



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 09:55 AM UTC

PDB ID : 8OYI / pdb_00008oyi
EMDB ID : EMD-17287
Title : particulate methane monooxygenase with 2,2,2-trifluoroethanol bound
Authors : Tucci, F.J.; Rosenzweig, A.C.
Deposited on : 2023-05-04
Resolution : 2.19 Å(reported)

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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:

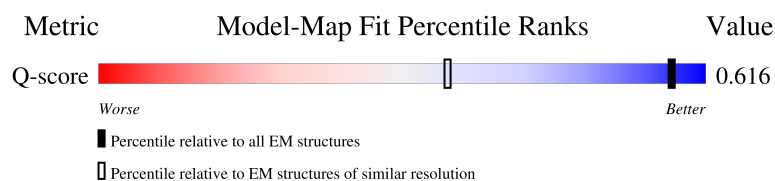
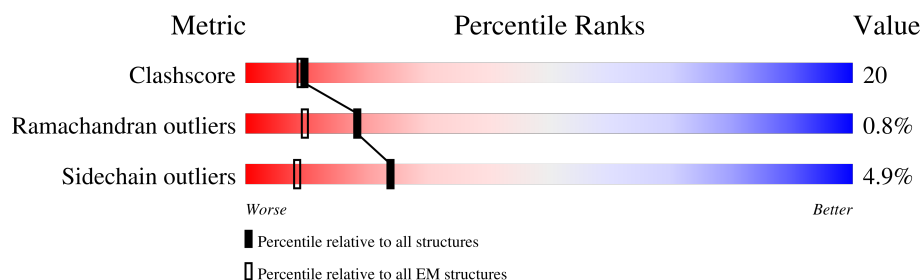
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 2.19 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.



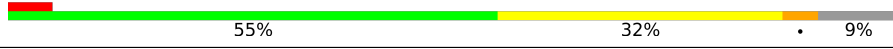
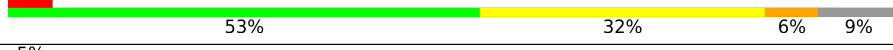
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2745 (1.70 - 2.69)

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	414	
1	E	414	
1	I	414	
2	B	247	

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Mol	Chain	Length	Quality of chain
2	F	247	
2	J	247	
3	C	260	
3	G	260	
3	K	260	

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	309	-	-	X	-
7	P1O	J	301	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 25318 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	C	N	O	S	0	0
			3017	1938	513	551	15		
1	E	382	Total	C	N	O	S	0	0
			3017	1938	513	551	15		
1	I	382	Total	C	N	O	S	0	0
			3017	1938	513	551	15		

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

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

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

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

- Molecule 4 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

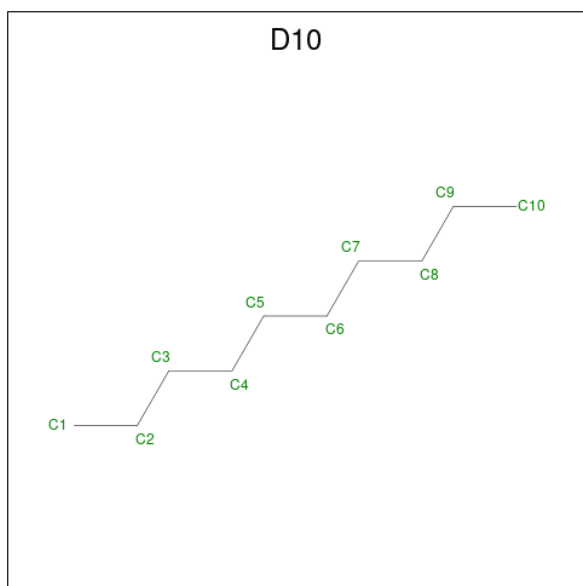
Mol	Chain	Residues	Atoms		AltConf
4	A	2	Total	Cu	0
			2	2	

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

- Molecule 5 is DECANE (CCD ID: D10) (formula: C₁₀H₂₂).



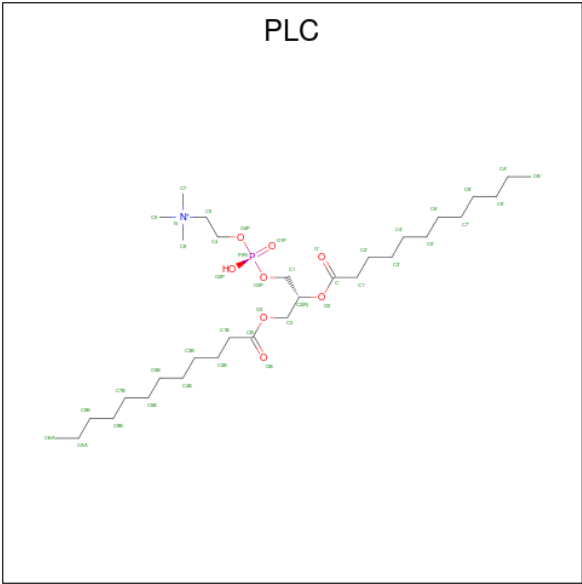
Mol	Chain	Residues	Atoms			AltConf
5	A	1	Total 32	C 10	H 22	0
5	B	1	Total 32	C 10	H 22	0
5	B	1	Total 32	C 10	H 22	0
5	B	1	Total 32	C 10	H 22	0
5	B	1	Total 32	C 10	H 22	0
5	C	1	Total 32	C 10	H 22	0

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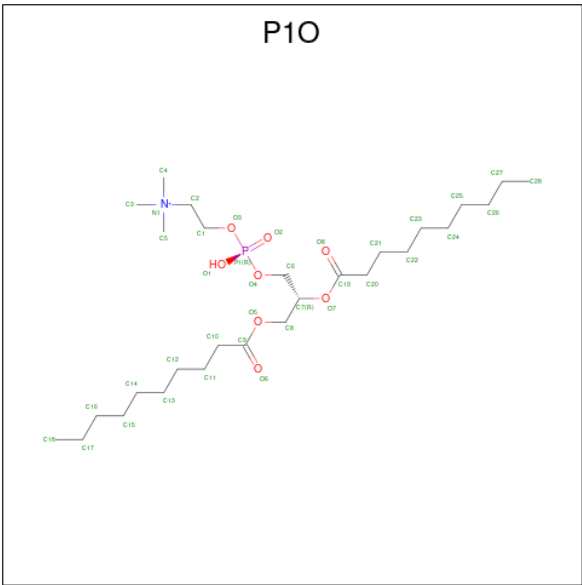
Mol	Chain	Residues	Atoms			AltConf
5	F	1	Total	C	H	0
			32	10	22	
5	F	1	Total	C	H	0
			32	10	22	
5	F	1	Total	C	H	0
			32	10	22	
5	F	1	Total	C	H	0
			32	10	22	
5	J	1	Total	C	H	0
			32	10	22	
5	J	1	Total	C	H	0
			32	10	22	
5	J	1	Total	C	H	0
			32	10	22	
5	J	1	Total	C	H	0
			32	10	22	
5	E	1	Total	C	H	0
			32	10	22	
5	I	1	Total	C	H	0
			32	10	22	
5	G	1	Total	C	H	0
			32	10	22	
5	K	1	Total	C	H	0
			32	10	22	

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



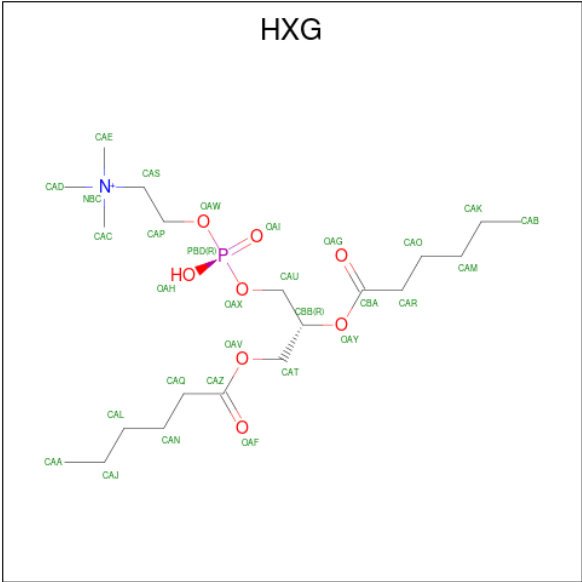
Mol	Chain	Residues	Atoms						AltConf
6	B	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	B	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	B	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	C	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	C	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	C	1	Total 106	C 32	H 64	N 1	O 8	P 1	0
6	F	1	Total 106	C 32	H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



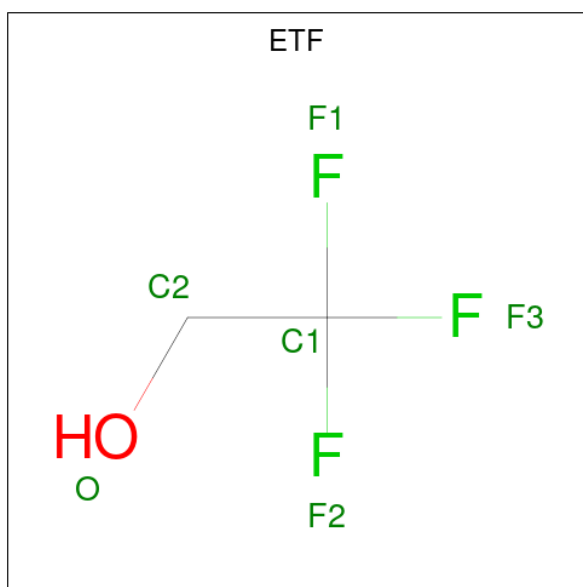
Mol	Chain	Residues	Atoms						AltConf
7	B	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	B	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	C	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	C	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	F	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	F	1	Total	C	H	N	O	P	0
			94	28	56	1	8	1	
7	J	1	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 70	C 20	H 40	N 1	O 8	P 1	0
8	C	1	Total 70	C 20	H 40	N 1	O 8	P 1	0
8	G	1	Total 70	C 20	H 40	N 1	O 8	P 1	0
8	G	1	Total 70	C 20	H 40	N 1	O 8	P 1	0
8	K	1	Total 70	C 20	H 40	N 1	O 8	P 1	0
8	K	1	Total 70	C 20	H 40	N 1	O 8	P 1	0

- 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	

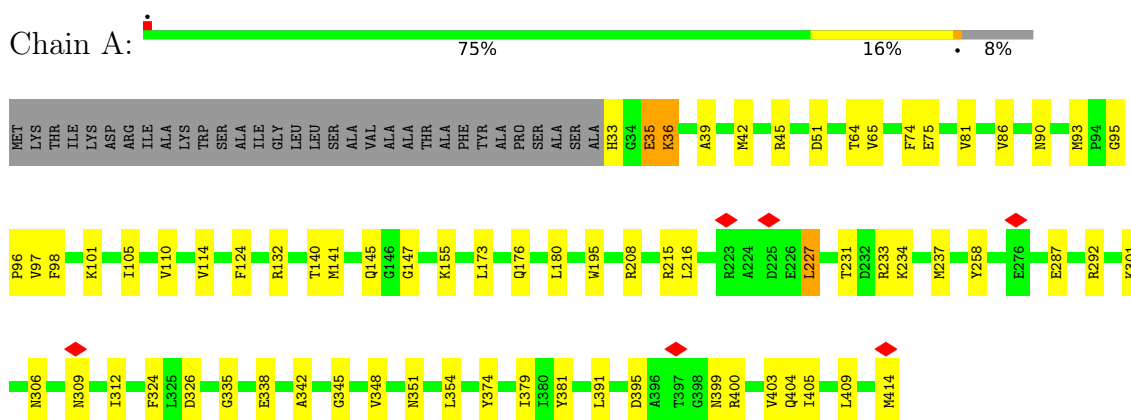
- Molecule 10 is water.

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

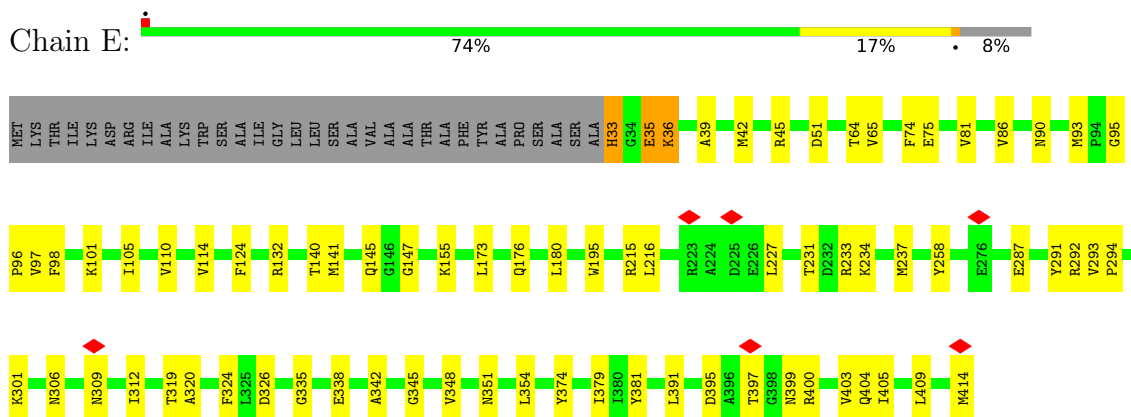
3 Residue-property plots [i](#)

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.

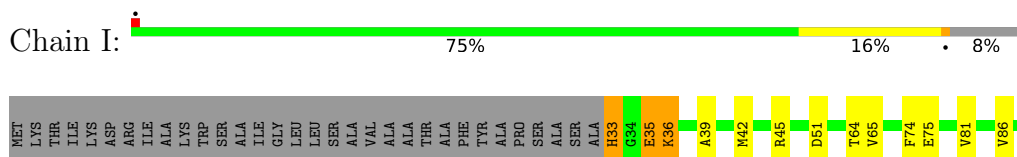
- Molecule 1: Particulate methane monooxygenase alpha subunit

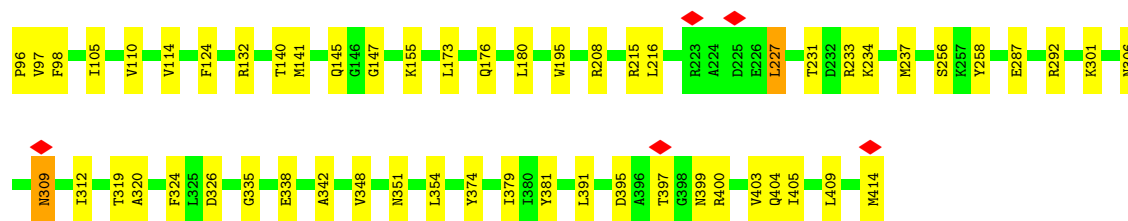


- Molecule 1: Particulate methane monooxygenase alpha subunit

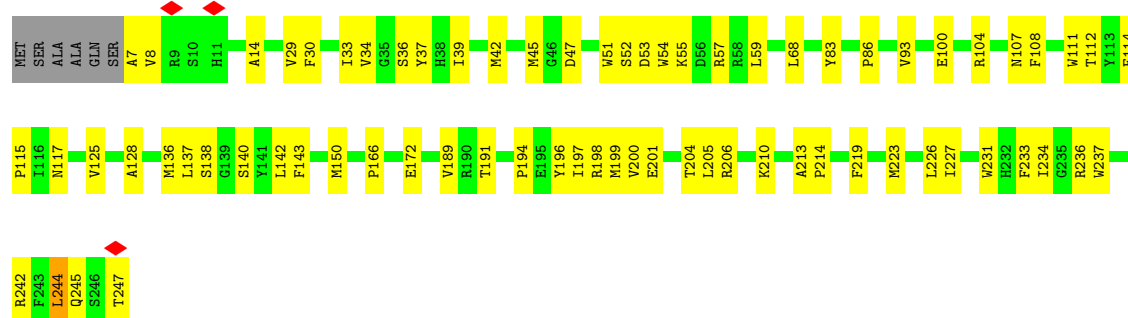


- Molecule 1: Particulate methane monooxygenase alpha subunit

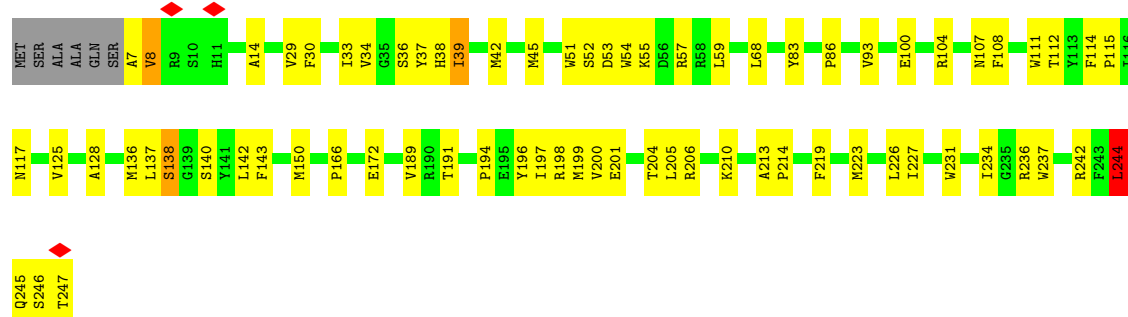




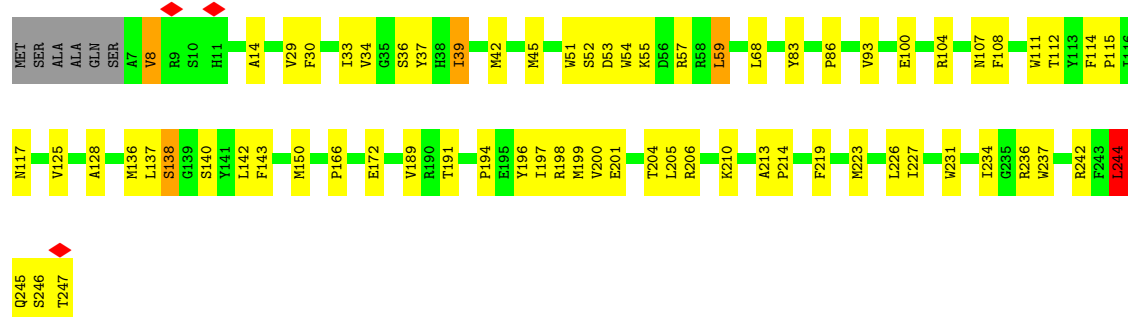
• Molecule 2: Particulate methane monooxygenase beta subunit



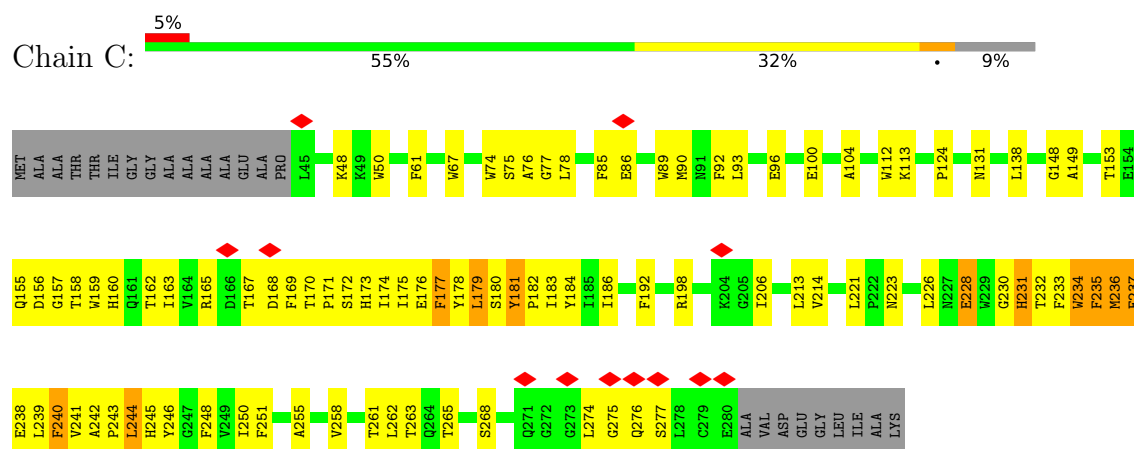
• Molecule 2: Particulate methane monooxygenase beta subunit



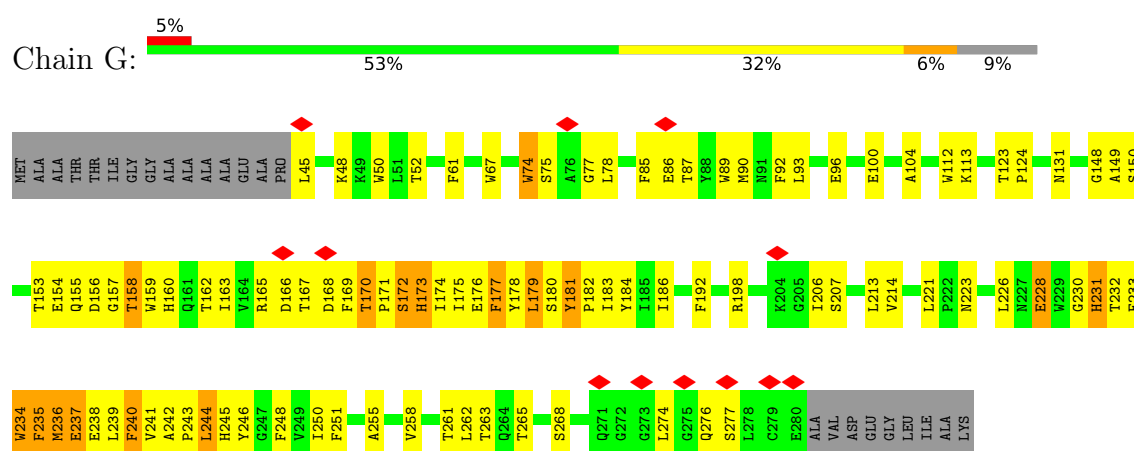
• Molecule 2: Particulate methane monooxygenase beta subunit



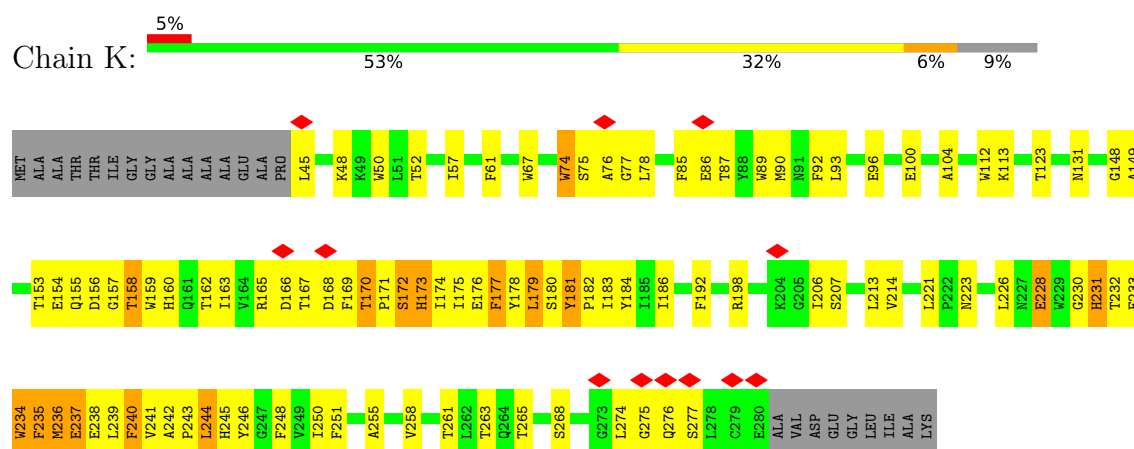
- Molecule 3: Ammonia monooxygenase/methane monooxygenase, subunit C family protein



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



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



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 ($e^-/\text{\AA}^2$)	53.56	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.319	Depositor
Minimum map value	-0.687	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.033	Depositor
Recommended contour level	0.197	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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.41	0/2051	0.65	2/2810 (0.1%)
3	G	0.41	0/2051	0.65	2/2810 (0.1%)
3	K	0.41	0/2051	0.65	2/2810 (0.1%)
All	All	0.29	0/21609	0.49	6/29505 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	C	231	HIS	N-CA-C	-5.72	104.45	112.12
3	G	231	HIS	N-CA-C	-5.72	104.46	112.12
3	K	231	HIS	N-CA-C	-5.71	104.47	112.12
3	K	177	PHE	CB-CA-C	-5.68	101.36	110.79
3	C	177	PHE	CB-CA-C	-5.66	101.40	110.79
3	G	177	PHE	CB-CA-C	-5.66	101.40	110.79

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	3017	0	2980	52	0
1	E	3017	0	2980	53	0
1	I	3017	0	2980	51	0
2	B	1977	0	1936	110	0
2	F	1977	0	1936	111	0
2	J	1977	0	1936	110	0
3	C	1972	0	1904	149	0
3	G	1972	0	1904	154	0
3	K	1972	0	1904	151	0
4	A	2	0	0	0	0
4	C	1	0	0	0	0
4	E	2	0	0	0	0
4	G	1	0	0	0	0
4	I	2	0	0	0	0
4	K	1	0	0	0	0
5	A	10	22	22	0	0
5	B	40	88	88	9	0
5	C	10	22	22	1	0
5	E	10	22	22	0	0
5	F	40	88	88	7	0
5	G	10	22	22	1	0
5	I	10	22	22	0	0
5	J	40	88	88	8	0
5	K	10	22	22	1	0
6	B	126	192	192	10	0
6	C	126	192	192	11	0
6	F	126	192	192	9	0
6	G	126	192	192	13	0
6	J	126	192	192	10	0
6	K	126	192	192	12	0
7	B	76	112	112	37	0
7	C	76	112	112	21	0
7	F	76	112	112	38	0
7	G	76	112	112	22	0
7	J	76	112	112	36	0
7	K	76	112	112	21	0
8	C	60	80	80	29	0
8	G	60	80	80	29	0
8	K	60	80	80	31	0
9	C	6	3	2	0	0
9	G	6	3	2	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	K	6	3	2	0	0
10	A	70	0	0	4	0
10	B	35	0	0	2	0
10	C	13	0	0	1	0
10	E	70	0	0	4	0
10	F	34	0	0	5	0
10	G	13	0	0	2	0
10	I	69	0	0	3	0
10	J	36	0	0	2	0
10	K	12	0	0	2	0
All	All	22849	2469	22926	901	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 20.

All (901) 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:306:HXG:H41	1.81	1.16
3:K:67:TRP:CD1	8:K:306:HXG:H41	1.81	1.16
2:J:206:ARG:HG3	3:G:236:MET:HE2	1.21	1.14
2:B:206:ARG:HG3	3:K:236:MET:HE2	1.25	1.11
6:K:302:PLC:H73	8:K:303:HXG:H36	1.33	1.11
3:K:171:PRO:HA	3:K:174:ILE:HD12	1.30	1.10
2:B:244:LEU:HD21	7:B:307:P1O:C8	1.82	1.09
3:G:171:PRO:HA	3:G:174:ILE:HD12	1.30	1.08
6:G:302:PLC:H73	8:G:303:HXG:H36	1.33	1.08
3:C:236:MET:HE2	2:F:206:ARG:HG3	1.22	1.08
2:B:244:LEU:HD21	7:B:307:P1O:H18	1.34	1.07
2:B:244:LEU:CD2	7:B:307:P1O:H18	1.83	1.07
3:C:67:TRP:HA	8:C:306:HXG:H39	1.07	1.07
2:F:244:LEU:HD21	7:F:309:P1O:C8	1.86	1.06
2:F:244:LEU:HD21	7:F:309:P1O:H18	1.37	1.06
3:K:67:TRP:HA	8:K:306:HXG:H39	1.07	1.05
3:G:67:TRP:HA	8:G:306:HXG:H39	1.07	1.05
3:C:171:PRO:HA	3:C:174:ILE:HD12	1.30	1.05
6:C:302:PLC:H73	8:C:303:HXG:H36	1.33	1.04
2:F:244:LEU:CD2	7:F:309:P1O:H18	1.87	1.04
2:B:244:LEU:HD23	7:B:307:P1O:C5	1.87	1.03
2:F:138:SER:HA	7:F:304:P1O:O8	1.60	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:244:LEU:HD23	7:J:301:P1O:C5	1.90	1.02
3:K:67:TRP:CD1	8:K:306:HXG:CAC	2.43	1.01
3:G:67:TRP:CD1	8:G:306:HXG:CAC	2.43	1.01
3:C:67:TRP:CA	8:C:306:HXG:H39	1.90	1.00
3:C:67:TRP:HA	8:C:306:HXG:CAC	1.91	1.00
3:G:67:TRP:CA	8:G:306:HXG:H39	1.90	1.00
2:B:244:LEU:HD23	7:B:307:P1O:H12	1.43	1.00
3:C:67:TRP:CD1	8:C:306:HXG:CAC	2.43	1.00
2:F:244:LEU:HD23	7:F:309:P1O:C5	1.90	1.00
2:J:244:LEU:HD21	7:J:301:P1O:C8	1.91	1.00
3:G:67:TRP:HA	8:G:306:HXG:CAC	1.91	1.00
2:J:138:SER:HA	7:J:303:P1O:O8	1.60	0.99
2:B:138:SER:HA	7:B:302:P1O:O8	1.60	0.99
3:K:67:TRP:CA	8:K:306:HXG:H39	1.90	0.99
2:J:244:LEU:CD2	7:J:301:P1O:H18	1.91	0.99
2:J:244:LEU:HD23	7:J:301:P1O:H12	1.45	0.99
3:K:67:TRP:HD1	8:K:306:HXG:H41	1.18	0.99
3:K:67:TRP:HA	8:K:306:HXG:CAC	1.91	0.98
2:J:244:LEU:HD21	7:J:301:P1O:H18	1.41	0.97
3:C:50:TRP:CE3	7:C:308:P1O:H39	2.03	0.94
3:K:50:TRP:CE3	7:K:308:P1O:H39	2.03	0.93
3:K:236:MET:HG2	8:K:303:HXG:CAE	1.98	0.93
3:C:236:MET:HG2	8:C:303:HXG:CAE	1.98	0.93
3:G:67:TRP:HD1	8:G:306:HXG:H41	1.18	0.93
3:G:236:MET:HG2	8:G:303:HXG:CAE	1.98	0.93
3:G:50:TRP:CE3	7:G:308:P1O:H39	2.03	0.93
2:F:244:LEU:HD23	7:F:309:P1O:H12	1.47	0.92
3:C:67:TRP:HD1	8:C:306:HXG:H41	1.18	0.92
2:B:107:ASN:ND2	3:C:155:GLN:HB2	1.84	0.92
10:C:404:HOH:O	2:F:150:MET:HE1	1.70	0.91
2:J:150:MET:HE1	10:G:405:HOH:O	1.71	0.91
2:F:107:ASN:ND2	3:G:155:GLN:HB2	1.86	0.91
2:B:244:LEU:HD11	7:B:307:P1O:H42	1.54	0.90
2:B:150:MET:HE1	10:K:405:HOH:O	1.71	0.88
2:F:244:LEU:HD11	7:F:309:P1O:H42	1.54	0.88
2:J:107:ASN:ND2	3:K:155:GLN:HB2	1.89	0.88
3:K:89:TRP:CE2	3:K:171:PRO:HB3	2.09	0.88
8:K:303:HXG:H9	5:K:304:D10:H42	1.56	0.88
3:C:89:TRP:CE2	3:C:171:PRO:HB3	2.09	0.88
8:C:303:HXG:H9	5:C:304:D10:H42	1.56	0.88
3:G:112:TRP:CH2	7:G:309:P1O:H2	2.10	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:206:ARG:CG	3:G:236:MET:HE2	2.04	0.87
3:C:112:TRP:CH2	7:C:309:P1O:H2	2.10	0.87
1:I:36:LYS:HG2	1:I:374:TYR:HA	1.56	0.87
3:G:89:TRP:CE2	3:G:171:PRO:HB3	2.09	0.86
1:E:36:LYS:HG2	1:E:374:TYR:HA	1.55	0.86
8:G:306:HXG:H37	8:G:306:HXG:H26	1.58	0.86
1:A:36:LYS:HG2	1:A:374:TYR:HA	1.55	0.86
8:G:303:HXG:H9	5:G:304:D10:H42	1.56	0.86
3:K:112:TRP:CH2	7:K:309:P1O:H2	2.10	0.86
3:C:232:THR:O	3:C:236:MET:HE1	1.77	0.85
3:C:67:TRP:HB2	8:C:306:HXG:H40	1.57	0.85
3:K:67:TRP:HD1	8:K:306:HXG:CAC	1.86	0.85
3:G:67:TRP:HB2	8:G:306:HXG:H40	1.57	0.85
3:K:232:THR:O	3:K:236:MET:HE1	1.77	0.85
8:C:306:HXG:H37	8:C:306:HXG:H26	1.58	0.85
2:B:244:LEU:HD11	7:B:307:P1O:C22	2.08	0.84
2:J:244:LEU:HD11	7:J:301:P1O:H42	1.58	0.84
3:G:232:THR:O	3:G:236:MET:HE1	1.77	0.84
8:K:306:HXG:H37	8:K:306:HXG:H26	1.58	0.83
3:K:67:TRP:HB2	8:K:306:HXG:H40	1.57	0.83
3:G:67:TRP:HD1	8:G:306:HXG:CAC	1.86	0.83
7:F:304:P1O:O8	7:F:304:P1O:H20	1.79	0.83
7:J:303:P1O:O8	7:J:303:P1O:H20	1.79	0.82
7:B:302:P1O:O8	7:B:302:P1O:H20	1.79	0.82
2:B:206:ARG:CG	3:K:236:MET:HE2	2.10	0.82
2:F:237:TRP:HE1	5:F:307:D10:H32	1.44	0.81
2:F:244:LEU:HD11	7:F:309:P1O:C22	2.10	0.81
2:B:237:TRP:HE1	5:B:305:D10:H32	1.44	0.80
2:J:237:TRP:HE1	5:J:306:D10:H32	1.44	0.80
3:G:171:PRO:CA	3:G:174:ILE:HD12	2.12	0.79
3:G:171:PRO:HA	3:G:174:ILE:CD1	2.12	0.79
1:A:81:VAL:HG13	1:A:147:GLY:HA3	1.65	0.79
3:G:67:TRP:CB	8:G:306:HXG:H40	2.13	0.79
3:C:67:TRP:CB	8:C:306:HXG:H40	2.13	0.79
3:K:171:PRO:HA	3:K:174:ILE:CD1	2.13	0.79
1:E:81:VAL:HG13	1:E:147:GLY:HA3	1.65	0.79
1:I:81:VAL:HG13	1:I:147:GLY:HA3	1.65	0.79
3:C:236:MET:HE2	2:F:206:ARG:CG	2.07	0.78
7:J:303:P1O:H48	7:J:303:P1O:H38	1.66	0.78
2:B:138:SER:CA	7:B:302:P1O:O8	2.32	0.77
3:G:112:TRP:CH2	7:G:309:P1O:C1	2.67	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:67:TRP:CB	8:K:306:HXG:H40	2.13	0.77
3:K:112:TRP:CH2	7:K:309:P1O:C1	2.67	0.77
3:C:67:TRP:HD1	8:C:306:HXG:CAC	1.86	0.77
3:G:236:MET:HE3	3:G:236:MET:HA	1.66	0.77
3:C:67:TRP:CB	8:C:306:HXG:CAC	2.63	0.77
3:G:67:TRP:CB	8:G:306:HXG:CAC	2.63	0.77
2:B:199:MET:SD	6:B:308:PLC:H2	2.24	0.77
2:F:138:SER:CA	7:F:304:P1O:O8	2.32	0.77
2:J:244:LEU:HD11	7:J:301:P1O:C22	2.14	0.77
7:B:302:P1O:H38	7:B:302:P1O:H48	1.66	0.76
3:K:67:TRP:CB	8:K:306:HXG:CAC	2.63	0.76
7:F:304:P1O:H48	7:F:304:P1O:H38	1.66	0.76
2:J:138:SER:CA	7:J:303:P1O:O8	2.32	0.76
3:K:171:PRO:CA	3:K:174:ILE:HD12	2.12	0.76
3:C:112:TRP:CH2	7:C:309:P1O:C1	2.67	0.76
3:K:112:TRP:CZ2	7:K:309:P1O:C1	2.69	0.76
3:C:48:LYS:HD3	3:C:50:TRP:HD1	1.50	0.76
3:C:112:TRP:CZ2	7:C:309:P1O:C1	2.69	0.76
3:K:48:LYS:HD3	3:K:50:TRP:HD1	1.51	0.76
2:F:7:ALA:CB	10:F:418:HOH:O	2.33	0.76
2:J:199:MET:SD	6:J:308:PLC:H2	2.25	0.76
3:K:236:MET:HA	3:K:236:MET:HE3	1.66	0.75
3:C:236:MET:HE3	3:C:236:MET:HA	1.66	0.75
3:G:112:TRP:CZ2	7:G:309:P1O:C1	2.69	0.75
3:G:148:GLY:HA2	3:G:180:SER:HB2	1.69	0.75
3:C:171:PRO:CA	3:C:174:ILE:HD12	2.12	0.75
3:G:48:LYS:HD3	3:G:50:TRP:HD1	1.50	0.75
2:F:138:SER:HB2	7:F:304:P1O:O8	1.87	0.74
3:C:171:PRO:HA	3:C:174:ILE:CD1	2.12	0.74
3:C:148:GLY:HA2	3:C:180:SER:HB2	1.69	0.74
7:F:309:P1O:O6	7:F:309:P1O:H43	1.88	0.74
3:C:85:PHE:HE1	3:C:171:PRO:HD3	1.53	0.74
3:K:85:PHE:HE1	3:K:171:PRO:HD3	1.53	0.74
3:G:85:PHE:HE1	3:G:171:PRO:HD3	1.53	0.74
2:B:138:SER:HB2	7:B:302:P1O:O8	1.87	0.73
7:B:307:P1O:O6	7:B:307:P1O:H43	1.88	0.73
3:K:148:GLY:HA2	3:K:180:SER:HB2	1.69	0.73
3:K:177:PHE:O	3:K:182:PRO:HD3	1.88	0.73
2:F:199:MET:SD	6:F:301:PLC:H2	2.28	0.73
2:J:138:SER:HB2	7:J:303:P1O:O8	1.87	0.73
3:G:177:PHE:O	3:G:182:PRO:HD3	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:177:PHE:O	3:C:182:PRO:HD3	1.88	0.72
7:J:301:P1O:O6	7:J:301:P1O:H43	1.88	0.72
3:K:236:MET:HA	3:K:236:MET:CE	2.20	0.72
3:G:236:MET:HG2	8:G:303:HXG:H37	1.70	0.72
2:B:112:THR:HG21	3:C:162:THR:HG21	1.72	0.71
3:C:236:MET:HA	3:C:236:MET:CE	2.20	0.71
3:C:236:MET:HG2	8:C:303:HXG:H37	1.70	0.71
3:G:236:MET:HA	3:G:236:MET:CE	2.20	0.71
3:K:236:MET:HG2	8:K:303:HXG:H37	1.71	0.71
2:B:244:LEU:HD21	7:B:307:P1O:H17	1.73	0.70
2:B:244:LEU:CD2	7:B:307:P1O:H12	2.21	0.70
2:J:138:SER:CB	7:J:303:P1O:O8	2.40	0.70
1:I:215:ARG:HG2	1:I:227:LEU:HD22	1.73	0.70
2:F:112:THR:HG21	3:G:162:THR:HG21	1.73	0.69
1:E:326:ASP:OD2	1:E:351:ASN:ND2	2.25	0.69
2:F:138:SER:CB	7:F:304:P1O:O8	2.40	0.69
1:A:93:MET:HE1	1:A:98:PHE:HB2	1.75	0.69
2:B:138:SER:CB	7:B:302:P1O:O8	2.40	0.69
1:E:93:MET:HE1	1:E:98:PHE:HB2	1.75	0.69
1:E:215:ARG:HG2	1:E:227:LEU:HD22	1.73	0.69
1:A:75:GLU:OE1	1:A:404:GLN:NE2	2.26	0.69
2:B:244:LEU:HD23	7:B:307:P1O:H18	1.74	0.69
3:C:112:TRP:CZ2	7:C:309:P1O:H2	2.28	0.69
3:G:112:TRP:CZ2	7:G:309:P1O:H2	2.28	0.69
1:A:215:ARG:HG2	1:A:227:LEU:HD22	1.73	0.69
1:I:93:MET:HE1	1:I:98:PHE:HB2	1.75	0.69
2:B:142:LEU:HD22	7:J:301:P1O:H48	1.76	0.68
2:J:112:THR:HG21	3:K:162:THR:HG21	1.74	0.68
1:I:326:ASP:OD2	1:I:351:ASN:ND2	2.25	0.68
3:C:67:TRP:CG	8:C:306:HXG:CAC	2.76	0.68
2:F:7:ALA:HB3	10:F:418:HOH:O	1.92	0.68
3:K:89:TRP:CD2	3:K:171:PRO:HB3	2.28	0.68
2:B:114:PHE:HE2	3:C:162:THR:HG22	1.57	0.68
3:C:67:TRP:CA	8:C:306:HXG:CAC	2.61	0.68
1:I:75:GLU:OE1	1:I:404:GLN:NE2	2.26	0.68
3:G:67:TRP:CG	8:G:306:HXG:CAC	2.76	0.68
2:J:108:PHE:CE2	3:K:158:THR:HG22	2.29	0.68
3:G:89:TRP:CD2	3:G:171:PRO:HB3	2.28	0.68
1:A:326:ASP:OD2	1:A:351:ASN:ND2	2.25	0.68
2:J:244:LEU:CD2	7:J:301:P1O:H12	2.23	0.68
3:C:89:TRP:CD2	3:C:171:PRO:HB3	2.28	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:112:TRP:CZ2	7:K:309:P1O:H2	2.28	0.68
3:K:67:TRP:CG	8:K:306:HXG:CAC	2.76	0.68
7:B:307:P1O:H48	2:F:142:LEU:HD22	1.75	0.67
1:E:75:GLU:OE1	1:E:404:GLN:NE2	2.25	0.67
2:B:108:PHE:CE2	3:C:158:THR:HG22	2.28	0.67
2:B:237:TRP:HE1	5:B:305:D10:C3	2.06	0.67
3:K:67:TRP:CA	8:K:306:HXG:CAC	2.61	0.67
2:F:244:LEU:HD23	7:F:309:P1O:H18	1.77	0.66
7:F:309:P1O:H48	2:J:142:LEU:HD22	1.75	0.66
1:A:176:GLN:NE2	10:A:601:HOH:O	2.29	0.66
2:F:108:PHE:CE2	3:G:158:THR:HG22	2.29	0.66
2:J:237:TRP:HE1	5:J:306:D10:C3	2.07	0.66
2:F:244:LEU:HD21	7:F:309:P1O:H17	1.75	0.66
3:C:242:ALA:HB3	3:C:244:LEU:HD23	1.78	0.66
3:G:242:ALA:HB3	3:G:244:LEU:HD23	1.78	0.66
2:F:237:TRP:HE1	5:F:307:D10:C3	2.07	0.66
6:F:301:PLC:H32	6:F:301:PLC:H1'2	1.78	0.66
3:K:242:ALA:HB3	3:K:244:LEU:HD23	1.78	0.65
2:F:114:PHE:HE2	3:G:162:THR:HG22	1.59	0.65
8:C:306:HXG:H37	8:C:306:HXG:CAU	2.26	0.65
8:G:306:HXG:H37	8:G:306:HXG:CAU	2.26	0.65
2:J:114:PHE:HE2	3:K:162:THR:HG22	1.60	0.65
6:J:308:PLC:H1'2	6:J:308:PLC:H32	1.78	0.65
8:K:306:HXG:H37	8:K:306:HXG:CAU	2.26	0.65
2:B:45:MET:HE3	2:B:68:LEU:HD11	1.77	0.65
2:F:45:MET:HE3	2:F:68:LEU:HD11	1.78	0.65
2:F:237:TRP:NE1	5:F:307:D10:H32	2.12	0.65
2:J:45:MET:HE3	2:J:68:LEU:HD11	1.77	0.65
6:B:308:PLC:H32	6:B:308:PLC:H1'2	1.78	0.64
7:J:301:P1O:H12	7:J:301:P1O:O3	1.98	0.64
7:B:307:P1O:H12	7:B:307:P1O:O3	1.98	0.64
3:C:261:THR:O	3:C:265:THR:HG23	1.98	0.64
7:F:309:P1O:H12	7:F:309:P1O:O3	1.98	0.64
3:K:261:THR:O	3:K:265:THR:HG23	1.98	0.64
2:B:237:TRP:NE1	5:B:305:D10:H32	2.12	0.64
2:J:237:TRP:NE1	5:J:306:D10:H32	2.12	0.63
3:G:234:TRP:HZ2	8:G:303:HXG:H18	1.63	0.63
3:K:234:TRP:HZ2	8:K:303:HXG:H18	1.64	0.63
2:J:244:LEU:HD21	7:J:301:P1O:H17	1.81	0.63
2:F:244:LEU:CD2	7:F:309:P1O:H12	2.24	0.63
2:B:226:LEU:HD21	3:K:251:PHE:HZ	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:234:TRP:HZ2	8:C:303:HXG:H18	1.63	0.63
1:E:176:GLN:NE2	10:E:601:HOH:O	2.32	0.62
7:G:308:P1O:H18	7:G:308:P1O:H24	1.82	0.62
3:G:67:TRP:CA	8:G:306:HXG:CAC	2.61	0.62
7:C:309:P1O:O8	7:C:309:P1O:H7	1.99	0.62
3:G:261:THR:O	3:G:265:THR:HG23	1.98	0.62
7:G:309:P1O:O8	7:G:309:P1O:H7	1.99	0.62
2:B:138:SER:O	2:B:138:SER:OG	2.17	0.62
7:C:308:P1O:H18	7:C:308:P1O:H24	1.82	0.62
3:G:192:PHE:HB2	3:G:214:VAL:HG21	1.82	0.62
2:F:237:TRP:NE1	5:F:307:D10:H51	2.15	0.61
2:J:226:LEU:HD21	3:G:251:PHE:HZ	1.64	0.61
7:K:309:P1O:O8	7:K:309:P1O:H7	1.99	0.61
3:C:192:PHE:HB2	3:C:214:VAL:HG21	1.82	0.61
3:G:50:TRP:CZ3	7:G:308:P1O:H39	2.35	0.61
7:K:308:P1O:H18	7:K:308:P1O:H24	1.82	0.61
2:J:237:TRP:NE1	5:J:306:D10:H51	2.15	0.61
3:K:192:PHE:HB2	3:K:214:VAL:HG21	1.82	0.61
3:C:50:TRP:CZ3	7:C:308:P1O:H39	2.35	0.60
2:B:237:TRP:NE1	5:B:305:D10:H51	2.15	0.60
2:J:8:VAL:HG11	2:J:14:ALA:HB2	1.83	0.60
3:K:50:TRP:CZ3	7:K:308:P1O:H39	2.35	0.60
2:F:138:SER:O	2:F:138:SER:OG	2.17	0.60
2:B:8:VAL:HG11	2:B:14:ALA:HB2	1.83	0.60
7:C:308:P1O:O2	7:C:308:P1O:H4	2.02	0.60
2:F:8:VAL:HG11	2:F:14:ALA:HB2	1.83	0.60
3:C:251:PHE:HZ	2:F:226:LEU:HD21	1.66	0.60
3:G:236:MET:H	8:G:303:HXG:H37	1.67	0.59
7:G:308:P1O:H4	7:G:308:P1O:O2	2.02	0.59
7:K:308:P1O:H4	7:K:308:P1O:O2	2.02	0.59
7:J:301:P1O:H15	7:J:301:P1O:H39	1.83	0.59
1:I:176:GLN:NE2	10:I:603:HOH:O	2.36	0.59
2:B:114:PHE:CE2	3:C:162:THR:HG22	2.38	0.59
7:B:307:P1O:H15	7:B:307:P1O:H39	1.83	0.59
3:C:112:TRP:CZ2	7:C:309:P1O:H7	2.38	0.59
3:K:112:TRP:CZ2	7:K:309:P1O:H7	2.38	0.59
7:F:309:P1O:H15	7:F:309:P1O:H39	1.83	0.59
2:J:242:ARG:HH21	7:J:301:P1O:C4	2.16	0.59
2:J:244:LEU:HD23	7:J:301:P1O:H18	1.82	0.59
2:J:199:MET:SD	6:J:308:PLC:C2	2.91	0.58
3:G:112:TRP:CZ2	7:G:309:P1O:H7	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:291:TYR:HA	10:E:650:HOH:O	2.03	0.58
3:C:236:MET:H	8:C:303:HXG:H37	1.67	0.58
2:J:242:ARG:HH21	7:J:301:P1O:H9	1.68	0.58
2:B:111:TRP:CH2	3:C:159:TRP:HE3	2.22	0.58
7:F:309:P1O:H17	7:F:309:P1O:H39	1.86	0.58
7:B:307:P1O:H17	7:B:307:P1O:H39	1.85	0.58
7:C:308:P1O:O2	7:C:308:P1O:H6	2.04	0.58
2:J:206:ARG:HB2	3:G:236:MET:HE1	1.86	0.58
3:K:236:MET:H	8:K:303:HXG:H37	1.67	0.58
2:B:242:ARG:HH21	7:B:307:P1O:C4	2.17	0.58
3:C:77:GLY:O	3:C:171:PRO:HG2	2.05	0.57
2:J:206:ARG:HB2	3:G:236:MET:CE	2.35	0.57
2:B:199:MET:SD	6:B:308:PLC:C2	2.92	0.57
3:C:112:TRP:CH2	7:C:309:P1O:H1	2.40	0.57
7:G:308:P1O:O2	7:G:308:P1O:H6	2.04	0.57
7:K:308:P1O:O2	7:K:308:P1O:H6	2.04	0.57
7:J:301:P1O:H17	7:J:301:P1O:H39	1.85	0.57
3:C:92:PHE:HA	6:C:307:PLC:OB	2.05	0.57
3:K:77:GLY:O	3:K:171:PRO:HG2	2.05	0.57
2:B:244:LEU:CD2	7:B:307:P1O:C8	2.57	0.57
3:C:100:GLU:OE2	3:C:179:LEU:HD12	2.05	0.57
1:E:231:THR:HG22	1:E:234:LYS:HE2	1.87	0.57
1:A:231:THR:HG22	1:A:234:LYS:HE2	1.87	0.56
3:G:92:PHE:HA	6:G:307:PLC:OB	2.05	0.56
1:A:309:ASN:OD1	1:A:309:ASN:N	2.39	0.56
3:C:112:TRP:CZ2	7:C:309:P1O:C3	2.89	0.56
3:G:77:GLY:O	3:G:171:PRO:HG2	2.05	0.56
3:K:92:PHE:HA	6:K:307:PLC:OB	2.05	0.56
2:F:114:PHE:CE2	3:G:162:THR:HG22	2.40	0.56
2:J:114:PHE:CE2	3:K:162:THR:HG22	2.41	0.56
1:I:231:THR:HG22	1:I:234:LYS:HE2	1.87	0.56
3:K:112:TRP:CZ2	7:K:309:P1O:C3	2.89	0.56
3:G:100:GLU:OE2	3:G:179:LEU:HD12	2.05	0.56
3:G:148:GLY:HA2	3:G:180:SER:CB	2.36	0.56
2:J:244:LEU:HD23	7:J:301:P1O:H13	1.86	0.56
3:K:148:GLY:HA2	3:K:180:SER:CB	2.36	0.56
2:B:213:ALA:HB3	6:B:301:PLC:H11	1.88	0.56
1:E:309:ASN:N	1:E:309:ASN:OD1	2.39	0.56
3:G:112:TRP:CZ2	7:G:309:P1O:C3	2.89	0.56
3:G:235:PHE:HA	8:G:303:HXG:H41	1.88	0.56
3:K:112:TRP:CH2	7:K:309:P1O:H1	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:236:MET:HE3	3:C:236:MET:CA	2.36	0.56
3:K:112:TRP:HB2	7:K:309:P1O:H40	1.88	0.56
2:B:244:LEU:HD11	7:B:307:P1O:H43	1.86	0.56
3:C:112:TRP:HB2	7:C:309:P1O:H40	1.88	0.56
2:J:244:LEU:CD2	7:J:301:P1O:C8	2.66	0.56
3:G:112:TRP:HB2	7:G:309:P1O:H40	1.88	0.55
2:J:138:SER:O	2:J:138:SER:OG	2.17	0.55
3:K:100:GLU:OE2	3:K:179:LEU:HD12	2.05	0.55
6:F:301:PLC:H63	6:F:301:PLC:O4P	2.06	0.55
1:E:45:ARG:HB3	1:E:74:PHE:CD2	2.42	0.55
3:G:85:PHE:CE1	3:G:171:PRO:HD3	2.40	0.55
6:B:308:PLC:H63	6:B:308:PLC:O4P	2.06	0.55
3:G:67:TRP:CD1	8:G:306:HXG:H40	2.39	0.55
2:F:111:TRP:CH2	3:G:159:TRP:HE3	2.24	0.55
2:F:199:MET:SD	6:F:301:PLC:C2	2.95	0.55
1:A:45:ARG:HB3	1:A:74:PHE:CD2	2.42	0.55
3:K:67:TRP:CD1	8:K:306:HXG:H40	2.39	0.55
3:K:235:PHE:HA	8:K:303:HXG:H41	1.88	0.55
2:F:242:ARG:HH21	7:F:309:P1O:C4	2.20	0.55
2:F:244:LEU:HD11	7:F:309:P1O:H43	1.89	0.54
6:J:308:PLC:H63	6:J:308:PLC:O4P	2.06	0.54
1:I:309:ASN:OD1	1:I:309:ASN:N	2.39	0.54
3:G:236:MET:HE3	3:G:236:MET:CA	2.36	0.54
2:B:242:ARG:HH21	7:B:307:P1O:H9	1.71	0.54
2:B:244:LEU:HD21	7:B:307:P1O:H15	1.88	0.54
1:I:45:ARG:HB3	1:I:74:PHE:CD2	2.42	0.54
3:G:159:TRP:HD1	3:G:160:HIS:HD1	1.56	0.54
3:C:235:PHE:HA	8:C:303:HXG:H41	1.88	0.54
2:B:114:PHE:HE2	3:C:162:THR:CG2	2.20	0.54
2:B:244:LEU:HD21	7:B:307:P1O:C6	2.37	0.54
3:C:235:PHE:CZ	3:C:243:PRO:HD2	2.43	0.54
2:J:206:ARG:O	3:G:232:THR:HG22	2.08	0.54
3:G:112:TRP:CH2	7:G:309:P1O:H1	2.40	0.54
3:K:159:TRP:HD1	3:K:160:HIS:HD1	1.56	0.54
2:J:213:ALA:HB3	6:J:302:PLC:H11	1.88	0.54
3:C:159:TRP:HD1	3:C:160:HIS:HD1	1.56	0.54
3:K:50:TRP:CD2	7:K:308:P1O:H16	2.43	0.54
3:K:155:GLN:O	3:K:158:THR:HG23	2.08	0.54
3:C:50:TRP:CD2	7:C:308:P1O:H16	2.43	0.54
2:F:213:ALA:HB3	6:F:303:PLC:H11	1.88	0.54
2:F:244:LEU:CD1	7:F:309:P1O:H42	2.35	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:206:ARG:O	3:K:232:THR:HG22	2.08	0.53
3:C:148:GLY:HA2	3:C:180:SER:CB	2.36	0.53
2:F:242:ARG:HH21	7:F:309:P1O:H9	1.73	0.53
2:J:244:LEU:HD21	7:J:301:P1O:H15	1.90	0.53
3:G:50:TRP:CD2	7:G:308:P1O:H16	2.43	0.53
3:K:235:PHE:CZ	3:K:243:PRO:HD2	2.43	0.53
3:C:232:THR:HG22	2:F:206:ARG:O	2.07	0.53
3:C:236:MET:HE1	2:F:206:ARG:HB2	1.90	0.53
2:J:111:TRP:CH2	3:K:159:TRP:HE3	2.26	0.53
1:A:292:ARG:NH1	1:A:414:MET:O	2.42	0.53
3:C:159:TRP:HD1	3:C:160:HIS:ND1	2.06	0.53
2:F:244:LEU:HD21	7:F:309:P1O:H15	1.91	0.53
3:G:235:PHE:CZ	3:G:243:PRO:HD2	2.43	0.53
2:B:111:TRP:CH2	3:C:159:TRP:CE3	2.97	0.53
3:G:153:THR:O	3:G:156:ASP:HB3	2.08	0.53
3:G:246:TYR:O	3:G:250:ILE:HG12	2.09	0.53
3:K:159:TRP:HD1	3:K:160:HIS:ND1	2.06	0.53
3:G:235:PHE:CE2	3:G:243:PRO:HD2	2.44	0.53
3:K:246:TYR:O	3:K:250:ILE:HG12	2.09	0.53
2:F:244:LEU:CD2	7:F:309:P1O:C8	2.61	0.53
2:J:242:ARG:NH2	7:J:301:P1O:C4	2.72	0.53
3:G:89:TRP:HB3	3:G:174:ILE:CD1	2.39	0.53
3:K:236:MET:HE3	3:K:236:MET:CA	2.36	0.53
3:C:246:TYR:O	3:C:250:ILE:HG12	2.09	0.53
3:K:235:PHE:CE2	3:K:243:PRO:HD2	2.44	0.53
2:B:244:LEU:CD1	7:B:307:P1O:H42	2.33	0.52
3:G:155:GLN:O	3:G:158:THR:HG23	2.08	0.52
3:C:236:MET:CE	3:C:236:MET:CA	2.84	0.52
2:F:136:MET:HE1	1:E:233:ARG:HA	1.92	0.52
3:G:159:TRP:HD1	3:G:160:HIS:ND1	2.06	0.52
3:C:153:THR:O	3:C:156:ASP:HB3	2.08	0.52
3:C:183:ILE:HG12	6:C:307:PLC:HEA3	1.92	0.52
3:G:236:MET:N	8:G:303:HGX:H37	2.25	0.52
3:C:155:GLN:O	3:C:158:THR:HG23	2.08	0.52
3:C:236:MET:N	8:C:303:HGX:H37	2.25	0.52
3:G:157:GLY:HA2	3:G:240:PHE:CE1	2.45	0.52
3:C:89:TRP:HB3	3:C:174:ILE:CD1	2.39	0.52
2:J:143:PHE:CD2	7:J:303:P1O:H26	2.45	0.52
3:G:236:MET:CE	3:G:236:MET:CA	2.85	0.52
2:F:114:PHE:HE2	3:G:162:THR:CG2	2.21	0.52
2:F:143:PHE:CD2	7:F:304:P1O:H26	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:89:TRP:HB3	3:K:174:ILE:CD1	2.39	0.52
8:K:303:HXG:H17	8:K:303:HXG:H10	1.92	0.52
2:J:136:MET:HE1	1:I:233:ARG:HA	1.91	0.52
1:E:155:LYS:NZ	1:E:335:GLY:O	2.35	0.52
1:I:39:ALA:HB3	1:I:42:MET:HG3	1.92	0.52
3:G:183:ILE:HG12	6:G:307:PLC:HEA3	1.92	0.52
1:A:233:ARG:HA	2:B:136:MET:HE1	1.91	0.52
3:C:67:TRP:CD1	8:C:306:HXG:H40	2.39	0.52
3:C:235:PHE:CE2	3:C:243:PRO:HD2	2.44	0.52
2:F:244:LEU:HD23	7:F:309:P1O:H13	1.87	0.52
3:K:153:THR:O	3:K:156:ASP:HB3	2.08	0.52
2:B:143:PHE:CD2	7:B:302:P1O:H26	2.45	0.51
6:B:308:PLC:H9'1	6:K:302:PLC:HTA1	1.92	0.51
3:C:243:PRO:C	3:C:245:HIS:H	2.18	0.51
3:C:236:MET:CE	2:F:206:ARG:HB2	2.40	0.51
2:F:111:TRP:CH2	3:G:159:TRP:CE3	2.99	0.51
1:E:39:ALA:HB3	1:E:42:MET:HG3	1.92	0.51
3:G:243:PRO:C	3:G:245:HIS:H	2.18	0.51
6:C:302:PLC:HTA1	6:F:301:PLC:H9'1	1.92	0.51
6:J:308:PLC:H9'1	6:G:302:PLC:HTA1	1.92	0.51
3:G:238:GLU:HG3	10:G:403:HOH:O	2.10	0.51
3:K:236:MET:N	8:K:303:HXG:H37	2.25	0.51
1:A:51:ASP:CG	1:A:400:ARG:HH22	2.18	0.51
3:C:157:GLY:HA2	3:C:240:PHE:CE1	2.45	0.51
3:G:112:TRP:CE2	7:G:309:P1O:C3	2.94	0.51
8:G:303:HXG:H17	8:G:303:HXG:H10	1.92	0.51
3:C:85:PHE:CE1	3:C:171:PRO:HD3	2.40	0.51
2:F:7:ALA:HB2	10:F:418:HOH:O	2.05	0.51
2:J:244:LEU:HD21	7:J:301:P1O:C6	2.41	0.51
3:K:157:GLY:HA2	3:K:240:PHE:CE1	2.45	0.51
1:E:292:ARG:NH1	1:E:414:MET:O	2.42	0.51
7:C:309:P1O:H12	7:C:309:P1O:H16	1.93	0.51
2:F:100:GLU:O	2:F:104:ARG:HG2	2.11	0.51
2:J:114:PHE:HE2	3:K:162:THR:CG2	2.22	0.51
2:J:100:GLU:O	2:J:104:ARG:HG2	2.11	0.51
3:K:112:TRP:CE2	7:K:309:P1O:C3	2.94	0.51
3:K:131:ASN:OD1	3:K:268:SER:OG	2.29	0.51
3:K:183:ILE:HG12	6:K:307:PLC:HEA3	1.92	0.51
3:K:243:PRO:C	3:K:245:HIS:H	2.18	0.51
3:C:112:TRP:CE2	7:C:309:P1O:C3	2.94	0.50
3:K:236:MET:CE	3:K:236:MET:CA	2.84	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:159:TRP:CE3	3:C:172:SER:HB3	2.46	0.50
2:B:206:ARG:HB2	3:K:236:MET:HE1	1.92	0.50
2:J:245:GLN:O	3:K:206:ILE:HD12	2.12	0.50
1:E:51:ASP:CG	1:E:400:ARG:HH22	2.18	0.50
3:K:159:TRP:NE1	3:K:163:ILE:HD11	2.26	0.50
3:C:90:MET:HE2	6:C:305:PLC:H61	1.93	0.50
3:C:93:LEU:HD13	3:C:174:ILE:HG23	1.94	0.50
3:C:170:THR:O	3:C:174:ILE:HG13	2.12	0.50
8:C:303:HXG:H17	8:C:303:HXG:H10	1.92	0.50
1:E:124:PHE:CE2	1:E:140:THR:HG21	2.47	0.50
1:I:155:LYS:NZ	1:I:335:GLY:O	2.35	0.50
3:G:112:TRP:CE2	7:G:309:P1O:H6	2.46	0.50
3:G:170:THR:O	3:G:174:ILE:HG13	2.12	0.50
3:K:192:PHE:HZ	3:K:206:ILE:HG23	1.75	0.50
3:C:112:TRP:CE2	7:C:309:P1O:H6	2.46	0.50
3:C:131:ASN:OD1	3:C:268:SER:OG	2.29	0.50
7:G:309:P1O:H12	7:G:309:P1O:H16	1.93	0.50
3:K:112:TRP:CE2	7:K:309:P1O:H6	2.46	0.50
1:A:39:ALA:HB3	1:A:42:MET:HG3	1.92	0.50
2:B:242:ARG:NH2	7:B:307:P1O:C4	2.75	0.50
2:F:244:LEU:HD21	7:F:309:P1O:C6	2.41	0.50
2:J:246:SER:HA	3:K:206:ILE:CD1	2.42	0.50
1:I:292:ARG:NH1	1:I:414:MET:O	2.42	0.50
3:G:192:PHE:HZ	3:G:206:ILE:HG23	1.75	0.50
3:K:238:GLU:HG3	10:K:403:HOH:O	2.12	0.50
1:A:155:LYS:NZ	1:A:335:GLY:O	2.35	0.50
2:B:30:PHE:O	2:B:34:VAL:HG23	2.12	0.50
2:B:86:PRO:HA	10:B:426:HOH:O	2.11	0.50
2:B:219:PHE:CD1	3:K:226:LEU:HA	2.47	0.50
2:J:30:PHE:O	2:J:34:VAL:HG23	2.12	0.50
1:I:51:ASP:CG	1:I:400:ARG:HH22	2.18	0.50
3:K:170:THR:O	3:K:174:ILE:HG13	2.12	0.50
2:B:206:ARG:HB2	3:K:236:MET:CE	2.41	0.50
3:C:192:PHE:HZ	3:C:206:ILE:HG23	1.75	0.50
3:G:90:MET:HE2	6:G:305:PLC:H61	1.93	0.50
3:G:93:LEU:HD13	3:G:174:ILE:HG23	1.94	0.49
3:K:90:MET:HE2	6:K:305:PLC:H61	1.93	0.49
3:K:159:TRP:CE3	3:K:172:SER:HB3	2.46	0.49
2:B:100:GLU:O	2:B:104:ARG:HG2	2.11	0.49
3:C:159:TRP:NE1	3:C:163:ILE:HD11	2.26	0.49
3:G:159:TRP:NE1	3:G:163:ILE:HD11	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:226:LEU:HD21	3:K:251:PHE:CZ	2.47	0.49
2:F:86:PRO:HA	10:F:423:HOH:O	2.11	0.49
3:K:75:SER:OG	3:K:75:SER:O	2.28	0.49
7:K:309:P1O:H16	7:K:309:P1O:H12	1.93	0.49
2:J:247:THR:HG23	3:K:206:ILE:HD13	1.94	0.49
6:F:301:PLC:OB	6:F:301:PLC:H81	2.12	0.49
3:G:159:TRP:CE3	3:G:172:SER:HB3	2.46	0.49
1:I:124:PHE:CE2	1:I:140:THR:HG21	2.47	0.49
3:G:131:ASN:OD1	3:G:268:SER:OG	2.29	0.49
6:B:308:PLC:OB	6:B:308:PLC:H81	2.12	0.49
1:A:90:ASN:HB3	1:A:141:MET:HG3	1.95	0.49
1:A:124:PHE:CE2	1:A:140:THR:HG21	2.47	0.49
1:I:105:ILE:HD12	1:I:114:VAL:HG21	1.94	0.49
1:A:105:ILE:HD12	1:A:114:VAL:HG21	1.94	0.49
3:C:275:GLY:HA3	3:C:276:GLN:HA	1.50	0.49
6:J:308:PLC:OB	6:J:308:PLC:H81	2.12	0.49
1:E:90:ASN:HB3	1:E:141:MET:HG3	1.95	0.49
6:C:302:PLC:C7	8:C:303:HXG:H36	2.24	0.49
2:J:111:TRP:CH2	3:K:159:TRP:CE3	3.01	0.49
2:F:30:PHE:O	2:F:34:VAL:HG23	2.12	0.48
2:J:206:ARG:HH12	3:G:238:GLU:CD	2.22	0.48
1:E:105:ILE:HD12	1:E:114:VAL:HG21	1.94	0.48
10:A:666:HOH:O	2:B:57:ARG:HG3	2.12	0.48
2:B:237:TRP:CD1	5:B:305:D10:H51	2.49	0.48
2:J:39:ILE:HG13	3:K:149:ALA:HB1	1.95	0.48
6:G:302:PLC:C7	8:G:303:HXG:H36	2.24	0.48
3:K:93:LEU:HD13	3:K:174:ILE:HG23	1.94	0.48
2:F:52:SER:HA	2:F:55:LYS:HG3	1.96	0.48
1:I:348:VAL:HB	1:I:351:ASN:HB2	1.96	0.48
3:K:104:ALA:HB2	3:K:186:ILE:HG12	1.96	0.48
1:A:173:LEU:HD21	2:B:172:GLU:HB2	1.96	0.48
3:K:85:PHE:CE1	3:K:171:PRO:HD3	2.40	0.48
1:A:348:VAL:HB	1:A:351:ASN:HB2	1.96	0.48
2:F:196:TYR:CE1	1:E:145:GLN:HA	2.49	0.48
3:K:275:GLY:HA3	3:K:276:GLN:HA	1.50	0.48
2:B:52:SER:HA	2:B:55:LYS:HG3	1.96	0.48
3:C:104:ALA:HB2	3:C:186:ILE:HG12	1.96	0.48
2:J:86:PRO:HA	10:J:423:HOH:O	2.12	0.48
1:I:93:MET:HG3	1:I:95:GLY:O	2.14	0.48
6:G:302:PLC:H2A2	6:G:302:PLC:H4'1	1.96	0.48
2:J:189:VAL:HG13	10:I:615:HOH:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:93:MET:HG3	1:E:95:GLY:O	2.14	0.48
2:F:138:SER:O	2:F:140:SER:N	2.47	0.48
2:F:242:ARG:NH2	7:F:309:P1O:C4	2.77	0.48
2:J:237:TRP:CD1	5:J:306:D10:H51	2.49	0.48
3:G:104:ALA:HB2	3:G:186:ILE:HG12	1.96	0.48
3:K:230:GLY:O	3:K:244:LEU:HD13	2.13	0.48
3:C:230:GLY:O	3:C:244:LEU:HD13	2.13	0.47
2:F:37:TYR:OH	2:F:68:LEU:O	2.29	0.47
2:J:52:SER:HA	2:J:55:LYS:HG3	1.96	0.47
2:J:138:SER:O	2:J:140:SER:N	2.47	0.47
2:J:244:LEU:HD11	7:J:301:P1O:H43	1.94	0.47
1:I:90:ASN:HB3	1:I:141:MET:HG3	1.95	0.47
10:A:615:HOH:O	2:B:189:VAL:HG13	2.14	0.47
2:F:237:TRP:CD1	5:F:307:D10:H51	2.49	0.47
1:A:93:MET:HG3	1:A:95:GLY:O	2.14	0.47
1:A:145:GLN:HA	2:B:196:TYR:CE1	2.49	0.47
2:B:244:LEU:HD23	7:B:307:P1O:H13	1.86	0.47
1:E:348:VAL:HB	1:E:351:ASN:HB2	1.96	0.47
2:B:138:SER:O	2:B:140:SER:N	2.47	0.47
2:J:219:PHE:CD1	3:G:226:LEU:HA	2.49	0.47
3:G:74:TRP:HA	3:G:78:LEU:HG	1.97	0.47
3:G:230:GLY:O	3:G:244:LEU:HD13	2.13	0.47
3:C:74:TRP:HA	3:C:78:LEU:HG	1.97	0.47
3:C:75:SER:O	3:C:75:SER:OG	2.28	0.47
2:J:172:GLU:HB2	1:I:173:LEU:HD21	1.96	0.47
3:G:75:SER:O	3:G:75:SER:OG	2.28	0.47
3:K:74:TRP:HA	3:K:78:LEU:HG	1.97	0.47
6:K:302:PLC:H2A2	6:K:302:PLC:H4'1	1.96	0.47
1:A:195:TRP:HB3	2:B:125:VAL:HB	1.97	0.47
2:B:57:ARG:HH21	2:B:201:GLU:CD	2.23	0.47
3:C:61:PHE:HA	6:C:307:PLC:H8'1	1.96	0.47
3:C:89:TRP:HB3	3:C:174:ILE:HD13	1.97	0.47
6:C:302:PLC:H4'1	6:C:302:PLC:H2A2	1.96	0.47
2:F:172:GLU:HB2	1:E:173:LEU:HD21	1.97	0.47
2:F:245:GLN:O	3:G:206:ILE:HD12	2.15	0.47
3:K:89:TRP:HB3	3:K:174:ILE:HD13	1.97	0.47
3:C:223:ASN:ND2	3:C:248:PHE:CD1	2.80	0.47
3:C:226:LEU:HA	2:F:219:PHE:CD1	2.49	0.47
2:J:242:ARG:NH2	7:J:301:P1O:H8	2.30	0.47
3:K:113:LYS:HB3	3:K:113:LYS:HE2	1.60	0.47
2:B:237:TRP:HB2	5:B:304:D10:H51	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:183:ILE:HG12	6:C:307:PLC:CBA	2.45	0.47
2:B:245:GLN:O	3:C:206:ILE:HD12	2.15	0.47
2:F:54:TRP:HH2	2:F:197:ILE:HG22	1.80	0.47
2:J:54:TRP:HH2	2:J:197:ILE:HG22	1.80	0.47
3:K:61:PHE:HA	6:K:307:PLC:H8'1	1.96	0.47
2:F:57:ARG:HH21	2:F:201:GLU:CD	2.23	0.46
2:J:205:LEU:HB2	1:E:147:GLY:O	2.15	0.46
10:J:406:HOH:O	3:K:263:THR:HG21	2.14	0.46
2:B:54:TRP:HH2	2:B:197:ILE:HG22	1.80	0.46
1:I:306:ASN:ND2	1:I:354:LEU:O	2.49	0.46
2:B:242:ARG:NH2	7:B:307:P1O:H8	2.30	0.46
2:J:57:ARG:HH21	2:J:201:GLU:CD	2.23	0.46
2:J:138:SER:HA	7:J:303:P1O:C19	2.42	0.46
3:G:112:TRP:CZ2	7:G:309:P1O:H1	2.50	0.46
1:A:258:TYR:CE2	2:B:166:PRO:HG3	2.51	0.46
1:A:306:ASN:ND2	1:A:354:LEU:O	2.49	0.46
2:B:247:THR:HG23	3:C:206:ILE:HD13	1.98	0.46
2:F:36:SER:HB3	3:G:255:ALA:HB3	1.98	0.46
2:F:39:ILE:HG13	3:G:149:ALA:HB1	1.98	0.46
2:F:237:TRP:NE1	5:F:307:D10:C3	2.76	0.46
10:F:407:HOH:O	3:G:263:THR:HG21	2.16	0.46
3:G:61:PHE:HA	6:G:307:PLC:H8'1	1.96	0.46
3:G:96:GLU:OE2	3:G:178:TYR:HB2	2.15	0.46
3:K:183:ILE:HG12	6:K:307:PLC:CBA	2.45	0.46
1:A:35:GLU:O	1:A:36:LYS:HB2	2.16	0.46
1:A:379:ILE:HG21	1:A:409:LEU:HD23	1.98	0.46
2:F:247:THR:HG23	3:G:206:ILE:HD13	1.96	0.46
3:G:183:ILE:HG12	6:G:307:PLC:CBA	2.45	0.46
6:K:302:PLC:C7	8:K:303:HXG:H36	2.24	0.46
2:J:104:ARG:HA	2:J:108:PHE:HB2	1.98	0.46
2:J:196:TYR:CE1	1:I:145:GLN:HA	2.50	0.46
2:J:237:TRP:HB2	5:J:305:D10:H51	1.97	0.46
1:I:105:ILE:HG13	1:I:110:VAL:HG21	1.98	0.46
1:A:237:MET:HG3	2:B:137:LEU:HD11	1.98	0.46
3:C:153:THR:HA	3:C:176:GLU:OE2	2.16	0.46
1:E:105:ILE:HG13	1:E:110:VAL:HG21	1.98	0.46
1:E:306:ASN:ND2	1:E:354:LEU:O	2.48	0.46
1:I:35:GLU:O	1:I:36:LYS:HB2	2.16	0.46
1:I:338:GLU:CD	1:I:338:GLU:H	2.24	0.46
2:B:36:SER:HB3	3:C:255:ALA:HB3	1.97	0.46
2:B:236:ARG:HA	2:B:236:ARG:HD3	1.65	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:234:TRP:CZ2	8:C:303:HXG:H18	2.48	0.46
2:F:237:TRP:HB2	5:F:306:D10:H51	1.97	0.46
1:E:35:GLU:O	1:E:36:LYS:HB2	2.16	0.46
3:G:239:LEU:HD21	3:G:241:VAL:HG23	1.98	0.46
3:K:96:GLU:OE2	3:K:178:TYR:HB2	2.15	0.46
3:K:153:THR:HA	3:K:176:GLU:OE2	2.16	0.46
3:K:221:LEU:HG	6:K:305:PLC:HE'3	1.98	0.46
3:C:181:TYR:HA	3:C:184:TYR:CE2	2.51	0.46
2:F:104:ARG:HA	2:F:108:PHE:HB2	1.98	0.46
3:G:155:GLN:HA	3:G:158:THR:HG23	1.98	0.46
3:G:221:LEU:HG	6:G:305:PLC:HE'3	1.98	0.46
3:G:234:TRP:CZ2	8:G:303:HXG:H18	2.48	0.46
1:A:381:TYR:CZ	2:F:210:LYS:HA	2.51	0.45
3:C:86:GLU:HA	3:C:90:MET:HB2	1.98	0.45
2:J:125:VAL:HB	1:I:195:TRP:HB3	1.97	0.45
1:E:338:GLU:CD	1:E:338:GLU:H	2.24	0.45
1:A:338:GLU:H	1:A:338:GLU:CD	2.24	0.45
3:C:170:THR:HG23	3:C:173:HIS:CG	2.51	0.45
2:J:36:SER:HB3	3:K:255:ALA:HB3	1.98	0.45
2:J:166:PRO:HG3	1:I:258:TYR:CE2	2.50	0.45
1:E:51:ASP:OD2	1:E:400:ARG:NH2	2.49	0.45
3:G:89:TRP:HB3	3:G:174:ILE:HD13	1.97	0.45
3:G:170:THR:HG23	3:G:173:HIS:CG	2.51	0.45
3:G:176:GLU:HG2	3:G:177:PHE:N	2.32	0.45
3:K:223:ASN:ND2	3:K:248:PHE:CD1	2.80	0.45
3:C:251:PHE:CZ	2:F:226:LEU:HD21	2.49	0.45
2:F:166:PRO:HG3	1:E:258:TYR:CE2	2.51	0.45
2:F:137:LEU:HD11	1:E:237:MET:HG3	1.99	0.45
1:E:379:ILE:HG21	1:E:409:LEU:HD23	1.98	0.45
1:I:379:ILE:HG21	1:I:409:LEU:HD23	1.98	0.45
3:G:153:THR:HA	3:G:176:GLU:OE2	2.16	0.45
2:J:83:TYR:CE2	1:I:216:LEU:HB2	2.52	0.45
2:J:231:TRP:HA	2:J:234:ILE:HB	1.99	0.45
3:G:181:TYR:HA	3:G:184:TYR:CE2	2.51	0.45
3:K:160:HIS:HE1	3:K:170:THR:HG21	1.82	0.45
3:C:96:GLU:OE2	3:C:178:TYR:HB2	2.15	0.45
2:J:236:ARG:HA	2:J:236:ARG:HD3	1.65	0.45
2:B:104:ARG:HA	2:B:108:PHE:HB2	1.98	0.45
3:C:239:LEU:HD21	3:C:241:VAL:HG23	1.98	0.45
2:F:125:VAL:HB	1:E:195:TRP:HB3	1.98	0.45
3:G:86:GLU:HA	3:G:90:MET:HB2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:181:TYR:HA	3:K:184:TYR:CE2	2.51	0.45
2:B:237:TRP:NE1	5:B:305:D10:C3	2.76	0.45
1:A:51:ASP:OD2	1:A:400:ARG:NH2	2.49	0.44
1:A:105:ILE:HG13	1:A:110:VAL:HG21	1.98	0.44
3:C:155:GLN:HA	3:C:158:THR:HG23	1.98	0.44
2:J:104:ARG:HA	2:J:104:ARG:HD3	1.87	0.44
2:J:226:LEU:HD21	3:G:251:PHE:CZ	2.48	0.44
3:K:155:GLN:HA	3:K:158:THR:HG23	1.98	0.44
3:K:170:THR:HG23	3:K:173:HIS:CG	2.51	0.44
2:B:7:ALA:N	3:C:124:PRO:HB2	2.32	0.44
2:B:231:TRP:HA	2:B:234:ILE:HB	1.99	0.44
2:F:83:TYR:CE2	1:E:216:LEU:HB2	2.52	0.44
2:F:138:SER:HA	7:F:304:P1O:C19	2.42	0.44
2:F:231:TRP:HA	2:F:234:ILE:HB	1.99	0.44
3:G:160:HIS:HE1	3:G:170:THR:HG21	1.82	0.44
3:K:239:LEU:HD21	3:K:241:VAL:HG23	1.98	0.44
2:J:191:THR:HB	3:K:162:THR:O	2.17	0.44
3:K:86:GLU:HA	3:K:90:MET:HB2	1.99	0.44
2:B:210:LYS:HA	1:I:381:TYR:CZ	2.53	0.44
3:C:221:LEU:HG	6:C:305:PLC:HE'3	1.98	0.44
2:J:137:LEU:HD11	1:I:237:MET:HG3	1.99	0.44
1:I:51:ASP:OD2	1:I:400:ARG:NH2	2.49	0.44
3:C:160:HIS:HE1	3:C:170:THR:HG21	1.82	0.44
3:C:176:GLU:HG2	3:C:177:PHE:N	2.32	0.44
2:J:219:PHE:CE1	3:G:226:LEU:HD13	2.52	0.44
2:B:219:PHE:CE1	3:K:226:LEU:HD13	2.53	0.44
2:J:237:TRP:NE1	5:J:306:D10:C3	2.76	0.44
3:G:174:ILE:HG13	3:G:174:ILE:H	1.65	0.44
3:K:67:TRP:CG	8:K:306:HXG:H40	2.51	0.44
2:B:206:ARG:HH12	3:K:238:GLU:CD	2.26	0.44
3:C:112:TRP:CZ2	7:C:309:P1O:H1	2.50	0.44
1:I:287:GLU:O	1:I:301:LYS:HB3	2.18	0.44
2:F:29:VAL:HG13	2:F:33:ILE:HD12	2.00	0.44
3:K:112:TRP:CZ2	7:K:309:P1O:H1	2.50	0.44
3:K:159:TRP:HE1	3:K:163:ILE:HD11	1.82	0.44
1:A:287:GLU:O	1:A:301:LYS:HB3	2.18	0.44
10:B:408:HOH:O	3:C:263:THR:HG21	2.17	0.44
2:J:37:TYR:OH	2:J:68:LEU:O	2.29	0.44
1:E:287:GLU:O	1:E:301:LYS:HB3	2.18	0.44
3:G:181:TYR:HB2	3:G:182:PRO:HD3	2.00	0.44
3:C:180:SER:O	3:C:184:TYR:HD2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:181:TYR:HB2	3:C:182:PRO:HD3	2.00	0.43
2:J:206:ARG:O	3:G:232:THR:HA	2.18	0.43
2:J:210:LYS:HA	1:E:381:TYR:CZ	2.52	0.43
3:G:113:LYS:HE2	3:G:113:LYS:HB3	1.60	0.43
3:K:176:GLU:HG2	3:K:177:PHE:N	2.32	0.43
1:A:306:ASN:HB2	1:A:354:LEU:HD23	2.00	0.43
2:B:29:VAL:HG13	2:B:33:ILE:HD12	2.00	0.43
3:G:155:GLN:C	3:G:158:THR:HG23	2.43	0.43
3:C:155:GLN:C	3:C:158:THR:HG23	2.44	0.43
2:J:244:LEU:CD1	7:J:301:P1O:H42	2.39	0.43
3:K:180:SER:O	3:K:184:TYR:HD2	2.01	0.43
3:K:213:LEU:HD22	3:K:258:VAL:HA	2.00	0.43
1:A:216:LEU:HB2	2:B:83:TYR:CE2	2.53	0.43
2:F:246:SER:HA	3:G:206:ILE:CD1	2.49	0.43
2:J:59:LEU:HD11	6:G:302:PLC:H1A1	2.00	0.43
1:E:312:ILE:HG12	1:E:354:LEU:HB3	2.01	0.43
3:C:238:GLU:CD	2:F:206:ARG:HH12	2.26	0.43
2:F:54:TRP:CH2	2:F:197:ILE:HG22	2.54	0.43
1:E:97:VAL:HA	1:E:132:ARG:HB2	2.00	0.43
3:K:235:PHE:H	8:K:303:HXG:H41	1.84	0.43
3:C:159:TRP:HE1	3:C:163:ILE:HD11	1.82	0.43
2:J:199:MET:SD	6:J:308:PLC:O2	2.77	0.43
1:E:324:PHE:HA	1:E:342:ALA:HB3	2.01	0.43
1:A:101:LYS:HD2	1:A:101:LYS:HA	1.79	0.43
2:B:115:PRO:HB2	2:B:117:ASN:OD1	2.19	0.43
2:F:242:ARG:NH2	7:F:309:P1O:H8	2.34	0.43
7:J:301:P1O:H15	7:J:301:P1O:C20	2.47	0.43
3:G:96:GLU:HA	6:G:307:PLC:H7A1	2.01	0.43
3:K:96:GLU:HA	6:K:307:PLC:H7A1	2.01	0.43
2:B:204:THR:C	1:I:42:MET:HE2	2.43	0.43
2:F:191:THR:HB	3:G:162:THR:O	2.19	0.43
2:J:54:TRP:CH2	2:J:197:ILE:HG22	2.54	0.43
1:E:306:ASN:HB2	1:E:354:LEU:HD23	2.01	0.43
1:I:312:ILE:HG12	1:I:354:LEU:HB3	2.01	0.43
3:G:96:GLU:OE1	3:G:179:LEU:HD13	2.19	0.43
3:G:198:ARG:HH21	7:G:308:P1O:H2	1.84	0.43
3:K:181:TYR:HB2	3:K:182:PRO:HD3	2.00	0.43
2:J:93:VAL:HG21	2:J:128:ALA:HB2	2.01	0.43
2:J:114:PHE:CE1	1:I:96:PRO:HG3	2.53	0.43
1:I:306:ASN:HB2	1:I:354:LEU:HD23	2.00	0.43
3:G:223:ASN:ND2	3:G:248:PHE:CD1	2.80	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:235:PHE:H	8:G:303:HXG:H41	1.83	0.43
3:K:57:ILE:HD12	3:K:57:ILE:HA	1.90	0.43
3:C:96:GLU:HA	6:C:307:PLC:H7A1	2.01	0.43
3:C:175:ILE:O	3:C:179:LEU:HB3	2.19	0.43
2:F:7:ALA:N	3:G:124:PRO:HB2	2.34	0.43
2:F:115:PRO:HB2	2:F:117:ASN:OD1	2.19	0.43
3:G:243:PRO:HG2	8:G:303:HXG:CAZ	2.49	0.43
3:K:160:HIS:CD2	3:K:165:ARG:NH2	2.87	0.43
1:A:96:PRO:HG3	2:B:114:PHE:CE1	2.54	0.42
1:A:97:VAL:HA	1:A:132:ARG:HB2	2.00	0.42
3:C:235:PHE:H	8:C:303:HXG:H41	1.84	0.42
3:C:243:PRO:HG2	8:C:303:HXG:CAZ	2.49	0.42
2:J:115:PRO:HB2	2:J:117:ASN:OD1	2.19	0.42
3:G:159:TRP:HE1	3:G:163:ILE:HD11	1.83	0.42
3:G:180:SER:O	3:G:184:TYR:HD2	2.01	0.42
3:K:155:GLN:C	3:K:158:THR:HG23	2.43	0.42
1:A:42:MET:HE2	2:F:204:THR:C	2.44	0.42
2:B:42:MET:SD	3:C:240:PHE:HE2	2.42	0.42
5:B:306:D10:H103	5:J:307:D10:H103	2.00	0.42
3:C:96:GLU:OE1	3:C:179:LEU:HD13	2.19	0.42
3:C:160:HIS:CD2	3:C:165:ARG:NH2	2.87	0.42
7:F:309:P1O:H43	7:F:309:P1O:H17	2.01	0.42
3:K:96:GLU:OE1	3:K:179:LEU:HD13	2.19	0.42
3:K:198:ARG:HH21	7:K:308:P1O:H2	1.84	0.42
1:A:312:ILE:HG12	1:A:354:LEU:HB3	2.01	0.42
2:B:93:VAL:HG21	2:B:128:ALA:HB2	2.01	0.42
7:B:307:P1O:H43	7:B:307:P1O:H17	2.01	0.42
2:J:29:VAL:HG13	2:J:33:ILE:HD12	2.00	0.42
2:J:214:PRO:HA	6:J:302:PLC:H2'1	2.02	0.42
3:G:173:HIS:O	3:G:174:ILE:C	2.62	0.42
3:G:175:ILE:O	3:G:179:LEU:HB3	2.19	0.42
3:K:243:PRO:HG2	8:K:303:HXG:CAZ	2.49	0.42
2:B:37:TYR:OH	2:B:68:LEU:O	2.29	0.42
3:C:173:HIS:O	3:C:174:ILE:C	2.62	0.42
3:C:213:LEU:HD22	3:C:258:VAL:HA	2.00	0.42
2:F:214:PRO:HA	6:F:303:PLC:H2'1	2.02	0.42
1:E:86:VAL:HB	1:E:145:GLN:HB2	2.01	0.42
1:I:324:PHE:HA	1:I:342:ALA:HB3	2.01	0.42
1:A:324:PHE:HA	1:A:342:ALA:HB3	2.01	0.42
3:C:243:PRO:C	3:C:245:HIS:N	2.76	0.42
2:F:38:HIS:ND1	3:G:150:SER:OG	2.24	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:213:ALA:HB3	6:J:302:PLC:C1	2.50	0.42
1:I:97:VAL:HA	1:I:132:ARG:HB2	2.00	0.42
2:F:189:VAL:HG13	10:E:613:HOH:O	2.19	0.42
2:F:198:ARG:HG2	2:F:200:VAL:HG13	2.01	0.42
2:J:198:ARG:HG2	2:J:200:VAL:HG13	2.01	0.42
7:J:301:P1O:H43	7:J:301:P1O:H17	2.01	0.42
1:E:101:LYS:HD2	1:E:101:LYS:HA	1.79	0.42
1:E:405:ILE:HA	10:E:634:HOH:O	2.20	0.42
1:I:42:MET:HE3	1:I:42:MET:HB3	1.91	0.42
3:G:213:LEU:HD22	3:G:258:VAL:HA	2.00	0.42
1:A:208:ARG:HE	1:A:208:ARG:HB2	1.51	0.42
2:B:213:ALA:HB3	6:B:301:PLC:C1	2.49	0.42
2:B:214:PRO:HA	6:B:301:PLC:H2'1	2.02	0.42
6:G:302:PLC:H1A1	6:G:302:PLC:H31	1.83	0.42
1:A:147:GLY:O	2:F:205:LEU:HB2	2.19	0.42
2:B:39:ILE:HG13	3:C:149:ALA:HB1	2.00	0.42
3:C:159:TRP:CD1	3:C:160:HIS:HD1	2.36	0.42
2:F:93:VAL:HG21	2:F:128:ALA:HB2	2.01	0.42
2:F:114:PHE:CE1	1:E:96:PRO:HG3	2.54	0.42
2:J:206:ARG:CB	3:G:236:MET:HE2	2.48	0.42
3:G:160:HIS:CD2	3:G:165:ARG:NH2	2.87	0.42
3:K:173:HIS:O	3:K:174:ILE:C	2.62	0.42
1:A:195:TRP:O	2:B:125:VAL:HG11	2.20	0.42
7:B:307:P1O:H15	7:B:307:P1O:C20	2.47	0.42
3:C:198:ARG:HH21	7:C:308:P1O:H2	1.84	0.42
2:F:136:MET:HE2	2:F:136:MET:HB3	1.91	0.42
2:F:194:PRO:HG2	2:F:197:ILE:HG13	2.02	0.42
1:I:319:THR:HG23	1:I:320:ALA:H	1.85	0.42
3:K:175:ILE:O	3:K:179:LEU:HB3	2.19	0.42
2:B:54:TRP:CH2	2:B:197:ILE:HG22	2.54	0.42
2:B:136:MET:HE2	2:B:136:MET:HB3	1.91	0.42
2:F:138:SER:C	2:F:140:SER:H	2.28	0.42
2:F:236:ARG:HA	2:F:236:ARG:HD3	1.65	0.42
1:E:391:LEU:HB2	1:E:403:VAL:HG22	2.01	0.42
3:G:243:PRO:C	3:G:245:HIS:N	2.76	0.42
1:A:86:VAL:HB	1:A:145:GLN:HB2	2.02	0.41
2:J:194:PRO:HG2	2:J:197:ILE:HG13	2.02	0.41
1:I:86:VAL:HB	1:I:145:GLN:HB2	2.01	0.41
3:G:112:TRP:CE2	7:G:309:P1O:H7	2.55	0.41
1:A:405:ILE:HA	10:A:631:HOH:O	2.19	0.41
3:C:168:ASP:OD1	3:C:228:GLU:OE1	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:125:VAL:HG11	1:E:195:TRP:O	2.20	0.41
2:J:244:LEU:O	3:K:206:ILE:HB	2.20	0.41
3:K:172:SER:O	3:K:175:ILE:HB	2.20	0.41
3:C:262:LEU:HD23	3:C:262:LEU:HA	1.91	0.41
2:J:42:MET:O	3:K:240:PHE:HD2	2.03	0.41
3:G:173:HIS:CE1	3:G:177:PHE:HE2	2.38	0.41
3:K:237:GLU:H	3:K:237:GLU:HG2	1.66	0.41
2:B:51:TRP:HB2	2:B:54:TRP:CD1	2.56	0.41
7:B:302:P1O:H2	7:B:302:P1O:H12	1.85	0.41
1:E:319:THR:HG23	1:E:320:ALA:H	1.85	0.41
1:I:208:ARG:HE	1:I:208:ARG:HB2	1.52	0.41
1:I:405:ILE:HA	10:I:632:HOH:O	2.20	0.41
3:G:172:SER:O	3:G:175:ILE:HB	2.20	0.41
3:K:174:ILE:HG13	3:K:174:ILE:H	1.65	0.41
3:C:167:THR:HG23	3:C:169:PHE:H	1.85	0.41
3:C:239:LEU:HG	3:C:241:VAL:H	1.86	0.41
3:K:168:ASP:OD1	3:K:228:GLU:OE1	2.38	0.41
3:K:173:HIS:CE1	3:K:177:PHE:HE2	2.38	0.41
3:K:236:MET:HG2	8:K:303:HXG:H38	1.95	0.41
1:A:42:MET:HE3	1:A:42:MET:HB3	1.91	0.41
2:B:197:ILE:HD13	3:C:237:GLU:HB2	2.03	0.41
3:C:242:ALA:CB	3:C:244:LEU:HD23	2.49	0.41
8:G:303:HXG:H40	8:G:303:HXG:OAW	2.21	0.41
1:A:391:LEU:HB2	1:A:403:VAL:HG22	2.02	0.41
2:B:191:THR:HB	3:C:162:THR:O	2.21	0.41
3:C:112:TRP:CZ2	7:C:309:P1O:C2	3.04	0.41
3:C:172:SER:O	3:C:175:ILE:HB	2.20	0.41
2:F:51:TRP:HB2	2:F:54:TRP:CD1	2.56	0.41
2:J:125:VAL:HG11	1:I:195:TRP:O	2.20	0.41
1:I:36:LYS:HG3	1:I:374:TYR:CD2	2.56	0.41
3:G:159:TRP:CD1	3:G:160:HIS:ND1	2.88	0.41
3:G:237:GLU:H	3:G:237:GLU:HG2	1.66	0.41
3:G:274:LEU:HD11	3:G:276:GLN:HE22	1.86	0.41
3:K:112:TRP:CE2	7:K:309:P1O:H7	2.56	0.41
3:K:234:TRP:CZ2	8:K:303:HXG:H18	2.48	0.41
2:B:138:SER:C	2:B:140:SER:H	2.28	0.41
2:B:194:PRO:HG2	2:B:197:ILE:HG13	2.02	0.41
3:C:274:LEU:HD11	3:C:276:GLN:HE22	1.86	0.41
2:J:51:TRP:HB2	2:J:54:TRP:CD1	2.56	0.41
2:J:223:MET:O	2:J:227:ILE:HG12	2.21	0.41
1:E:395:ASP:OD2	1:E:399:ASN:ND2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:167:THR:HG23	3:K:169:PHE:H	1.85	0.41
1:A:36:LYS:HG3	1:A:374:TYR:CD2	2.56	0.41
1:A:42:MET:HE3	2:F:205:LEU:HG	2.03	0.41
2:B:138:SER:HA	7:B:302:P1O:C19	2.42	0.41
2:B:198:ARG:HG2	2:B:200:VAL:HG13	2.01	0.41
2:B:223:MET:O	2:B:227:ILE:HG12	2.21	0.41
3:C:173:HIS:CE1	3:C:177:PHE:HE2	2.38	0.41
8:C:303:HXG:OAW	8:C:303:HXG:H40	2.21	0.41
2:F:53:ASP:O	2:F:198:ARG:HD2	2.21	0.41
7:F:304:P1O:H2	7:F:304:P1O:H12	1.85	0.41
7:F:309:P1O:H17	7:F:309:P1O:C20	2.51	0.41
7:F:309:P1O:H15	7:F:309:P1O:C20	2.47	0.41
7:F:309:P1O:C6	7:F:309:P1O:C20	2.99	0.41
2:J:53:ASP:O	2:J:198:ARG:HD2	2.21	0.41
2:J:59:LEU:HD13	2:J:59:LEU:HA	1.90	0.41
2:J:136:MET:HE2	2:J:136:MET:HB3	1.91	0.41
2:J:138:SER:C	2:J:140:SER:H	2.28	0.41
2:J:204:THR:C	1:E:42:MET:HE2	2.45	0.41
1:E:33:HIS:HD2	3:G:78:LEU:HB3	1.86	0.41
1:I:33:HIS:HD2	3:K:78:LEU:HB3	1.86	0.41
1:I:391:LEU:HB2	1:I:403:VAL:HG22	2.01	0.41
3:G:167:THR:HG23	3:G:169:PHE:H	1.85	0.41
3:K:48:LYS:HD3	3:K:50:TRP:CD1	2.41	0.41
3:C:75:SER:O	3:C:76:ALA:HB3	2.21	0.41
3:C:226:LEU:HD13	2:F:219:PHE:CE1	2.55	0.41
3:G:168:ASP:OD1	3:G:228:GLU:OE1	2.38	0.41
3:G:239:LEU:HG	3:G:241:VAL:H	1.86	0.41
2:B:59:LEU:HD11	6:K:302:PLC:H1A1	2.03	0.40
2:F:42:MET:SD	3:G:240:PHE:HE2	2.44	0.40
3:G:159:TRP:CD1	3:G:160:HIS:HD1	2.37	0.40
3:G:262:LEU:HD23	3:G:262:LEU:HA	1.91	0.40
3:K:159:TRP:CD1	3:K:160:HIS:ND1	2.88	0.40
3:K:274:LEU:HD11	3:K:276:GLN:HE22	1.86	0.40
8:K:303:HXG:OAW	8:K:303:HXG:H40	2.21	0.40
1:A:342:ALA:HB1	1:A:345:GLY:HA3	2.04	0.40
2:B:53:ASP:O	2:B:198:ARG:HD2	2.21	0.40
2:B:205:LEU:HG	1:I:42:MET:HE3	2.02	0.40
2:F:38:HIS:CD2	2:F:42:MET:HE2	2.57	0.40
1:E:293:VAL:HA	1:E:294:PRO:HA	1.92	0.40
3:G:112:TRP:CZ2	7:G:309:P1O:C2	3.04	0.40
1:A:395:ASP:OD2	1:A:399:ASN:ND2	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:42:MET:O	3:G:240:PHE:HD2	2.04	0.40
3:K:89:TRP:CZ2	3:K:171:PRO:HB3	2.56	0.40
2:B:47:ASP:CG	3:C:239:LEU:HD12	2.47	0.40
2:B:233:PHE:CG	5:B:303:D10:H51	2.57	0.40
3:C:113:LYS:HE2	3:C:113:LYS:HB3	1.60	0.40
2:F:213:ALA:HB3	6:F:303:PLC:C1	2.50	0.40
2:F:223:MET:O	2:F:227:ILE:HG12	2.21	0.40
3:G:242:ALA:CB	3:G:244:LEU:HD23	2.49	0.40
3:K:75:SER:O	3:K:76:ALA:HB3	2.21	0.40
2:B:42:MET:O	3:C:240:PHE:HD2	2.04	0.40
2:B:199:MET:SD	6:B:308:PLC:O2	2.79	0.40
7:B:307:P1O:H17	7:B:307:P1O:C20	2.51	0.40
3:C:138:LEU:HD23	3:C:138:LEU:HA	1.91	0.40
7:J:301:P1O:C6	7:J:301:P1O:C20	2.99	0.40
1:E:36:LYS:HG3	1:E:374:TYR:CD2	2.56	0.40
1:E:342:ALA:HB1	1:E:345:GLY:HA3	2.04	0.40
1:I:395:ASP:OD2	1:I:399:ASN:ND2	2.54	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	380/414 (92%)	364 (96%)	16 (4%)	0	100	100
1	E	380/414 (92%)	364 (96%)	16 (4%)	0	100	100
1	I	380/414 (92%)	364 (96%)	16 (4%)	0	100	100
2	B	239/247 (97%)	227 (95%)	11 (5%)	1 (0%)	30	34
2	F	239/247 (97%)	227 (95%)	11 (5%)	1 (0%)	30	34
2	J	239/247 (97%)	228 (95%)	10 (4%)	1 (0%)	30	34
3	C	234/260 (90%)	207 (88%)	21 (9%)	6 (3%)	4	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	G	234/260 (90%)	208 (89%)	20 (8%)	6 (3%)	4	2
3	K	234/260 (90%)	207 (88%)	21 (9%)	6 (3%)	4	2
All	All	2559/2763 (93%)	2396 (94%)	142 (6%)	21 (1%)	18	16

All (21) 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
2	F	244	LEU
2	J	244	LEU
3	G	233	PHE
3	K	233	PHE
3	C	235	PHE
3	G	235	PHE
3	K	235	PHE
3	C	234	TRP
3	G	234	TRP
3	K	234	TRP
3	C	179	LEU
3	G	179	LEU
3	K	179	LEU
3	C	181	TYR
3	G	181	TYR
3	K	181	TYR

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	169/345 (49%)	162 (96%)	7 (4%)	27	37
1	E	212/345 (61%)	205 (97%)	7 (3%)	33	45

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	I	323/345 (94%)	313 (97%)	10 (3%)	35	48
2	F	206/210 (98%)	201 (98%)	5 (2%)	43	58
2	J	206/210 (98%)	201 (98%)	5 (2%)	43	58
3	C	50/212 (24%)	44 (88%)	6 (12%)	5	4
3	G	200/212 (94%)	182 (91%)	18 (9%)	9	9
3	K	200/212 (94%)	182 (91%)	18 (9%)	9	9
All	All	1566/2091 (75%)	1490 (95%)	76 (5%)	24	29

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	HIS
1	A	35	GLU
1	A	36	LYS
1	A	64	THR
1	A	65	VAL
1	A	180	LEU
1	A	227	LEU
3	C	228	GLU
3	C	231	HIS
3	C	236	MET
3	C	237	GLU
3	C	240	PHE
3	C	244	LEU
2	F	8	VAL
2	F	39	ILE
2	F	59	LEU
2	F	138	SER
2	F	244	LEU
2	J	8	VAL
2	J	39	ILE
2	J	59	LEU
2	J	138	SER
2	J	244	LEU
1	E	33	HIS
1	E	35	GLU
1	E	36	LYS
1	E	64	THR
1	E	65	VAL
1	E	180	LEU

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Mol	Chain	Res	Type
1	E	397	THR
1	I	33	HIS
1	I	35	GLU
1	I	36	LYS
1	I	64	THR
1	I	65	VAL
1	I	180	LEU
1	I	227	LEU
1	I	256	SER
1	I	309	ASN
1	I	397	THR
3	G	45	LEU
3	G	52	THR
3	G	74	TRP
3	G	87	THR
3	G	123	THR
3	G	154	GLU
3	G	158	THR
3	G	166	ASP
3	G	170	THR
3	G	172	SER
3	G	173	HIS
3	G	207	SER
3	G	228	GLU
3	G	231	HIS
3	G	236	MET
3	G	237	GLU
3	G	240	PHE
3	G	244	LEU
3	K	45	LEU
3	K	52	THR
3	K	74	TRP
3	K	87	THR
3	K	123	THR
3	K	154	GLU
3	K	158	THR
3	K	166	ASP
3	K	170	THR
3	K	172	SER
3	K	173	HIS
3	K	207	SER
3	K	228	GLU

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Mol	Chain	Res	Type
3	K	231	HIS
3	K	236	MET
3	K	237	GLU
3	K	240	PHE
3	K	244	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	ASN
2	F	174	ASN
2	F	187	ASN
2	J	174	ASN
2	J	187	ASN
1	E	168	ASN
1	I	168	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 [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	G	307	-	41,41,41	0.89	0	47,49,49	0.69	2 (4%)
7	P1O	C	309	-	37,37,37	1.13	2 (5%)	43,45,45	1.14	3 (6%)
6	PLC	J	308	-	41,41,41	1.05	2 (4%)	47,49,49	1.05	4 (8%)
9	ETF	C	310	4	5,5,5	0.76	0	7,7,7	0.31	0
7	P1O	J	303	-	37,37,37	1.12	2 (5%)	43,45,45	1.12	4 (9%)
5	D10	F	308	-	9,9,9	0.19	0	8,8,8	0.55	0
5	D10	B	305	-	9,9,9	0.21	0	8,8,8	0.55	0
6	PLC	G	305	-	41,41,41	0.90	0	47,49,49	0.69	1 (2%)
6	PLC	G	302	-	41,41,41	1.06	2 (4%)	47,49,49	1.11	3 (6%)
7	P1O	G	309	-	37,37,37	1.13	2 (5%)	43,45,45	1.13	3 (6%)
6	PLC	F	302	-	41,41,41	0.84	0	47,49,49	0.75	1 (2%)
5	D10	F	305	-	9,9,9	0.20	0	8,8,8	0.56	0
7	P1O	F	304	-	37,37,37	1.11	2 (5%)	43,45,45	1.12	3 (6%)
5	D10	E	501	-	9,9,9	0.20	0	8,8,8	0.56	0
6	PLC	B	301	-	41,41,41	0.91	0	47,49,49	0.85	1 (2%)
6	PLC	K	302	-	41,41,41	1.06	2 (4%)	47,49,49	1.11	3 (6%)
6	PLC	K	307	-	41,41,41	0.89	0	47,49,49	0.68	2 (4%)
9	ETF	G	310	4	5,5,5	0.76	0	7,7,7	0.31	0
7	P1O	J	301	-	37,37,37	1.11	2 (5%)	43,45,45	1.11	3 (6%)
8	HXG	K	303	-	29,29,29	0.35	0	35,37,37	0.36	0
5	D10	F	306	-	9,9,9	0.21	0	8,8,8	0.57	0
7	P1O	C	308	-	37,37,37	1.10	2 (5%)	43,45,45	1.13	4 (9%)
6	PLC	C	307	-	41,41,41	0.89	0	47,49,49	0.69	2 (4%)
6	PLC	K	305	-	41,41,41	0.90	0	47,49,49	0.68	1 (2%)
8	HXG	C	303	-	29,29,29	0.35	0	35,37,37	0.36	0
7	P1O	B	307	-	37,37,37	1.13	2 (5%)	43,45,45	1.10	3 (6%)
5	D10	J	307	-	9,9,9	0.20	0	8,8,8	0.56	0
6	PLC	J	302	-	41,41,41	0.91	0	47,49,49	0.85	1 (2%)
6	PLC	B	309	-	41,41,41	0.84	0	47,49,49	0.75	1 (2%)
8	HXG	K	306	-	29,29,29	0.35	0	35,37,37	0.37	0
6	PLC	C	302	-	41,41,41	1.06	2 (4%)	47,49,49	1.11	3 (6%)
5	D10	J	306	-	9,9,9	0.21	0	8,8,8	0.55	0
8	HXG	C	306	-	29,29,29	0.36	0	35,37,37	0.37	0
5	D10	K	304	-	9,9,9	0.20	0	8,8,8	0.56	0
5	D10	J	305	-	9,9,9	0.20	0	8,8,8	0.56	0
5	D10	C	304	-	9,9,9	0.20	0	8,8,8	0.56	0
6	PLC	F	301	-	41,41,41	1.05	2 (4%)	47,49,49	1.05	4 (8%)
5	D10	I	501	-	9,9,9	0.20	0	8,8,8	0.56	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	D10	B	304	-	9,9,9	0.21	0	8,8,8	0.56	0
7	P1O	K	308	-	37,37,37	1.10	2 (5%)	43,45,45	1.13	4 (9%)
7	P1O	F	309	-	37,37,37	1.12	2 (5%)	43,45,45	1.10	3 (6%)
5	D10	A	503	-	9,9,9	0.20	0	8,8,8	0.55	0
7	P1O	G	308	-	37,37,37	1.10	2 (5%)	43,45,45	1.13	4 (9%)
7	P1O	K	309	-	37,37,37	1.13	2 (5%)	43,45,45	1.14	3 (6%)
5	D10	B	303	-	9,9,9	0.20	0	8,8,8	0.57	0
6	PLC	C	305	-	41,41,41	0.90	0	47,49,49	0.69	1 (2%)
8	HXG	G	303	-	29,29,29	0.35	0	35,37,37	0.37	0
5	D10	F	307	-	9,9,9	0.21	0	8,8,8	0.55	0
6	PLC	F	303	-	41,41,41	0.91	0	47,49,49	0.84	1 (2%)
8	HXG	G	306	-	29,29,29	0.35	0	35,37,37	0.37	0
6	PLC	J	309	-	41,41,41	0.84	0	47,49,49	0.75	1 (2%)
5	D10	B	306	-	9,9,9	0.20	0	8,8,8	0.55	0
5	D10	J	304	-	9,9,9	0.20	0	8,8,8	0.56	0
5	D10	G	304	-	9,9,9	0.20	0	8,8,8	0.56	0
6	PLC	B	308	-	41,41,41	1.05	2 (4%)	47,49,49	1.05	4 (8%)
9	ETF	K	310	4	5,5,5	0.76	0	7,7,7	0.31	0
7	P1O	B	302	-	37,37,37	1.11	2 (5%)	43,45,45	1.12	3 (6%)

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	G	307	-	-	16/45/45/45	-
7	P1O	C	309	-	-	26/41/41/41	-
6	PLC	J	308	-	-	25/45/45/45	-
9	ETF	C	310	4	-	0/3/3/3	-
7	P1O	J	303	-	-	26/41/41/41	-
5	D10	F	308	-	-	5/7/7/7	-
5	D10	B	305	-	-	6/7/7/7	-
6	PLC	G	305	-	-	14/45/45/45	-
6	PLC	G	302	-	-	22/45/45/45	-
7	P1O	G	309	-	-	26/41/41/41	-
6	PLC	F	302	-	-	15/45/45/45	-
5	D10	F	305	-	-	0/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	P1O	F	304	-	-	26/41/41/41	-
5	D10	E	501	-	-	0/7/7/7	-
6	PLC	B	301	-	-	15/45/45/45	-
6	PLC	K	302	-	-	22/45/45/45	-
6	PLC	K	307	-	-	16/45/45/45	-
9	ETF	G	310	4	-	0/3/3/3	-
7	P1O	J	301	-	-	20/41/41/41	-
8	HXG	K	303	-	-	9/33/33/33	-
5	D10	F	306	-	-	2/7/7/7	-
7	P1O	C	308	-	-	19/41/41/41	-
6	PLC	C	307	-	-	16/45/45/45	-
6	PLC	K	305	-	-	14/45/45/45	-
8	HXG	C	303	-	-	9/33/33/33	-
7	P1O	B	307	-	-	20/41/41/41	-
5	D10	J	307	-	-	5/7/7/7	-
6	PLC	J	302	-	-	15/45/45/45	-
6	PLC	B	309	-	-	15/45/45/45	-
8	HXG	K	306	-	-	7/33/33/33	-
6	PLC	C	302	-	-	22/45/45/45	-
5	D10	J	306	-	-	6/7/7/7	-
8	HXG	C	306	-	-	7/33/33/33	-
5	D10	K	304	-	-	0/7/7/7	-
5	D10	J	305	-	-	2/7/7/7	-
5	D10	C	304	-	-	0/7/7/7	-
6	PLC	F	301	-	-	25/45/45/45	-
5	D10	I	501	-	-	0/7/7/7	-
5	D10	B	304	-	-	2/7/7/7	-
7	P1O	K	308	-	-	19/41/41/41	-
7	P1O	F	309	-	-	20/41/41/41	-
5	D10	A	503	-	-	0/7/7/7	-
7	P1O	G	308	-	-	19/41/41/41	-
7	P1O	K	309	-	-	26/41/41/41	-
5	D10	B	303	-	-	0/7/7/7	-
6	PLC	C	305	-	-	14/45/45/45	-
8	HXG	G	303	-	-	9/33/33/33	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	D10	F	307	-	-	6/7/7/7	-
6	PLC	F	303	-	-	15/45/45/45	-
8	HXG	G	306	-	-	7/33/33/33	-
6	PLC	J	309	-	-	15/45/45/45	-
5	D10	B	306	-	-	5/7/7/7	-
5	D10	J	304	-	-	0/7/7/7	-
5	D10	G	304	-	-	0/7/7/7	-
6	PLC	B	308	-	-	25/45/45/45	-
9	ETF	K	310	4	-	0/3/3/3	-
7	P1O	B	302	-	-	26/41/41/41	-

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	307	P1O	O5-C9	4.36	1.46	1.33
7	F	309	P1O	O5-C9	4.34	1.46	1.33
7	J	303	P1O	O5-C9	4.32	1.45	1.33
7	B	302	P1O	O5-C9	4.30	1.45	1.33
7	C	309	P1O	O5-C9	4.30	1.45	1.33
7	K	309	P1O	O5-C9	4.29	1.45	1.33
7	G	309	P1O	O5-C9	4.29	1.45	1.33
7	F	304	P1O	O5-C9	4.28	1.45	1.33
7	J	301	P1O	O5-C9	4.28	1.45	1.33
6	C	302	PLC	O3-CB	4.25	1.45	1.33
6	G	302	PLC	O3-CB	4.25	1.45	1.33
6	K	302	PLC	O3-CB	4.25	1.45	1.33
7	G	308	P1O	O5-C9	4.23	1.45	1.33
7	C	308	P1O	O5-C9	4.23	1.45	1.33
7	K	308	P1O	O5-C9	4.23	1.45	1.33
6	F	301	PLC	O3-CB	4.22	1.45	1.33
6	B	308	PLC	O3-CB	4.22	1.45	1.33
6	J	308	PLC	O3-CB	4.21	1.45	1.33
7	F	309	P1O	O7-C19	4.16	1.46	1.34
7	B	307	P1O	O7-C19	4.16	1.46	1.34
7	J	301	P1O	O7-C19	4.15	1.46	1.34
7	C	309	P1O	O7-C19	4.14	1.46	1.34
7	K	309	P1O	O7-C19	4.14	1.46	1.34
6	J	308	PLC	O2-C'	4.14	1.46	1.34
7	G	309	P1O	O7-C19	4.13	1.46	1.34
6	B	308	PLC	O2-C'	4.12	1.45	1.34
6	F	301	PLC	O2-C'	4.11	1.45	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	J	303	P1O	O7-C19	4.10	1.45	1.34
6	G	302	PLC	O2-C'	4.09	1.45	1.34
7	B	302	P1O	O7-C19	4.08	1.45	1.34
7	F	304	P1O	O7-C19	4.07	1.45	1.34
6	K	302	PLC	O2-C'	4.06	1.45	1.34
6	C	302	PLC	O2-C'	4.05	1.45	1.34
7	G	308	P1O	O7-C19	4.02	1.45	1.34
7	C	308	P1O	O7-C19	4.02	1.45	1.34
7	K	308	P1O	O7-C19	4.01	1.45	1.34

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	J	301	P1O	O7-C19-C20	4.26	120.69	111.48
7	F	309	P1O	O7-C19-C20	4.24	120.65	111.48
7	B	307	P1O	O7-C19-C20	4.21	120.58	111.48
7	K	309	P1O	O7-C19-C20	4.17	120.51	111.48
7	C	309	P1O	O7-C19-C20	4.17	120.49	111.48
7	G	309	P1O	O7-C19-C20	4.16	120.48	111.48
7	C	308	P1O	O7-C19-C20	4.15	120.47	111.48
7	K	308	P1O	O7-C19-C20	4.14	120.45	111.48
7	G	308	P1O	O7-C19-C20	4.14	120.44	111.48
7	B	302	P1O	O7-C19-C20	4.14	120.43	111.48
7	J	303	P1O	O7-C19-C20	4.12	120.40	111.48
7	F	304	P1O	O7-C19-C20	4.11	120.37	111.48
6	C	302	PLC	O2-C'-C1'	4.09	120.34	111.48
6	K	302	PLC	O2-C'-C1'	4.09	120.32	111.48
6	G	302	PLC	O2-C'-C1'	4.08	120.30	111.48
6	J	308	PLC	O2-C'-C1'	4.06	120.27	111.48
6	B	308	PLC	O2-C'-C1'	4.06	120.27	111.48
6	F	301	PLC	O2-C'-C1'	4.06	120.27	111.48
6	B	309	PLC	O2-C'-C1'	3.80	119.70	111.48
6	F	302	PLC	O2-C'-C1'	3.80	119.70	111.48
6	J	309	PLC	O2-C'-C1'	3.79	119.67	111.48
6	J	302	PLC	O2-C'-C1'	3.72	119.53	111.48
6	B	301	PLC	O2-C'-C1'	3.72	119.53	111.48
6	F	303	PLC	O2-C'-C1'	3.68	119.45	111.48
6	C	307	PLC	O2-C'-C1'	3.11	118.20	111.48
6	G	307	PLC	O2-C'-C1'	3.10	118.18	111.48
6	K	307	PLC	O2-C'-C1'	3.09	118.17	111.48
6	K	302	PLC	O3-CB-C1B	2.88	120.62	111.83
6	C	302	PLC	O3-CB-C1B	2.88	120.60	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	G	302	PLC	O3-CB-C1B	2.87	120.58	111.83
7	B	307	P1O	O5-C9-C10	2.82	120.43	111.83
7	J	301	P1O	O5-C9-C10	2.81	120.40	111.83
7	F	309	P1O	O5-C9-C10	2.80	120.37	111.83
6	B	308	PLC	O3-CB-C1B	2.80	120.37	111.83
7	K	308	P1O	O5-C9-C10	2.79	120.34	111.83
7	G	308	P1O	O5-C9-C10	2.79	120.34	111.83
6	F	301	PLC	O3-CB-C1B	2.79	120.33	111.83
7	C	308	P1O	O5-C9-C10	2.78	120.32	111.83
7	F	304	P1O	O5-C9-C10	2.78	120.32	111.83
6	J	308	PLC	O3-CB-C1B	2.78	120.31	111.83
7	J	303	P1O	O5-C9-C10	2.77	120.29	111.83
7	B	302	P1O	O5-C9-C10	2.77	120.28	111.83
7	K	309	P1O	O5-C9-C10	2.71	120.09	111.83
7	C	309	P1O	O5-C9-C10	2.71	120.09	111.83
7	G	309	P1O	O5-C9-C10	2.71	120.09	111.83
6	G	302	PLC	C2-O2-C'	-2.50	111.81	117.80
6	K	302	PLC	C2-O2-C'	-2.48	111.85	117.80
6	C	302	PLC	C2-O2-C'	-2.48	111.85	117.80
7	C	309	P1O	C7-O7-C19	-2.46	111.90	117.80
7	K	309	P1O	C7-O7-C19	-2.46	111.91	117.80
7	G	309	P1O	C7-O7-C19	-2.45	111.93	117.80
6	C	305	PLC	O2-C'-C1'	2.42	116.72	111.48
6	G	305	PLC	O2-C'-C1'	2.42	116.72	111.48
6	K	305	PLC	O2-C'-C1'	2.41	116.70	111.48
7	B	302	P1O	C7-O7-C19	-2.37	112.13	117.80
7	J	303	P1O	C7-O7-C19	-2.35	112.18	117.80
7	F	304	P1O	C7-O7-C19	-2.34	112.20	117.80
7	G	308	P1O	C7-O7-C19	-2.30	112.30	117.80
7	C	308	P1O	C7-O7-C19	-2.30	112.30	117.80
7	K	308	P1O	C7-O7-C19	-2.28	112.35	117.80
6	G	307	PLC	O2-C'-O'	-2.25	118.44	123.70
6	K	307	PLC	O2-C'-O'	-2.24	118.47	123.70
6	C	307	PLC	O2-C'-O'	-2.24	118.47	123.70
6	J	308	PLC	C2-O2-C'	-2.23	112.46	117.80
6	B	308	PLC	C2-O2-C'	-2.22	112.47	117.80
6	F	301	PLC	C2-O2-C'	-2.22	112.49	117.80
7	F	309	P1O	O7-C19-O8	-2.14	118.69	123.70
7	J	301	P1O	O7-C19-O8	-2.12	118.75	123.70
7	B	307	P1O	O7-C19-O8	-2.09	118.81	123.70
6	J	308	PLC	O2-C'-O'	-2.02	118.97	123.70
7	G	308	P1O	C1-C2-N1	-2.02	109.35	115.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	308	PLC	O2-C'-O'	-2.02	118.99	123.70
7	J	303	P1O	O7-C19-O8	-2.01	119.00	123.70
7	C	308	P1O	C1-C2-N1	-2.01	109.36	115.82
7	K	308	P1O	C1-C2-N1	-2.01	109.37	115.82
6	F	301	PLC	O2-C'-O'	-2.00	119.02	123.70

There are no chirality outliers.

All (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	309	PLC	C1'-C'-O2-C2
6	B	309	PLC	C1-O3P-P-O1P
6	B	309	PLC	C1-O3P-P-O4P
6	B	309	PLC	C4-O4P-P-O1P
6	C	302	PLC	O4P-C4-C5-N
6	C	302	PLC	C1'-C'-O2-C2
6	C	302	PLC	C1B-CB-O3-C3
6	C	302	PLC	OB-CB-O3-C3
6	C	302	PLC	C1-O3P-P-O2P
6	C	302	PLC	C1-O3P-P-O4P
6	C	302	PLC	C4-O4P-P-O2P
6	C	302	PLC	C4-O4P-P-O3P
6	C	305	PLC	C1'-C'-O2-C2
6	C	305	PLC	C4-O4P-P-O1P
6	C	305	PLC	C4-O4P-P-O3P
6	C	307	PLC	C1'-C'-O2-C2
6	C	307	PLC	O'-C'-O2-C2
6	C	307	PLC	C1-O3P-P-O1P
6	F	301	PLC	C1'-C'-O2-C2
6	F	302	PLC	C1'-C'-O2-C2
6	F	302	PLC	C1-O3P-P-O1P
6	F	302	PLC	C1-O3P-P-O4P
6	F	302	PLC	C4-O4P-P-O1P
6	F	303	PLC	C1'-C'-O2-C2
6	F	303	PLC	C1-O3P-P-O2P
6	F	303	PLC	C1-O3P-P-O4P
6	J	302	PLC	C1'-C'-O2-C2
6	J	302	PLC	C1-O3P-P-O2P

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Mol	Chain	Res	Type	Atoms
6	J	302	PLC	C1-O3P-P-O4P
6	J	308	PLC	C1'-C'-O2-C2
6	J	309	PLC	C1'-C'-O2-C2
6	J	309	PLC	C1-O3P-P-O1P
6	J	309	PLC	C1-O3P-P-O4P
6	J	309	PLC	C4-O4P-P-O1P
6	G	302	PLC	O4P-C4-C5-N
6	G	302	PLC	C1'-C'-O2-C2
6	G	302	PLC	C1B-CB-O3-C3
6	G	302	PLC	OB-CB-O3-C3
6	G	302	PLC	C1-O3P-P-O2P
6	G	302	PLC	C1-O3P-P-O4P
6	G	302	PLC	C4-O4P-P-O2P
6	G	302	PLC	C4-O4P-P-O3P
6	G	305	PLC	C1'-C'-O2-C2
6	G	305	PLC	C4-O4P-P-O1P
6	G	305	PLC	C4-O4P-P-O3P
6	G	307	PLC	C1'-C'-O2-C2
6	G	307	PLC	O'-C'-O2-C2
6	G	307	PLC	C1-O3P-P-O1P
6	K	302	PLC	O4P-C4-C5-N
6	K	302	PLC	C1'-C'-O2-C2
6	K	302	PLC	C1B-CB-O3-C3
6	K	302	PLC	OB-CB-O3-C3
6	K	302	PLC	C1-O3P-P-O2P
6	K	302	PLC	C1-O3P-P-O4P
6	K	302	PLC	C4-O4P-P-O2P
6	K	302	PLC	C4-O4P-P-O3P
6	K	305	PLC	C1'-C'-O2-C2
6	K	305	PLC	C4-O4P-P-O1P
6	K	305	PLC	C4-O4P-P-O3P
6	K	307	PLC	C1'-C'-O2-C2
6	K	307	PLC	O'-C'-O2-C2
6	K	307	PLC	C1-O3P-P-O1P
7	B	302	P1O	C1-O3-P1-O1
7	B	302	P1O	C1-O3-P1-O2
7	B	302	P1O	C1-O3-P1-O4
7	B	302	P1O	C6-O4-P1-O1
7	B	302	P1O	C6-O4-P1-O3
7	B	307	P1O	C6-O4-P1-O1
7	B	307	P1O	C6-O4-P1-O2
7	B	307	P1O	C6-O4-P1-O3

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Mol	Chain	Res	Type	Atoms
7	B	307	P1O	O8-C19-O7-C7
7	C	308	P1O	C1-O3-P1-O1
7	C	308	P1O	C1-O3-P1-O4
7	C	308	P1O	C6-O4-P1-O3
7	C	308	P1O	C2-C1-O3-P1
7	C	308	P1O	O3-C1-C2-N1
7	C	308	P1O	O6-C9-O5-C8
7	C	308	P1O	C10-C9-O5-C8
7	C	308	P1O	O8-C19-O7-C7
7	C	309	P1O	C1-O3-P1-O1
7	C	309	P1O	C1-O3-P1-O2
7	C	309	P1O	C6-O4-P1-O1
7	C	309	P1O	C6-O4-P1-O2
7	C	309	P1O	C6-O4-P1-O3
7	F	304	P1O	C1-O3-P1-O1
7	F	304	P1O	C1-O3-P1-O2
7	F	304	P1O	C1-O3-P1-O4
7	F	304	P1O	C6-O4-P1-O1
7	F	304	P1O	C6-O4-P1-O3
7	F	309	P1O	C6-O4-P1-O1
7	F	309	P1O	C6-O4-P1-O2
7	F	309	P1O	C6-O4-P1-O3
7	F	309	P1O	O8-C19-O7-C7
7	J	301	P1O	C6-O4-P1-O1
7	J	301	P1O	C6-O4-P1-O2
7	J	301	P1O	C6-O4-P1-O3
7	J	301	P1O	O8-C19-O7-C7
7	J	303	P1O	C1-O3-P1-O1
7	J	303	P1O	C1-O3-P1-O2
7	J	303	P1O	C1-O3-P1-O4
7	J	303	P1O	C6-O4-P1-O1
7	J	303	P1O	C6-O4-P1-O3
7	G	308	P1O	C1-O3-P1-O1
7	G	308	P1O	C1-O3-P1-O4
7	G	308	P1O	C6-O4-P1-O3
7	G	308	P1O	C2-C1-O3-P1
7	G	308	P1O	O3-C1-C2-N1
7	G	308	P1O	O6-C9-O5-C8
7	G	308	P1O	C10-C9-O5-C8
7	G	308	P1O	O8-C19-O7-C7
7	G	309	P1O	C1-O3-P1-O1
7	G	309	P1O	C1-O3-P1-O2

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Mol	Chain	Res	Type	Atoms
7	G	309	P1O	C6-O4-P1-O1
7	G	309	P1O	C6-O4-P1-O2
7	G	309	P1O	C6-O4-P1-O3
7	K	308	P1O	C1-O3-P1-O1
7	K	308	P1O	C1-O3-P1-O4
7	K	308	P1O	C6-O4-P1-O3
7	K	308	P1O	C2-C1-O3-P1
7	K	308	P1O	O3-C1-C2-N1
7	K	308	P1O	O6-C9-O5-C8
7	K	308	P1O	C10-C9-O5-C8
7	K	308	P1O	O8-C19-O7-C7
7	K	309	P1O	C1-O3-P1-O1
7	K	309	P1O	C1-O3-P1-O2
7	K	309	P1O	C6-O4-P1-O1
7	K	309	P1O	C6-O4-P1-O2
7	K	309	P1O	C6-O4-P1-O3
8	C	303	HXG	CAU-OAX-PBD-OAH
8	C	303	HXG	CAU-OAX-PBD-OAW
8	C	303	HXG	CAP-OAW-PBD-OAX
8	C	306	HXG	OAW-CAP-CAS-NBC
8	G	303	HXG	CAU-OAX-PBD-OAH
8	G	303	HXG	CAU-OAX-PBD-OAW
8	G	303	HXG	CAP-OAW-PBD-OAX
8	G	306	HXG	OAW-CAP-CAS-NBC
8	K	303	HXG	CAU-OAX-PBD-OAH
8	K	303	HXG	CAU-OAX-PBD-OAW
8	K	303	HXG	CAP-OAW-PBD-OAX
8	K	306	HXG	OAW-CAP-CAS-NBC
6	B	301	PLC	OB-CB-O3-C3
6	F	303	PLC	OB-CB-O3-C3
6	J	302	PLC	OB-CB-O3-C3
6	C	305	PLC	C2-C3-O3-CB
6	G	305	PLC	C2-C3-O3-CB
6	K	305	PLC	C2-C3-O3-CB
6	B	301	PLC	C1B-CB-O3-C3
6	F	303	PLC	C1B-CB-O3-C3
6	J	302	PLC	C1B-CB-O3-C3
7	B	302	P1O	C10-C9-O5-C8
7	F	304	P1O	C10-C9-O5-C8
7	J	303	P1O	C10-C9-O5-C8
6	B	309	PLC	OB-CB-O3-C3
6	C	305	PLC	OB-CB-O3-C3

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Mol	Chain	Res	Type	Atoms
6	C	307	PLC	OB-CB-O3-C3
6	F	302	PLC	OB-CB-O3-C3
6	J	309	PLC	OB-CB-O3-C3
6	G	305	PLC	OB-CB-O3-C3
6	G	307	PLC	OB-CB-O3-C3
6	K	305	PLC	OB-CB-O3-C3
6	K	307	PLC	OB-CB-O3-C3
7	B	302	P1O	O6-C9-O5-C8
7	F	304	P1O	O6-C9-O5-C8
7	J	303	P1O	O6-C9-O5-C8
6	B	301	PLC	O'-C'-O2-C2
6	B	308	PLC	O'-C'-O2-C2
6	B	309	PLC	O'-C'-O2-C2
6	C	302	PLC	O'-C'-O2-C2
6	C	305	PLC	O'-C'-O2-C2
6	F	301	PLC	O'-C'-O2-C2
6	F	302	PLC	O'-C'-O2-C2
6	F	303	PLC	O'-C'-O2-C2
6	J	302	PLC	O'-C'-O2-C2
6	J	308	PLC	O'-C'-O2-C2
6	J	309	PLC	O'-C'-O2-C2
6	G	302	PLC	O'-C'-O2-C2
6	G	305	PLC	O'-C'-O2-C2
6	K	302	PLC	O'-C'-O2-C2
6	K	305	PLC	O'-C'-O2-C2
6	B	301	PLC	C2-C3-O3-CB
6	F	303	PLC	C2-C3-O3-CB
6	J	302	PLC	C2-C3-O3-CB
6	C	305	PLC	C1B-CB-O3-C3
6	C	307	PLC	C1B-CB-O3-C3
6	G	305	PLC	C1B-CB-O3-C3
6	G	307	PLC	C1B-CB-O3-C3
6	K	305	PLC	C1B-CB-O3-C3
6	K	307	PLC	C1B-CB-O3-C3
7	B	307	P1O	C20-C19-O7-C7
7	C	308	P1O	C20-C19-O7-C7
7	F	309	P1O	C20-C19-O7-C7
7	J	301	P1O	C20-C19-O7-C7
7	G	308	P1O	C20-C19-O7-C7
7	K	308	P1O	C20-C19-O7-C7
6	B	309	PLC	C1B-CB-O3-C3
6	F	302	PLC	C1B-CB-O3-C3

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Mol	Chain	Res	Type	Atoms
6	J	309	PLC	C1B-CB-O3-C3
6	B	308	PLC	C4-C5-N-C6
6	F	301	PLC	C4-C5-N-C6
6	J	308	PLC	C4-C5-N-C6
6	B	308	PLC	C1B-CB-O3-C3
6	F	301	PLC	C1B-CB-O3-C3
6	J	308	PLC	C1B-CB-O3-C3
7	C	308	P1O	C9-C10-C11-C12
7	G	308	P1O	C9-C10-C11-C12
7	K	308	P1O	C9-C10-C11-C12
7	C	309	P1O	C10-C9-O5-C8
7	G	309	P1O	C10-C9-O5-C8
7	K	309	P1O	C10-C9-O5-C8
7	B	307	P1O	C19-C20-C21-C22
8	C	303	HXG	CAL-CAN-CAQ-CAZ
8	G	303	HXG	CAL-CAN-CAQ-CAZ
8	K	303	HXG	CAL-CAN-CAQ-CAZ
6	B	308	PLC	OB-CB-O3-C3
6	F	301	PLC	OB-CB-O3-C3
6	J	308	PLC	OB-CB-O3-C3
7	C	309	P1O	C7-C6-O4-P1
7	G	309	P1O	C7-C6-O4-P1
7	K	309	P1O	C7-C6-O4-P1
7	F	309	P1O	C19-C20-C21-C22
7	J	301	P1O	C19-C20-C21-C22
8	C	306	HXG	CAM-CAO-CAR-CBA
8	G	306	HXG	CAM-CAO-CAR-CBA
8	K	306	HXG	CAM-CAO-CAR-CBA
6	B	308	PLC	C4-C5-N-C7
6	F	301	PLC	C4-C5-N-C7
6	J	308	PLC	C4-C5-N-C7
7	B	302	P1O	C20-C19-O7-C7
7	F	304	P1O	C20-C19-O7-C7
7	J	303	P1O	C20-C19-O7-C7
7	B	302	P1O	O8-C19-O7-C7
7	F	304	P1O	O8-C19-O7-C7
7	J	303	P1O	O8-C19-O7-C7
6	B	308	PLC	CB-C1B-C2B-C3B
6	F	301	PLC	CB-C1B-C2B-C3B
6	J	308	PLC	CB-C1B-C2B-C3B
7	C	309	P1O	O6-C9-O5-C8
7	G	309	P1O	O6-C9-O5-C8

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Mol	Chain	Res	Type	Atoms
7	K	309	P1O	O6-C9-O5-C8
6	C	305	PLC	C3-C2-O2-C'
6	G	305	PLC	C3-C2-O2-C'
6	K	305	PLC	C3-C2-O2-C'
6	B	308	PLC	C4-C5-N-C8
6	F	301	PLC	C4-C5-N-C8
6	J	308	PLC	C4-C5-N-C8
6	J	308	PLC	C2B-C3B-C4B-C5B
7	C	308	P1O	C21-C22-C23-C24
7	G	308	P1O	C21-C22-C23-C24
7	K	308	P1O	C21-C22-C23-C24
6	B	308	PLC	C2B-C3B-C4B-C5B
7	C	309	P1O	C13-C14-C15-C16
7	K	309	P1O	C13-C14-C15-C16
6	F	301	PLC	C2B-C3B-C4B-C5B
7	B	302	P1O	C11-C12-C13-C14
7	J	303	P1O	C11-C12-C13-C14
7	G	309	P1O	C13-C14-C15-C16
6	C	302	PLC	C6B-C7B-C8B-C9B
6	G	302	PLC	C6B-C7B-C8B-C9B
7	F	304	P1O	C11-C12-C13-C14
6	K	302	PLC	C6B-C7B-C8B-C9B
8	C	303	HXG	CAR-CBA-OAY-CBB
8	G	303	HXG	CAR-CBA-OAY-CBB
8	K	303	HXG	CAR-CBA-OAY-CBB
6	F	303	PLC	C5'-C6'-C7'-C8'
6	J	302	PLC	C5'-C6'-C7'-C8'
6	J	308	PLC	C4B-C5B-C6B-C7B
7	B	307	P1O	C23-C24-C25-C26
7	F	309	P1O	C23-C24-C25-C26
7	J	301	P1O	C23-C24-C25-C26
6	B	301	PLC	C5'-C6'-C7'-C8'
6	B	308	PLC	C4B-C5B-C6B-C7B
6	F	301	PLC	C4B-C5B-C6B-C7B
5	B	305	D10	C5-C6-C7-C8
6	B	308	PLC	C3B-C4B-C5B-C6B
6	F	301	PLC	C3B-C4B-C5B-C6B
6	J	308	PLC	C3B-C4B-C5B-C6B
7	B	307	P1O	C9-C10-C11-C12
7	F	309	P1O	C9-C10-C11-C12
7	J	301	P1O	C9-C10-C11-C12
6	B	309	PLC	C1'-C2'-C3'-C4'

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Mol	Chain	Res	Type	Atoms
6	J	309	PLC	C1'-C2'-C3'-C4'
7	B	307	P1O	C14-C15-C16-C17
7	F	309	P1O	C14-C15-C16-C17
7	C	309	P1O	C20-C19-O7-C7
7	G	309	P1O	C20-C19-O7-C7
7	K	309	P1O	C20-C19-O7-C7
5	F	307	D10	C5-C6-C7-C8
5	J	306	D10	C5-C6-C7-C8
6	F	302	PLC	C1'-C2'-C3'-C4'
7	J	301	P1O	C14-C15-C16-C17
6	C	302	PLC	C1'-C2'-C3'-C4'
6	G	302	PLC	C1'-C2'-C3'-C4'
6	K	302	PLC	C1'-C2'-C3'-C4'
7	F	304	P1O	C21-C22-C23-C24
7	B	302	P1O	C21-C22-C23-C24
7	J	303	P1O	C21-C22-C23-C24
7	K	309	P1O	C24-C25-C26-C27
7	C	309	P1O	C24-C25-C26-C27
7	G	309	P1O	C24-C25-C26-C27
7	B	302	P1O	C13-C14-C15-C16
7	F	304	P1O	C13-C14-C15-C16
7	C	309	P1O	C21-C22-C23-C24
7	G	309	P1O	C21-C22-C23-C24
7	K	309	P1O	C21-C22-C23-C24
7	J	303	P1O	C13-C14-C15-C16
6	C	305	PLC	CB-C1B-C2B-C3B
6	G	305	PLC	CB-C1B-C2B-C3B
6	K	305	PLC	CB-C1B-C2B-C3B
8	C	303	HXG	OAG-CBA-OAY-CBB
8	G	303	HXG	OAG-CBA-OAY-CBB
8	K	303	HXG	OAG-CBA-OAY-CBB
6	B	308	PLC	C4'-C5'-C6'-C7'
6	F	301	PLC	C4'-C5'-C6'-C7'
6	J	308	PLC	C4'-C5'-C6'-C7'
7	J	301	P1O	C13-C14-C15-C16
7	B	307	P1O	C13-C14-C15-C16
7	F	309	P1O	C13-C14-C15-C16
6	C	302	PLC	CB-C1B-C2B-C3B
6	G	302	PLC	CB-C1B-C2B-C3B
6	K	302	PLC	CB-C1B-C2B-C3B
6	B	308	PLC	C6'-C7'-C8'-C9'
6	F	301	PLC	C6'-C7'-C8'-C9'

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Mol	Chain	Res	Type	Atoms
6	J	308	PLC	C6'-C7'-C8'-C9'
6	G	302	PLC	C3B-C4B-C5B-C6B
6	C	302	PLC	C3B-C4B-C5B-C6B
6	K	302	PLC	C3B-C4B-C5B-C6B
5	F	308	D10	C5-C6-C7-C8
5	B	306	D10	C5-C6-C7-C8
5	J	307	D10	C5-C6-C7-C8
6	C	302	PLC	C4B-C5B-C6B-C7B
7	C	309	P1O	C12-C13-C14-C15
7	G	309	P1O	C12-C13-C14-C15
7	K	309	P1O	C12-C13-C14-C15
6	G	302	PLC	C4B-C5B-C6B-C7B
6	K	302	PLC	C4B-C5B-C6B-C7B
6	K	302	PLC	C6'-C7'-C8'-C9'
6	C	302	PLC	C6'-C7'-C8'-C9'
6	G	302	PLC	C6'-C7'-C8'-C9'
7	B	302	P1O	C23-C24-C25-C26
7	F	304	P1O	C23-C24-C25-C26
7	J	303	P1O	C23-C24-C25-C26
6	J	302	PLC	C4'-C5'-C6'-C7'
6	F	303	PLC	C4'-C5'-C6'-C7'
8	C	303	HXG	OAX-CAU-CBB-CAT
8	G	303	HXG	OAX-CAU-CBB-CAT
8	K	303	HXG	OAX-CAU-CBB-CAT
7	C	309	P1O	O8-C19-O7-C7
7	G	309	P1O	O8-C19-O7-C7
7	K	309	P1O	O8-C19-O7-C7
6	B	301	PLC	C4'-C5'-C6'-C7'
7	C	309	P1O	C10-C11-C12-C13
7	G	309	P1O	C10-C11-C12-C13
7	K	309	P1O	C10-C11-C12-C13
7	B	307	P1O	C12-C13-C14-C15
7	F	309	P1O	C12-C13-C14-C15
7	J	301	P1O	C12-C13-C14-C15
7	B	307	P1O	C22-C23-C24-C25
7	C	308	P1O	C22-C23-C24-C25
7	F	304	P1O	C20-C21-C22-C23
7	G	308	P1O	C22-C23-C24-C25
7	K	308	P1O	C22-C23-C24-C25
7	B	302	P1O	C20-C21-C22-C23
7	J	301	P1O	C22-C23-C24-C25
7	J	303	P1O	C20-C21-C22-C23

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Mol	Chain	Res	Type	Atoms
7	F	309	P1O	C22-C23-C24-C25
6	C	307	PLC	C4'-C5'-C6'-C7'
6	G	307	PLC	C4'-C5'-C6'-C7'
6	K	307	PLC	C4'-C5'-C6'-C7'
7	B	307	P1O	C6-C7-O7-C19
7	F	309	P1O	C6-C7-O7-C19
7	J	301	P1O	C6-C7-O7-C19
5	J	305	D10	C5-C6-C7-C8
5	B	305	D10	C4-C5-C6-C7
5	B	306	D10	C4-C5-C6-C7
5	F	307	D10	C4-C5-C6-C7
5	F	308	D10	C4-C5-C6-C7
5	J	306	D10	C4-C5-C6-C7
5	J	307	D10	C4-C5-C6-C7
5	B	304	D10	C5-C6-C7-C8
5	F	306	D10	C5-C6-C7-C8
7	B	307	P1O	C15-C16-C17-C18
7	F	309	P1O	C15-C16-C17-C18
7	J	301	P1O	C15-C16-C17-C18
5	B	306	D10	C6-C7-C8-C9
5	F	308	D10	C6-C7-C8-C9
5	J	307	D10	C6-C7-C8-C9
6	C	302	PLC	O2-C2-C3-O3
6	G	302	PLC	O2-C2-C3-O3
6	K	302	PLC	O2-C2-C3-O3
6	B	308	PLC	C6B-C7B-C8B-C9B
6	F	301	PLC	C6B-C7B-C8B-C9B
6	J	308	PLC	C6B-C7B-C8B-C9B
5	F	307	D10	C2-C3-C4-C5
5	J	306	D10	C2-C3-C4-C5
5	B	305	D10	C2-C3-C4-C5
5	J	307	D10	C2-C3-C4-C5
5	B	306	D10	C2-C3-C4-C5
5	F	308	D10	C2-C3-C4-C5
6	C	307	PLC	C7'-C8'-C9'-CA'
6	G	307	PLC	C7'-C8'-C9'-CA'
6	K	307	PLC	C7'-C8'-C9'-CA'
6	B	308	PLC	C7B-C8B-C9B-CAA
6	F	301	PLC	C7B-C8B-C9B-CAA
6	J	308	PLC	C7B-C8B-C9B-CAA
5	B	305	D10	C6-C7-C8-C9
7	C	309	P1O	O4-C6-C7-C8

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Mol	Chain	Res	Type	Atoms
7	G	309	P1O	O4-C6-C7-C8
7	K	309	P1O	O4-C6-C7-C8
5	F	307	D10	C6-C7-C8-C9
5	J	306	D10	C6-C7-C8-C9
6	C	302	PLC	C8B-C9B-CAA-CBA
6	G	302	PLC	C8B-C9B-CAA-CBA
6	K	302	PLC	C8B-C9B-CAA-CBA
6	B	301	PLC	C1-C2-C3-O3
6	F	303	PLC	C1-C2-C3-O3
6	J	302	PLC	C1-C2-C3-O3
6	B	309	PLC	C'-C1'-C2'-C3'
6	F	302	PLC	C'-C1'-C2'-C3'
6	C	307	PLC	C1B-C2B-C3B-C4B
6	K	307	PLC	C1B-C2B-C3B-C4B
6	G	307	PLC	C1B-C2B-C3B-C4B
7	B	302	P1O	C12-C13-C14-C15
7	F	304	P1O	C12-C13-C14-C15
7	J	303	P1O	C12-C13-C14-C15
6	J	309	PLC	C'-C1'-C2'-C3'
6	K	302	PLC	C8'-C9'-CA'-CB'
6	C	302	PLC	C8'-C9'-CA'-CB'
6	G	302	PLC	C8'-C9'-CA'-CB'
6	J	309	PLC	C5B-C6B-C7B-C8B
6	B	309	PLC	C5B-C6B-C7B-C8B
6	F	302	PLC	C5B-C6B-C7B-C8B
7	C	309	P1O	C19-C20-C21-C22
7	G	309	P1O	C19-C20-C21-C22
7	K	309	P1O	C19-C20-C21-C22
7	C	309	P1O	C15-C16-C17-C18
7	G	309	P1O	C15-C16-C17-C18
7	K	309	P1O	C15-C16-C17-C18
5	B	305	D10	C3-C4-C5-C6
5	J	306	D10	C3-C4-C5-C6
6	K	302	PLC	C3'-C4'-C5'-C6'
6	C	302	PLC	C3'-C4'-C5'-C6'
5	F	307	D10	C3-C4-C5-C6
6	C	302	PLC	C1B-C2B-C3B-C4B
6	G	302	PLC	C3'-C4'-C5'-C6'
6	G	302	PLC	C1B-C2B-C3B-C4B
6	K	302	PLC	C1B-C2B-C3B-C4B
6	B	308	PLC	O3P-C1-C2-C3
6	F	301	PLC	O3P-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
6	J	308	PLC	O3P-C1-C2-C3
6	B	308	PLC	C3-C2-O2-C'
6	B	309	PLC	C1-C2-O2-C'
6	F	301	PLC	C3-C2-O2-C'
6	F	302	PLC	C1-C2-O2-C'
6	J	308	PLC	C3-C2-O2-C'
6	J	309	PLC	C1-C2-O2-C'
8	C	306	HXG	CAT-CBB-OAY-CBA
8	G	306	HXG	CAT-CBB-OAY-CBA
8	K	306	HXG	CAT-CBB-OAY-CBA
6	B	308	PLC	O3P-C1-C2-O2
6	F	301	PLC	O3P-C1-C2-O2
6	J	308	PLC	O3P-C1-C2-O2
7	C	309	P1O	O4-C6-C7-O7
7	G	309	P1O	O4-C6-C7-O7
7	K	309	P1O	O4-C6-C7-O7
8	C	303	HXG	OAX-CAU-CBB-OAY
8	G	303	HXG	OAX-CAU-CBB-OAY
8	K	303	HXG	OAX-CAU-CBB-OAY
6	B	309	PLC	C5-C4-O4P-P
6	C	307	PLC	C5-C4-O4P-P
6	F	302	PLC	C5-C4-O4P-P
6	J	309	PLC	C5-C4-O4P-P
6	G	307	PLC	C5-C4-O4P-P
6	K	307	PLC	C5-C4-O4P-P
7	C	309	P1O	C22-C23-C24-C25
7	G	309	P1O	C14-C15-C16-C17
7	K	309	P1O	C14-C15-C16-C17
7	K	309	P1O	C22-C23-C24-C25
7	C	309	P1O	O7-C7-C8-O5
7	G	309	P1O	O7-C7-C8-O5
7	K	309	P1O	O7-C7-C8-O5
7	C	309	P1O	C14-C15-C16-C17
7	G	309	P1O	C22-C23-C24-C25
7	C	308	P1O	C11-C12-C13-C14
7	G	308	P1O	C11-C12-C13-C14
7	K	308	P1O	C11-C12-C13-C14
7	B	302	P1O	C1-C2-N1-C5
7	F	304	P1O	C1-C2-N1-C5
6	B	309	PLC	O4P-C4-C5-N
6	F	302	PLC	O4P-C4-C5-N
6	J	309	PLC	O4P-C4-C5-N

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Mol	Chain	Res	Type	Atoms
7	F	304	P1O	C24-C25-C26-C27
7	J	303	P1O	C24-C25-C26-C27
7	F	304	P1O	C9-C10-C11-C12
7	B	302	P1O	C24-C25-C26-C27
7	B	302	P1O	C1-C2-N1-C4
7	F	304	P1O	C1-C2-N1-C4
7	J	303	P1O	C1-C2-N1-C4
7	J	303	P1O	C1-C2-N1-C5
7	B	302	P1O	C9-C10-C11-C12
7	J	303	P1O	C9-C10-C11-C12
7	B	302	P1O	C15-C16-C17-C18
7	J	303	P1O	C15-C16-C17-C18
7	F	304	P1O	C15-C16-C17-C18
6	K	305	PLC	C7B-C8B-C9B-CAA
6	G	305	PLC	C7B-C8B-C9B-CAA
6	C	305	PLC	C7B-C8B-C9B-CAA
7	B	307	P1O	C21-C22-C23-C24
7	F	304	P1O	C10-C11-C12-C13
7	F	309	P1O	C21-C22-C23-C24
7	J	301	P1O	C21-C22-C23-C24
6	B	301	PLC	O2-C2-C3-O3
6	F	303	PLC	O2-C2-C3-O3
6	J	302	PLC	O2-C2-C3-O3
5	B	306	D10	C3-C4-C5-C6
5	F	308	D10	C3-C4-C5-C6
5	J	307	D10	C3-C4-C5-C6
7	B	302	P1O	C10-C11-C12-C13
7	J	303	P1O	C10-C11-C12-C13
6	C	302	PLC	C1-C2-C3-O3
6	G	302	PLC	C1-C2-C3-O3
6	K	302	PLC	C1-C2-C3-O3
6	J	308	PLC	C3'-C4'-C5'-C6'
6	B	301	PLC	C1-O3P-P-O1P
6	B	308	PLC	C4-O4P-P-O2P
6	B	308	PLC	C4-O4P-P-O3P
6	B	309	PLC	C1-O3P-P-O2P
6	F	301	PLC	C4-O4P-P-O2P
6	F	301	PLC	C4-O4P-P-O3P
6	F	302	PLC	C1-O3P-P-O2P
6	F	303	PLC	C1-O3P-P-O1P
6	J	302	PLC	C1-O3P-P-O1P
6	J	308	PLC	C4-O4P-P-O2P

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Mol	Chain	Res	Type	Atoms
6	J	308	PLC	C4-O4P-P-O3P
6	J	309	PLC	C1-O3P-P-O2P
7	B	302	P1O	C6-O4-P1-O2
7	B	307	P1O	C1-O3-P1-O1
7	C	308	P1O	C6-O4-P1-O2
7	C	309	P1O	C1-O3-P1-O4
7	F	304	P1O	C6-O4-P1-O2
7	F	309	P1O	C1-O3-P1-O1
7	J	301	P1O	C1-O3-P1-O1
7	J	303	P1O	C6-O4-P1-O2
7	G	308	P1O	C6-O4-P1-O2
7	G	309	P1O	C1-O3-P1-O4
7	K	308	P1O	C6-O4-P1-O2
7	K	309	P1O	C1-O3-P1-O4
8	C	303	HXG	CAP-OAW-PBD-OAI
8	C	306	HXG	CAU-OAX-PBD-OAI
8	G	303	HXG	CAP-OAW-PBD-OAI
8	G	306	HXG	CAU-OAX-PBD-OAI
8	K	303	HXG	CAP-OAW-PBD-OAI
8	K	306	HXG	CAU-OAX-PBD-OAI
6	F	301	PLC	C3'-C4'-C5'-C6'
6	B	308	PLC	C3'-C4'-C5'-C6'
6	G	305	PLC	C3B-C4B-C5B-C6B
6	K	305	PLC	C3B-C4B-C5B-C6B
6	C	305	PLC	C3B-C4B-C5B-C6B
7	K	308	P1O	C24-C25-C26-C27
7	C	308	P1O	C24-C25-C26-C27
7	G	308	P1O	C24-C25-C26-C27
6	G	307	PLC	C'-C1'-C2'-C3'
6	K	307	PLC	C'-C1'-C2'-C3'
6	C	307	PLC	C'-C1'-C2'-C3'
6	C	307	PLC	C2-C1-O3P-P
6	G	307	PLC	C2-C1-O3P-P
6	K	307	PLC	C2-C1-O3P-P
6	K	302	PLC	C5B-C6B-C7B-C8B
6	C	302	PLC	C5B-C6B-C7B-C8B
6	G	302	PLC	C5B-C6B-C7B-C8B
7	F	304	P1O	O7-C7-C8-O5
7	J	303	P1O	O7-C7-C8-O5
8	C	306	HXG	OAV-CAT-CBB-OAY
8	G	306	HXG	OAV-CAT-CBB-OAY
8	K	306	HXG	OAV-CAT-CBB-OAY

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Mol	Chain	Res	Type	Atoms
6	C	305	PLC	C5B-C6B-C7B-C8B
6	G	305	PLC	C5B-C6B-C7B-C8B
6	K	305	PLC	C5B-C6B-C7B-C8B
6	B	308	PLC	C5'-C6'-C7'-C8'
6	F	301	PLC	C5'-C6'-C7'-C8'
6	J	308	PLC	C5'-C6'-C7'-C8'
8	C	306	HXG	CBB-CAU-OAX-PBD
8	G	306	HXG	CBB-CAU-OAX-PBD
8	K	306	HXG	CBB-CAU-OAX-PBD
7	B	302	P1O	C1-C2-N1-C3
7	F	304	P1O	C1-C2-N1-C3
7	J	303	P1O	C1-C2-N1-C3
5	B	305	D10	C7-C8-C9-C10
7	C	309	P1O	C6-C7-C8-O5
7	G	309	P1O	C6-C7-C8-O5
7	K	309	P1O	C6-C7-C8-O5
5	F	307	D10	C7-C8-C9-C10
5	J	306	D10	C7-C8-C9-C10
6	C	307	PLC	C3-C2-O2-C'
6	G	307	PLC	C3-C2-O2-C'
6	K	307	PLC	C3-C2-O2-C'
7	C	309	P1O	C23-C24-C25-C26
7	K	309	P1O	C23-C24-C25-C26
7	G	309	P1O	C23-C24-C25-C26
6	G	305	PLC	C8B-C9B-CAA-CBA
6	C	305	PLC	C8B-C9B-CAA-CBA
7	B	302	P1O	O7-C7-C8-O5
6	K	305	PLC	C8B-C9B-CAA-CBA
6	K	307	PLC	C6B-C7B-C8B-C9B
6	G	307	PLC	C6B-C7B-C8B-C9B
6	C	307	PLC	C6B-C7B-C8B-C9B
6	J	309	PLC	C7B-C8B-C9B-CAA
6	B	309	PLC	C7B-C8B-C9B-CAA
6	F	302	PLC	C7B-C8B-C9B-CAA
7	C	309	P1O	C9-C10-C11-C12
7	G	309	P1O	C9-C10-C11-C12
7	K	309	P1O	C9-C10-C11-C12
5	F	306	D10	C7-C8-C9-C10
5	J	305	D10	C7-C8-C9-C10
5	B	304	D10	C7-C8-C9-C10
7	F	309	P1O	O7-C19-C20-C21
7	J	301	P1O	O7-C19-C20-C21

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Mol	Chain	Res	Type	Atoms
7	B	307	P1O	O7-C19-C20-C21
6	B	308	PLC	C2B-C1B-CB-O3
6	F	301	PLC	C2B-C1B-CB-O3
6	J	308	PLC	C2B-C1B-CB-O3
7	F	309	P1O	C25-C26-C27-C28
7	J	301	P1O	C25-C26-C27-C28
7	K	308	P1O	C10-C11-C12-C13
7	C	308	P1O	C10-C11-C12-C13
7	G	308	P1O	C10-C11-C12-C13
6	C	307	PLC	O2-C'-C1'-C2'
6	G	307	PLC	O2-C'-C1'-C2'
6	K	307	PLC	O2-C'-C1'-C2'
7	B	307	P1O	C25-C26-C27-C28
6	B	301	PLC	C1-C2-O2-C'
6	B	301	PLC	C3-C2-O2-C'
6	C	307	PLC	C1-C2-O2-C'
6	F	303	PLC	C1-C2-O2-C'
6	F	303	PLC	C3-C2-O2-C'
6	J	302	PLC	C1-C2-O2-C'
6	J	302	PLC	C3-C2-O2-C'
6	G	307	PLC	C1-C2-O2-C'
6	K	307	PLC	C1-C2-O2-C'
8	K	306	HXG	OAG-CBA-OAY-CBB
7	C	308	P1O	C7-C6-O4-P1
7	G	308	P1O	C7-C6-O4-P1
7	K	308	P1O	C7-C6-O4-P1
8	G	306	HXG	OAG-CBA-OAY-CBB
7	C	308	P1O	C11-C10-C9-O5
7	G	308	P1O	C11-C10-C9-O5
7	K	308	P1O	C11-C10-C9-O5
6	J	302	PLC	C6'-C7'-C8'-C9'
6	C	307	PLC	O'-C'-C1'-C2'
6	G	307	PLC	O'-C'-C1'-C2'
6	K	307	PLC	O'-C'-C1'-C2'
8	C	306	HXG	OAG-CBA-OAY-CBB
6	F	303	PLC	C6'-C7'-C8'-C9'
6	B	301	PLC	C6'-C7'-C8'-C9'
7	B	302	P1O	C6-C7-C8-O5
7	F	304	P1O	C6-C7-C8-O5
7	J	303	P1O	C6-C7-C8-O5
6	B	308	PLC	C2B-C1B-CB-OB
6	F	301	PLC	C2B-C1B-CB-OB

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Mol	Chain	Res	Type	Atoms
6	J	308	PLC	C2B-C1B-CB-OB
7	B	307	P1O	O4-C6-C7-C8
7	F	309	P1O	O4-C6-C7-C8
7	J	301	P1O	O4-C6-C7-C8
7	G	308	P1O	C11-C10-C9-O6
7	B	307	P1O	O8-C19-C20-C21
7	C	308	P1O	C11-C10-C9-O6
7	F	309	P1O	O8-C19-C20-C21
7	K	308	P1O	C11-C10-C9-O6
6	B	308	PLC	O2-C'-C1'-C2'
6	C	305	PLC	C2B-C1B-CB-O3
6	F	301	PLC	O2-C'-C1'-C2'
6	J	308	PLC	O2-C'-C1'-C2'
6	G	305	PLC	C2B-C1B-CB-O3
7	B	302	P1O	C11-C10-C9-O5
7	F	304	P1O	C11-C10-C9-O5
7	J	303	P1O	C11-C10-C9-O5
7	J	301	P1O	O8-C19-C20-C21
6	K	305	PLC	C2B-C1B-CB-O3

There are no ring outliers.

45 monomers are involved in 343 short contacts:

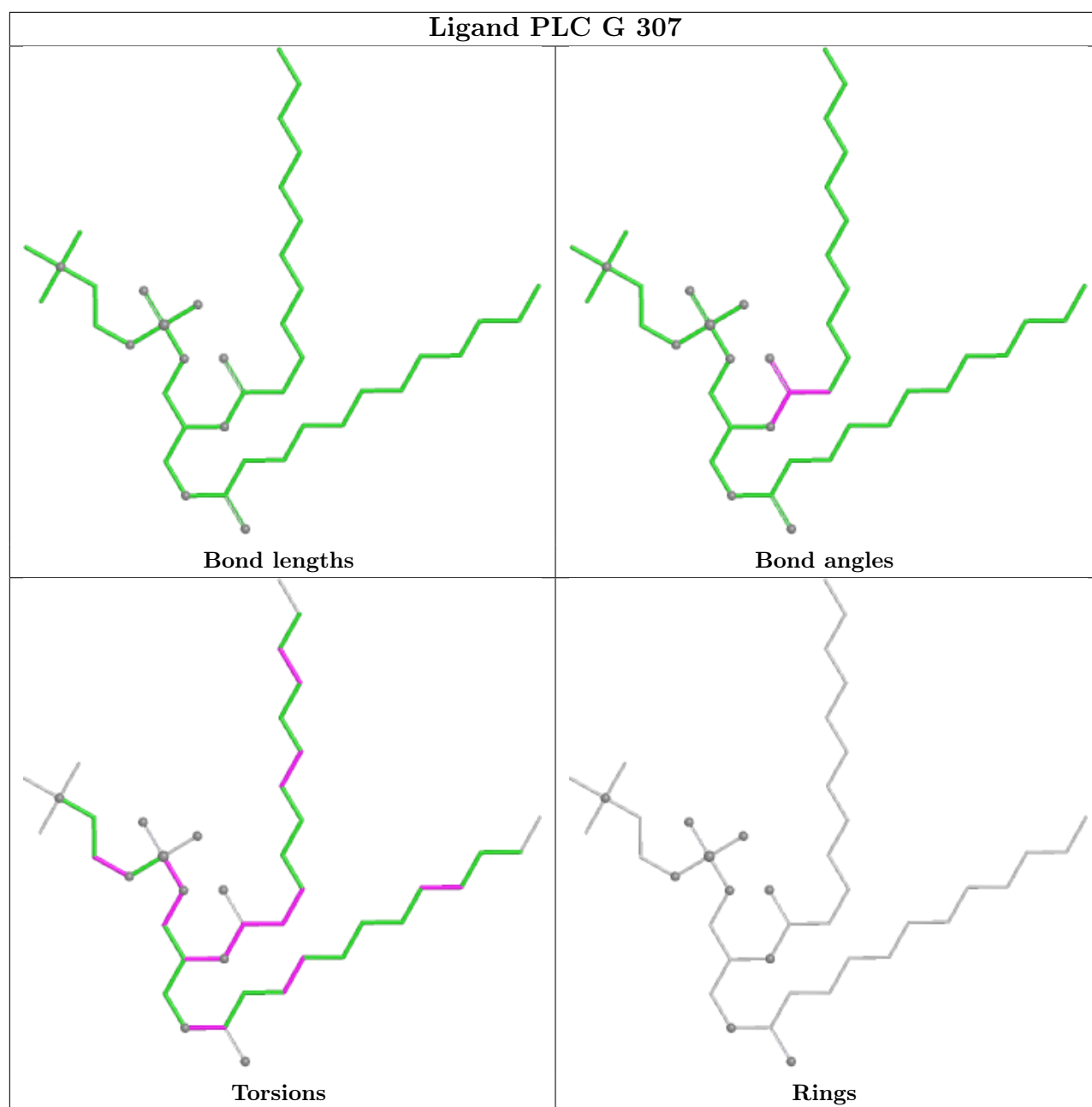
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	G	307	PLC	5	0
7	C	309	P1O	14	0
6	J	308	PLC	7	0
7	J	303	P1O	8	0
5	B	305	D10	6	0
6	G	305	PLC	2	0
6	G	302	PLC	6	0
7	G	309	P1O	15	0
7	F	304	P1O	9	0
6	B	301	PLC	3	0
6	K	302	PLC	5	0
6	K	307	PLC	5	0
7	J	301	P1O	28	0
8	K	303	HXG	15	0
5	F	306	D10	1	0
7	C	308	P1O	7	0
6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

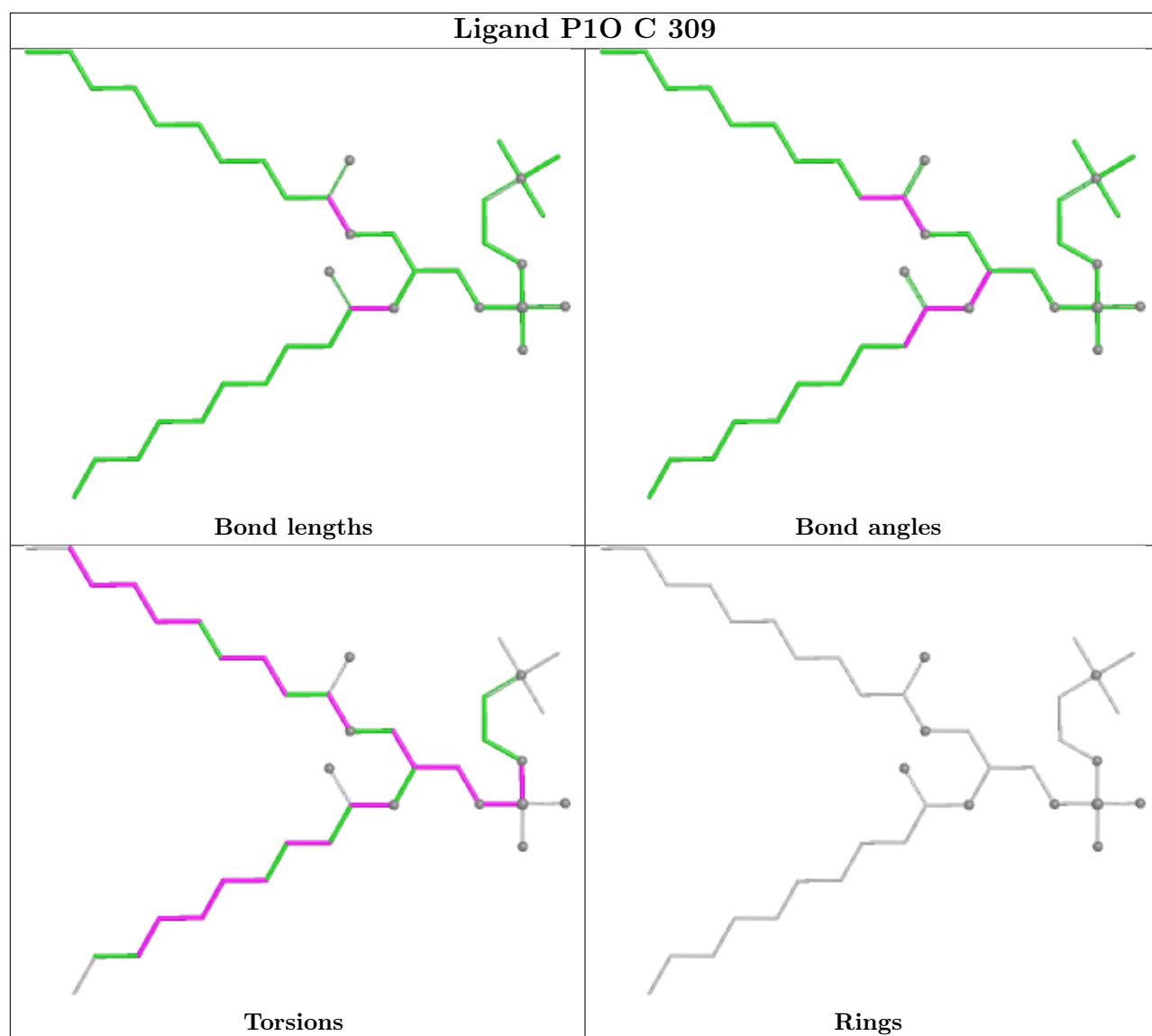
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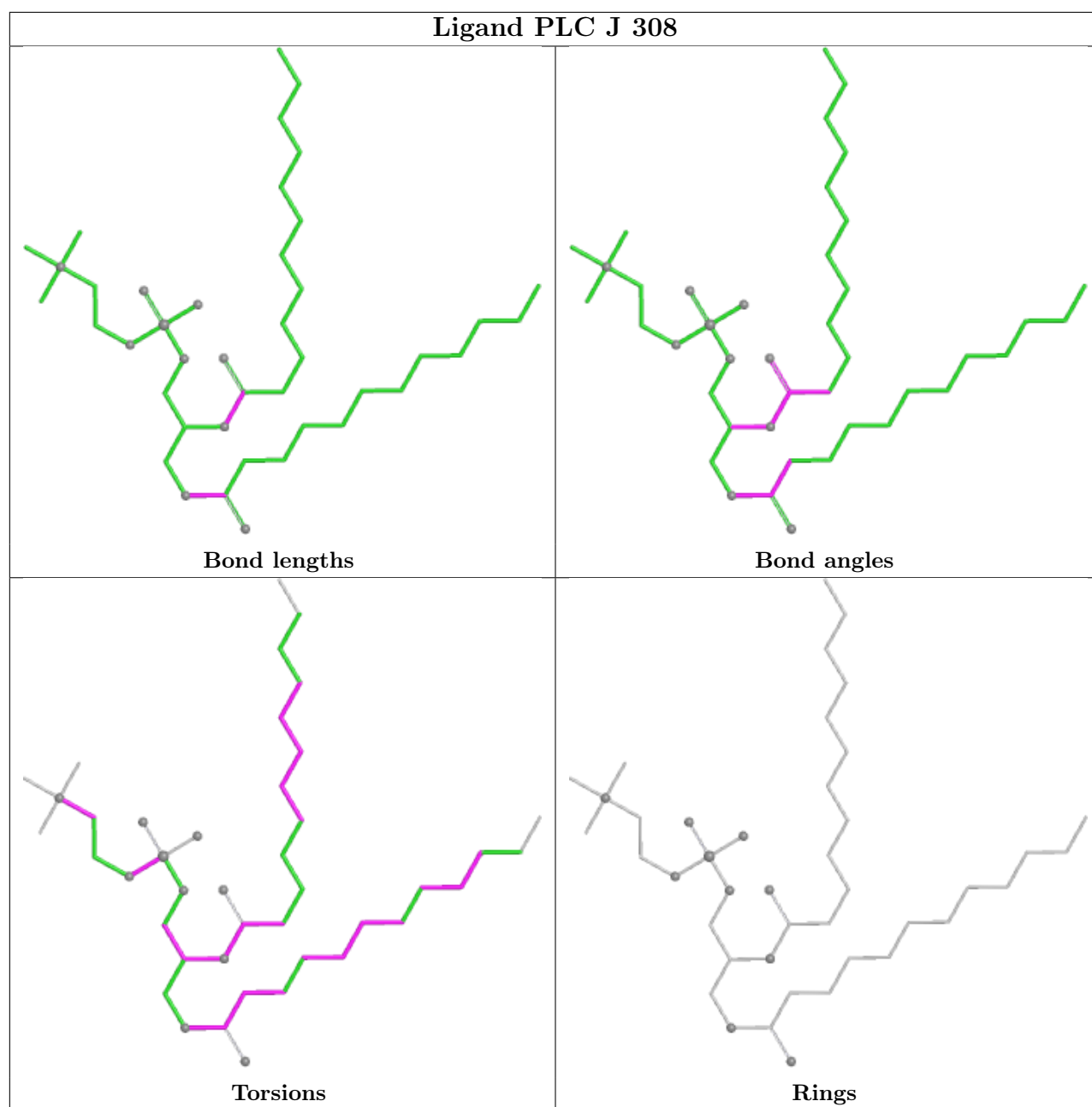
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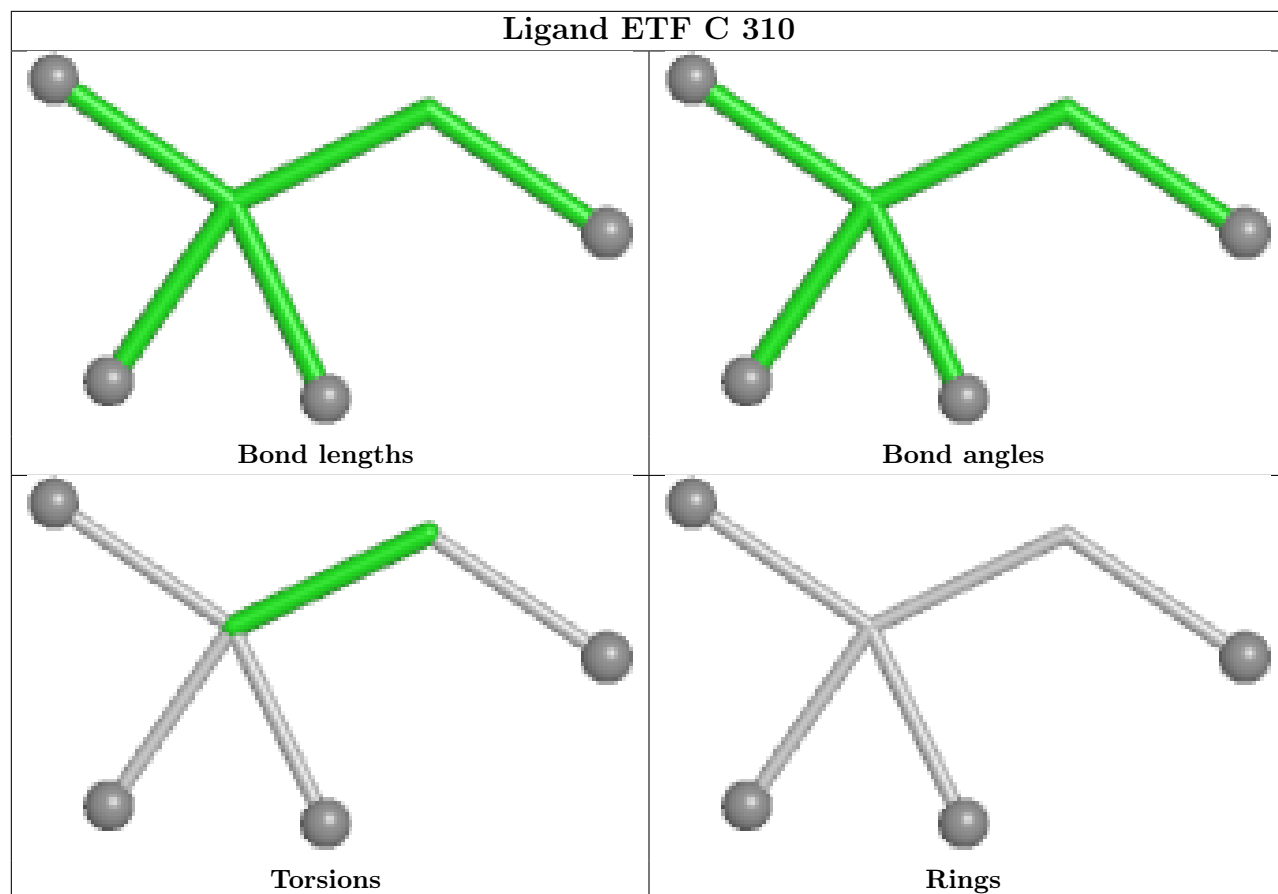
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	C	303	HXG	14	0
7	B	307	P1O	28	0
5	J	307	D10	1	0
6	J	302	PLC	3	0
8	K	306	HXG	16	0
6	C	302	PLC	4	0
5	J	306	D10	6	0
8	C	306	HXG	15	0
5	K	304	D10	1	0
5	J	305	D10	1	0
5	C	304	D10	1	0
6	F	301	PLC	6	0
5	B	304	D10	1	0
7	K	308	P1O	7	0
7	F	309	P1O	29	0
7	G	308	P1O	7	0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

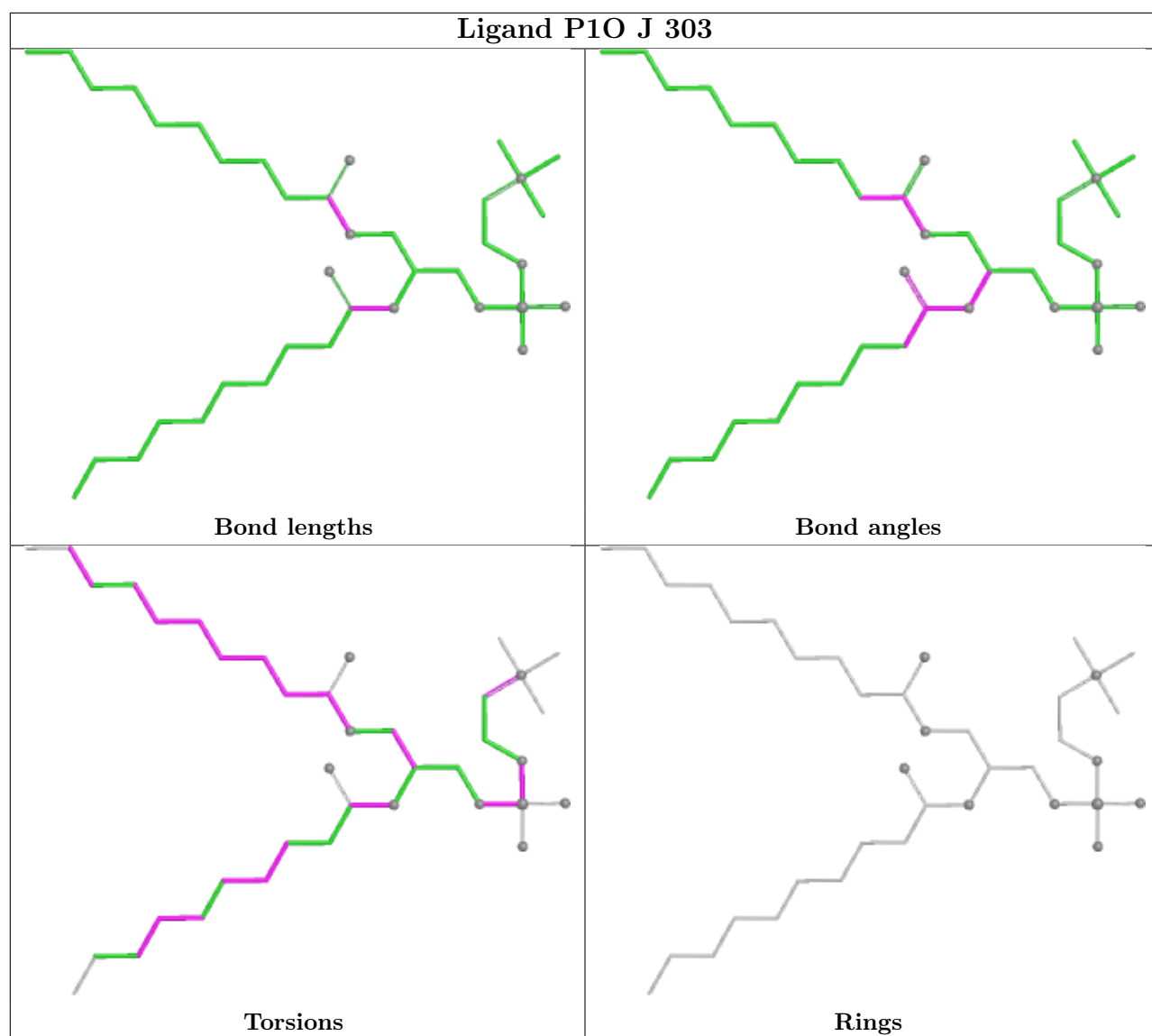
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

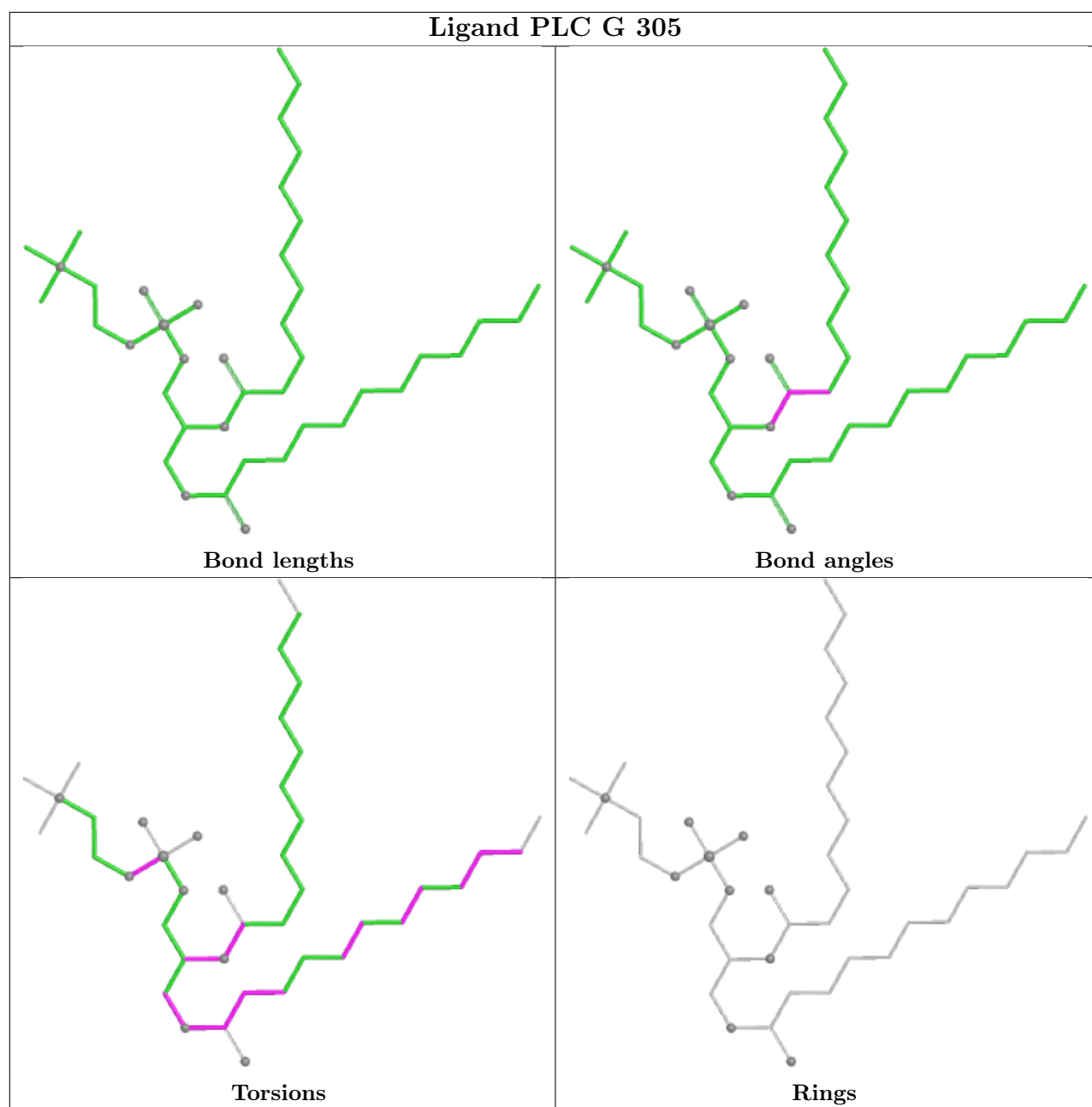


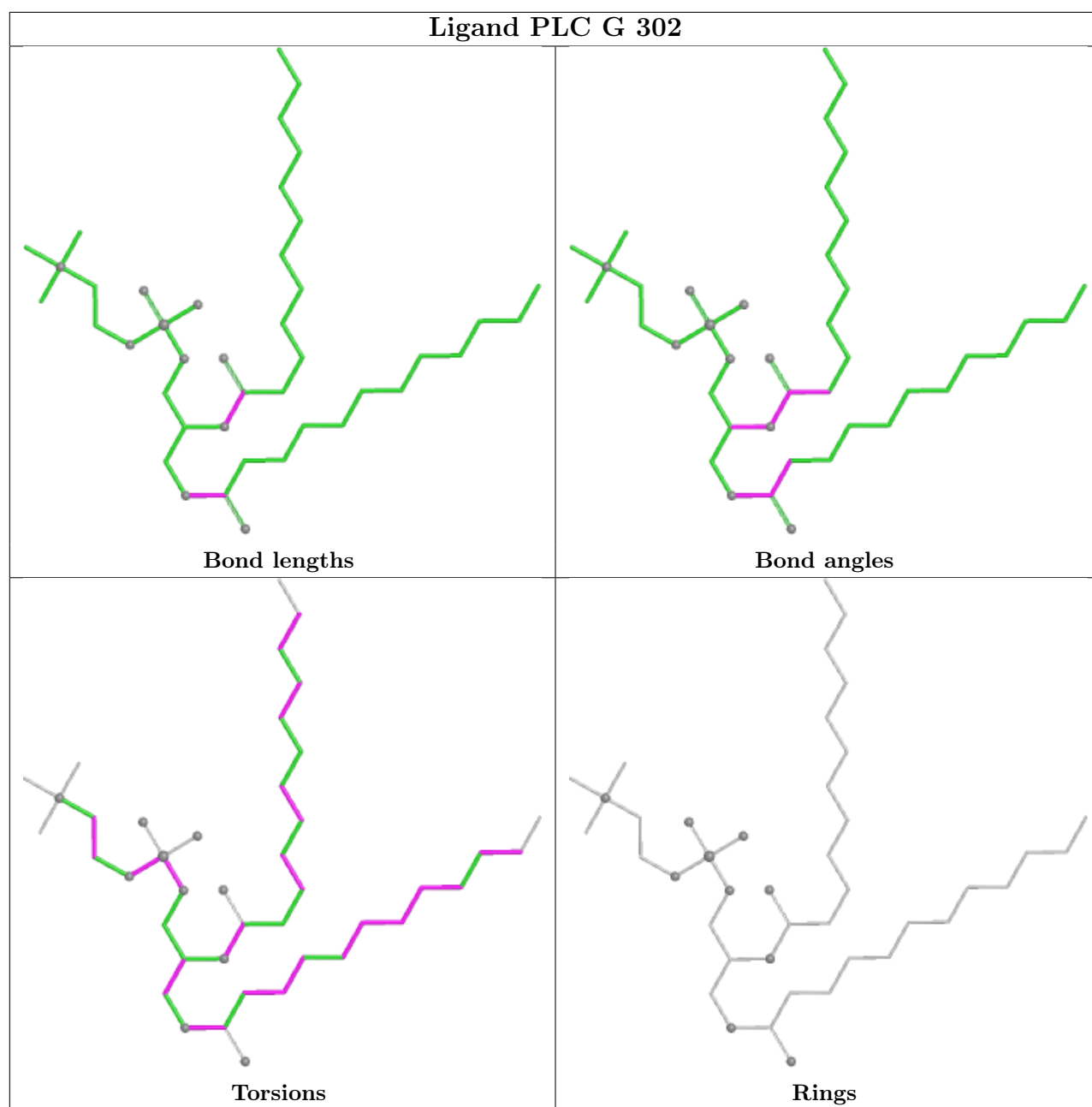


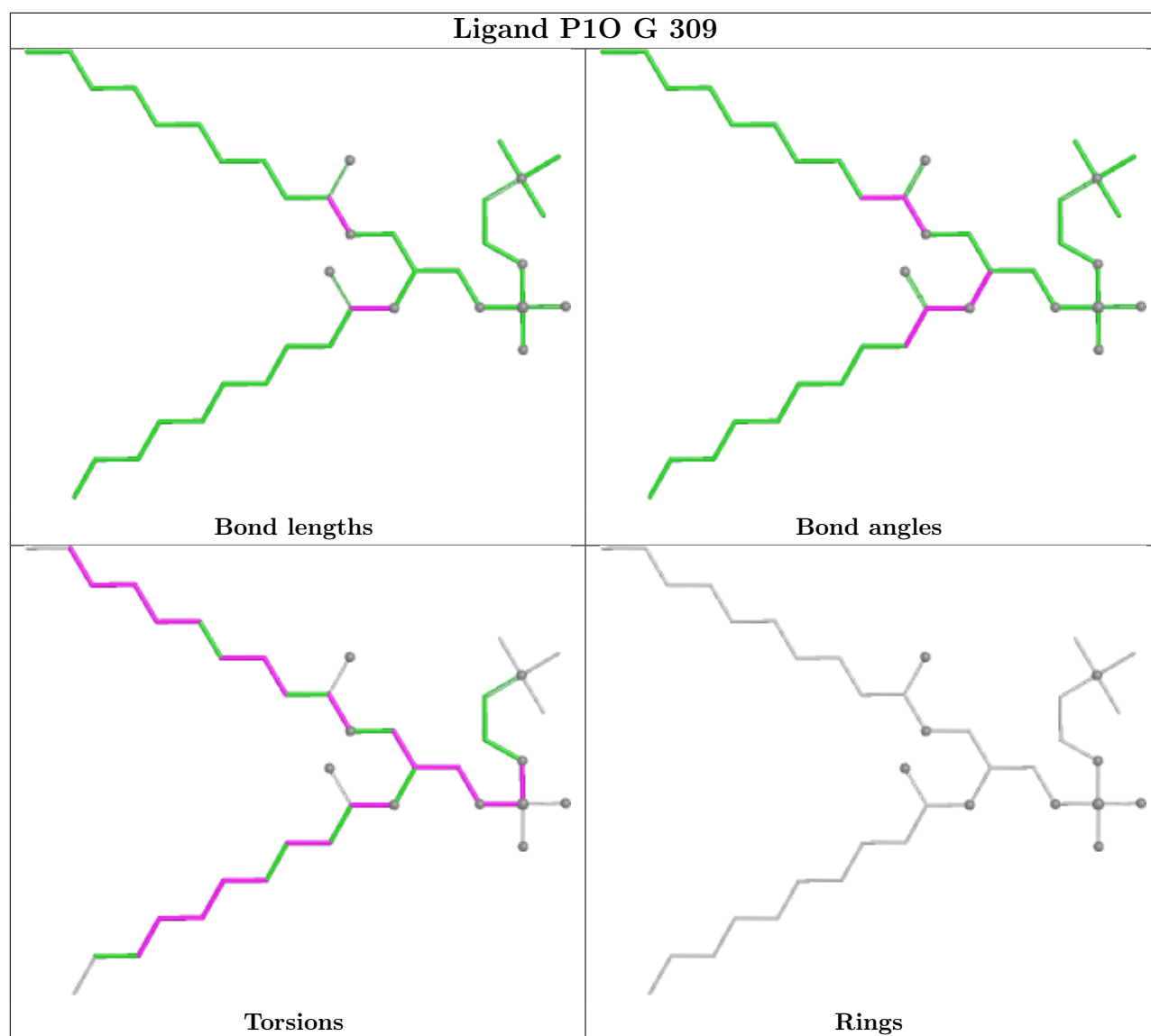


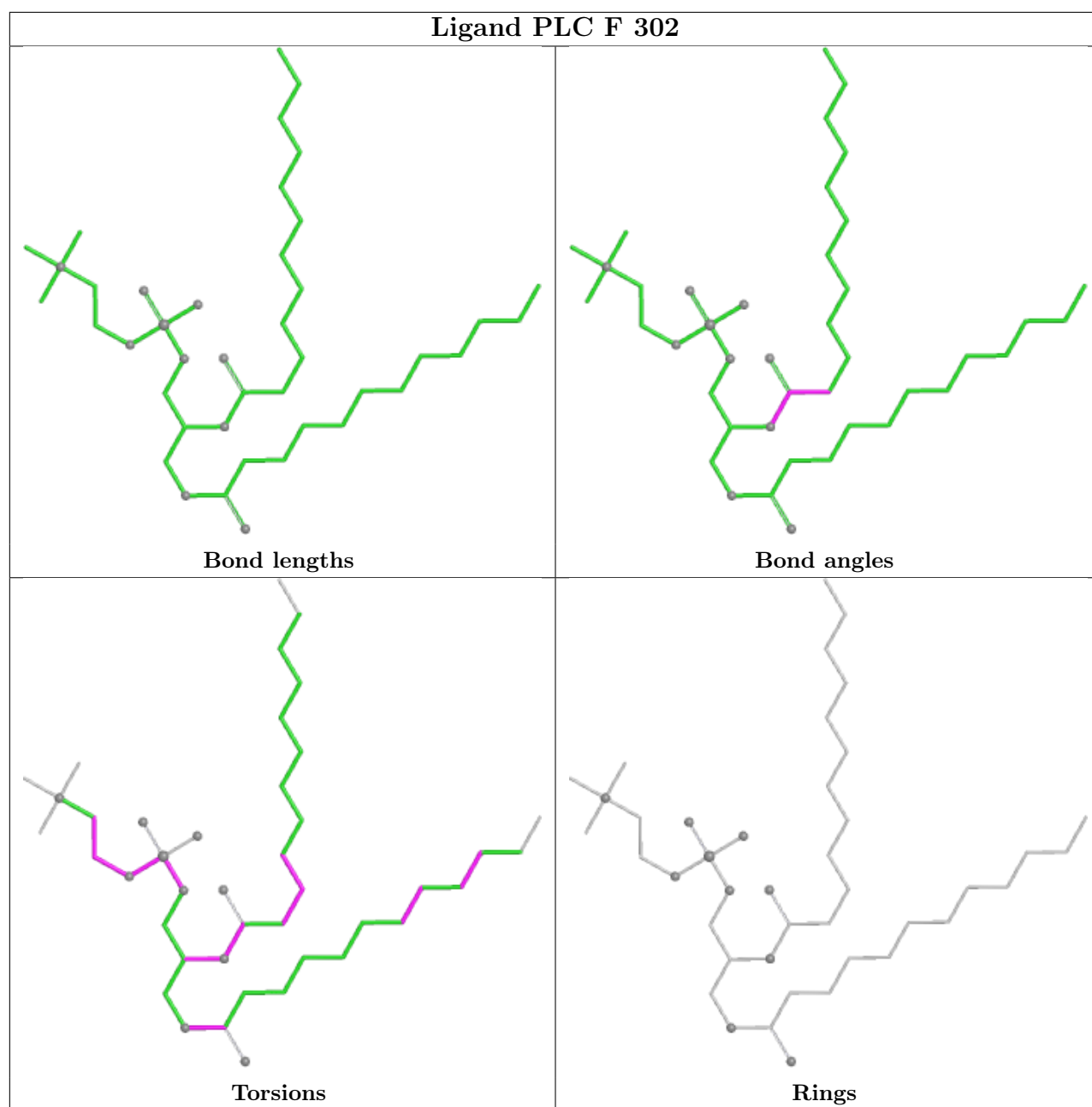


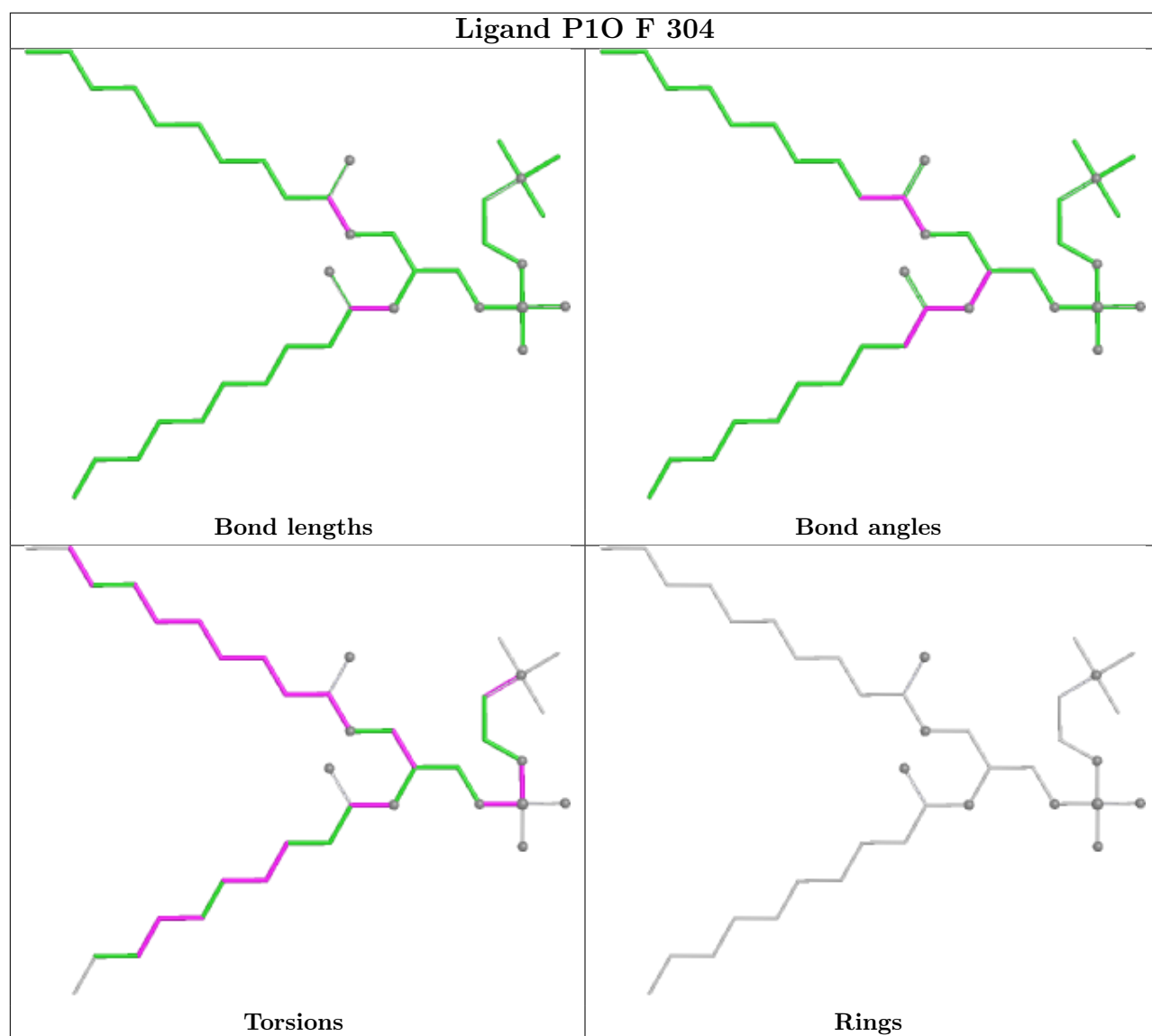


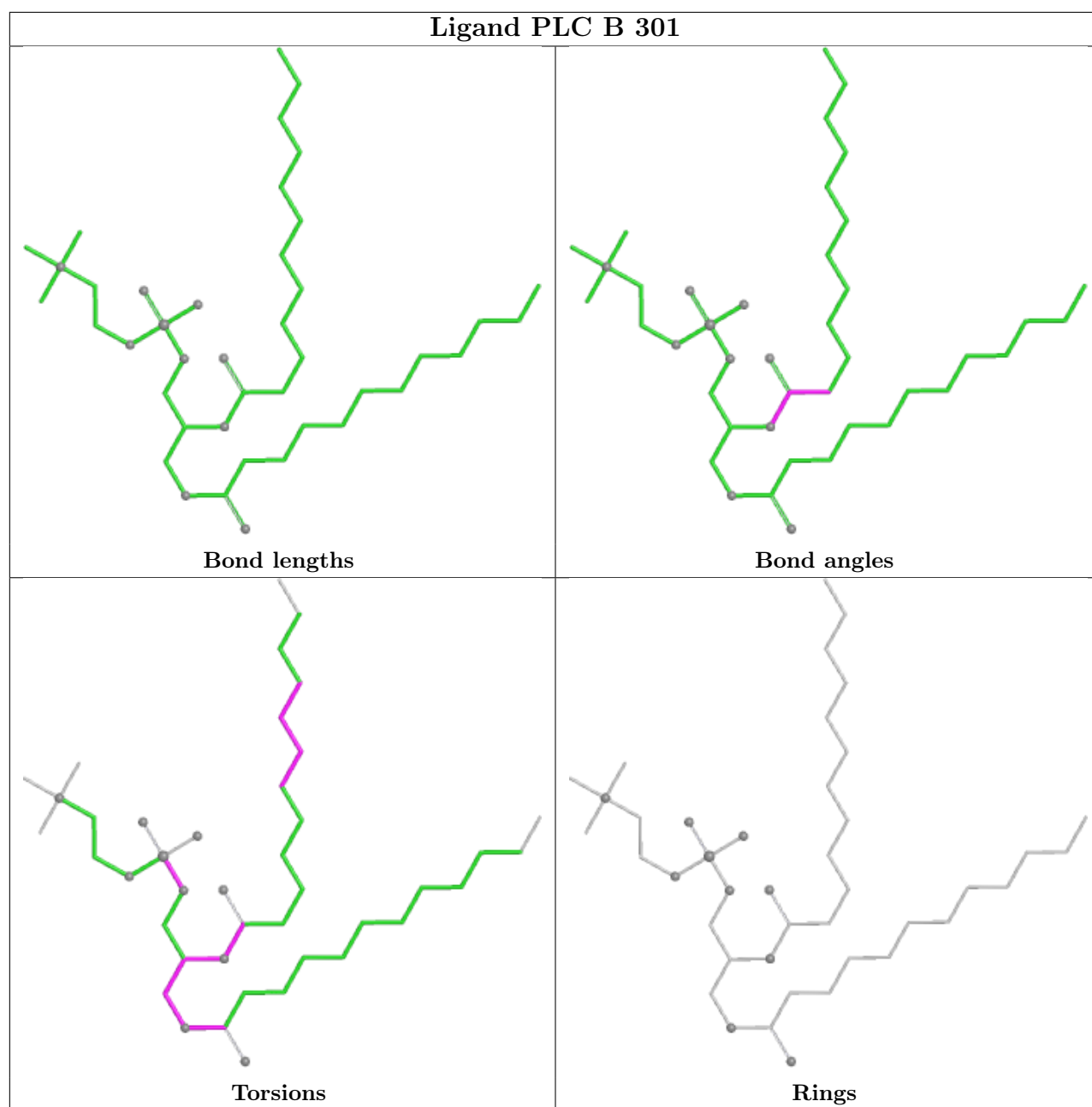


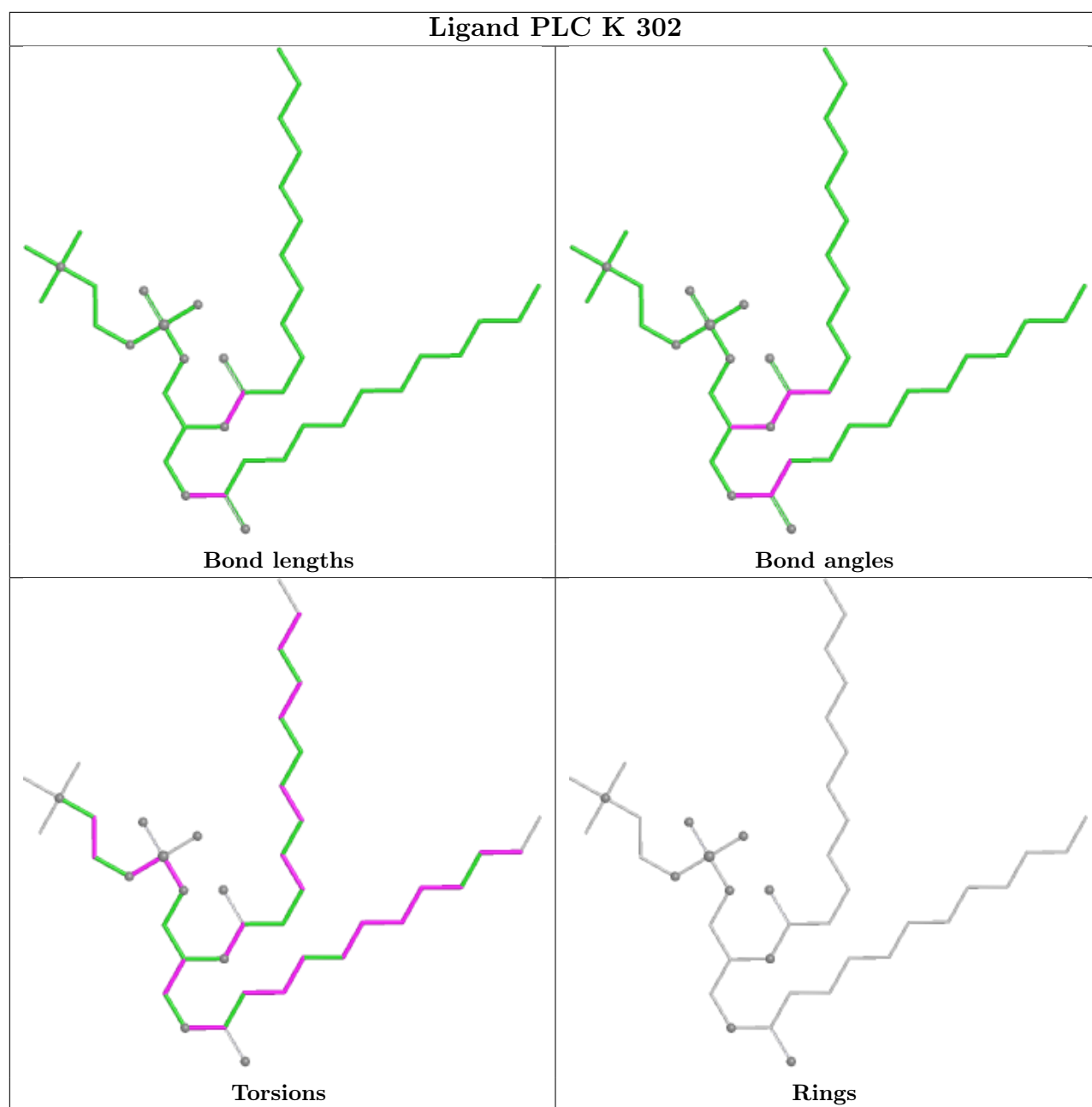


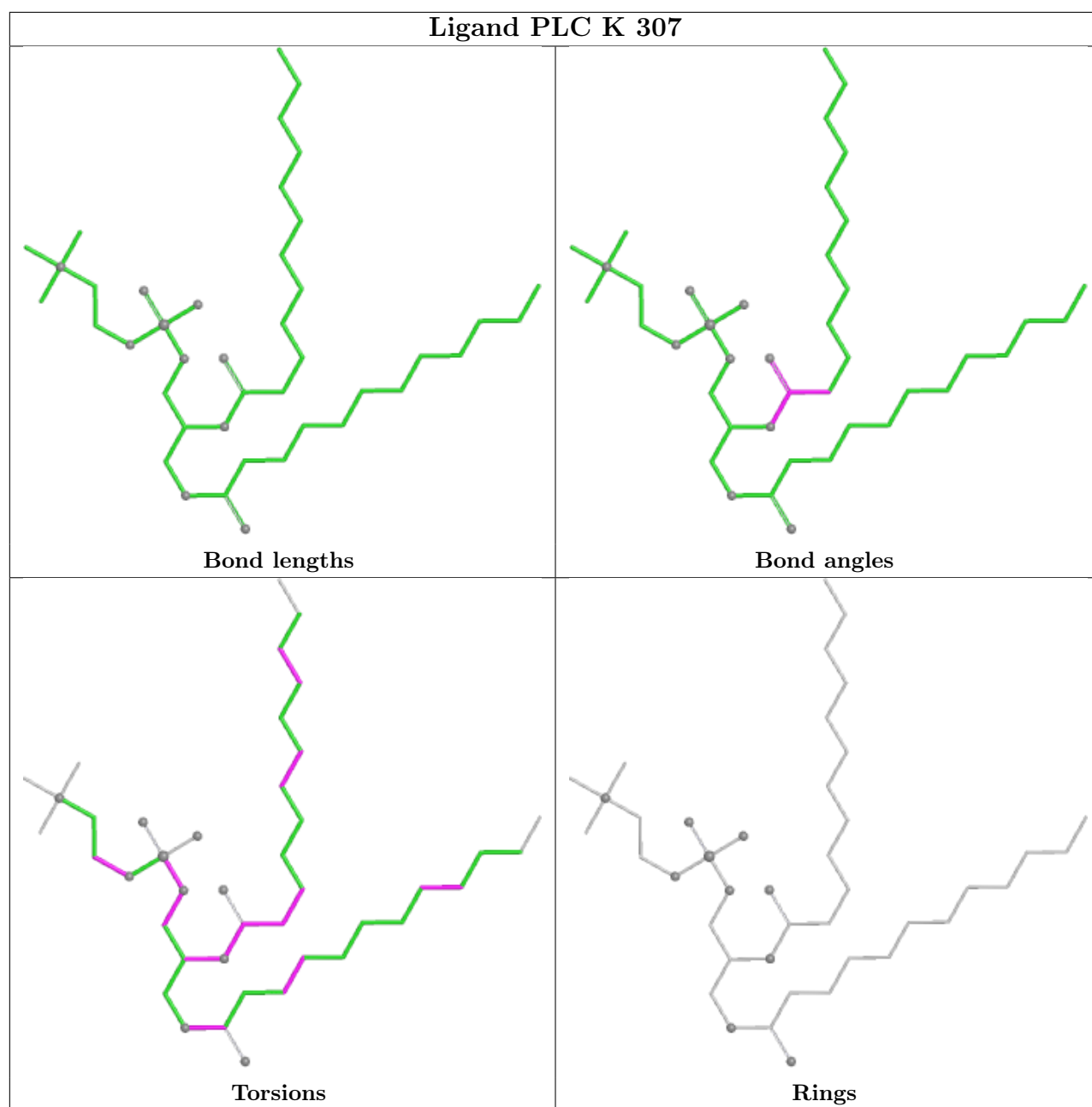


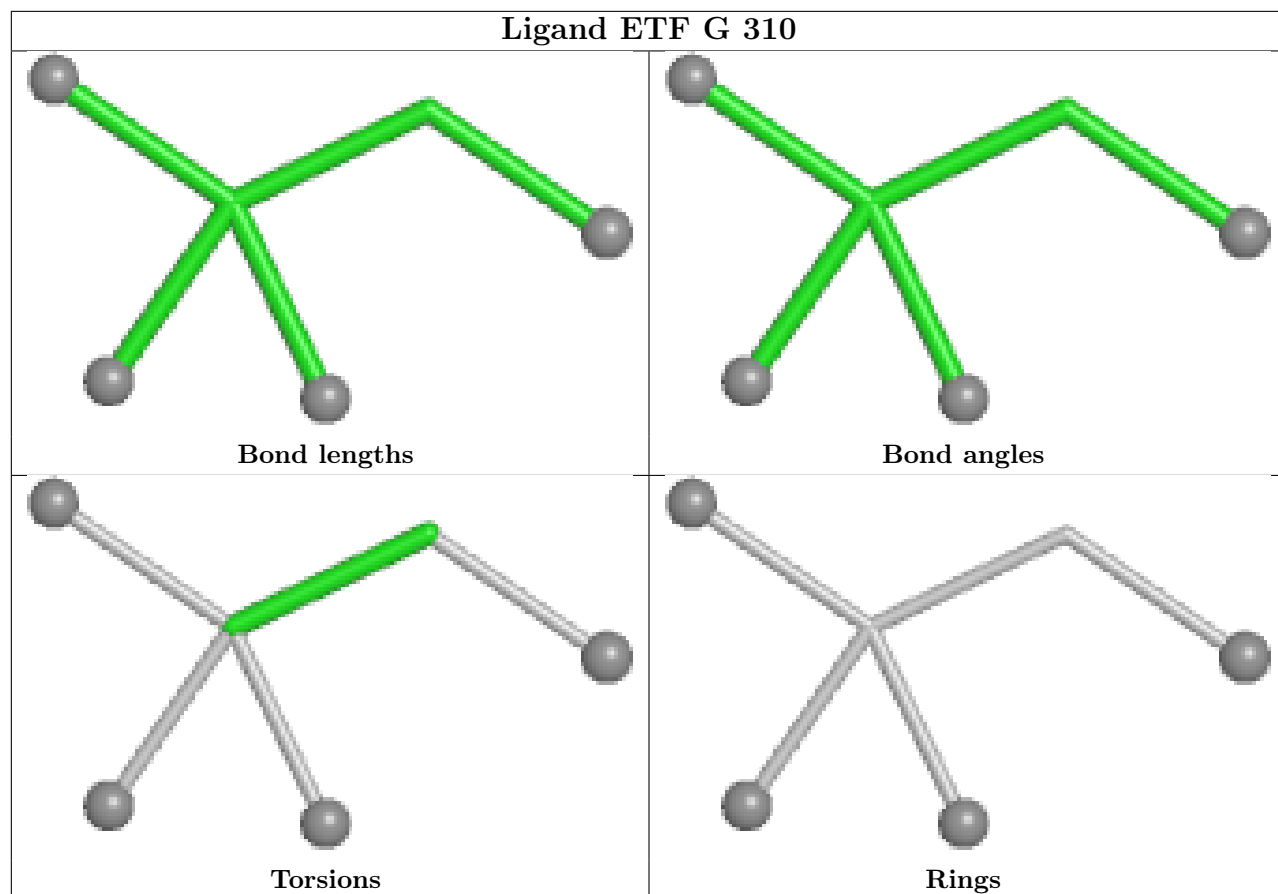


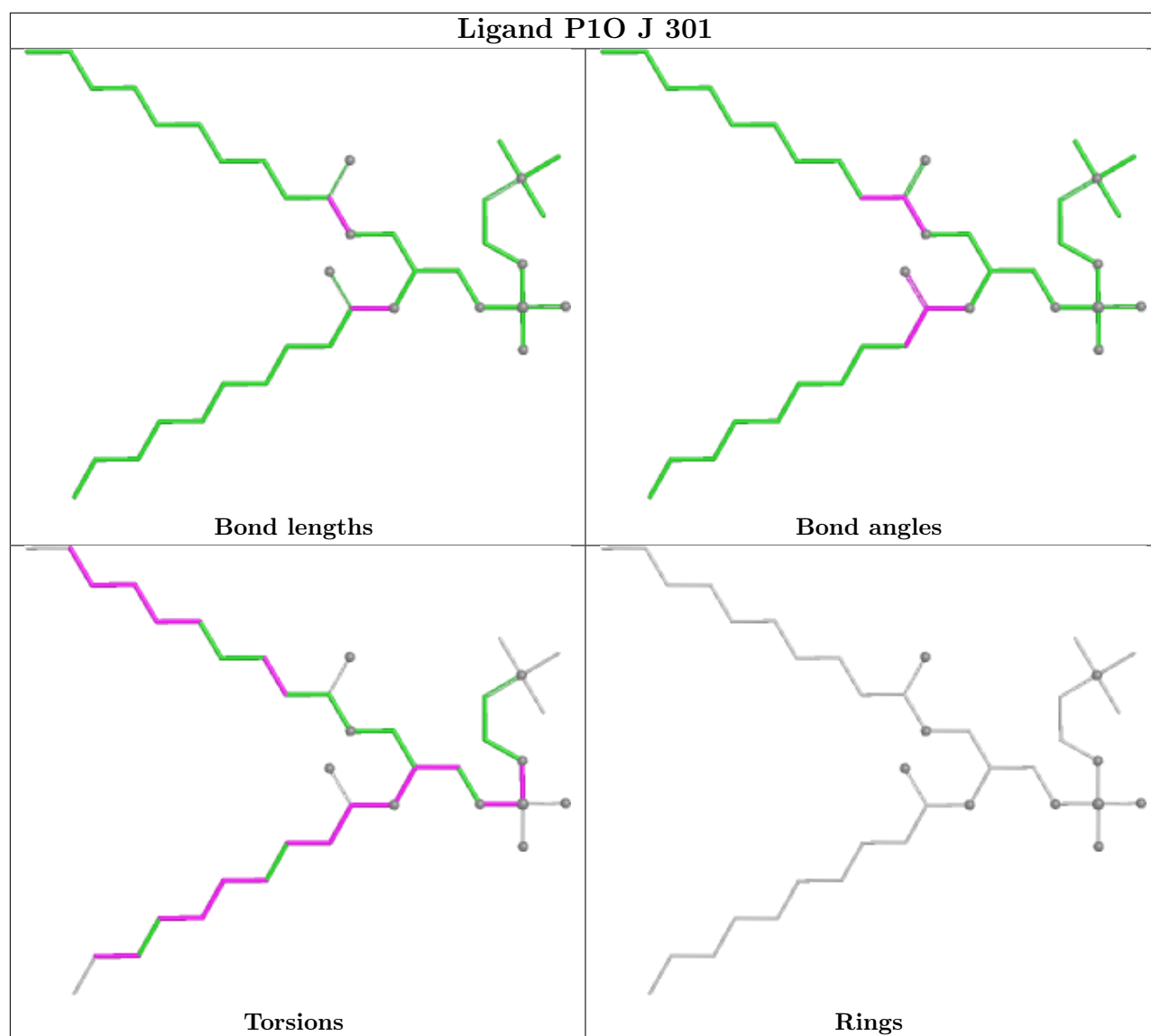


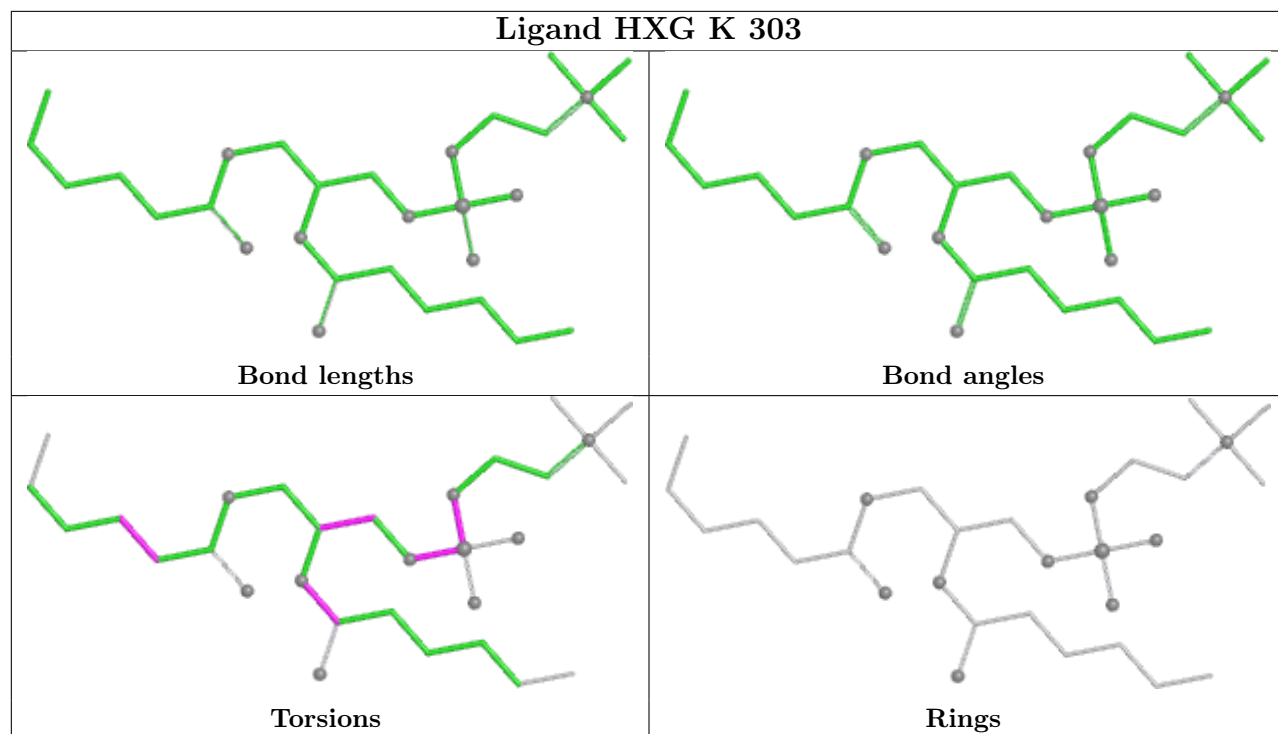


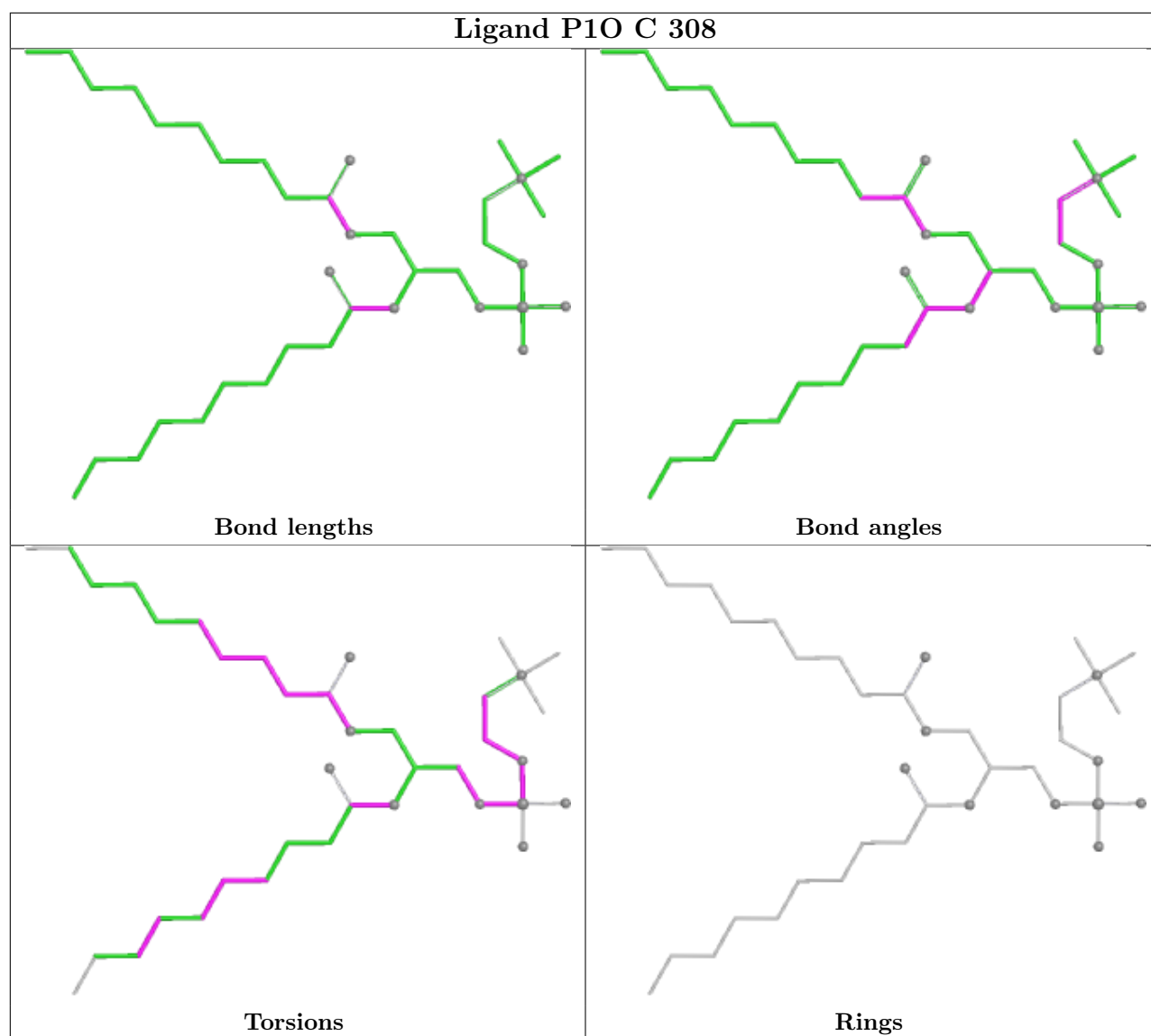


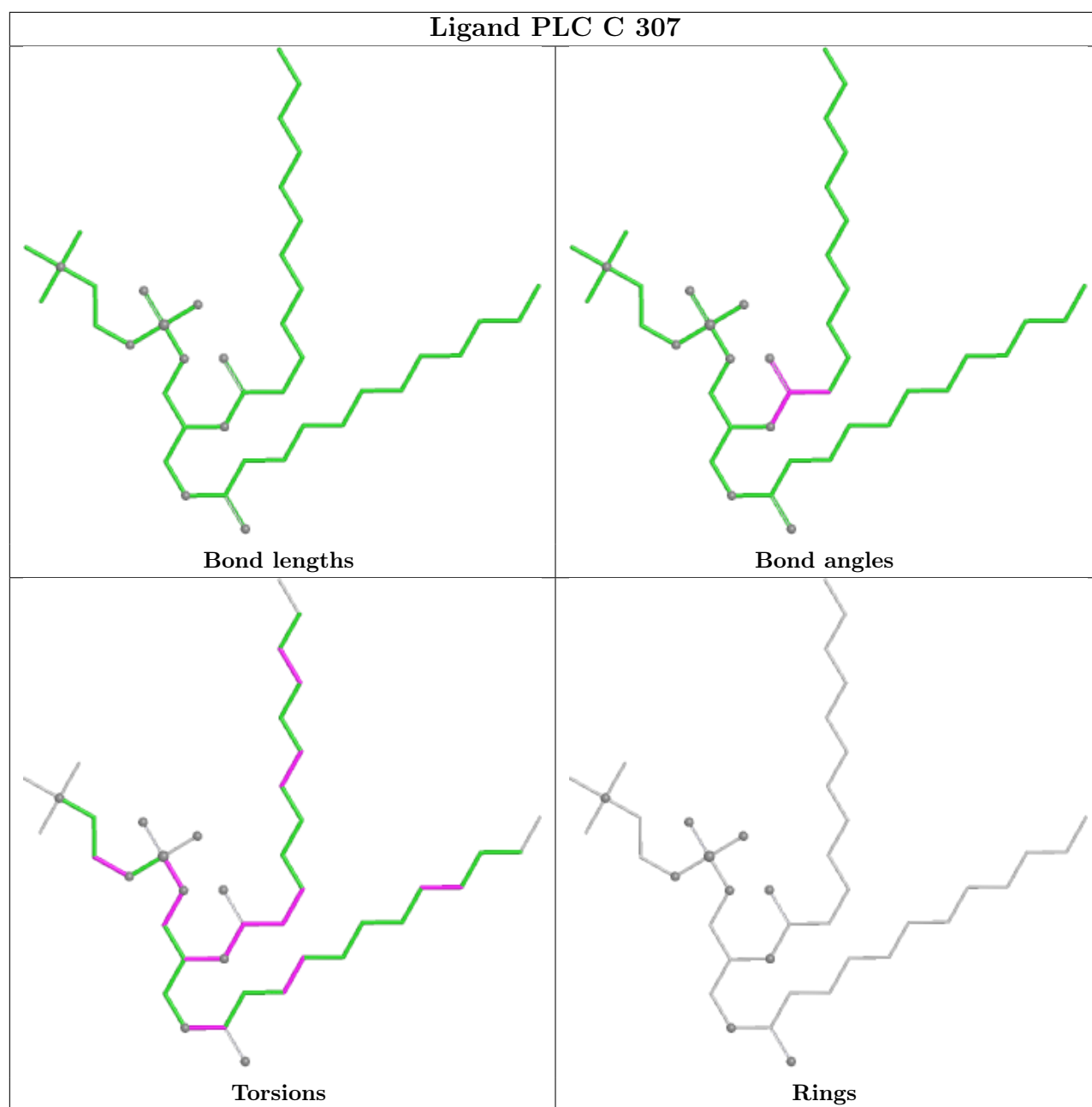


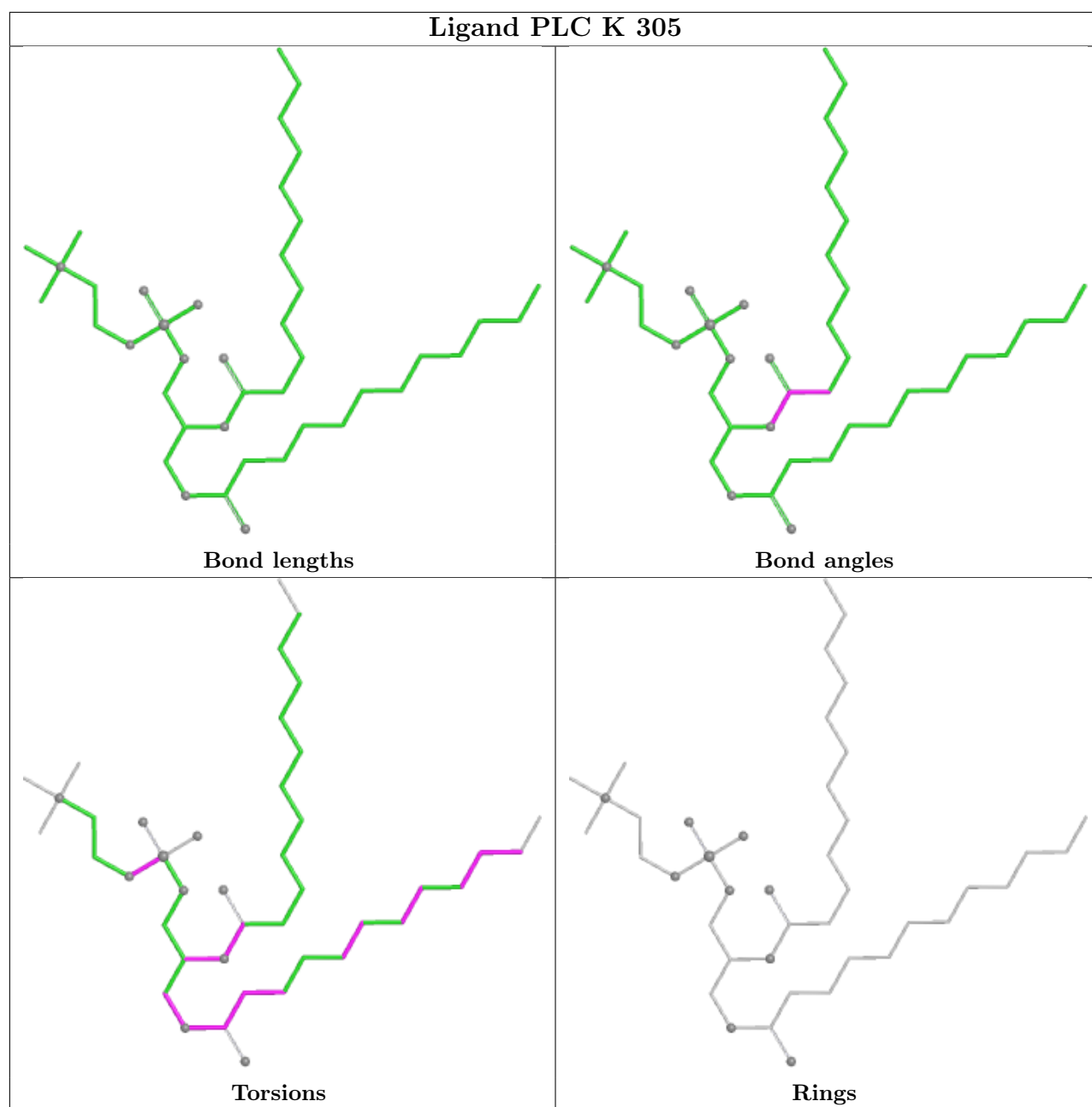


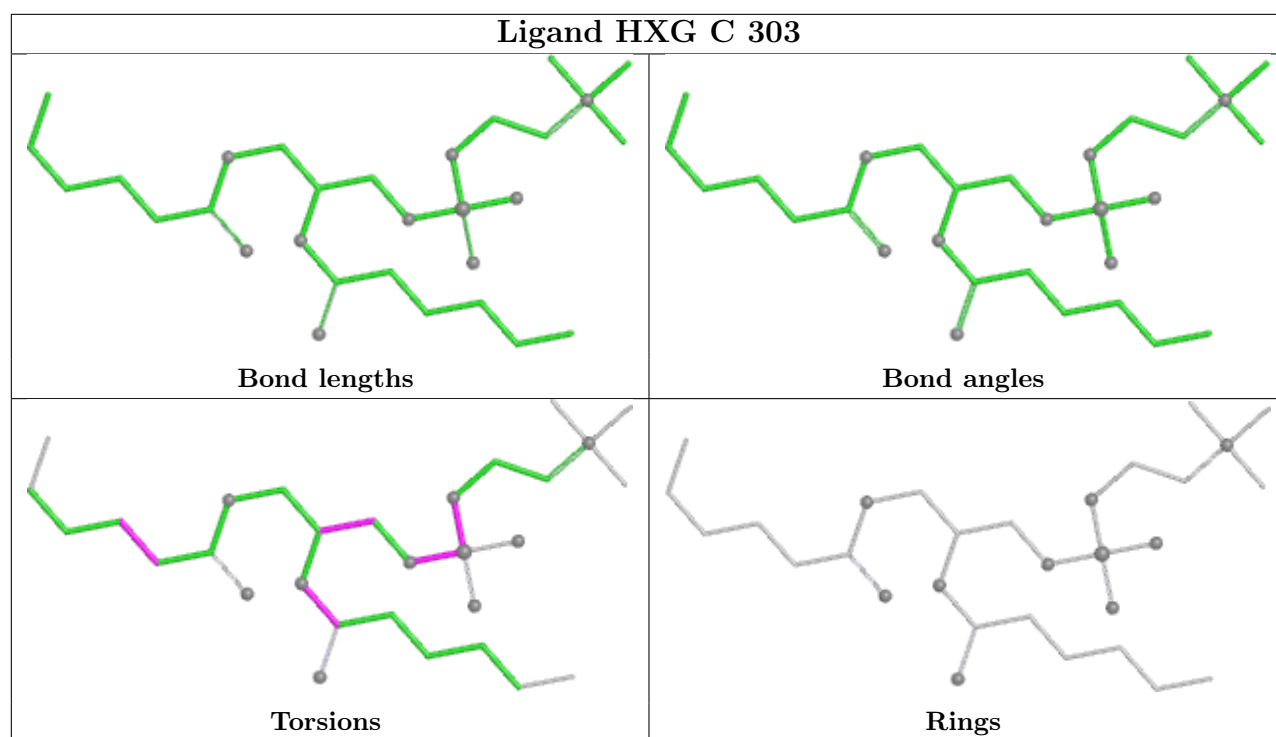


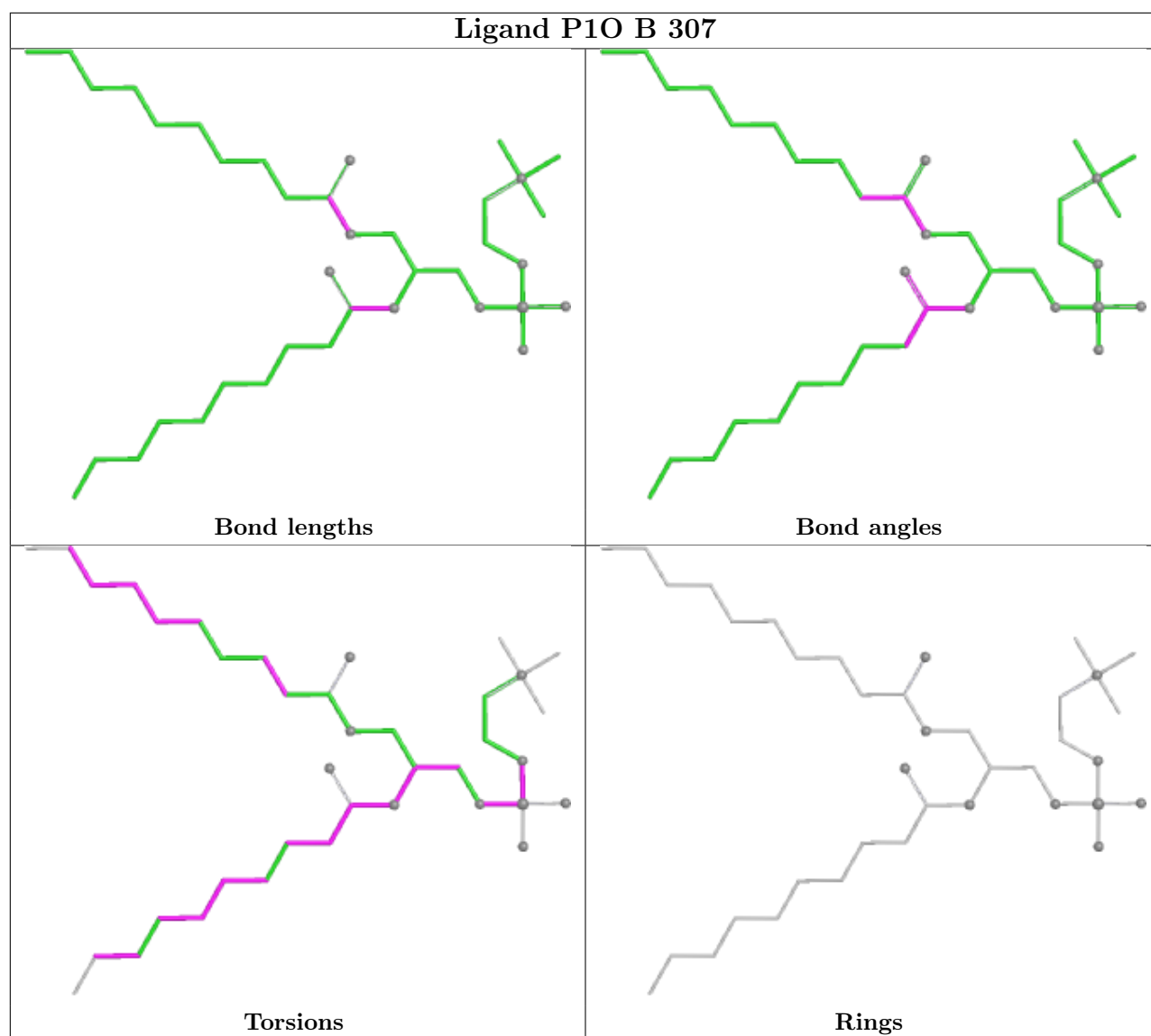


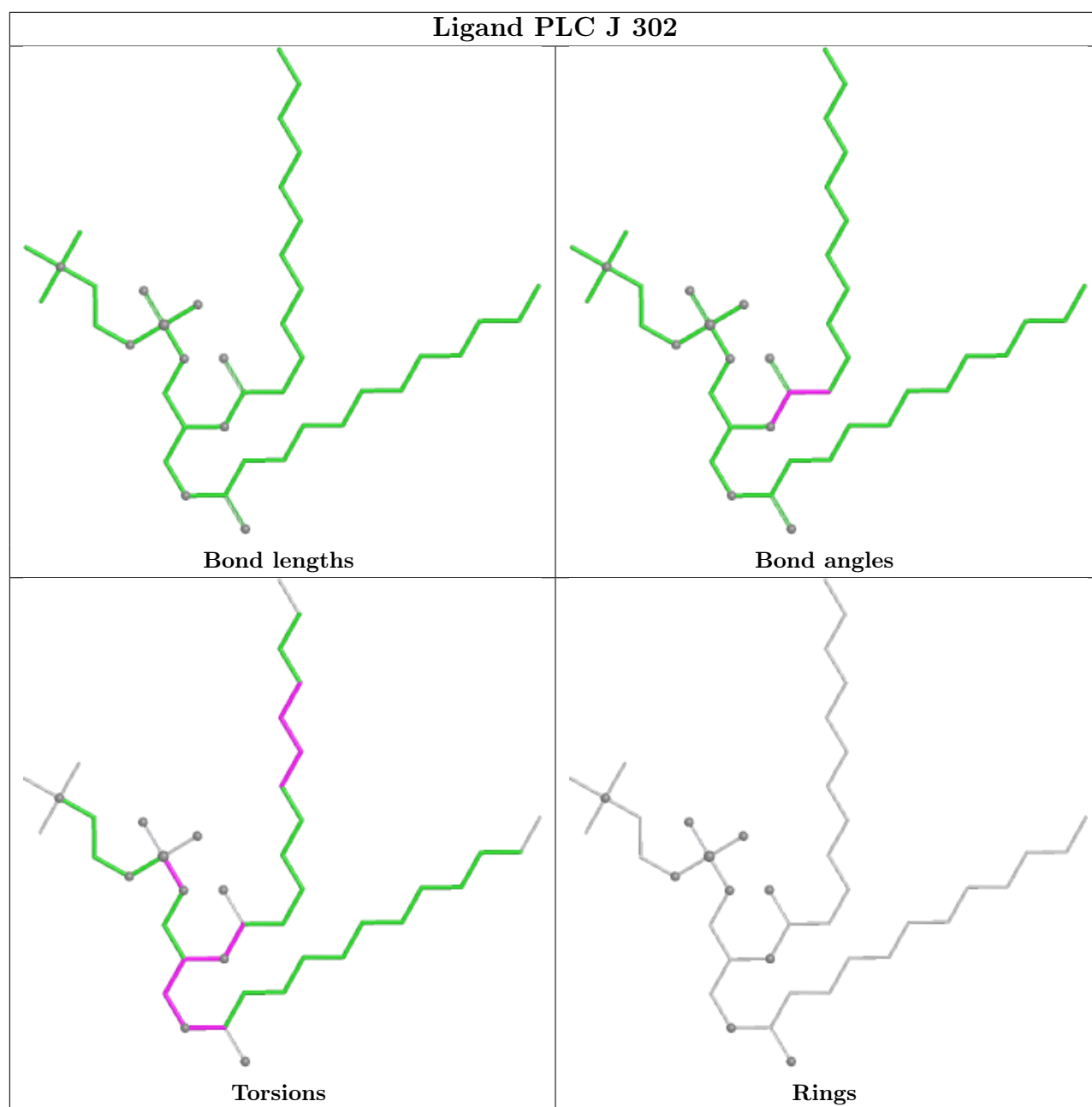


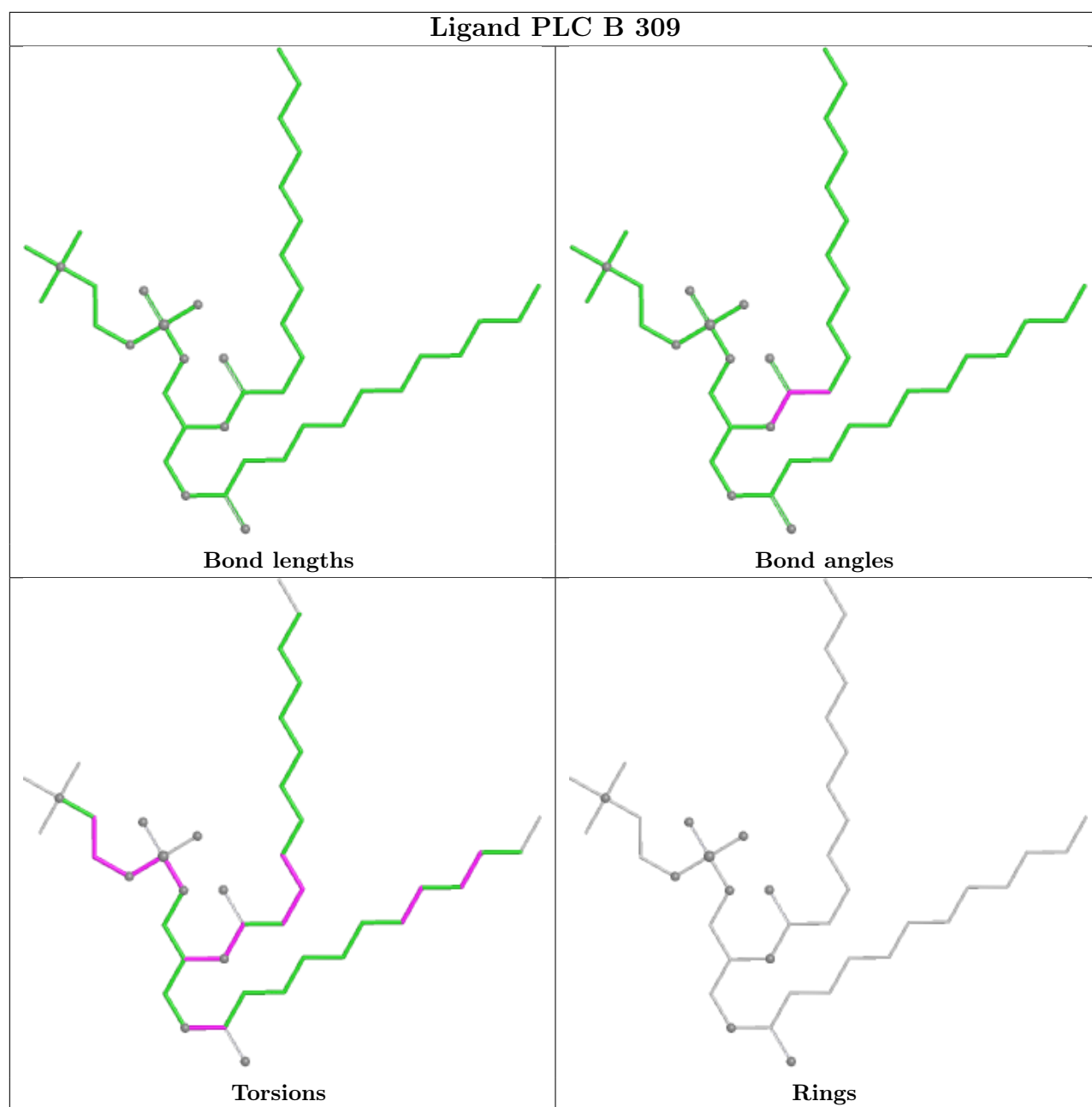


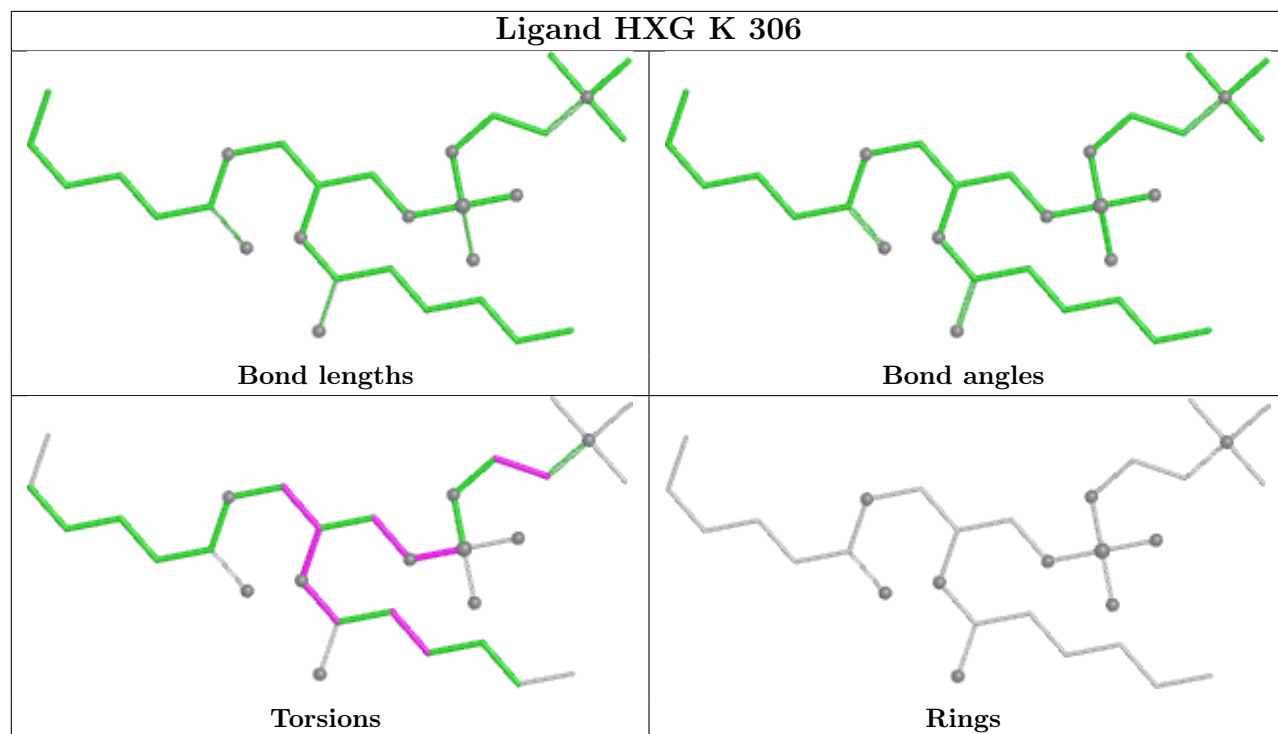


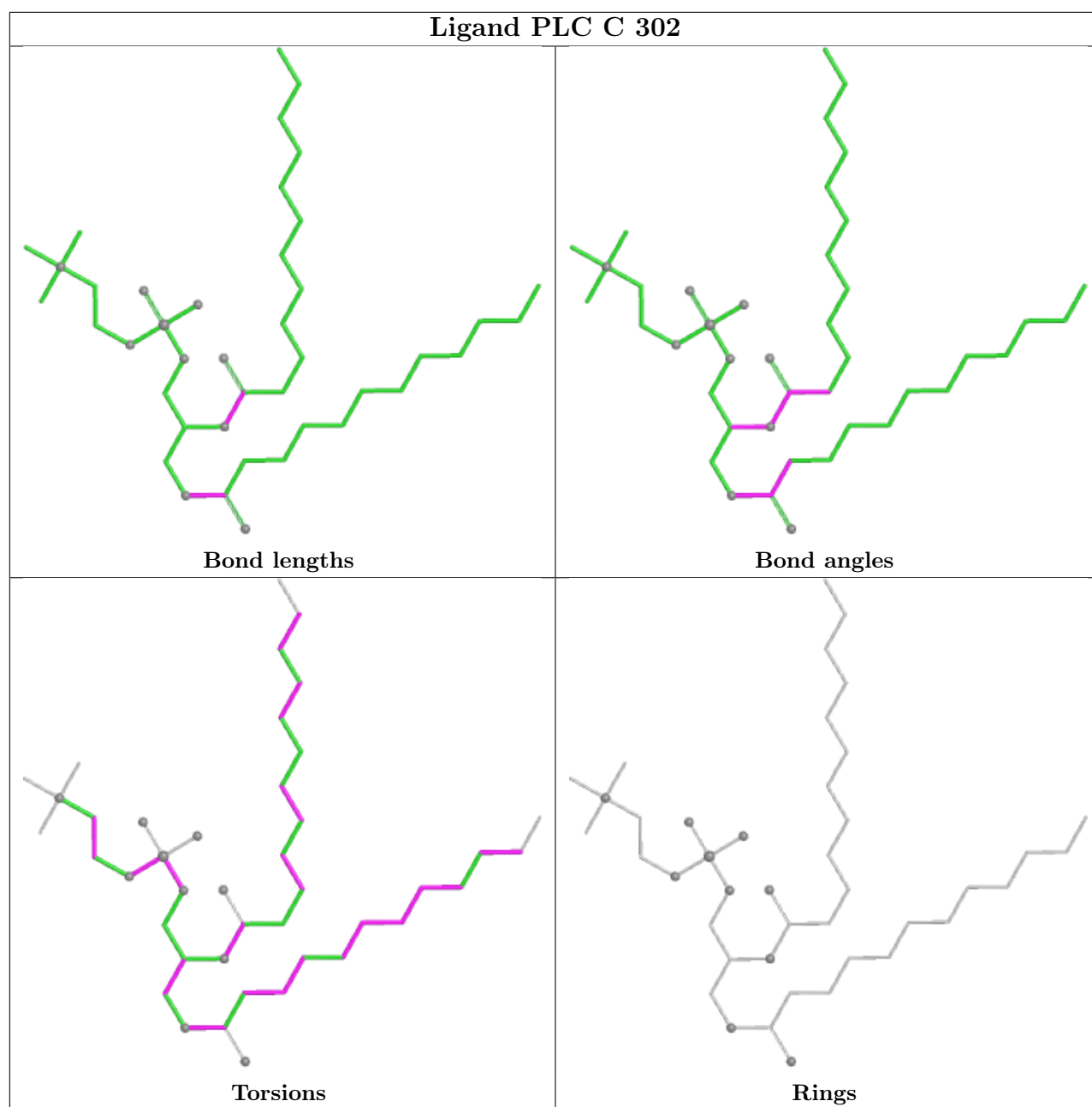


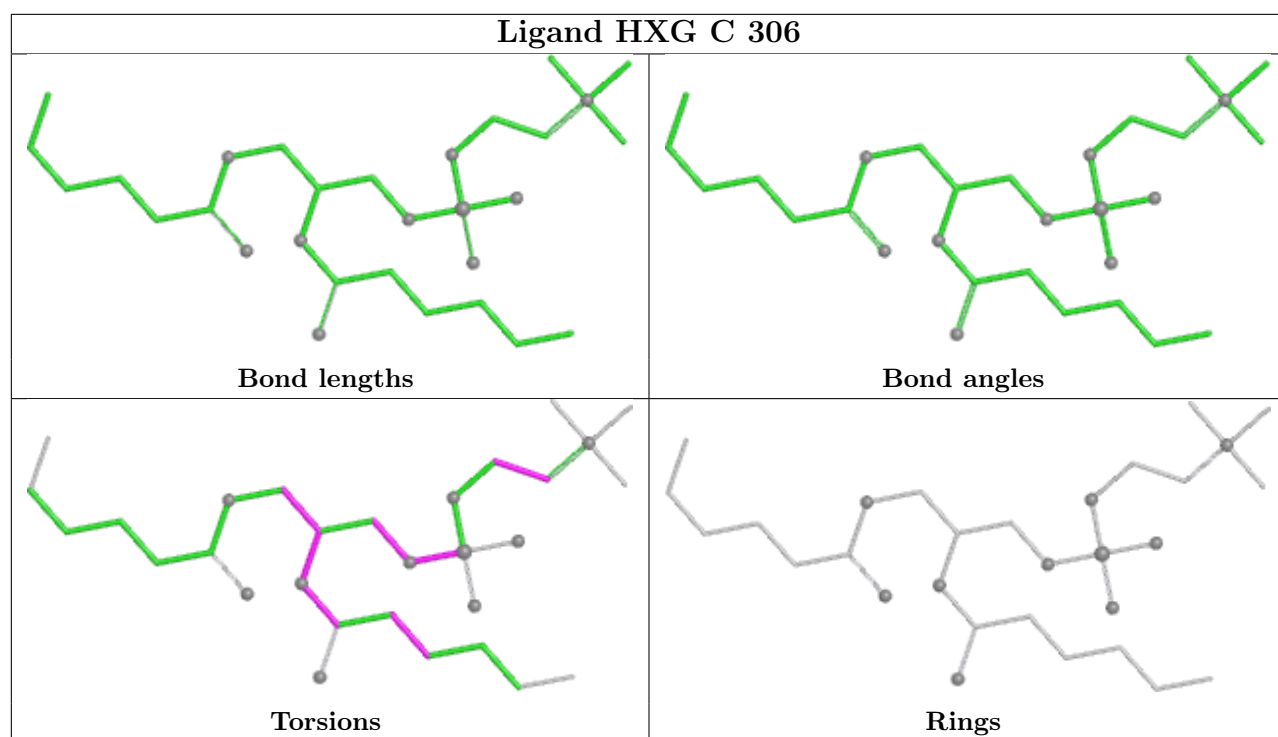


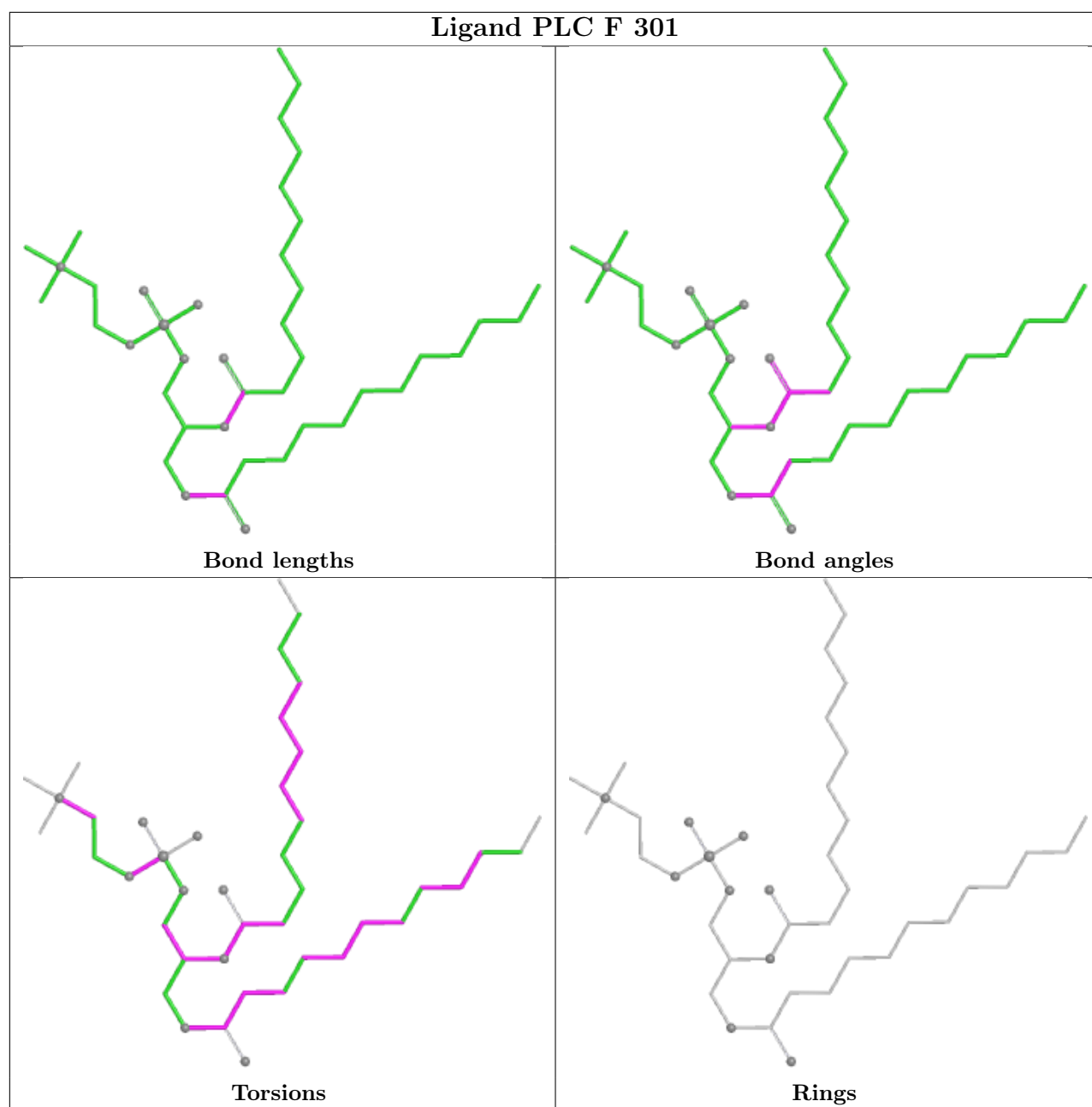


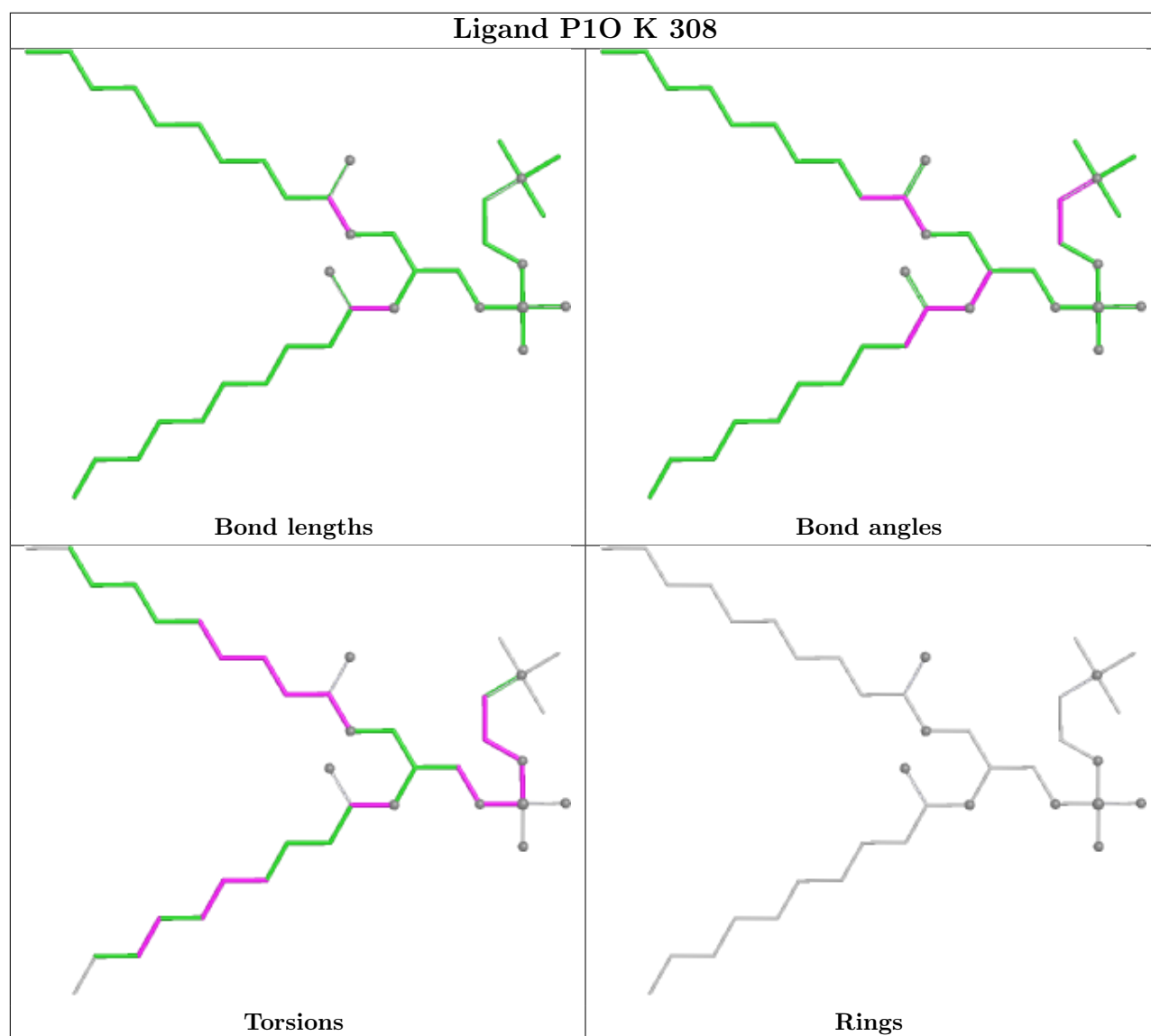


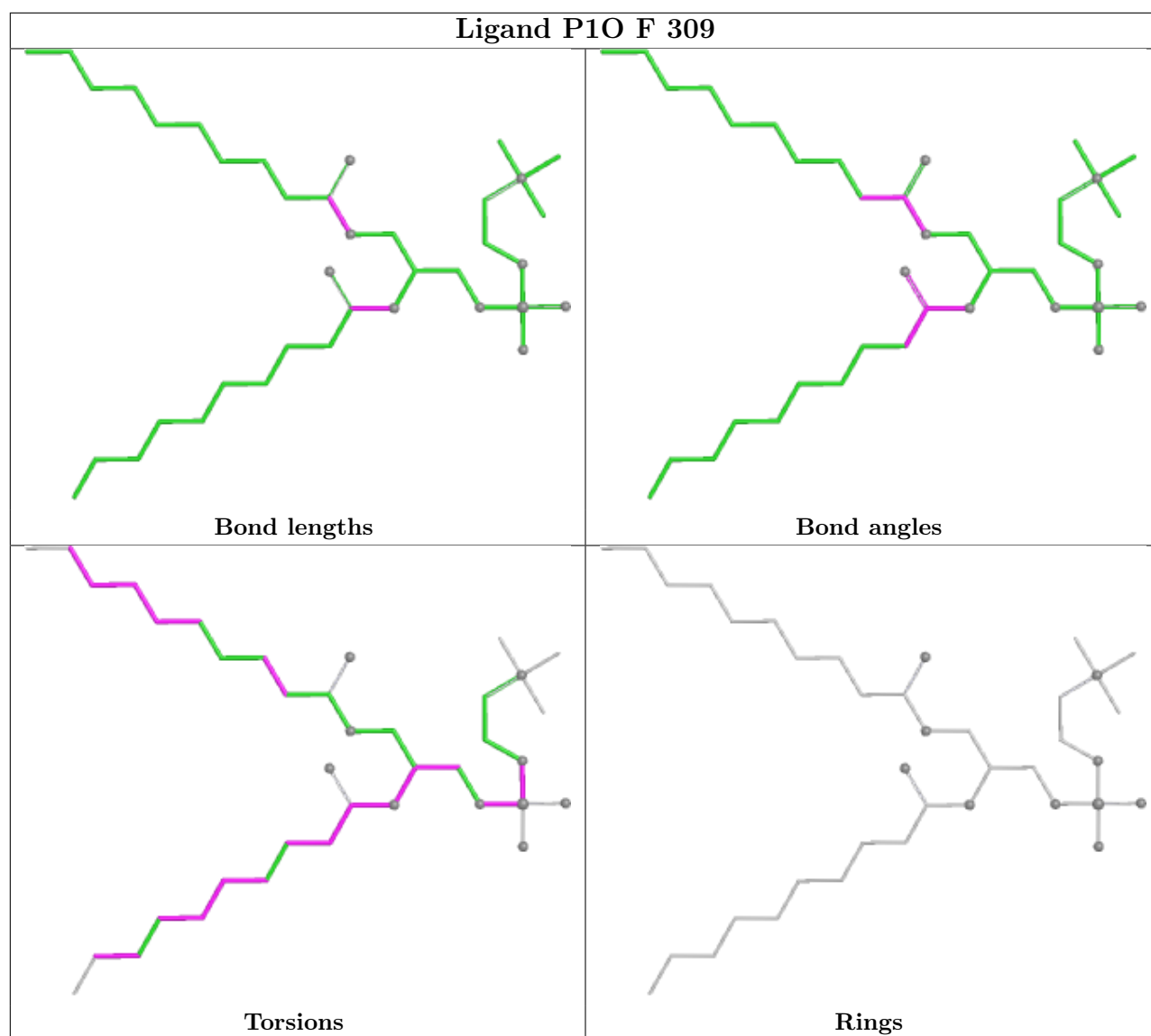


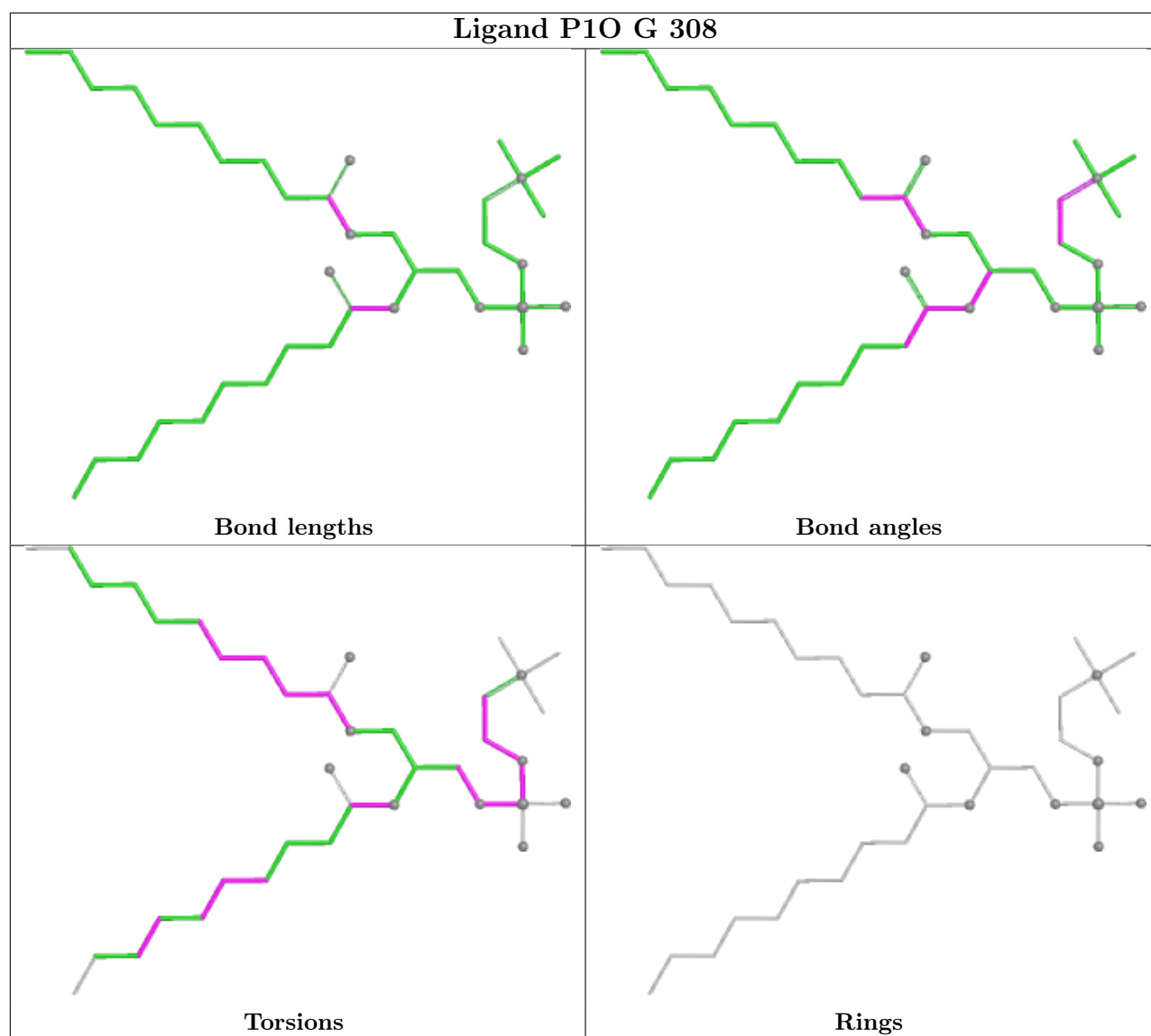


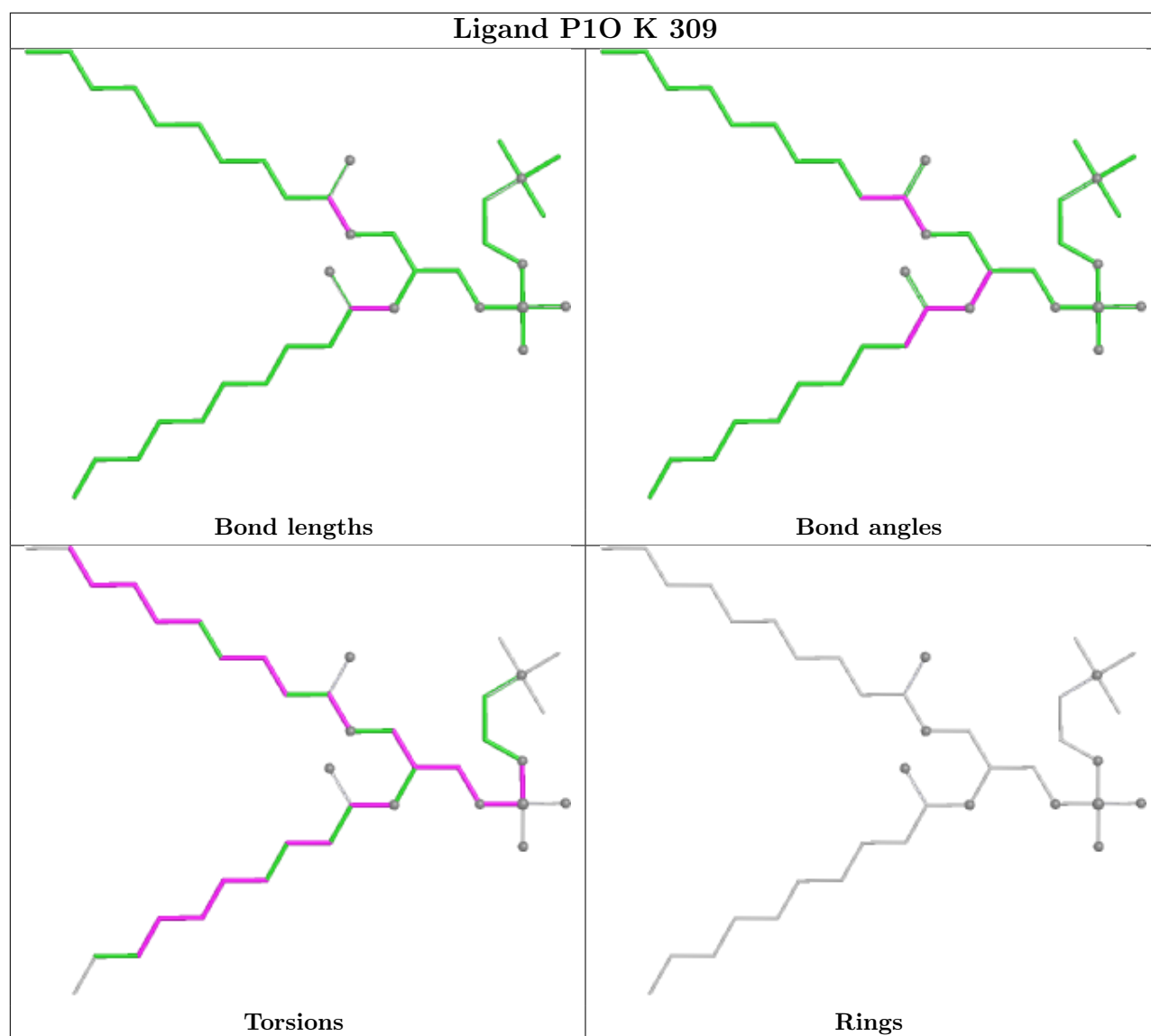


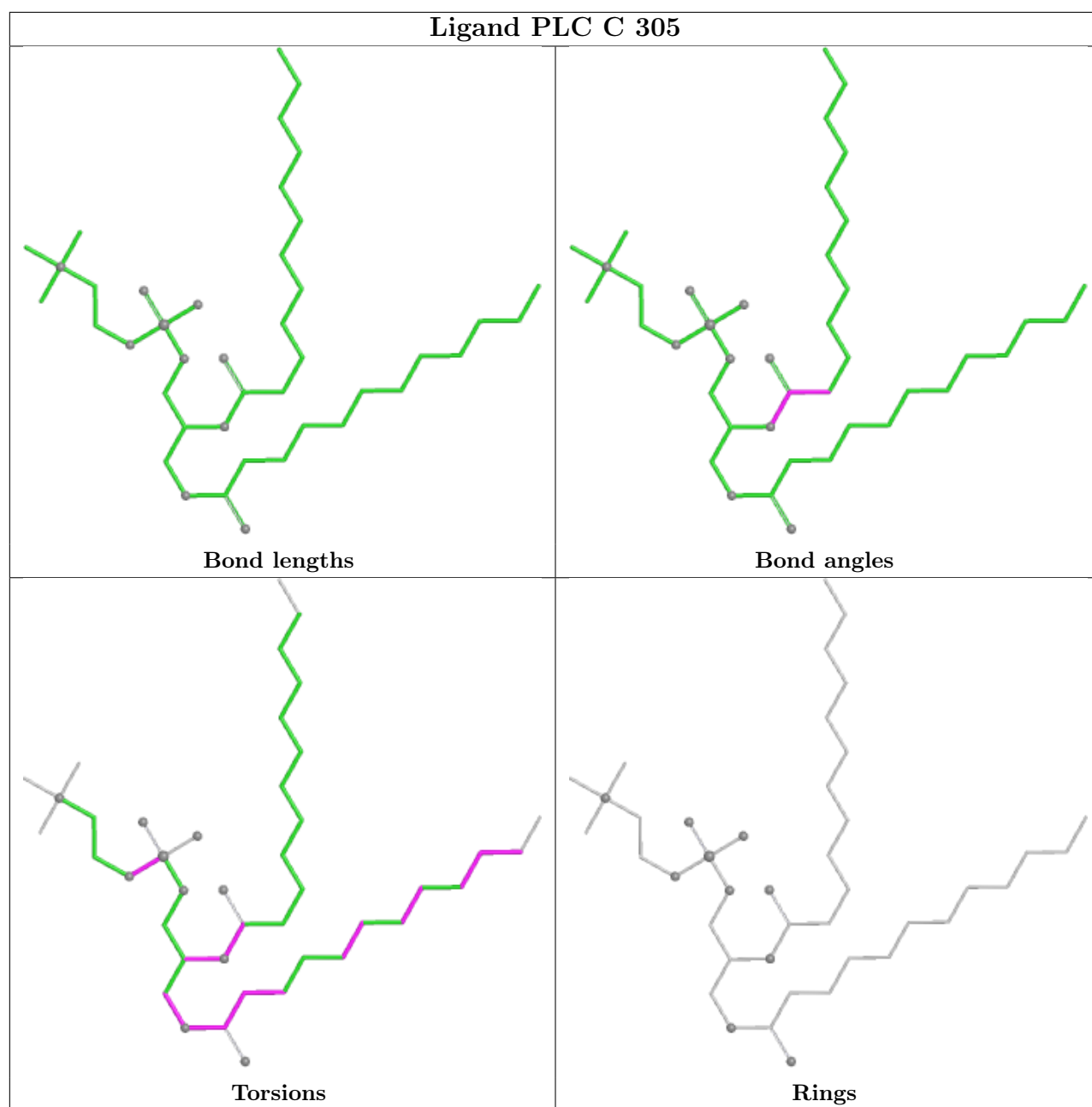


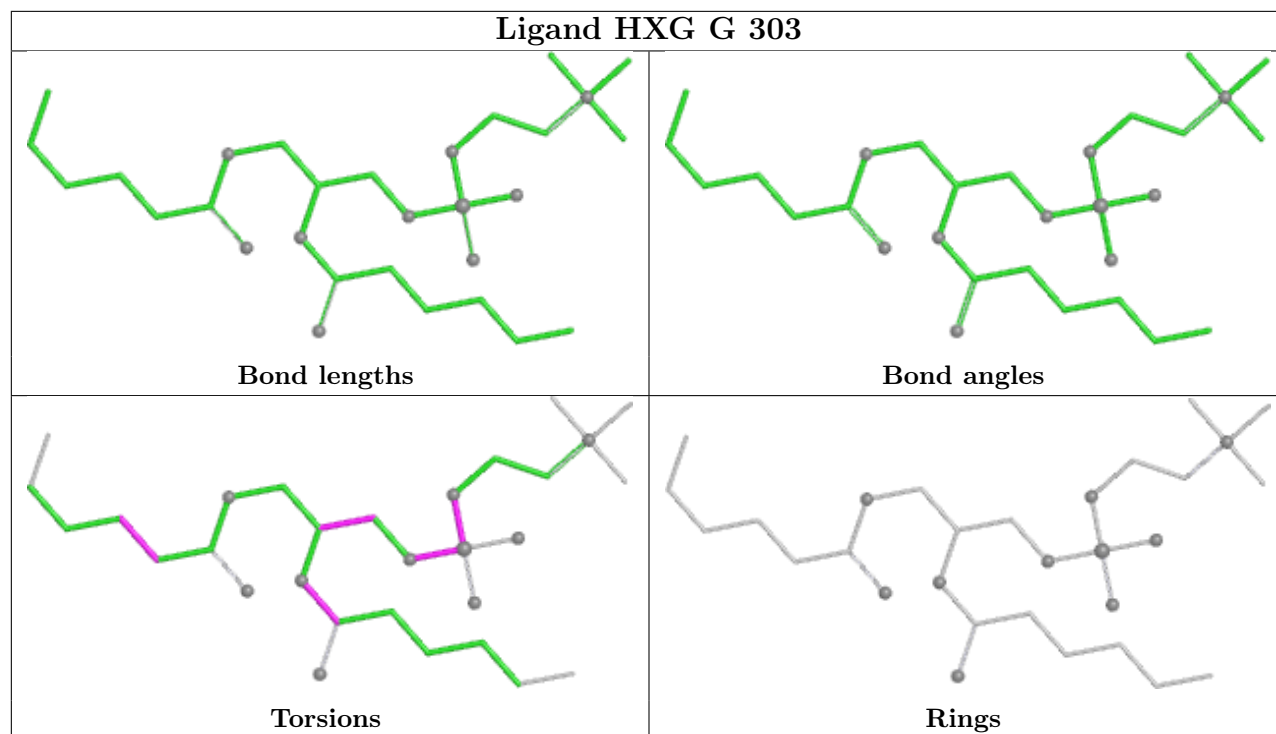


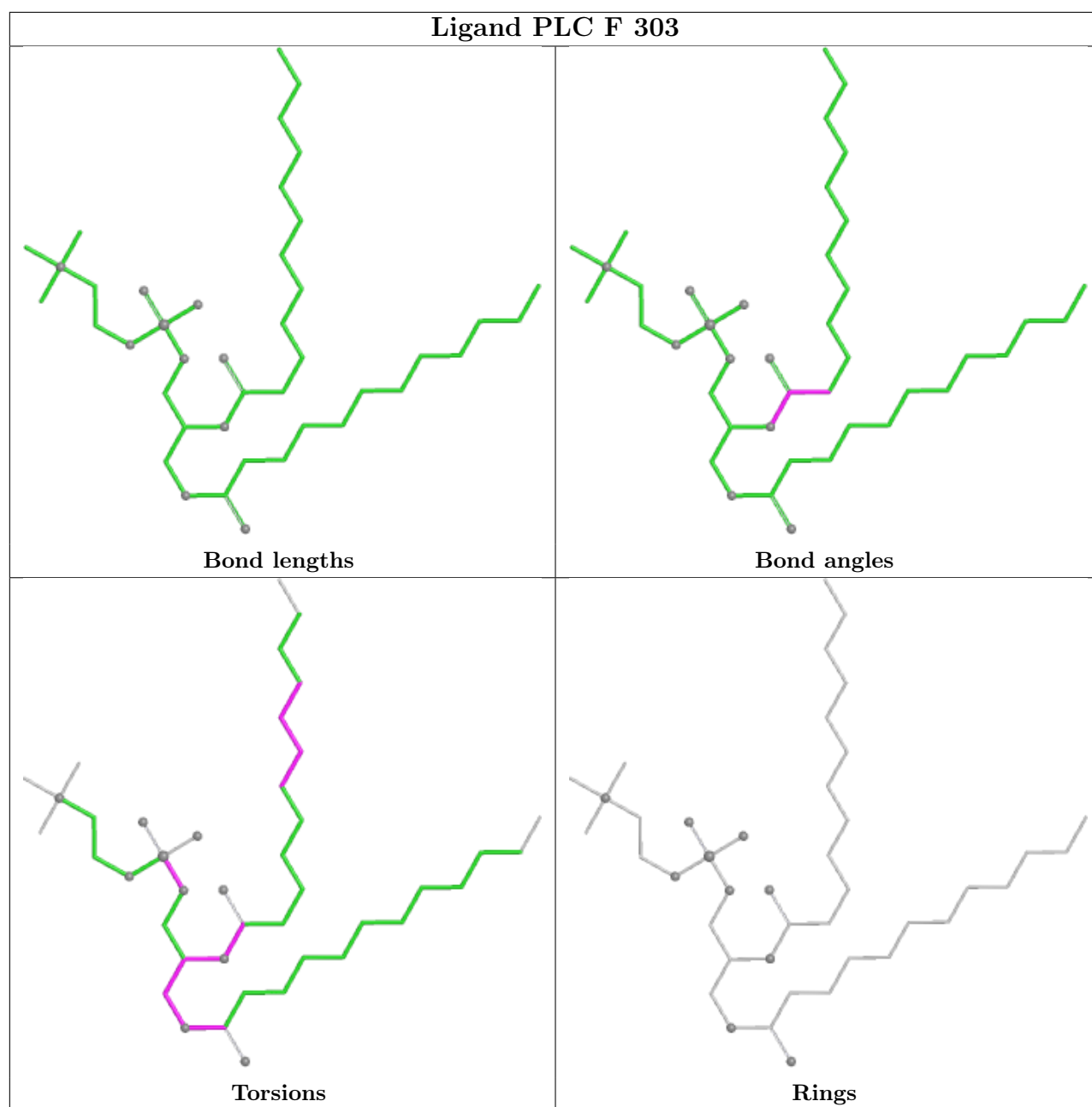


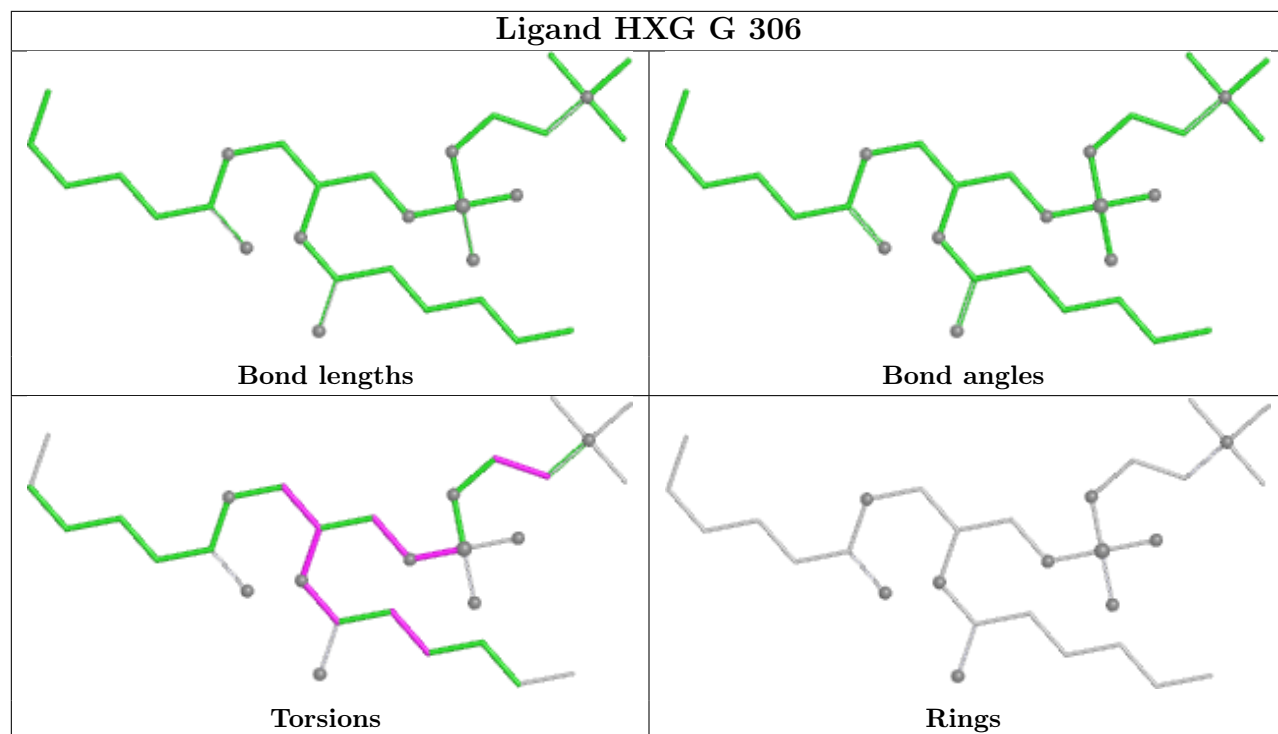


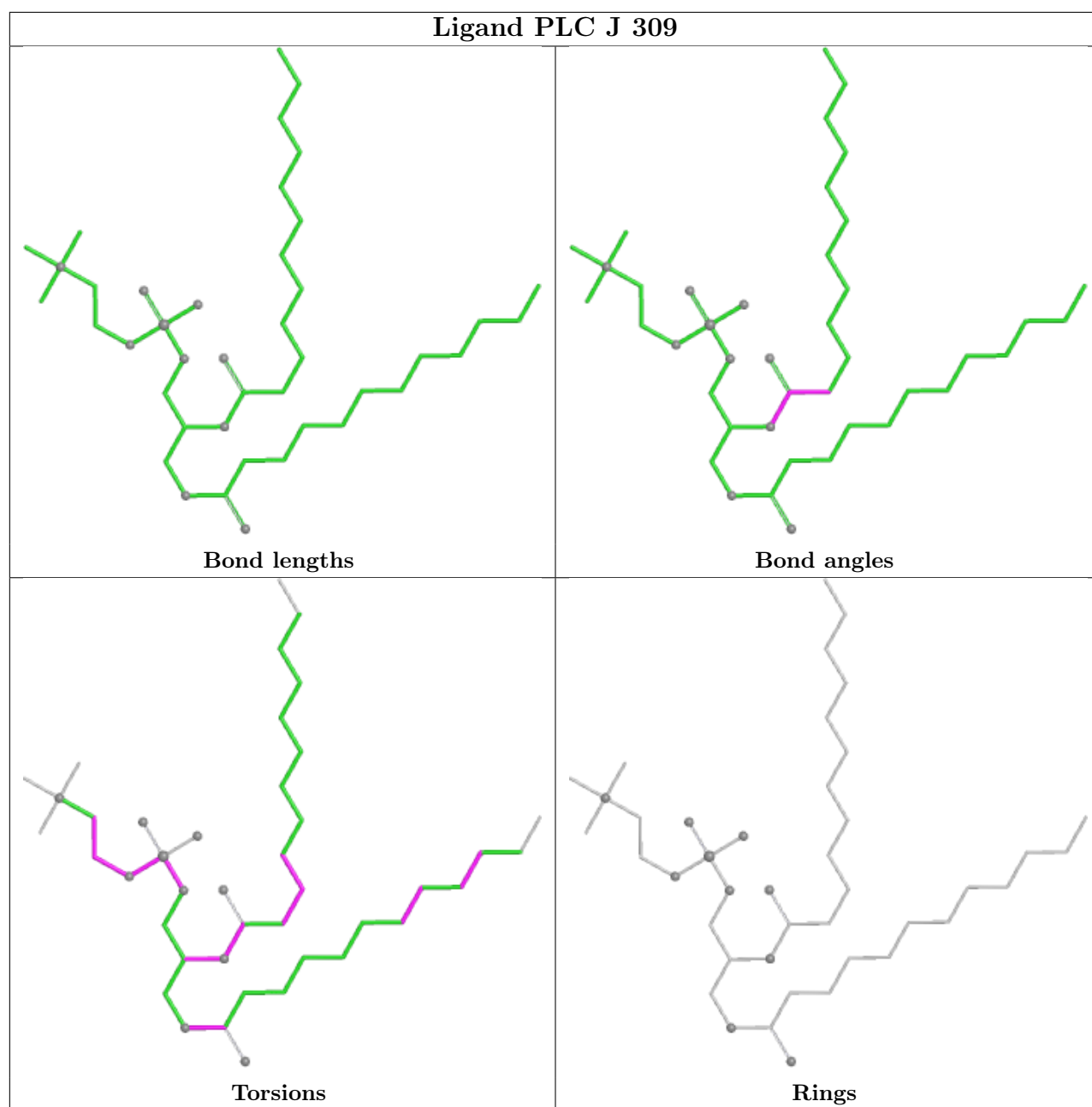


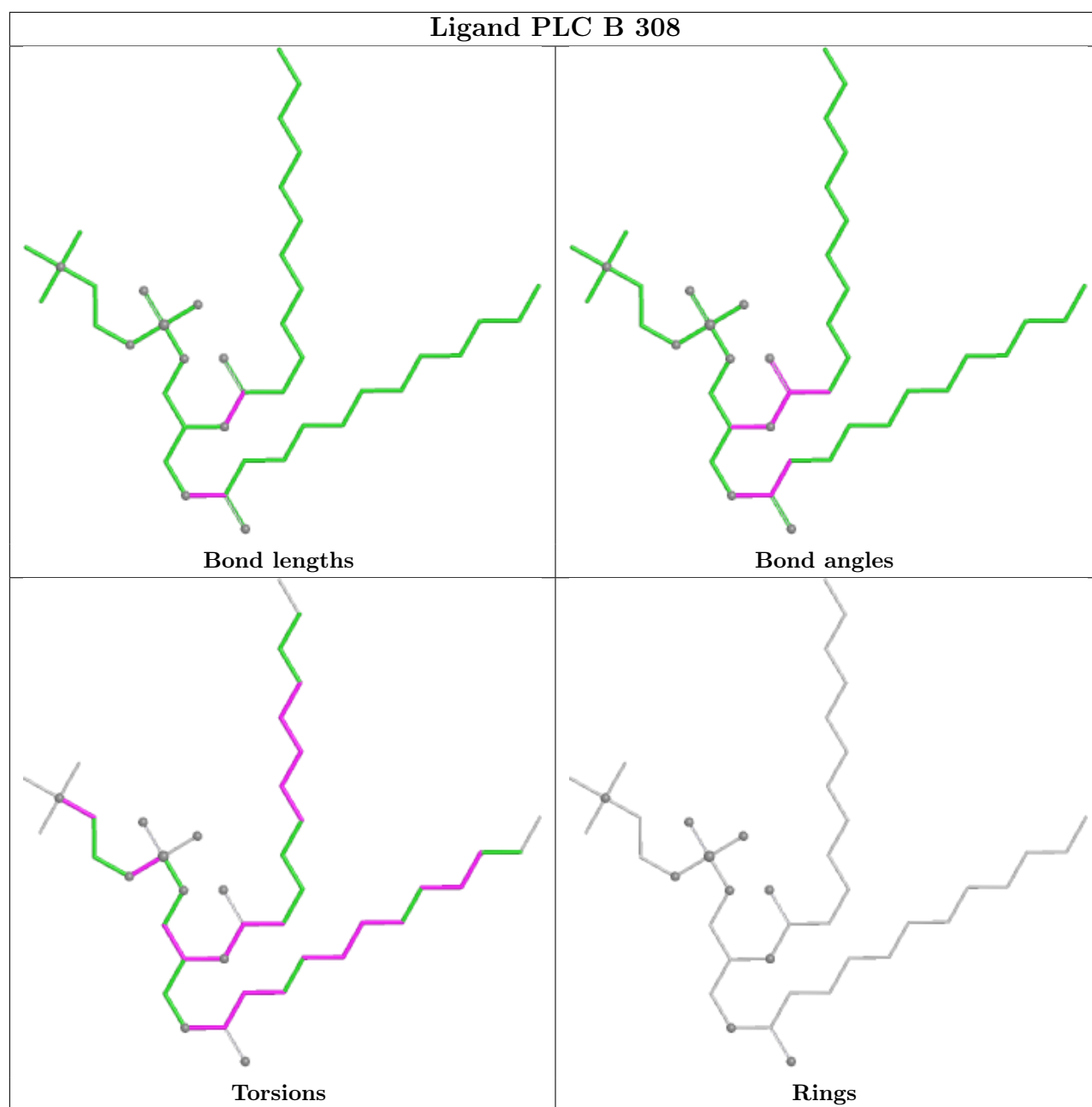


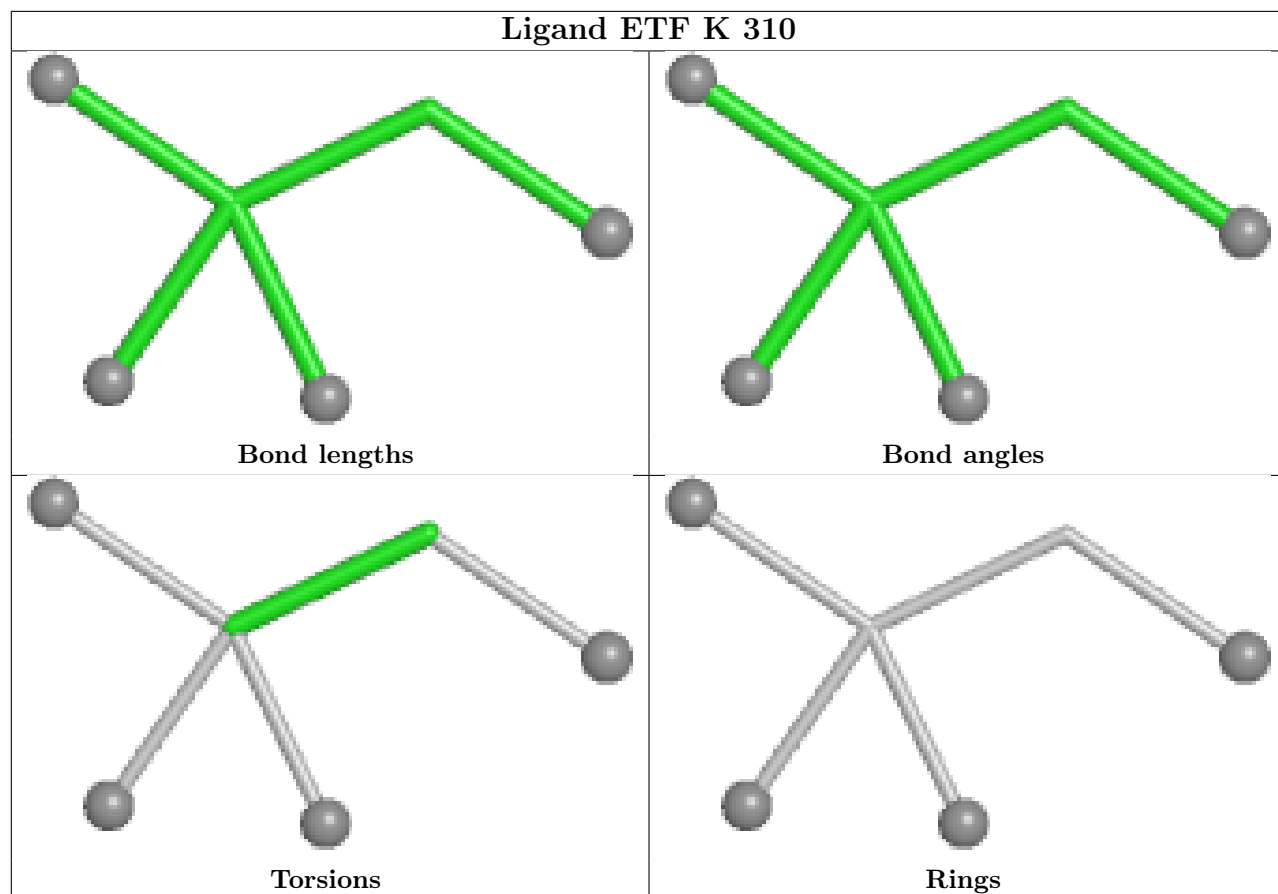


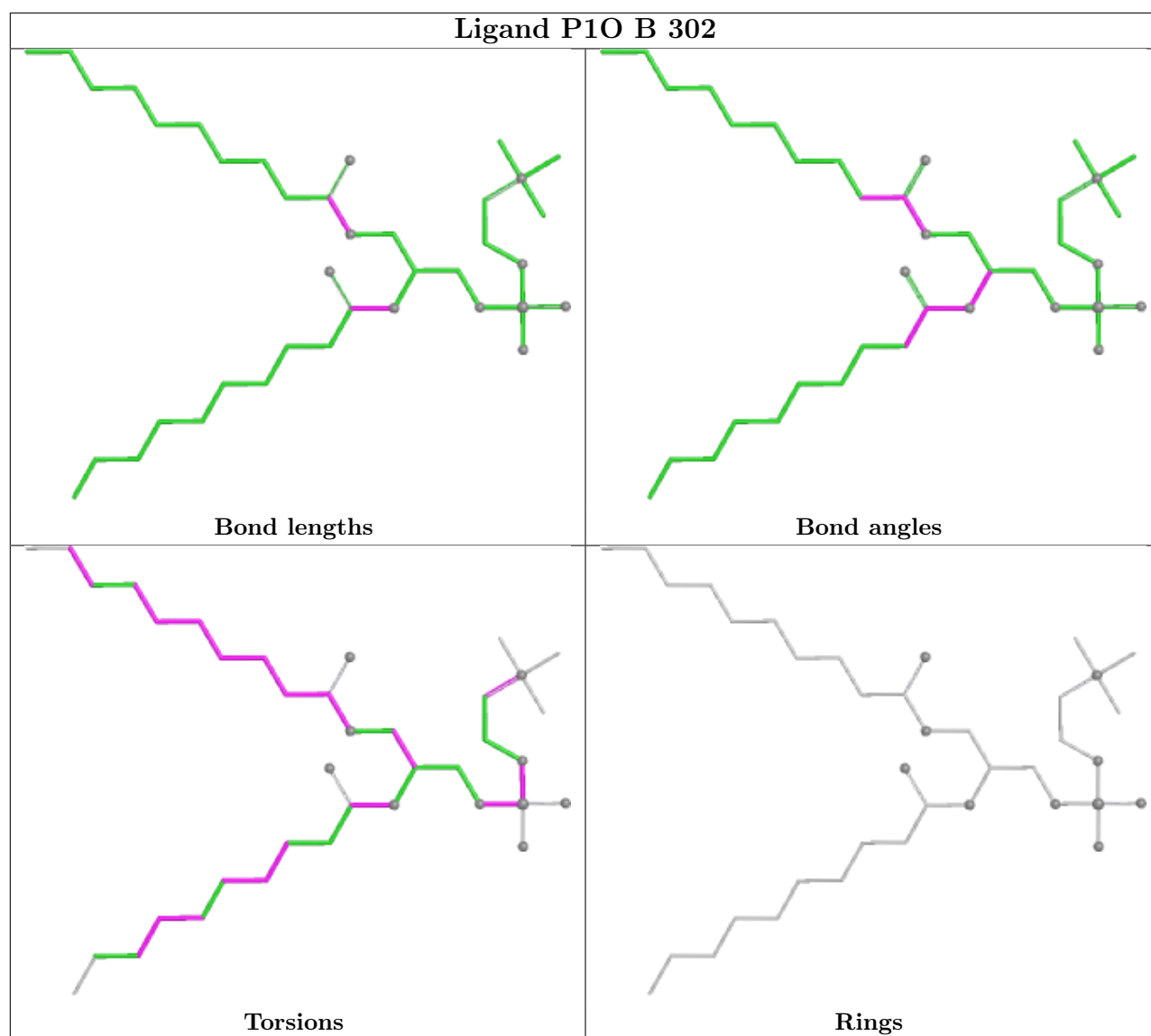












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

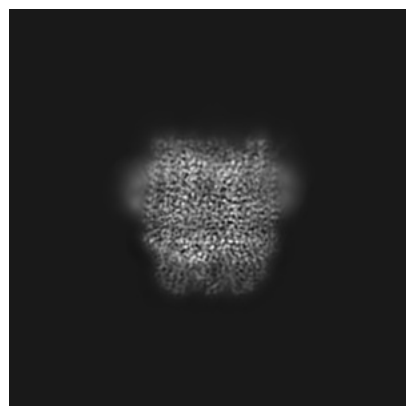
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17287. These allow visual inspection of the internal detail of the map and identification of artifacts.

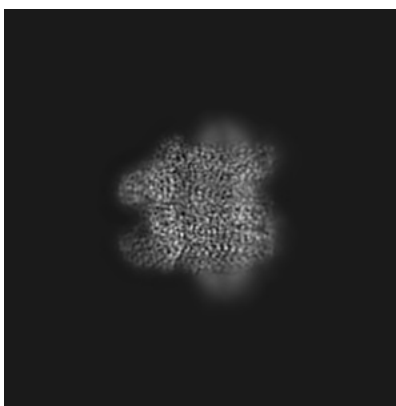
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

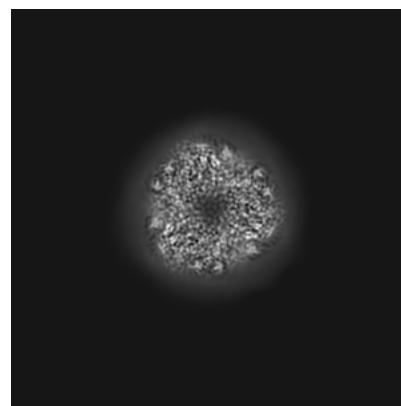
6.1.1 Primary map



X

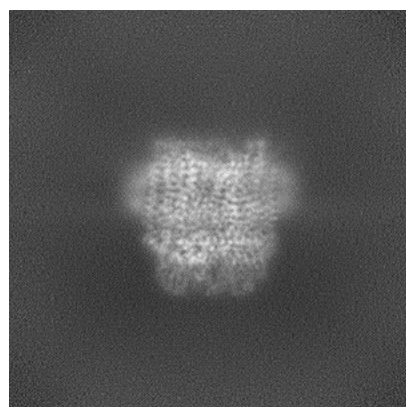


Y

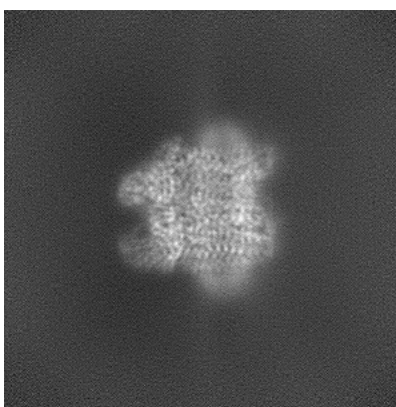


Z

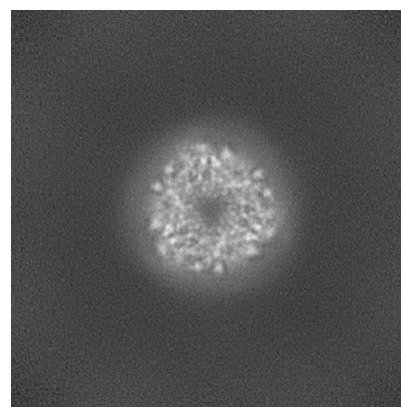
6.1.2 Raw map



X



Y

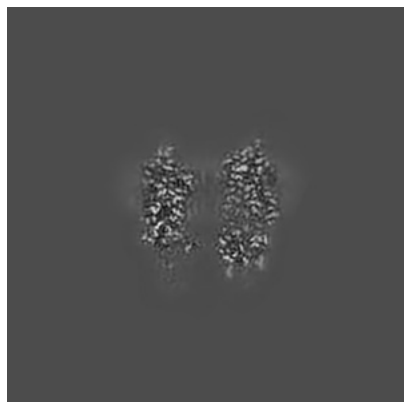


Z

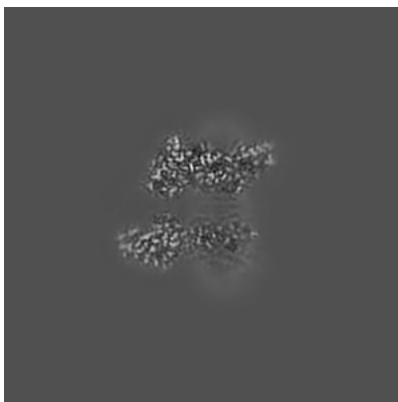
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

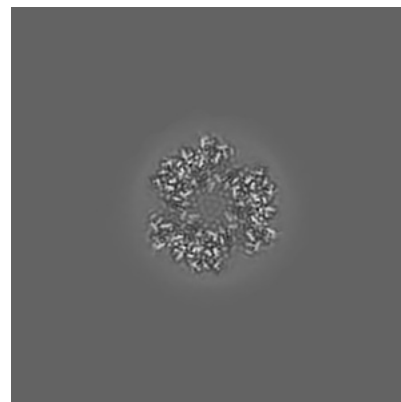
6.2.1 Primary map



X Index: 256

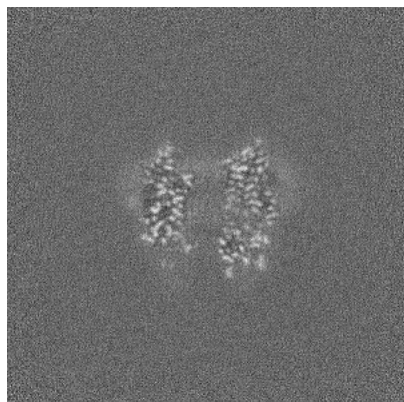


Y Index: 256

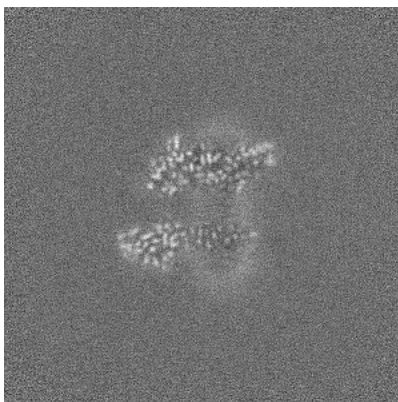


Z Index: 256

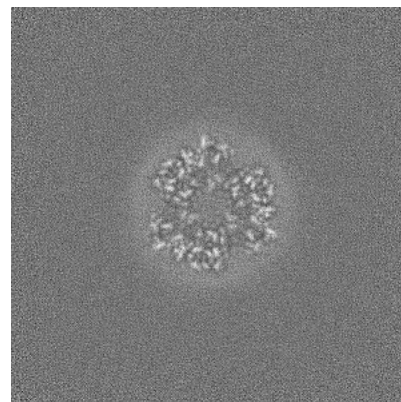
6.2.2 Raw map



X Index: 256



Y Index: 256

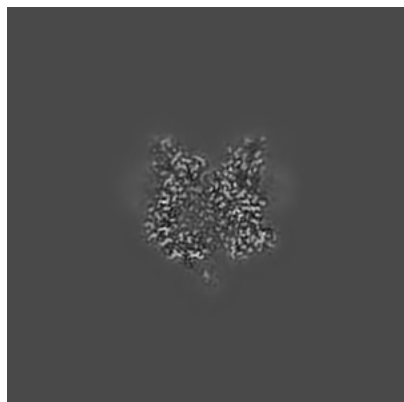


Z Index: 256

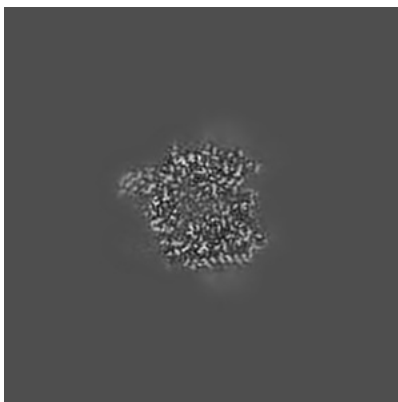
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

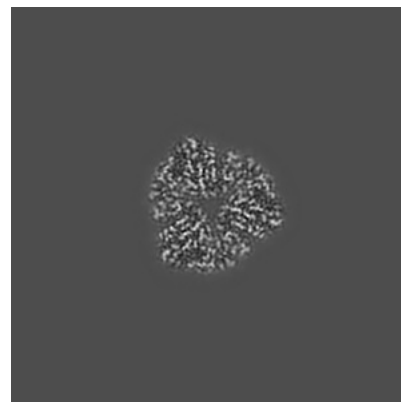
6.3.1 Primary map



X Index: 226

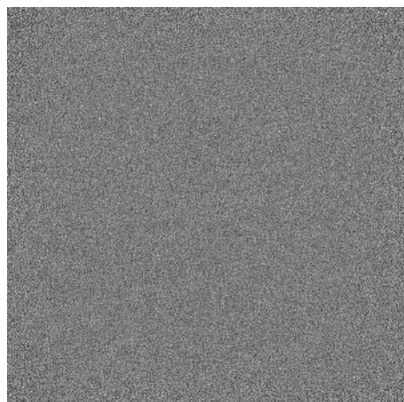


Y Index: 288

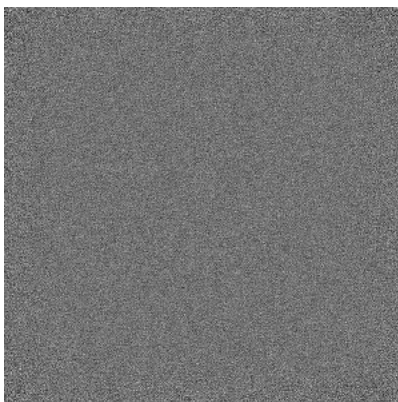


Z Index: 214

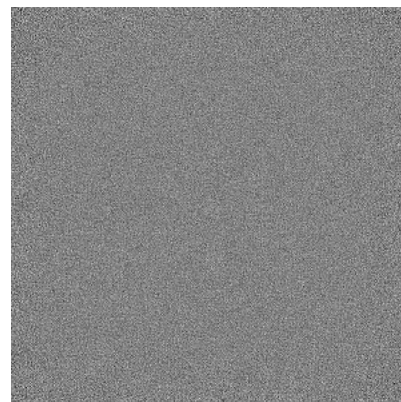
6.3.2 Raw map



X Index: 0



Y Index: 0

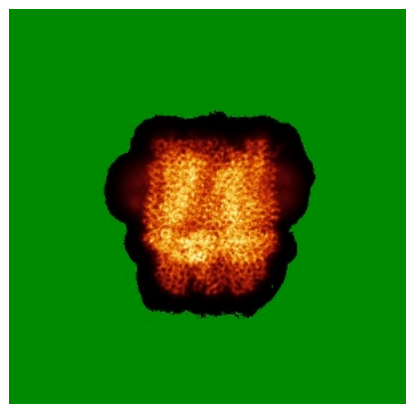


Z Index: 0

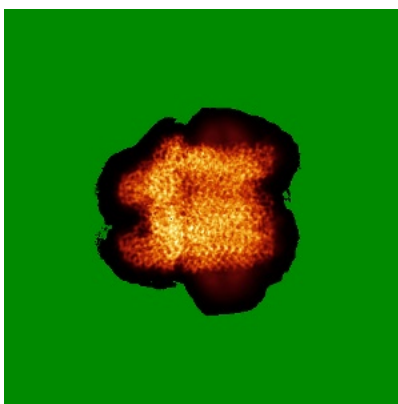
The images above show the largest variance slices of the map in three orthogonal directions.

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

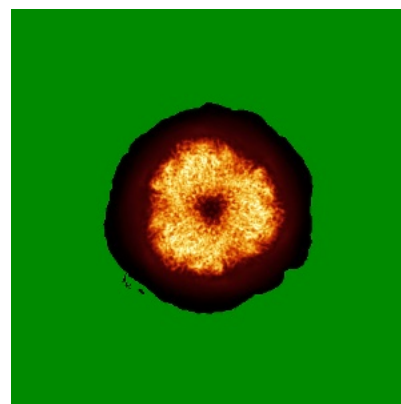
6.4.1 Primary map



X

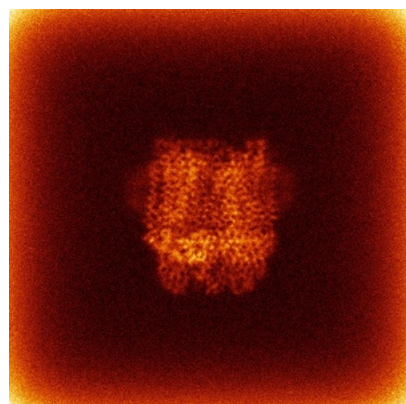


Y

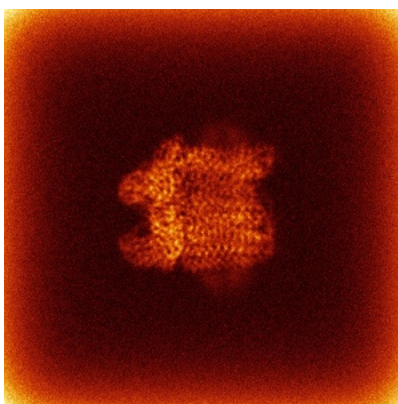


Z

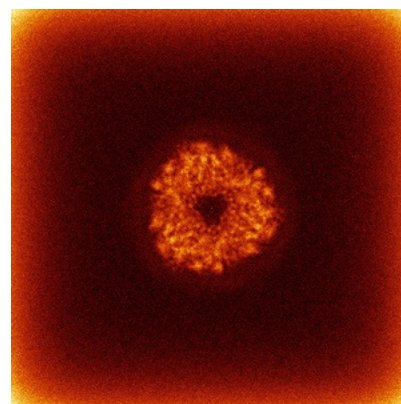
6.4.2 Raw map



X



Y

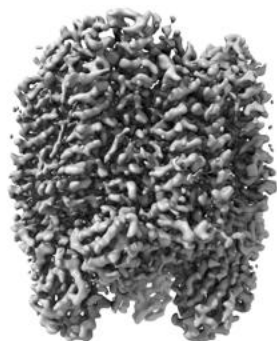


Z

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



X



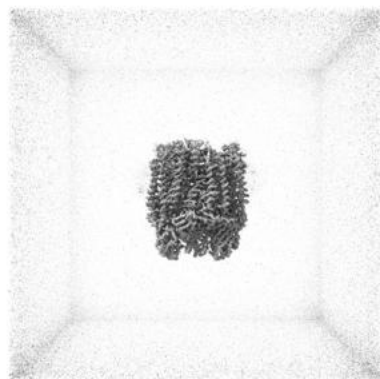
Y



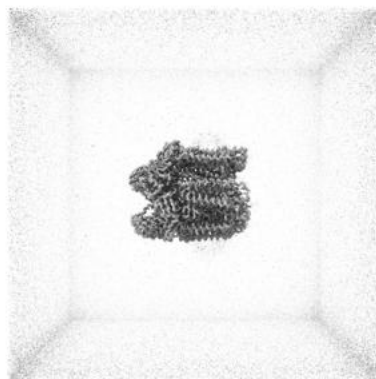
Z

The images above show the 3D surface view of the map at the recommended contour level 0.197. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

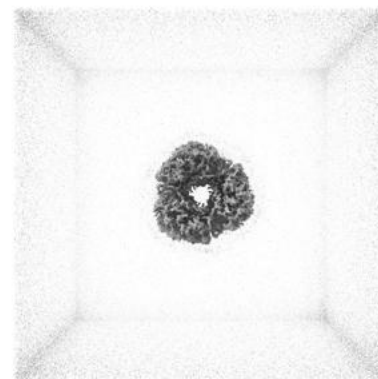
6.5.2 Raw map



X



Y



Z

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.

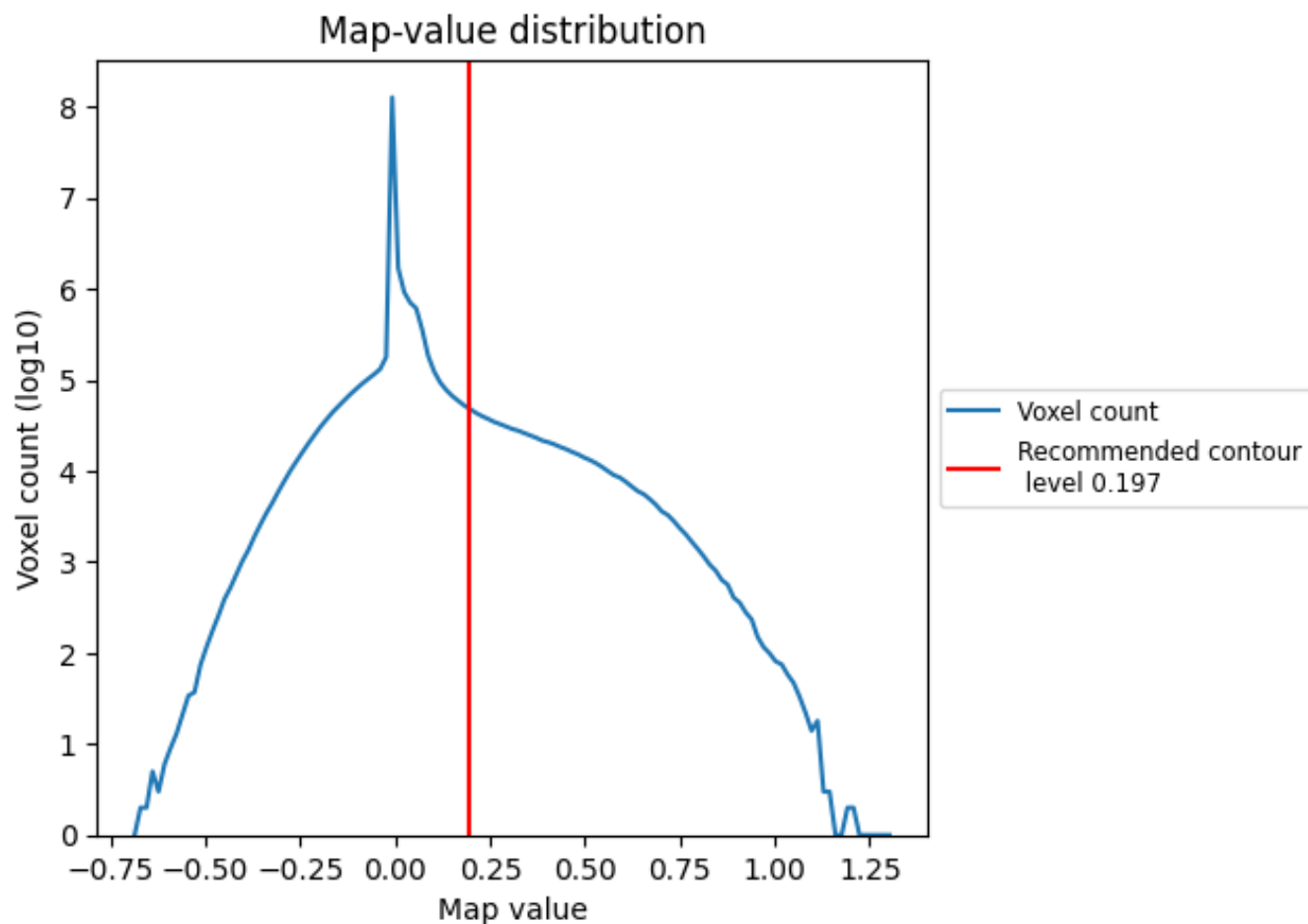
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

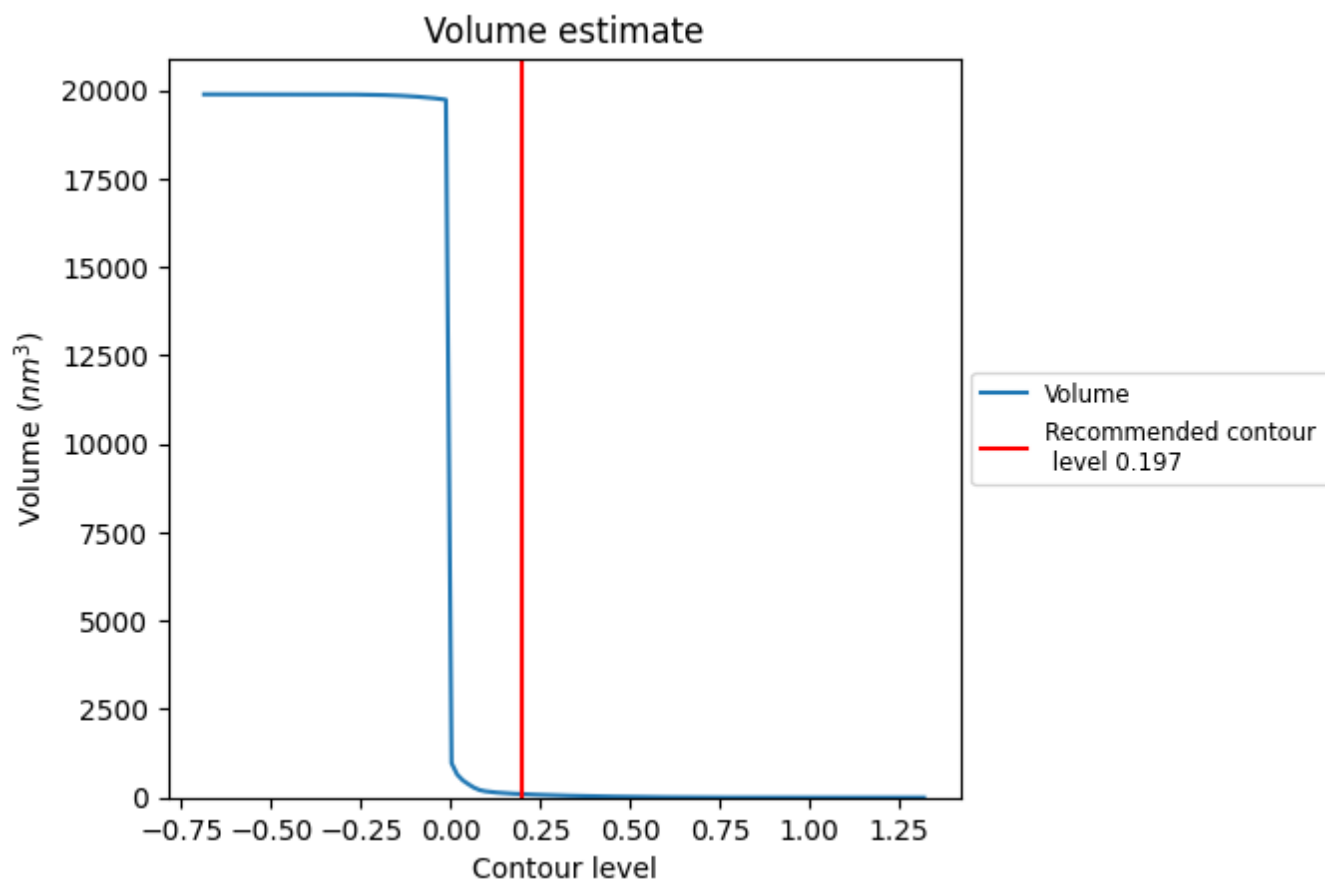
This section contains the results of statistical analysis of the map.

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.

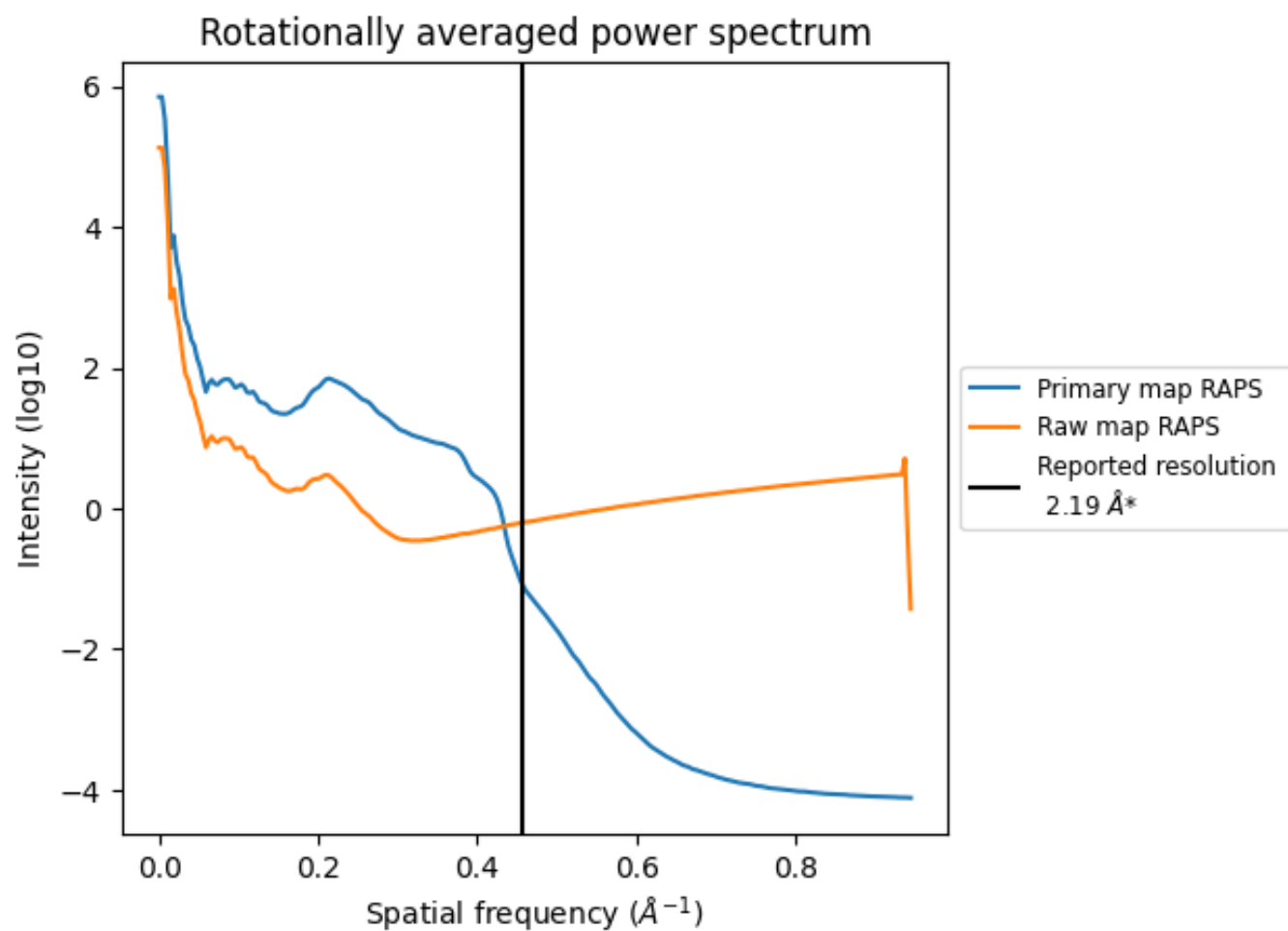
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 99 nm^3 ; this corresponds to an approximate mass of 89 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 ⓘ

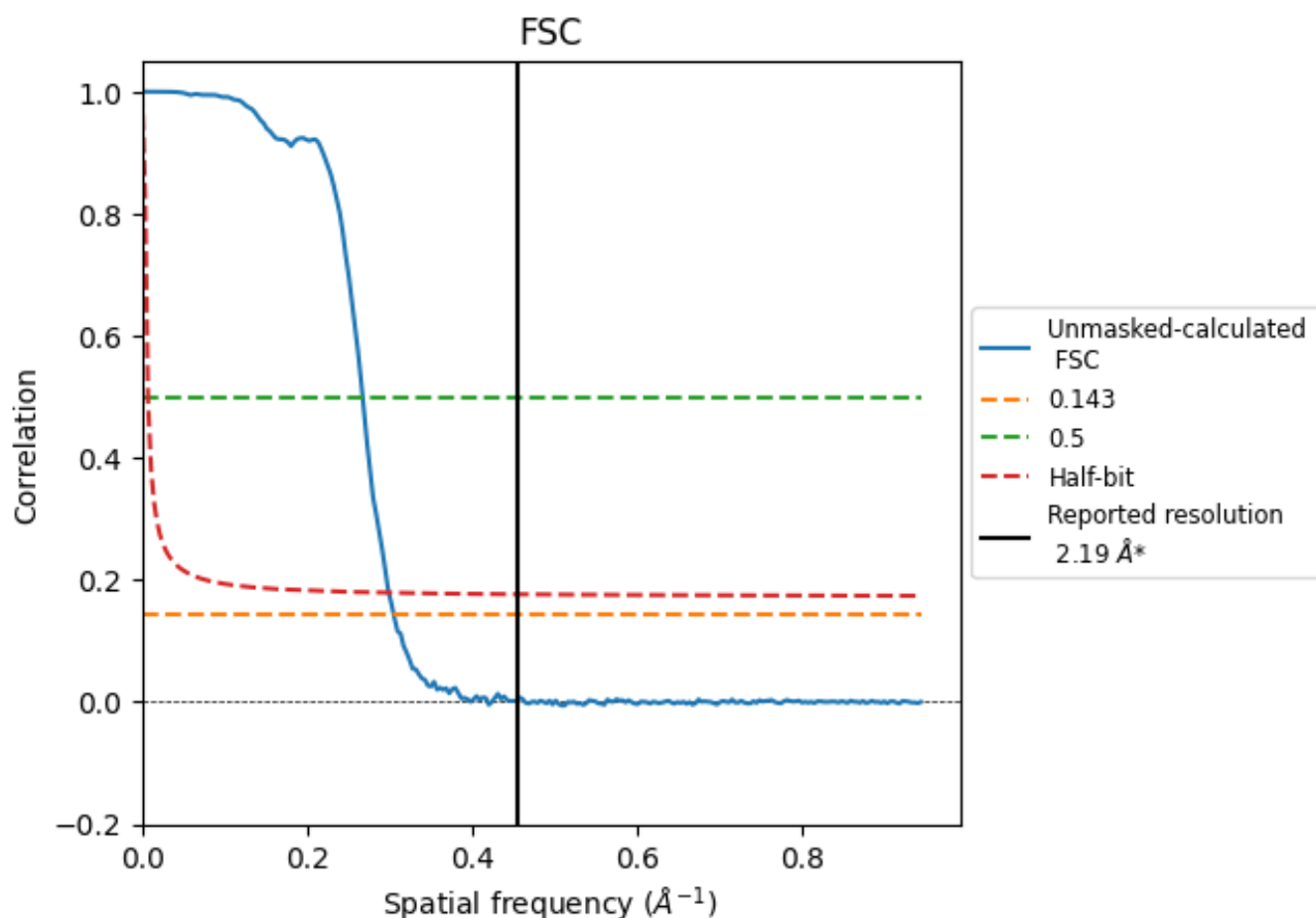


*Reported resolution corresponds to spatial frequency of 0.457 Å⁻¹

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 0.457 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.19	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.27	3.74	3.34

*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 3.27 differs from the reported value 2.19 by more than 10 %

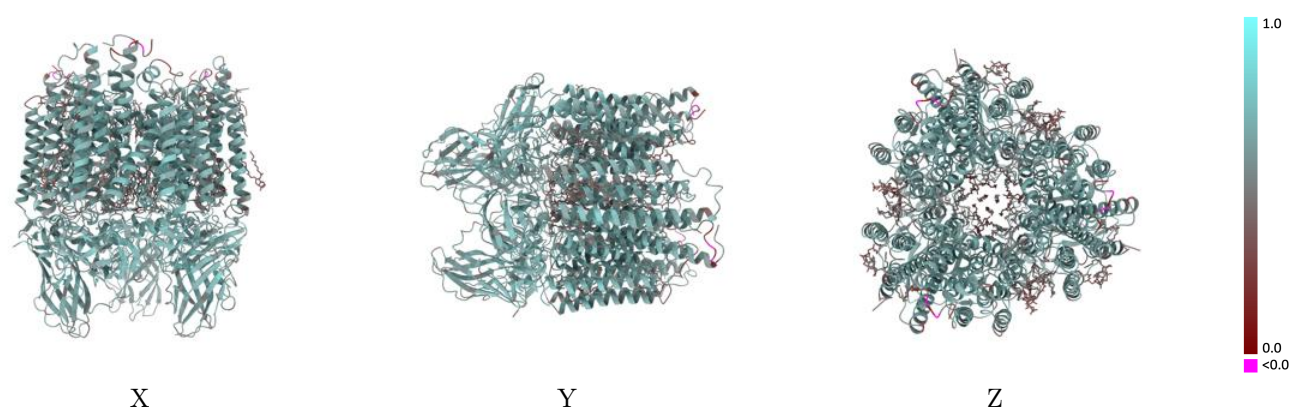
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-17287 and PDB model 8OYI. Per-residue inclusion information can be found in section 3 on page 11.

9.1 Map-model overlay [i](#)

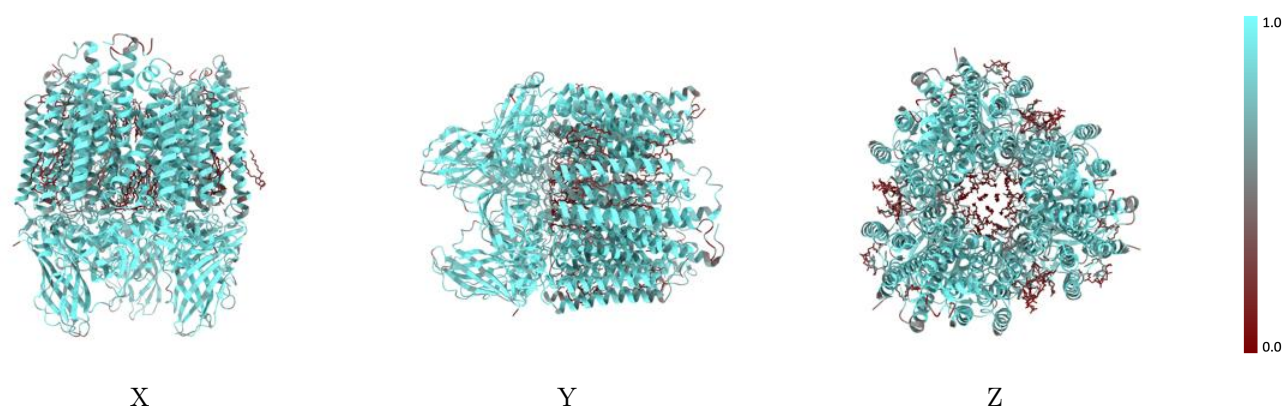
This section was not generated.

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



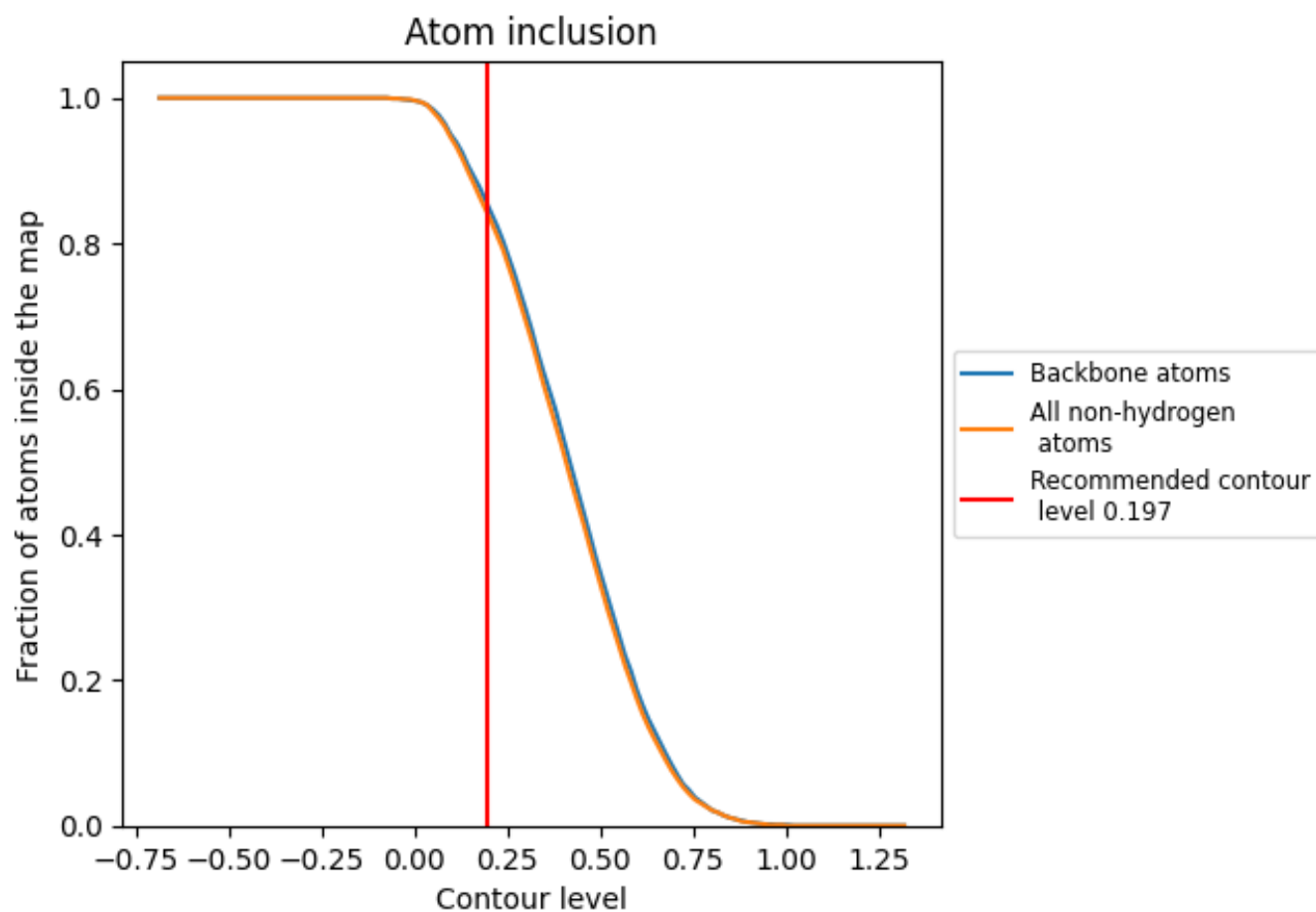
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.197).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 84% 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 (0.197) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8400	0.6160
A	0.8810	0.6240
B	0.8470	0.6360
C	0.7800	0.5870
E	0.8810	0.6240
F	0.8500	0.6330
G	0.7790	0.5870
I	0.8810	0.6240
J	0.8490	0.6330
K	0.7820	0.5860

1.0

0.0

<0.0