



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 09:37 PM UTC

PDB ID : 7RSH / pdb_00007rsh
EMDB ID : EMD-24670
Title : SthK Y26F Closed State
Authors : Gao, X.; Nimigean, C.
Deposited on : 2021-08-11
Resolution : 3.00 Å(reported)
Based on initial model : 6CJU

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

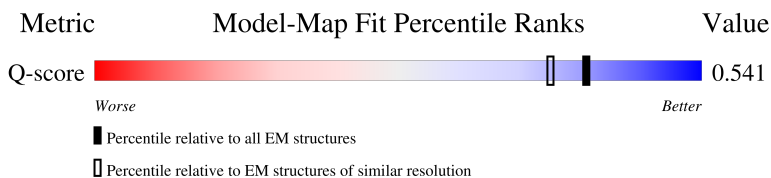
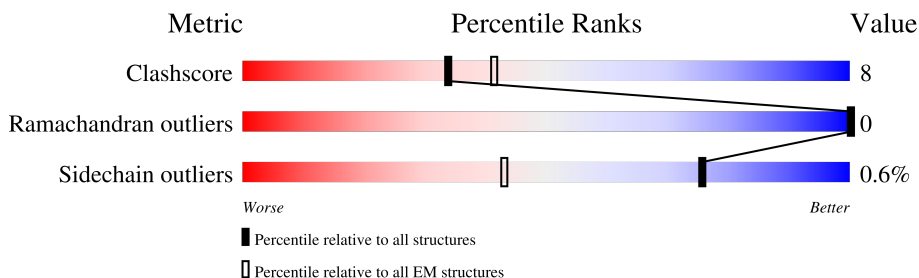
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14081 (2.50 - 3.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	456	15% 71% 17% 12%
1	B	456	15% 71% 16% 12%
1	C	456	15% 71% 17% 12%
1	D	456	15% 71% 17% 12%

2 Entry composition [i](#)

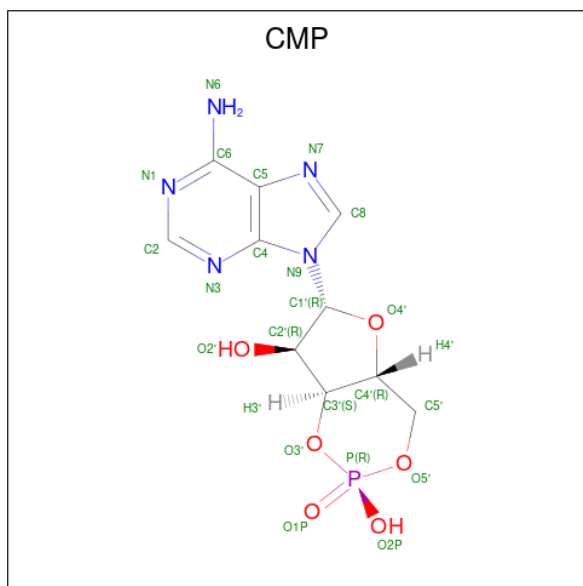
There are 3 unique types of molecules in this entry. The entry contains 13744 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called SthK.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	401	Total	C	N	O	S	0	0
			3191	2093	532	559	7		
1	B	401	Total	C	N	O	S	0	0
			3191	2093	532	559	7		
1	C	401	Total	C	N	O	S	0	0
			3191	2093	532	559	7		
1	D	401	Total	C	N	O	S	0	0
			3191	2093	532	559	7		

- Molecule 2 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (CCD ID: CMP) (formula: $C_{10}H_{12}N_5O_6P$).



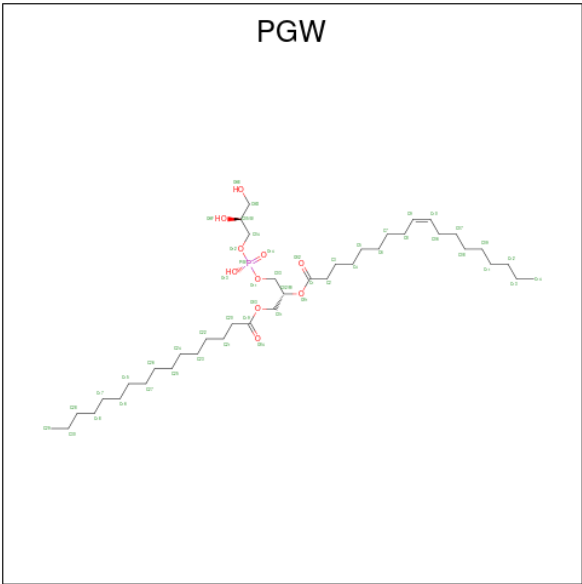
Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	N	O	P	0
			22	10	5	6	1	
2	B	1	Total	C	N	O	P	0
			22	10	5	6	1	

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Mol	Chain	Residues	Atoms					AltConf
2	C	1	Total	C	N	O	P	0
			22	10	5	6	1	
2	D	1	Total	C	N	O	P	0
			22	10	5	6	1	

- Molecule 3 is (1R)-2-{[(S)-{[(2S)-2,3-dihydroxypropyl]oxy}(hydroxy)phosphoryl]oxy}-1-[(hexadecanoyloxy)methyl]ethyl (9Z)-octadec-9-enoate (CCD ID: PGW) (formula: C₄₀H₇₇O₁₀P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
3	A	1	Total 43	C 32	O 10	P 1	0
3	A	1	Total 15	C 15			0
3	A	1	Total 13	C 13			0
3	A	1	Total 24	C 20	O 4		0
3	A	1	Total 13	C 13			0
3	A	1	Total 9	C 9			0
3	A	1	Total 15	C 15			0
3	A	1	Total 15	C 15			0

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Mol	Chain	Residues	Atoms				AltConf
3	A	1	Total	C			0
			15	15			
3	A	1	Total	C			0
			13	13			
3	A	1	Total	C			0
			8	8			
3	A	1	Total	C			0
			12	12			
3	A	1	Total	C			0
			15	15			
3	A	1	Total	C			0
			13	13			
3	B	1	Total	C	O	P	0
			43	32	10	1	
3	B	1	Total	C			0
			15	15			
3	B	1	Total	C			0
			13	13			
3	B	1	Total	C	O		0
			24	20	4		
3	B	1	Total	C			0
			13	13			
3	B	1	Total	C			0
			9	9			
3	B	1	Total	C			0
			15	15			
3	B	1	Total	C			0
			15	15			
3	B	1	Total	C			0
			15	15			
3	B	1	Total	C			0
			13	13			
3	B	1	Total	C			0
			8	8			
3	B	1	Total	C			0
			12	12			
3	B	1	Total	C			0
			15	15			
3	B	1	Total	C			0
			13	13			
3	C	1	Total	C	O	P	0
			43	32	10	1	

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Mol	Chain	Residues	Atoms	AltConf
3	C	1	Total C 15 15	0
3	C	1	Total C 13 13	0
3	C	1	Total C O 24 20 4	0
3	C	1	Total C 13 13	0
3	C	1	Total C 9 9	0
3	C	1	Total C 15 15	0
3	C	1	Total C 15 15	0
3	C	1	Total C 15 15	0
3	C	1	Total C 13 13	0
3	C	1	Total C 8 8	0
3	C	1	Total C 12 12	0
3	C	1	Total C 15 15	0
3	C	1	Total C 13 13	0
3	D	1	Total C 13 13	0
3	D	1	Total C O P 43 32 10 1	0
3	D	1	Total C 15 15	0
3	D	1	Total C 13 13	0
3	D	1	Total C O 24 20 4	0
3	D	1	Total C 13 13	0
3	D	1	Total C 9 9	0
3	D	1	Total C 15 15	0

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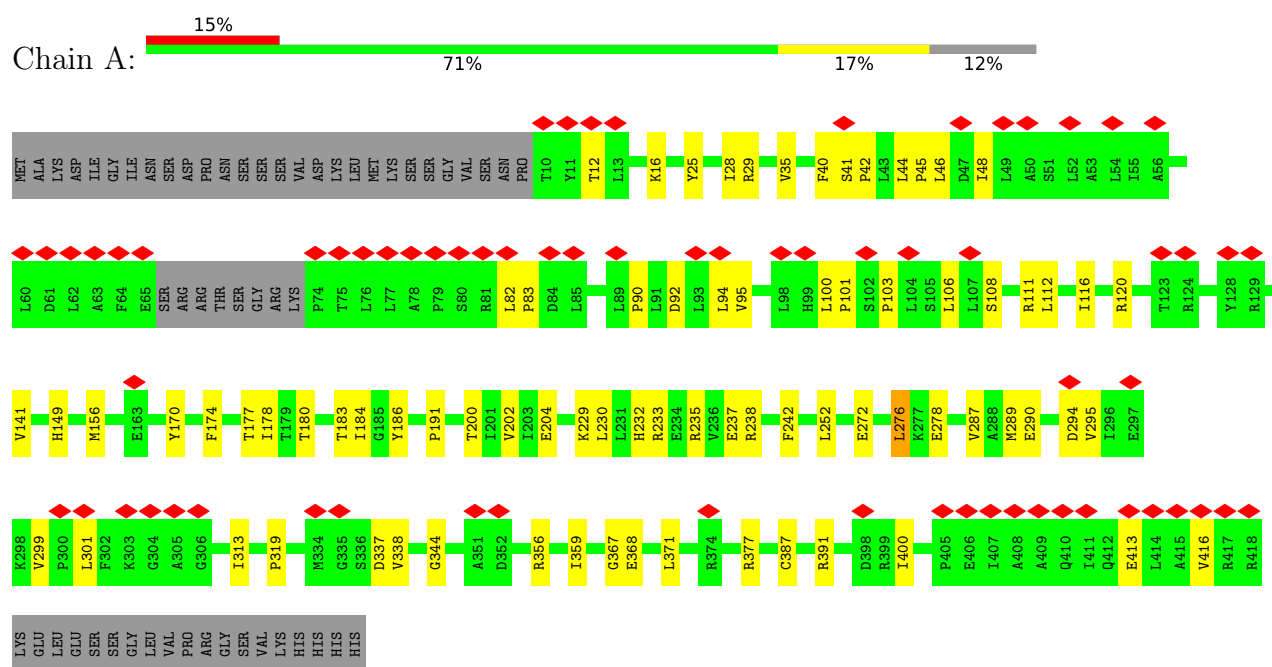
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Mol	Chain	Residues	Atoms		AltConf
3	D	1	Total	C	0
			15	15	
3	D	1	Total	C	0
			15	15	
3	D	1	Total	C	0
			13	13	
3	D	1	Total	C	0
			8	8	
3	D	1	Total	C	0
			12	12	
3	D	1	Total	C	0
			15	15	

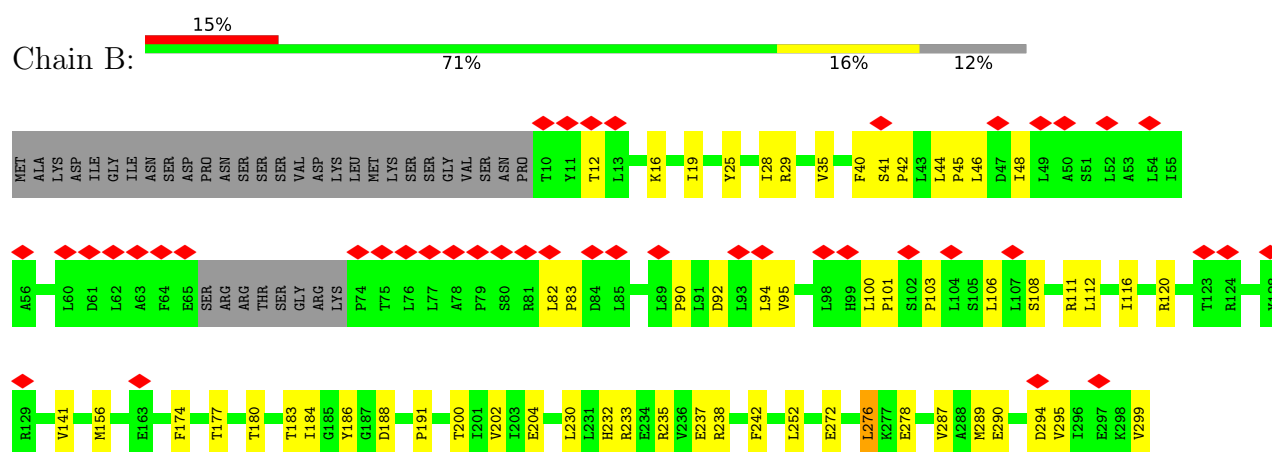
3 Residue-property plots [i](#)

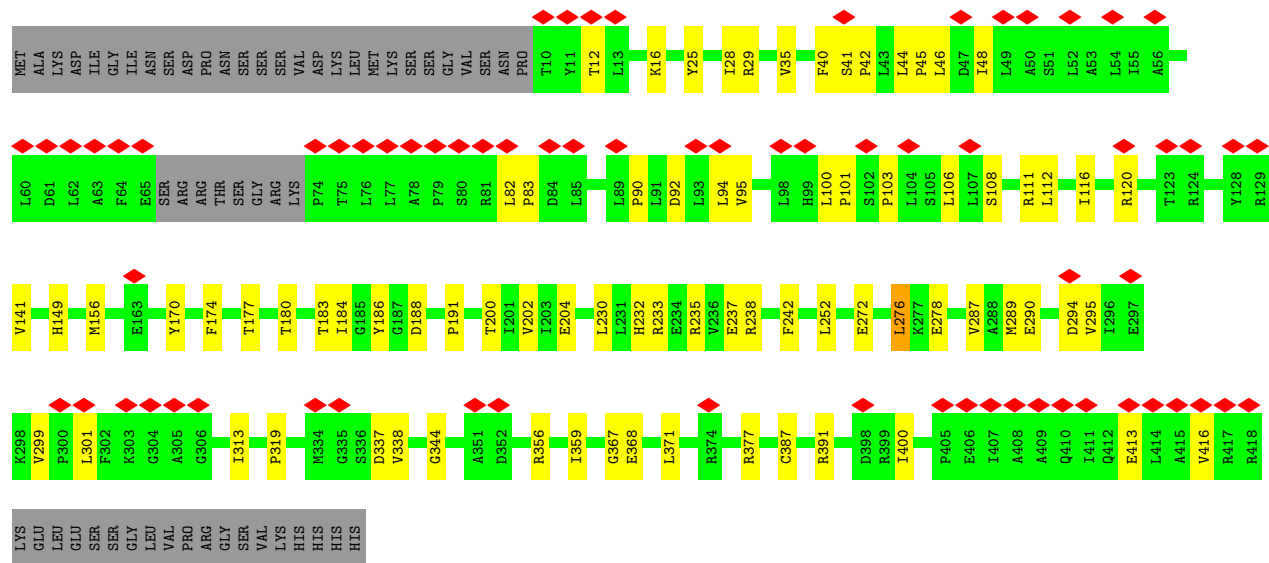
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: SthK



• Molecule 1: SthK





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C4	Depositor
Number of particles used	46786	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 ($e^-/\text{\AA}^2$)	1.365	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	130000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.262	Depositor
Minimum map value	-0.201	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.03	Depositor
Map size (Å)	254.1072, 254.1072, 254.1072	wwPDB
Map dimensions	240, 240, 240	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.05878, 1.05878, 1.05878	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: PGW, CMP

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	0.20	0/3264	0.43	1/4447 (0.0%)
1	B	0.20	0/3264	0.43	1/4447 (0.0%)
1	C	0.20	0/3264	0.43	1/4447 (0.0%)
1	D	0.20	0/3264	0.43	1/4447 (0.0%)
All	All	0.20	0/13056	0.43	4/17788 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	28	ILE	N-CA-C	-5.60	108.04	113.53
1	D	28	ILE	N-CA-C	-5.59	108.05	113.53
1	A	28	ILE	N-CA-C	-5.57	108.08	113.53
1	C	28	ILE	N-CA-C	-5.54	108.10	113.53

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	3191	0	3331	64	0
1	B	3191	0	3331	63	0
1	C	3191	0	3331	64	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3191	0	3331	62	0
2	A	22	0	11	4	0
2	B	22	0	11	3	0
2	C	22	0	11	4	0
2	D	22	0	11	4	0
3	A	223	0	341	7	0
3	B	223	0	341	7	0
3	C	223	0	341	8	0
3	D	223	0	341	7	0
All	All	13744	0	14732	233	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 8.

All (233) 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)
2:A:501:CMP:H2	2:A:501:CMP:C2	0.97	1.48
2:B:501:CMP:H2	2:B:501:CMP:C2	0.97	1.48
2:C:501:CMP:H2	2:C:501:CMP:C2	0.97	1.48
2:D:502:CMP:C2	2:D:502:CMP:H2	0.97	1.47
1:C:276:LEU:HD11	1:C:287:VAL:HG11	1.72	0.72
1:A:276:LEU:HD11	1:A:287:VAL:HG11	1.72	0.72
1:A:186:TYR:HH	1:B:180:THR:HG1	1.36	0.72
1:B:276:LEU:HD11	1:B:287:VAL:HG11	1.72	0.72
1:D:276:LEU:HD11	1:D:287:VAL:HG11	1.72	0.72
1:C:29:ARG:HH22	1:C:40:PHE:HD1	1.39	0.70
1:B:29:ARG:HH22	1:B:40:PHE:HD1	1.39	0.70
1:D:29:ARG:HH22	1:D:40:PHE:HD1	1.40	0.70
1:A:29:ARG:HH22	1:A:40:PHE:HD1	1.39	0.69
1:B:186:TYR:HH	1:C:180:THR:HG1	1.37	0.69
1:C:35:VAL:HG11	1:C:156:MET:HE3	1.75	0.68
1:B:35:VAL:HG11	1:B:156:MET:HE3	1.75	0.67
1:A:35:VAL:HG11	1:A:156:MET:HE3	1.75	0.67
1:D:35:VAL:HG11	1:D:156:MET:HE3	1.75	0.67
1:B:242:PHE:CZ	1:C:272:GLU:HG2	2.32	0.65
1:C:186:TYR:HH	1:D:180:THR:HG1	1.40	0.65
1:A:242:PHE:CZ	1:B:272:GLU:HG2	2.32	0.65
1:C:242:PHE:CZ	1:D:272:GLU:HG2	2.32	0.65
1:B:95:VAL:HG12	1:B:100:LEU:HD22	1.79	0.64
1:A:272:GLU:HG2	1:D:242:PHE:CZ	2.32	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:95:VAL:HG12	1:D:100:LEU:HD22	1.79	0.64
1:B:356:ARG:HH11	1:B:359:ILE:HD11	1.62	0.64
1:C:356:ARG:HH11	1:C:359:ILE:HD11	1.62	0.64
1:A:95:VAL:HG12	1:A:100:LEU:HD22	1.79	0.64
1:C:95:VAL:HG12	1:C:100:LEU:HD22	1.79	0.62
1:A:356:ARG:HH11	1:A:359:ILE:HD11	1.62	0.62
1:D:356:ARG:HH11	1:D:359:ILE:HD11	1.62	0.62
1:B:184:ILE:HG12	1:C:183:THR:HA	1.82	0.61
1:A:183:THR:HA	1:D:184:ILE:HG12	1.83	0.60
1:D:82:LEU:HD23	1:D:83:PRO:HD3	1.83	0.60
1:A:184:ILE:HG12	1:B:183:THR:HA	1.83	0.60
1:D:95:VAL:HG21	1:D:108:SER:HB3	1.84	0.60
1:C:82:LEU:HD23	1:C:83:PRO:HD3	1.83	0.60
1:A:82:LEU:HD23	1:A:83:PRO:HD3	1.83	0.60
1:A:95:VAL:HG21	1:A:108:SER:HB3	1.84	0.60
1:C:95:VAL:HG21	1:C:108:SER:HB3	1.84	0.60
1:B:82:LEU:HD23	1:B:83:PRO:HD3	1.83	0.59
1:B:95:VAL:HG21	1:B:108:SER:HB3	1.84	0.59
1:C:184:ILE:HG12	1:D:183:THR:HA	1.85	0.59
1:A:238:ARG:NH2	1:B:278:GLU:OE2	2.27	0.57
1:B:252:LEU:HD13	1:C:290:GLU:HG3	1.86	0.57
1:A:338:VAL:HG11	1:A:371:LEU:HD11	1.88	0.56
1:A:278:GLU:OE2	1:D:238:ARG:NH2	2.27	0.56
1:A:252:LEU:HD13	1:B:290:GLU:HG3	1.87	0.56
1:D:413:GLU:O	1:D:416:VAL:HG12	2.06	0.56
1:A:413:GLU:O	1:A:416:VAL:HG12	2.06	0.56
1:B:413:GLU:O	1:B:416:VAL:HG12	2.06	0.55
1:C:413:GLU:O	1:C:416:VAL:HG12	2.06	0.55
1:D:338:VAL:HG11	1:D:371:LEU:HD11	1.88	0.55
1:A:233:ARG:O	1:A:237:GLU:HG2	2.07	0.55
1:A:290:GLU:HG3	1:D:252:LEU:HD13	1.87	0.55
1:B:233:ARG:O	1:B:237:GLU:HG2	2.07	0.55
1:B:338:VAL:HG11	1:B:371:LEU:HD11	1.88	0.55
1:C:41:SER:OG	1:C:42:PRO:HD3	2.07	0.55
1:C:233:ARG:O	1:C:237:GLU:HG2	2.07	0.55
1:B:238:ARG:NH2	1:C:278:GLU:OE2	2.27	0.54
1:C:252:LEU:HD13	1:D:290:GLU:HG3	1.88	0.54
1:D:41:SER:OG	1:D:42:PRO:HD3	2.07	0.54
1:B:41:SER:OG	1:B:42:PRO:HD3	2.07	0.54
1:C:338:VAL:HG11	1:C:371:LEU:HD11	1.88	0.54
1:B:25:TYR:CZ	1:B:29:ARG:HD2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:25:TYR:CZ	1:D:29:ARG:HD2	2.43	0.54
1:A:141:VAL:HG11	3:A:509:PGW:H08	1.89	0.54
1:B:272:GLU:OE1	1:B:319:PRO:HG3	2.08	0.54
1:C:141:VAL:HG11	3:C:509:PGW:H08	1.90	0.54
1:A:272:GLU:OE1	1:A:319:PRO:HG3	2.08	0.53
1:C:25:TYR:CZ	1:C:29:ARG:HD2	2.43	0.53
1:A:41:SER:OG	1:A:42:PRO:HD3	2.07	0.53
1:D:299:VAL:HG12	1:D:301:LEU:HD23	1.91	0.53
1:D:272:GLU:OE1	1:D:319:PRO:HG3	2.08	0.53
1:B:141:VAL:HG11	3:B:509:PGW:H08	1.89	0.53
1:C:299:VAL:HG12	1:C:301:LEU:HD23	1.91	0.53
1:A:242:PHE:HZ	1:B:272:GLU:HG2	1.72	0.53
1:A:299:VAL:HG12	1:A:301:LEU:HD23	1.91	0.53
1:D:141:VAL:HG11	3:D:510:PGW:H08	1.89	0.53
1:D:233:ARG:O	1:D:237:GLU:HG2	2.07	0.53
1:A:25:TYR:CZ	1:A:29:ARG:HD2	2.43	0.53
1:B:299:VAL:HG12	1:B:301:LEU:HD23	1.91	0.53
1:C:272:GLU:OE1	1:C:319:PRO:HG3	2.08	0.52
1:A:276:LEU:HD11	1:A:287:VAL:CG1	2.39	0.52
1:B:276:LEU:HD11	1:B:287:VAL:CG1	2.39	0.52
1:C:242:PHE:HZ	1:D:272:GLU:HG2	1.72	0.51
3:D:507:PGW:H24A	3:D:508:PGW:H22A	1.92	0.51
1:B:242:PHE:HZ	1:C:272:GLU:HG2	1.72	0.51
1:A:272:GLU:HG2	1:D:242:PHE:HZ	1.72	0.51
1:C:112:LEU:HB2	3:C:514:PGW:H06	1.93	0.51
1:D:112:LEU:HB2	3:D:515:PGW:H06	1.93	0.51
3:A:506:PGW:H24A	3:A:507:PGW:H22A	1.92	0.51
3:B:506:PGW:H24A	3:B:507:PGW:H22A	1.92	0.51
1:C:276:LEU:HD11	1:C:287:VAL:CG1	2.39	0.50
1:D:276:LEU:HD11	1:D:287:VAL:CG1	2.39	0.50
1:A:112:LEU:HB2	3:A:514:PGW:H06	1.93	0.50
1:B:112:LEU:HB2	3:B:514:PGW:H06	1.93	0.50
1:C:45:PRO:O	1:C:48:ILE:HG22	2.12	0.50
1:C:202:VAL:HG11	3:C:512:PGW:H26	1.94	0.50
1:A:191:PRO:HG3	1:A:200:THR:OG1	2.12	0.50
3:C:506:PGW:H24A	3:C:507:PGW:H22A	1.92	0.50
1:D:313:ILE:HG22	1:D:400:ILE:HD11	1.94	0.50
1:A:202:VAL:HG11	3:A:512:PGW:H26	1.94	0.50
1:C:191:PRO:HG3	1:C:200:THR:OG1	2.12	0.49
1:D:202:VAL:HG11	3:D:513:PGW:H26	1.94	0.49
1:B:202:VAL:HG11	3:B:512:PGW:H26	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:PRO:O	1:A:48:ILE:HG22	2.12	0.49
1:C:313:ILE:HG22	1:C:400:ILE:HD11	1.94	0.49
1:D:45:PRO:O	1:D:48:ILE:HG22	2.12	0.49
1:D:92:ASP:C	1:D:94:LEU:H	2.20	0.49
1:D:368:GLU:OE1	2:D:502:CMP:O2'	2.31	0.49
1:B:45:PRO:O	1:B:48:ILE:HG22	2.12	0.49
1:D:191:PRO:HG3	1:D:200:THR:OG1	2.12	0.49
1:A:92:ASP:C	1:A:94:LEU:H	2.20	0.48
1:C:238:ARG:NH2	1:D:278:GLU:OE2	2.29	0.48
1:C:92:ASP:C	1:C:94:LEU:H	2.20	0.48
1:B:191:PRO:HG3	1:B:200:THR:OG1	2.12	0.48
1:B:313:ILE:HG22	1:B:400:ILE:HD11	1.94	0.48
1:B:188:ASP:OD1	1:B:188:ASP:N	2.46	0.48
1:A:184:ILE:HD12	1:B:204:GLU:HG2	1.96	0.48
1:A:313:ILE:HG22	1:A:400:ILE:HD11	1.94	0.48
1:B:92:ASP:C	1:B:94:LEU:H	2.20	0.48
1:B:368:GLU:OE1	2:B:501:CMP:O2'	2.31	0.47
1:C:184:ILE:HD12	1:D:204:GLU:HG2	1.96	0.47
1:B:92:ASP:OD1	1:B:108:SER:HB2	2.15	0.47
1:A:106:LEU:HD23	3:A:511:PGW:H22	1.96	0.47
1:C:106:LEU:HD23	3:C:511:PGW:H22	1.96	0.47
1:D:337:ASP:OD2	1:D:391:ARG:NH2	2.37	0.47
1:D:188:ASP:OD1	1:D:188:ASP:N	2.46	0.47
1:A:92:ASP:OD1	1:A:108:SER:HB2	2.15	0.47
1:C:92:ASP:OD1	1:C:108:SER:HB2	2.15	0.47
1:C:368:GLU:OE1	2:C:501:CMP:O2'	2.31	0.47
1:A:204:GLU:HG2	1:D:184:ILE:HD12	1.96	0.47
1:D:106:LEU:HD23	3:D:512:PGW:H22	1.96	0.47
1:C:90:PRO:HA	1:C:111:ARG:HH11	1.80	0.46
1:B:184:ILE:HD12	1:C:204:GLU:HG2	1.98	0.46
1:B:106:LEU:HD23	3:B:511:PGW:H22	1.96	0.46
1:B:90:PRO:HA	1:B:111:ARG:HH11	1.80	0.46
1:C:188:ASP:OD1	1:C:188:ASP:N	2.46	0.46
1:D:92:ASP:OD1	1:D:108:SER:HB2	2.15	0.46
1:A:337:ASP:OD2	1:A:391:ARG:NH2	2.37	0.45
1:B:337:ASP:OD2	1:B:391:ARG:NH2	2.37	0.45
1:D:289:MET:HE3	1:D:289:MET:HB3	1.87	0.45
1:A:25:TYR:CZ	1:A:46:LEU:HD22	2.51	0.45
1:B:25:TYR:CZ	1:B:46:LEU:HD22	2.51	0.45
1:C:25:TYR:CZ	1:C:46:LEU:HD22	2.51	0.45
1:D:25:TYR:CZ	1:D:46:LEU:HD22	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:PRO:HA	1:A:111:ARG:HH11	1.80	0.45
1:A:368:GLU:OE1	2:A:501:CMP:O2'	2.31	0.45
1:D:90:PRO:HA	1:D:111:ARG:HH11	1.81	0.45
1:A:90:PRO:HA	1:A:111:ARG:NH1	2.32	0.44
1:B:232:HIS:O	1:B:235:ARG:HG2	2.18	0.44
1:C:233:ARG:NH1	3:C:515:PGW:H21	2.32	0.44
1:D:101:PRO:HB2	1:D:103:PRO:HD2	1.99	0.44
1:A:233:ARG:NH1	3:A:515:PGW:H21	2.33	0.44
1:D:90:PRO:HA	1:D:111:ARG:NH1	2.32	0.44
1:D:233:ARG:NH1	3:D:501:PGW:H21	2.33	0.44
1:A:232:HIS:O	1:A:235:ARG:HG2	2.18	0.44
1:C:44:LEU:N	1:C:45:PRO:HD2	2.33	0.44
1:C:344:GLY:HA3	1:C:387:CYS:HA	2.00	0.44
1:B:233:ARG:NH1	3:B:515:PGW:H21	2.32	0.44
1:A:101:PRO:HB2	1:A:103:PRO:HD2	1.99	0.43
1:A:344:GLY:HA3	1:A:387:CYS:HA	2.00	0.43
1:B:101:PRO:HB2	1:B:103:PRO:HD2	1.99	0.43
1:B:289:MET:HE3	1:B:289:MET:HB3	1.86	0.43
1:C:232:HIS:O	1:C:235:ARG:HG2	2.18	0.43
1:D:44:LEU:N	1:D:45:PRO:HD2	2.33	0.43
1:C:90:PRO:HA	1:C:111:ARG:NH1	2.32	0.43
1:C:377:ARG:NH1	2:C:501:CMP:O2P	2.44	0.43
1:A:44:LEU:N	1:A:45:PRO:HD2	2.33	0.43
1:D:344:GLY:HA3	1:D:387:CYS:HA	2.00	0.43
1:C:12:THR:O	1:C:16:LYS:HG2	2.18	0.43
1:D:12:THR:O	1:D:16:LYS:HG2	2.18	0.43
1:D:232:HIS:O	1:D:235:ARG:HG2	2.18	0.43
1:D:149:HIS:ND1	1:D:170:TYR:OH	2.42	0.43
1:B:344:GLY:HA3	1:B:387:CYS:HA	2.00	0.43
1:A:289:MET:HE3	1:A:289:MET:HB3	1.87	0.43
1:B:90:PRO:HA	1:B:111:ARG:NH1	2.32	0.43
1:B:186:TYR:OH	1:C:180:THR:OG1	2.22	0.43
3:A:502:PGW:H16A	3:A:502:PGW:H8A	2.01	0.43
1:B:12:THR:O	1:B:16:LYS:HG2	2.18	0.43
1:C:92:ASP:C	1:C:94:LEU:N	2.78	0.42
1:A:278:GLU:HB3	1:D:235:ARG:HH21	1.84	0.42
1:B:44:LEU:N	1:B:45:PRO:HD2	2.33	0.42
1:A:12:THR:O	1:A:16:LYS:HG2	2.18	0.42
1:A:180:THR:OG1	1:D:186:TYR:OH	2.23	0.42
1:B:92:ASP:C	1:B:94:LEU:N	2.78	0.42
1:C:101:PRO:HB2	1:C:103:PRO:HD2	1.99	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:174:PHE:HA	1:C:177:THR:HG22	2.01	0.42
1:D:92:ASP:C	1:D:94:LEU:N	2.78	0.42
1:C:235:ARG:HH21	1:D:278:GLU:HB3	1.84	0.42
1:B:116:ILE:HG22	1:B:120:ARG:NH1	2.35	0.42
3:B:502:PGW:H16A	3:B:502:PGW:H8A	2.01	0.42
3:C:515:PGW:H24A	3:C:515:PGW:H21A	1.88	0.42
1:A:116:ILE:HG22	1:A:120:ARG:NH1	2.35	0.42
1:A:367:GLY:HA2	2:A:501:CMP:O2P	2.19	0.42
1:B:235:ARG:HH21	1:C:278:GLU:HB3	1.83	0.42
1:B:367:GLY:HA2	2:B:501:CMP:O2P	2.19	0.42
1:A:235:ARG:HH21	1:B:278:GLU:HB3	1.84	0.42
1:C:356:ARG:NH1	1:C:359:ILE:HD11	2.33	0.42
1:D:116:ILE:HG22	1:D:120:ARG:NH1	2.35	0.42
1:A:92:ASP:C	1:A:94:LEU:N	2.78	0.42
1:A:294:ASP:OD1	1:A:295:VAL:HG23	2.20	0.42
1:B:29:ARG:HH21	1:B:40:PHE:HA	1.85	0.42
1:D:29:ARG:HH21	1:D:40:PHE:HA	1.85	0.42
1:D:377:ARG:NH1	2:D:502:CMP:O2P	2.44	0.42
1:C:29:ARG:HH21	1:C:40:PHE:HA	1.85	0.41
1:D:174:PHE:HA	1:D:177:THR:HG22	2.01	0.41
1:D:367:GLY:HA2	2:D:502:CMP:O2P	2.20	0.41
3:D:503:PGW:H8A	3:D:503:PGW:H16A	2.01	0.41
1:A:377:ARG:NH1	2:A:501:CMP:O2P	2.44	0.41
1:B:356:ARG:NH1	1:B:359:ILE:HD11	2.33	0.41
1:D:294:ASP:OD1	1:D:295:VAL:HG23	2.20	0.41
1:B:25:TYR:OH	1:B:29:ARG:HD2	2.21	0.41
1:C:367:GLY:HA2	2:C:501:CMP:O2P	2.20	0.41
1:B:174:PHE:HA	1:B:177:THR:HG22	2.01	0.41
1:C:116:ILE:HG22	1:C:120:ARG:NH1	2.35	0.41
1:C:294:ASP:OD1	1:C:295:VAL:HG23	2.20	0.41
1:A:174:PHE:HA	1:A:177:THR:HG22	2.01	0.41
1:B:294:ASP:OD1	1:B:295:VAL:HG23	2.20	0.41
3:C:502:PGW:H8A	3:C:502:PGW:H16A	2.01	0.41
1:A:149:HIS:ND1	1:A:170:TYR:OH	2.42	0.41
1:A:229:LYS:HE3	1:A:229:LYS:HB2	1.91	0.41
1:C:25:TYR:OH	1:C:29:ARG:HD2	2.21	0.41
1:C:178:ILE:HD13	1:C:178:ILE:HA	1.93	0.41
1:A:25:TYR:OH	1:A:29:ARG:HD2	2.21	0.40
1:C:64:PHE:O	1:C:65:GLU:HG3	2.21	0.40
1:A:178:ILE:HD13	1:A:178:ILE:HA	1.93	0.40
1:A:29:ARG:HH21	1:A:40:PHE:HA	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:ILE:HD13	1:B:19:ILE:HA	1.96	0.40
1:D:356:ARG:NH1	1:D:359:ILE:HD11	2.33	0.40
1:A:356:ARG:NH1	1:A:359:ILE:HD11	2.33	0.40
1:C:176:TRP:CH2	1:C:203:ILE:HD11	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	397/456 (87%)	365 (92%)	32 (8%)	0	100	100
1	B	397/456 (87%)	366 (92%)	31 (8%)	0	100	100
1	C	397/456 (87%)	365 (92%)	32 (8%)	0	100	100
1	D	397/456 (87%)	365 (92%)	32 (8%)	0	100	100
All	All	1588/1824 (87%)	1461 (92%)	127 (8%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	344/393 (88%)	342 (99%)	2 (1%)	78	88
1	B	344/393 (88%)	342 (99%)	2 (1%)	78	88

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	344/393 (88%)	342 (99%)	2 (1%)	78	88
1	D	344/393 (88%)	342 (99%)	2 (1%)	78	88
All	All	1376/1572 (88%)	1368 (99%)	8 (1%)	76	88

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	230	LEU
1	A	276	LEU
1	B	230	LEU
1	B	276	LEU
1	C	230	LEU
1	C	276	LEU
1	D	230	LEU
1	D	276	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	GLN
1	B	159	GLN
1	C	159	GLN
1	D	159	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

60 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PGW	D	509	-	14,14,50	1.05	1 (7%)	13,13,56	1.09	1 (7%)
3	PGW	B	510	-	14,14,50	1.04	1 (7%)	13,13,56	1.14	1 (7%)
3	PGW	C	502	-	42,42,50	0.95	4 (9%)	45,48,56	1.14	2 (4%)
3	PGW	A	515	-	12,12,50	0.31	0	11,11,56	0.87	0
2	CMP	D	502	-	25,25,25	1.71	8 (32%)	37,39,39	1.93	8 (21%)
3	PGW	C	515	-	12,12,50	0.31	0	11,11,56	0.87	0
3	PGW	C	509	-	14,14,50	1.07	1 (7%)	13,13,56	1.07	1 (7%)
3	PGW	A	511	-	12,12,50	0.29	0	11,11,56	0.84	0
3	PGW	C	506	-	12,12,50	0.30	0	11,11,56	0.88	0
3	PGW	C	514	-	14,14,50	1.07	1 (7%)	13,13,56	1.10	2 (15%)
3	PGW	C	504	-	12,12,50	0.26	0	11,11,56	0.93	0
3	PGW	D	506	-	23,23,50	1.02	2 (8%)	24,24,56	1.15	2 (8%)
3	PGW	C	503	-	14,14,50	1.06	1 (7%)	13,13,56	1.09	1 (7%)
3	PGW	B	505	-	23,23,50	1.03	2 (8%)	24,24,56	1.15	2 (8%)
3	PGW	B	509	-	14,14,50	1.07	1 (7%)	13,13,56	1.07	1 (7%)
3	PGW	B	512	-	7,7,50	0.31	0	6,6,56	0.75	0
3	PGW	D	513	-	7,7,50	0.31	0	6,6,56	0.75	0
3	PGW	D	514	-	11,11,50	0.30	0	10,10,56	0.86	0
3	PGW	C	510	-	14,14,50	1.04	1 (7%)	13,13,56	1.14	1 (7%)
3	PGW	B	513	-	11,11,50	0.30	0	10,10,56	0.86	0
3	PGW	A	514	-	14,14,50	1.07	1 (7%)	13,13,56	1.10	2 (15%)
3	PGW	B	504	-	12,12,50	0.26	0	11,11,56	0.93	0
3	PGW	B	503	-	14,14,50	1.06	1 (7%)	13,13,56	1.09	1 (7%)
3	PGW	A	503	-	14,14,50	1.06	1 (7%)	13,13,56	1.09	1 (7%)
3	PGW	D	511	-	14,14,50	1.04	1 (7%)	13,13,56	1.14	1 (7%)
3	PGW	A	510	-	14,14,50	1.04	1 (7%)	13,13,56	1.14	1 (7%)
3	PGW	D	508	-	8,8,50	0.31	0	7,7,56	0.79	0
3	PGW	D	507	-	12,12,50	0.30	0	11,11,56	0.88	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PGW	B	514	-	14,14,50	1.07	1 (7%)	13,13,56	1.10	2 (15%)
2	CMP	B	501	-	25,25,25	1.71	8 (32%)	37,39,39	1.93	8 (21%)
3	PGW	D	515	-	14,14,50	1.07	1 (7%)	13,13,56	1.10	2 (15%)
3	PGW	A	512	-	7,7,50	0.31	0	6,6,56	0.75	0
3	PGW	D	504	-	14,14,50	1.06	1 (7%)	13,13,56	1.09	1 (7%)
3	PGW	C	505	-	23,23,50	1.03	2 (8%)	24,24,56	1.15	2 (8%)
3	PGW	C	508	-	14,14,50	1.05	1 (7%)	13,13,56	1.09	1 (7%)
3	PGW	C	512	-	7,7,50	0.31	0	6,6,56	0.75	0
3	PGW	A	509	-	14,14,50	1.07	1 (7%)	13,13,56	1.07	1 (7%)
3	PGW	A	504	-	12,12,50	0.26	0	11,11,56	0.93	0
3	PGW	B	502	-	42,42,50	0.95	4 (9%)	45,48,56	1.14	2 (4%)
3	PGW	C	513	-	11,11,50	0.30	0	10,10,56	0.86	0
3	PGW	B	507	-	8,8,50	0.31	0	7,7,56	0.79	0
3	PGW	B	508	-	14,14,50	1.05	1 (7%)	13,13,56	1.09	1 (7%)
3	PGW	A	505	-	23,23,50	1.03	2 (8%)	24,24,56	1.15	2 (8%)
3	PGW	A	508	-	14,14,50	1.05	1 (7%)	13,13,56	1.09	1 (7%)
3	PGW	C	507	-	8,8,50	0.31	0	7,7,56	0.79	0
3	PGW	B	515	-	12,12,50	0.30	0	11,11,56	0.86	0
3	PGW	A	513	-	11,11,50	0.30	0	10,10,56	0.86	0
3	PGW	D	501	-	12,12,50	0.30	0	11,11,56	0.87	0
3	PGW	A	507	-	8,8,50	0.31	0	7,7,56	0.79	0
2	CMP	A	501	-	25,25,25	1.71	8 (32%)	37,39,39	1.93	8 (21%)
3	PGW	D	510	-	14,14,50	1.07	1 (7%)	13,13,56	1.07	1 (7%)
2	CMP	C	501	-	25,25,25	1.71	8 (32%)	37,39,39	1.93	8 (21%)
3	PGW	D	503	-	42,42,50	0.95	4 (9%)	45,48,56	1.14	2 (4%)
3	PGW	B	511	-	12,12,50	0.29	0	11,11,56	0.84	0
3	PGW	B	506	-	12,12,50	0.30	0	11,11,56	0.88	0
3	PGW	D	512	-	12,12,50	0.29	0	11,11,56	0.84	0
3	PGW	A	506	-	12,12,50	0.30	0	11,11,56	0.88	0
3	PGW	D	505	-	12,12,50	0.26	0	11,11,56	0.93	0
3	PGW	C	511	-	12,12,50	0.29	0	11,11,56	0.84	0
3	PGW	A	502	-	42,42,50	0.95	4 (9%)	45,48,56	1.14	2 (4%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PGW	D	509	-	-	2/12/12/55	-
3	PGW	B	510	-	-	3/12/12/55	-
3	PGW	C	502	-	-	23/47/47/55	-
3	PGW	A	515	-	-	3/10/10/55	-
2	CMP	D	502	-	-	0/4/31/31	0/4/4/4
3	PGW	C	515	-	-	3/10/10/55	-
3	PGW	C	509	-	-	5/12/12/55	-
3	PGW	A	511	-	-	4/10/10/55	-
3	PGW	C	506	-	-	2/10/10/55	-
3	PGW	C	514	-	-	6/12/12/55	-
3	PGW	C	504	-	-	2/10/10/55	-
3	PGW	D	506	-	-	7/23/23/55	-
3	PGW	C	503	-	-	4/12/12/55	-
3	PGW	B	505	-	-	7/23/23/55	-
3	PGW	B	509	-	-	5/12/12/55	-
3	PGW	B	512	-	-	0/5/5/55	-
3	PGW	D	513	-	-	0/5/5/55	-
3	PGW	D	514	-	-	4/9/9/55	-
3	PGW	C	510	-	-	3/12/12/55	-
3	PGW	B	513	-	-	4/9/9/55	-
3	PGW	A	514	-	-	6/12/12/55	-
3	PGW	B	504	-	-	2/10/10/55	-
3	PGW	B	503	-	-	4/12/12/55	-
3	PGW	A	503	-	-	4/12/12/55	-
3	PGW	D	511	-	-	3/12/12/55	-
3	PGW	A	510	-	-	3/12/12/55	-
3	PGW	D	508	-	-	1/6/6/55	-
3	PGW	D	507	-	-	2/10/10/55	-
3	PGW	B	514	-	-	6/12/12/55	-
2	CMP	B	501	-	-	0/4/31/31	0/4/4/4
3	PGW	D	515	-	-	6/12/12/55	-
3	PGW	A	512	-	-	0/5/5/55	-
3	PGW	D	504	-	-	4/12/12/55	-
3	PGW	C	505	-	-	7/23/23/55	-
3	PGW	C	508	-	-	2/12/12/55	-
3	PGW	C	512	-	-	0/5/5/55	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PGW	A	509	-	-	5/12/12/55	-
3	PGW	A	504	-	-	2/10/10/55	-
3	PGW	B	502	-	-	23/47/47/55	-
3	PGW	C	513	-	-	4/9/9/55	-
3	PGW	B	507	-	-	1/6/6/55	-
3	PGW	B	508	-	-	2/12/12/55	-
3	PGW	A	505	-	-	7/23/23/55	-
3	PGW	A	508	-	-	2/12/12/55	-
3	PGW	C	507	-	-	1/6/6/55	-
3	PGW	B	515	-	-	3/10/10/55	-
3	PGW	A	513	-	-	4/9/9/55	-
3	PGW	D	501	-	-	3/10/10/55	-
3	PGW	A	507	-	-	1/6/6/55	-
2	CMP	A	501	-	-	0/4/31/31	0/4/4/4
3	PGW	D	510	-	-	5/12/12/55	-
2	CMP	C	501	-	-	0/4/31/31	0/4/4/4
3	PGW	D	503	-	-	23/47/47/55	-
3	PGW	B	511	-	-	4/10/10/55	-
3	PGW	B	506	-	-	2/10/10/55	-
3	PGW	D	512	-	-	4/10/10/55	-
3	PGW	A	506	-	-	2/10/10/55	-
3	PGW	D	505	-	-	2/10/10/55	-
3	PGW	C	511	-	-	4/10/10/55	-
3	PGW	A	502	-	-	23/47/47/55	-

All (76) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	502	CMP	C5-C4	4.70	1.47	1.39
2	C	501	CMP	C5-C4	4.66	1.47	1.39
2	A	501	CMP	C5-C4	4.65	1.47	1.39
2	B	501	CMP	C5-C4	4.65	1.47	1.39
3	D	515	PGW	C10-C9	3.82	1.53	1.31
3	A	514	PGW	C10-C9	3.81	1.53	1.31
3	B	514	PGW	C10-C9	3.81	1.53	1.31
3	C	514	PGW	C10-C9	3.80	1.53	1.31
3	A	509	PGW	C10-C9	3.80	1.53	1.31
3	B	509	PGW	C10-C9	3.80	1.53	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	510	PGW	C10-C9	3.80	1.53	1.31
3	C	509	PGW	C10-C9	3.80	1.53	1.31
3	C	503	PGW	C10-C9	3.77	1.53	1.31
3	A	503	PGW	C10-C9	3.76	1.53	1.31
3	B	503	PGW	C10-C9	3.76	1.53	1.31
3	D	504	PGW	C10-C9	3.76	1.53	1.31
3	A	508	PGW	C10-C9	3.74	1.53	1.31
3	B	508	PGW	C10-C9	3.74	1.53	1.31
3	D	509	PGW	C10-C9	3.74	1.52	1.31
3	C	508	PGW	C10-C9	3.73	1.52	1.31
3	A	510	PGW	C10-C9	3.72	1.52	1.31
3	B	510	PGW	C10-C9	3.72	1.52	1.31
3	D	511	PGW	C10-C9	3.72	1.52	1.31
3	C	510	PGW	C10-C9	3.71	1.52	1.31
2	A	501	CMP	P-O3'	2.75	1.62	1.57
2	B	501	CMP	P-O3'	2.75	1.62	1.57
2	C	501	CMP	P-O3'	2.75	1.62	1.57
2	D	502	CMP	P-O3'	2.75	1.62	1.57
2	C	501	CMP	C5-C6	2.65	1.48	1.41
2	A	501	CMP	C5-C6	2.63	1.48	1.41
2	B	501	CMP	C5-C6	2.62	1.48	1.41
2	D	502	CMP	C5-C6	2.61	1.48	1.41
3	A	502	PGW	O01-C02	-2.60	1.40	1.46
3	B	502	PGW	O01-C02	-2.60	1.40	1.46
3	C	502	PGW	O01-C02	-2.57	1.40	1.46
3	D	503	PGW	O01-C02	-2.57	1.40	1.46
2	C	501	CMP	P-O5'	2.48	1.60	1.57
2	A	501	CMP	P-O5'	2.45	1.60	1.57
2	B	501	CMP	P-O5'	2.45	1.60	1.57
2	D	502	CMP	P-O5'	2.45	1.60	1.57
2	A	501	CMP	O5'-C5'	-2.42	1.42	1.46
2	B	501	CMP	O5'-C5'	-2.42	1.42	1.46
2	C	501	CMP	O5'-C5'	-2.42	1.42	1.46
2	D	502	CMP	O5'-C5'	-2.42	1.42	1.46
2	A	501	CMP	C5-N7	-2.42	1.34	1.39
2	B	501	CMP	C5-N7	-2.42	1.34	1.39
2	C	501	CMP	C5-N7	-2.42	1.34	1.39
2	D	502	CMP	C5-N7	-2.42	1.34	1.39
3	C	505	PGW	O03-C19	2.42	1.40	1.33
3	D	506	PGW	O01-C1	2.41	1.40	1.33
3	A	505	PGW	O01-C1	2.40	1.40	1.33
3	C	505	PGW	O01-C1	2.40	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	505	PGW	O01-C1	2.39	1.40	1.33
3	A	505	PGW	O03-C19	2.39	1.40	1.33
3	B	505	PGW	O03-C19	2.39	1.40	1.33
3	D	506	PGW	O03-C19	2.37	1.40	1.33
3	C	502	PGW	O03-C19	2.36	1.40	1.33
3	A	502	PGW	O03-C19	2.35	1.40	1.33
3	B	502	PGW	O03-C19	2.35	1.40	1.33
3	D	503	PGW	O03-C19	2.35	1.40	1.33
2	A	501	CMP	C8-N7	2.30	1.36	1.31
2	B	501	CMP	C8-N7	2.30	1.36	1.31
2	C	501	CMP	C8-N7	2.30	1.36	1.31
2	D	502	CMP	C8-N7	2.30	1.36	1.31
3	A	502	PGW	O03-C01	-2.26	1.40	1.45
3	B	502	PGW	O03-C01	-2.26	1.40	1.45
3	C	502	PGW	O03-C01	-2.26	1.40	1.45
3	D	503	PGW	O03-C01	-2.26	1.40	1.45
3	A	502	PGW	O01-C1	2.11	1.40	1.34
3	B	502	PGW	O01-C1	2.11	1.40	1.34
3	C	502	PGW	O01-C1	2.11	1.40	1.34
3	D	503	PGW	O01-C1	2.11	1.40	1.34
2	A	501	CMP	O3'-C3'	-2.03	1.41	1.44
2	B	501	CMP	O3'-C3'	-2.03	1.41	1.44
2	C	501	CMP	O3'-C3'	-2.03	1.41	1.44
2	D	502	CMP	O3'-C3'	-2.03	1.41	1.44

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	502	CMP	C5-C4-N3	-5.92	118.57	126.72
2	B	501	CMP	C5-C4-N3	-5.91	118.58	126.72
2	A	501	CMP	C5-C4-N3	-5.91	118.58	126.72
2	C	501	CMP	C5-C4-N3	-5.86	118.64	126.72
2	B	501	CMP	N3-C4-N9	4.64	135.05	127.17
2	A	501	CMP	N3-C4-N9	4.63	135.04	127.17
2	D	502	CMP	N3-C4-N9	4.63	135.04	127.17
2	C	501	CMP	N3-C4-N9	4.61	135.01	127.17
3	A	502	PGW	O01-C1-C2	4.24	120.66	111.48
3	B	502	PGW	O01-C1-C2	4.24	120.66	111.48
3	C	502	PGW	O01-C1-C2	4.24	120.66	111.48
3	D	503	PGW	O01-C1-C2	4.24	120.66	111.48
2	B	501	CMP	C2-N3-C4	3.75	120.99	111.83
2	A	501	CMP	C2-N3-C4	3.75	120.98	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	502	CMP	C2-N3-C4	3.75	120.98	111.83
2	C	501	CMP	C2-N3-C4	3.74	120.98	111.83
2	C	501	CMP	O2P-P-O1P	3.45	119.17	108.56
2	D	502	CMP	O2P-P-O1P	3.44	119.14	108.56
2	A	501	CMP	O2P-P-O1P	3.44	119.14	108.56
2	B	501	CMP	O2P-P-O1P	3.44	119.13	108.56
2	D	502	CMP	C4-C5-N7	-3.32	106.79	110.58
2	A	501	CMP	C4-C5-N7	-3.30	106.81	110.58
2	B	501	CMP	C4-C5-N7	-3.30	106.81	110.58
2	C	501	CMP	C4-C5-N7	-3.28	106.83	110.58
2	C	501	CMP	N3-C2-N1	-3.27	123.63	128.58
2	A	501	CMP	N3-C2-N1	-3.25	123.66	128.58
2	B	501	CMP	N3-C2-N1	-3.25	123.66	128.58
2	D	502	CMP	N3-C2-N1	-3.23	123.69	128.58
3	A	502	PGW	O03-C19-C20	2.62	119.83	111.83
3	B	502	PGW	O03-C19-C20	2.62	119.83	111.83
3	C	502	PGW	O03-C19-C20	2.62	119.81	111.83
3	D	503	PGW	O03-C19-C20	2.61	119.80	111.83
3	A	505	PGW	O01-C1-C2	2.56	119.66	111.83
3	B	505	PGW	O01-C1-C2	2.56	119.64	111.83
3	D	506	PGW	O01-C1-C2	2.56	119.63	111.83
3	C	505	PGW	O01-C1-C2	2.56	119.63	111.83
3	D	506	PGW	O03-C19-C20	2.51	119.48	111.83
3	C	505	PGW	O03-C19-C20	2.50	119.46	111.83
3	A	505	PGW	O03-C19-C20	2.50	119.44	111.83
3	B	505	PGW	O03-C19-C20	2.50	119.44	111.83
2	A	501	CMP	C5-N7-C8	2.40	107.22	103.45
2	B	501	CMP	C5-N7-C8	2.40	107.22	103.45
2	C	501	CMP	C5-N7-C8	2.40	107.22	103.45
2	D	502	CMP	C5-N7-C8	2.40	107.22	103.45
2	C	501	CMP	C4-N9-C8	2.38	108.24	105.74
2	A	501	CMP	C4-N9-C8	2.36	108.22	105.74
2	B	501	CMP	C4-N9-C8	2.36	108.22	105.74
3	A	510	PGW	C06-C10-C9	-2.34	107.33	124.83
3	B	510	PGW	C06-C10-C9	-2.34	107.33	124.83
3	D	511	PGW	C06-C10-C9	-2.34	107.33	124.83
3	C	510	PGW	C06-C10-C9	-2.33	107.36	124.83
2	D	502	CMP	C4-N9-C8	2.33	108.18	105.74
3	D	504	PGW	C06-C10-C9	-2.19	108.41	124.83
3	C	503	PGW	C06-C10-C9	-2.19	108.42	124.83
3	A	503	PGW	C06-C10-C9	-2.19	108.42	124.83
3	B	503	PGW	C06-C10-C9	-2.19	108.42	124.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	508	PGW	C06-C10-C9	-2.15	108.70	124.83
3	A	508	PGW	C06-C10-C9	-2.15	108.70	124.83
3	B	508	PGW	C06-C10-C9	-2.15	108.70	124.83
3	D	509	PGW	C06-C10-C9	-2.15	108.72	124.83
3	A	514	PGW	C06-C10-C9	-2.14	108.76	124.83
3	B	514	PGW	C06-C10-C9	-2.14	108.76	124.83
3	C	514	PGW	C06-C10-C9	-2.14	108.77	124.83
3	D	515	PGW	C06-C10-C9	-2.14	108.78	124.83
3	A	509	PGW	C06-C10-C9	-2.14	108.79	124.83
3	B	509	PGW	C06-C10-C9	-2.14	108.79	124.83
3	D	510	PGW	C06-C10-C9	-2.14	108.79	124.83
3	C	509	PGW	C06-C10-C9	-2.14	108.83	124.83
3	D	515	PGW	C8-C9-C10	-2.04	109.51	124.83
3	A	514	PGW	C8-C9-C10	-2.04	109.52	124.83
3	B	514	PGW	C8-C9-C10	-2.04	109.52	124.83
3	C	514	PGW	C8-C9-C10	-2.04	109.56	124.83

There are no chirality outliers.

All (264) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	PGW	C03-O11-P-O12
3	A	502	PGW	C03-O11-P-O13
3	A	502	PGW	C03-O11-P-O14
3	A	502	PGW	O02-C1-O01-C02
3	B	502	PGW	C03-O11-P-O12
3	B	502	PGW	C03-O11-P-O13
3	B	502	PGW	C03-O11-P-O14
3	B	502	PGW	O02-C1-O01-C02
3	C	502	PGW	C03-O11-P-O12
3	C	502	PGW	C03-O11-P-O13
3	C	502	PGW	C03-O11-P-O14
3	C	502	PGW	O02-C1-O01-C02
3	D	503	PGW	C03-O11-P-O12
3	D	503	PGW	C03-O11-P-O13
3	D	503	PGW	C03-O11-P-O14
3	D	503	PGW	O02-C1-O01-C02
3	A	502	PGW	C2-C1-O01-C02
3	B	502	PGW	C2-C1-O01-C02
3	C	502	PGW	C2-C1-O01-C02
3	D	503	PGW	C2-C1-O01-C02
3	A	502	PGW	C20-C19-O03-C01

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Mol	Chain	Res	Type	Atoms
3	B	502	PGW	C20-C19-O03-C01
3	C	502	PGW	C20-C19-O03-C01
3	D	503	PGW	C20-C19-O03-C01
3	A	503	PGW	C06-C10-C9-C8
3	B	503	PGW	C06-C10-C9-C8
3	C	503	PGW	C06-C10-C9-C8
3	D	504	PGW	C06-C10-C9-C8
3	A	502	PGW	O04-C19-O03-C01
3	B	502	PGW	O04-C19-O03-C01
3	C	502	PGW	O04-C19-O03-C01
3	D	503	PGW	O04-C19-O03-C01
3	A	505	PGW	C2-C1-O01-C02
3	B	505	PGW	C2-C1-O01-C02
3	C	505	PGW	C2-C1-O01-C02
3	D	506	PGW	C2-C1-O01-C02
3	A	502	PGW	OAF-C05-CAD-OAE
3	B	502	PGW	OAF-C05-CAD-OAE
3	C	502	PGW	OAF-C05-CAD-OAE
3	D	503	PGW	OAF-C05-CAD-OAE
3	A	505	PGW	C19-C20-C21-C22
3	B	505	PGW	C19-C20-C21-C22
3	C	505	PGW	C19-C20-C21-C22
3	D	506	PGW	C19-C20-C21-C22
3	A	505	PGW	O02-C1-O01-C02
3	B	505	PGW	O02-C1-O01-C02
3	C	505	PGW	O02-C1-O01-C02
3	D	506	PGW	O02-C1-O01-C02
3	A	502	PGW	C04-C05-CAD-OAE
3	B	502	PGW	C04-C05-CAD-OAE
3	C	502	PGW	C04-C05-CAD-OAE
3	D	503	PGW	C04-C05-CAD-OAE
3	A	502	PGW	C4-C5-C6-C7
3	B	502	PGW	C4-C5-C6-C7
3	C	502	PGW	C4-C5-C6-C7
3	D	503	PGW	C4-C5-C6-C7
3	C	515	PGW	C24-C25-C26-C27
3	A	515	PGW	C24-C25-C26-C27
3	B	515	PGW	C24-C25-C26-C27
3	D	501	PGW	C24-C25-C26-C27
3	A	511	PGW	C22-C23-C24-C25
3	B	511	PGW	C22-C23-C24-C25
3	C	511	PGW	C22-C23-C24-C25

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Mol	Chain	Res	Type	Atoms
3	D	512	PGW	C22-C23-C24-C25
3	A	511	PGW	C20-C21-C22-C23
3	B	511	PGW	C20-C21-C22-C23
3	C	511	PGW	C20-C21-C22-C23
3	D	512	PGW	C20-C21-C22-C23
3	D	509	PGW	C08-C09-C11-C12
3	A	503	PGW	C6-C7-C8-C9
3	B	503	PGW	C6-C7-C8-C9
3	C	503	PGW	C6-C7-C8-C9
3	D	504	PGW	C6-C7-C8-C9
3	A	505	PGW	C21-C22-C23-C24
3	A	508	PGW	C08-C09-C11-C12
3	B	505	PGW	C21-C22-C23-C24
3	B	508	PGW	C08-C09-C11-C12
3	C	502	PGW	C2-C3-C4-C5
3	C	505	PGW	C21-C22-C23-C24
3	C	508	PGW	C08-C09-C11-C12
3	D	506	PGW	C21-C22-C23-C24
3	A	502	PGW	C2-C3-C4-C5
3	B	502	PGW	C2-C3-C4-C5
3	D	503	PGW	C2-C3-C4-C5
3	B	502	PGW	C20-C21-C22-C23
3	A	502	PGW	C20-C21-C22-C23
3	C	502	PGW	C20-C21-C22-C23
3	D	503	PGW	C20-C21-C22-C23
3	A	505	PGW	O03-C01-C02-O01
3	B	505	PGW	O03-C01-C02-O01
3	C	505	PGW	O03-C01-C02-O01
3	D	506	PGW	O03-C01-C02-O01
3	A	510	PGW	C06-C07-C08-C09
3	B	510	PGW	C06-C07-C08-C09
3	C	510	PGW	C06-C07-C08-C09
3	D	511	PGW	C06-C07-C08-C09
3	A	505	PGW	C20-C19-O03-C01
3	B	505	PGW	C20-C19-O03-C01
3	C	505	PGW	C20-C19-O03-C01
3	D	506	PGW	C20-C19-O03-C01
3	A	508	PGW	C6-C7-C8-C9
3	A	514	PGW	C10-C06-C07-C08
3	B	508	PGW	C6-C7-C8-C9
3	B	514	PGW	C10-C06-C07-C08
3	C	508	PGW	C6-C7-C8-C9

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Mol	Chain	Res	Type	Atoms
3	C	514	PGW	C10-C06-C07-C08
3	D	509	PGW	C6-C7-C8-C9
3	D	515	PGW	C10-C06-C07-C08
3	A	514	PGW	C06-C07-C08-C09
3	B	514	PGW	C06-C07-C08-C09
3	C	514	PGW	C06-C07-C08-C09
3	D	515	PGW	C06-C07-C08-C09
3	A	504	PGW	C27-C15-C16-C17
3	B	504	PGW	C27-C15-C16-C17
3	C	504	PGW	C27-C15-C16-C17
3	D	505	PGW	C27-C15-C16-C17
3	A	506	PGW	C24-C25-C26-C27
3	C	506	PGW	C24-C25-C26-C27
3	D	507	PGW	C24-C25-C26-C27
3	B	506	PGW	C24-C25-C26-C27
3	A	505	PGW	O04-C19-O03-C01
3	B	505	PGW	O04-C19-O03-C01
3	C	505	PGW	O04-C19-O03-C01
3	C	503	PGW	C3-C4-C5-C6
3	A	503	PGW	C3-C4-C5-C6
3	B	503	PGW	C3-C4-C5-C6
3	D	504	PGW	C3-C4-C5-C6
3	D	506	PGW	O04-C19-O03-C01
3	A	502	PGW	O03-C01-C02-C03
3	B	502	PGW	O03-C01-C02-C03
3	C	502	PGW	O03-C01-C02-C03
3	D	503	PGW	O03-C01-C02-C03
3	A	502	PGW	C21-C22-C23-C24
3	B	502	PGW	C21-C22-C23-C24
3	C	502	PGW	C21-C22-C23-C24
3	D	503	PGW	C21-C22-C23-C24
3	A	514	PGW	C4-C5-C6-C7
3	B	514	PGW	C4-C5-C6-C7
3	C	514	PGW	C4-C5-C6-C7
3	D	515	PGW	C4-C5-C6-C7
3	A	504	PGW	C23-C24-C25-C26
3	C	504	PGW	C23-C24-C25-C26
3	D	505	PGW	C23-C24-C25-C26
3	B	504	PGW	C23-C24-C25-C26
3	A	509	PGW	C06-C07-C08-C09
3	B	509	PGW	C06-C07-C08-C09
3	C	509	PGW	C06-C07-C08-C09

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Mol	Chain	Res	Type	Atoms
3	D	510	PGW	C06-C07-C08-C09
3	A	502	PGW	C3-C4-C5-C6
3	B	502	PGW	C3-C4-C5-C6
3	C	502	PGW	C3-C4-C5-C6
3	D	503	PGW	C3-C4-C5-C6
3	A	502	PGW	O03-C01-C02-O01
3	B	502	PGW	O03-C01-C02-O01
3	C	502	PGW	O03-C01-C02-O01
3	D	503	PGW	O03-C01-C02-O01
3	B	513	PGW	C24-C25-C26-C27
3	D	514	PGW	C24-C25-C26-C27
3	A	513	PGW	C24-C25-C26-C27
3	C	513	PGW	C24-C25-C26-C27
3	A	515	PGW	C16-C15-C27-C26
3	D	501	PGW	C16-C15-C27-C26
3	B	515	PGW	C16-C15-C27-C26
3	C	515	PGW	C16-C15-C27-C26
3	A	502	PGW	C7-C8-C9-C10
3	C	502	PGW	C7-C8-C9-C10
3	D	503	PGW	C7-C8-C9-C10
3	A	502	PGW	C04-O12-P-O14
3	B	502	PGW	C04-O12-P-O14
3	C	502	PGW	C04-O12-P-O14
3	D	503	PGW	C04-O12-P-O14
3	B	502	PGW	C7-C8-C9-C10
3	A	509	PGW	C2-C3-C4-C5
3	B	509	PGW	C2-C3-C4-C5
3	C	509	PGW	C2-C3-C4-C5
3	D	510	PGW	C2-C3-C4-C5
3	A	502	PGW	C07-C06-C10-C9
3	B	502	PGW	C07-C06-C10-C9
3	C	502	PGW	C07-C06-C10-C9
3	D	503	PGW	C07-C06-C10-C9
3	A	513	PGW	C23-C24-C25-C26
3	B	513	PGW	C23-C24-C25-C26
3	C	513	PGW	C23-C24-C25-C26
3	D	514	PGW	C23-C24-C25-C26
3	A	514	PGW	C7-C8-C9-C10
3	B	514	PGW	C7-C8-C9-C10
3	C	514	PGW	C7-C8-C9-C10
3	D	515	PGW	C7-C8-C9-C10
3	A	514	PGW	C07-C08-C09-C11

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Mol	Chain	Res	Type	Atoms
3	C	511	PGW	C23-C24-C25-C26
3	D	515	PGW	C07-C08-C09-C11
3	A	511	PGW	C23-C24-C25-C26
3	B	511	PGW	C23-C24-C25-C26
3	C	511	PGW	C25-C26-C27-C15
3	D	512	PGW	C23-C24-C25-C26
3	D	512	PGW	C25-C26-C27-C15
3	B	514	PGW	C07-C08-C09-C11
3	C	514	PGW	C07-C08-C09-C11
3	A	511	PGW	C25-C26-C27-C15
3	B	511	PGW	C25-C26-C27-C15
3	B	506	PGW	C22-C23-C24-C25
3	A	506	PGW	C22-C23-C24-C25
3	D	507	PGW	C22-C23-C24-C25
3	C	506	PGW	C22-C23-C24-C25
3	A	509	PGW	C7-C8-C9-C10
3	A	510	PGW	C07-C06-C10-C9
3	B	509	PGW	C7-C8-C9-C10
3	B	510	PGW	C07-C06-C10-C9
3	C	509	PGW	C7-C8-C9-C10
3	C	510	PGW	C07-C06-C10-C9
3	D	510	PGW	C7-C8-C9-C10
3	D	511	PGW	C07-C06-C10-C9
3	A	515	PGW	C20-C21-C22-C23
3	C	515	PGW	C20-C21-C22-C23
3	D	501	PGW	C20-C21-C22-C23
3	B	515	PGW	C20-C21-C22-C23
3	A	509	PGW	C10-C06-C07-C08
3	B	509	PGW	C10-C06-C07-C08
3	C	509	PGW	C10-C06-C07-C08
3	D	510	PGW	C10-C06-C07-C08
3	A	510	PGW	C7-C8-C9-C10
3	A	514	PGW	C07-C06-C10-C9
3	B	510	PGW	C7-C8-C9-C10
3	B	514	PGW	C07-C06-C10-C9
3	C	510	PGW	C7-C8-C9-C10
3	C	514	PGW	C07-C06-C10-C9
3	D	511	PGW	C7-C8-C9-C10
3	A	513	PGW	C22-C23-C24-C25
3	B	513	PGW	C22-C23-C24-C25
3	D	514	PGW	C22-C23-C24-C25
3	C	513	PGW	C22-C23-C24-C25

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Mol	Chain	Res	Type	Atoms
3	C	502	PGW	O03-C19-C20-C21
3	D	503	PGW	O03-C19-C20-C21
3	A	503	PGW	C7-C8-C9-C10
3	B	503	PGW	C7-C8-C9-C10
3	C	503	PGW	C7-C8-C9-C10
3	D	504	PGW	C7-C8-C9-C10
3	D	515	PGW	C07-C06-C10-C9
3	A	502	PGW	O03-C19-C20-C21
3	B	502	PGW	O03-C19-C20-C21
3	A	507	PGW	C24-C25-C26-C27
3	B	507	PGW	C24-C25-C26-C27
3	C	507	PGW	C24-C25-C26-C27
3	D	508	PGW	C24-C25-C26-C27
3	A	513	PGW	C16-C17-C18-C28
3	B	513	PGW	C16-C17-C18-C28
3	C	513	PGW	C16-C17-C18-C28
3	D	514	PGW	C16-C17-C18-C28
3	C	509	PGW	C3-C4-C5-C6
3	D	510	PGW	C3-C4-C5-C6
3	A	509	PGW	C3-C4-C5-C6
3	B	509	PGW	C3-C4-C5-C6
3	A	502	PGW	C16-C17-C18-C28
3	D	503	PGW	C16-C17-C18-C28
3	B	502	PGW	C16-C17-C18-C28
3	C	502	PGW	C16-C17-C18-C28
3	B	502	PGW	O04-C19-C20-C21
3	A	502	PGW	O04-C19-C20-C21
3	D	503	PGW	O04-C19-C20-C21
3	A	502	PGW	C24-C25-C26-C27
3	B	502	PGW	C24-C25-C26-C27
3	C	502	PGW	C24-C25-C26-C27
3	C	502	PGW	O04-C19-C20-C21
3	D	503	PGW	C24-C25-C26-C27

There are no ring outliers.

36 monomers are involved in 44 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	502	PGW	1	0
3	A	515	PGW	1	0
2	D	502	CMP	4	0
3	C	515	PGW	2	0

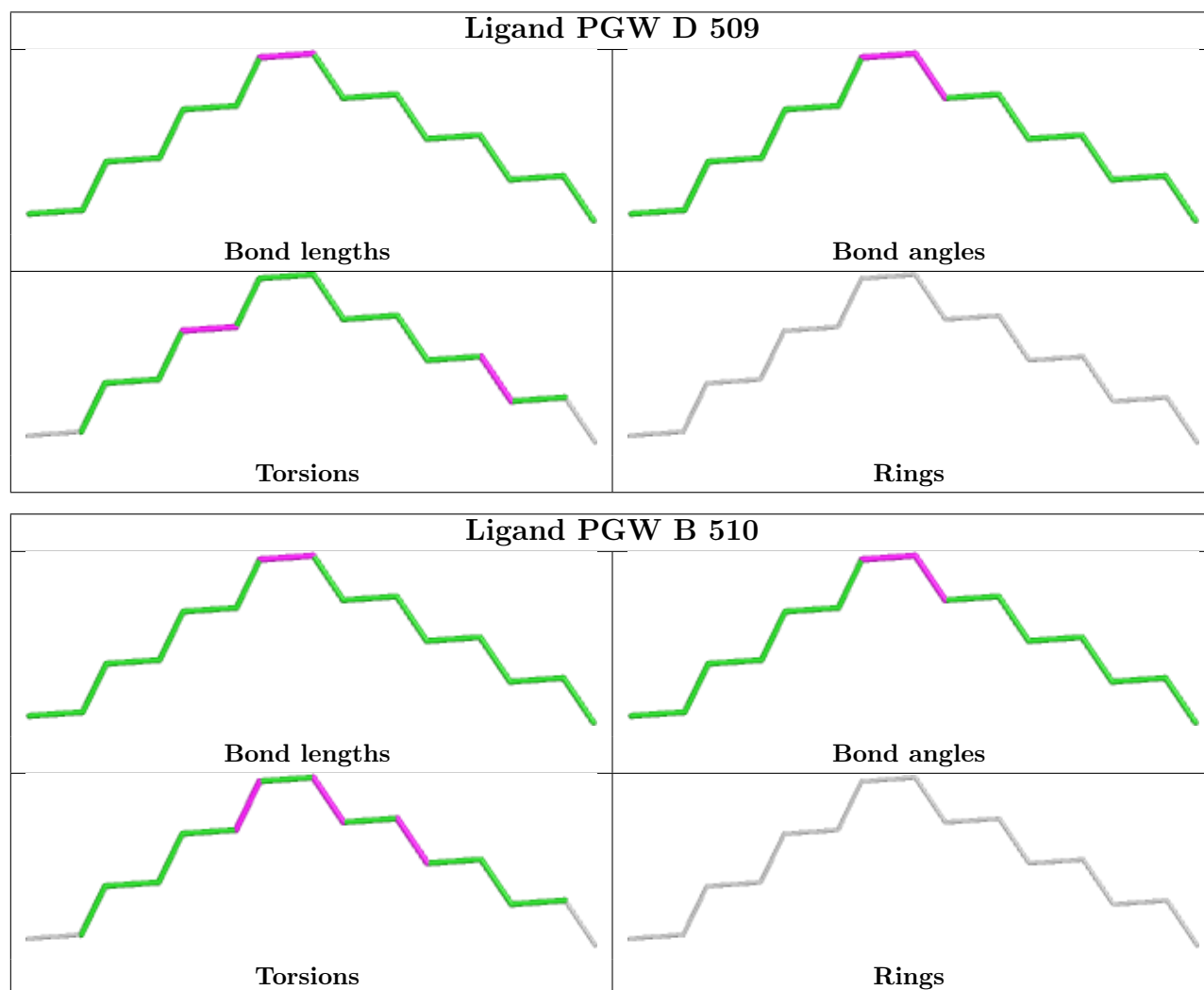
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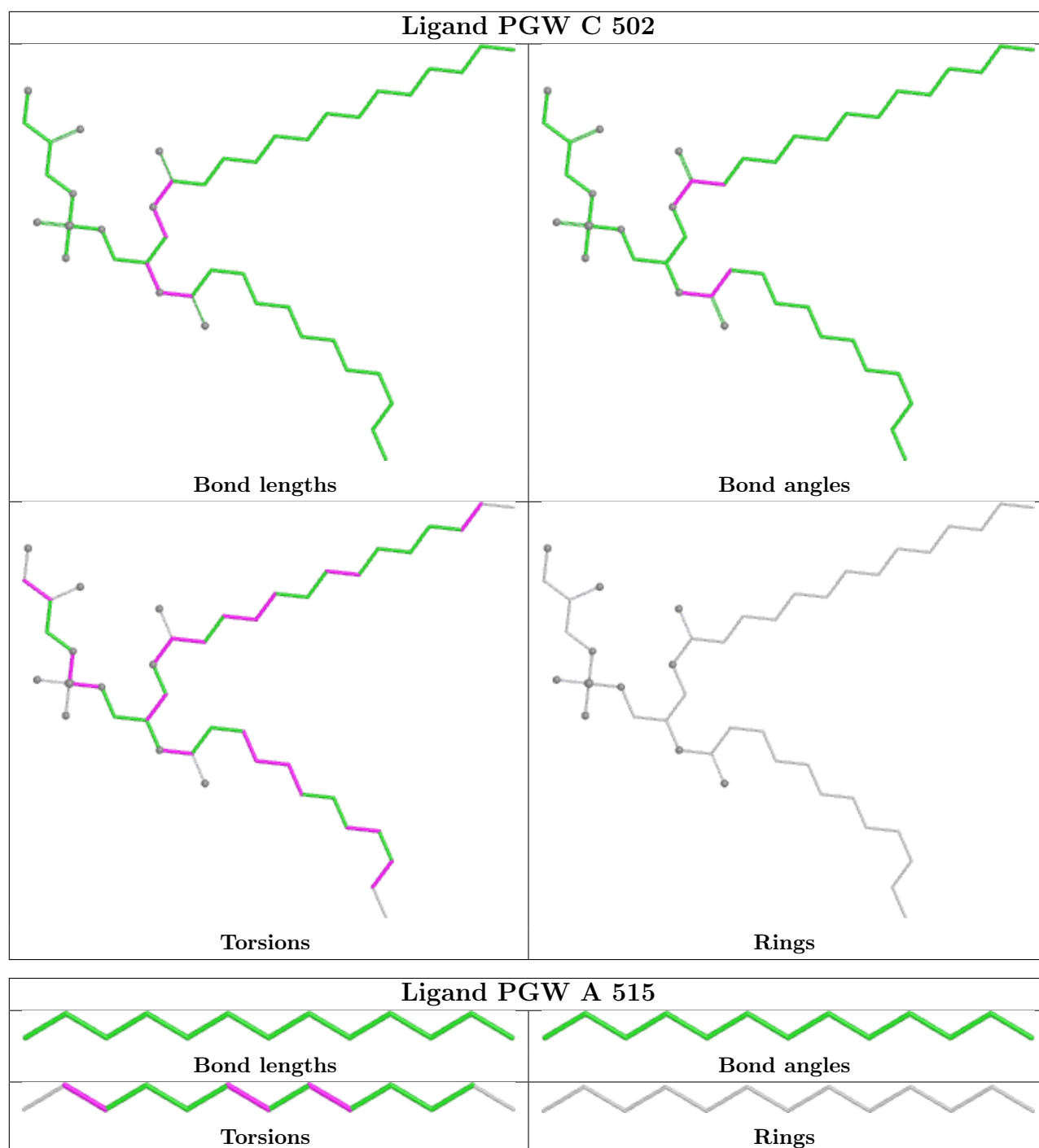
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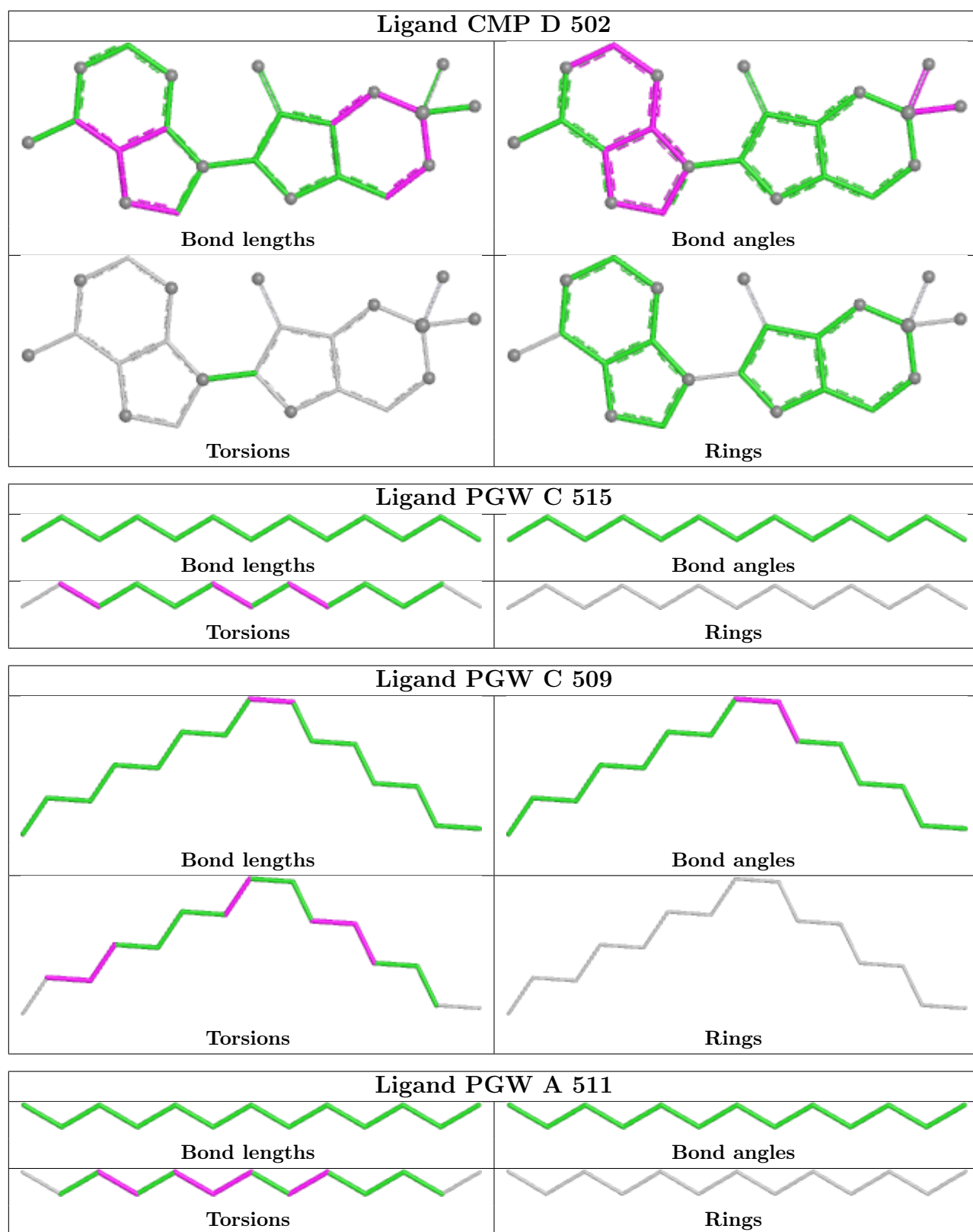
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	509	PGW	1	0
3	A	511	PGW	1	0
3	C	506	PGW	1	0
3	C	514	PGW	1	0
3	B	509	PGW	1	0
3	B	512	PGW	1	0
3	D	513	PGW	1	0
3	A	514	PGW	1	0
3	D	508	PGW	1	0
3	D	507	PGW	1	0
3	B	514	PGW	1	0
2	B	501	CMP	3	0
3	D	515	PGW	1	0
3	A	512	PGW	1	0
3	C	512	PGW	1	0
3	A	509	PGW	1	0
3	B	502	PGW	1	0
3	B	507	PGW	1	0
3	C	507	PGW	1	0
3	B	515	PGW	1	0
3	D	501	PGW	1	0
3	A	507	PGW	1	0
2	A	501	CMP	4	0
3	D	510	PGW	1	0
2	C	501	CMP	4	0
3	D	503	PGW	1	0
3	B	511	PGW	1	0
3	B	506	PGW	1	0
3	D	512	PGW	1	0
3	A	506	PGW	1	0
3	C	511	PGW	1	0
3	A	502	PGW	1	0

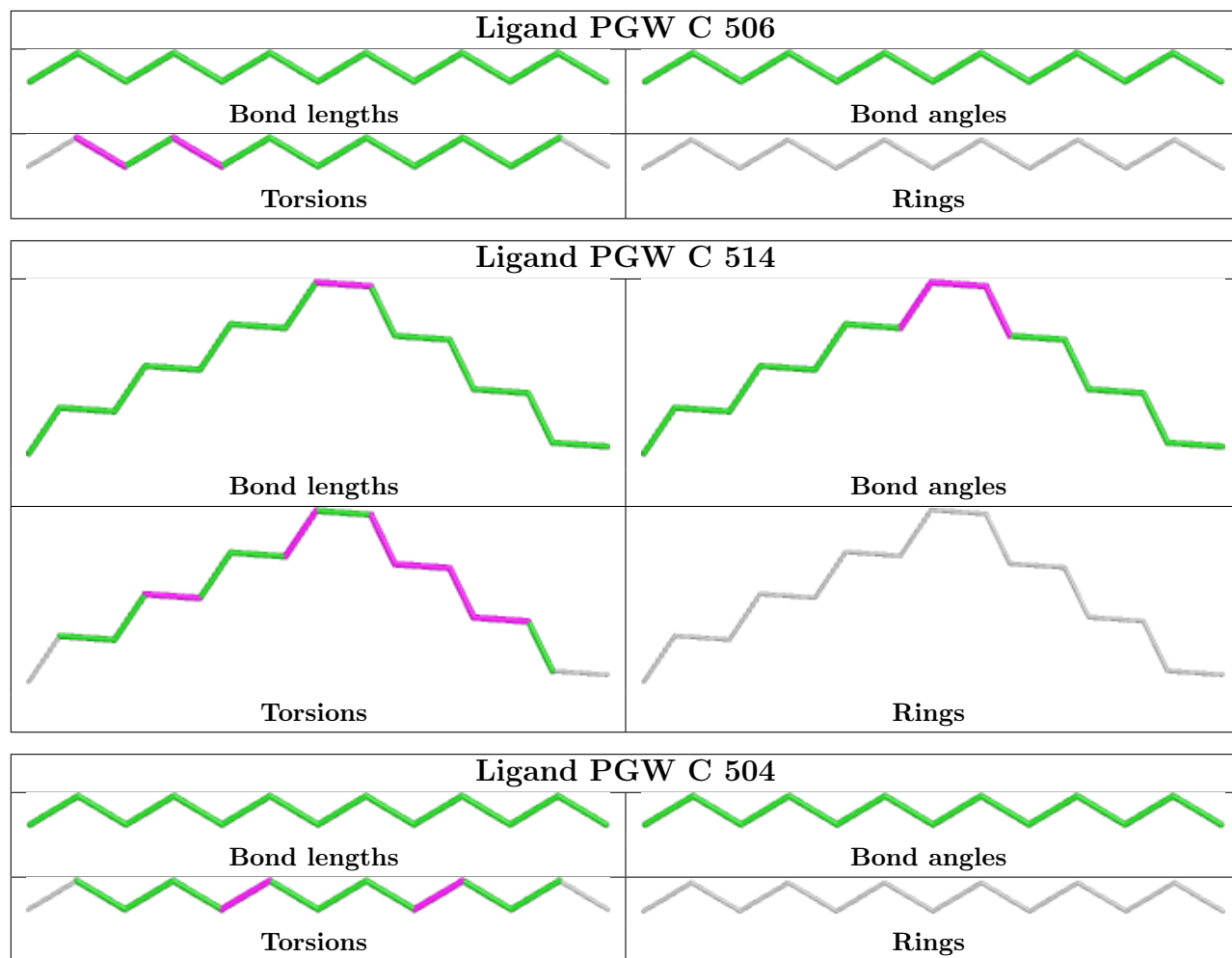
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

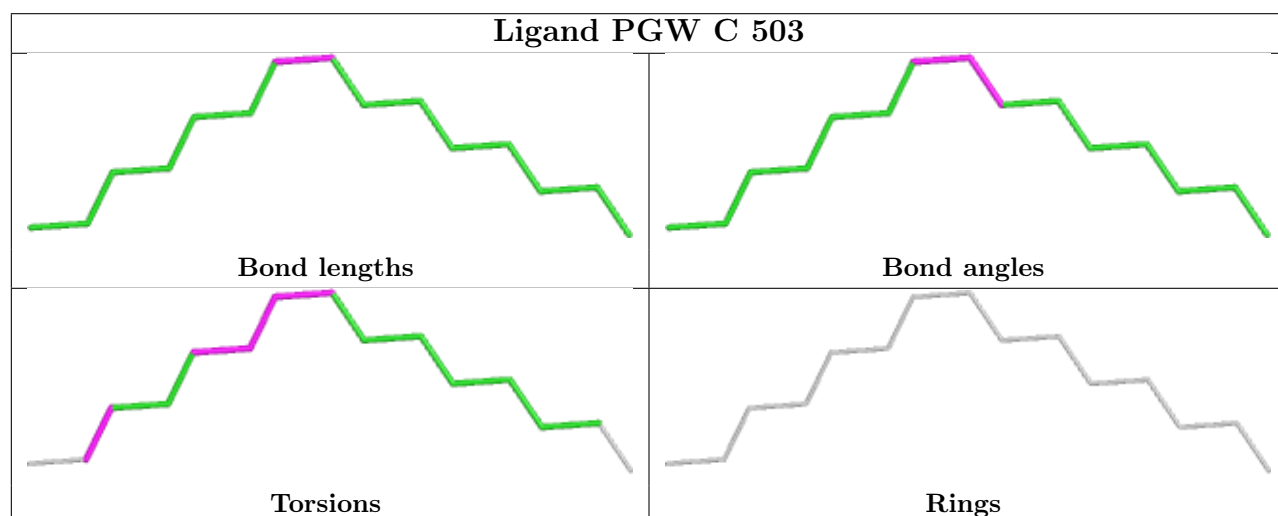
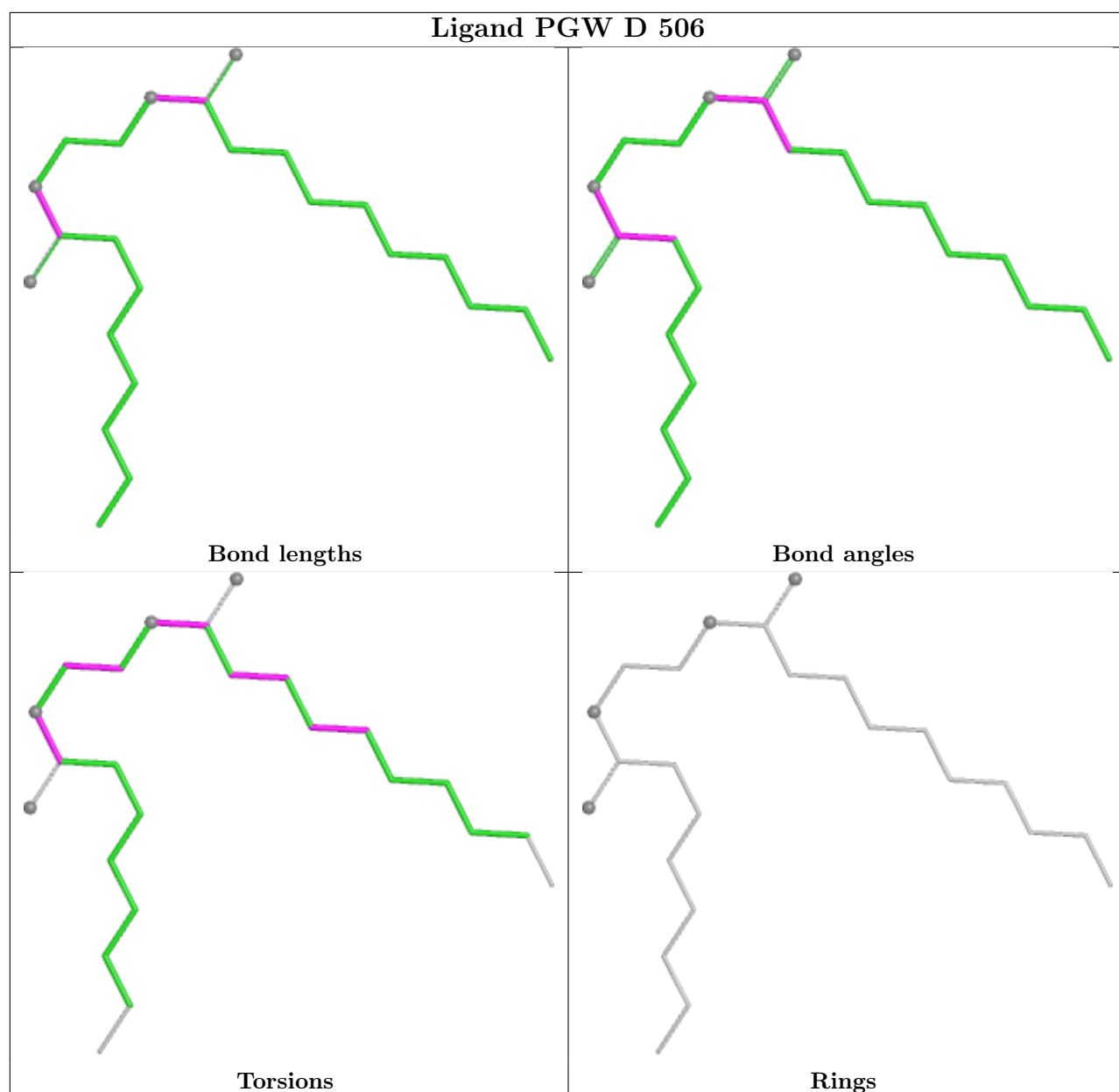
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

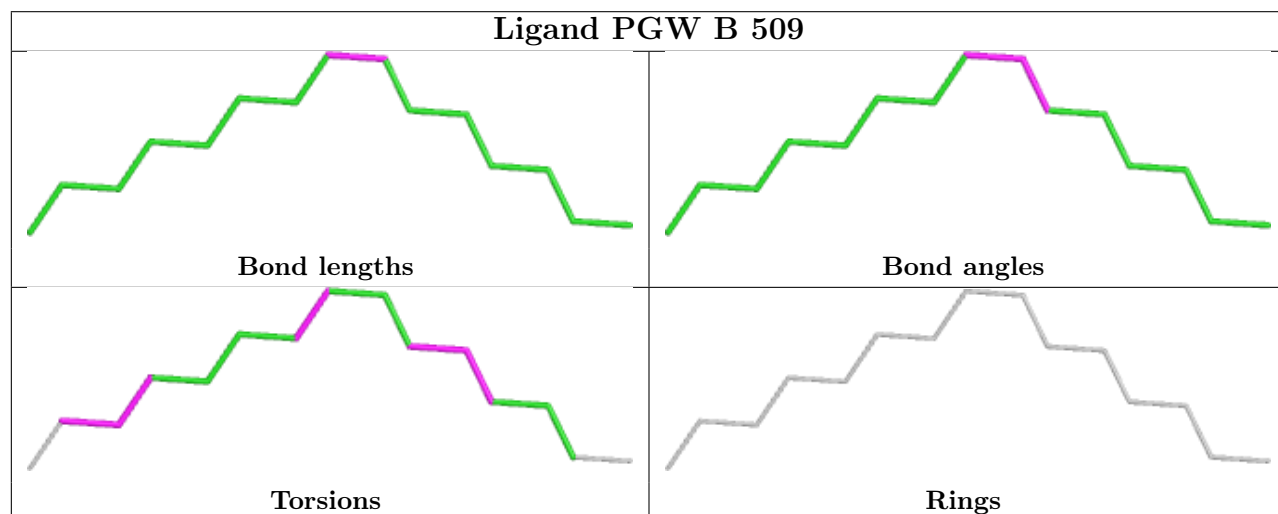
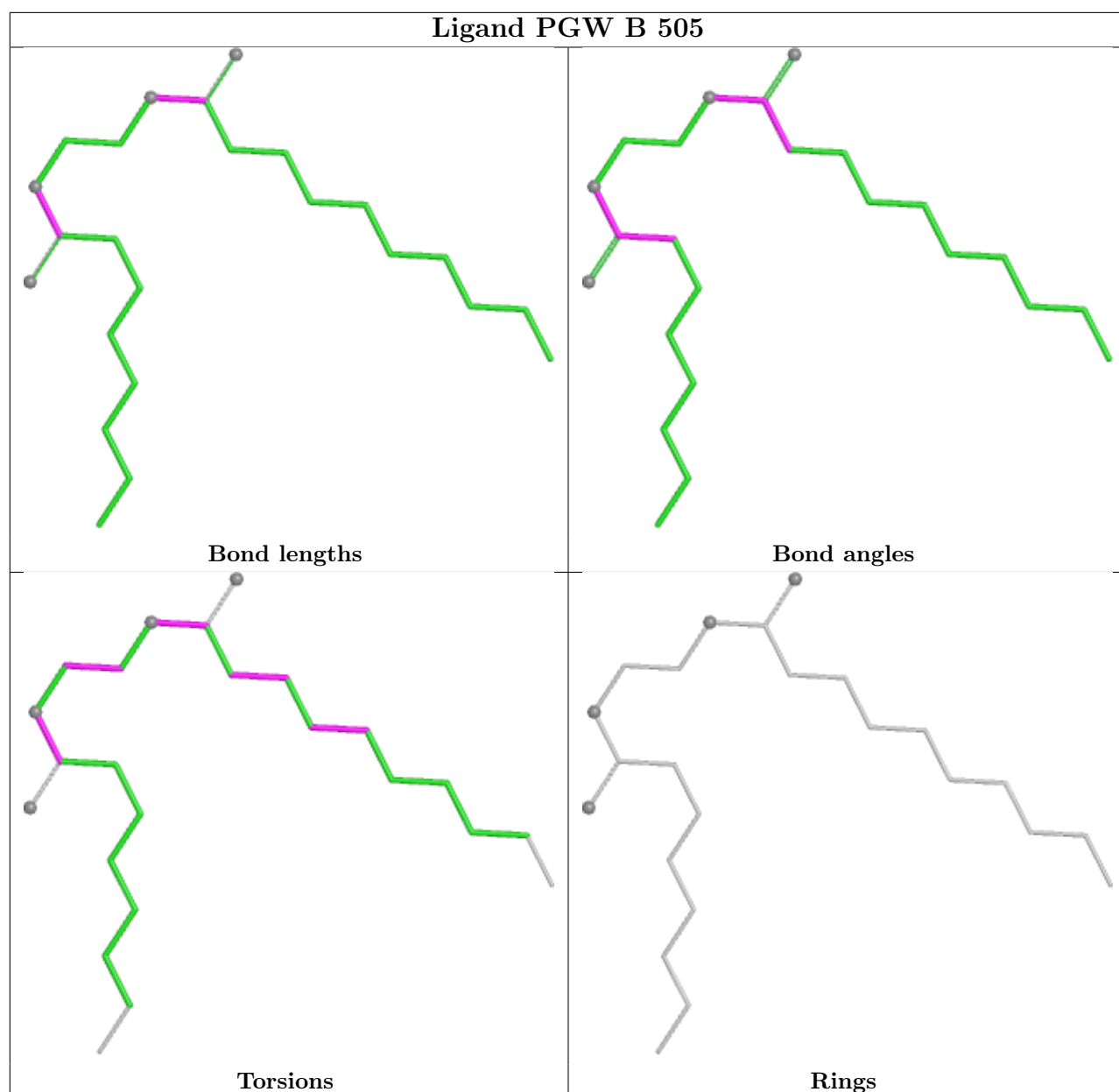


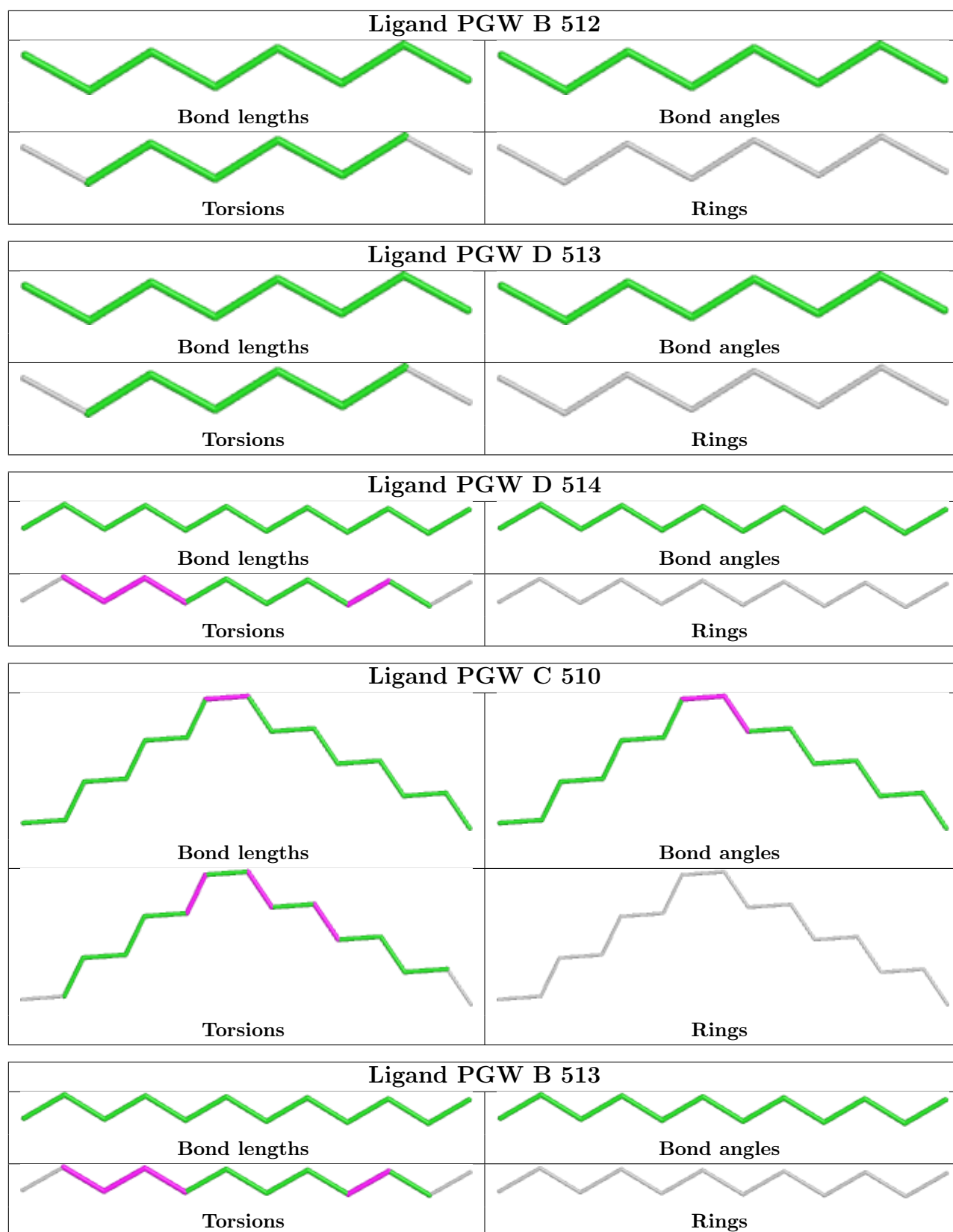


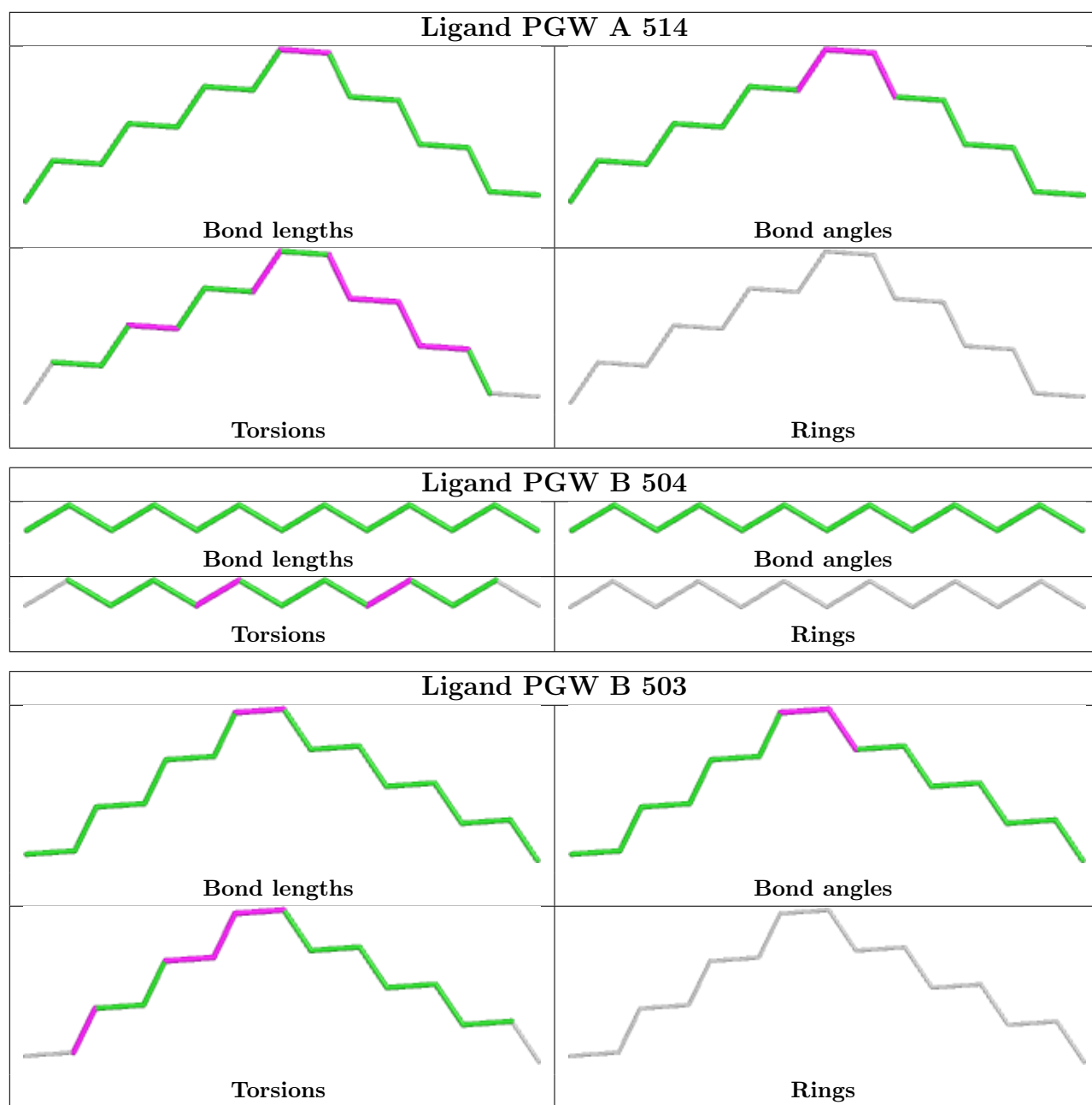


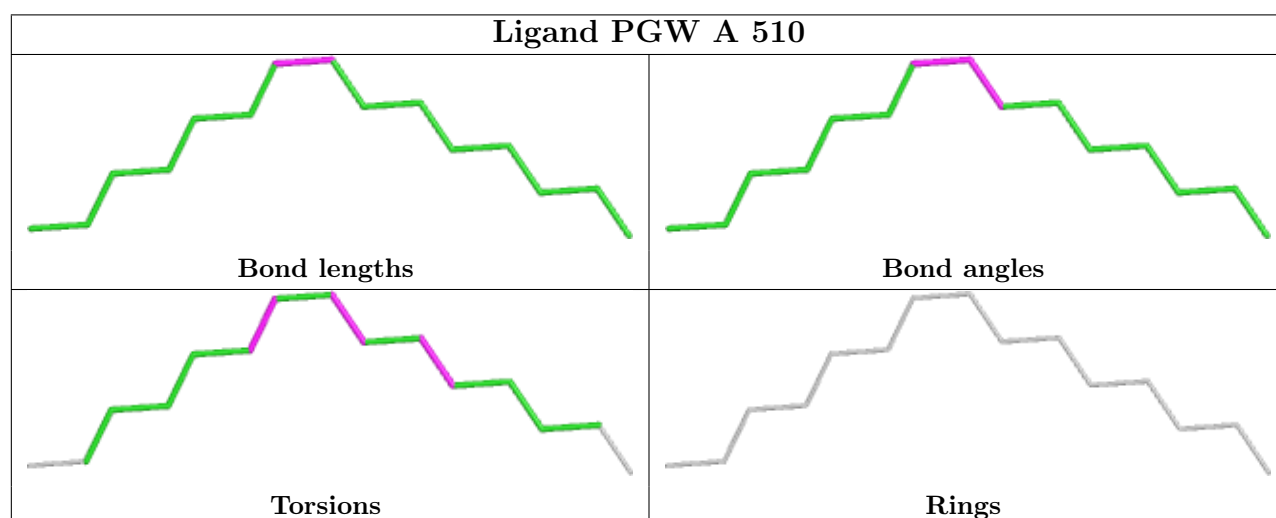
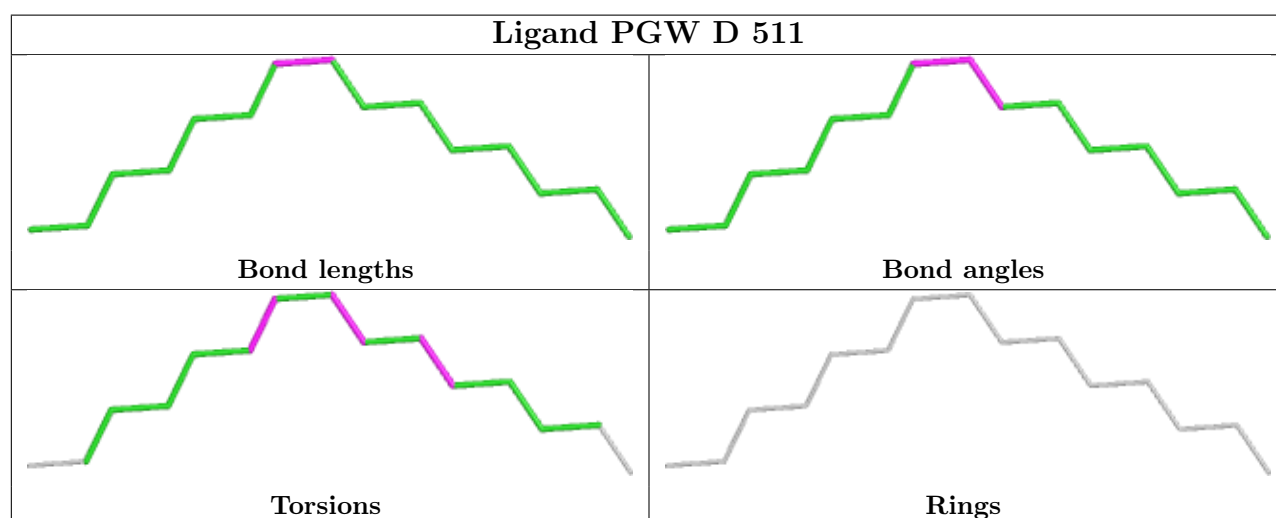
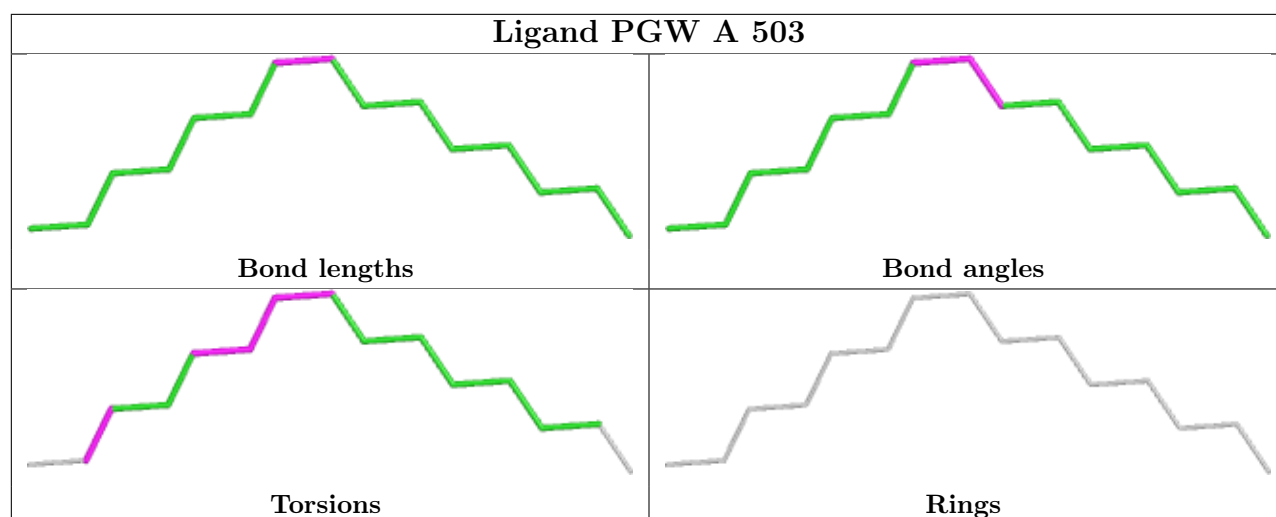


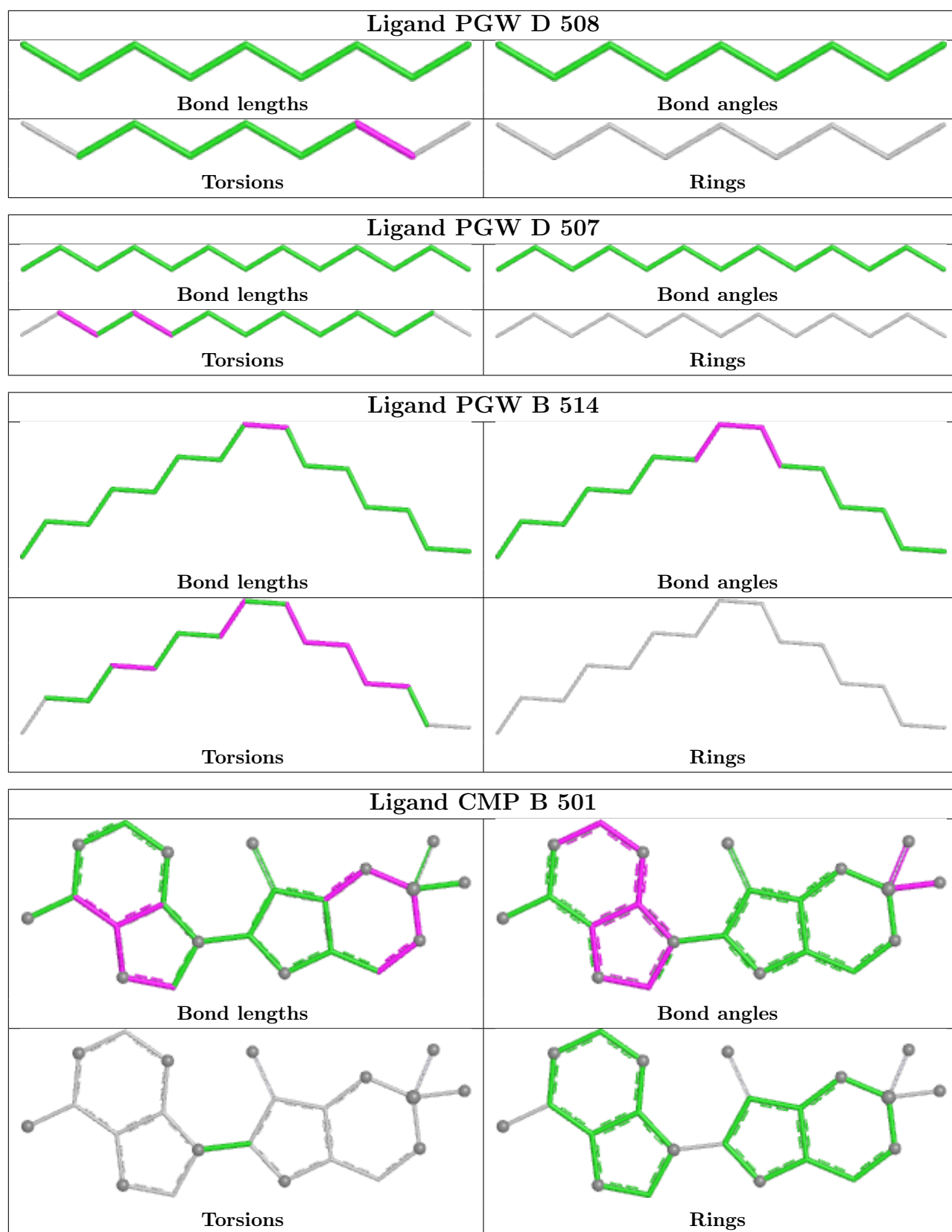


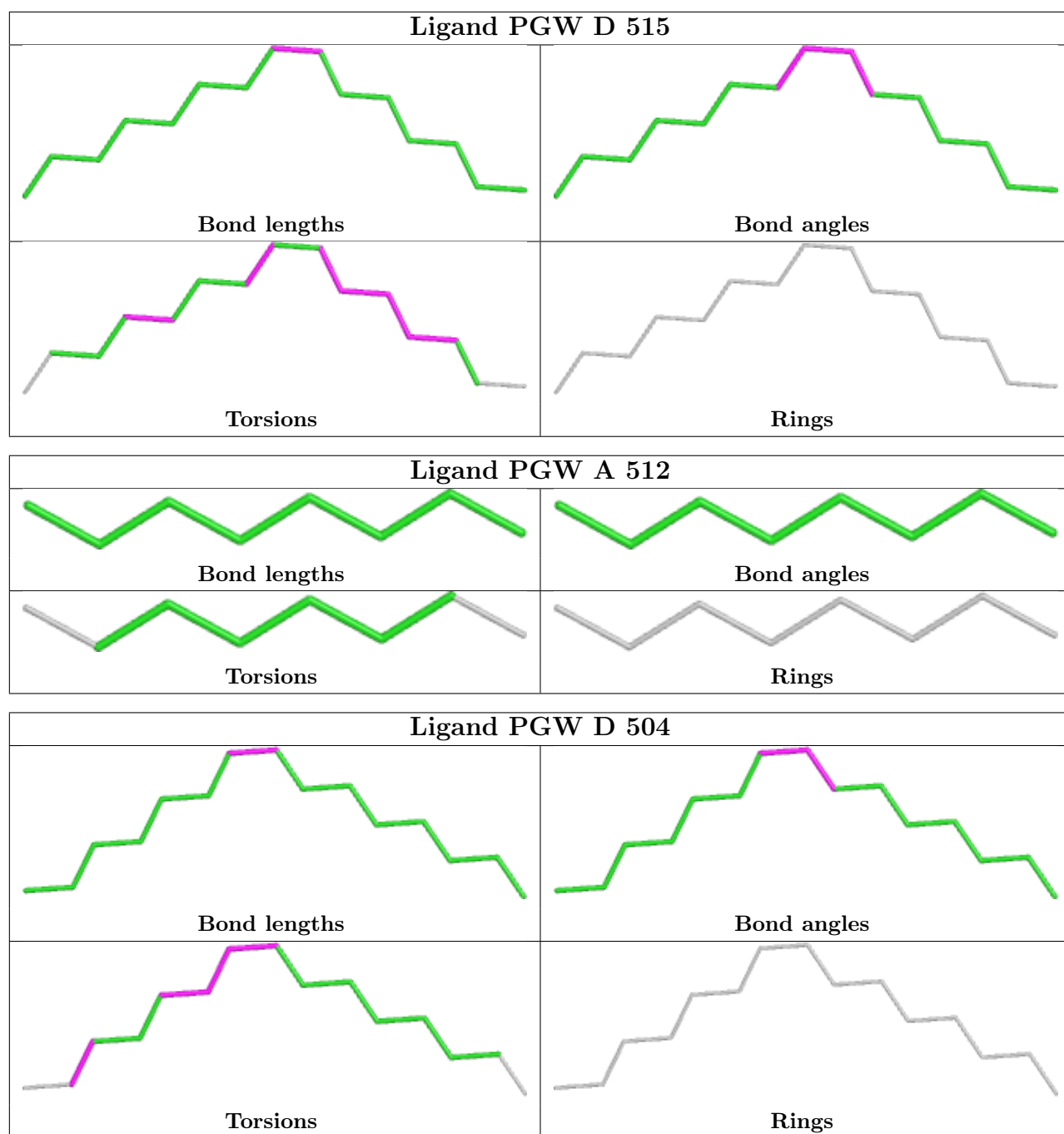


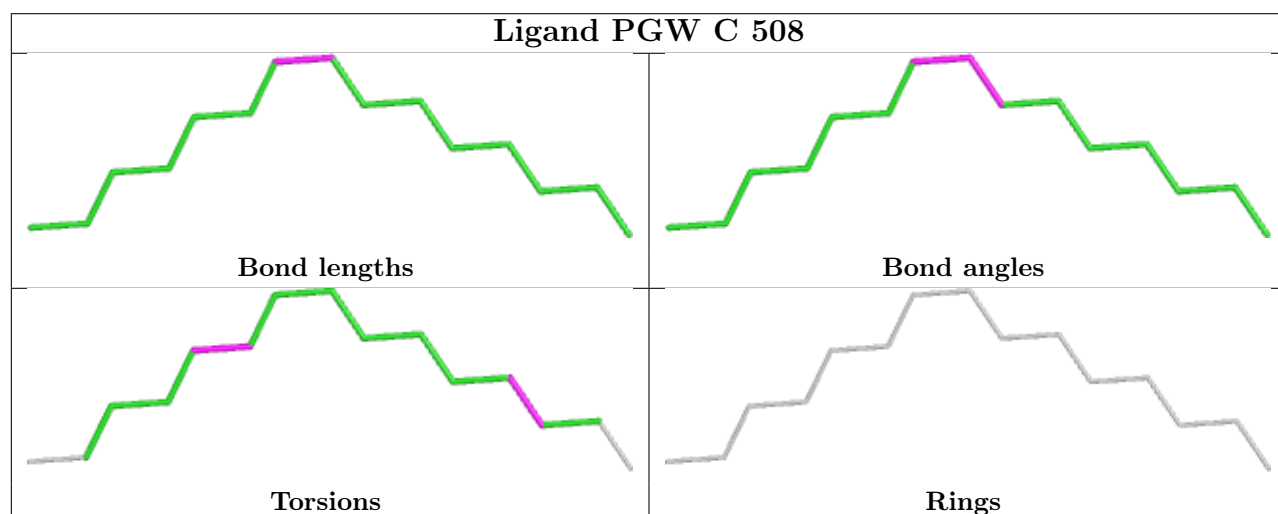
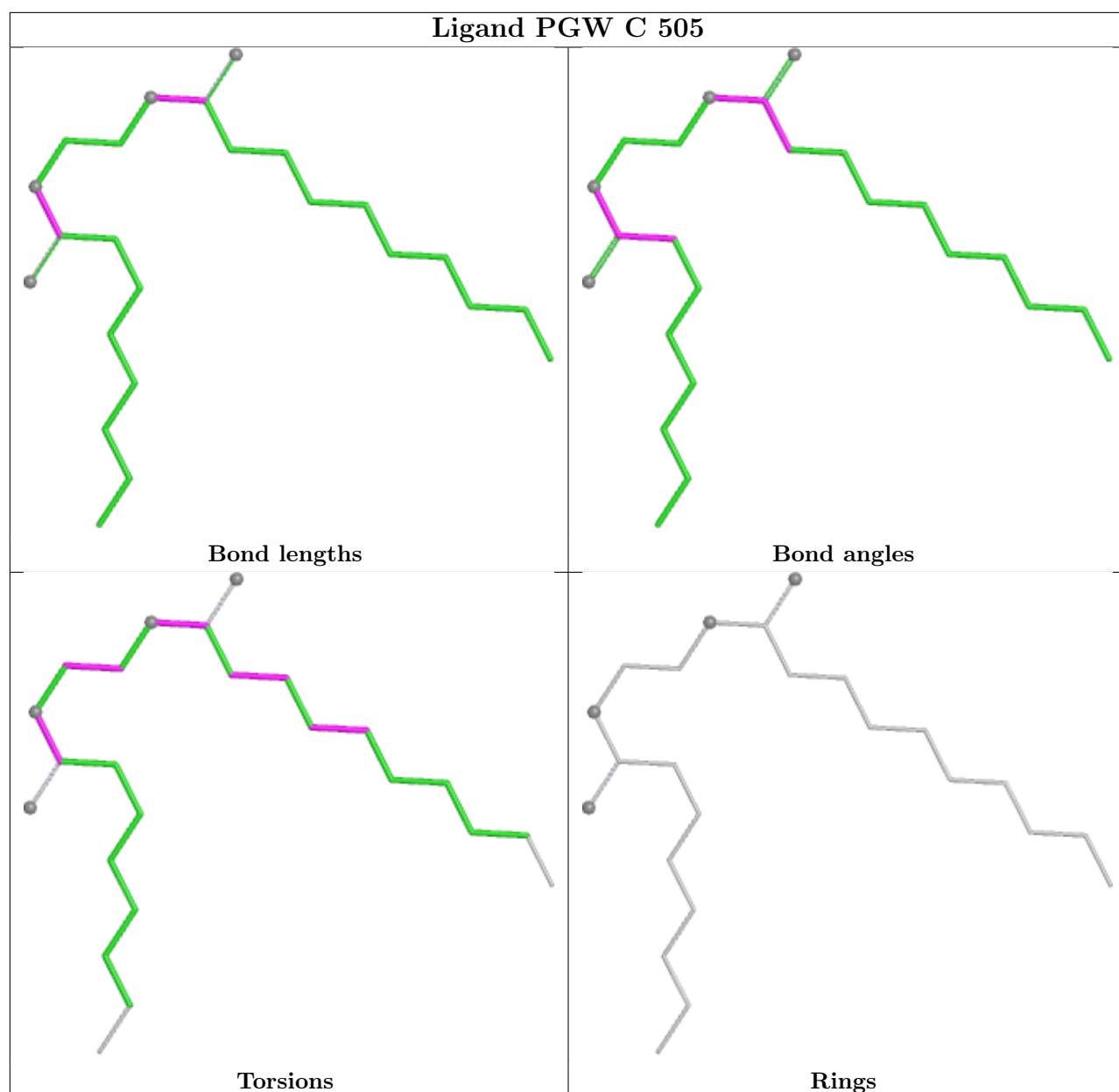


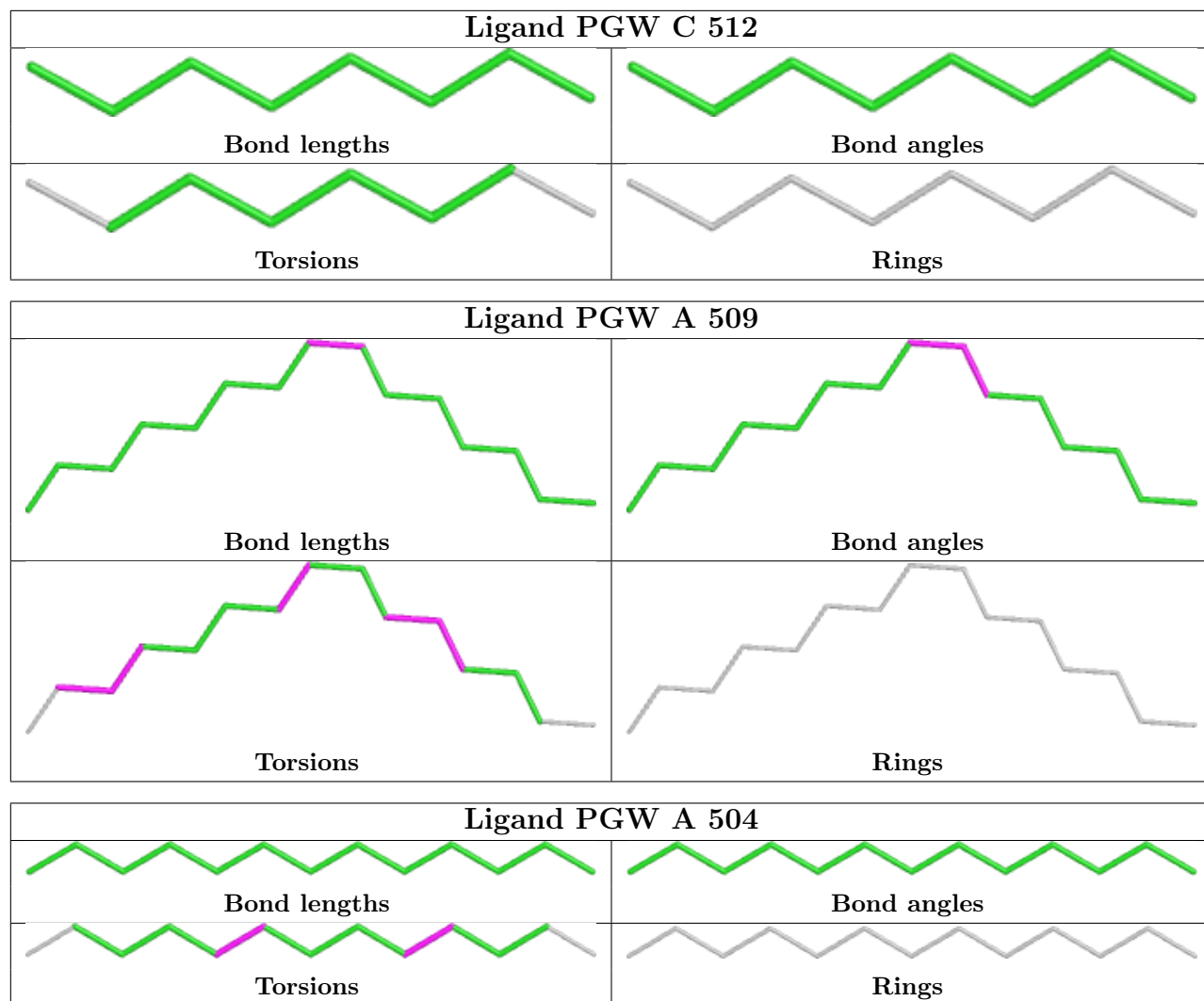


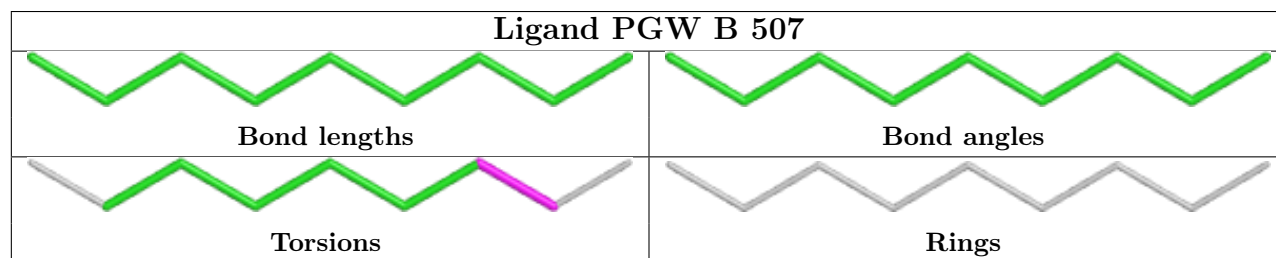
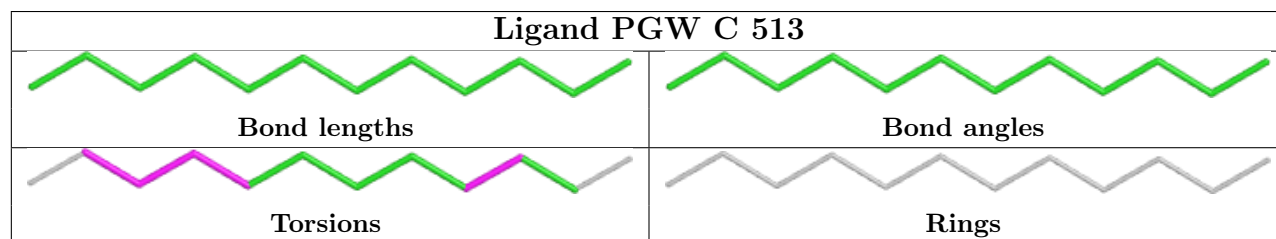
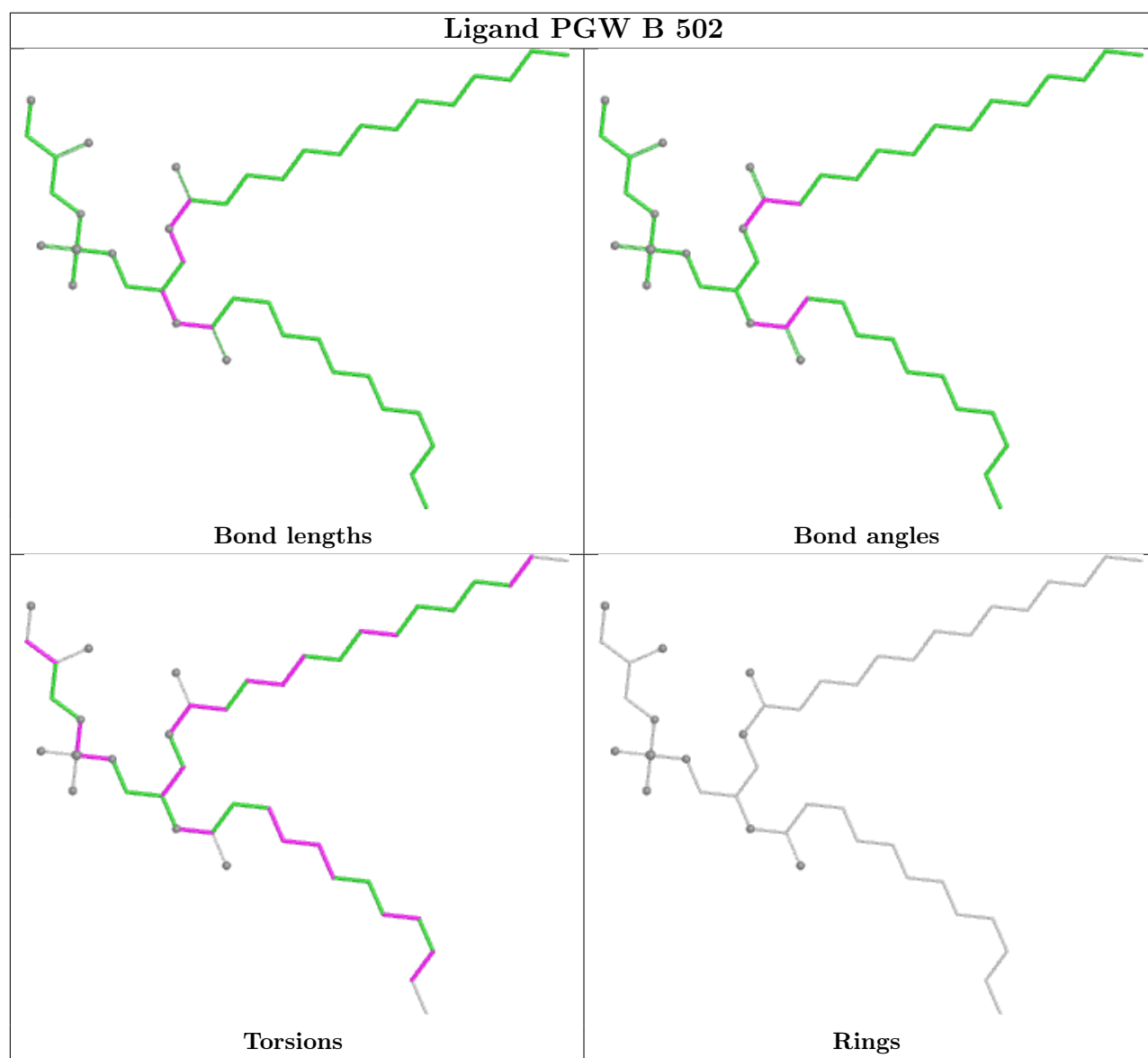


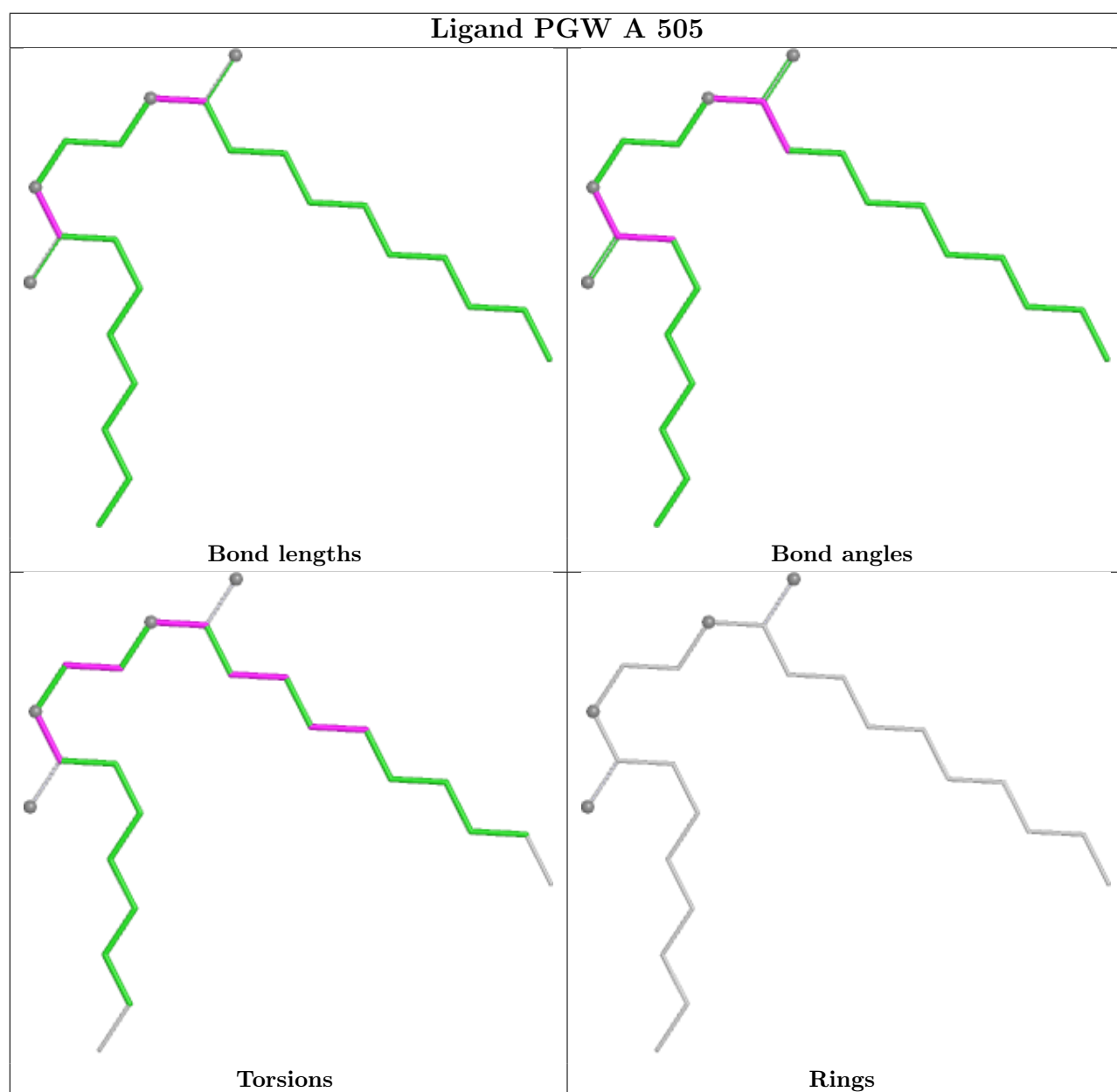
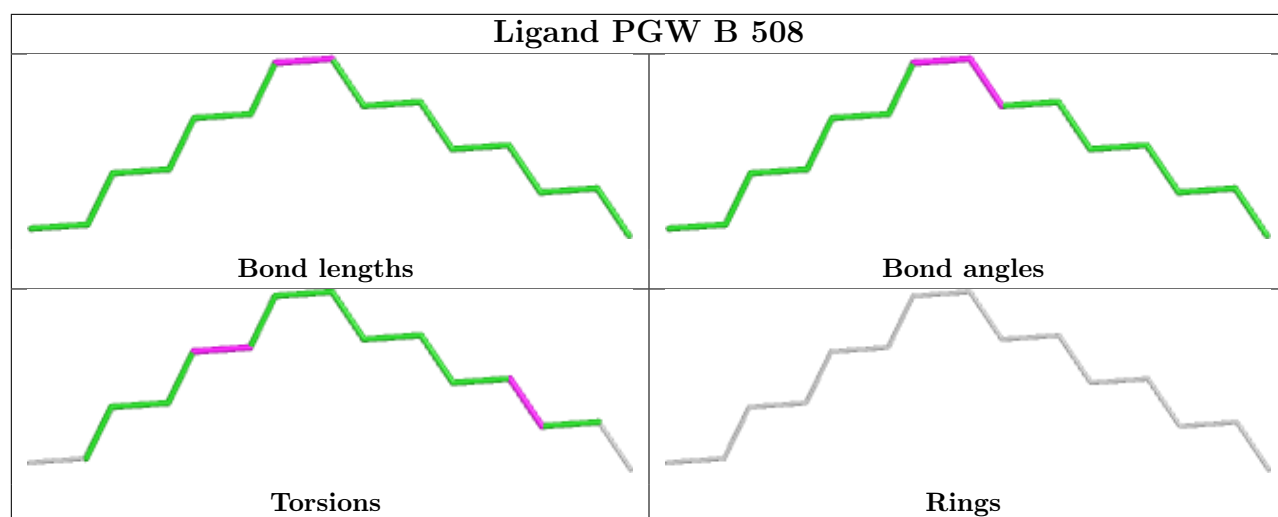


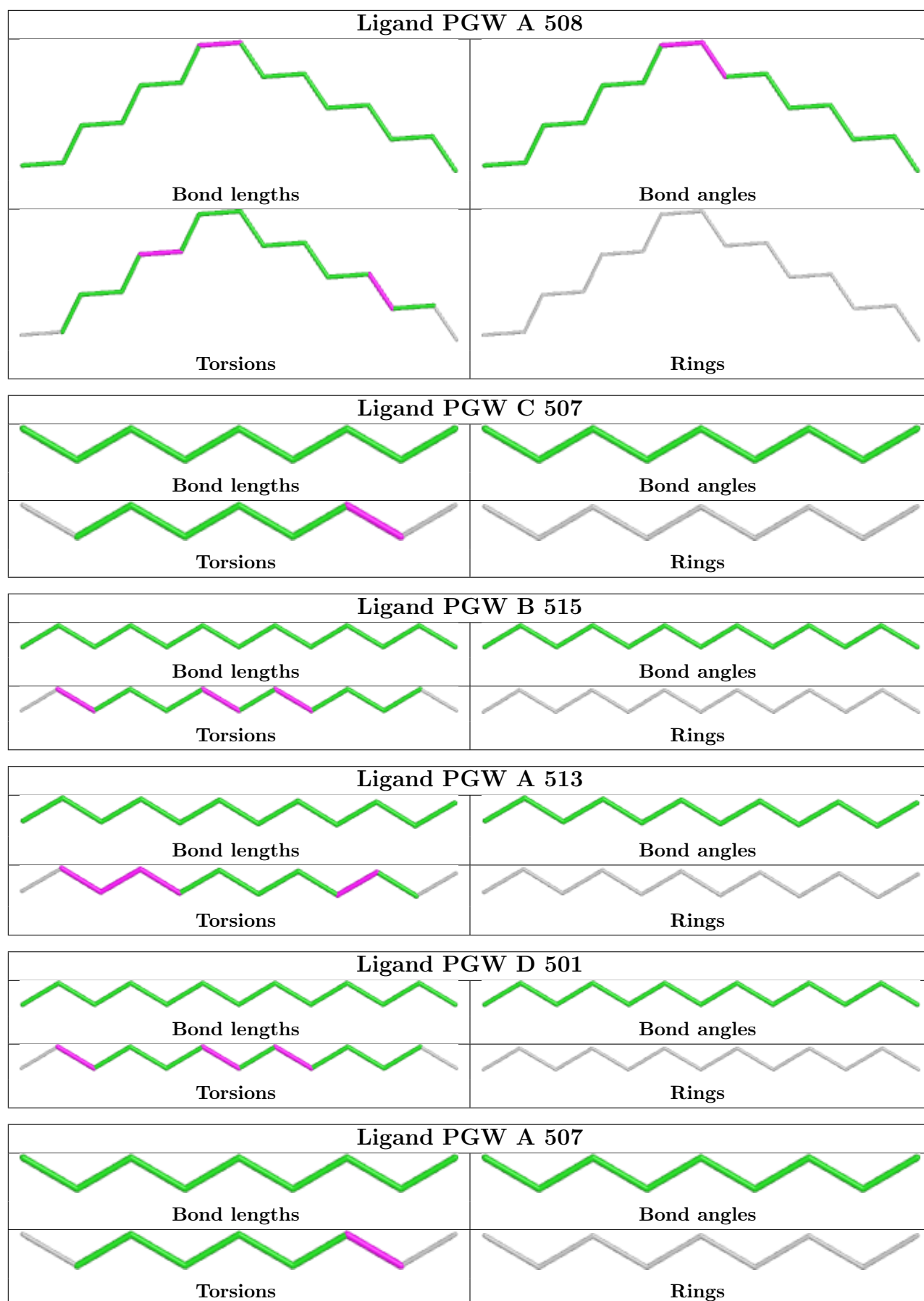


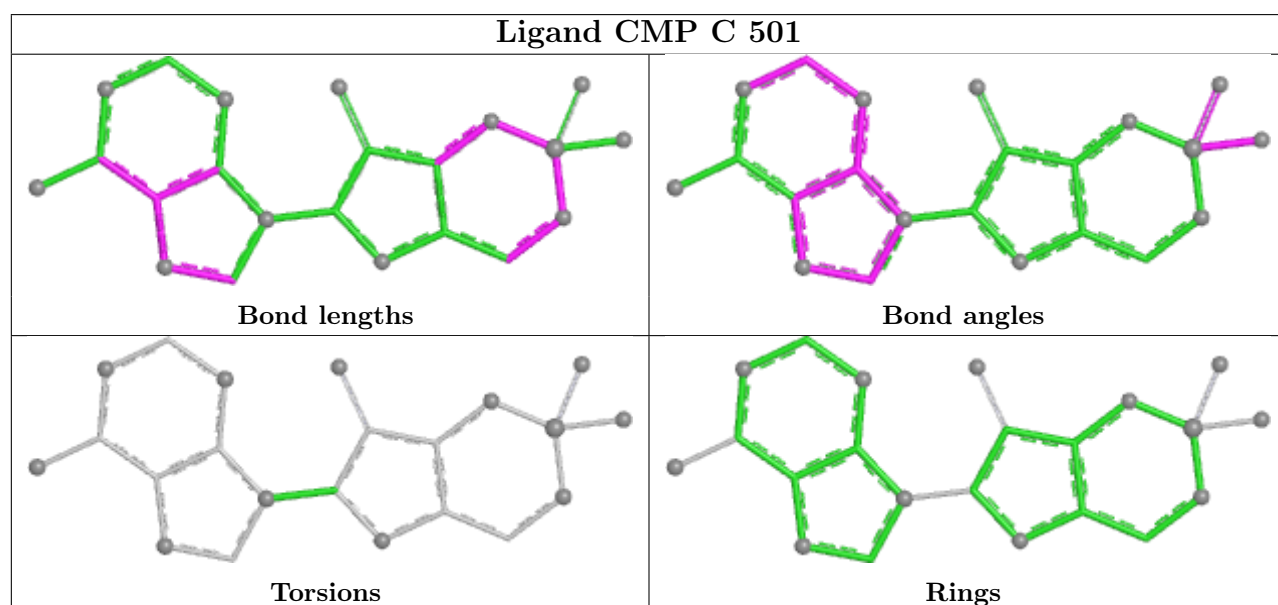
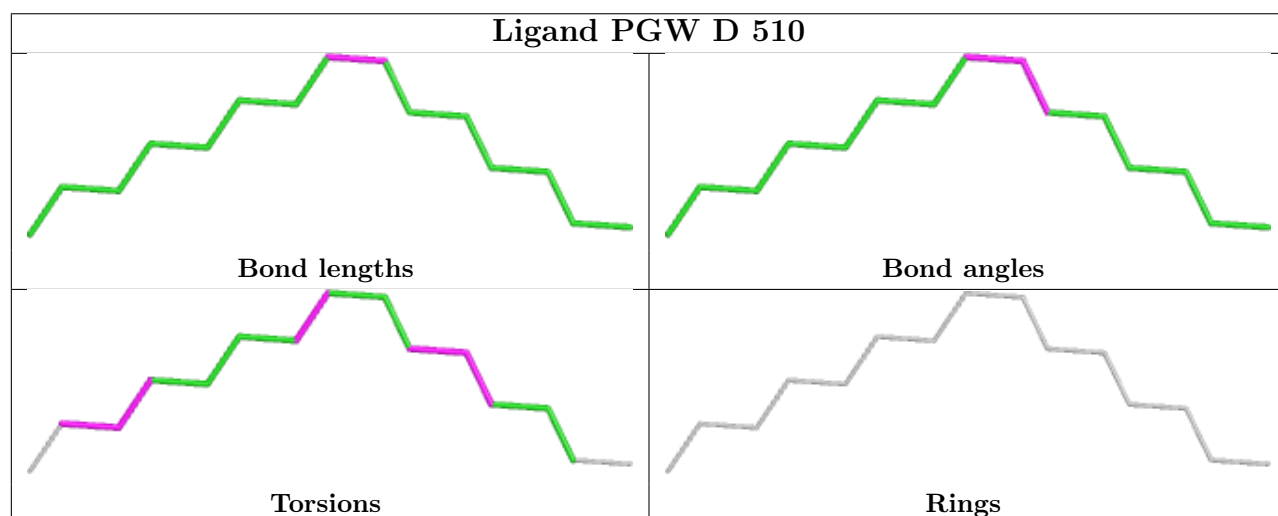
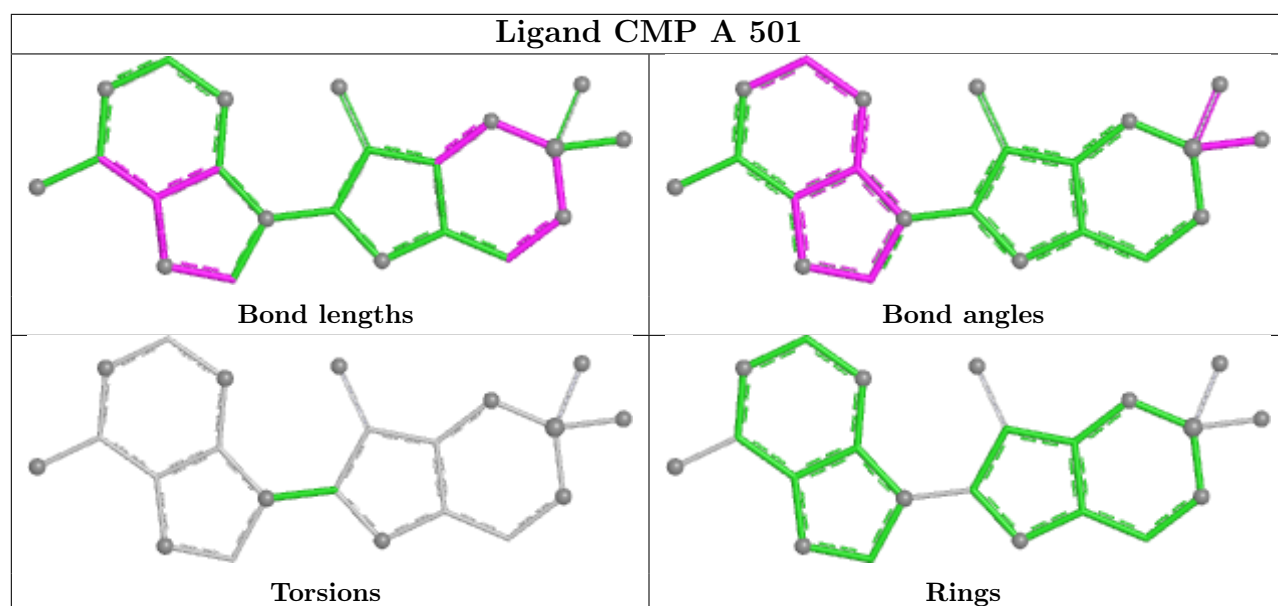


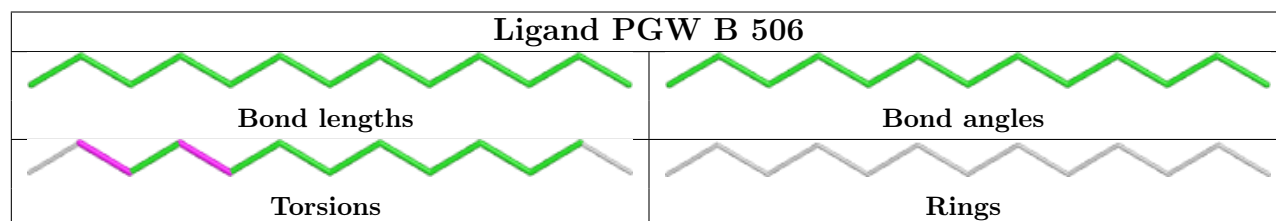
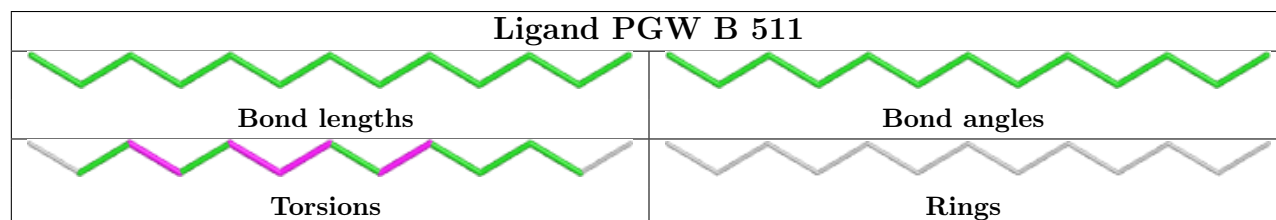
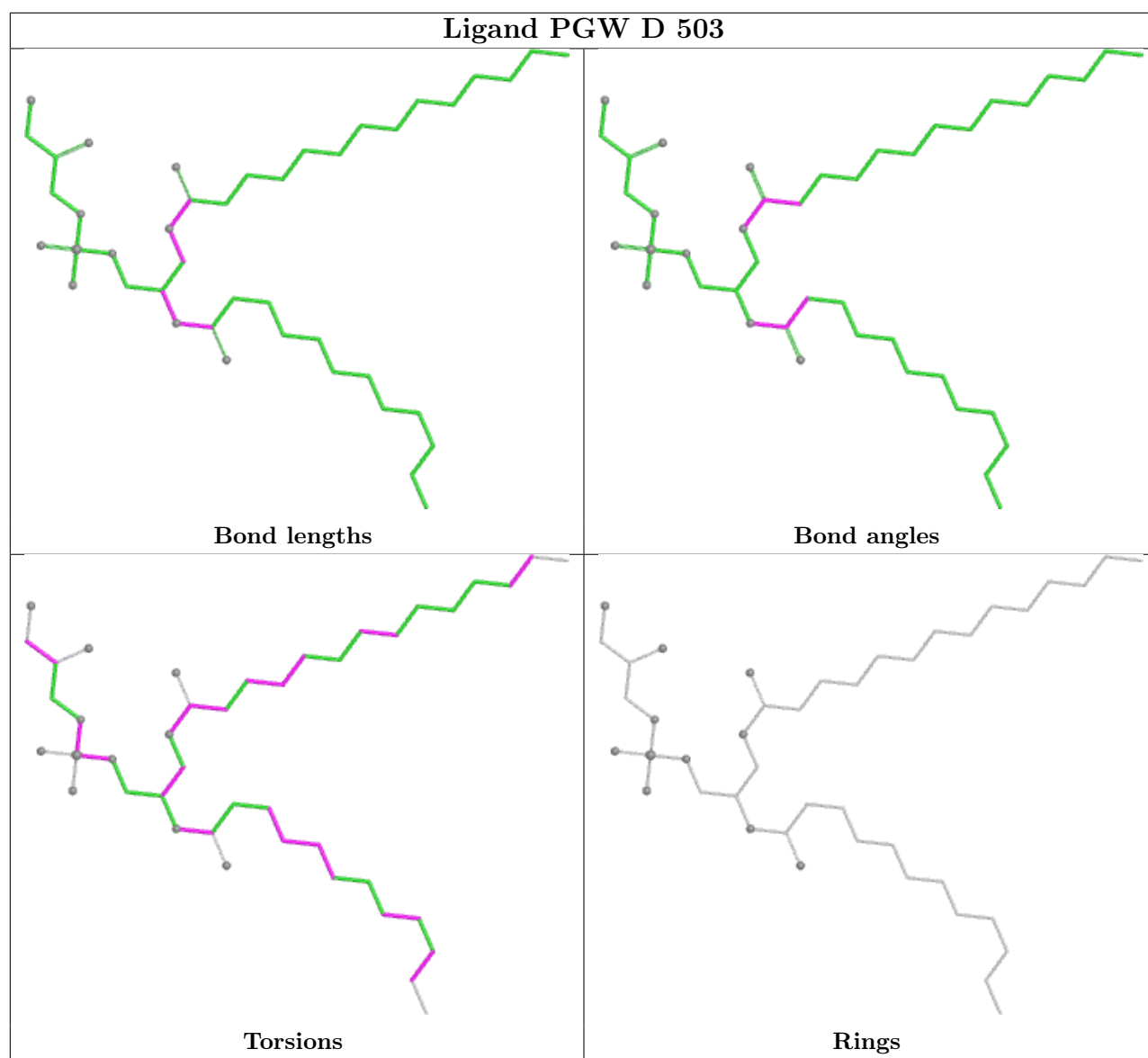


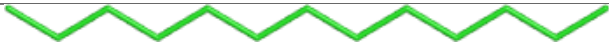
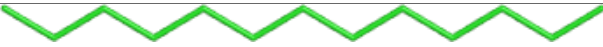
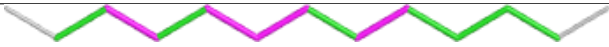








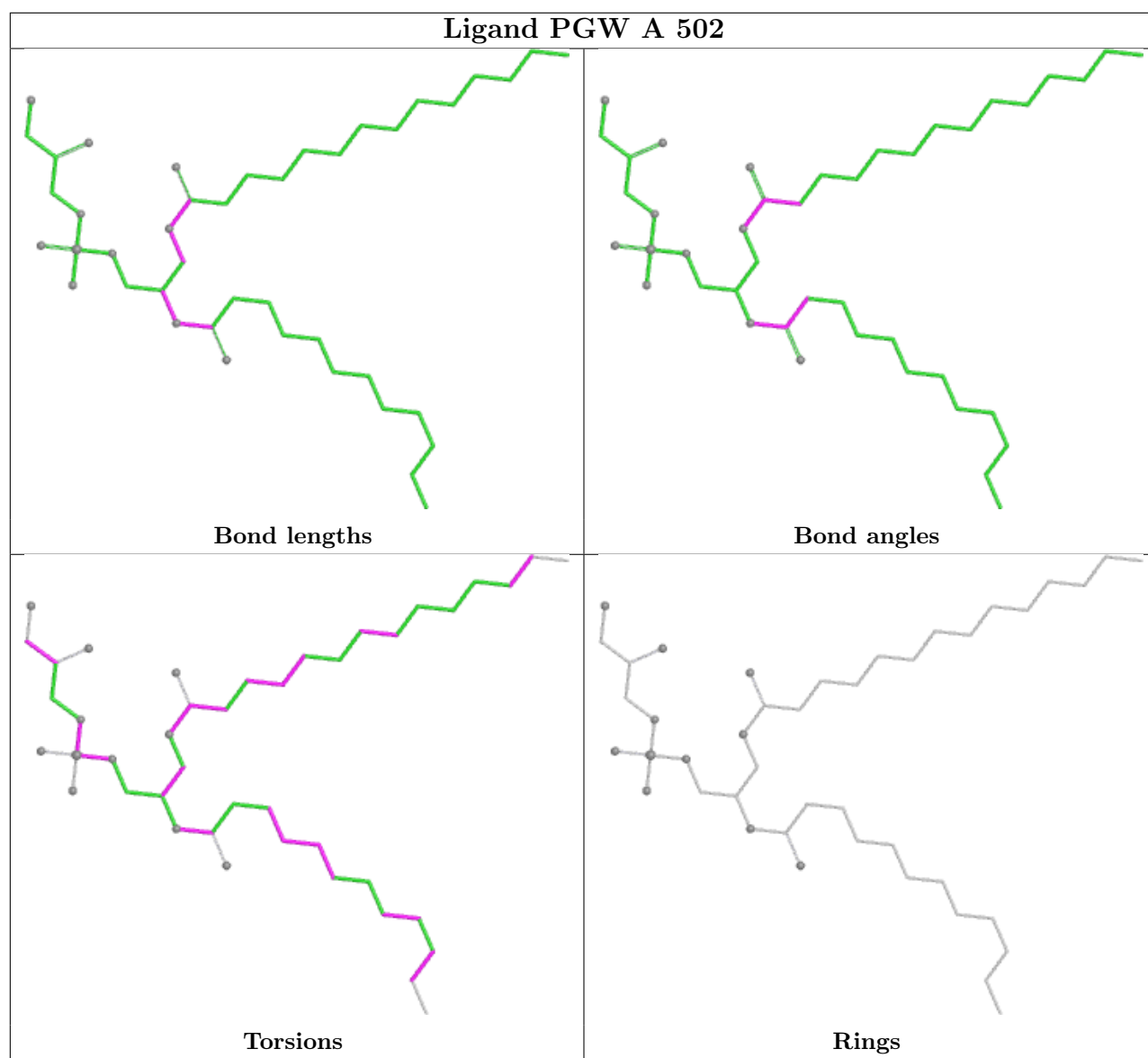


Ligand PGW D 512	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand PGW A 506	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand PGW D 505	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand PGW C 511	
 Bond lengths	 Bond angles
 Torsions	 Rings



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.

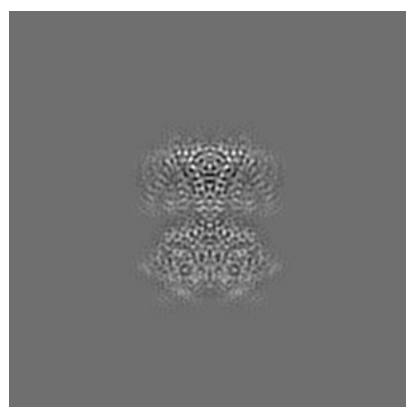
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24670. These allow visual inspection of the internal detail of the map and identification of artifacts.

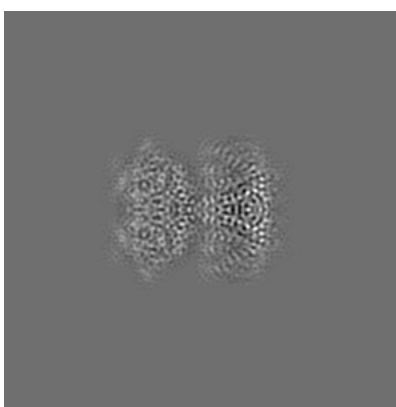
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

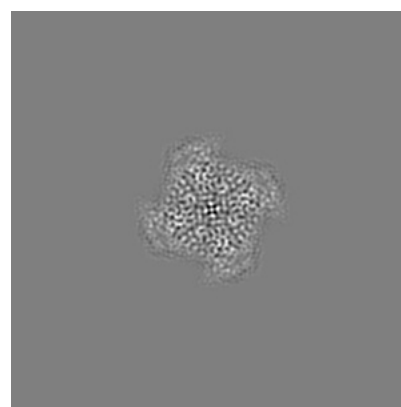
6.1.1 Primary map



X



Y

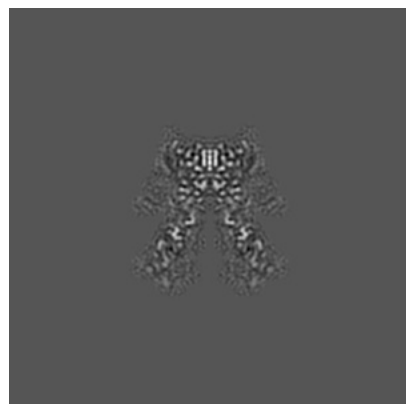


Z

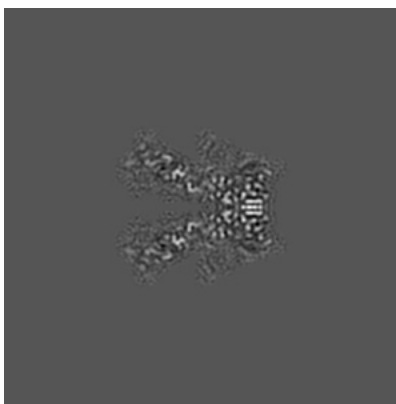
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

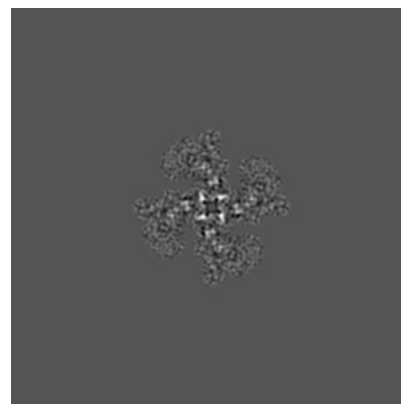
6.2.1 Primary map



X Index: 120



Y Index: 120

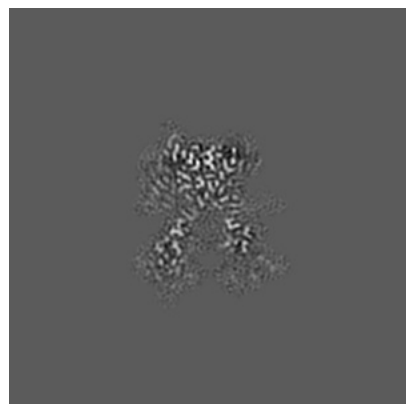


Z Index: 120

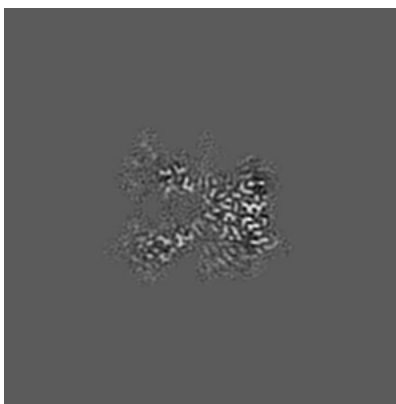
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

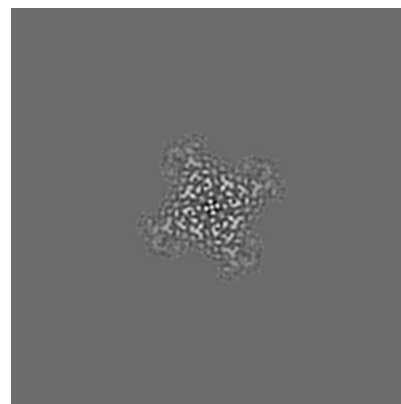
6.3.1 Primary map



X Index: 124



Y Index: 116

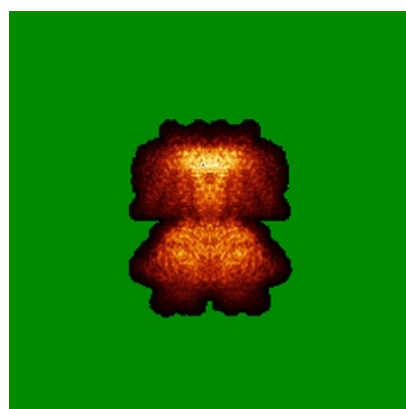


Z Index: 150

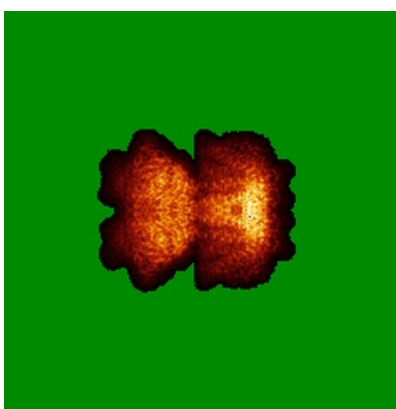
The images above show the largest variance slices of the map in three orthogonal directions.

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

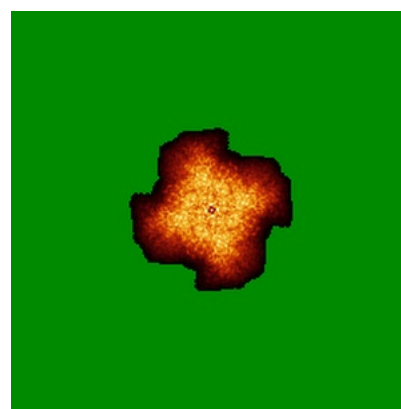
6.4.1 Primary map



X



Y



Z

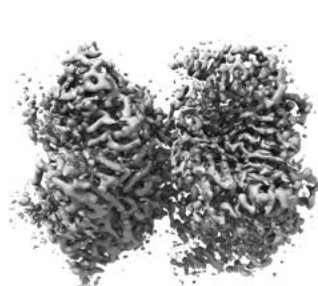
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

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 0.03. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

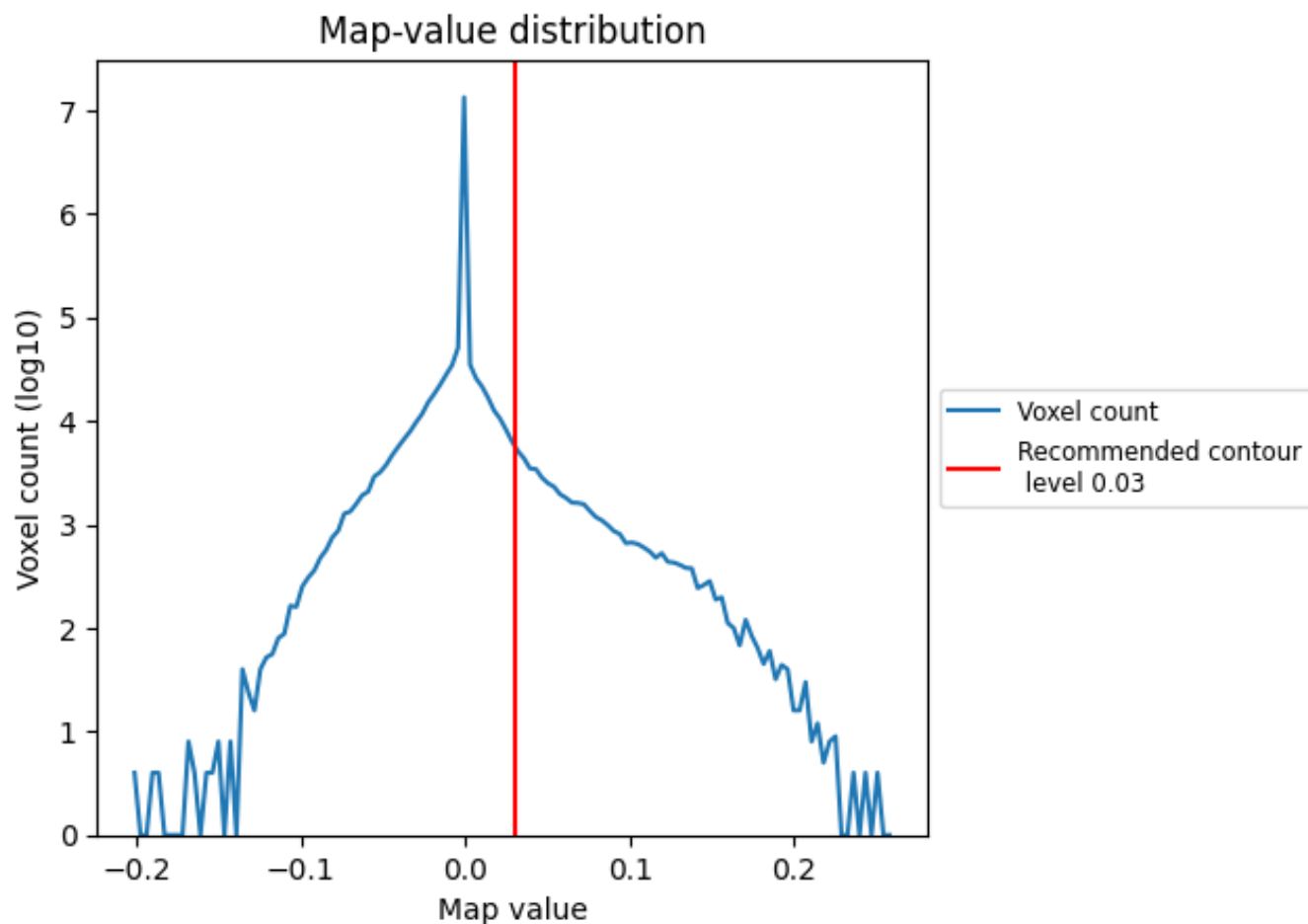
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

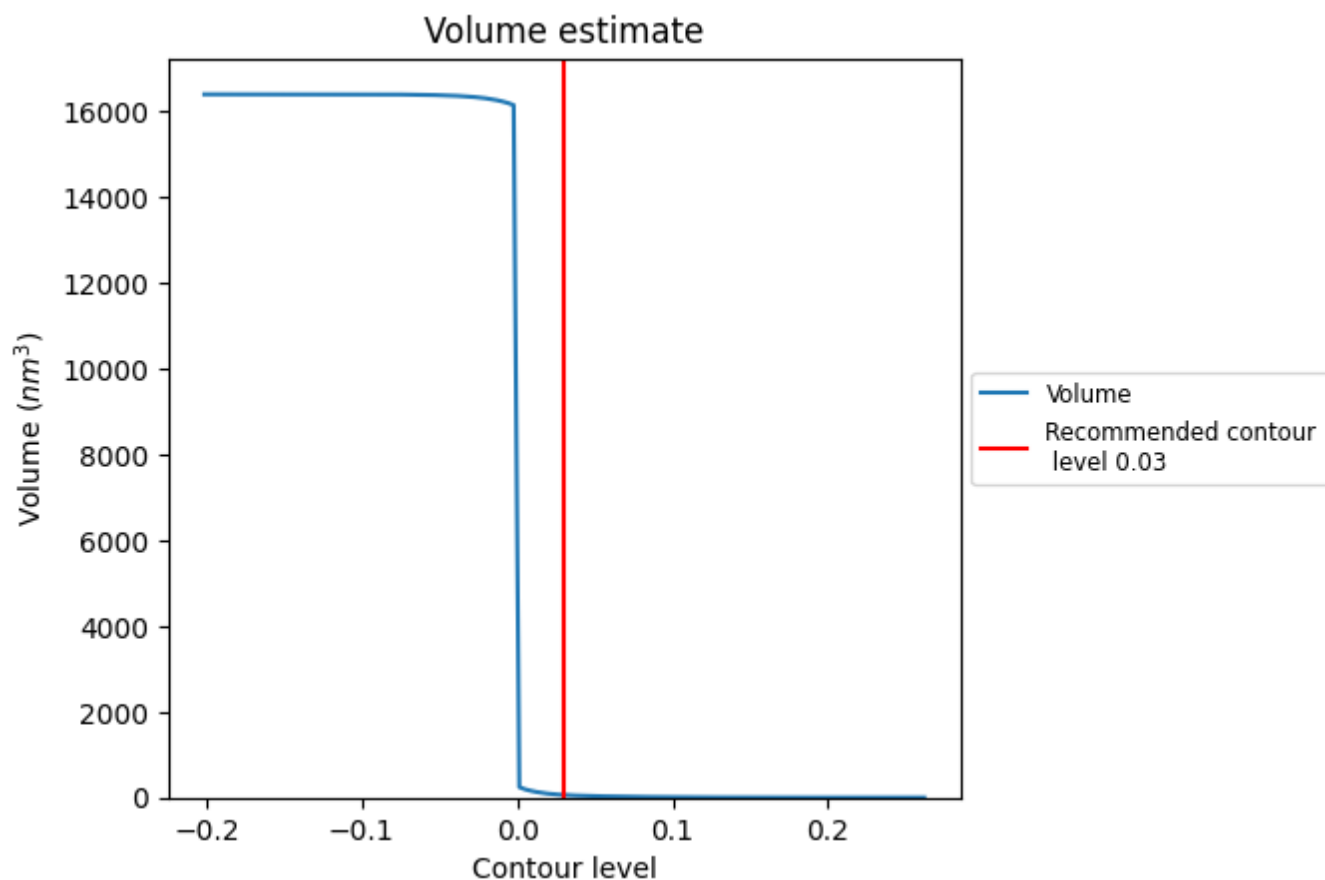
This section contains the results of statistical analysis of the map.

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.

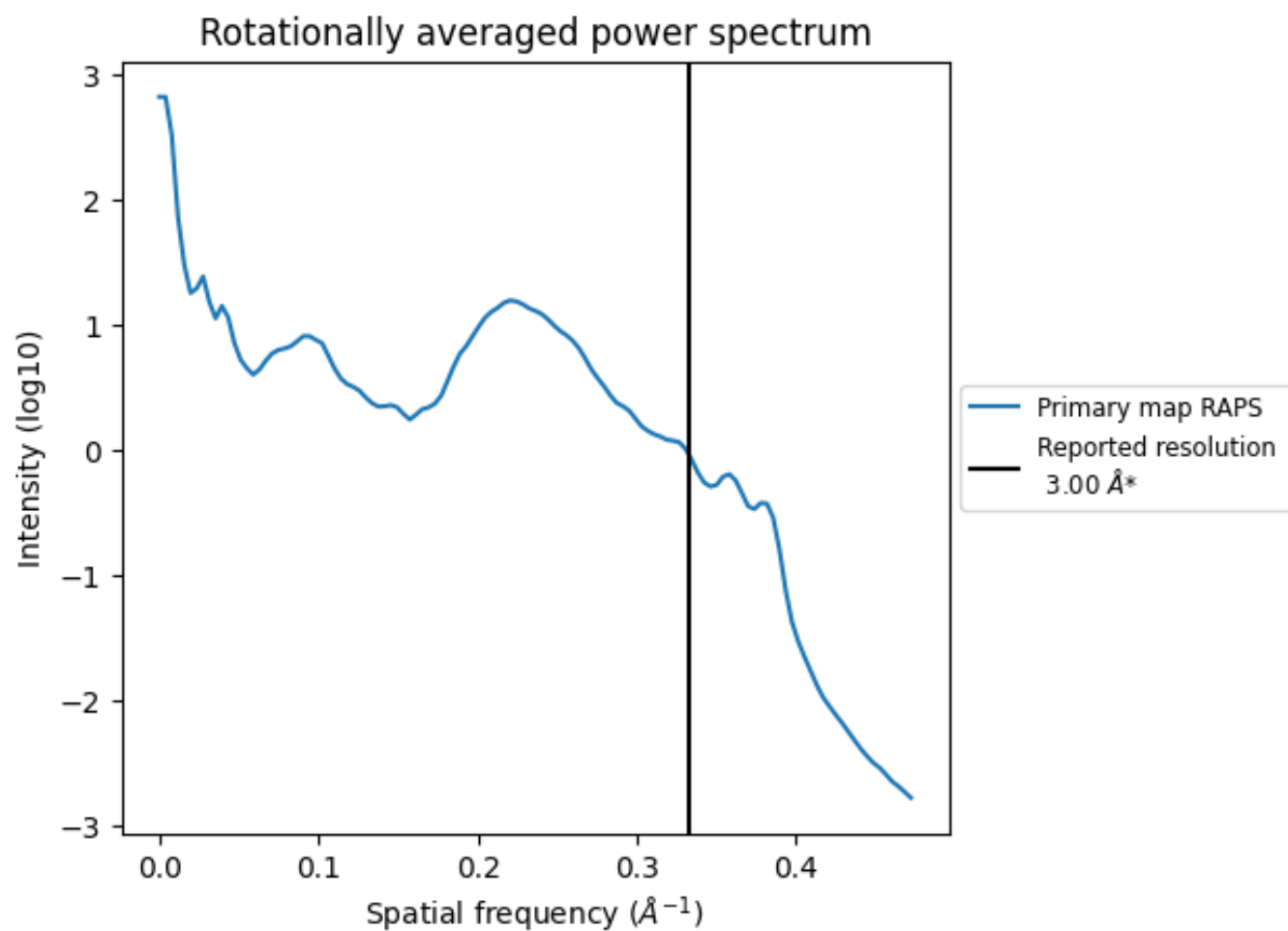
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 60 nm³; this corresponds to an approximate mass of 54 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ

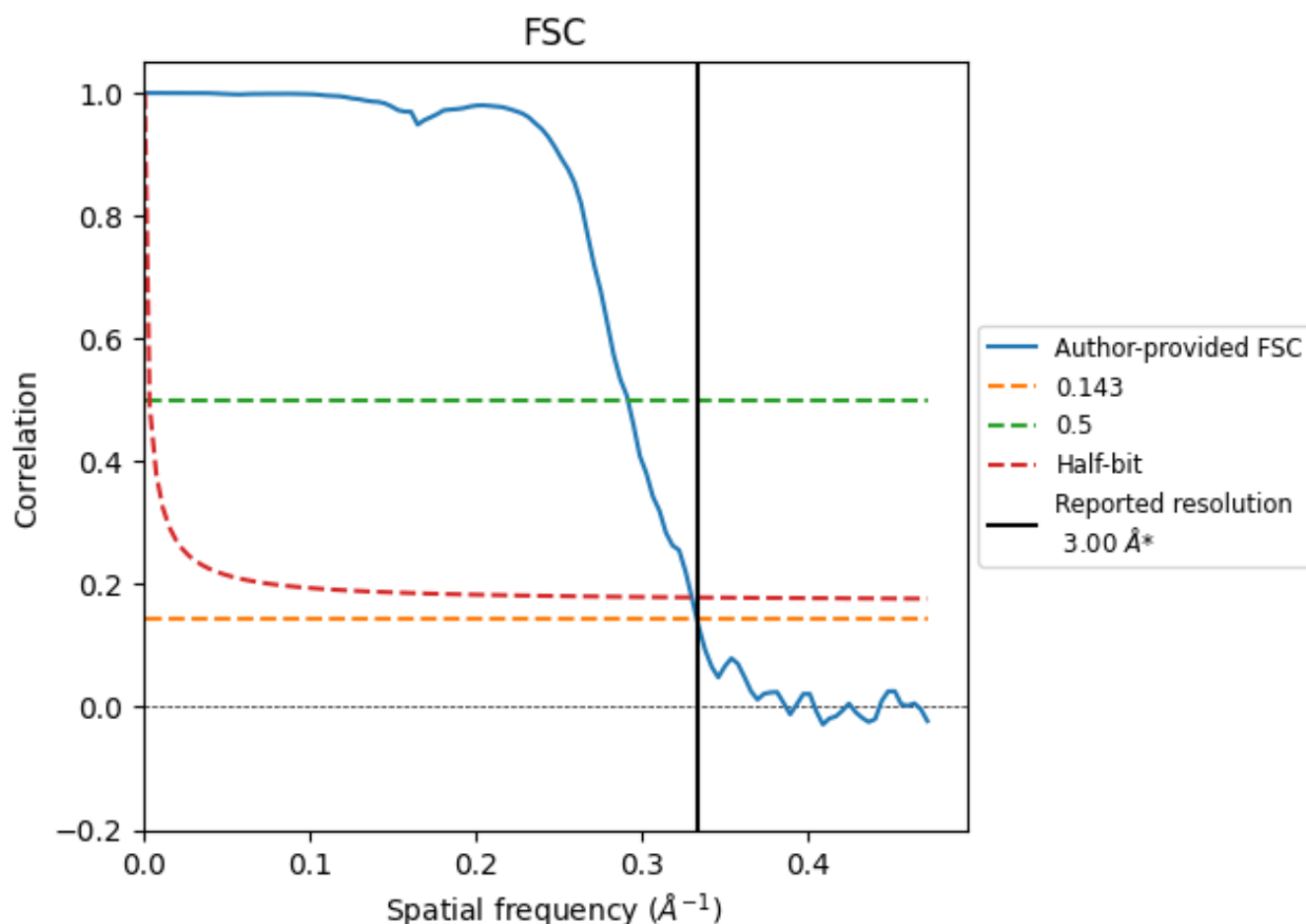


*Reported resolution corresponds to spatial frequency of 0.333 Å⁻¹

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 0.333 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	3.00	3.43	3.03
Unmasked-calculated*	-	-	-

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

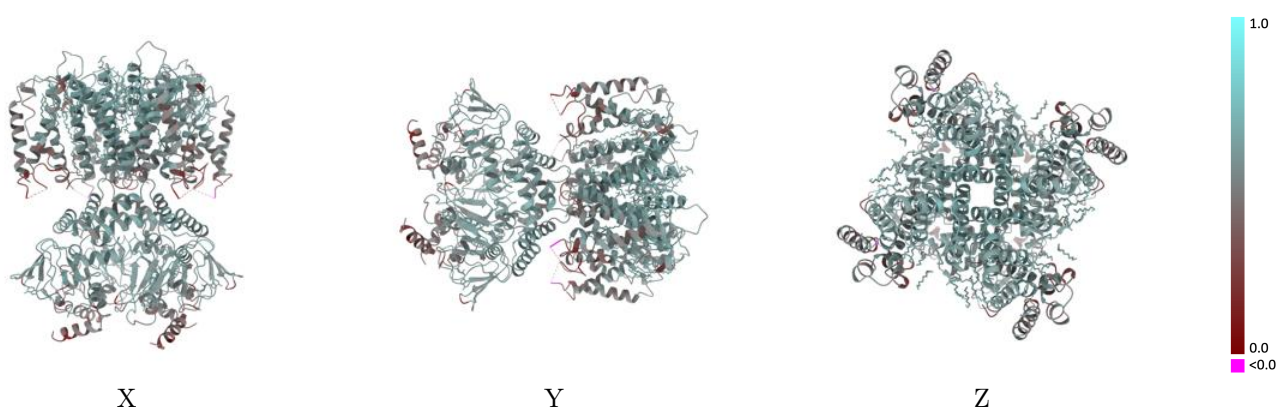
9 Map-model fit [i](#)

This section contains information regarding the fit between EMD map EMD-24670 and PDB model 7RSH. Per-residue inclusion information can be found in section 3 on page 8.

9.1 Map-model overlay [i](#)

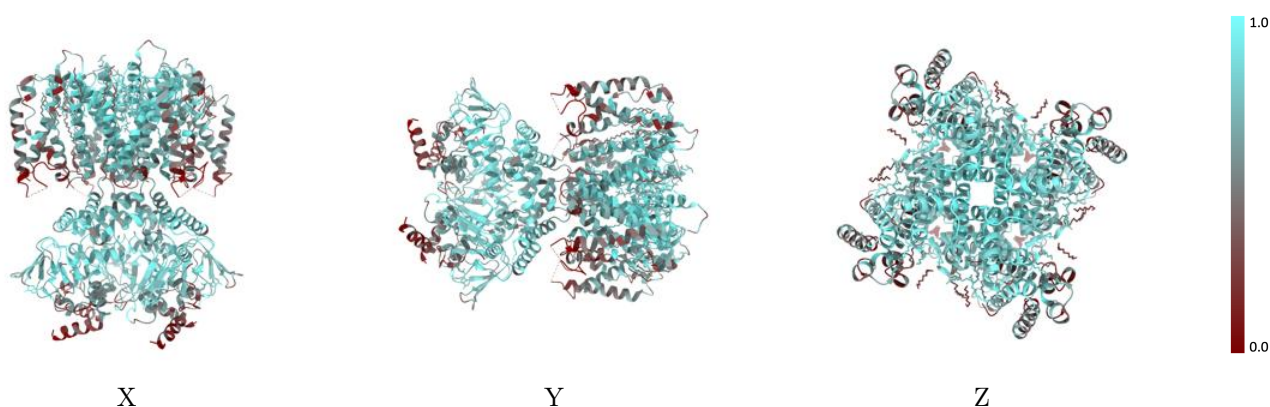
This section was not generated.

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



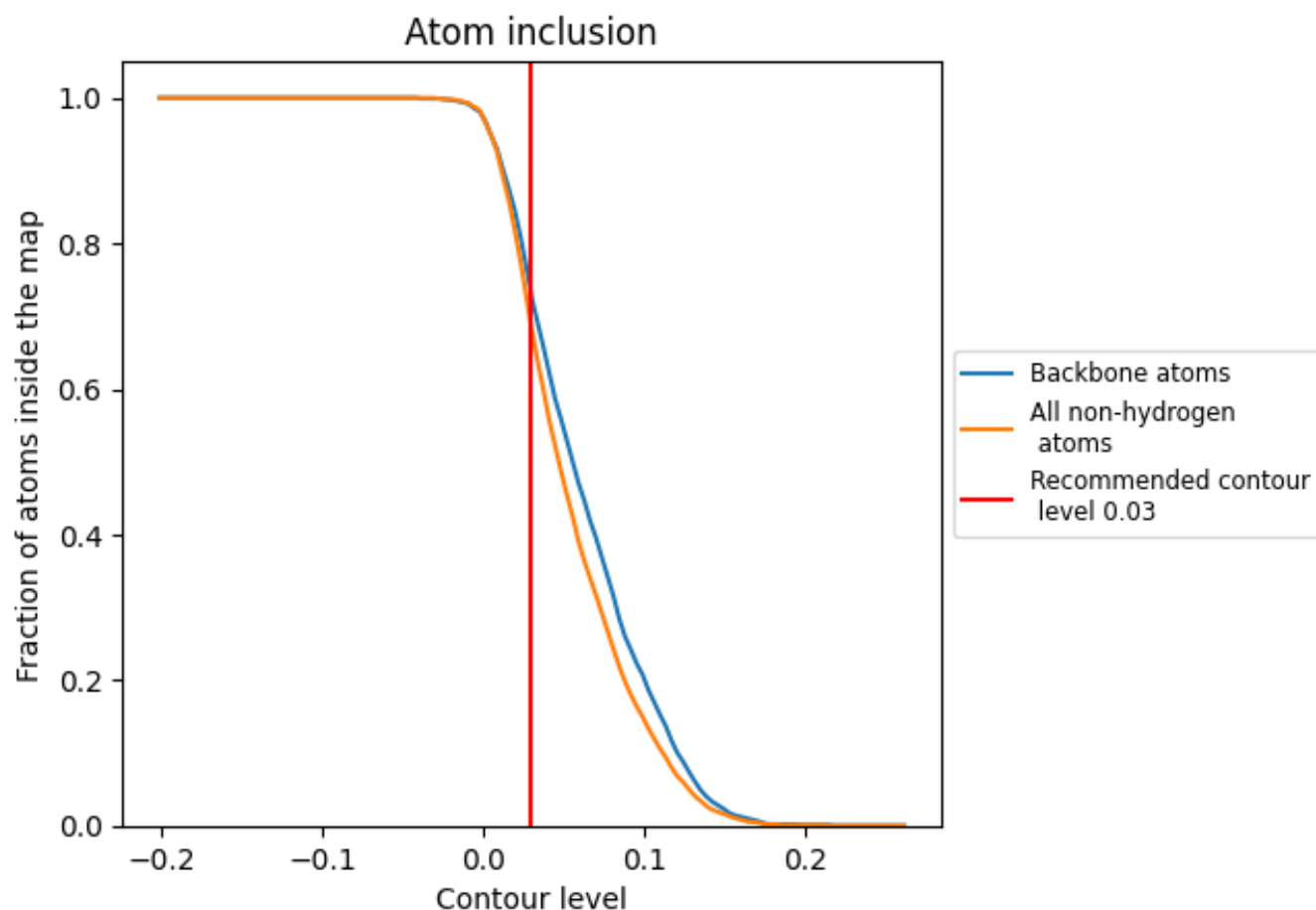
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.03).

9.4 Atom inclusion [i](#)



At the recommended contour level, 73% of all backbone atoms, 69% 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 (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6860	0.5410
A	0.6860	0.5420
B	0.6860	0.5410
C	0.6870	0.5400
D	0.6840	0.5410

