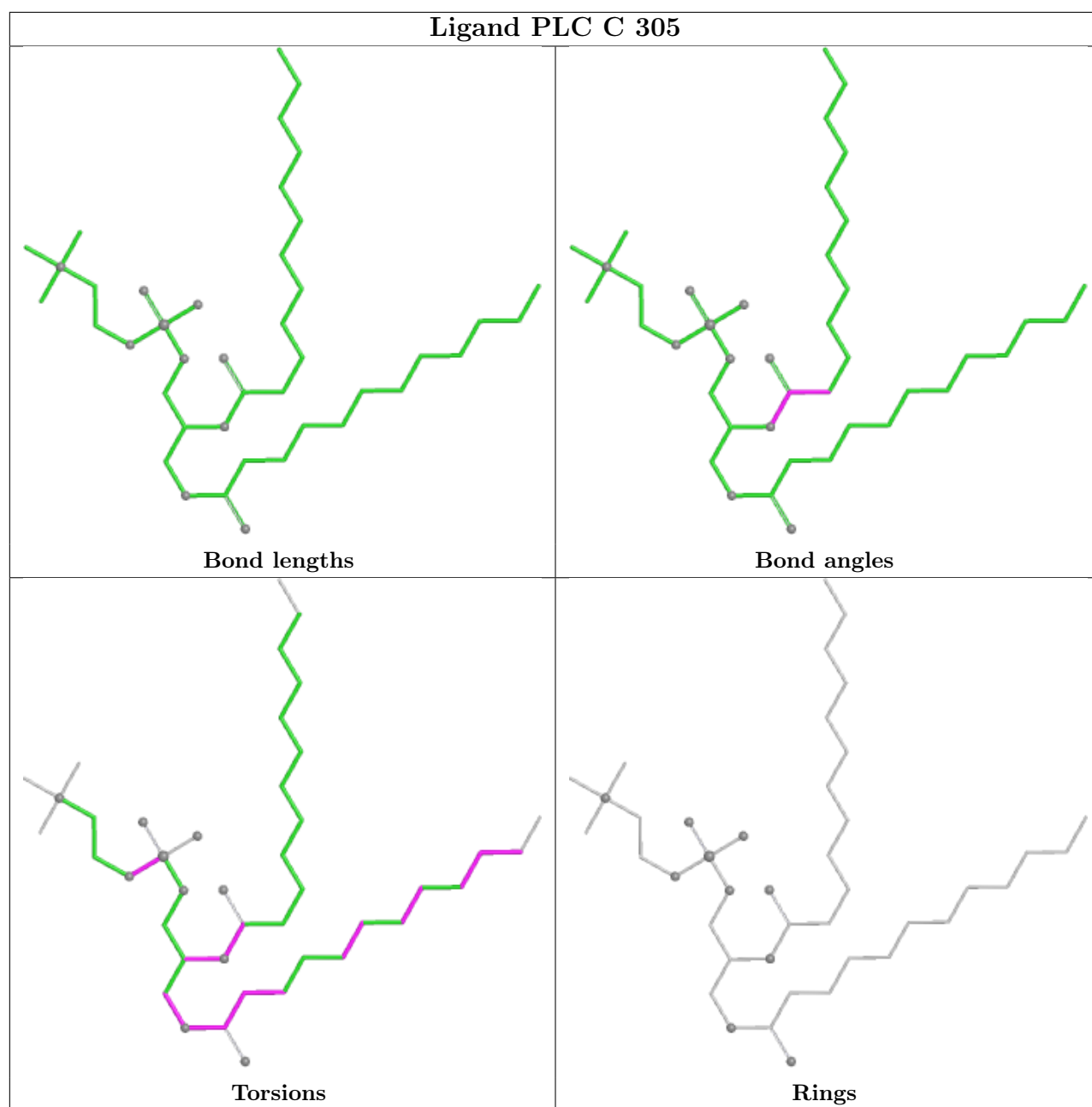
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	305	PLC	2	0
6	K	306	PLC	2	0
8	K	304	HXG	8	0
6	G	301	PLC	6	0
7	B	302	P1O	7	0
8	K	307	HXG	16	0
6	K	301	PLC	7	0
7	J	303	P1O	8	0
4	B	303	D10	1	0
8	C	306	HXG	15	0
7	C	309	P1O	14	0
6	C	311	PLC	7	0
7	F	308	P1O	29	0
7	B	307	P1O	28	0
7	C	308	P1O	7	0
6	B	301	PLC	3	0
6	J	302	PLC	3	0
7	J	301	P1O	26	0
7	K	309	P1O	7	0
4	C	304	D10	3	0
4	G	305	D10	3	0
6	K	303	PLC	7	0
8	G	304	HXG	8	0
9	G	311	ETF	1	0
4	K	305	D10	3	0
9	K	311	ETF	1	0
6	C	302	PLC	5	0
8	C	303	HXG	8	0
4	F	306	D10	6	0
4	B	305	D10	6	0

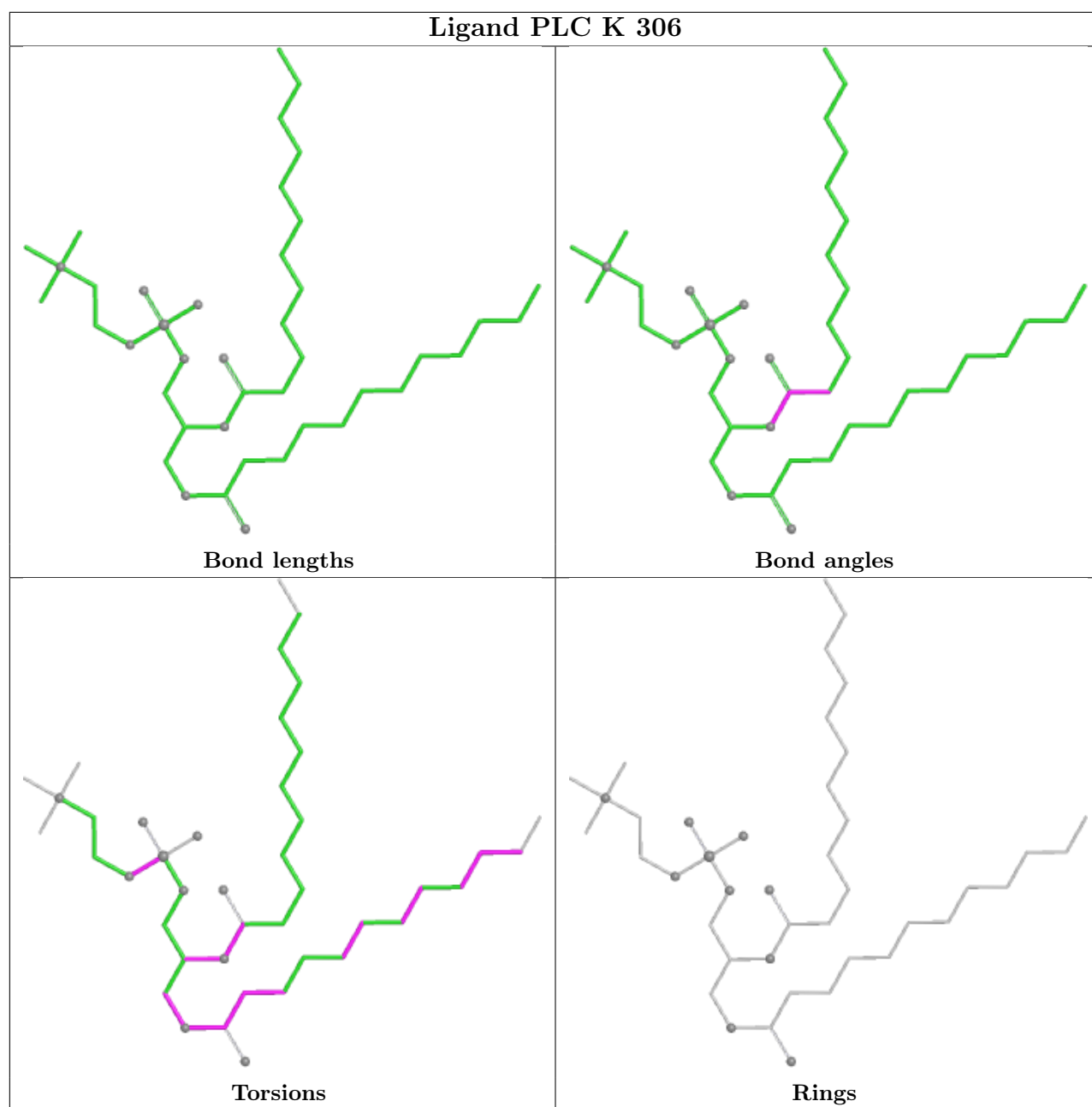
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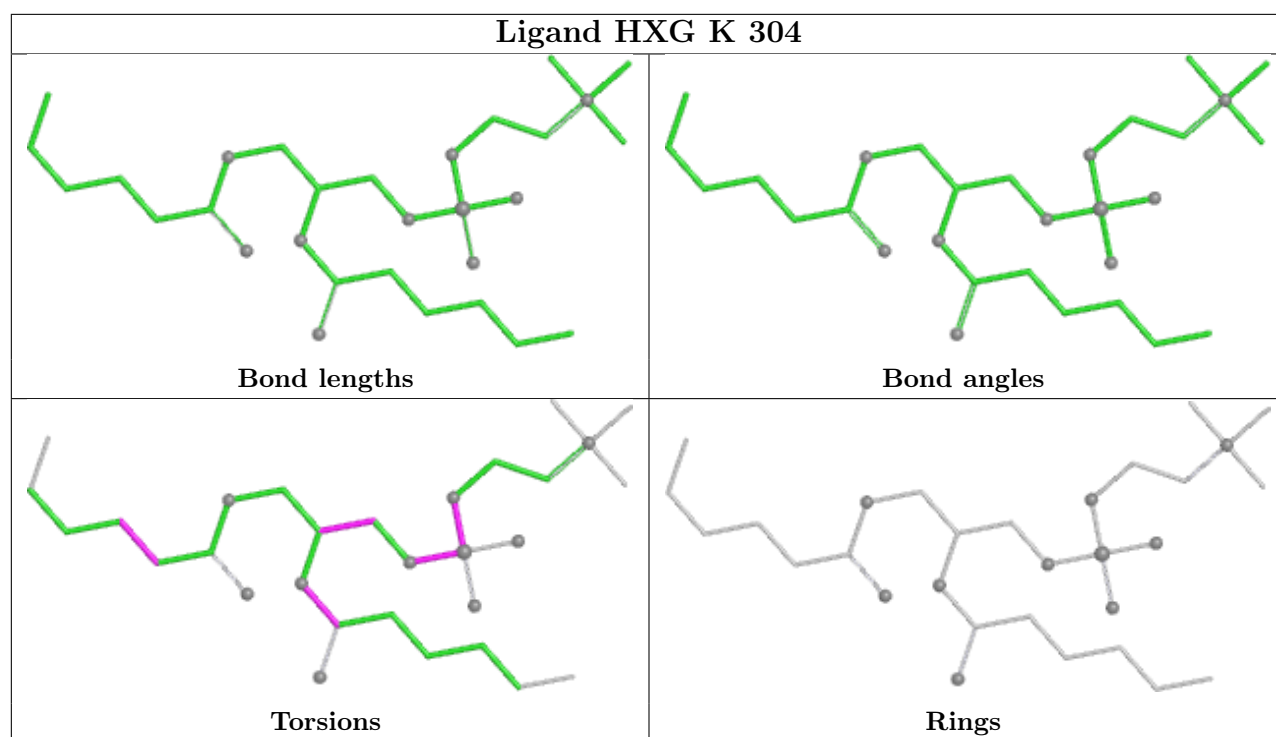
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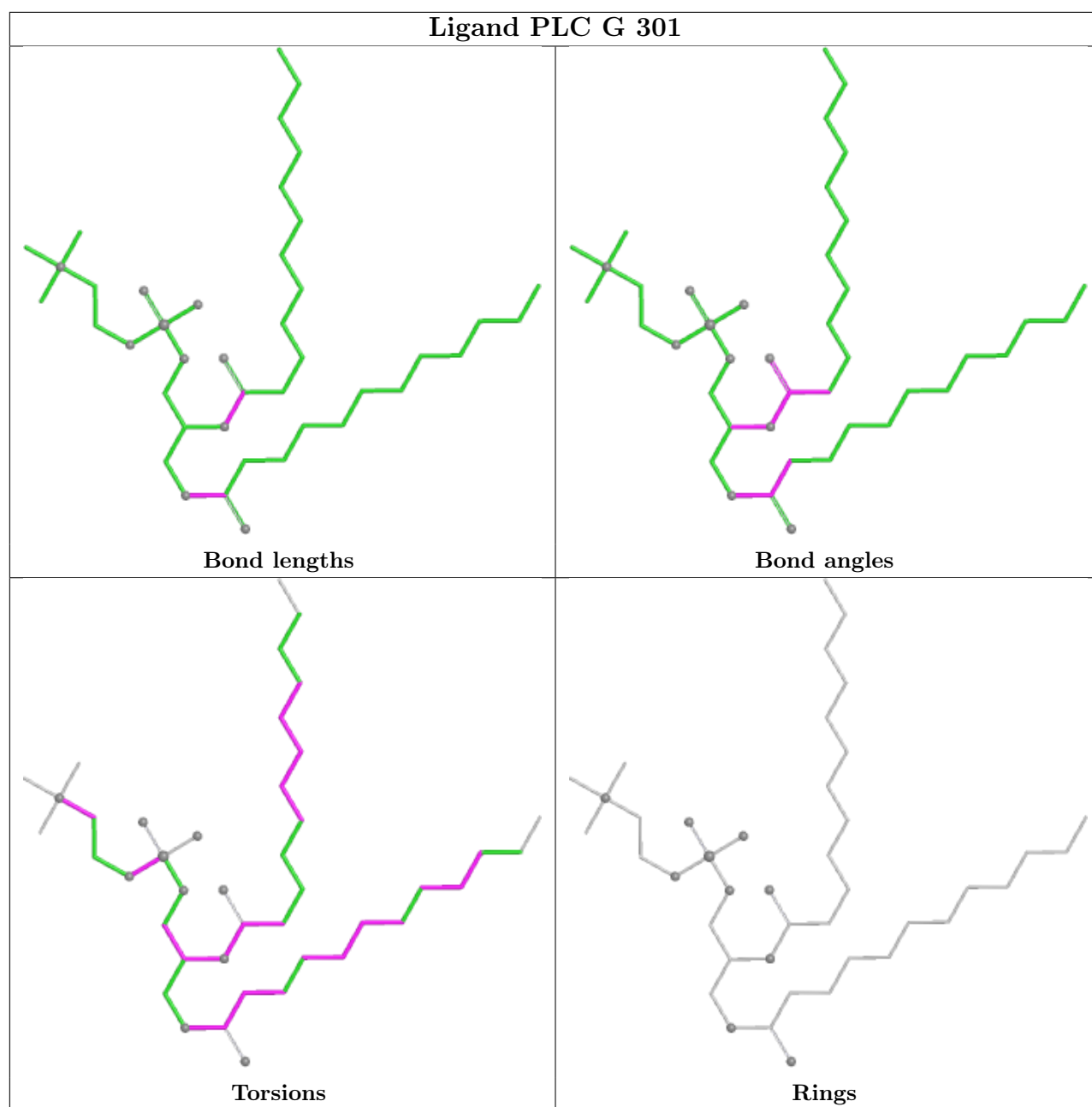
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	307	D10	1	0
4	J	305	D10	1	0
6	G	303	PLC	6	0
4	F	305	D10	1	0
4	B	304	D10	1	0
4	J	306	D10	6	0
6	G	306	PLC	2	0
7	G	310	P1O	16	0
7	F	303	P1O	8	0
7	G	309	P1O	7	0
4	B	306	D10	1	0
6	G	308	PLC	5	0
7	K	310	P1O	15	0
8	G	307	HXG	15	0
6	K	308	PLC	5	0
6	C	307	PLC	5	0
6	F	302	PLC	3	0
9	C	310	ETF	4	0

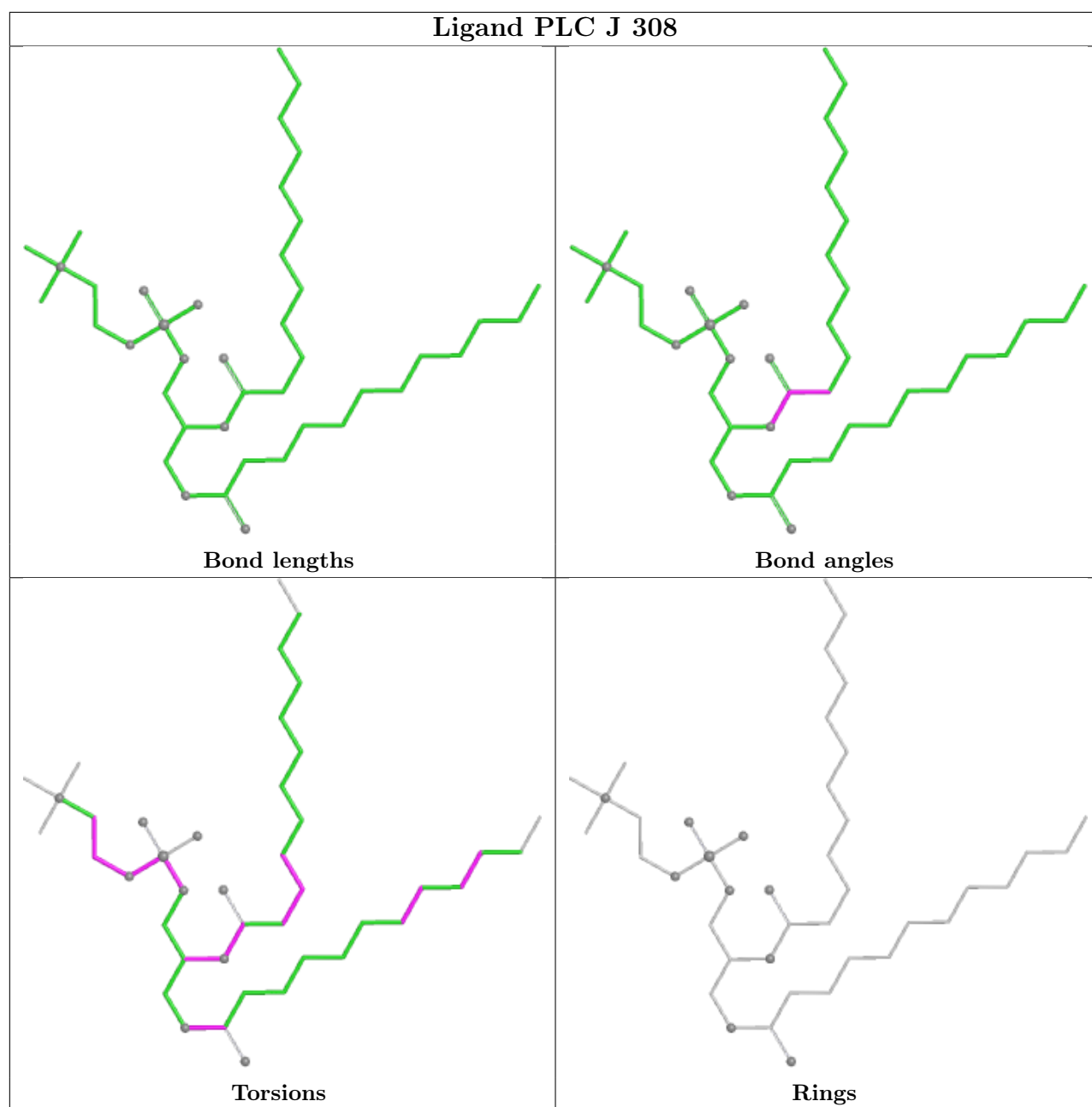
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

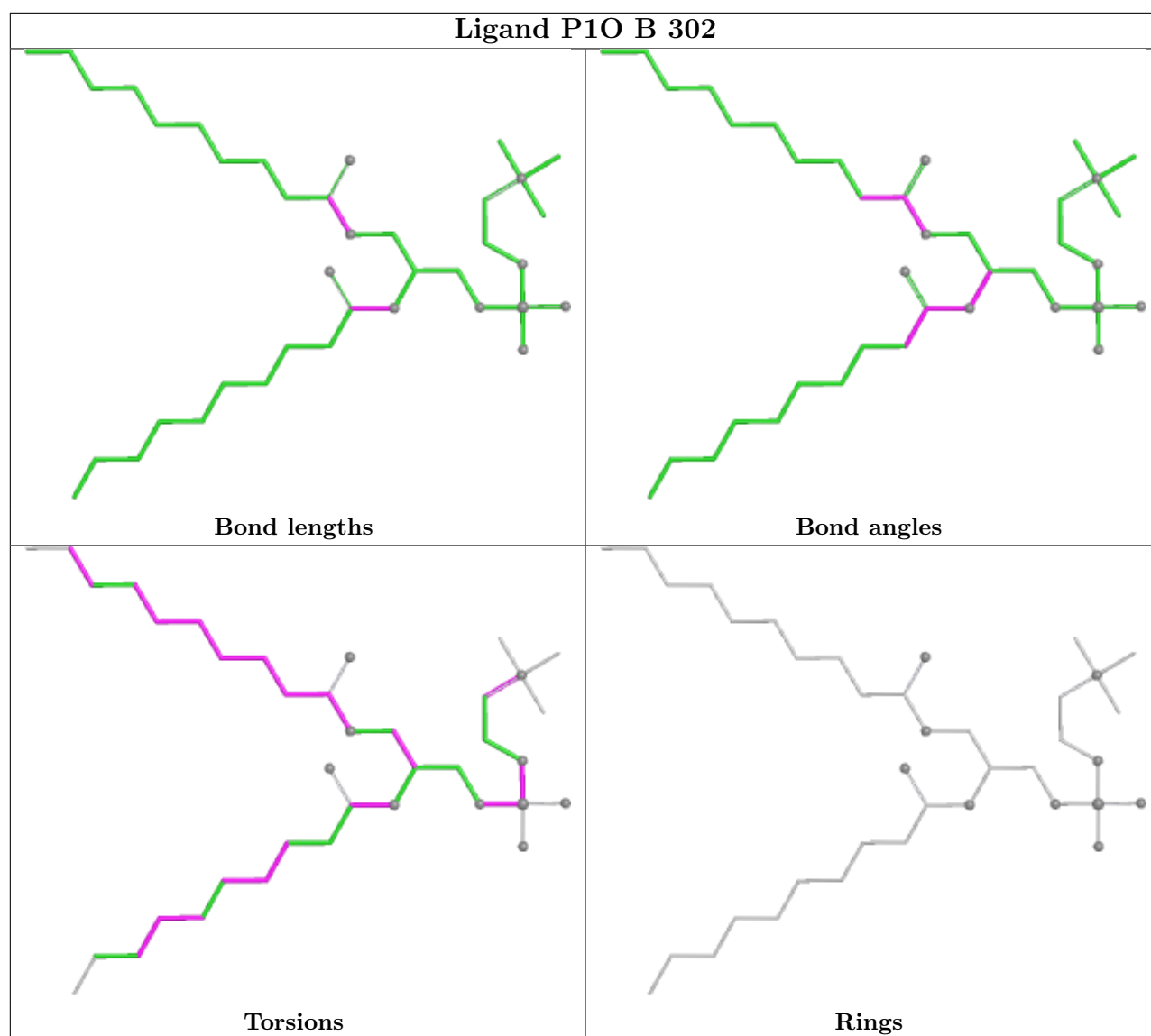


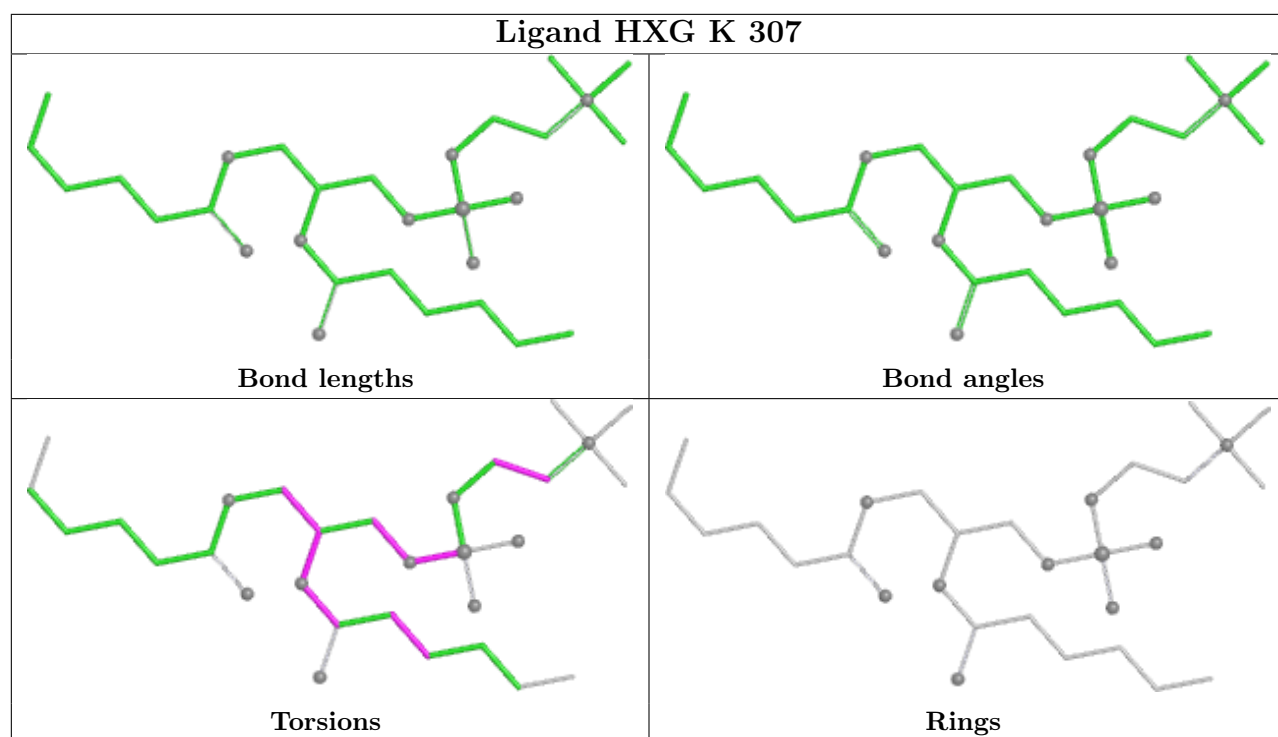


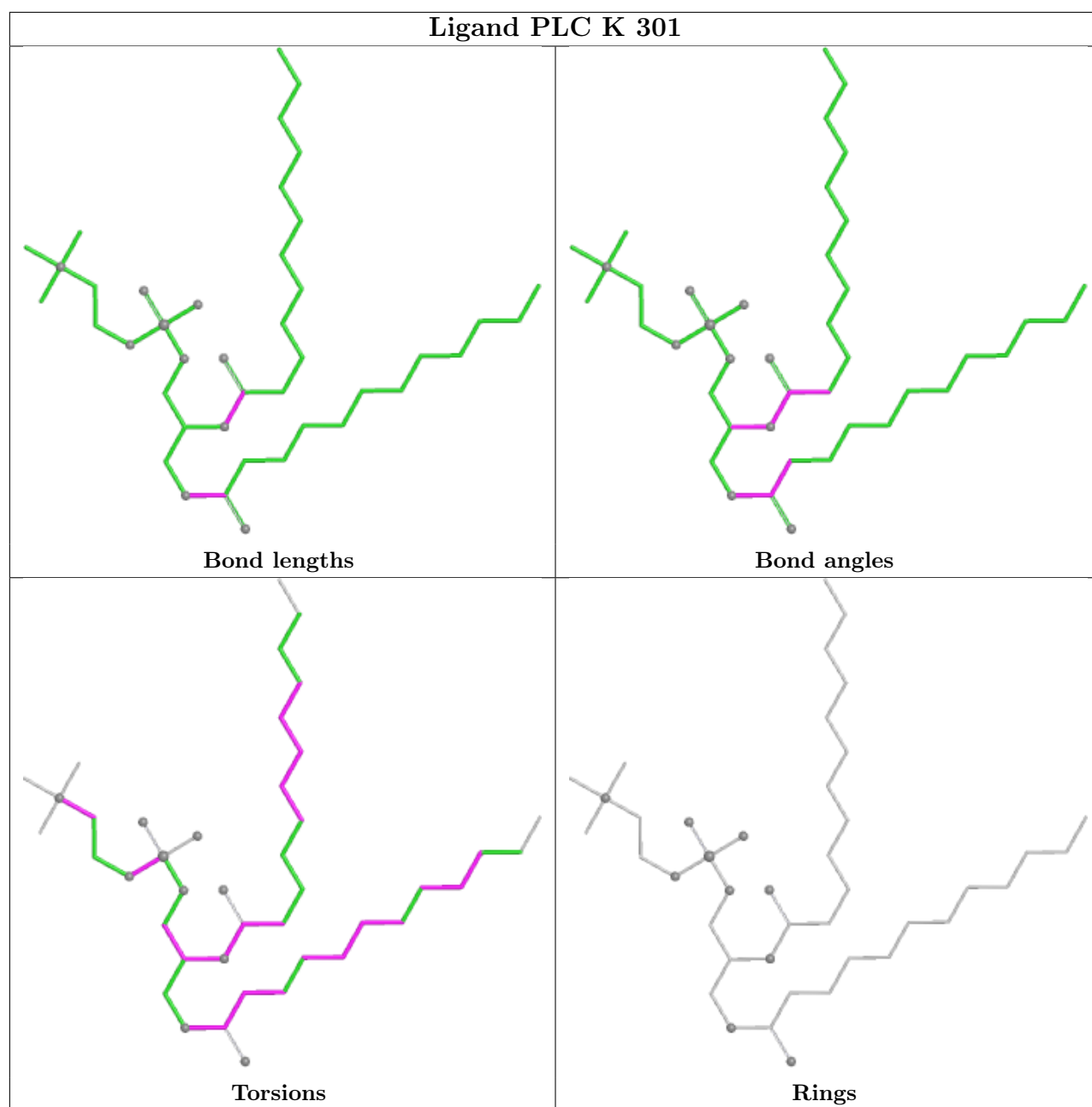


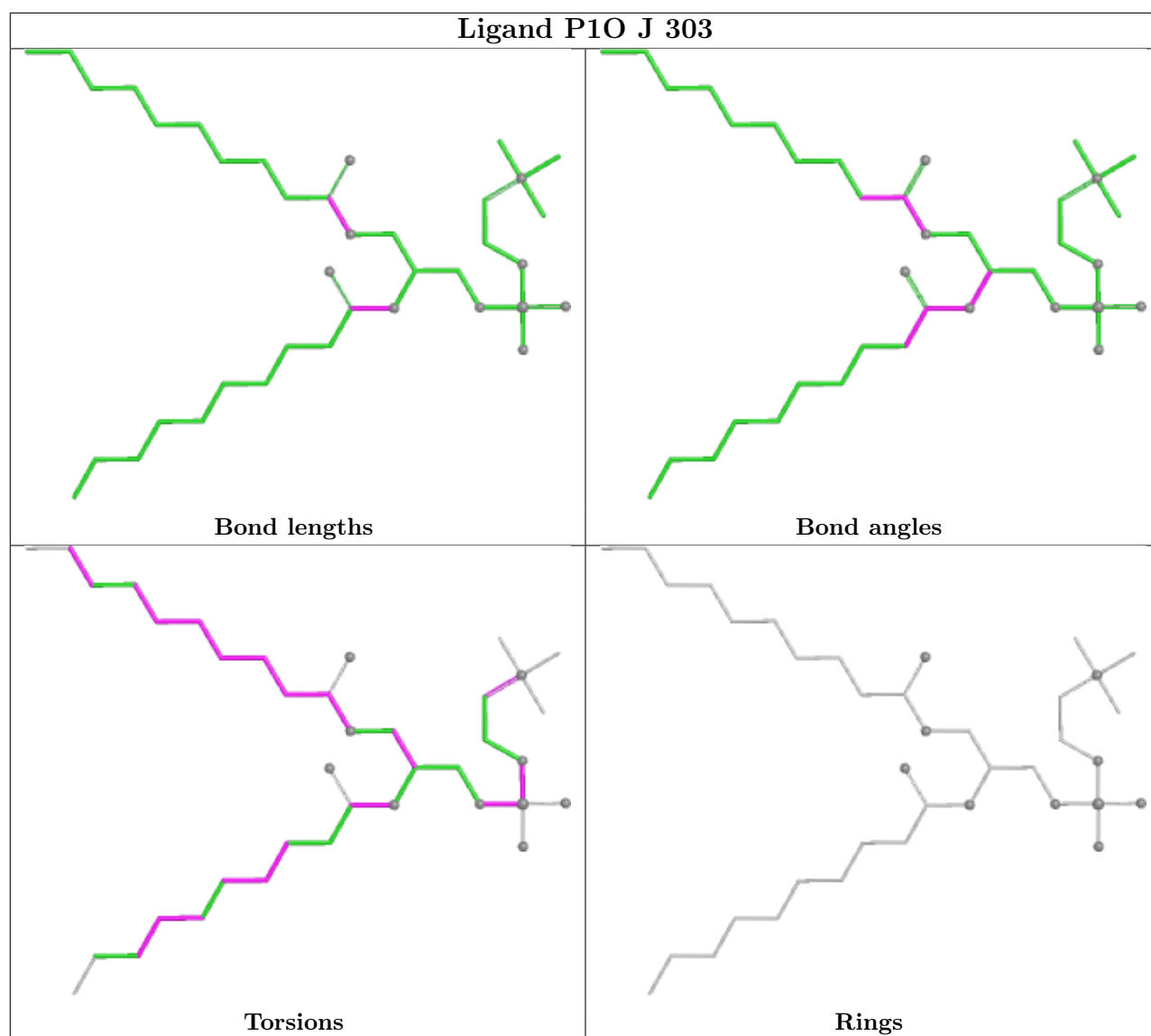


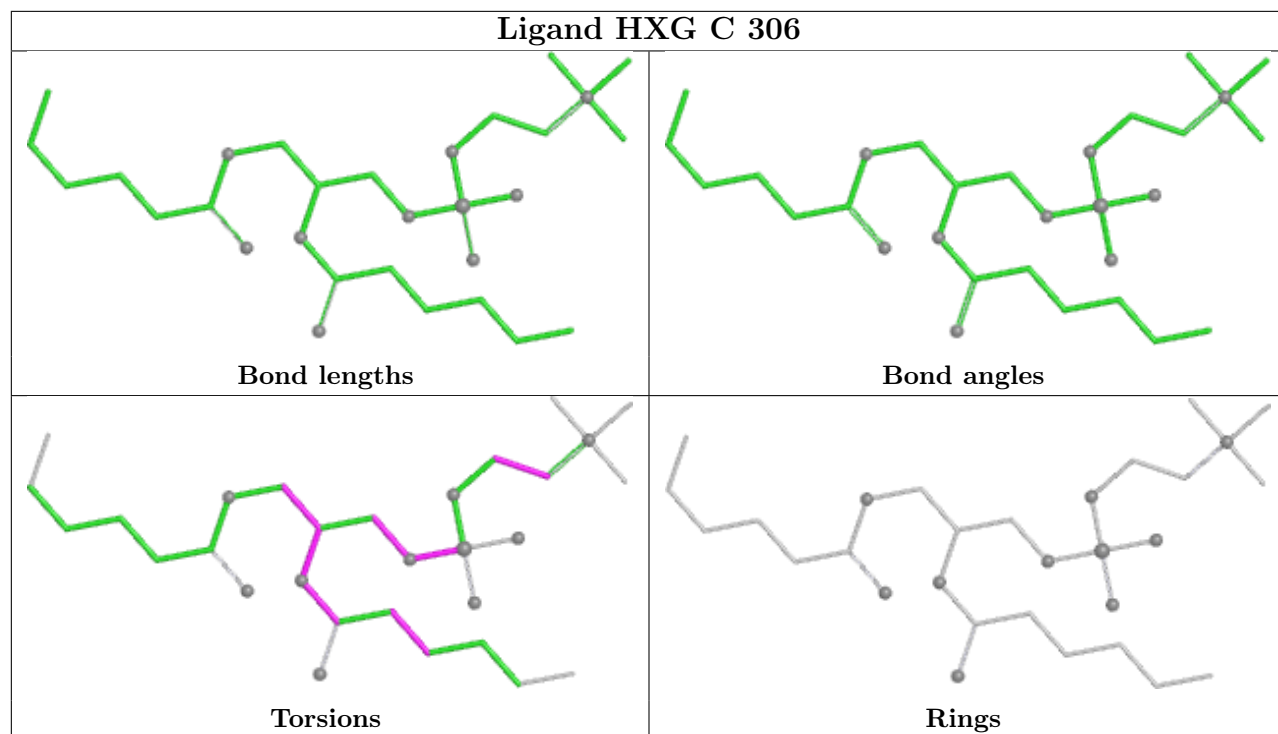


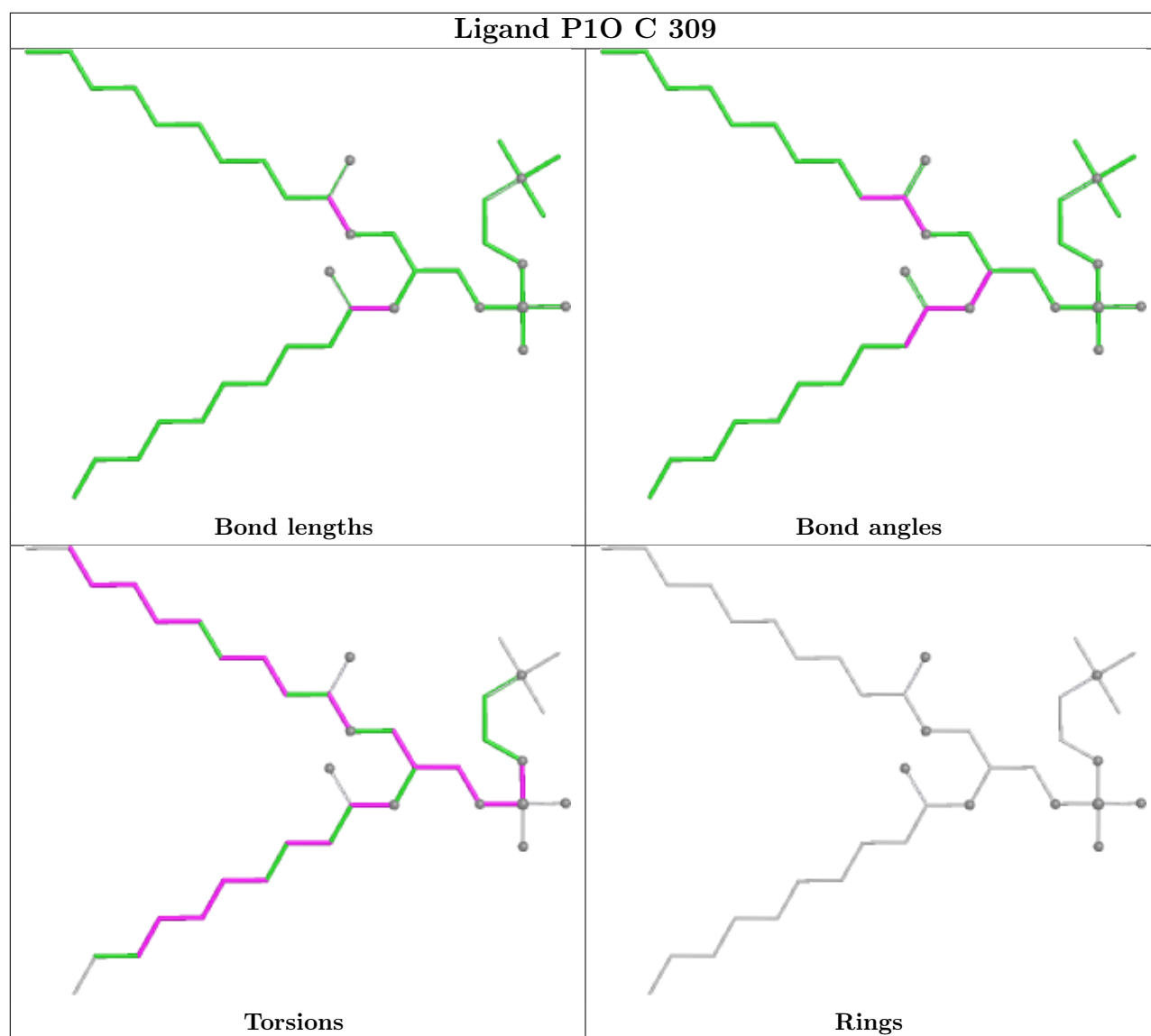


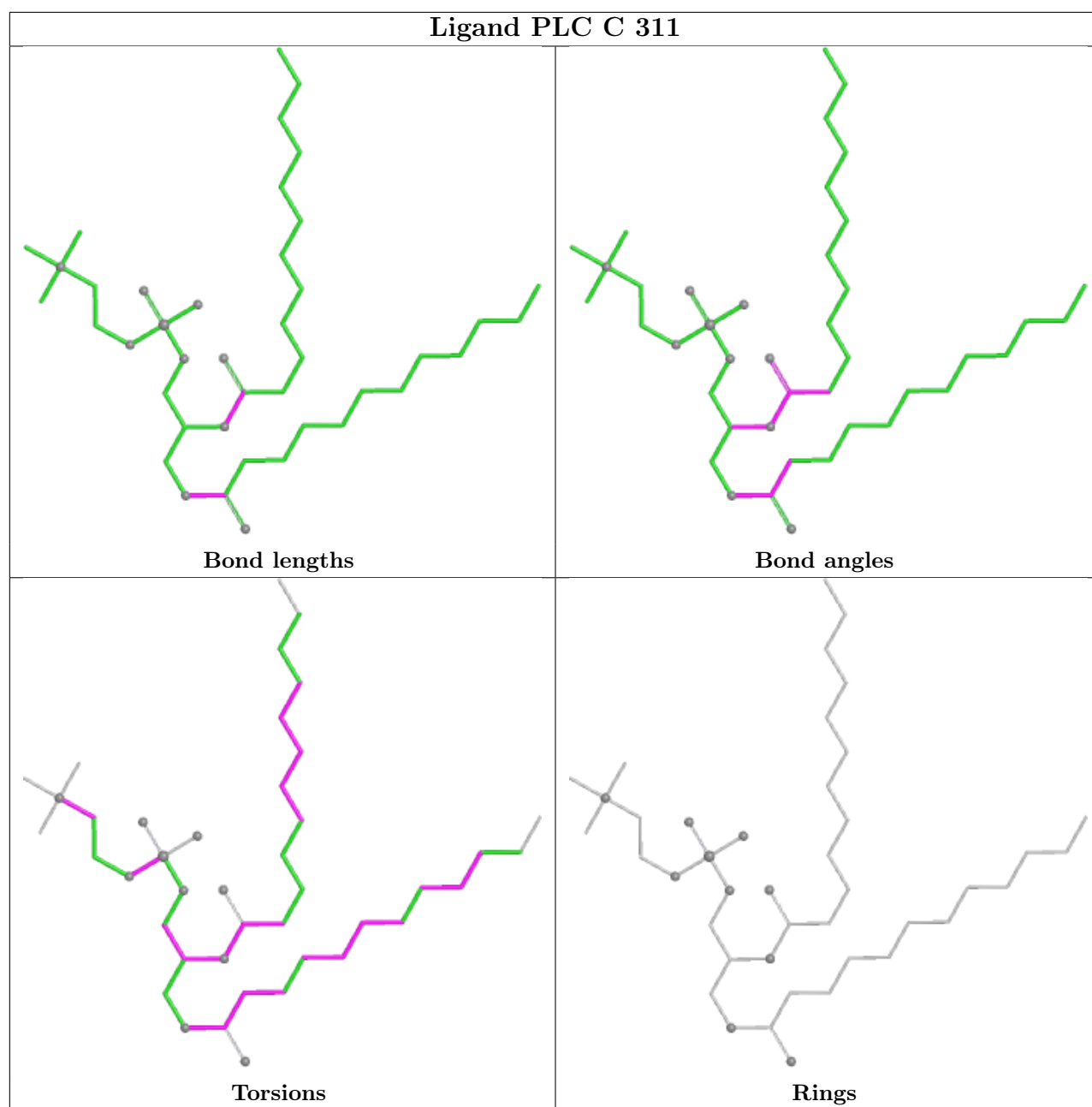


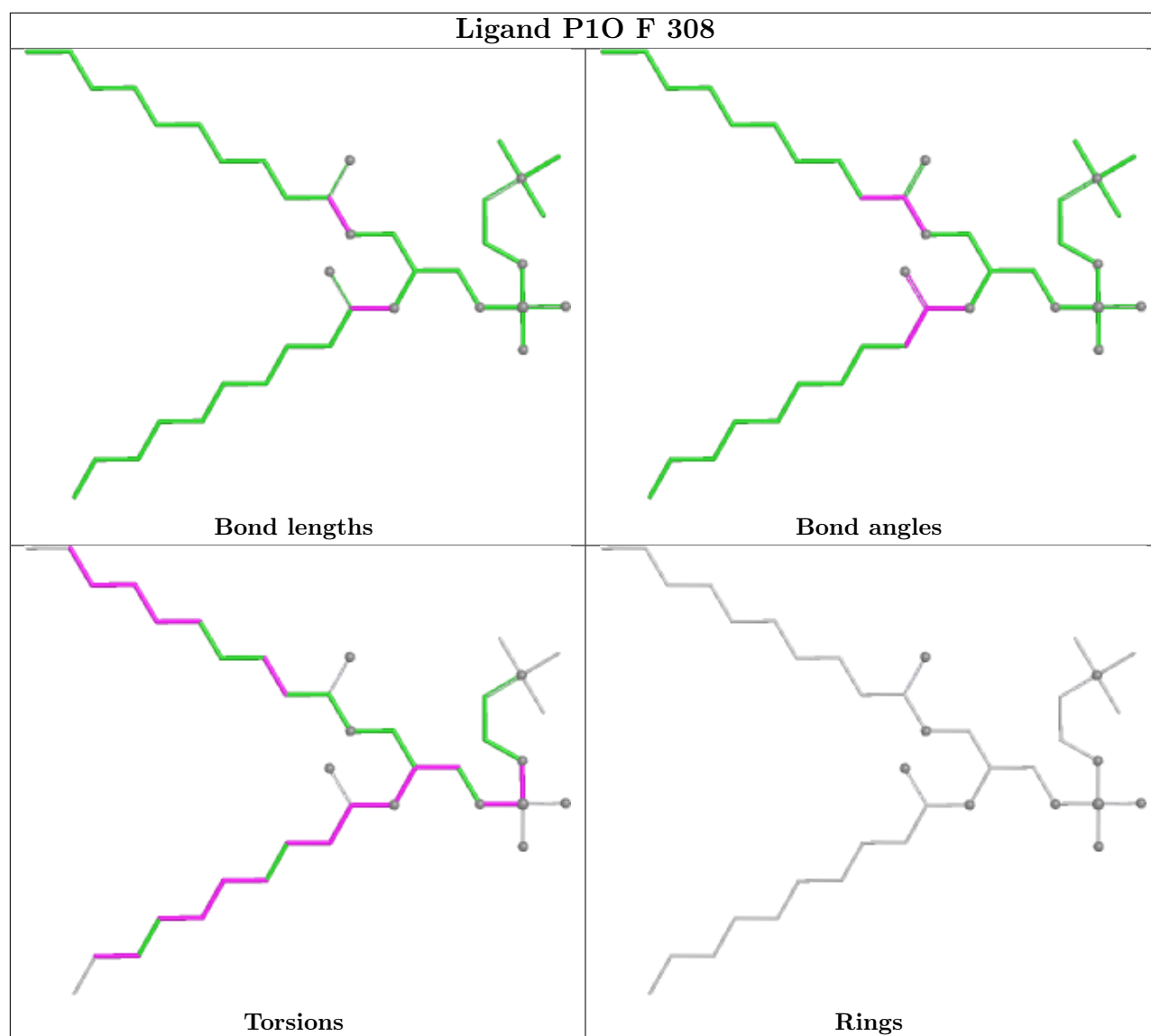


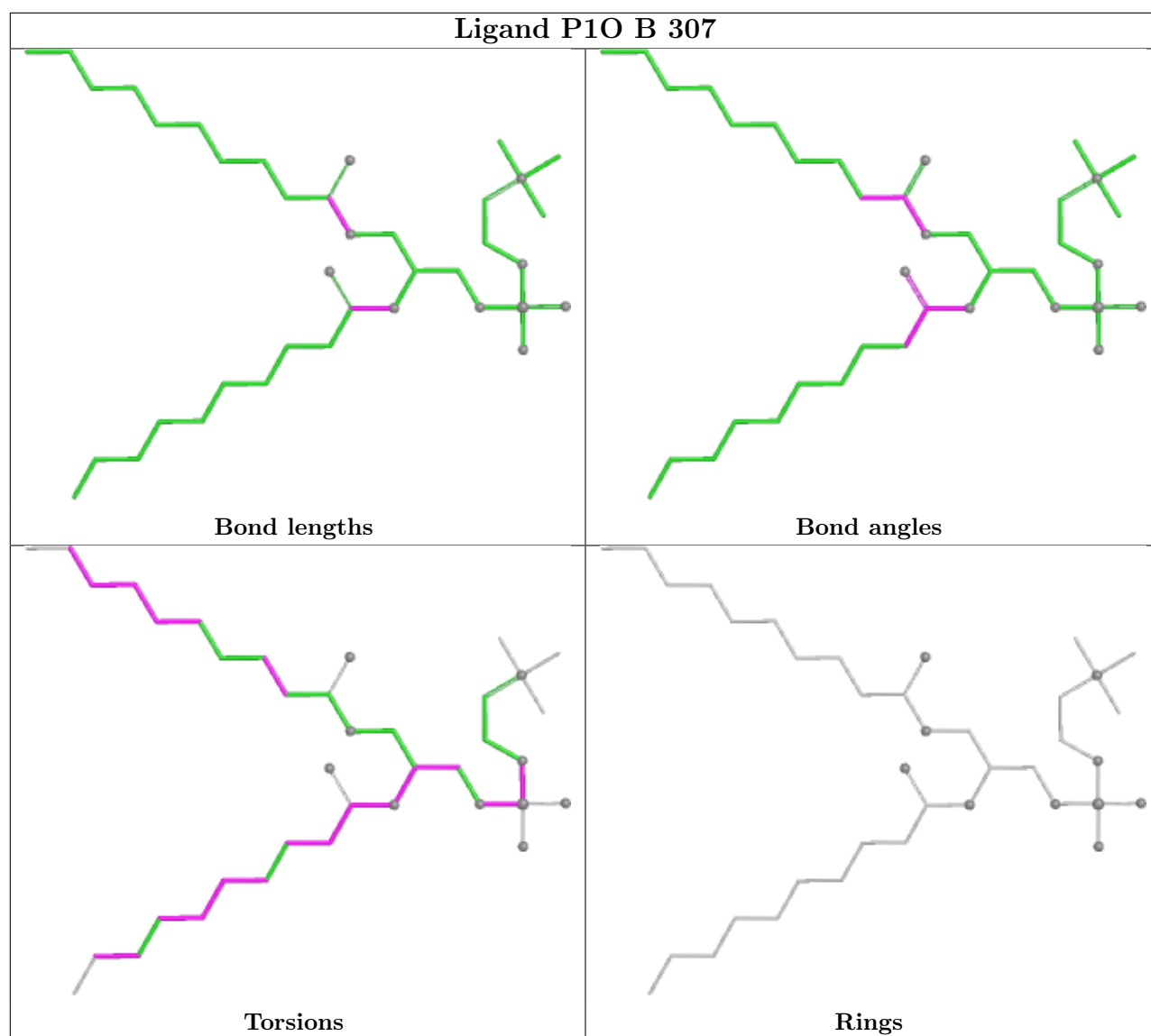


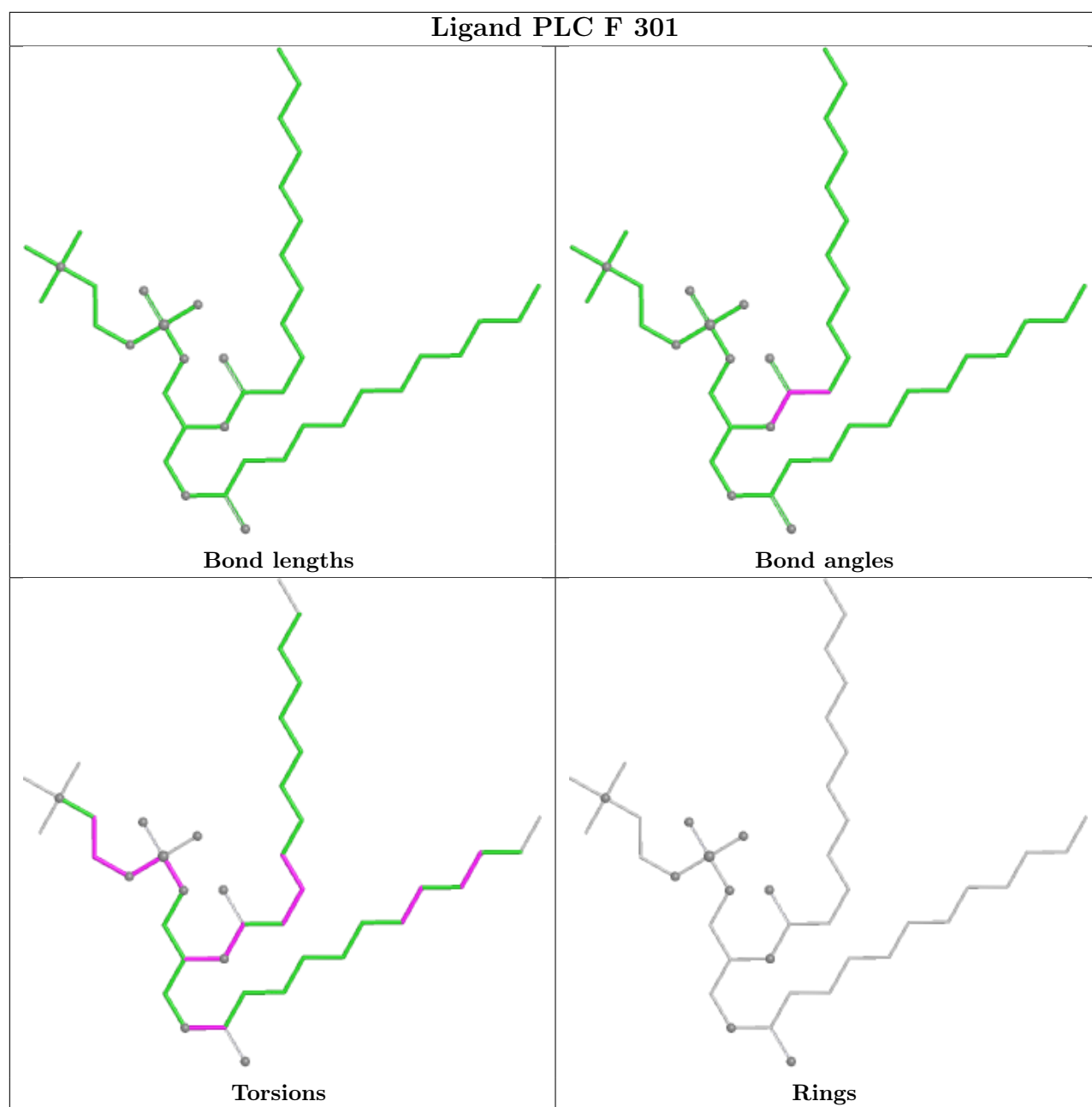


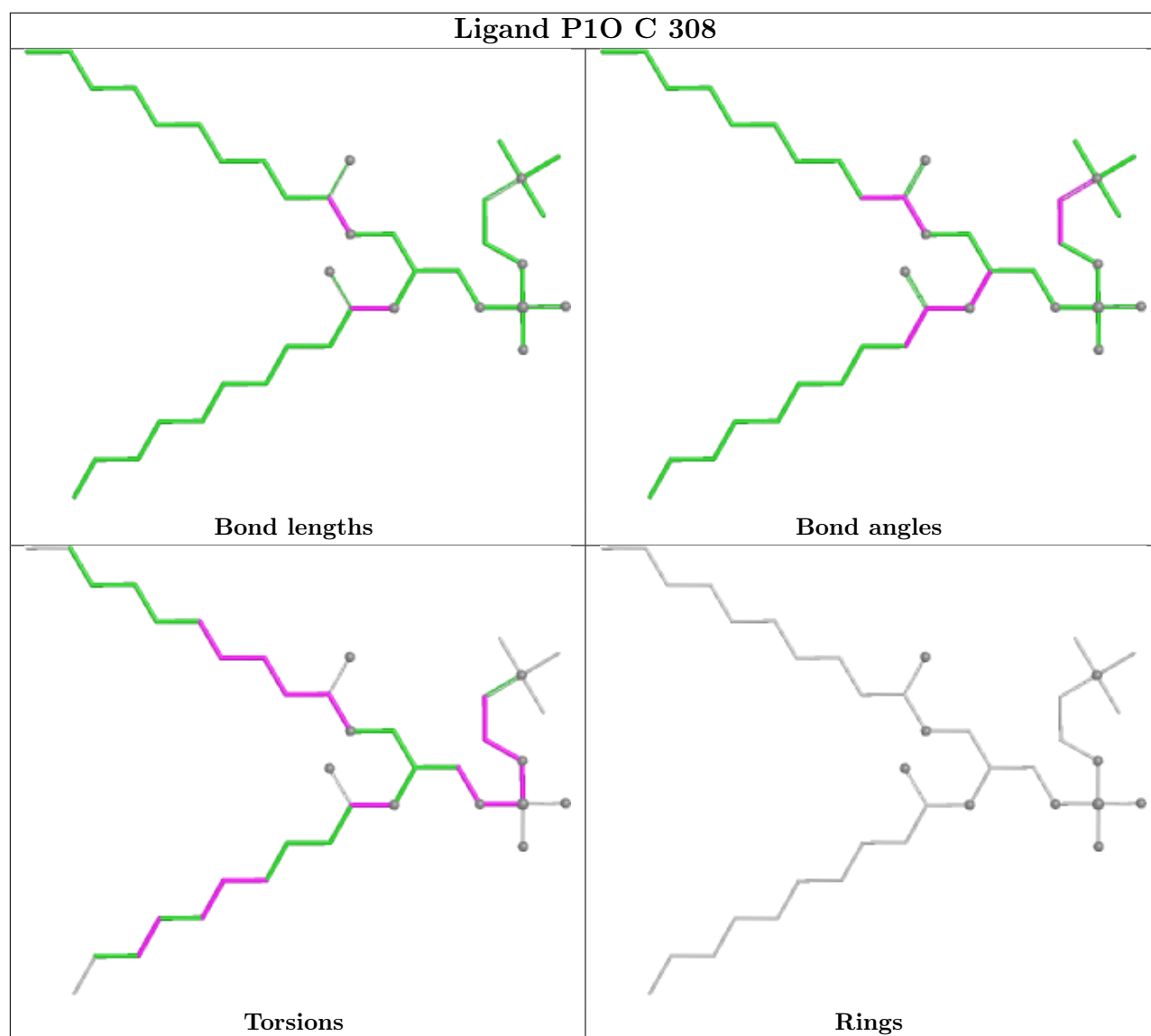


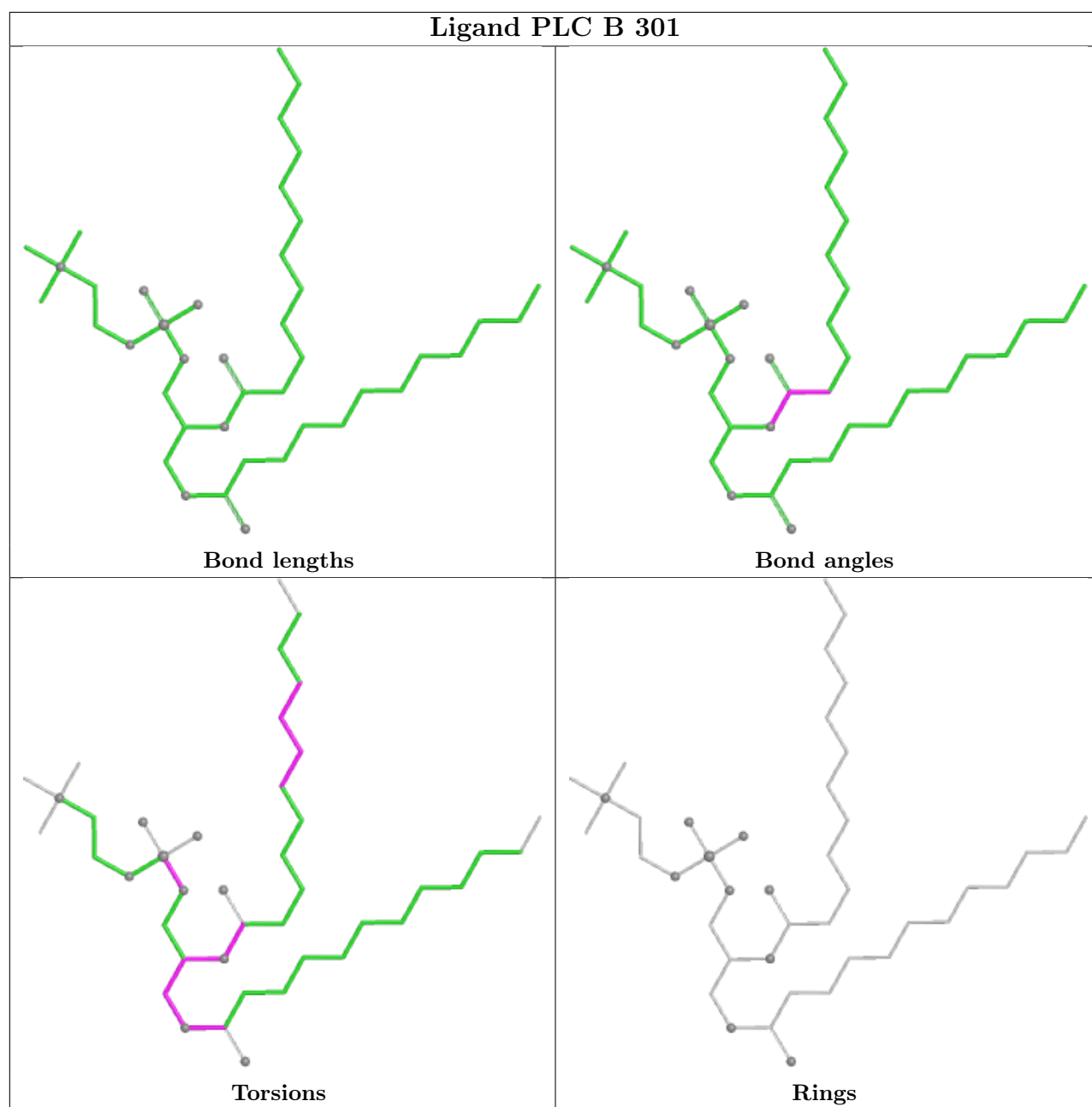


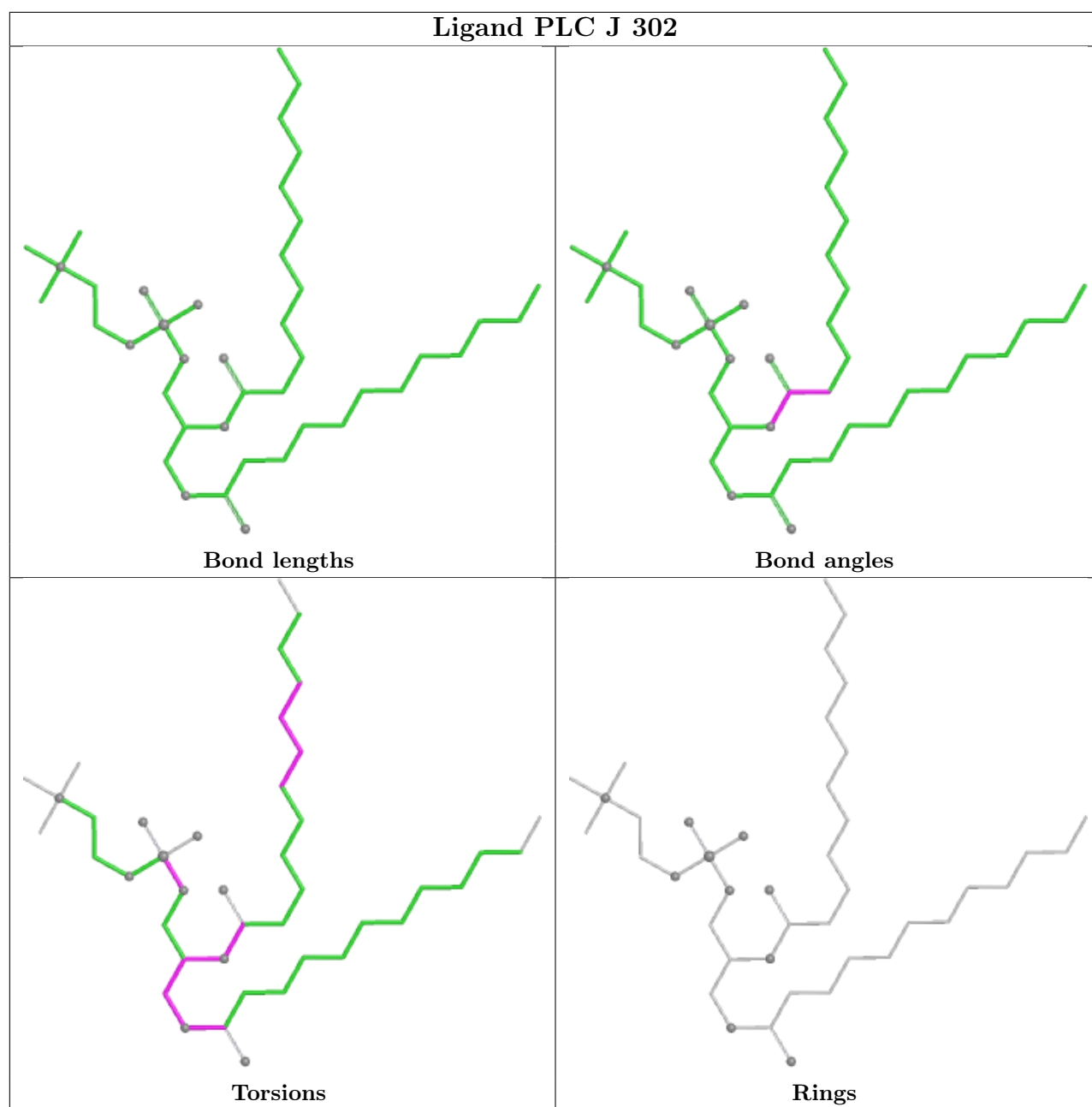


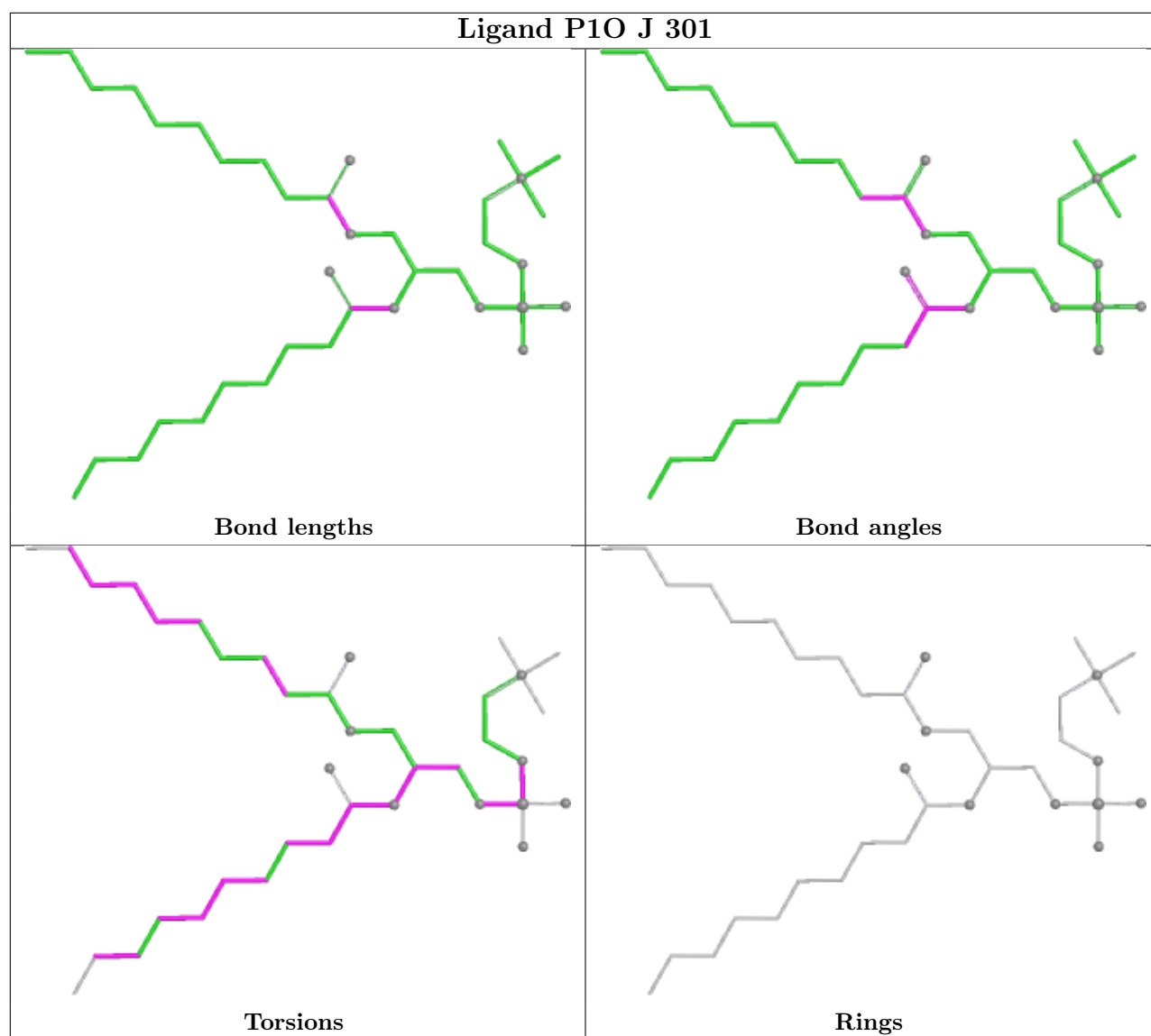


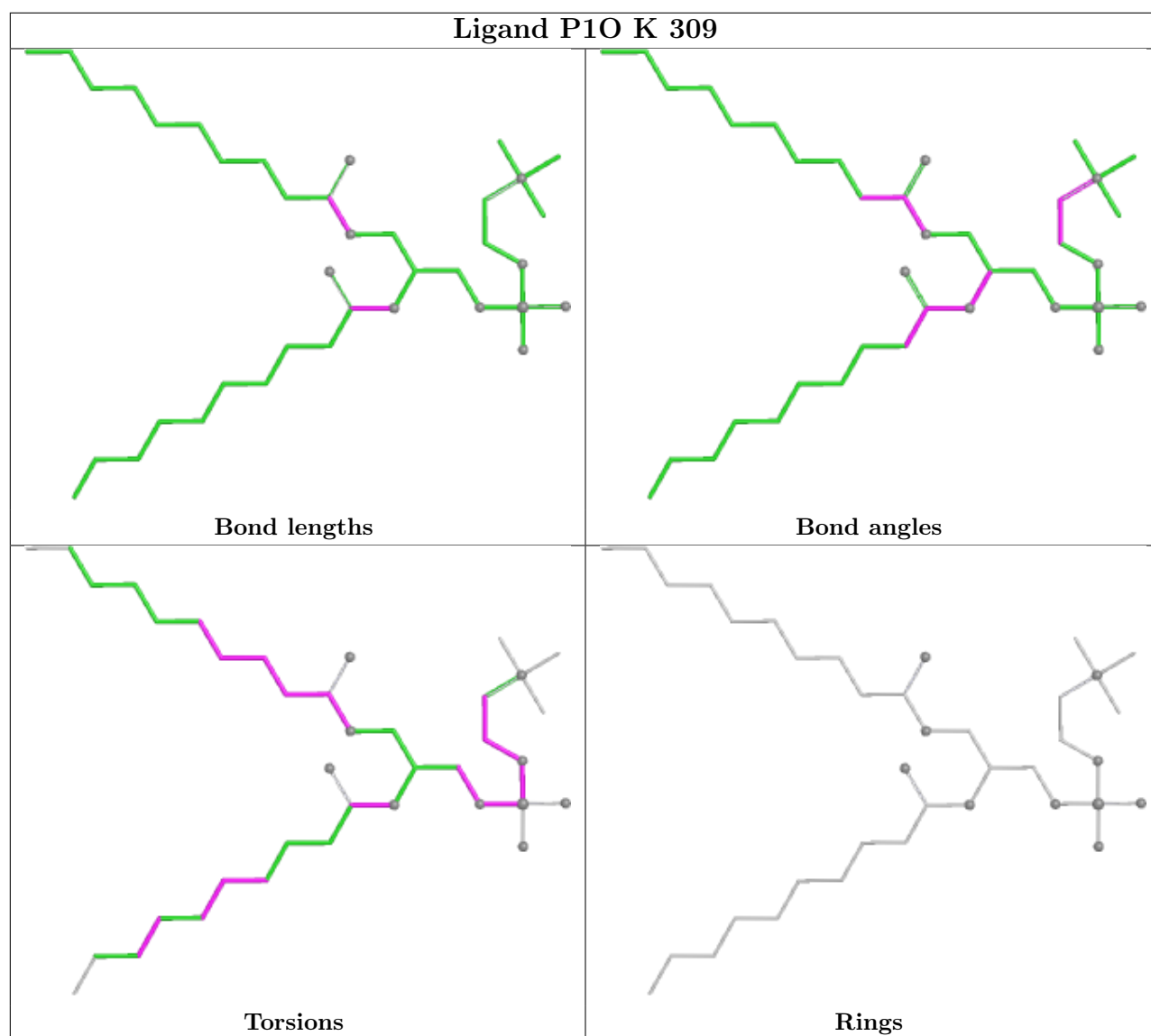


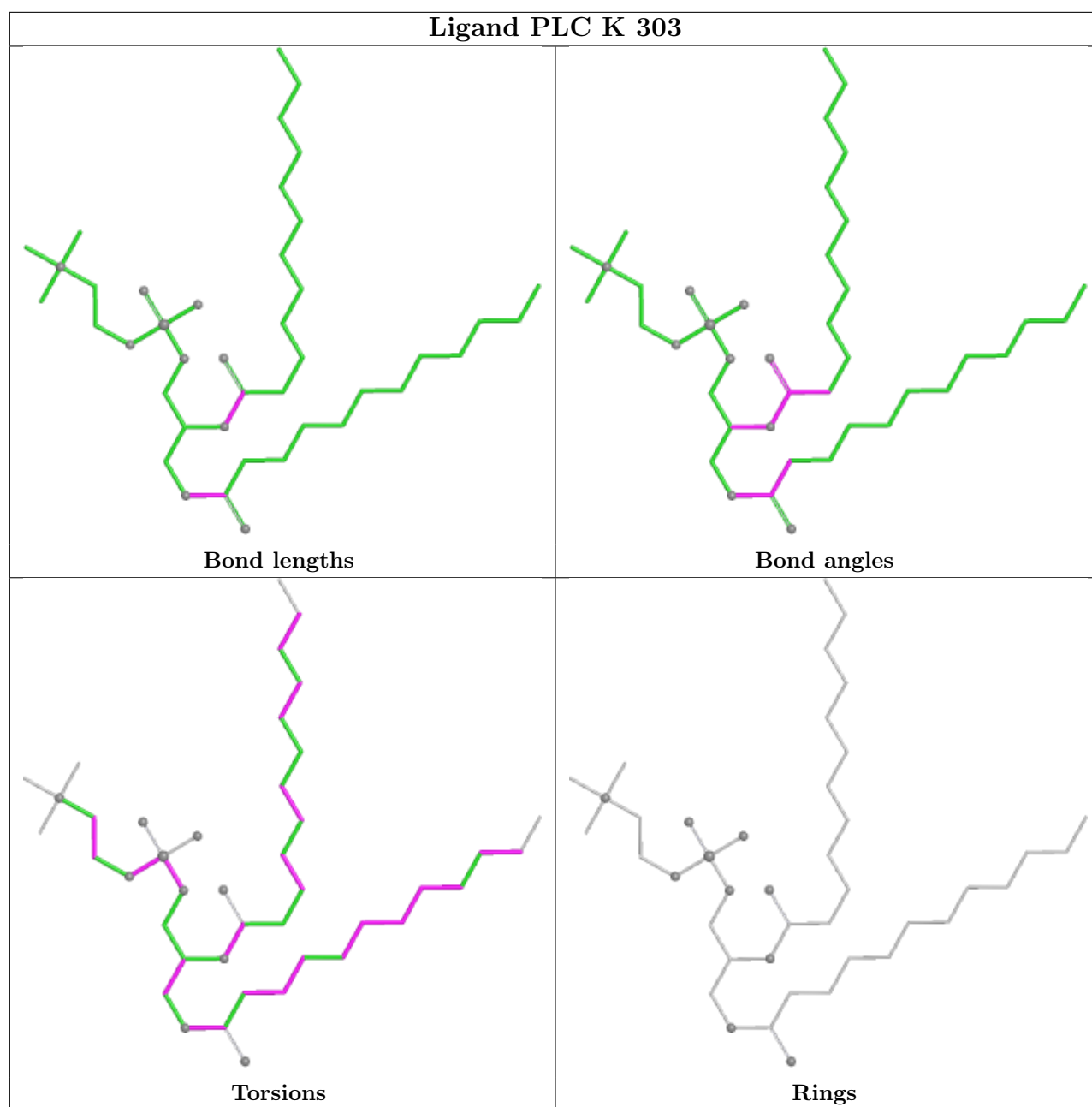


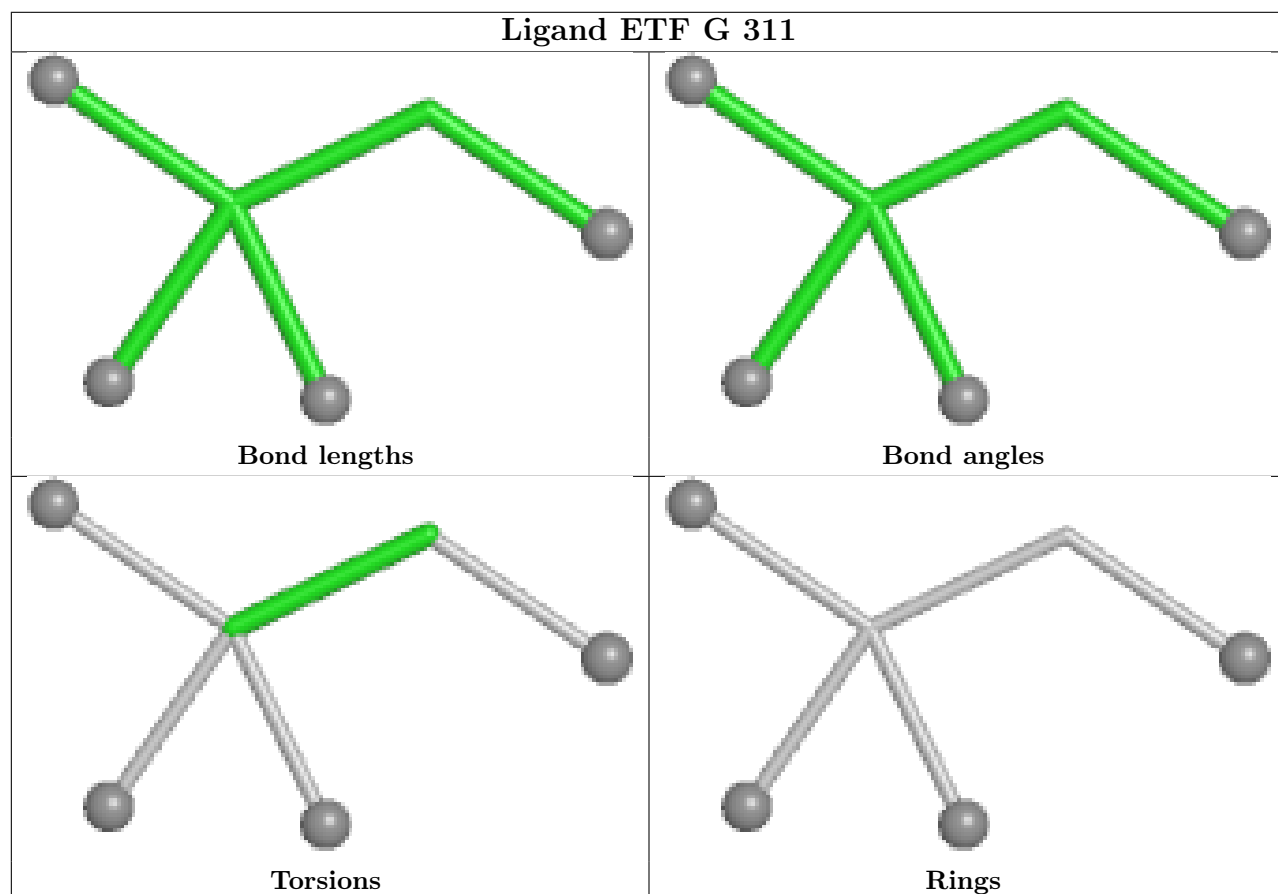
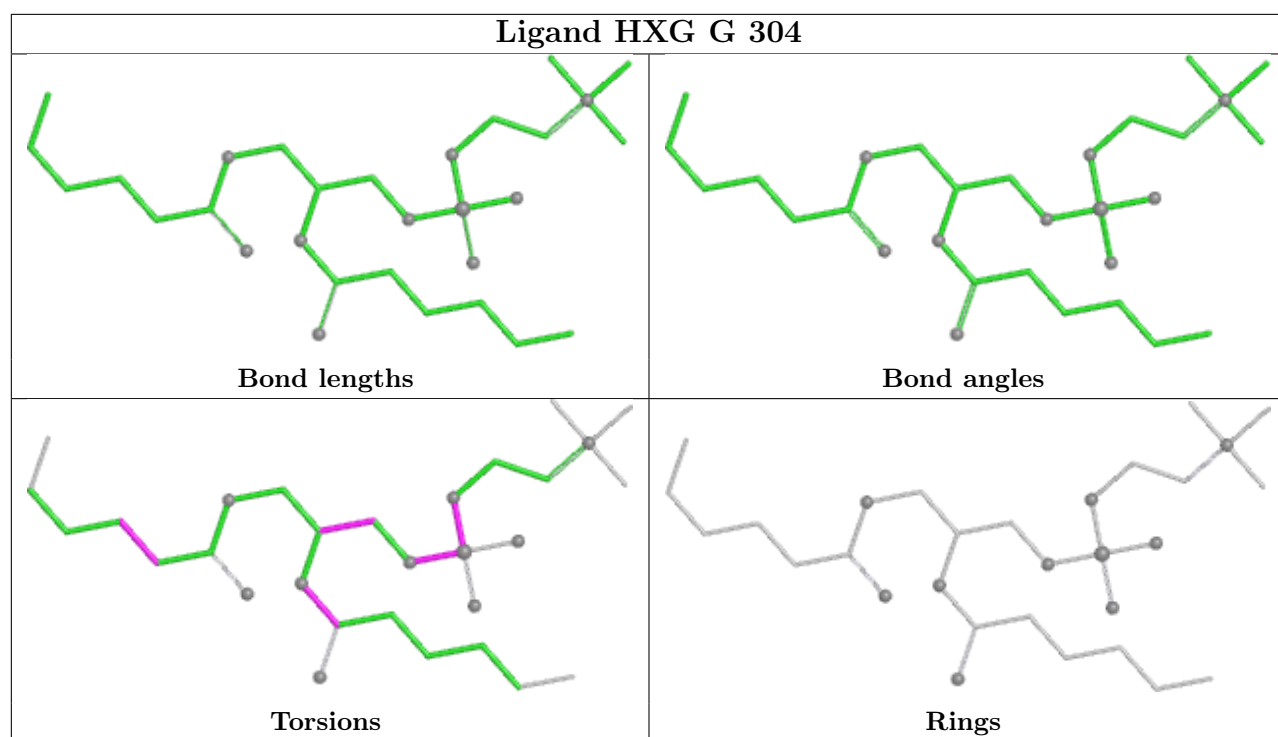


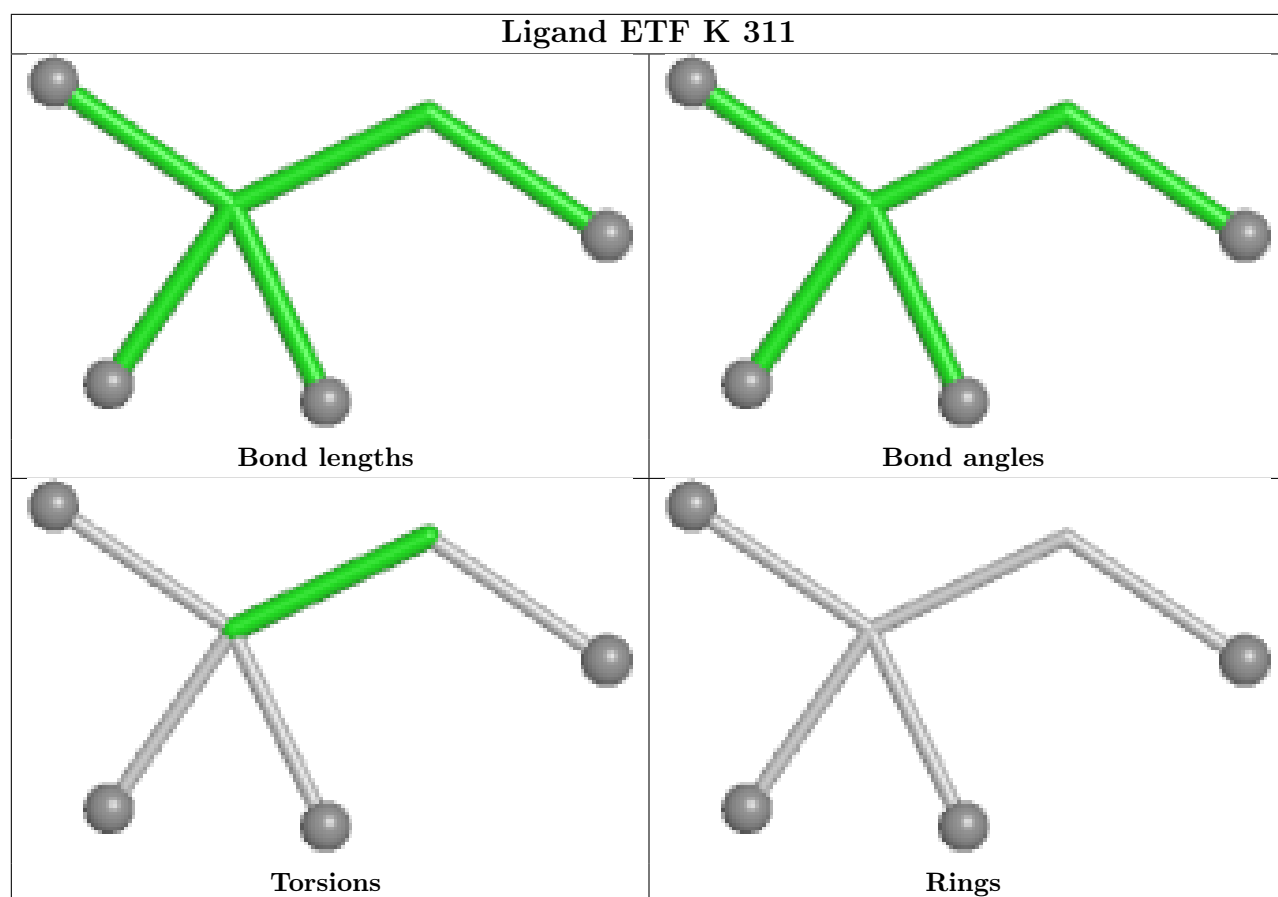


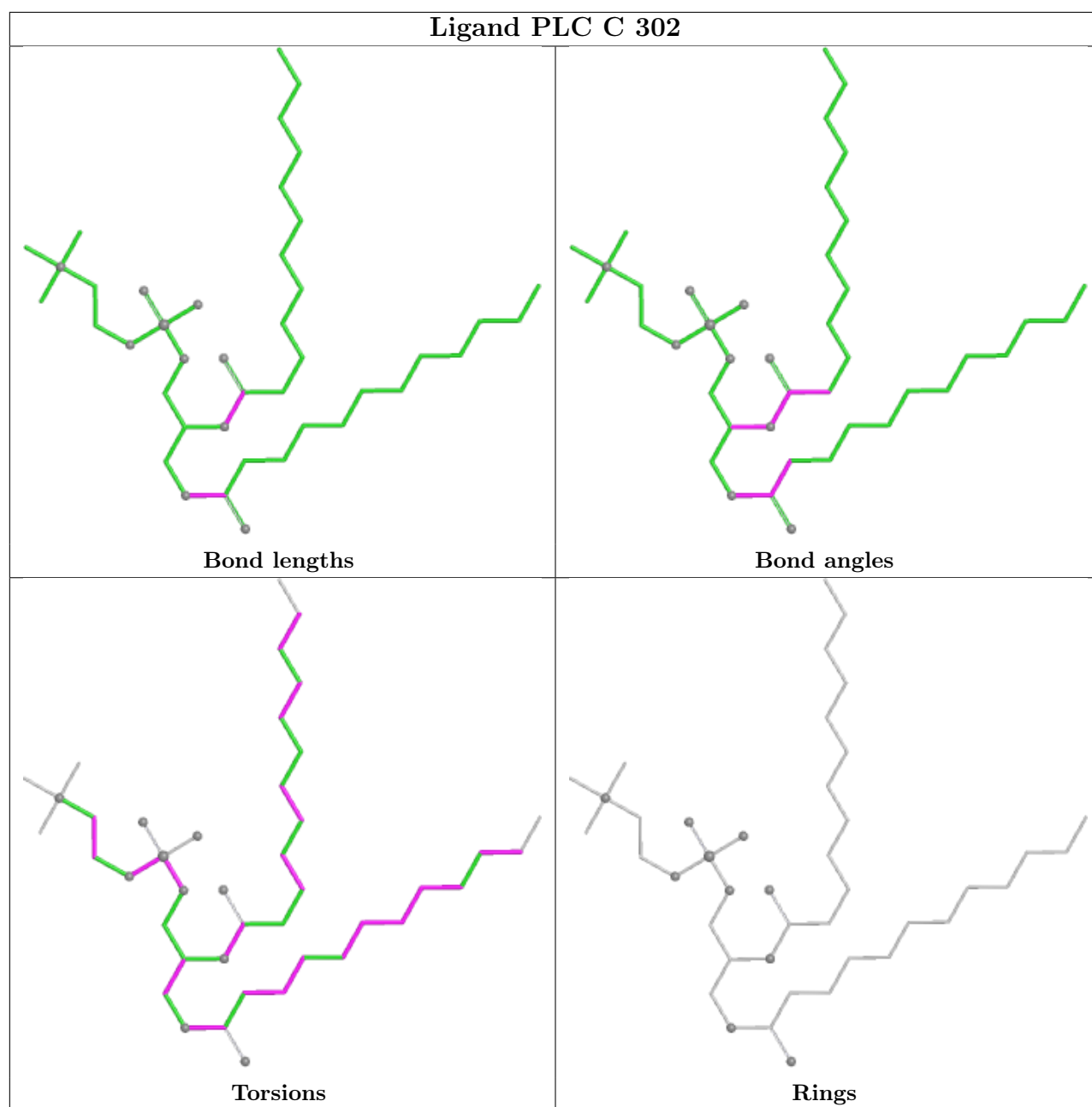


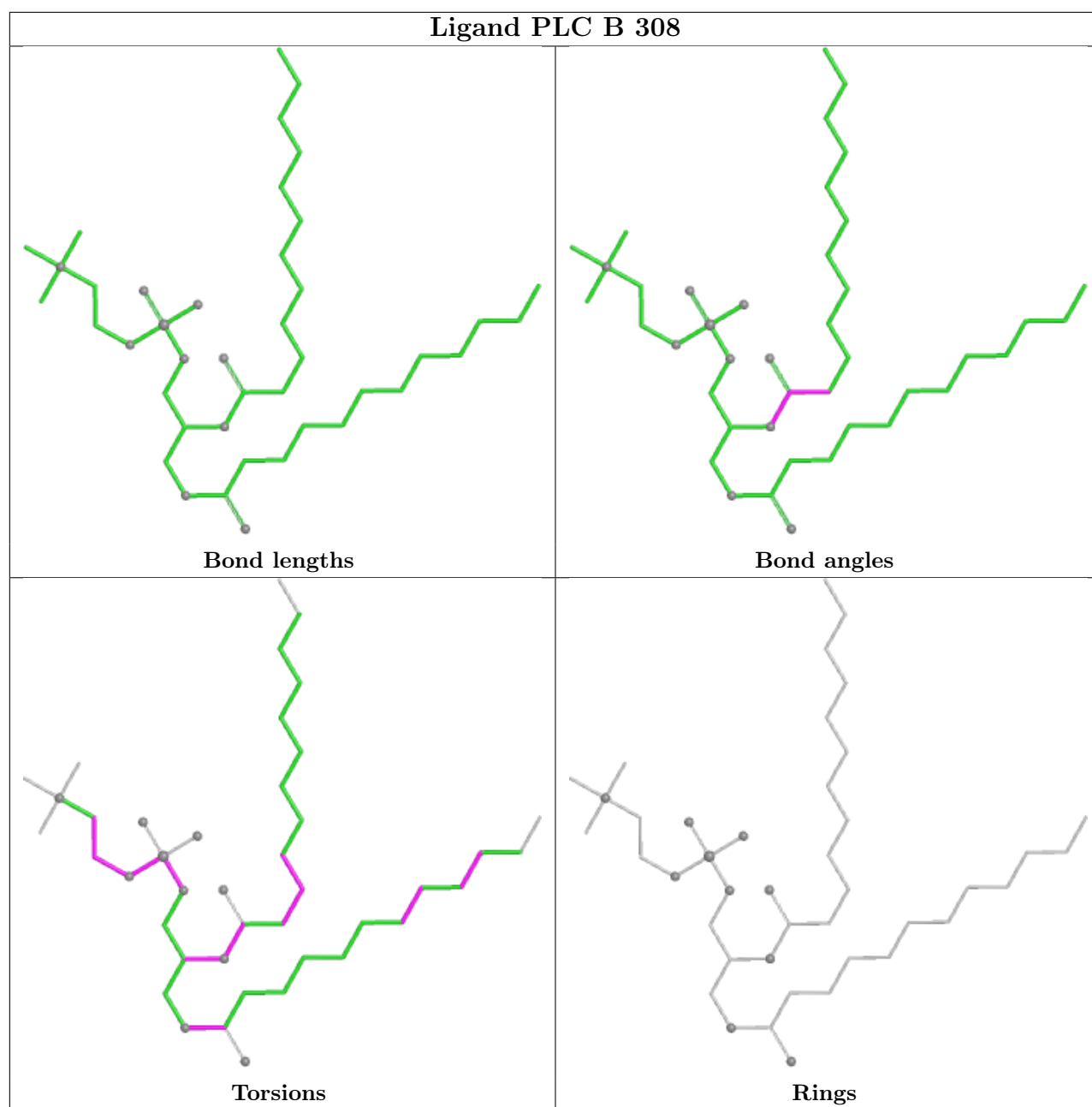


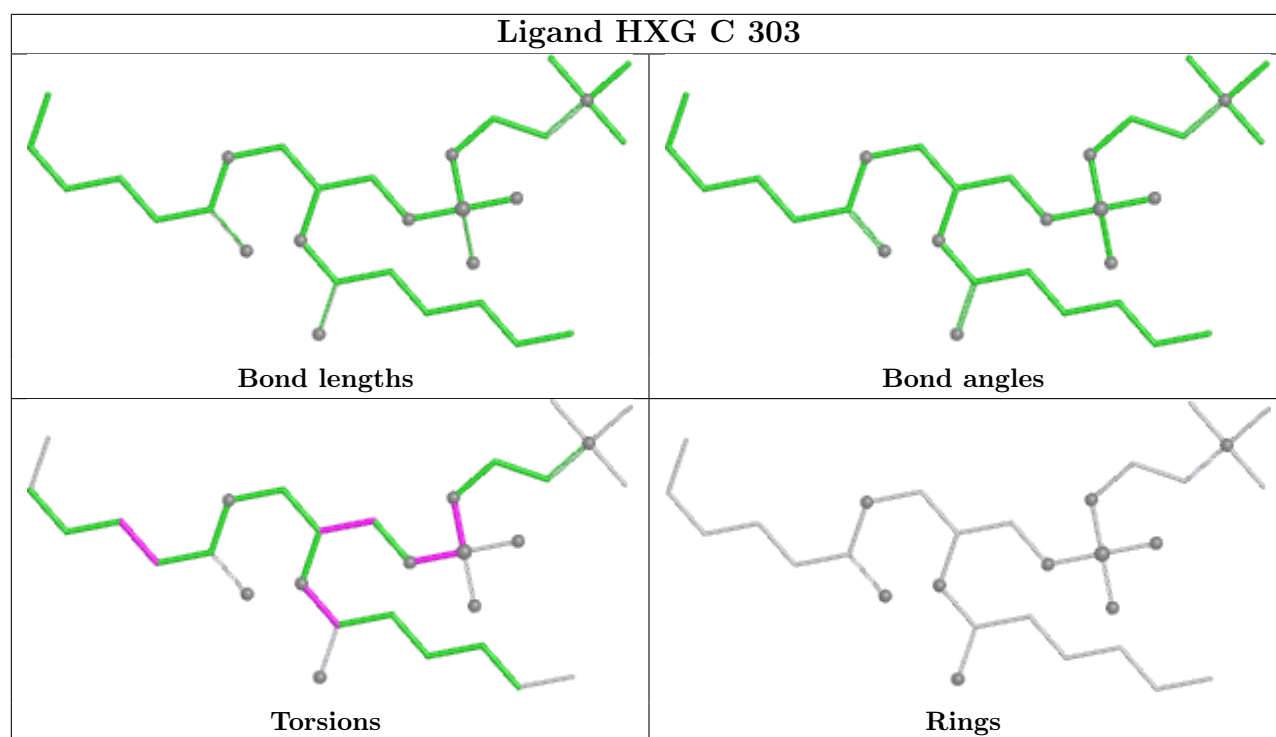


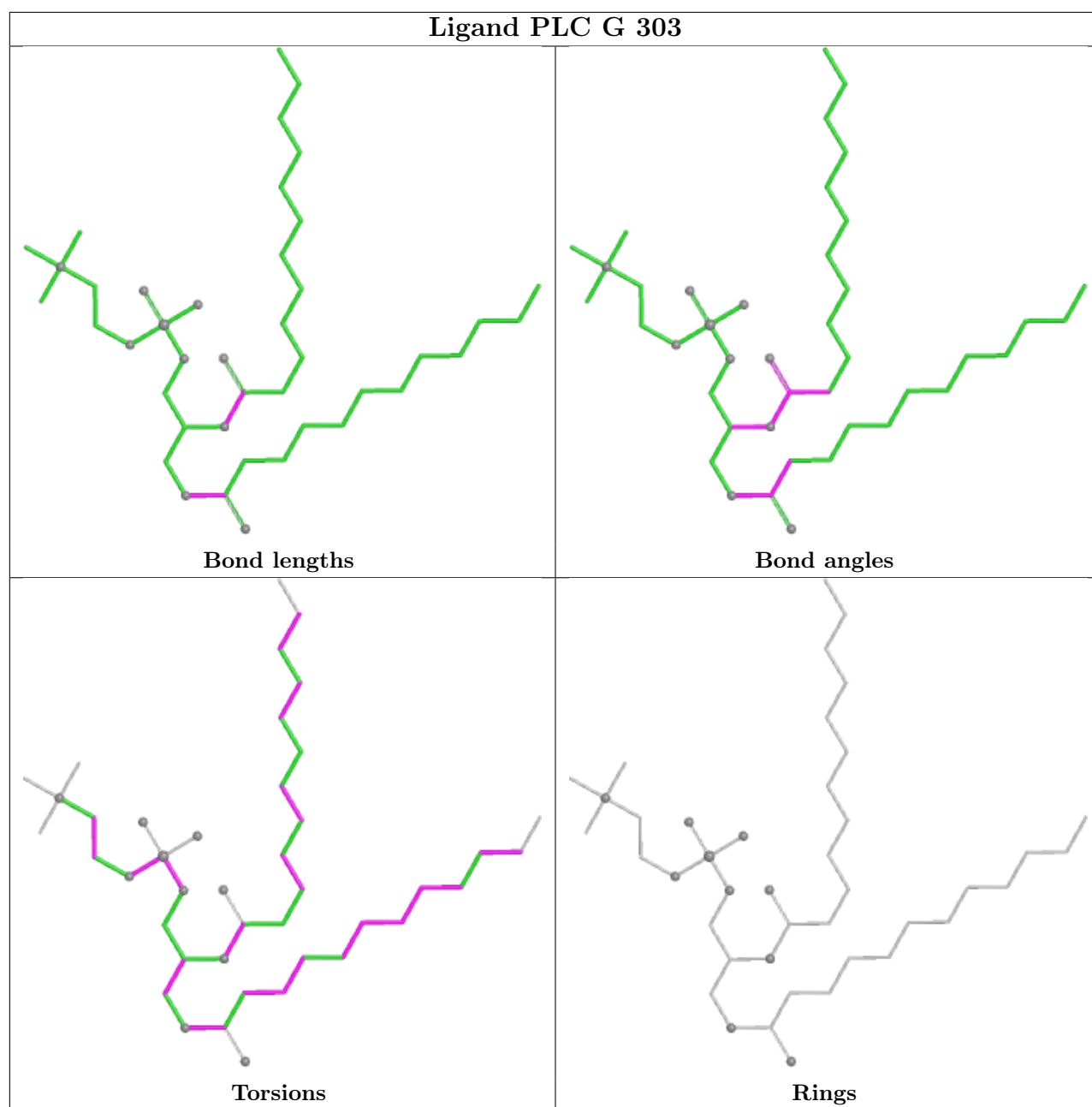


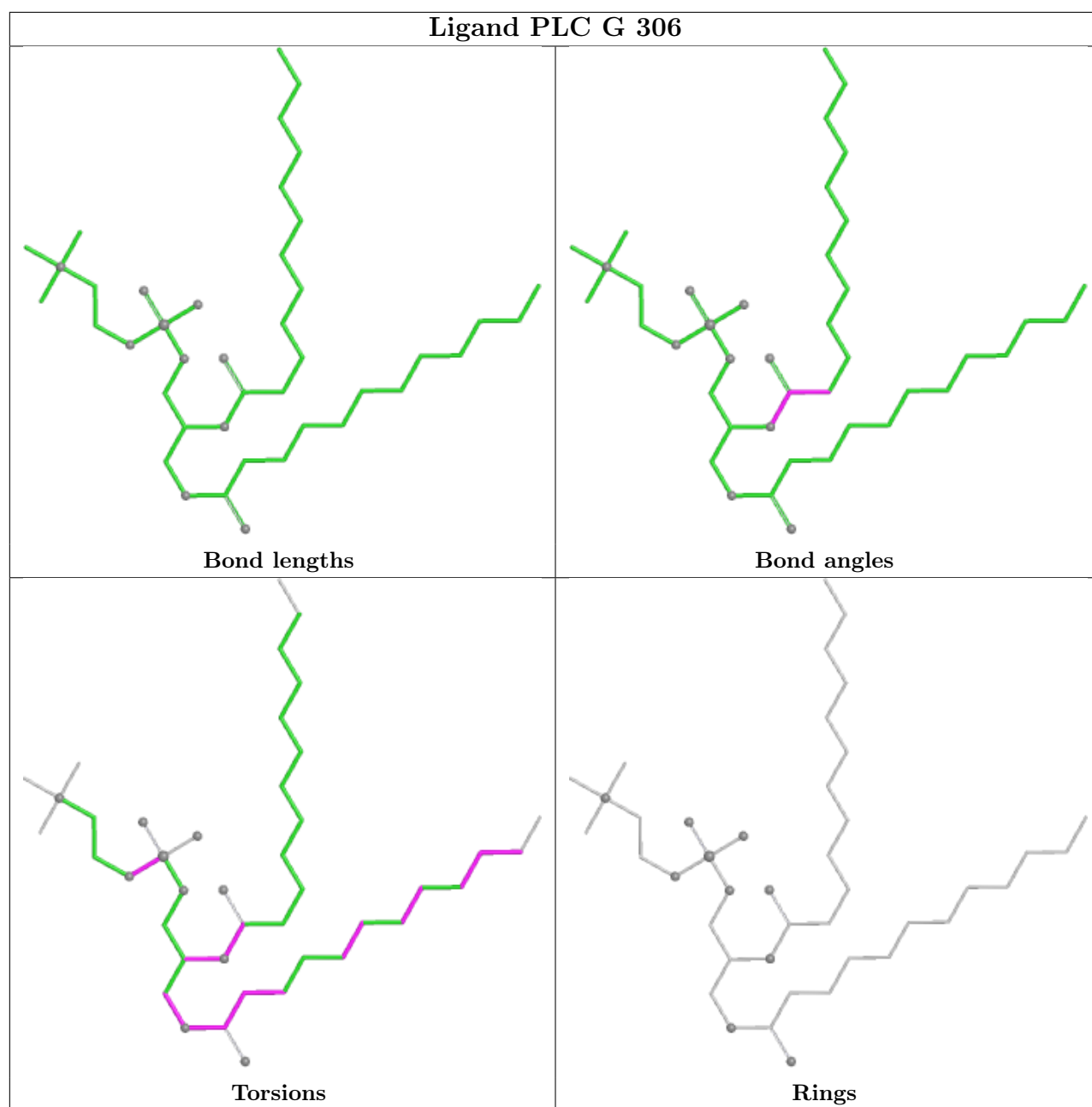


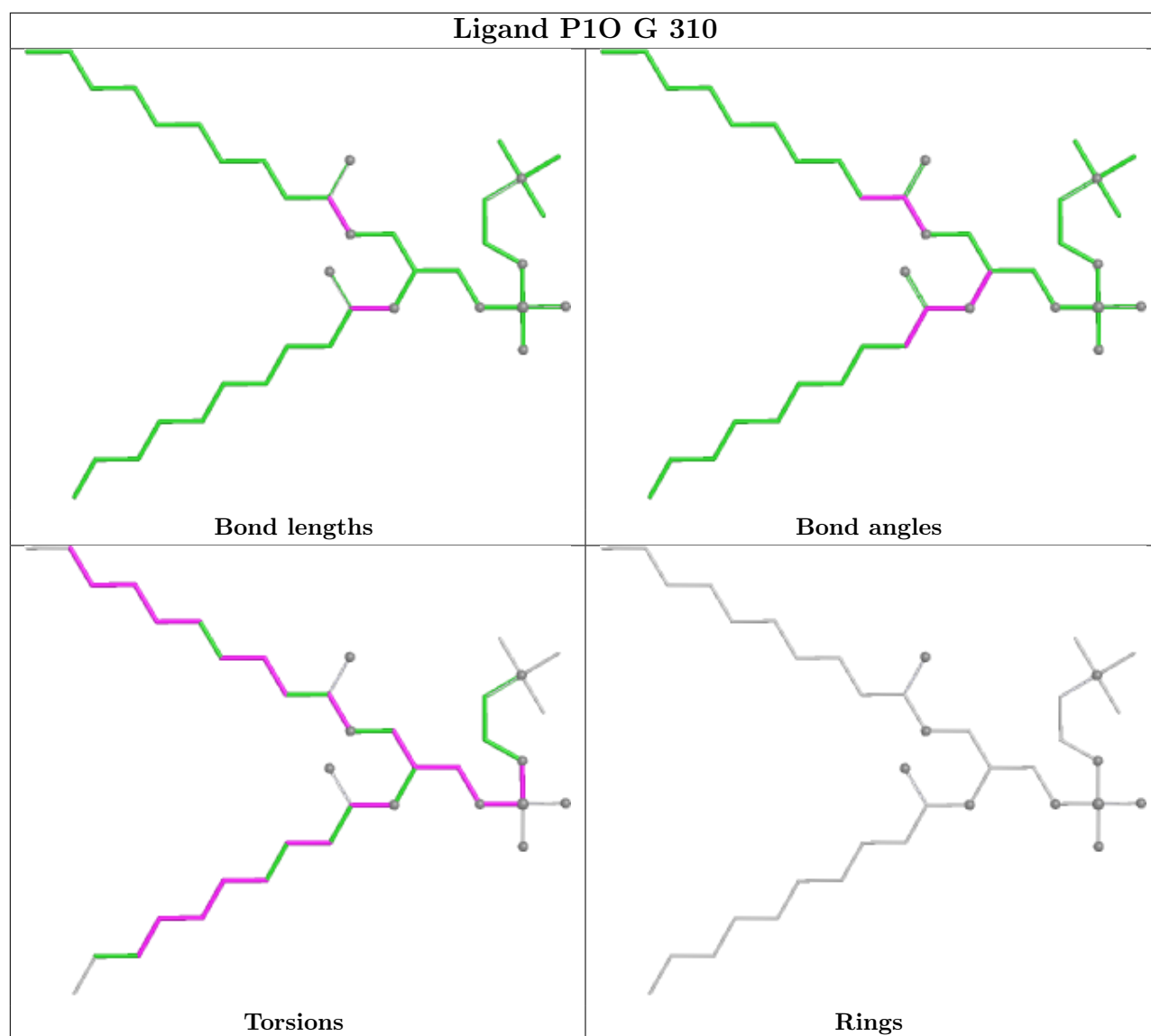


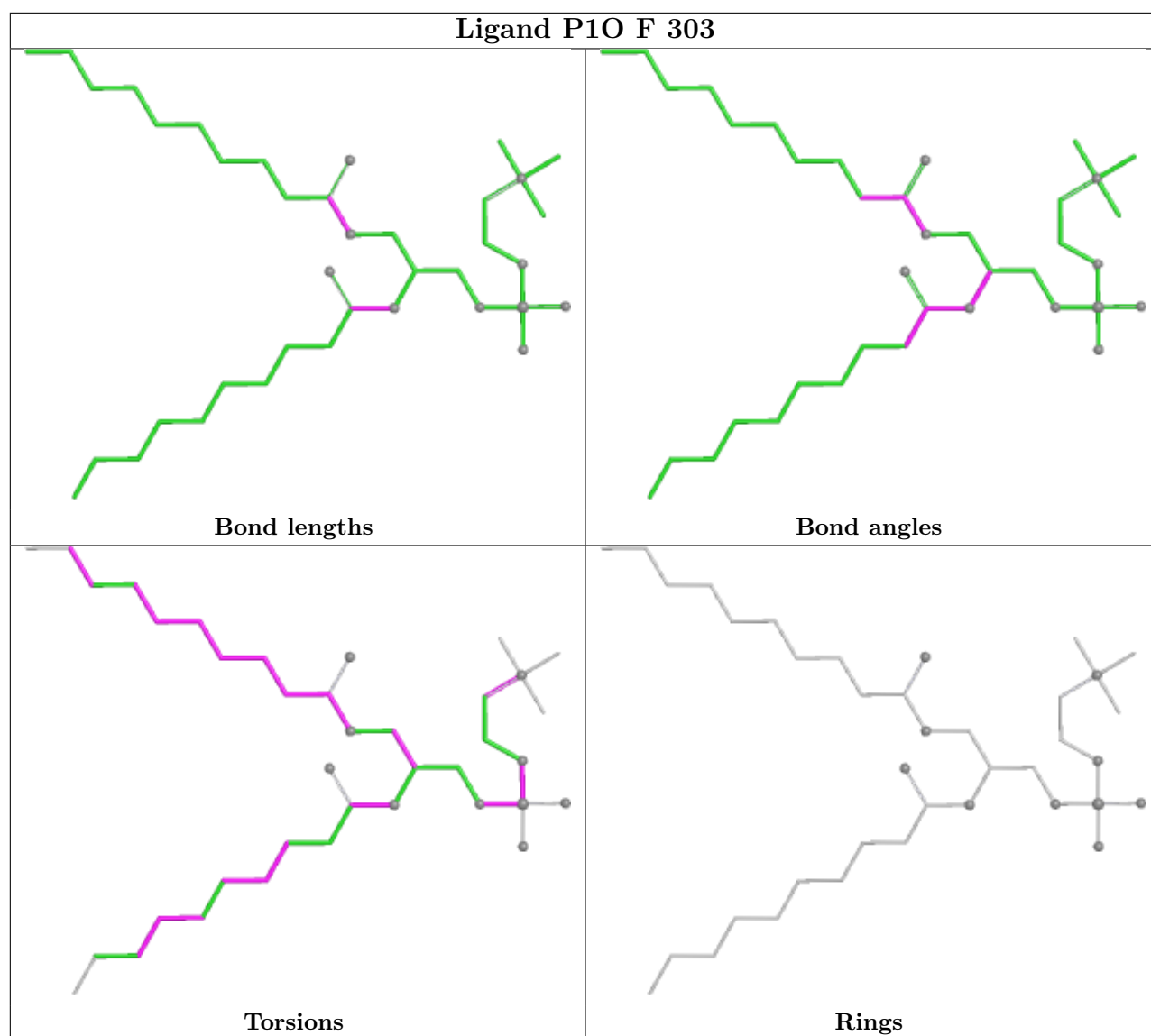


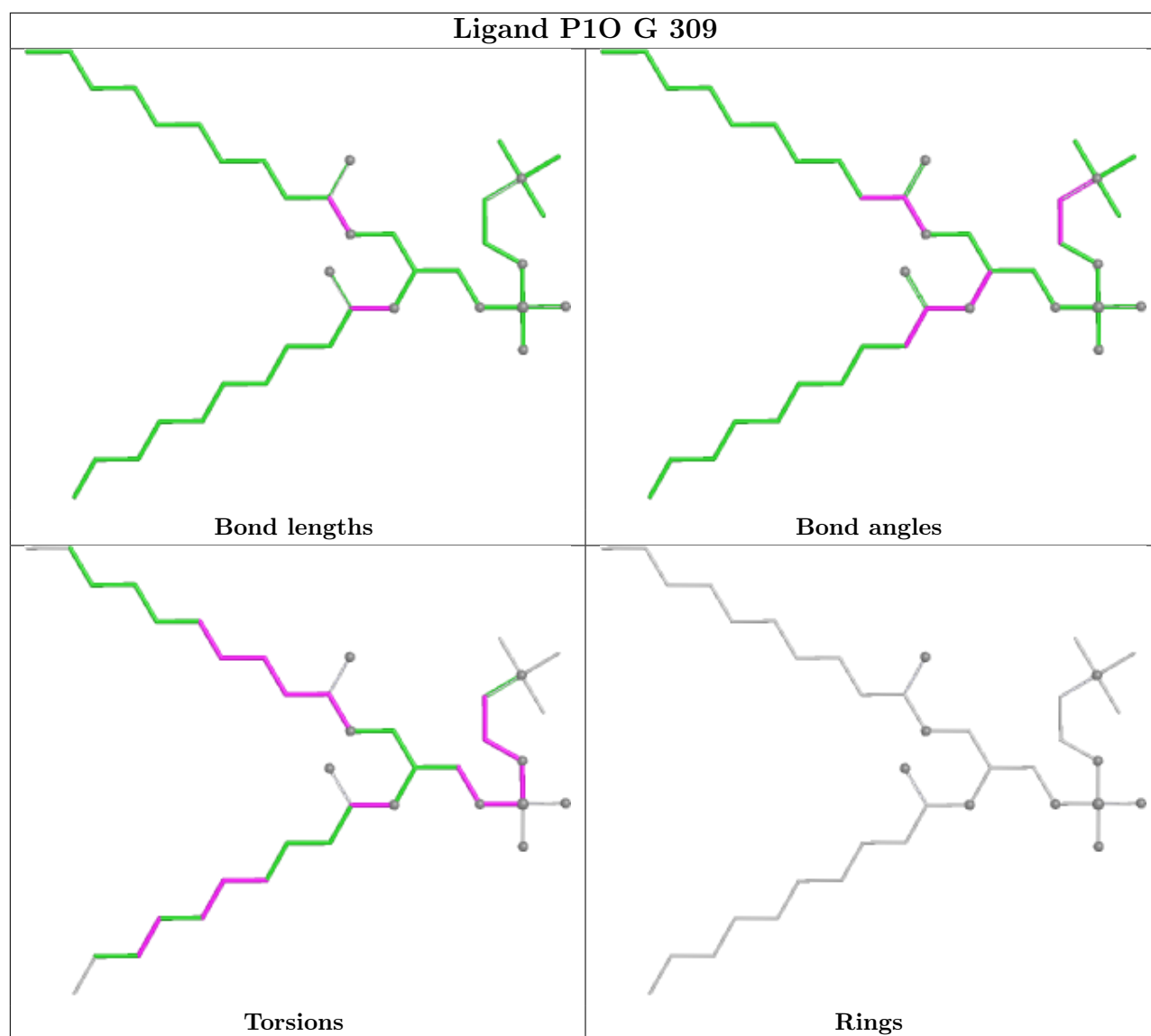


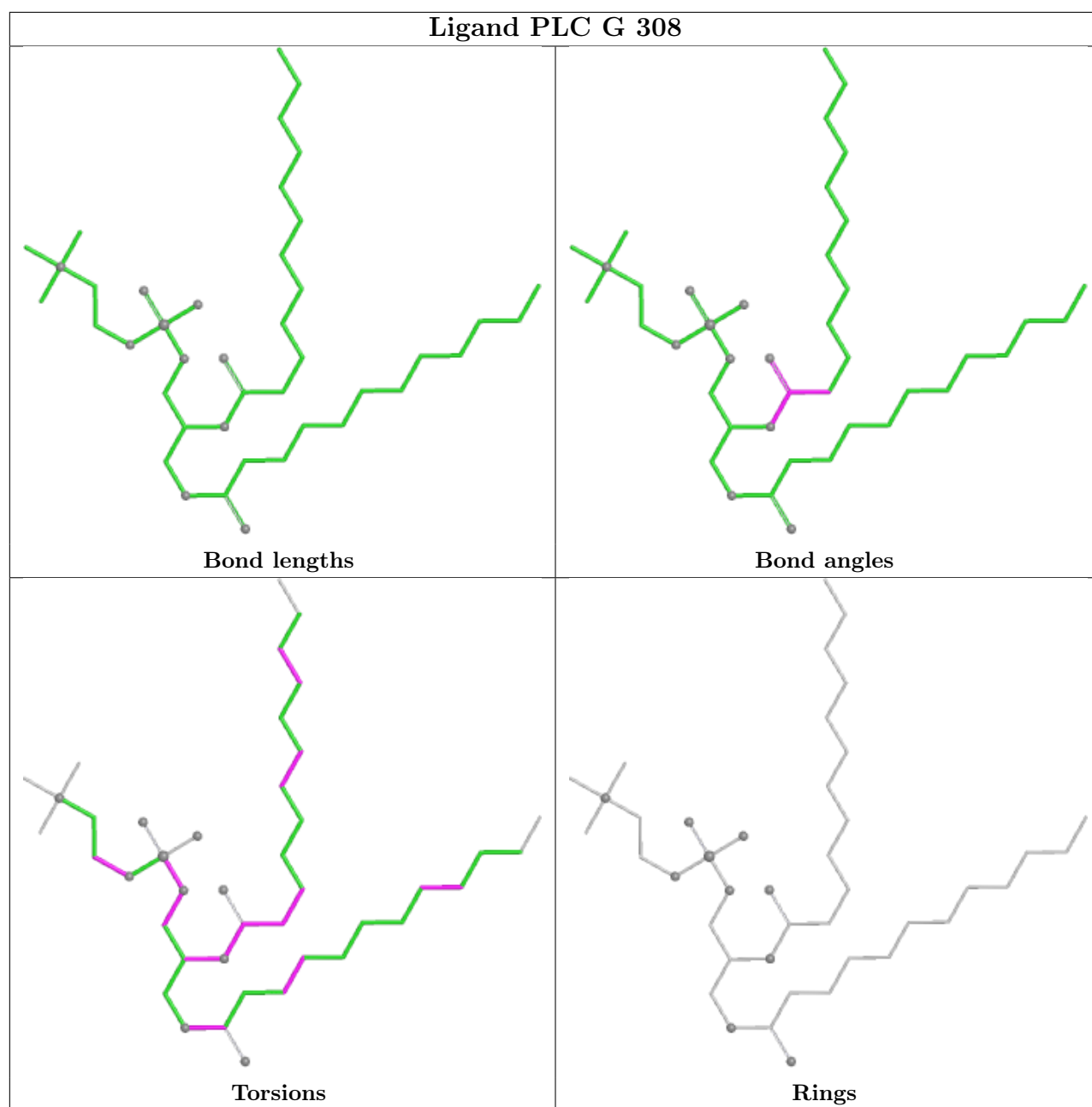


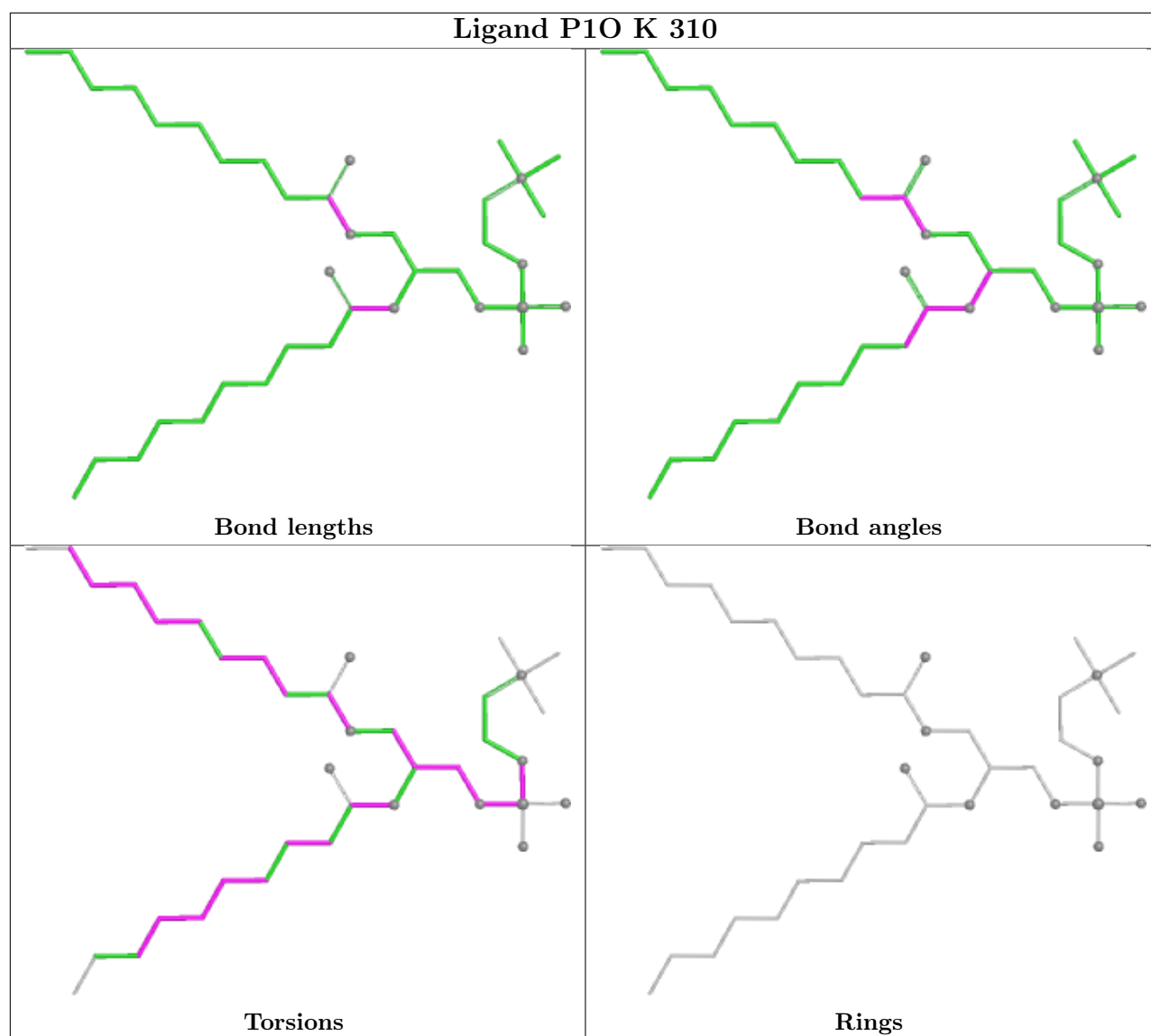


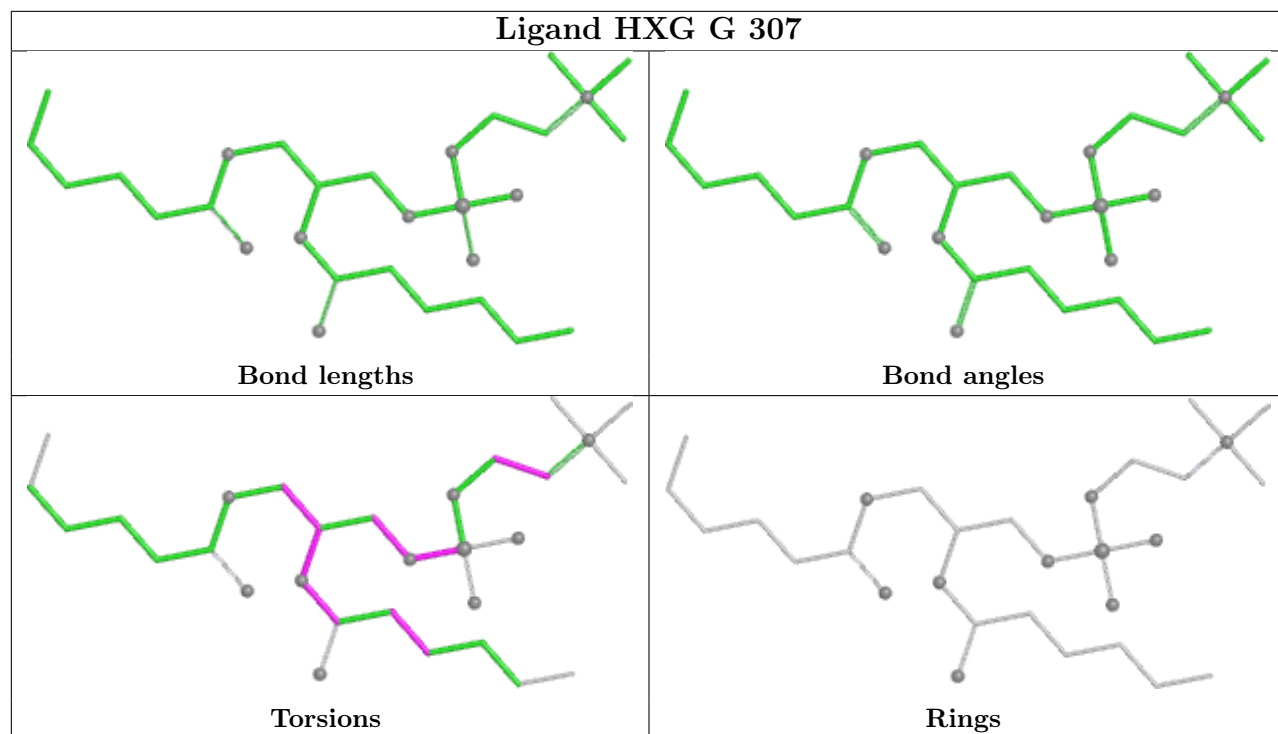


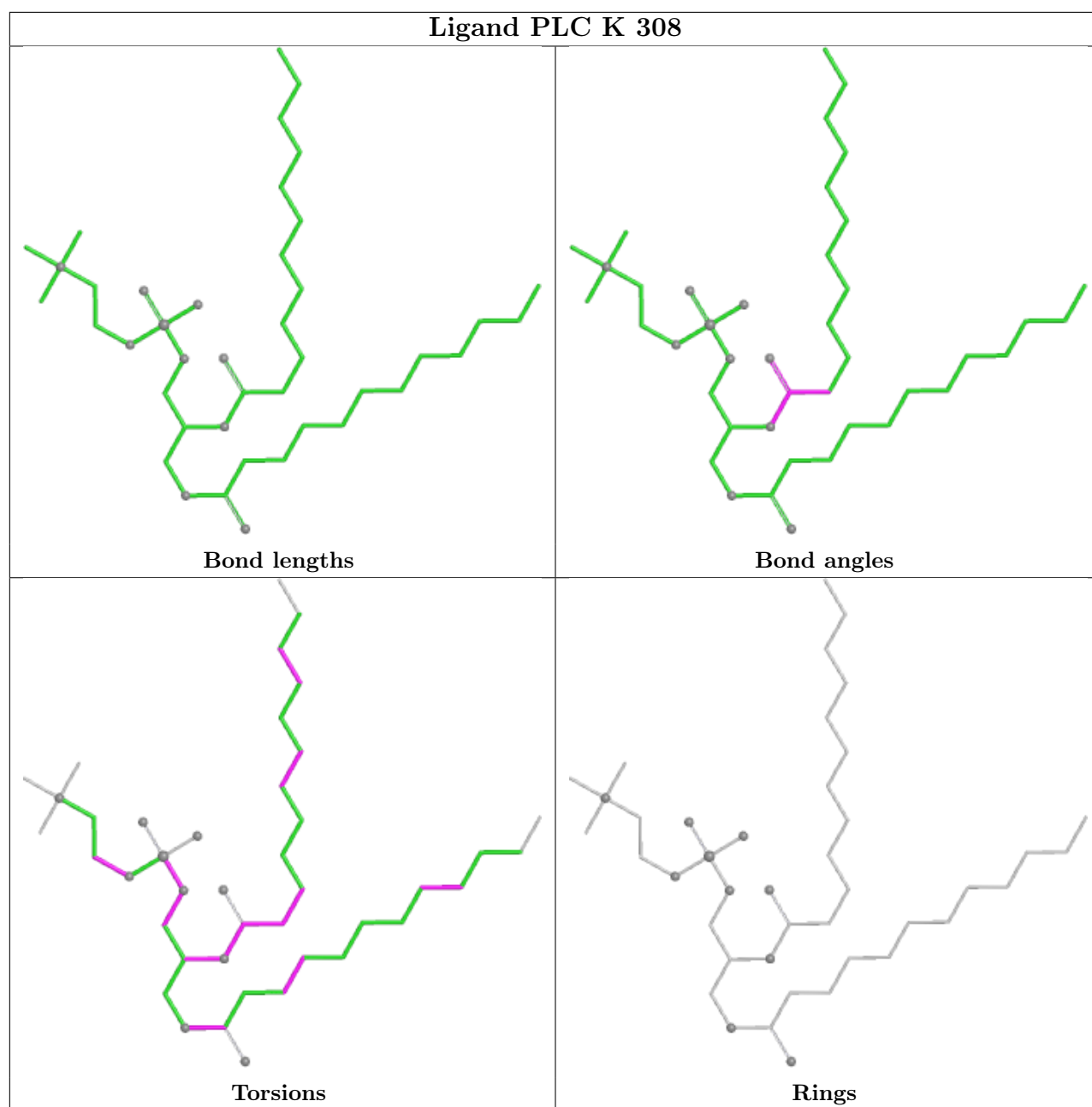


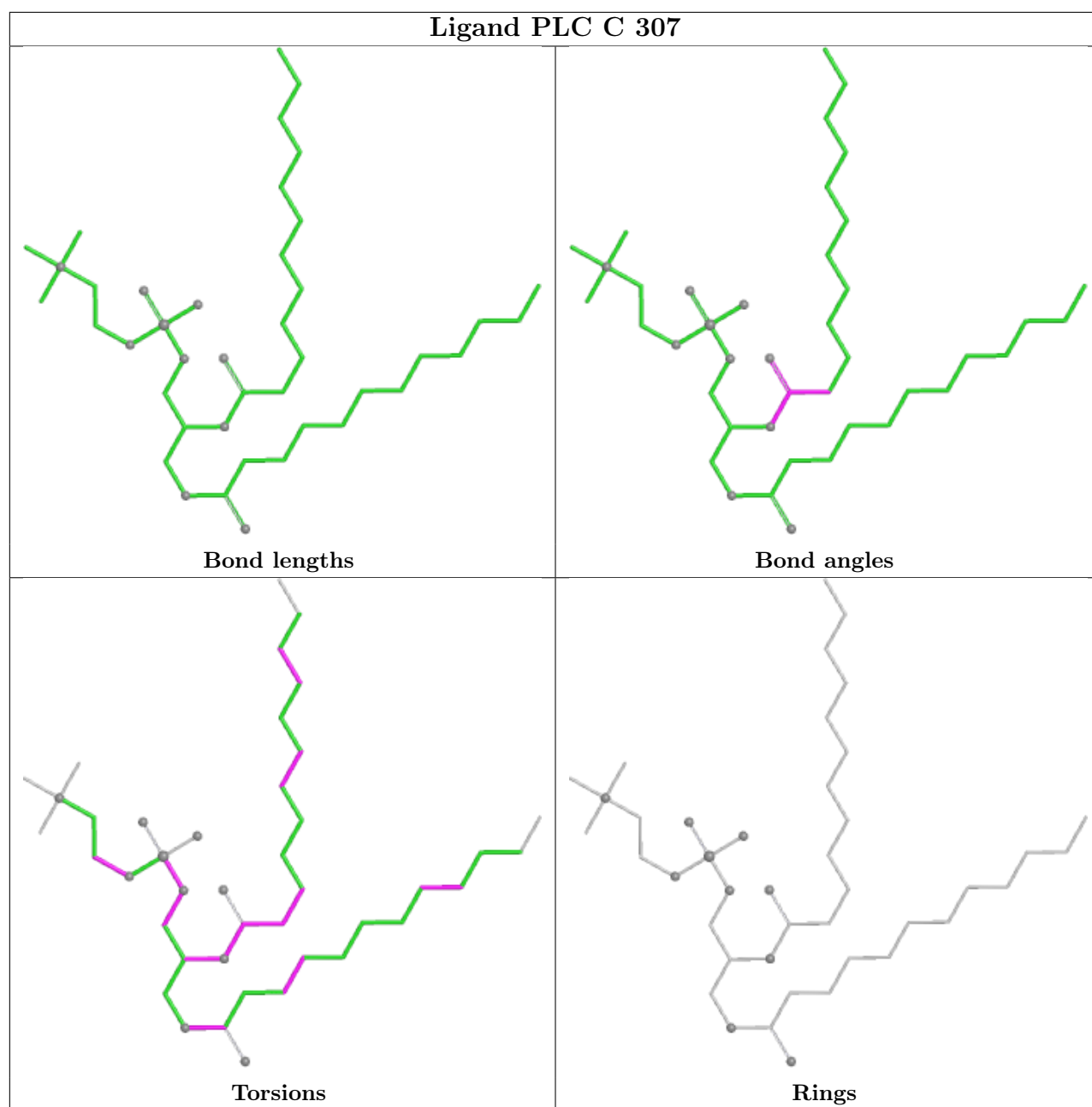


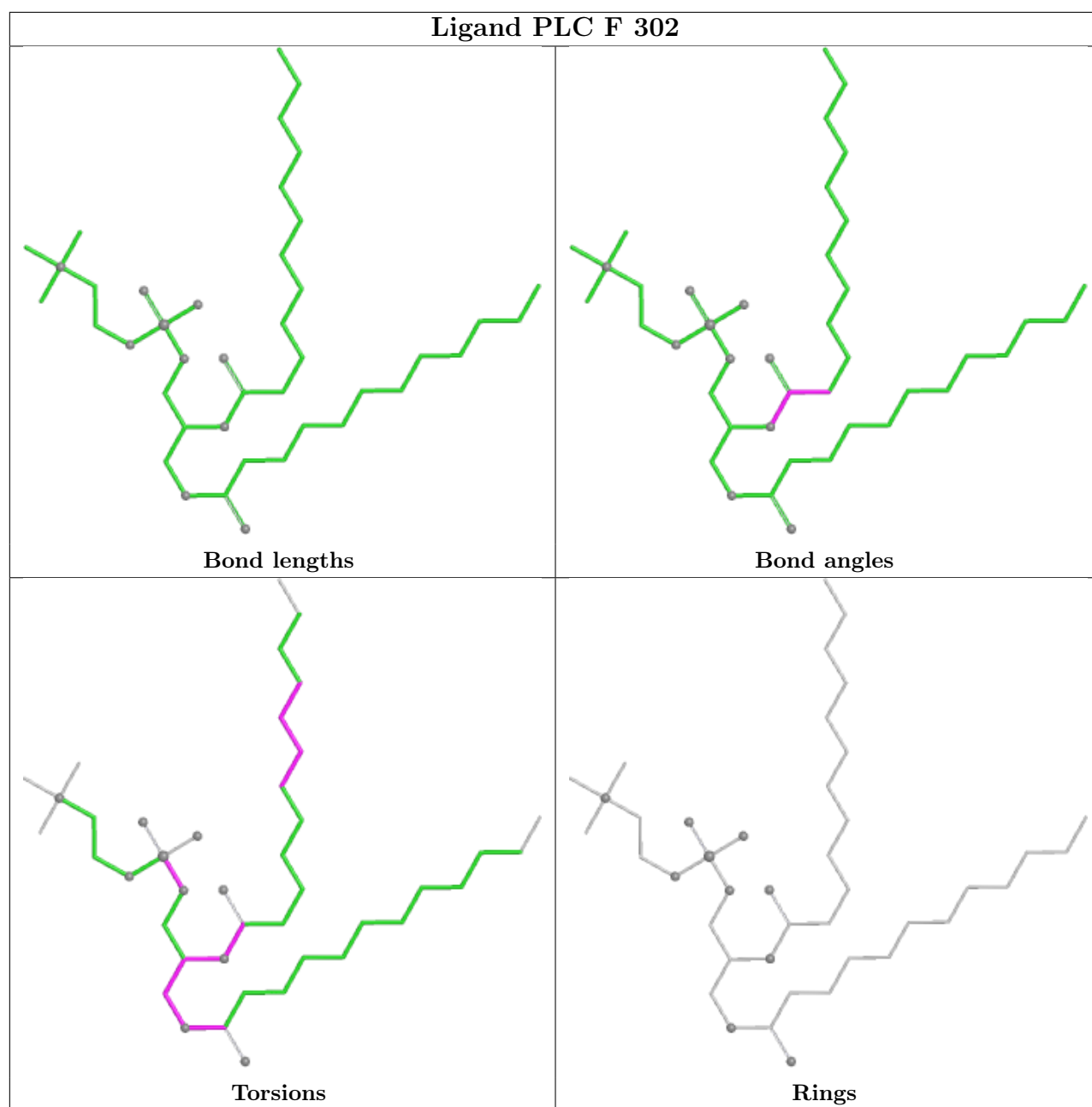


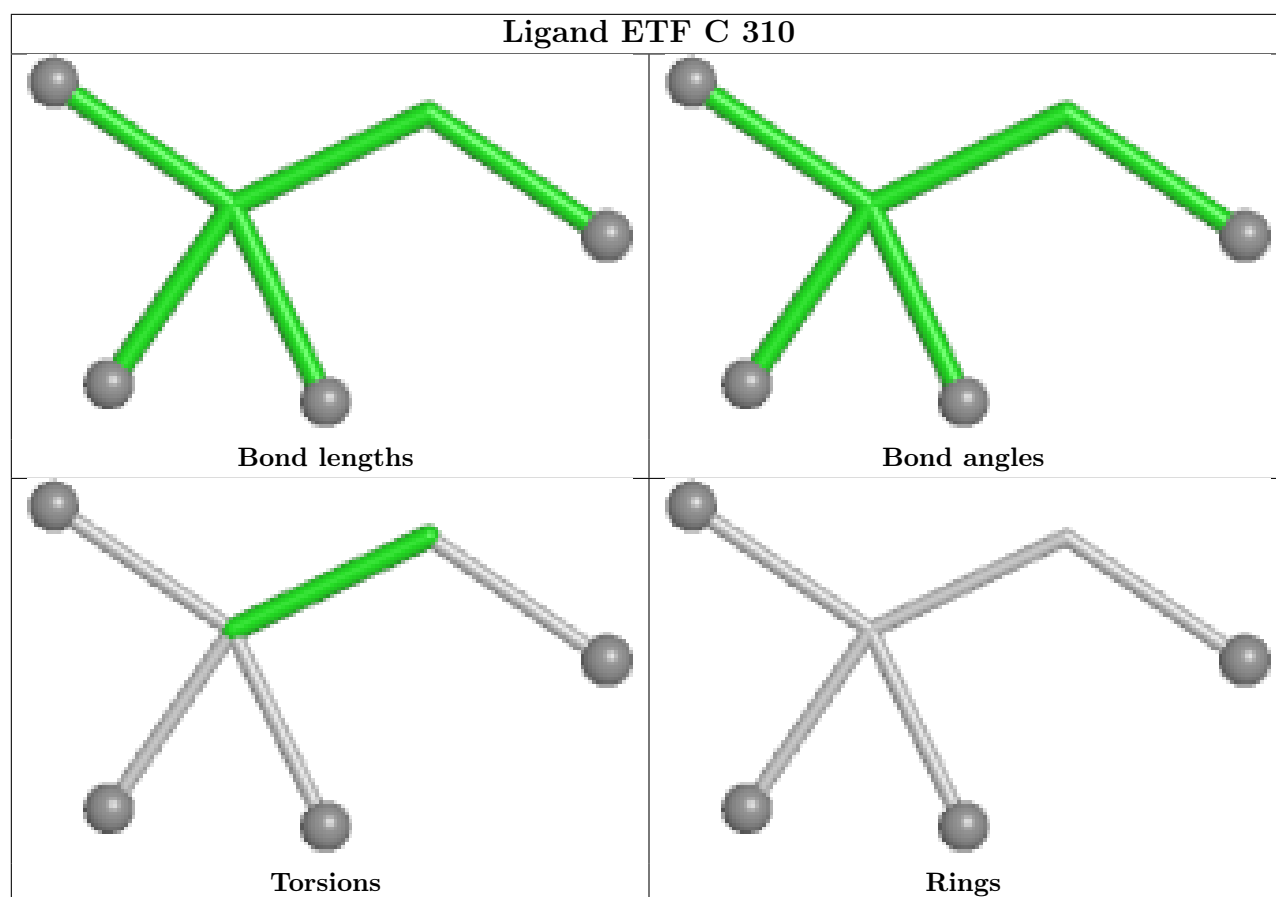












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

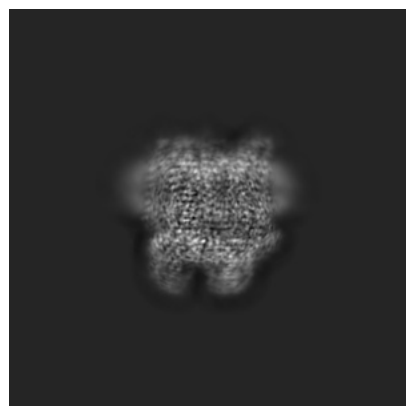
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40714. These allow visual inspection of the internal detail of the map and identification of artifacts.

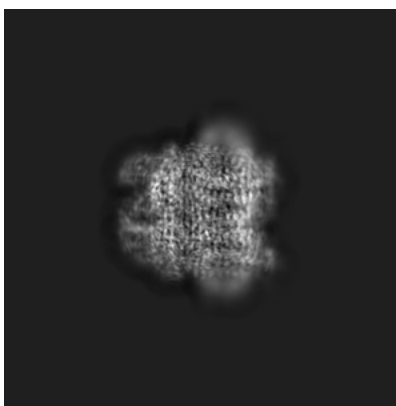
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

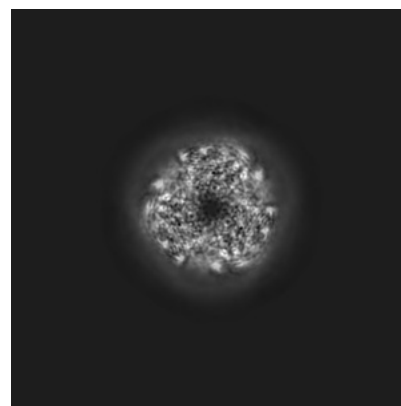
6.1.1 Primary map



X

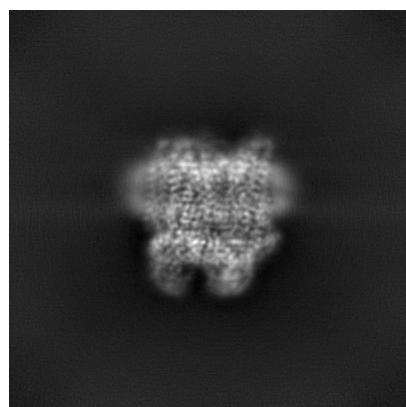


Y

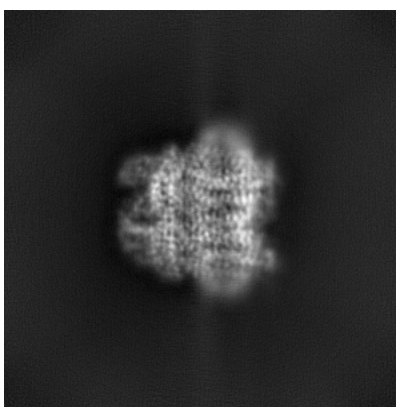


Z

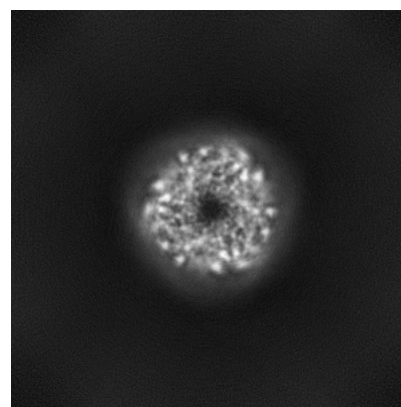
6.1.2 Raw map



X



Y

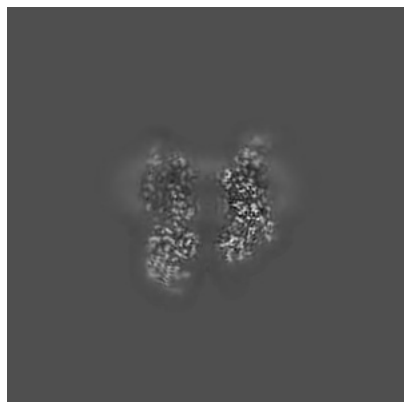


Z

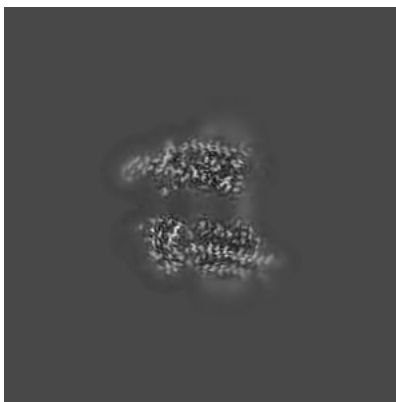
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

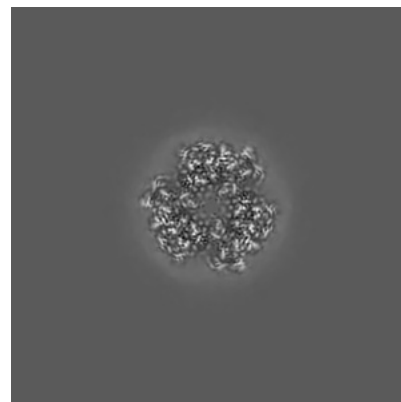
6.2.1 Primary map



X Index: 256

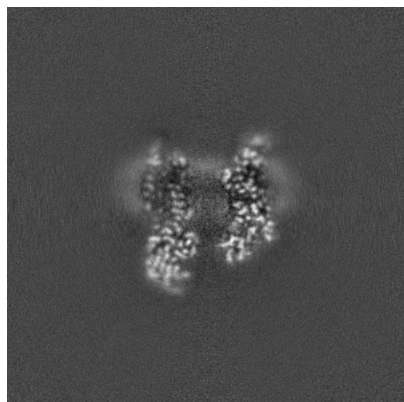


Y Index: 256

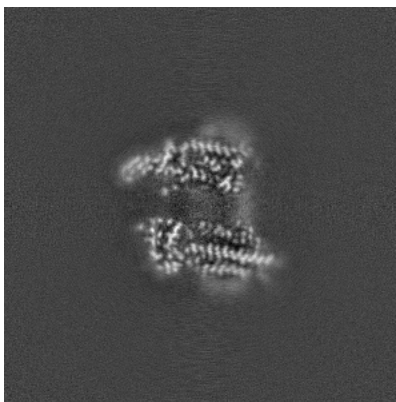


Z Index: 256

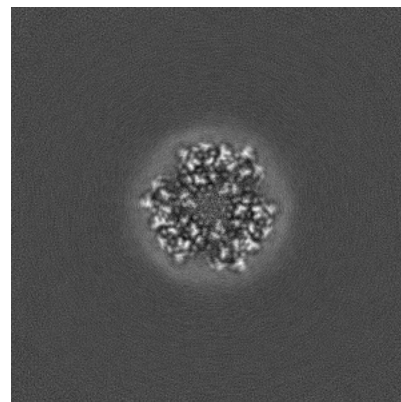
6.2.2 Raw map



X Index: 256



Y Index: 256

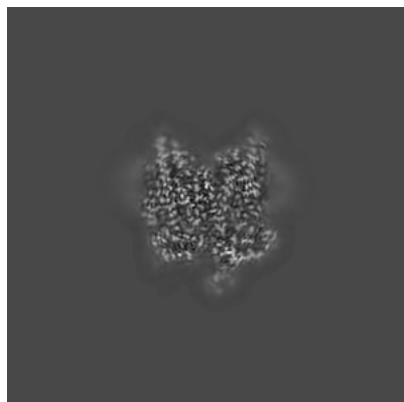


Z Index: 256

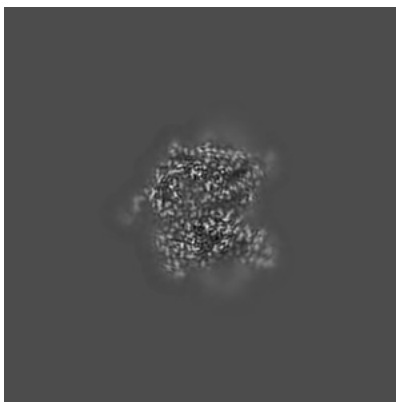
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

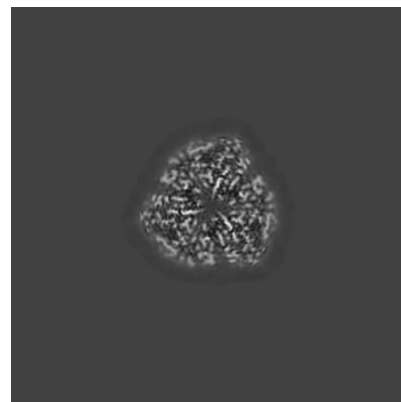
6.3.1 Primary map



X Index: 287

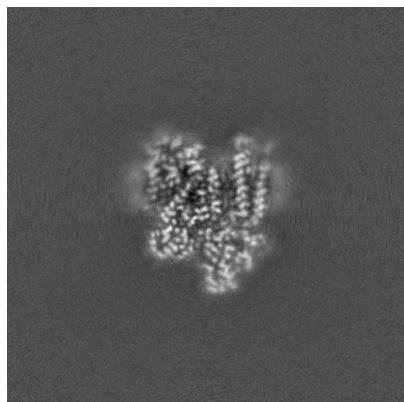


Y Index: 229

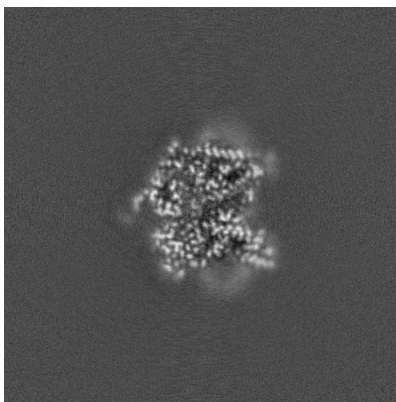


Z Index: 215

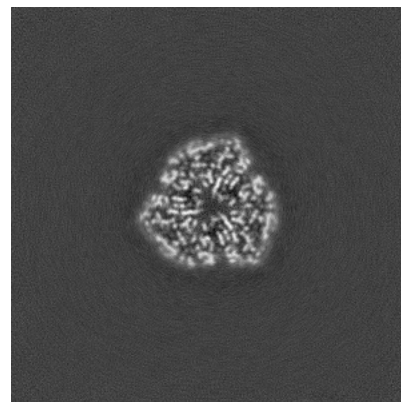
6.3.2 Raw map



X Index: 300



Y Index: 231

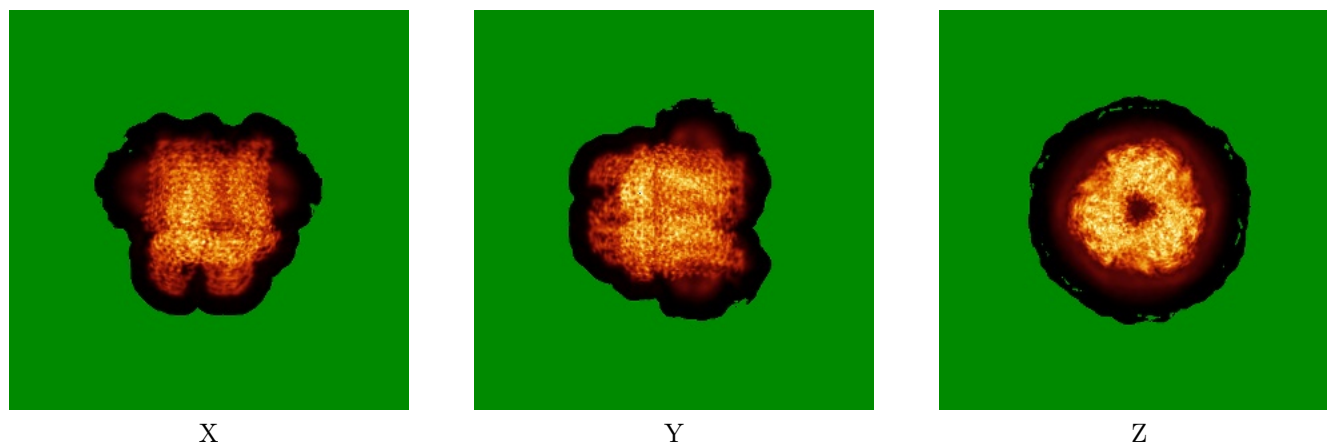


Z Index: 215

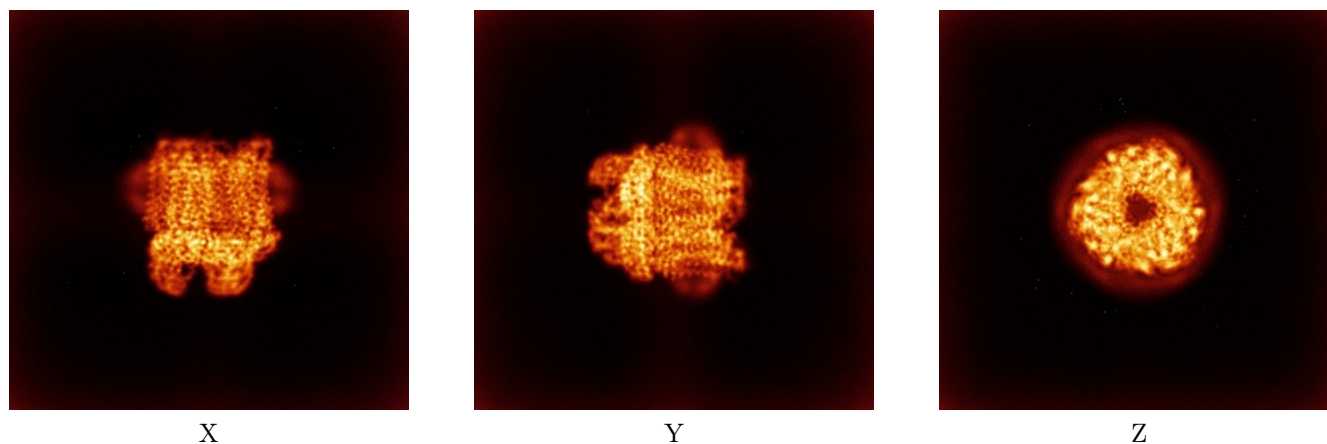
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

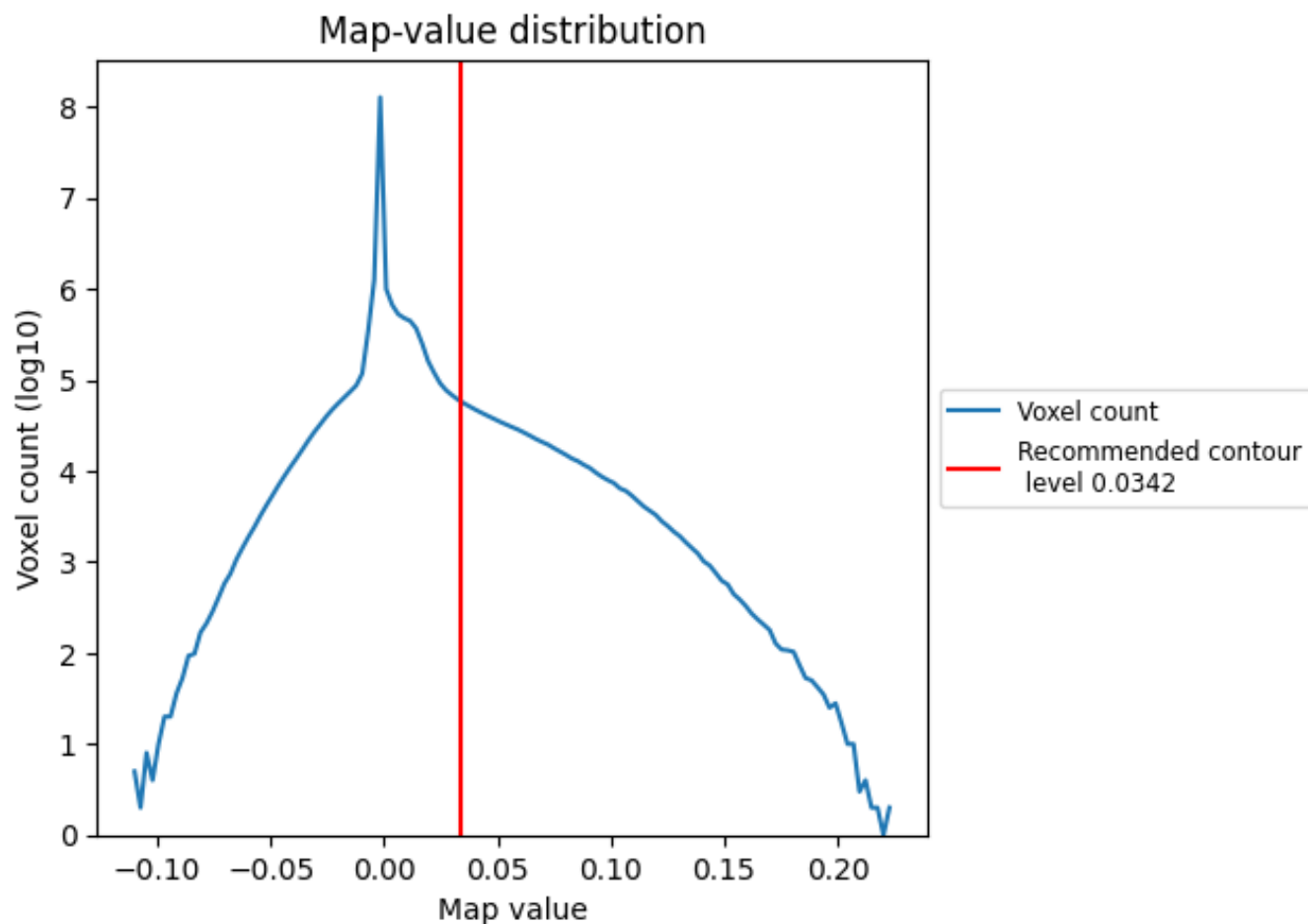
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

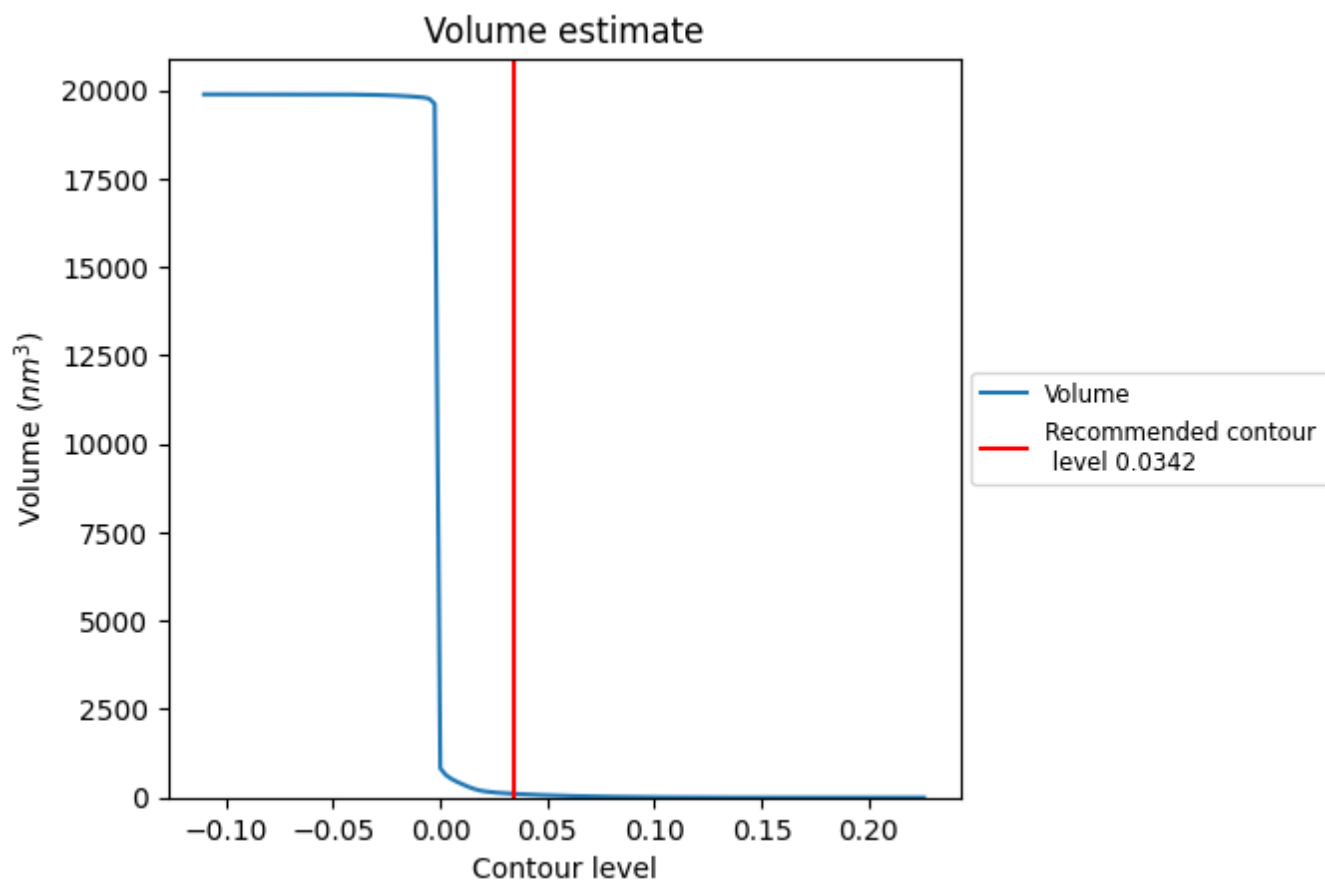
This section contains the results of statistical analysis of the map.

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.

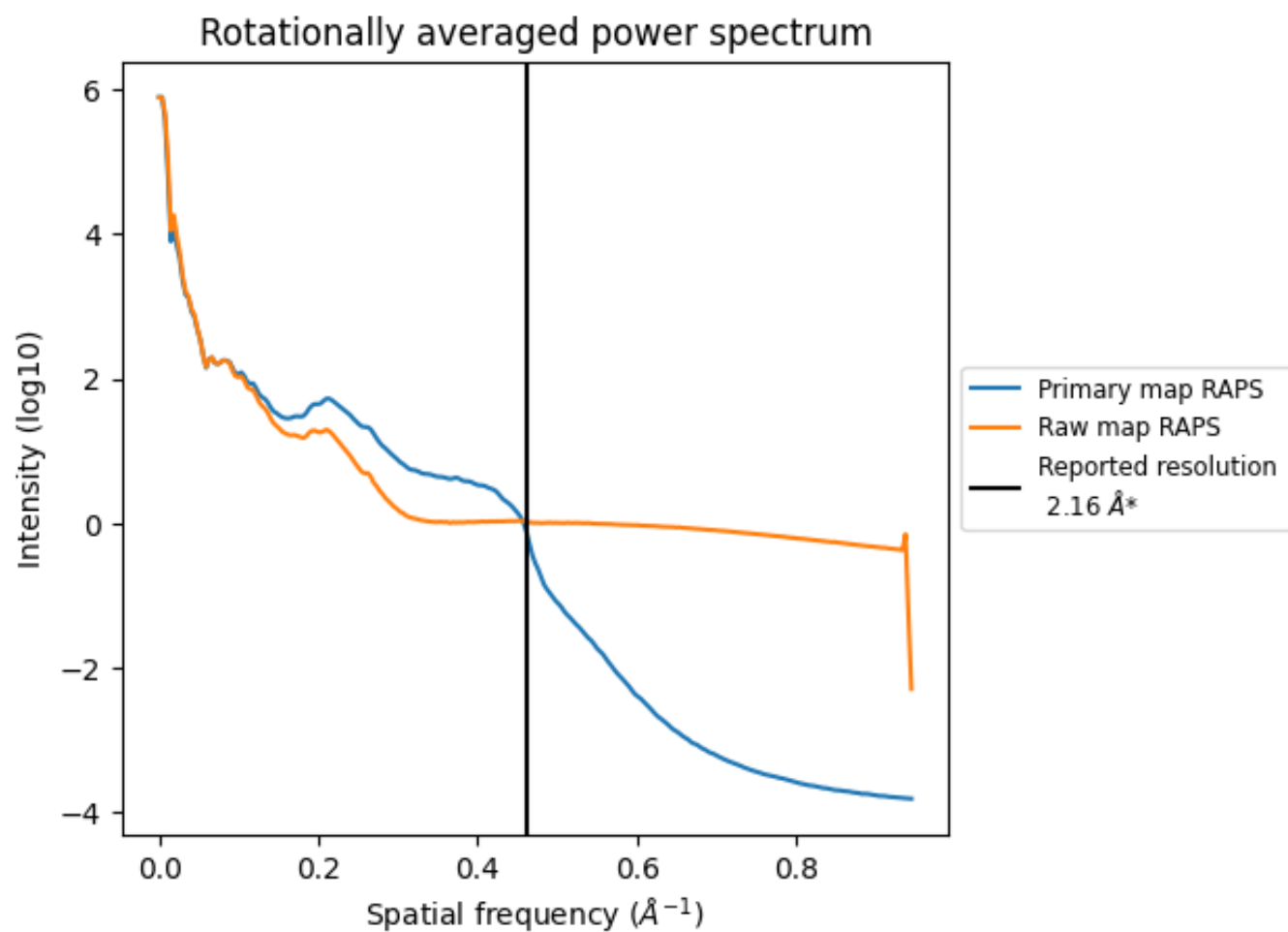
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 107 nm^3 ; this corresponds to an approximate mass of 97 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 ⓘ

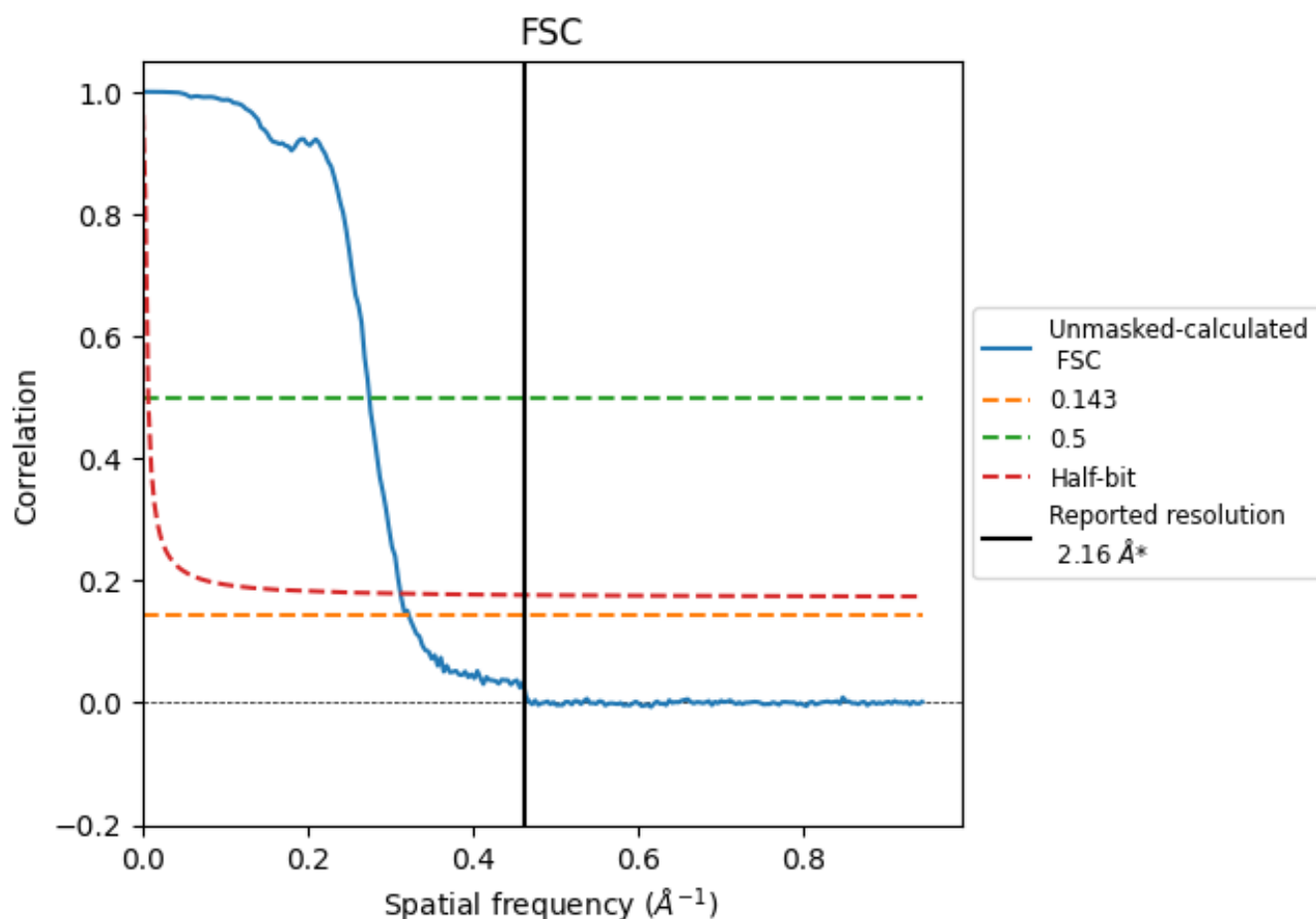


*Reported resolution corresponds to spatial frequency of $0.463 \text{ \AA}^{-1}$

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 0.463 Å⁻¹

8.2 Resolution estimates [i](#)

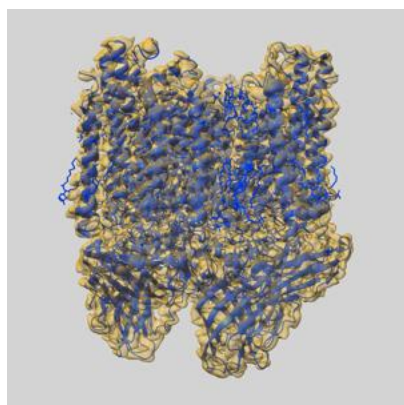
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.16	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.10	3.64	3.19

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.10 differs from the reported value 2.16 by more than 10 %

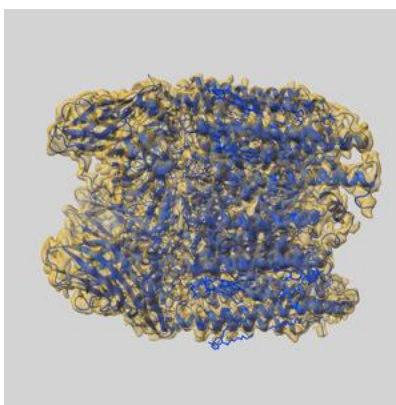
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-40714 and PDB model 8SQW. Per-residue inclusion information can be found in section [3](#) on page [11](#).

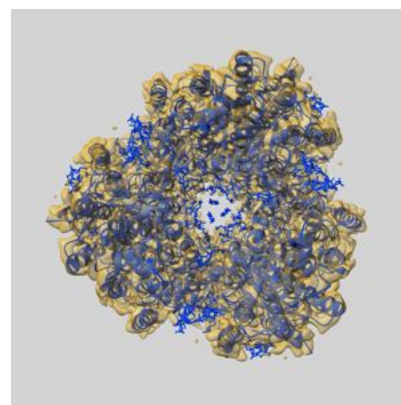
9.1 Map-model overlay [i](#)



X



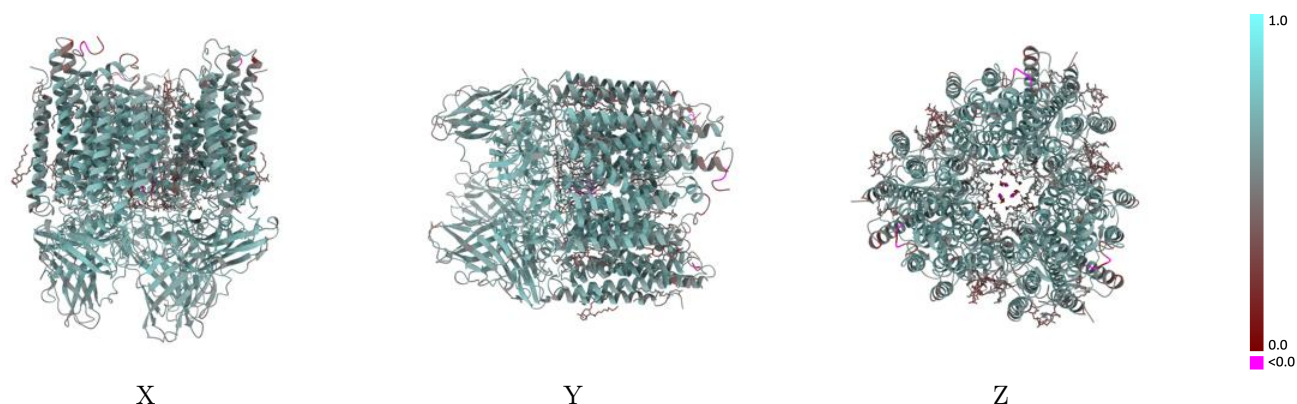
Y



Z

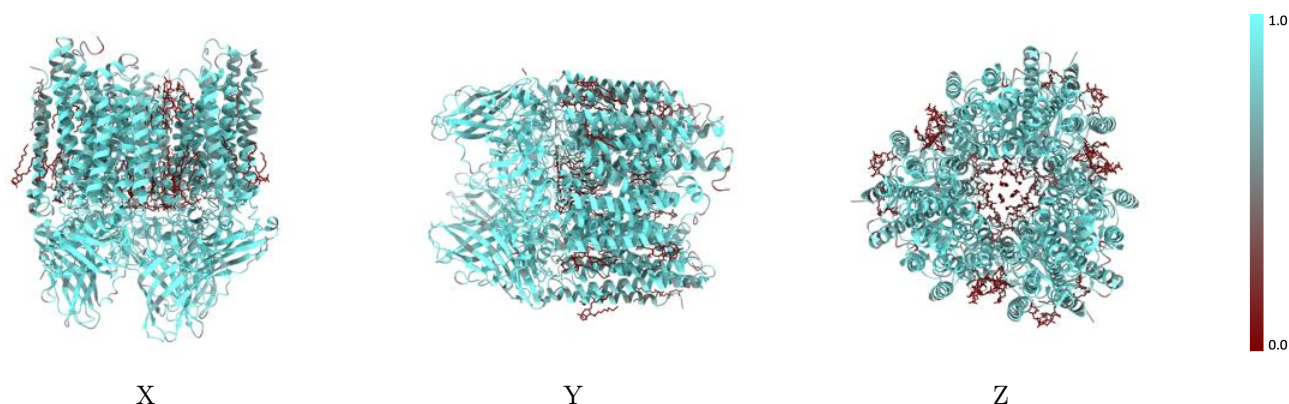
The images above show the 3D surface view of the map at the recommended contour level 0.0342 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



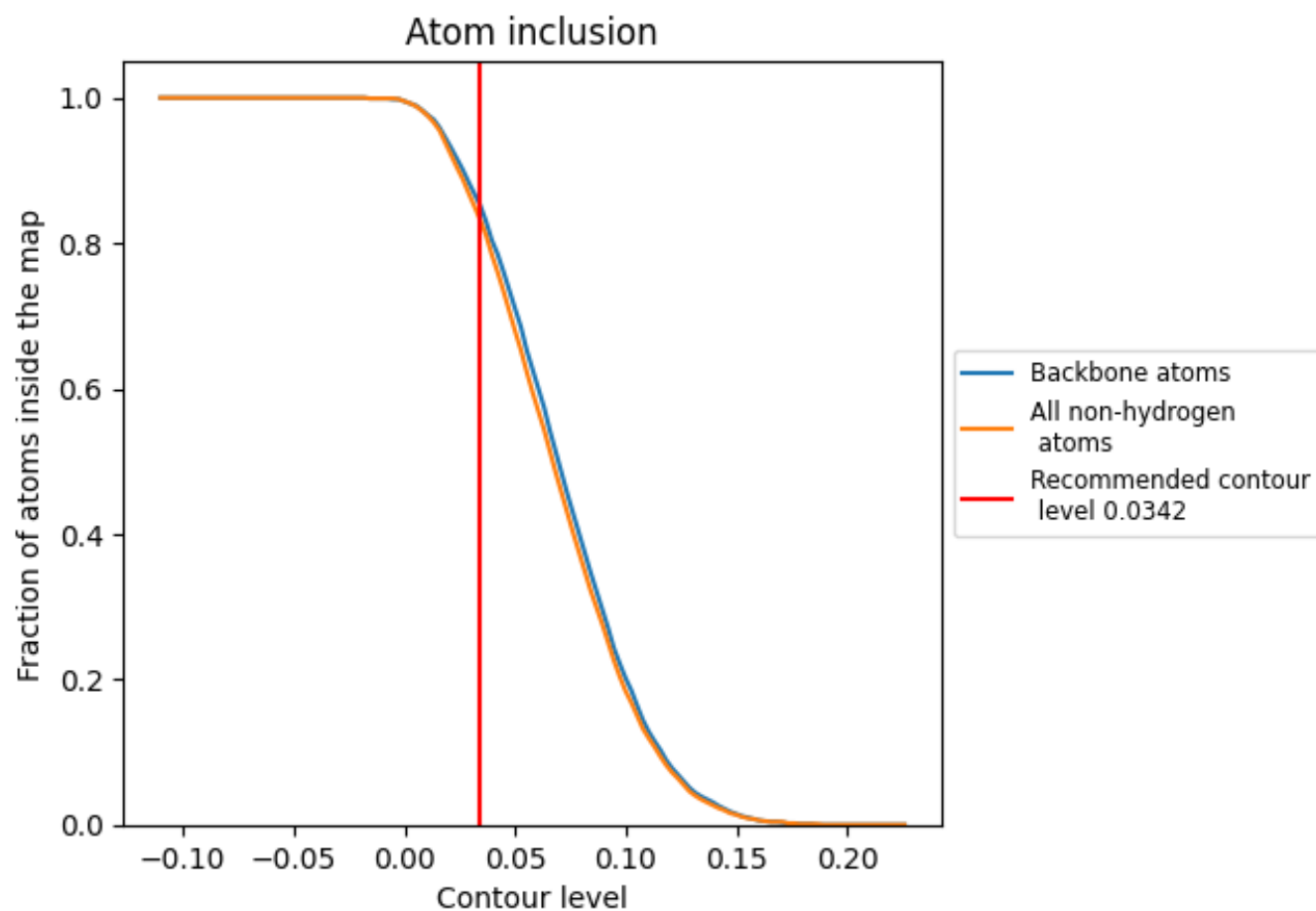
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0342).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 83% 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 (0.0342) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8320	0.6090
A	0.8790	0.6140
B	0.8630	0.6440
C	0.7380	0.5690
E	0.8780	0.6130
F	0.8630	0.6440
G	0.7350	0.5700
I	0.8790	0.6140
J	0.8640	0.6430
K	0.7350	0.5700

1.0

0.0

<0.0