



wwPDB EM Validation Summary Report ⓘ

May 25, 2025 – 01:27 PM EDT

PDB ID : 9CUI / pdb_00009cui
EMDB ID : EMD-45934
Title : Structure of human full-length ancestral TRPV6 channel in Calmodulin-bound state
Authors : Neuberger, A.; Nadezhdin, K.D.; Sobolevsky, A.I.
Deposited on : 2024-07-26
Resolution : 3.42 Å(reported)

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<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.dev118
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0rc1
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

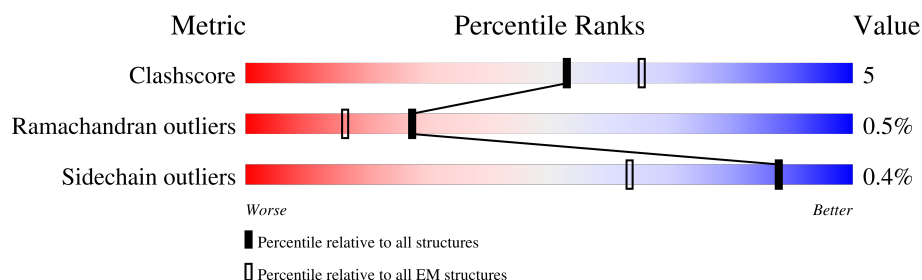
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.42 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	765	
1	B	765	
1	C	765	
1	D	765	
2	E	149	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 22713 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 Transient receptor potential cation channel subfamily V member 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	612	Total	C	N	O	S	0	0
			4916	3176	834	868	38		
1	B	620	Total	C	N	O	S	0	0
			4991	3220	854	879	38		
1	C	653	Total	C	N	O	S	0	0
			5286	3400	922	926	38		
1	D	617	Total	C	N	O	S	0	0
			4972	3213	847	874	38		

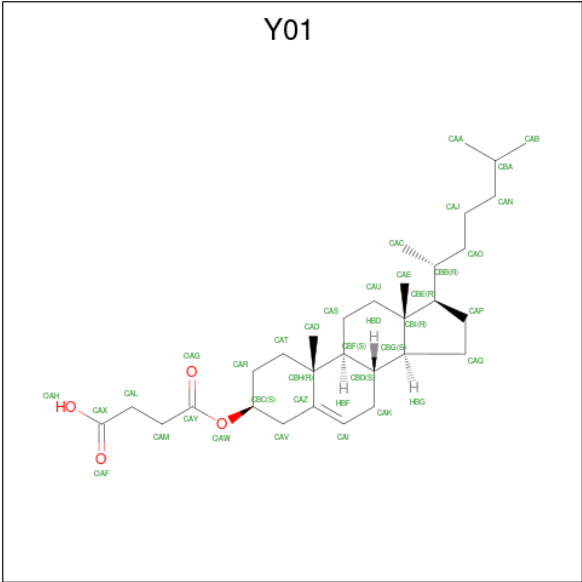
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	157	ARG	CYS	variant	UNP Q9H1D0
A	378	VAL	MET	variant	UNP Q9H1D0
A	681	THR	MET	variant	UNP Q9H1D0
B	157	ARG	CYS	variant	UNP Q9H1D0
B	378	VAL	MET	variant	UNP Q9H1D0
B	681	THR	MET	variant	UNP Q9H1D0
C	157	ARG	CYS	variant	UNP Q9H1D0
C	378	VAL	MET	variant	UNP Q9H1D0
C	681	THR	MET	variant	UNP Q9H1D0
D	157	ARG	CYS	variant	UNP Q9H1D0
D	378	VAL	MET	variant	UNP Q9H1D0
D	681	THR	MET	variant	UNP Q9H1D0

- Molecule 2 is a protein called Calmodulin-1.

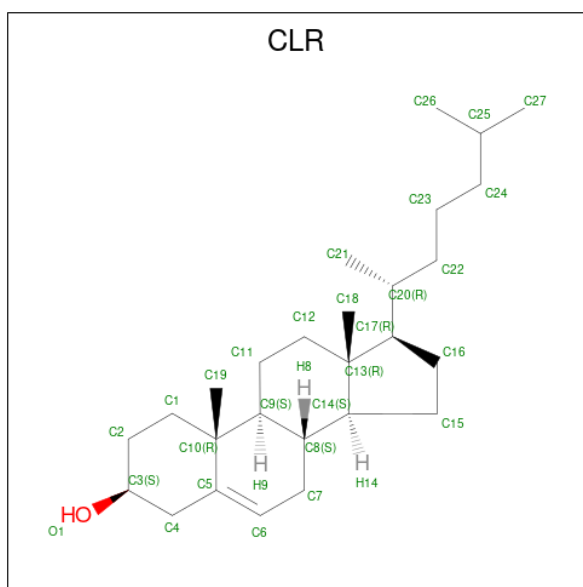
Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	148	Total	C	N	O	S	0	0
			1165	714	188	254	9		

- Molecule 3 is CHOLESTEROL HEMISUCCINATE (CCD ID: Y01) (formula: C₃₁H₅₀O₄).



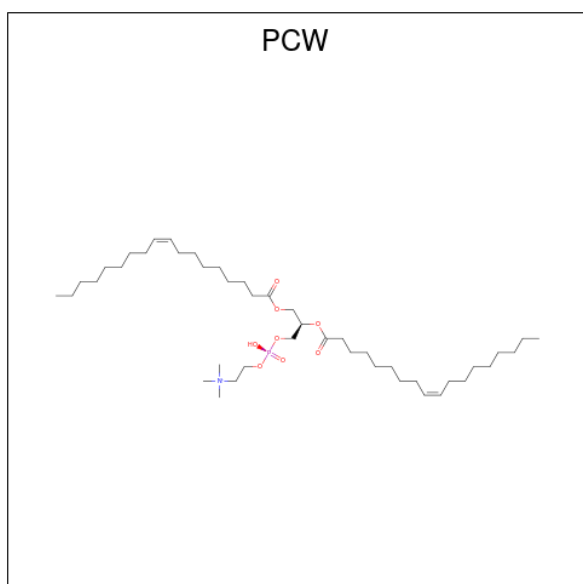
Mol	Chain	Residues	Atoms			AltConf
3	A	1	Total	C	O	0
			35	31	4	
3	A	1	Total	C	O	0
			35	31	4	
3	B	1	Total	C	O	0
			35	31	4	
3	B	1	Total	C	O	0
			35	31	4	
3	B	1	Total	C	O	0
			35	31	4	
3	C	1	Total	C	O	0
			35	31	4	
3	C	1	Total	C	O	0
			35	31	4	
3	D	1	Total	C	O	0
			35	31	4	

- Molecule 4 is CHOLESTEROL (CCD ID: CLR) (formula: C₂₇H₄₆O).



Mol	Chain	Residues	Atoms			AltConf
4	A	1	Total	C	O	0
			28	27	1	
4	B	1	Total	C	O	0
			28	27	1	
4	C	1	Total	C	O	0
			28	27	1	
4	D	1	Total	C	O	0
			28	27	1	

- Molecule 5 is 1,2-DIOLEOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PCW) (formula: $C_{44}H_{85}NO_8P$).



Mol	Chain	Residues	Atoms					AltConf
5	A	1	Total 51	C 41	N 1	O 8	P 1	0
5	A	1	Total 16	C 16				0
5	A	1	Total 16	C 16				0
5	A	1	Total 11	C 11				0
5	A	1	Total 8	C 8				0
5	A	1	Total 8	C 8				0
5	A	1	Total 16	C 16				0
5	A	1	Total 13	C 13				0
5	A	1	Total 43	C 33	N 1	O 8	P 1	0
5	A	1	Total 51	C 41	N 1	O 8	P 1	0
5	A	1	Total 8	C 8				0
5	A	1	Total 15	C 15				0
5	B	1	Total 15	C 15				0
5	B	1	Total 51	C 41	N 1	O 8	P 1	0
5	B	1	Total 16	C 16				0
5	B	1	Total 11	C 11				0
5	B	1	Total 8	C 8				0
5	B	1	Total 16	C 16				0
5	B	1	Total 13	C 13				0
5	B	1	Total 43	C 33	N 1	O 8	P 1	0
5	B	1	Total 51	C 41	N 1	O 8	P 1	0
5	C	1	Total 15	C 15				0

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Mol	Chain	Residues	Atoms	AltConf
5	C	1	Total C 16 16	0
5	C	1	Total C N O P 51 41 1 8 1	0
5	C	1	Total C 16 16	0
5	C	1	Total C 11 11	0
5	C	1	Total C 8 8	0
5	C	1	Total C 16 16	0
5	C	1	Total C 13 13	0
5	C	1	Total C N O P 43 33 1 8 1	0
5	C	1	Total C N O P 51 41 1 8 1	0
5	D	1	Total C 13 13	0
5	D	1	Total C N O P 43 33 1 8 1	0
5	D	1	Total C N O P 51 41 1 8 1	0
5	D	1	Total C 15 15	0
5	D	1	Total C 16 16	0
5	D	1	Total C N O P 51 41 1 8 1	0
5	D	1	Total C 16 16	0
5	D	1	Total C 16 16	0
5	D	1	Total C 11 11	0
5	D	1	Total C 8 8	0
5	D	1	Total C 8 8	0
5	D	1	Total C 16 16	0

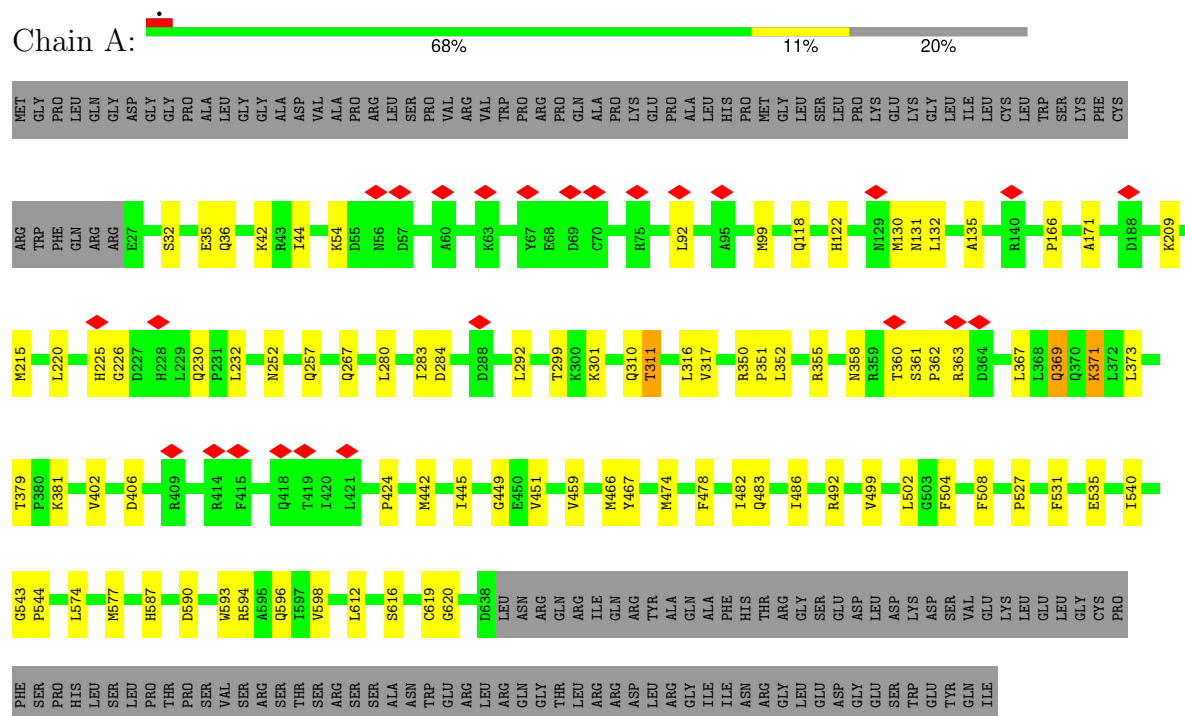
- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
6	A	2	Total 2	Ca 2	0
6	B	1	Total 1	Ca 1	0
6	E	4	Total 4	Ca 4	0

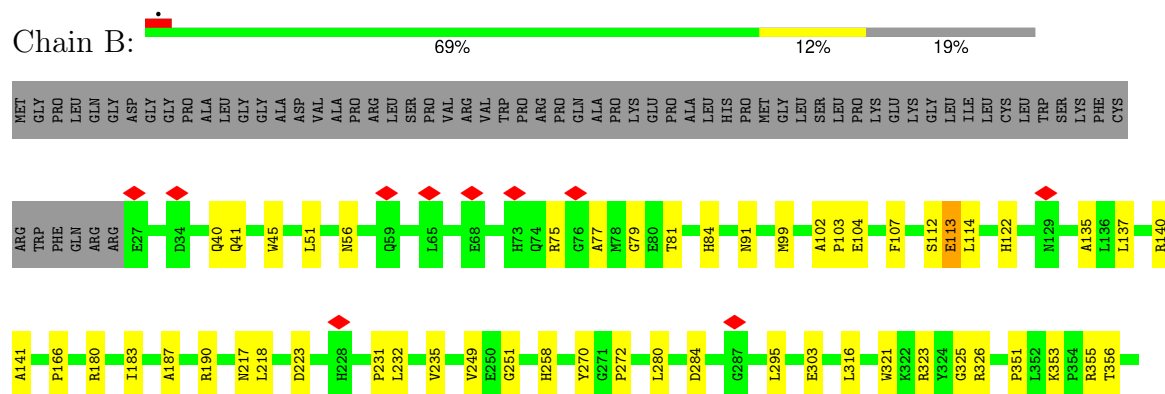
3 Residue-property plots

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

- Molecule 1: Transient receptor potential cation channel subfamily V member 6




- Molecule 1: Transient receptor potential cation channel subfamily V member 6

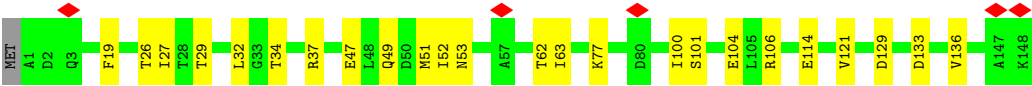


VAL
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ASN
ARG
GLY
LEU
GLU
ASP
GLY
SER
TRP
GLU
TYR

GLN
ILE

● Molecule 2: Calmodulin-1

Chain E:  83% 16%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	232058	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	55	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.351	Depositor
Minimum map value	-2.180	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.103	Depositor
Recommended contour level	0.326	Depositor
Map size (\AA)	273.152, 273.152, 273.152	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.067, 1.067, 1.067	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, CLR, PCW, Y01

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	0.26	0/5029	0.73	4/6824 (0.1%)
1	B	0.28	0/5104	0.77	1/6923 (0.0%)
1	C	0.28	0/5406	0.76	4/7326 (0.1%)
1	D	0.27	0/5088	0.73	4/6903 (0.1%)
2	E	0.21	0/1177	0.70	0/1580
All	All	0.27	0/21804	0.74	13/29556 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	5
1	C	0	1
1	D	0	4
All	All	0	11

There are no bond length outliers.

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	515	GLU	CA-C-N	7.14	132.75	121.17
1	D	515	GLU	C-N-CA	7.14	132.75	121.17
1	D	571	LEU	N-CA-C	-6.23	106.90	114.75
1	A	406	ASP	N-CA-C	-6.20	106.69	114.56
1	B	540	ILE	N-CA-C	-6.14	106.78	112.43

There are no chirality outliers.

5 of 11 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	230	GLN	Peptide
1	B	102	ALA	Peptide
1	B	113	GLU	Peptide
1	B	223	ASP	Peptide
1	B	231	PRO	Peptide

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	4916	0	4972	53	0
1	B	4991	0	5055	58	0
1	C	5286	0	5342	59	0
1	D	4972	0	5025	44	0
2	E	1165	0	1096	14	0
3	A	70	0	98	6	0
3	B	105	0	147	10	0
3	C	70	0	98	3	0
3	D	35	0	49	3	0
4	A	28	0	46	5	0
4	B	28	0	46	5	0
4	C	28	0	46	5	0
4	D	28	0	46	4	0
5	A	256	0	380	9	0
5	B	224	0	330	8	0
5	C	240	0	356	7	0
5	D	264	0	394	7	0
6	A	2	0	0	0	0
6	B	1	0	0	0	0
6	E	4	0	0	0	0
All	All	22713	0	23526	235	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 5.

The worst 5 of 235 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:426:HIS:CE1	3:B:803:Y01:HAL1	2.29	0.67
1:A:351:PRO:HB2	1:A:373:LEU:HB2	1.81	0.63
2:E:101:SER:H	2:E:104:GLU:HB2	1.65	0.62
1:A:531:PHE:HB2	3:B:801:Y01:HAS1	1.81	0.62
1:D:130:MET:HB3	1:D:132:LEU:H	1.65	0.61

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	610/765 (80%)	547 (90%)	60 (10%)	3 (0%)	25	55
1	B	618/765 (81%)	539 (87%)	75 (12%)	4 (1%)	22	51
1	C	649/765 (85%)	566 (87%)	80 (12%)	3 (0%)	25	55
1	D	615/765 (80%)	543 (88%)	68 (11%)	4 (1%)	19	48
2	E	146/149 (98%)	131 (90%)	15 (10%)	0	100	100
All	All	2638/3209 (82%)	2326 (88%)	298 (11%)	14 (0%)	27	55

5 of 14 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	371	LYS
1	B	114	LEU
1	B	104	GLU
1	B	639	LEU
1	C	114	LEU

5.3.2 Protein sidechains [i](#)

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

entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	531/662 (80%)	527 (99%)	4 (1%)	79	87
1	B	539/662 (81%)	537 (100%)	2 (0%)	89	93
1	C	568/662 (86%)	567 (100%)	1 (0%)	92	96
1	D	536/662 (81%)	534 (100%)	2 (0%)	89	93
2	E	126/127 (99%)	126 (100%)	0	100	100
All	All	2300/2775 (83%)	2291 (100%)	9 (0%)	88	93

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	183	ILE
1	D	367	LEU
1	A	612	LEU
1	B	367	LEU
1	B	397	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 60 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	365	ASN
1	D	572	ASN
1	C	91	ASN
1	D	546	ASN
2	E	143	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 62 ligands modelled in this entry, 7 are monoatomic - leaving 55 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
5	PCW	C	811	-	12,12,53	0.85	0	11,11,61	0.34	0
5	PCW	A	809	-	15,15,53	1.53	2 (13%)	14,14,61	0.73	0
5	PCW	A	804	-	15,15,53	1.49	2 (13%)	14,14,61	0.72	0
5	PCW	A	813	-	7,7,53	0.91	0	6,6,61	0.30	0
4	CLR	D	807	-	31,31,31	0.38	0	48,48,48	0.57	0
5	PCW	A	810	-	12,12,53	0.84	0	11,11,61	0.40	0
5	PCW	B	809	-	7,7,53	0.92	0	6,6,61	0.29	0
5	PCW	C	813	-	50,50,53	1.33	7 (14%)	56,58,61	0.95	3 (5%)
5	PCW	C	802	-	15,15,53	1.52	2 (13%)	14,14,61	0.70	0
4	CLR	A	802	-	31,31,31	0.38	0	48,48,48	0.54	0
5	PCW	D	802	-	42,42,53	1.42	7 (16%)	48,50,61	1.01	3 (6%)
5	PCW	A	808	-	7,7,53	0.92	0	6,6,61	0.31	0
5	PCW	B	808	-	10,10,53	0.95	0	9,9,61	0.45	0
5	PCW	D	809	-	15,15,53	1.52	2 (13%)	14,14,61	0.70	0
5	PCW	D	811	-	10,10,53	0.95	0	9,9,61	0.43	0
5	PCW	B	802	-	14,14,53	0.87	0	13,13,61	0.36	0
5	PCW	B	807	-	15,15,53	1.51	2 (13%)	14,14,61	0.70	0
5	PCW	C	810	-	15,15,53	1.52	2 (13%)	14,14,61	0.67	0
5	PCW	A	805	-	15,15,53	1.52	2 (13%)	14,14,61	0.67	0
5	PCW	A	812	-	50,50,53	1.34	7 (14%)	56,58,61	0.91	3 (5%)
3	Y01	A	814	-	38,38,38	0.53	0	57,57,57	0.79	2 (3%)
3	Y01	B	805	-	38,38,38	0.53	0	57,57,57	0.78	1 (1%)
3	Y01	B	801	-	38,38,38	0.51	0	57,57,57	0.97	5 (8%)
5	PCW	C	809	-	7,7,53	0.92	0	6,6,61	0.29	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PCW	D	804	-	14,14,53	0.88	0	13,13,61	0.35	0
5	PCW	C	806	-	50,50,53	1.33	6 (12%)	56,58,61	0.97	3 (5%)
5	PCW	D	808	-	50,50,53	1.31	6 (12%)	56,58,61	0.99	3 (5%)
5	PCW	A	811	-	42,42,53	1.42	7 (16%)	48,50,61	1.01	3 (6%)
5	PCW	C	808	-	10,10,53	0.95	0	9,9,61	0.43	0
5	PCW	A	806	-	10,10,53	0.95	0	9,9,61	0.49	0
4	CLR	B	804	-	31,31,31	0.41	0	48,48,48	0.81	0
4	CLR	C	804	-	31,31,31	0.45	0	48,48,48	0.71	0
5	PCW	B	806	-	50,50,53	1.33	6 (12%)	56,58,61	0.93	3 (5%)
5	PCW	B	811	-	12,12,53	0.82	0	11,11,61	0.40	0
5	PCW	D	801	-	12,12,53	0.85	0	11,11,61	0.34	0
3	Y01	D	806	-	38,38,38	0.51	0	57,57,57	0.72	0
3	Y01	C	805	-	38,38,38	0.54	0	57,57,57	0.95	4 (7%)
5	PCW	A	815	-	14,14,53	0.88	0	13,13,61	0.33	0
5	PCW	D	812	-	7,7,53	0.90	0	6,6,61	0.34	0
5	PCW	A	803	-	50,50,53	1.34	6 (12%)	56,58,61	0.96	3 (5%)
5	PCW	B	813	-	50,50,53	1.34	6 (12%)	56,58,61	0.94	3 (5%)
3	Y01	C	803	-	38,38,38	0.54	0	57,57,57	0.73	0
3	Y01	A	801	-	38,38,38	0.54	0	57,57,57	0.63	0
3	Y01	B	803	-	38,38,38	0.52	0	57,57,57	0.70	1 (1%)
5	PCW	B	810	-	15,15,53	1.51	2 (13%)	14,14,61	0.69	0
5	PCW	D	805	-	15,15,53	1.52	2 (13%)	14,14,61	0.68	0
5	PCW	D	814	-	15,15,53	1.53	2 (13%)	14,14,61	0.69	0
5	PCW	C	807	-	15,15,53	1.51	2 (13%)	14,14,61	0.69	0
5	PCW	B	812	-	42,42,53	1.43	7 (16%)	48,50,61	0.98	3 (6%)
5	PCW	D	813	-	7,7,53	0.92	0	6,6,61	0.27	0
5	PCW	C	812	-	42,42,53	1.40	6 (14%)	48,50,61	0.99	3 (6%)
5	PCW	A	807	-	7,7,53	0.90	0	6,6,61	0.34	0
5	PCW	D	810	-	15,15,53	1.53	2 (13%)	14,14,61	0.69	0
5	PCW	D	803	-	50,50,53	1.34	6 (12%)	56,58,61	0.93	3 (5%)
5	PCW	C	801	-	14,14,53	0.87	0	13,13,61	0.41	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PCW	C	811	-	-	4/10/10/57	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PCW	A	809	-	-	8/13/13/57	-
5	PCW	A	804	-	-	6/13/13/57	-
5	PCW	A	813	-	-	1/5/5/57	-
4	CLR	D	807	-	-	7/10/68/68	0/4/4/4
5	PCW	A	810	-	-	4/10/10/57	-
5	PCW	B	809	-	-	2/5/5/57	-
5	PCW	C	813	-	-	28/54/54/57	-
5	PCW	C	802	-	-	9/13/13/57	-
4	CLR	A	802	-	-	5/10/68/68	0/4/4/4
5	PCW	D	802	-	-	20/46/46/57	-
5	PCW	A	808	-	-	1/5/5/57	-
5	PCW	B	808	-	-	4/8/8/57	-
5	PCW	D	809	-	-	7/13/13/57	-
5	PCW	D	811	-	-	4/8/8/57	-
5	PCW	B	802	-	-	6/12/12/57	-
5	PCW	B	807	-	-	10/13/13/57	-
5	PCW	C	810	-	-	8/13/13/57	-
5	PCW	A	805	-	-	5/13/13/57	-
5	PCW	A	812	-	-	24/54/54/57	-
3	Y01	A	814	-	-	14/19/77/77	0/4/4/4
3	Y01	B	805	-	-	14/19/77/77	0/4/4/4
3	Y01	B	801	-	-	6/19/77/77	0/4/4/4
5	PCW	C	809	-	-	2/5/5/57	-
5	PCW	D	804	-	-	6/12/12/57	-
5	PCW	C	806	-	-	29/54/54/57	-
5	PCW	D	808	-	-	24/54/54/57	-
5	PCW	A	811	-	-	22/46/46/57	-
5	PCW	C	808	-	-	1/8/8/57	-
5	PCW	A	806	-	-	3/8/8/57	-
4	CLR	B	804	-	-	8/10/68/68	0/4/4/4
4	CLR	C	804	-	-	7/10/68/68	0/4/4/4
5	PCW	B	806	-	-	24/54/54/57	-
5	PCW	B	811	-	-	4/10/10/57	-
5	PCW	D	801	-	-	5/10/10/57	-
3	Y01	D	806	-	-	10/19/77/77	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	Y01	C	805	-	-	12/19/77/77	0/4/4/4
5	PCW	A	815	-	-	7/12/12/57	-
5	PCW	D	812	-	-	3/5/5/57	-
5	PCW	A	803	-	-	23/54/54/57	-
5	PCW	B	813	-	-	29/54/54/57	-
3	Y01	C	803	-	-	9/19/77/77	0/4/4/4
3	Y01	A	801	-	-	10/19/77/77	0/4/4/4
3	Y01	B	803	-	-	14/19/77/77	0/4/4/4
5	PCW	B	810	-	-	8/13/13/57	-
5	PCW	D	805	-	-	6/13/13/57	-
5	PCW	D	814	-	-	6/13/13/57	-
5	PCW	C	807	-	-	9/13/13/57	-
5	PCW	B	812	-	-	18/46/46/57	-
5	PCW	D	813	-	-	2/5/5/57	-
5	PCW	C	812	-	-	20/46/46/57	-
5	PCW	A	807	-	-	2/5/5/57	-
5	PCW	D	810	-	-	11/13/13/57	-
5	PCW	D	803	-	-	24/54/54/57	-
5	PCW	C	801	-	-	6/12/12/57	-

The worst 5 of 101 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	810	PCW	C20-C19	4.15	1.55	1.31
5	A	809	PCW	C20-C19	4.15	1.55	1.31
5	D	809	PCW	C20-C19	4.14	1.55	1.31
5	D	810	PCW	C20-C19	4.14	1.55	1.31
5	D	814	PCW	C20-C19	4.13	1.55	1.31

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	808	PCW	O2-C31-C32	4.22	120.61	111.48
5	A	803	PCW	O2-C31-C32	4.21	120.59	111.48
5	C	806	PCW	O2-C31-C32	4.16	120.48	111.48
5	D	802	PCW	O2-C31-C32	4.06	120.26	111.48
5	A	811	PCW	O2-C31-C32	3.99	120.12	111.48

There are no chirality outliers.

5 of 561 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	814	Y01	CAM-CAY-OAW-CBC
3	B	803	Y01	CAO-CBB-CBE-CAP
3	B	803	Y01	CAO-CBB-CBE-CBI
3	B	803	Y01	CAC-CBB-CBE-CBI
3	B	805	Y01	CAR-CBC-OAW-CAY

There are no ring outliers.

33 monomers are involved in 67 short contacts:

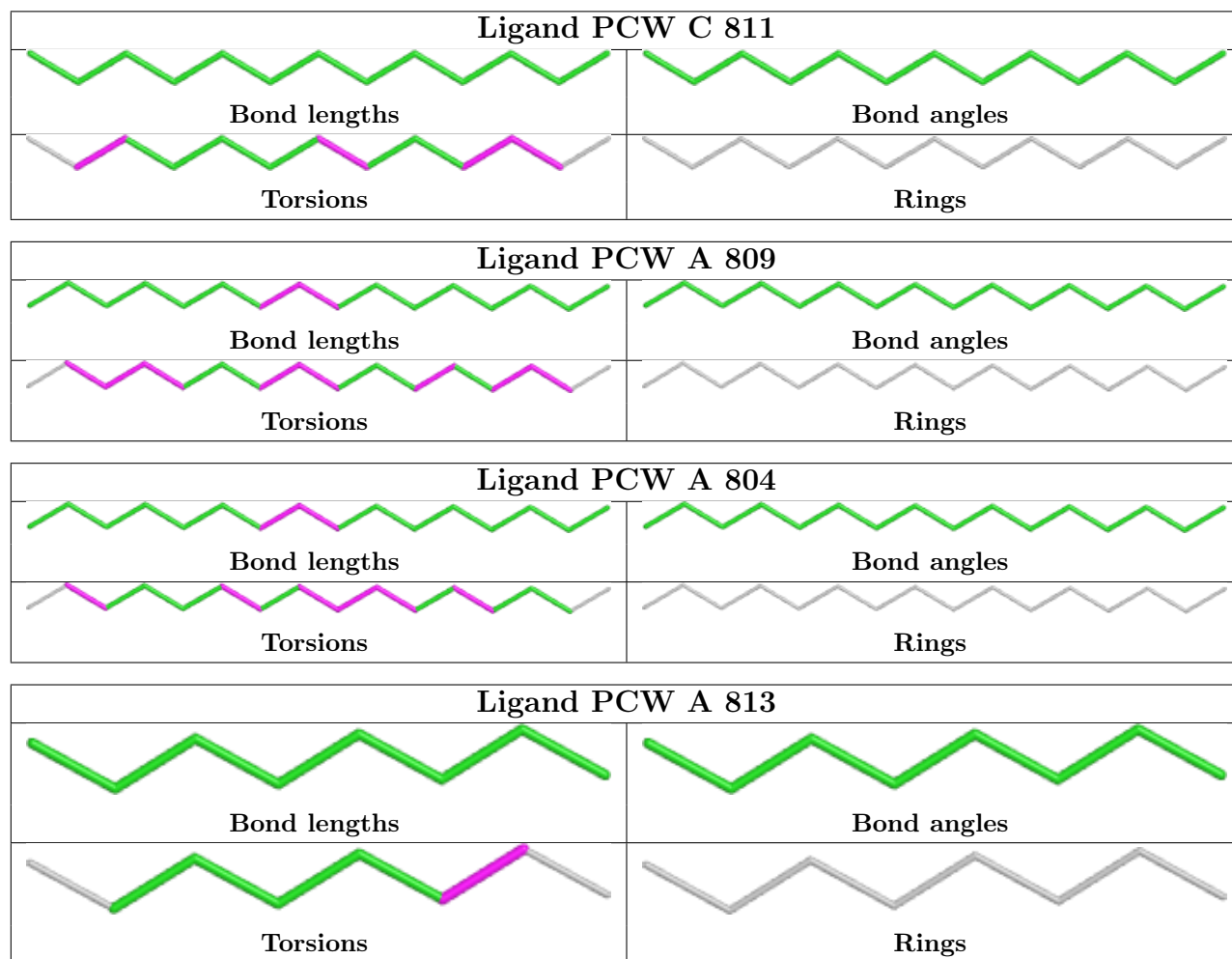
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	804	PCW	1	0
4	D	807	CLR	4	0
5	A	810	PCW	1	0
5	C	813	PCW	1	0
5	C	802	PCW	1	0
4	A	802	CLR	5	0
5	B	808	PCW	1	0
5	D	809	PCW	1	0
5	B	802	PCW	1	0
5	C	810	PCW	1	0
5	A	812	PCW	2	0
3	A	814	Y01	3	0
3	B	805	Y01	3	0
3	B	801	Y01	4	0
5	D	804	PCW	3	0
5	C	806	PCW	2	0
5	D	808	PCW	3	0
5	A	811	PCW	1	0
5	C	808	PCW	1	0
5	A	806	PCW	2	0
4	B	804	CLR	5	0
4	C	804	CLR	5	0
5	B	806	PCW	4	0
3	D	806	Y01	3	0
3	C	805	Y01	1	0
5	A	815	PCW	1	0
5	A	803	PCW	2	0
3	C	803	Y01	2	0
3	A	801	Y01	3	0
3	B	803	Y01	3	0

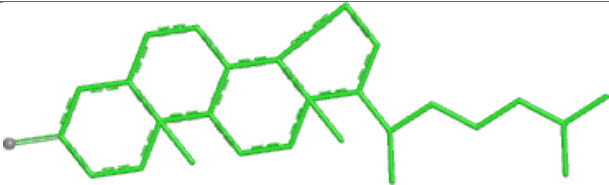
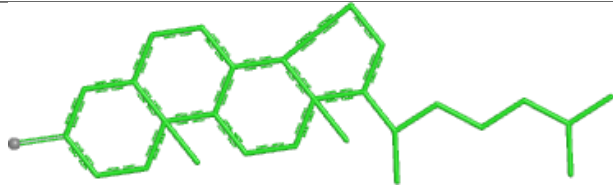
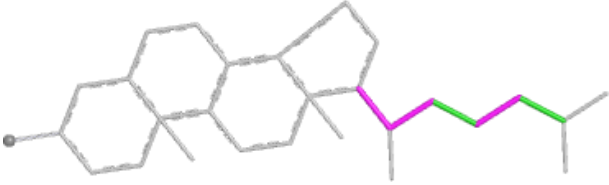
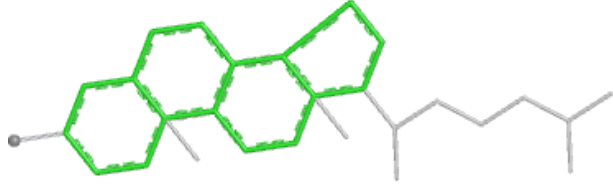
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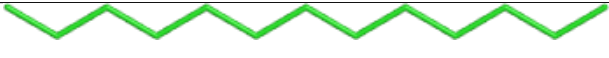
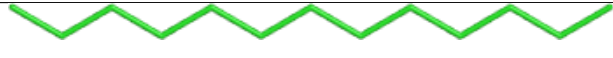


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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	810	PCW	1	0
5	B	812	PCW	1	0
5	C	801	PCW	1	0

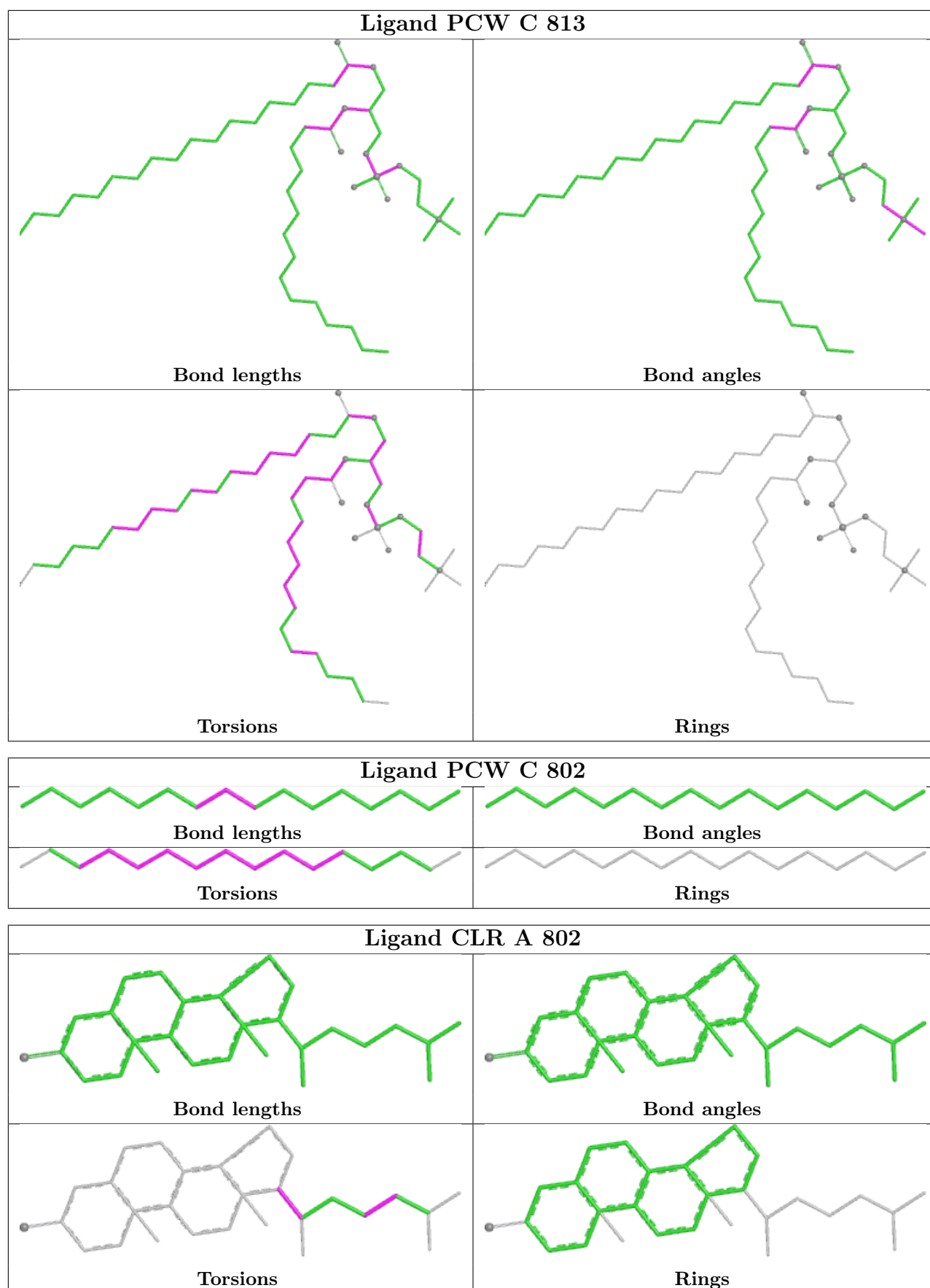
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

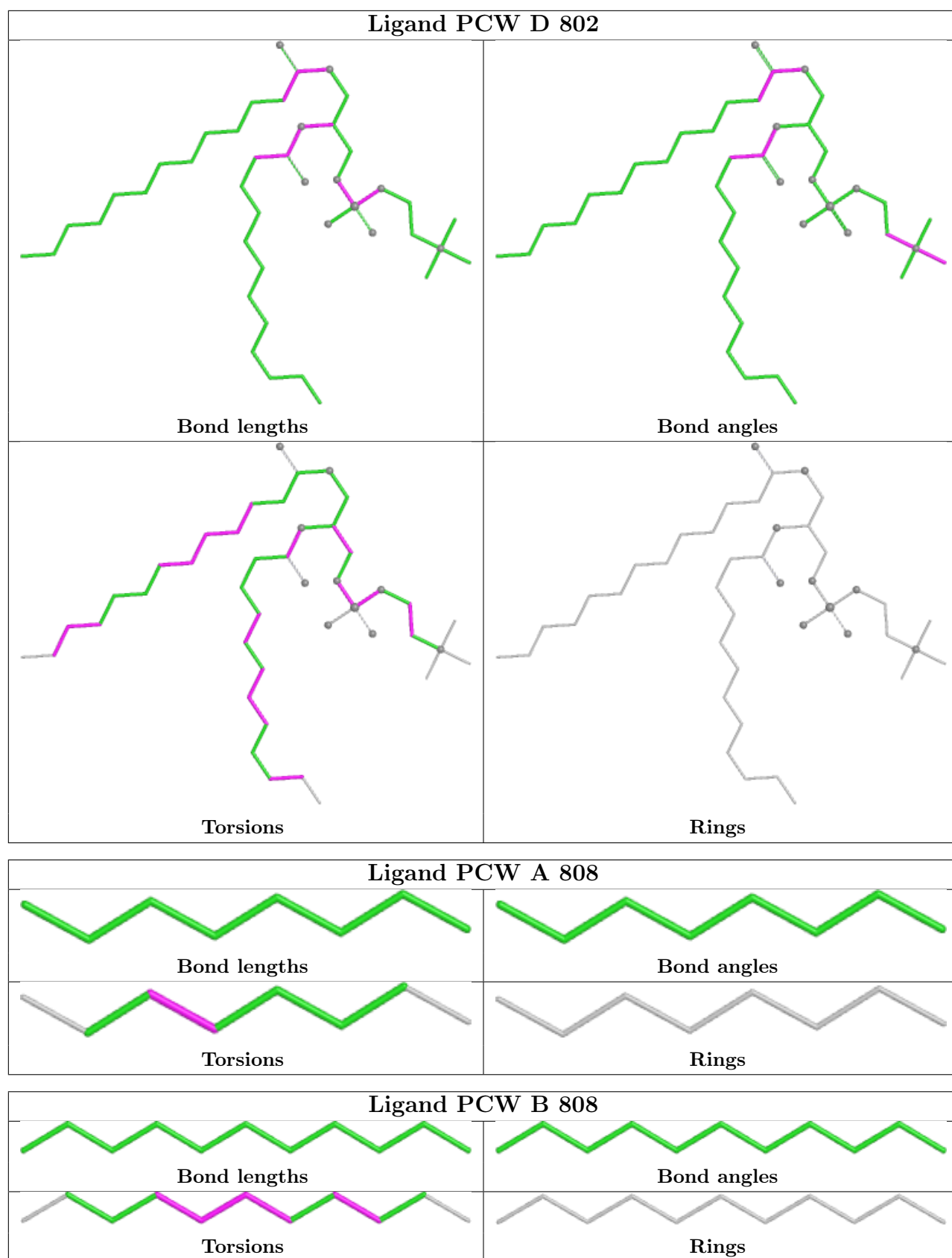



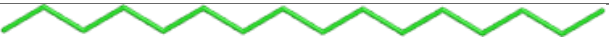


Ligand CLR D 807	
	
Bond lengths	Bond angles
	
Torsions	Rings





Ligand PCW A 810	
	
Bond lengths	Bond angles
	
Torsions	Rings





Ligand PCW B 809	
	
Bond lengths	Bond angles
	
Torsions	Rings

















Ligand PCW D 809			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

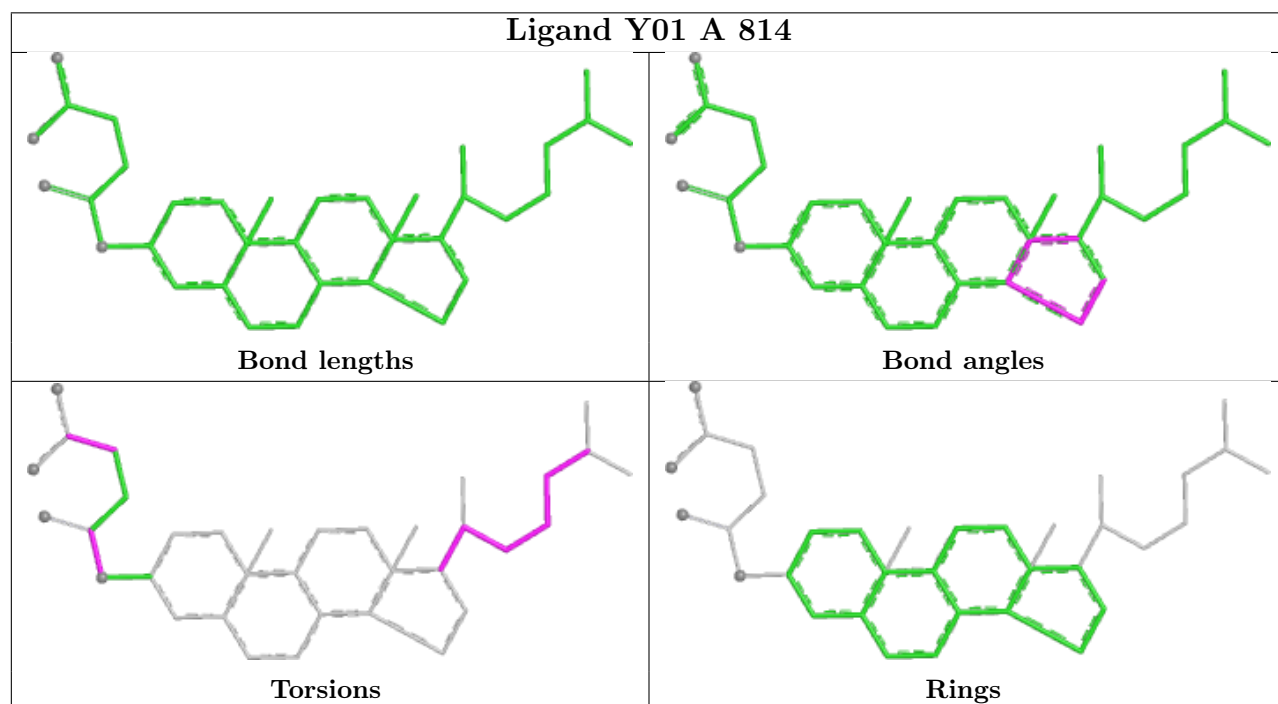
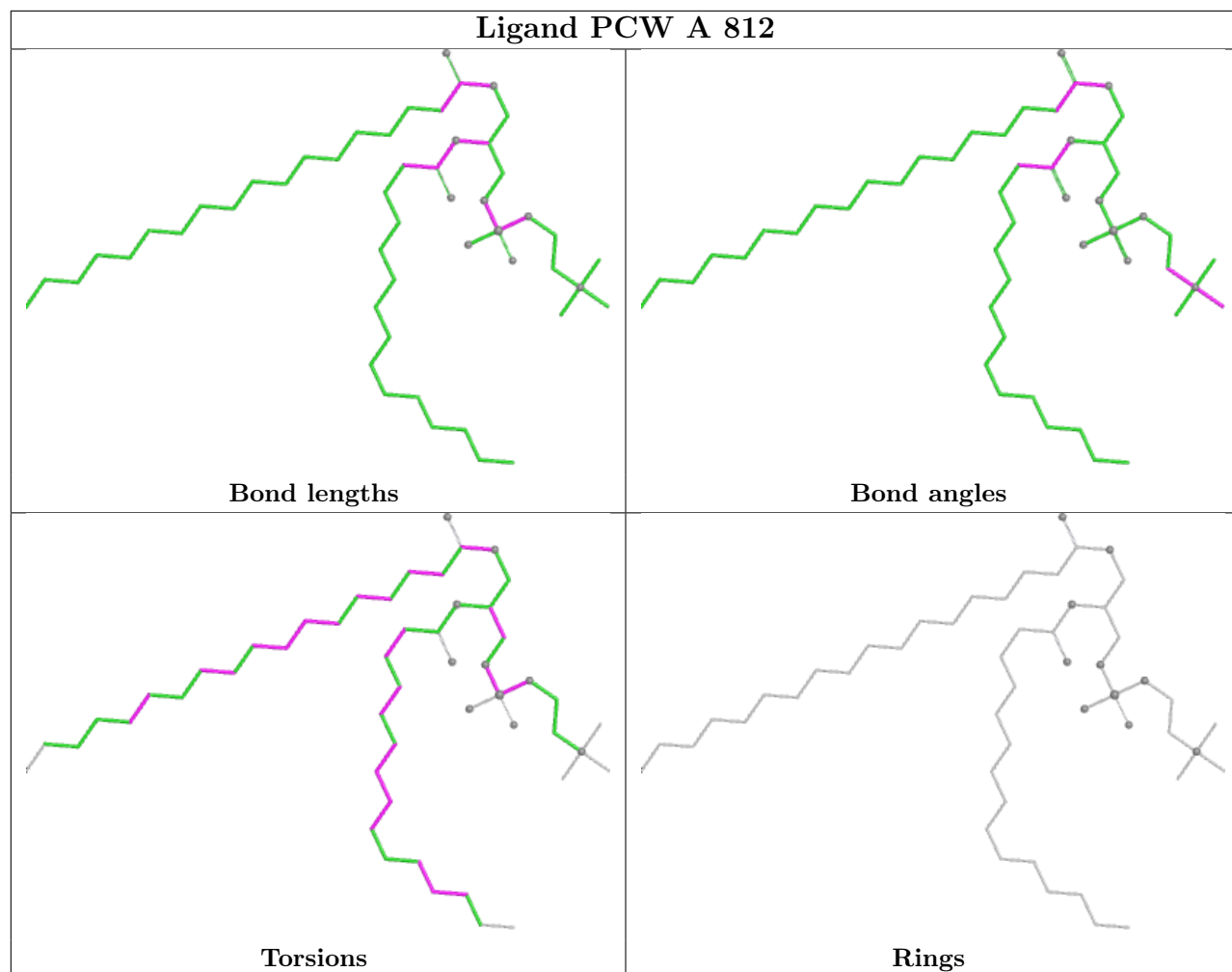
Ligand PCW D 811			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

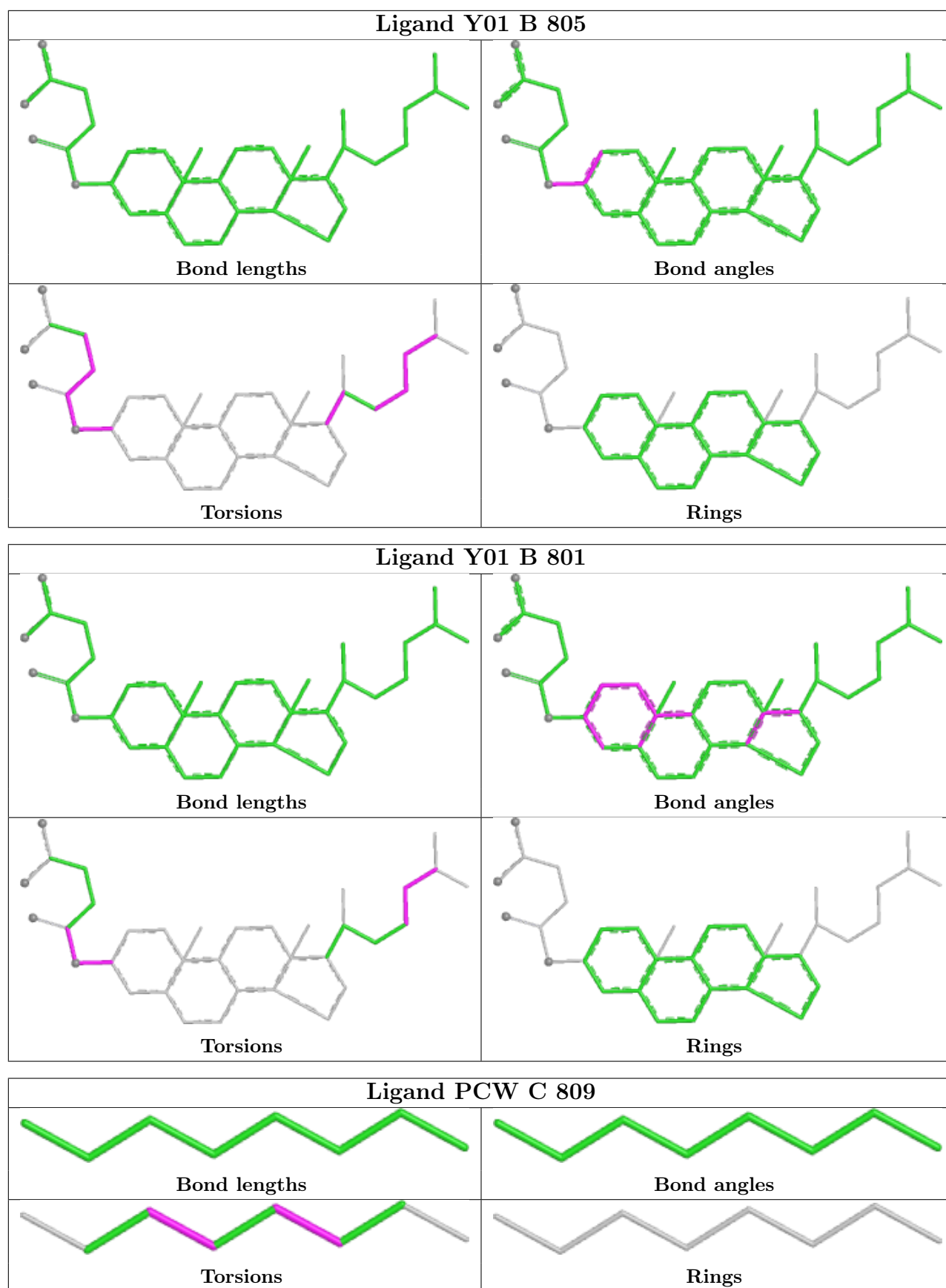
Ligand PCW B 802			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

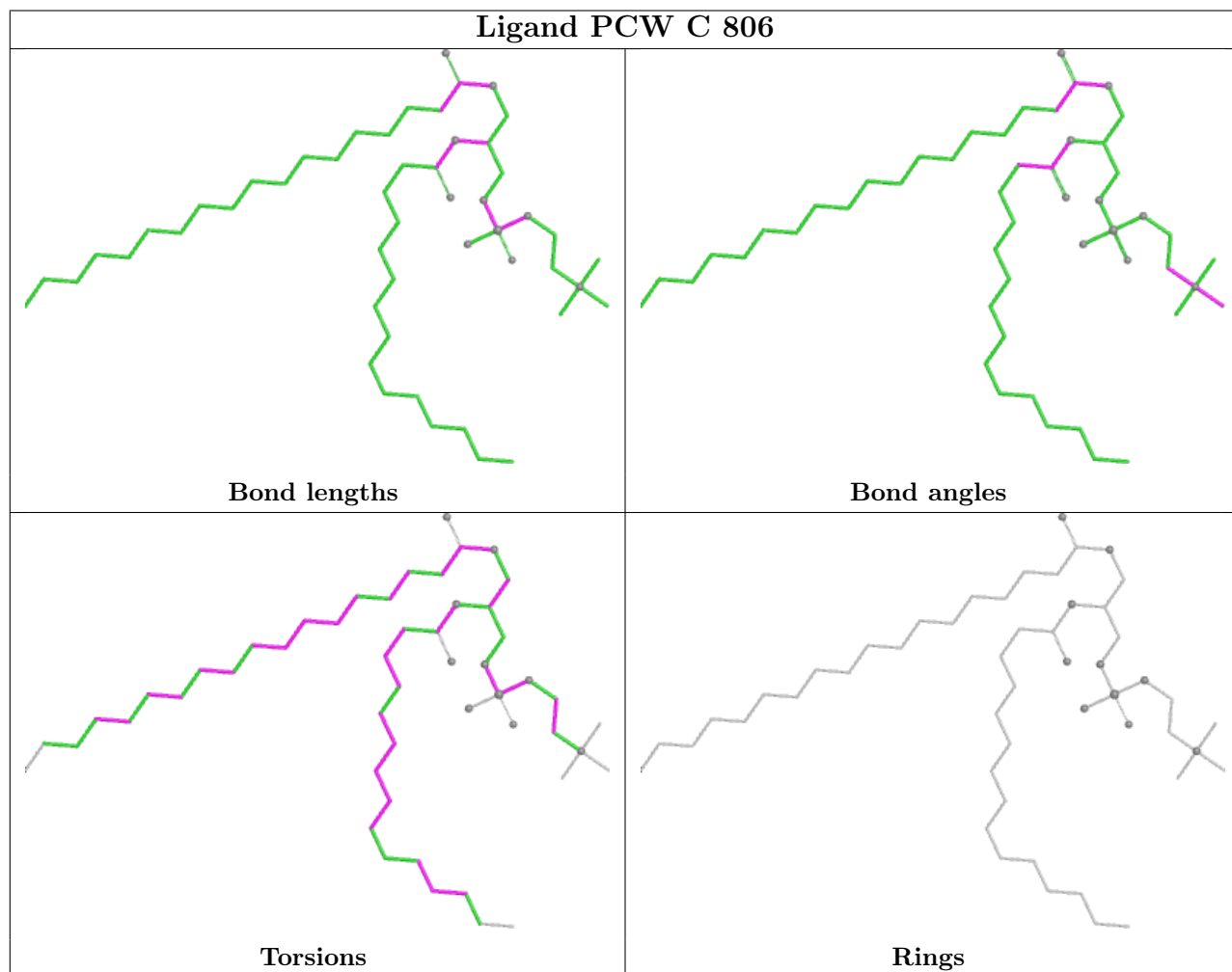
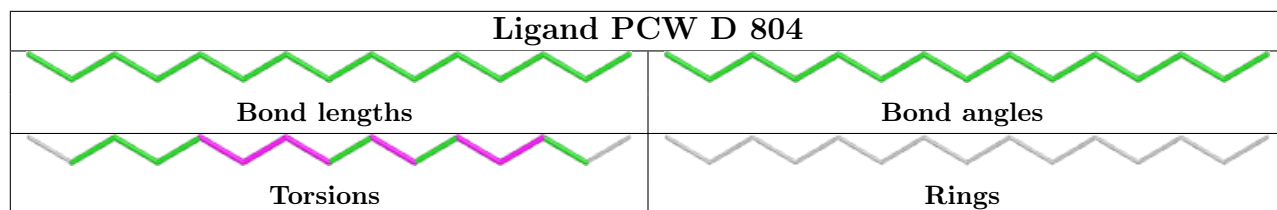
Ligand PCW B 807			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

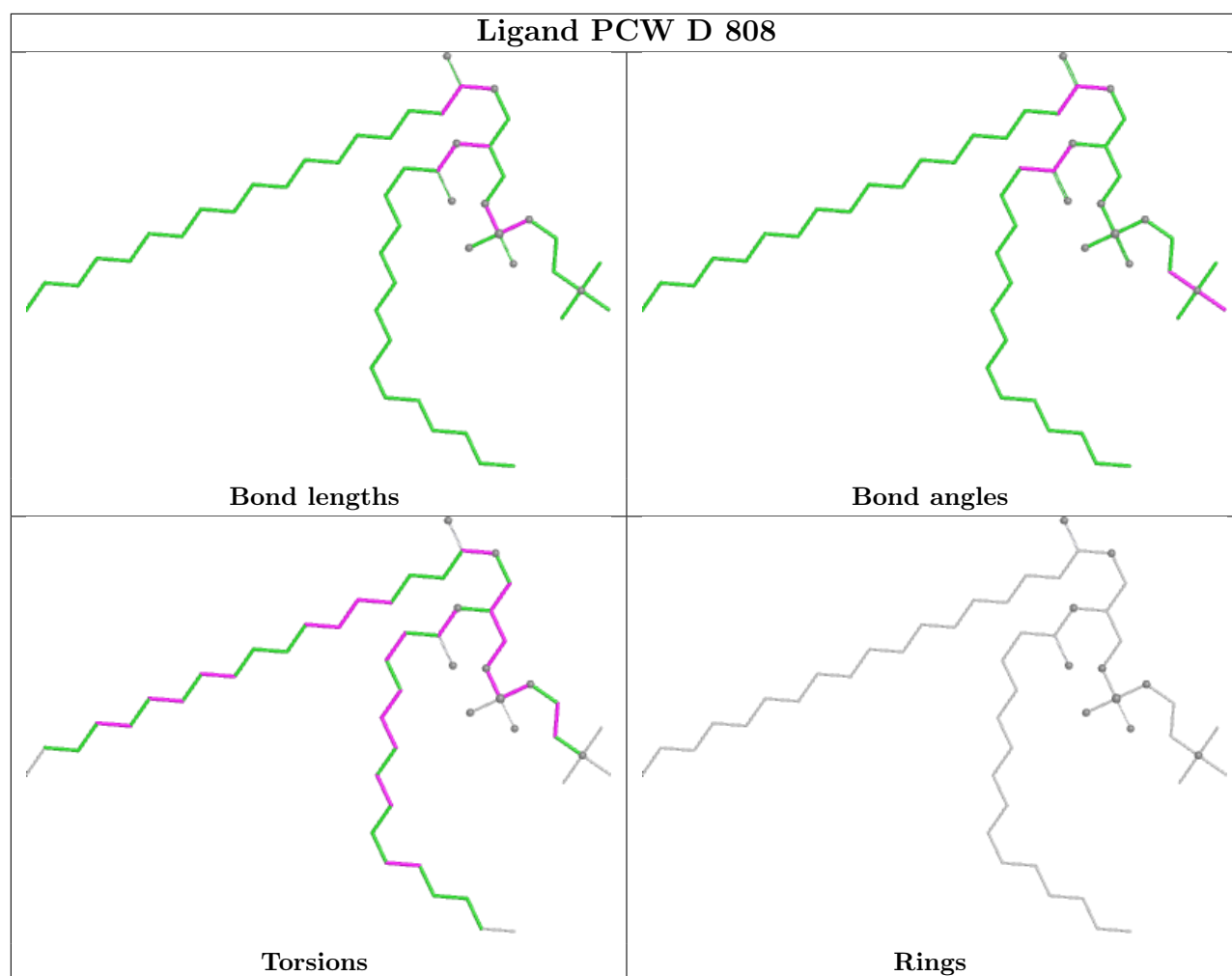
Ligand PCW C 810			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

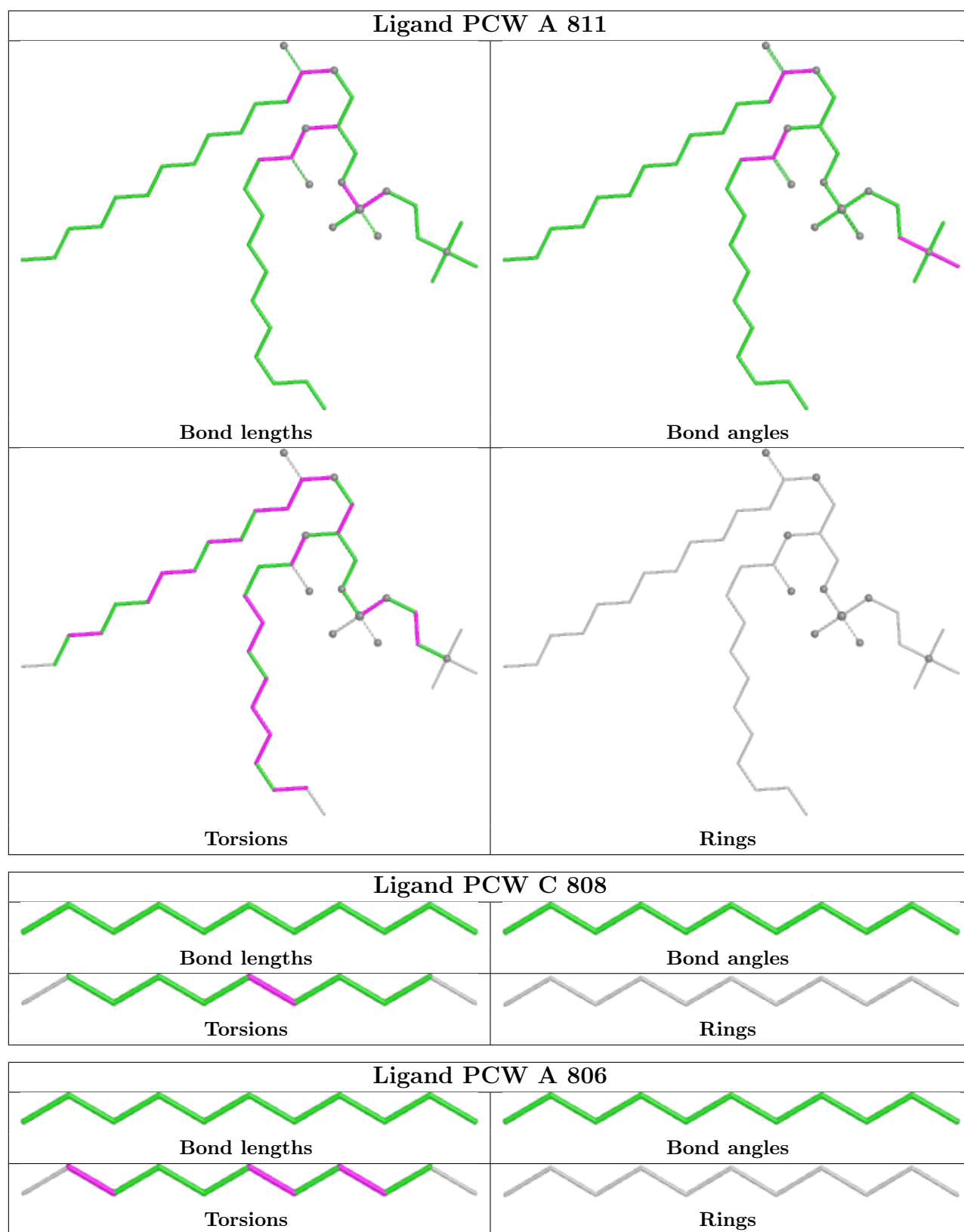
Ligand PCW A 805			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

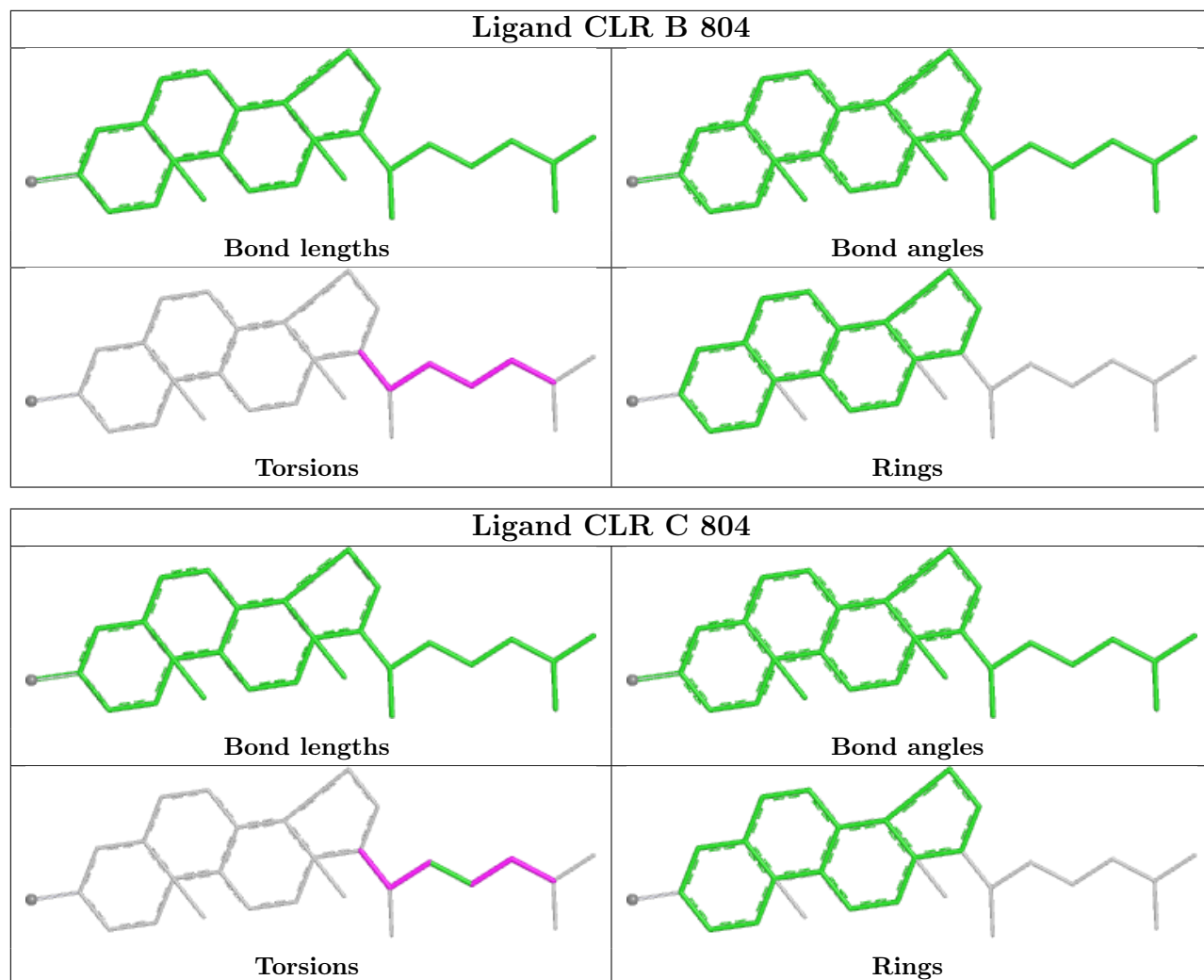


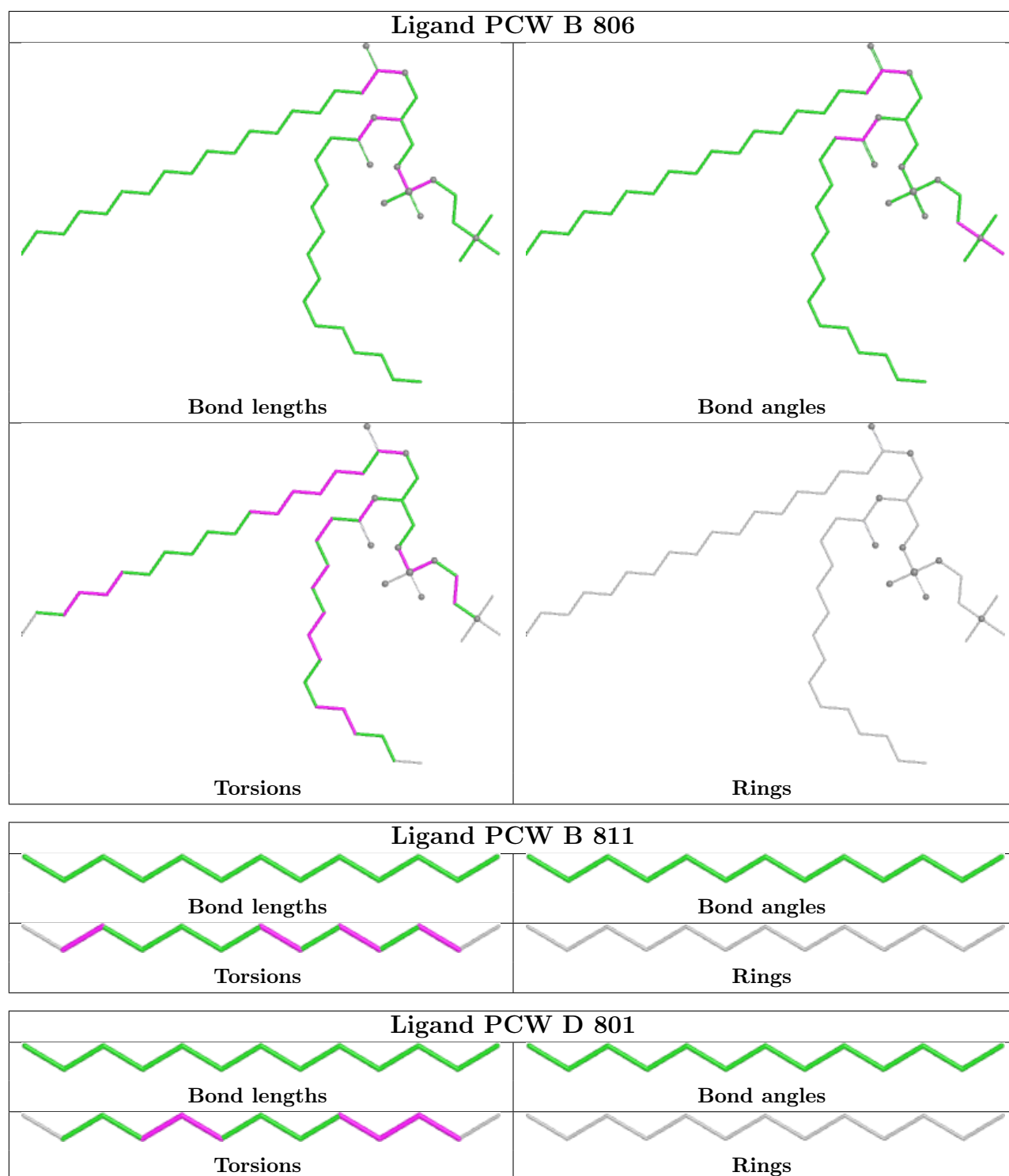


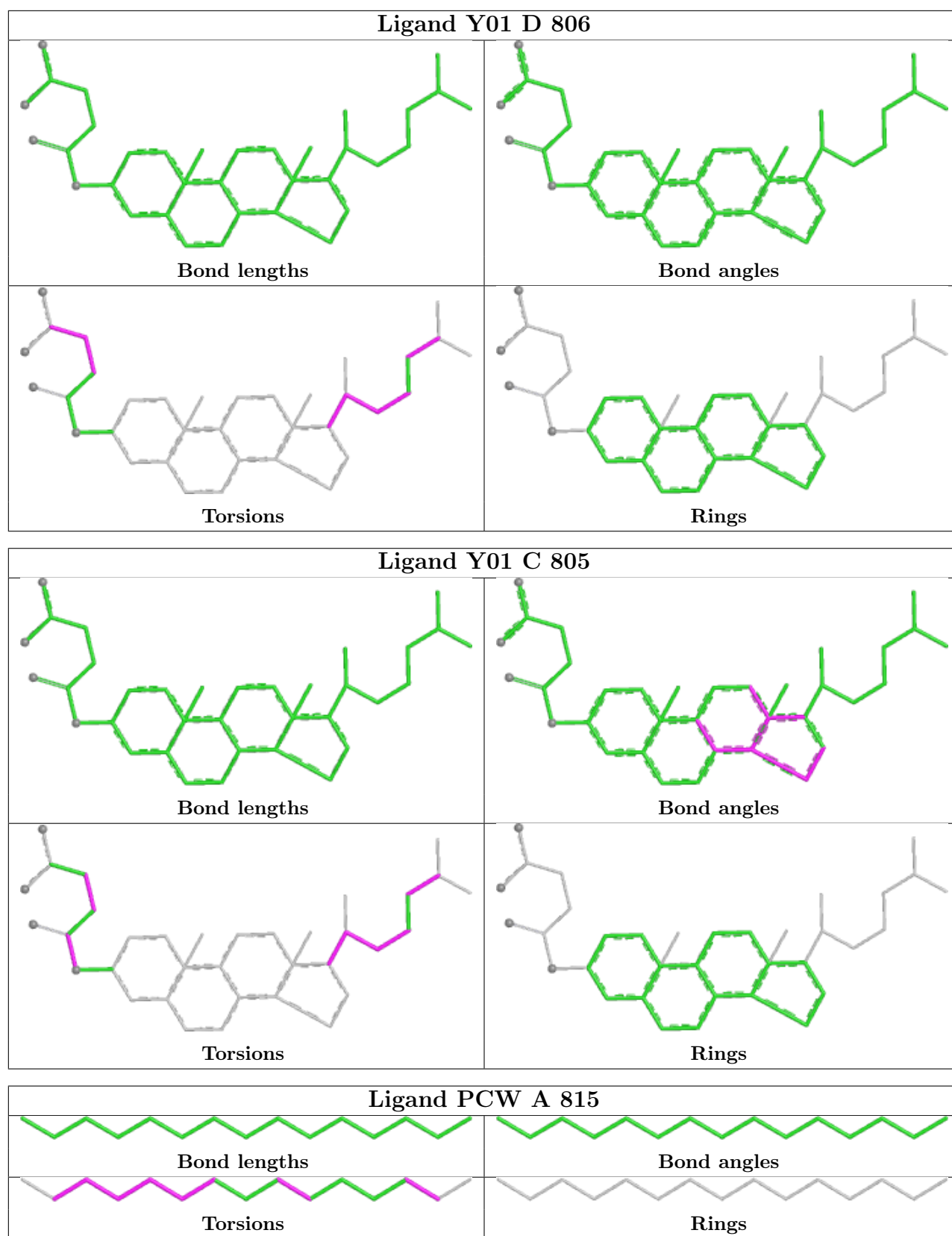


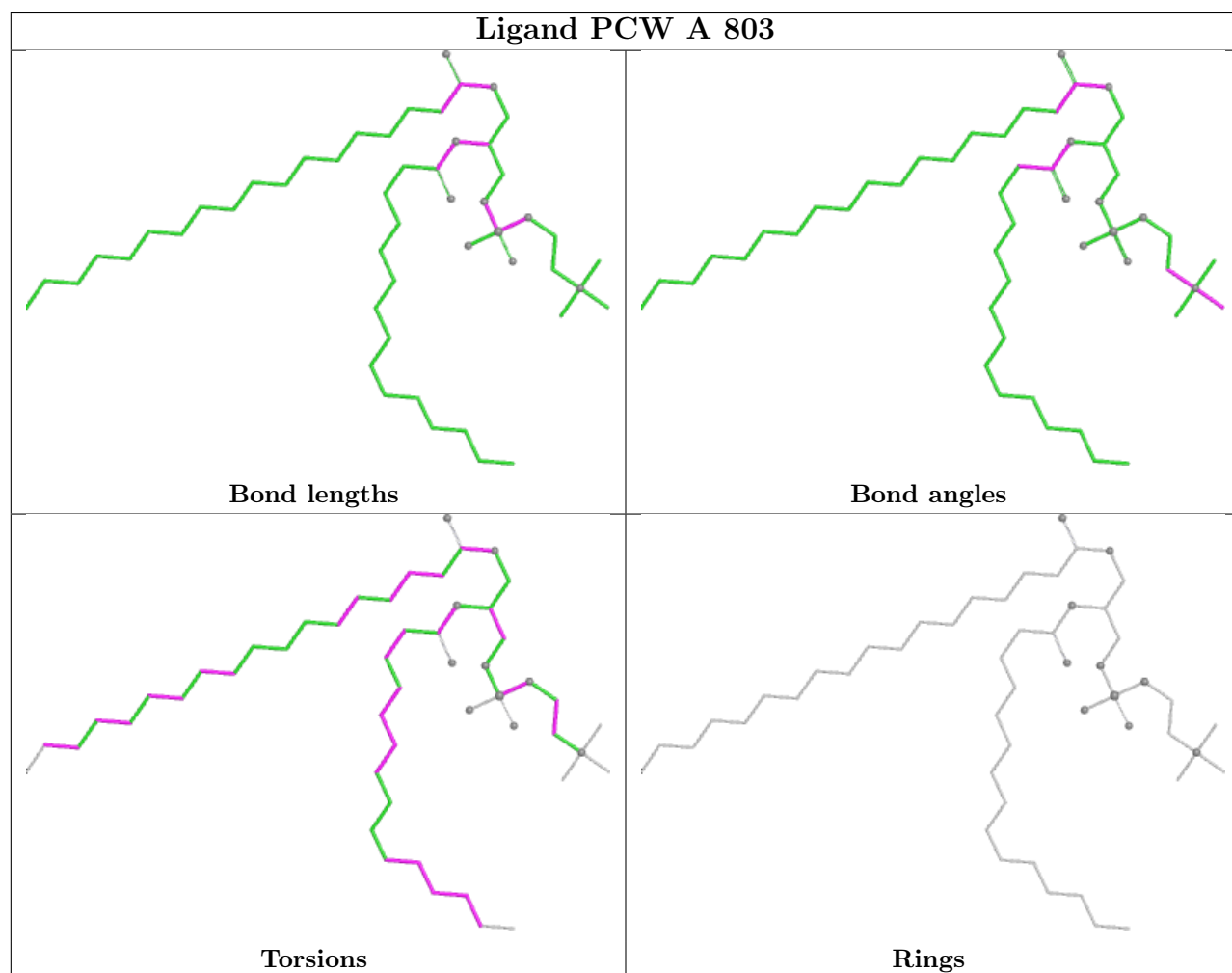
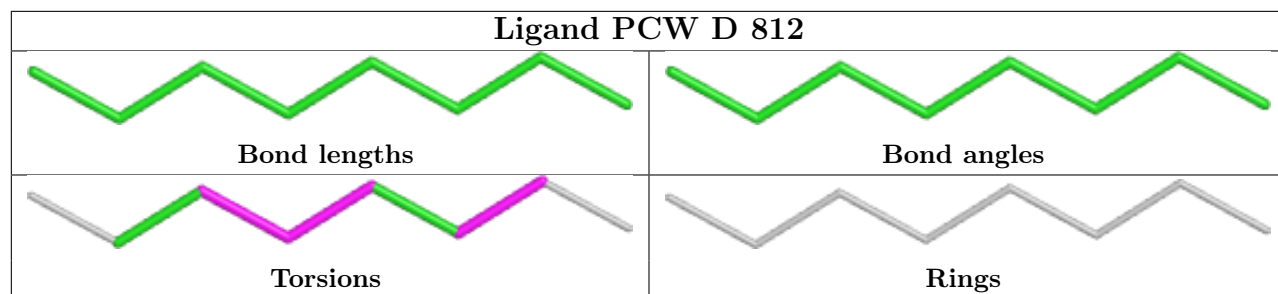


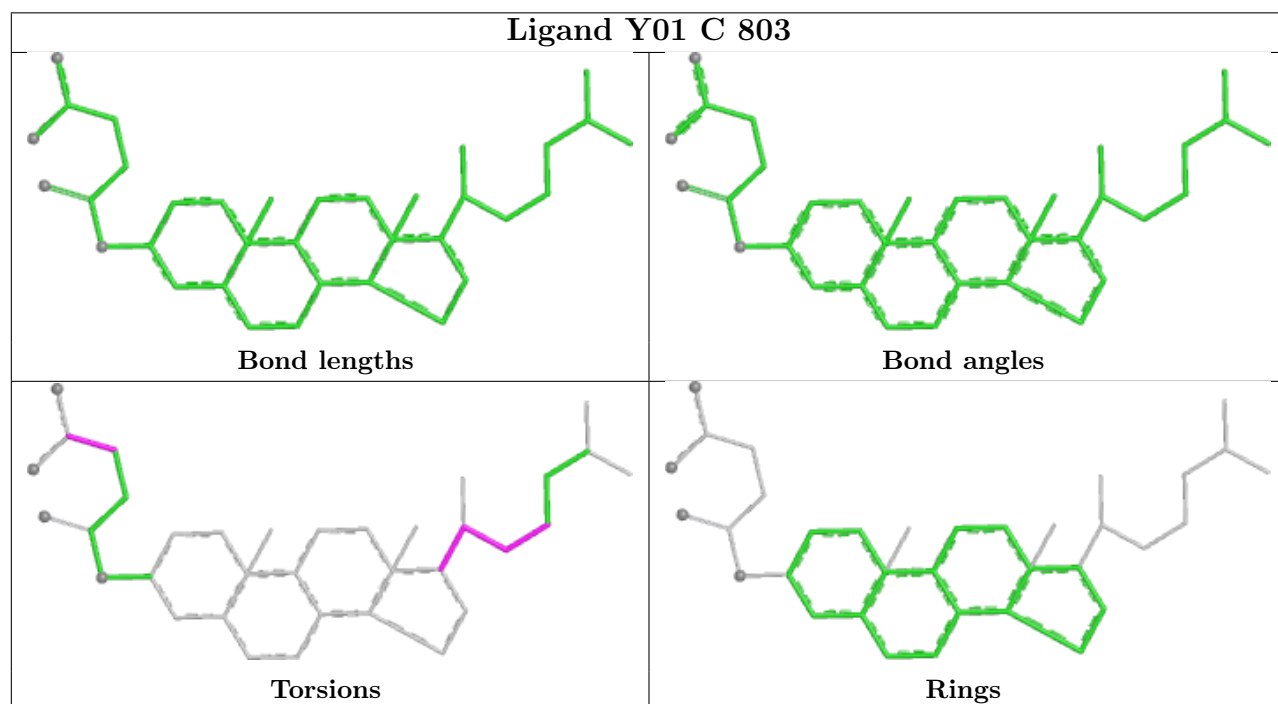
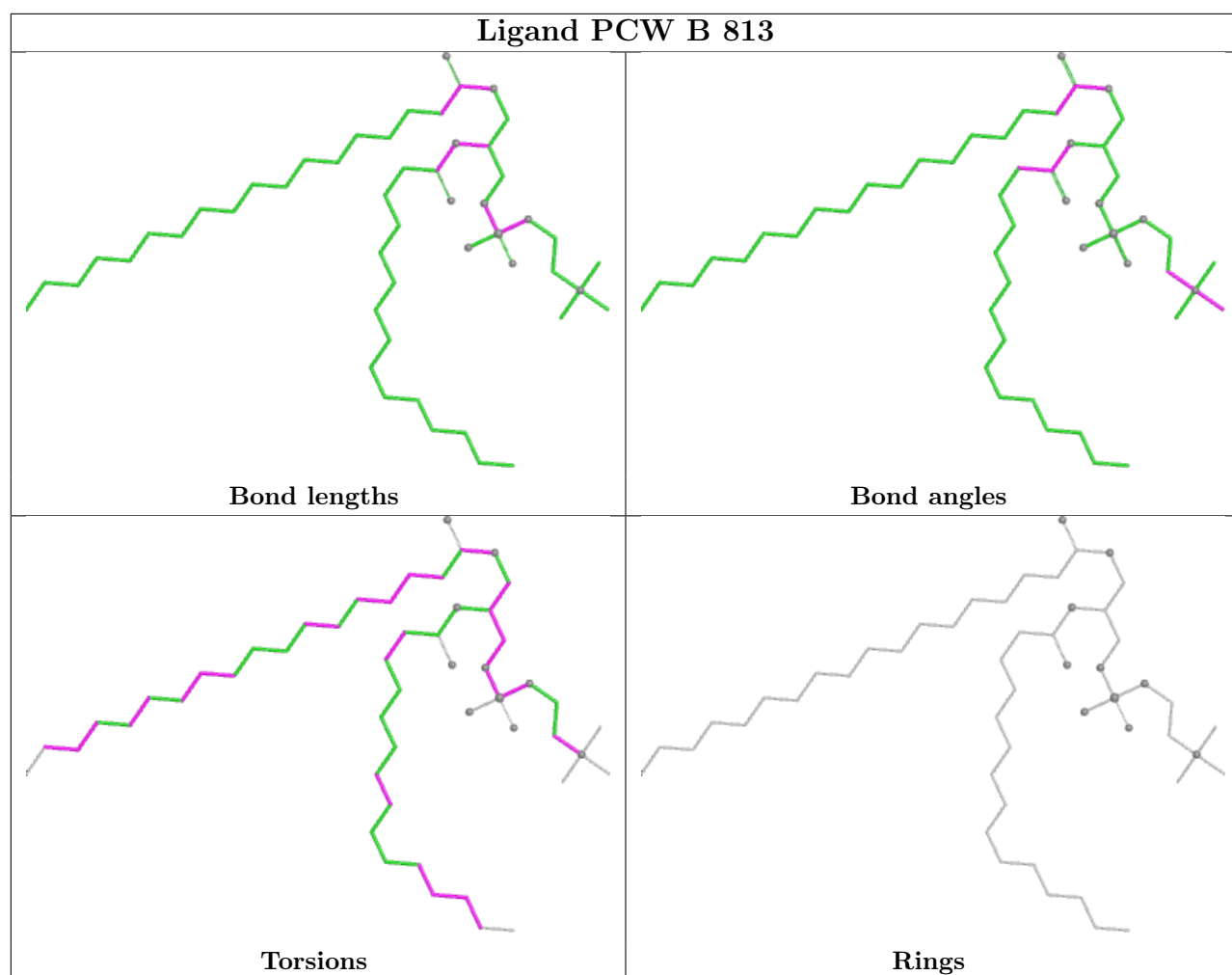


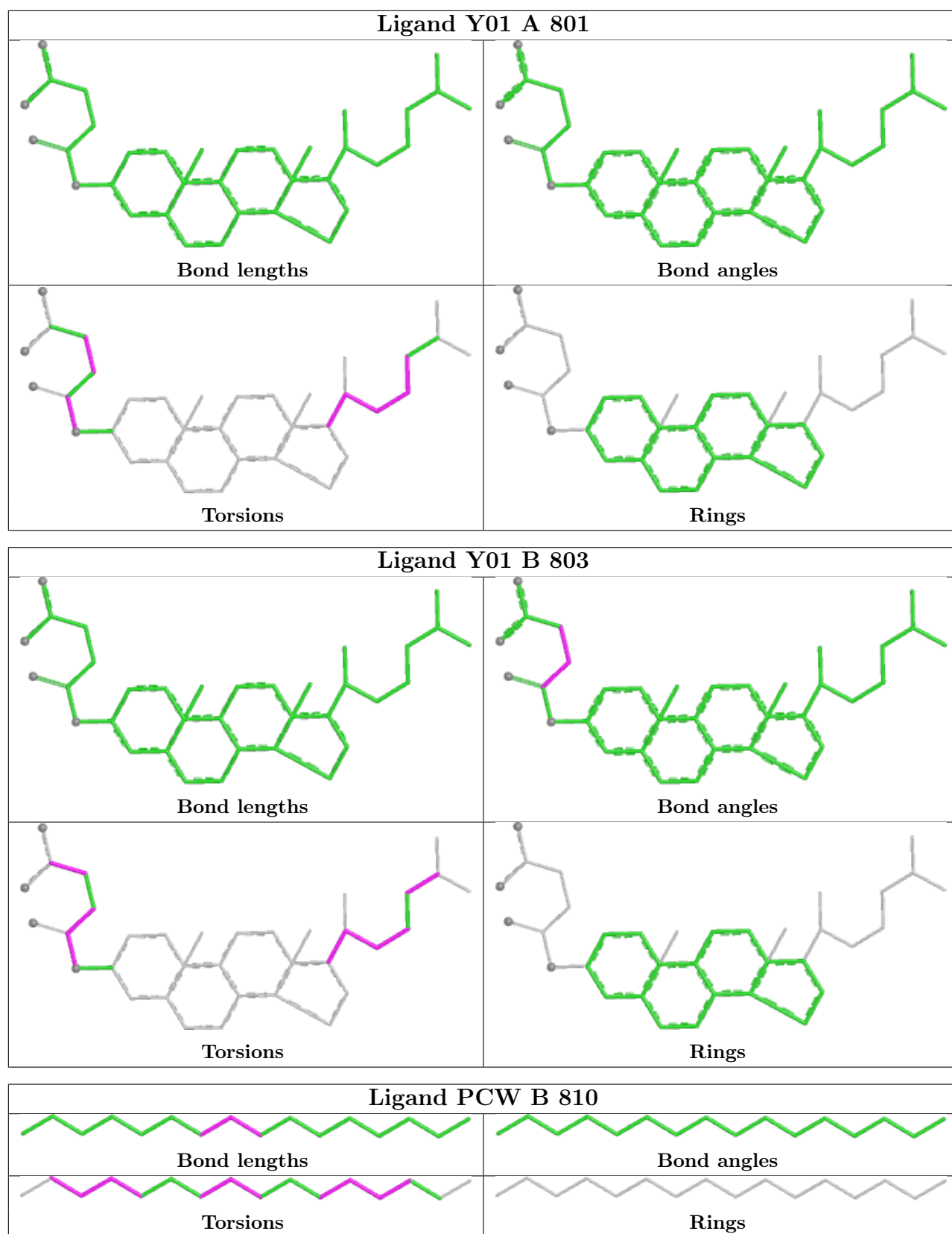


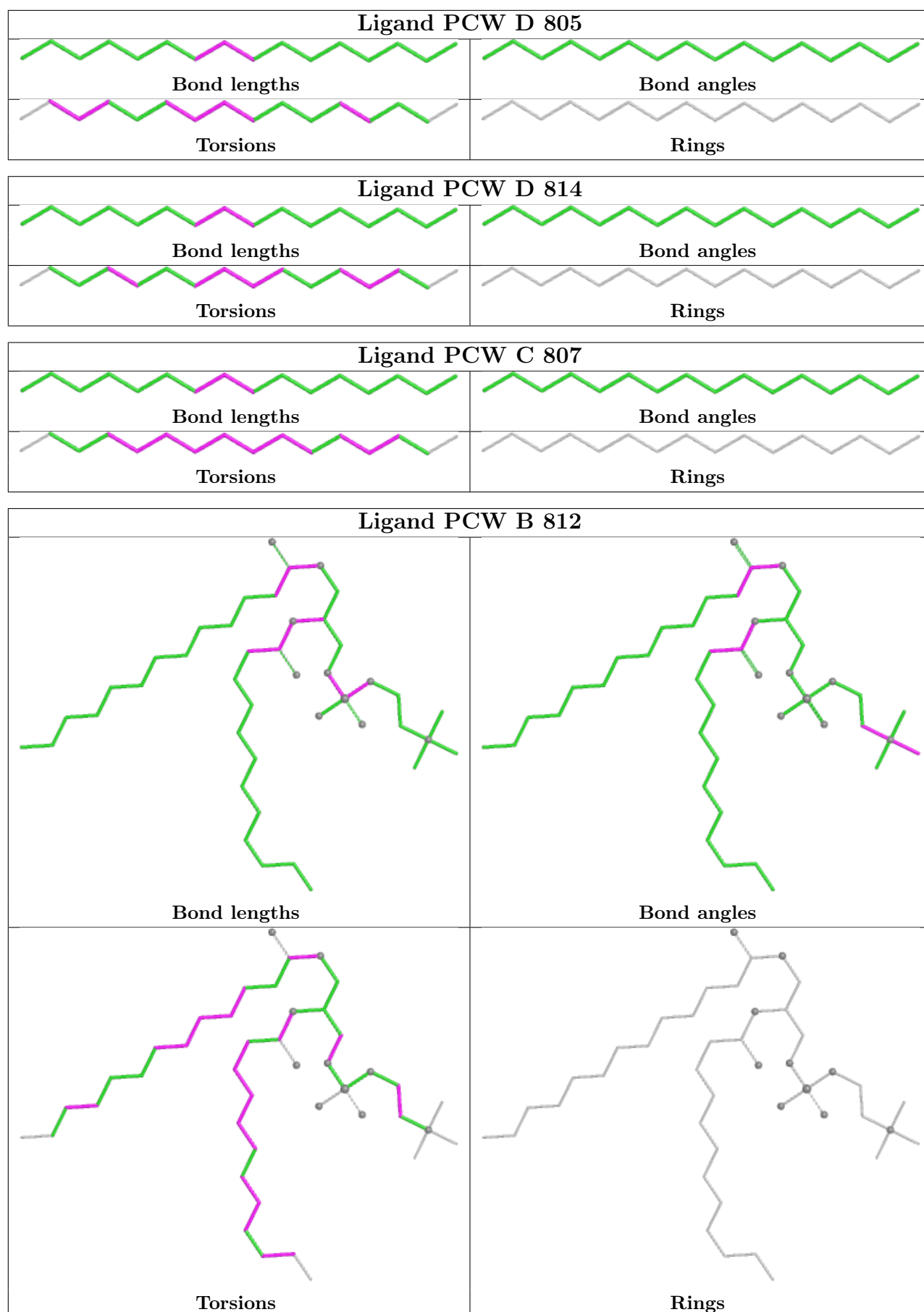


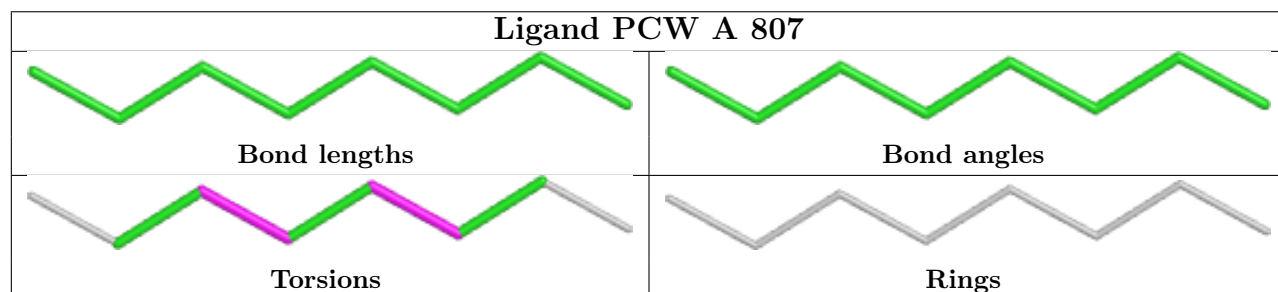
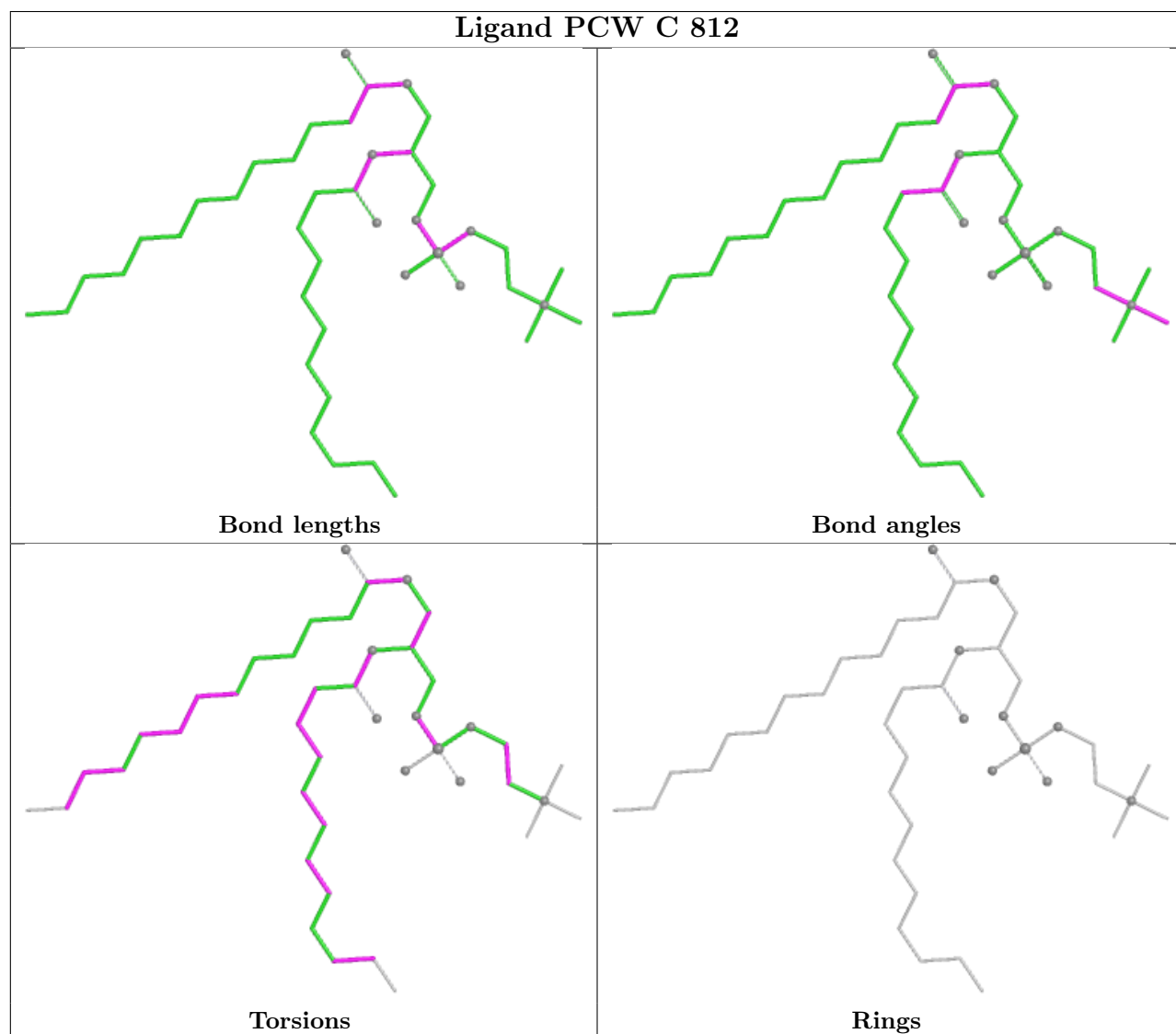
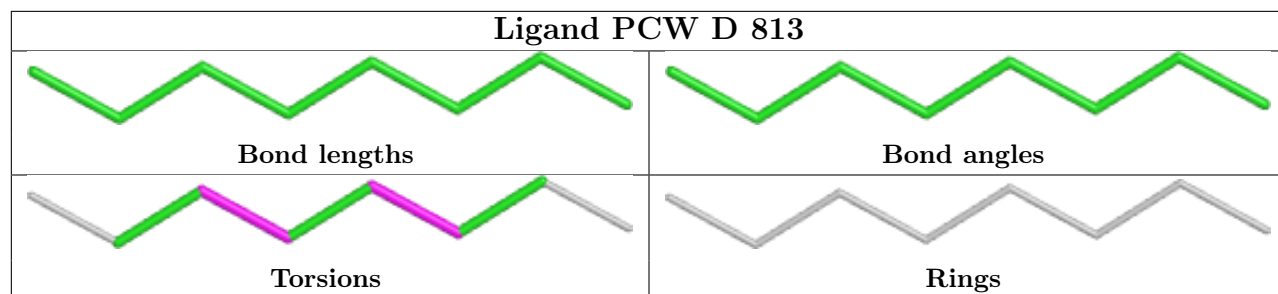


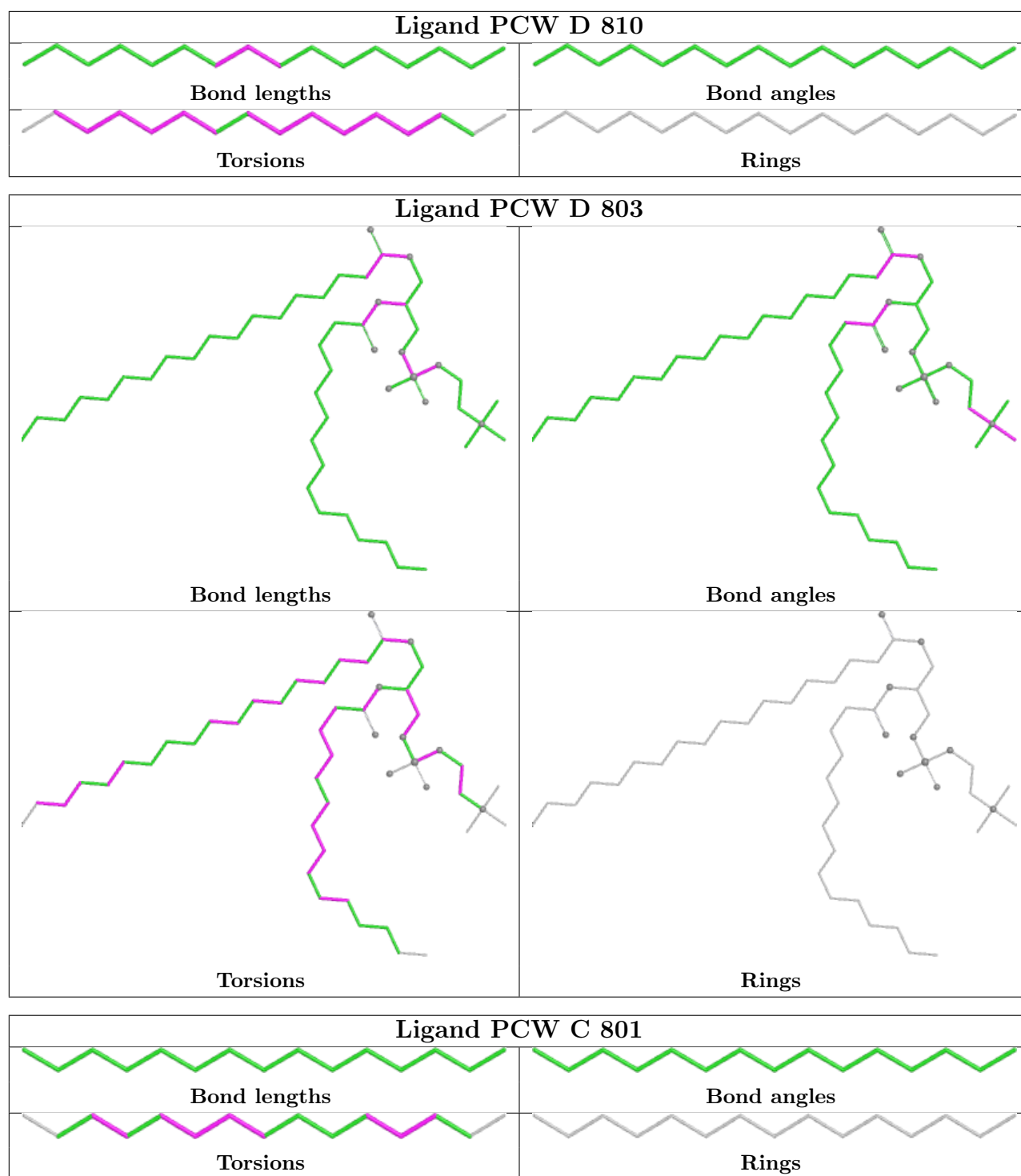












5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

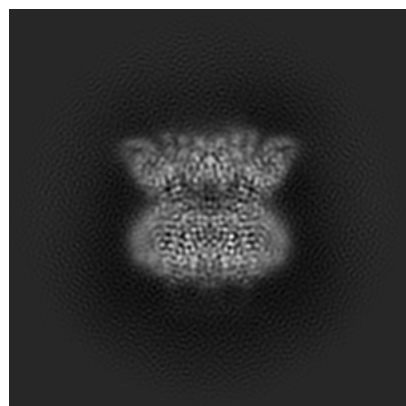
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-45934. These allow visual inspection of the internal detail of the map and identification of artifacts.

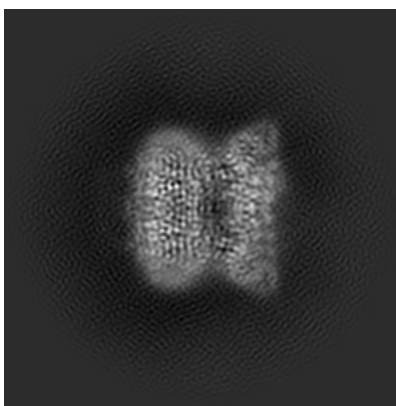
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

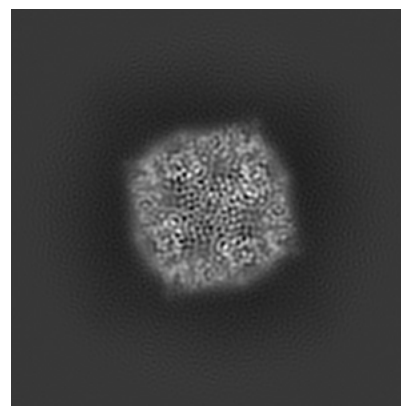
6.1.1 Primary map



X

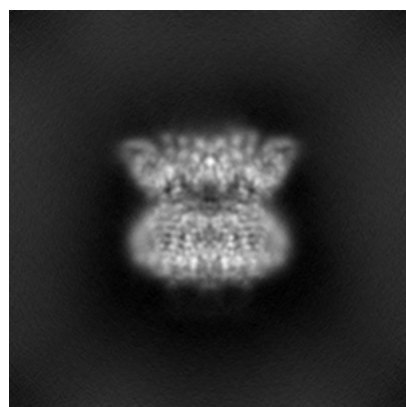


Y

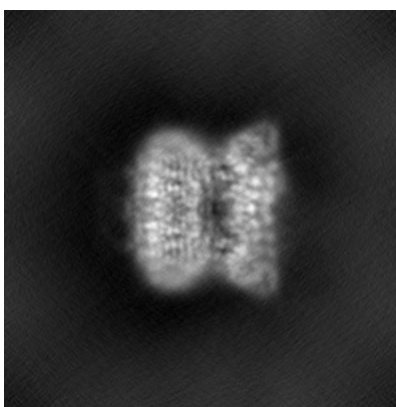


Z

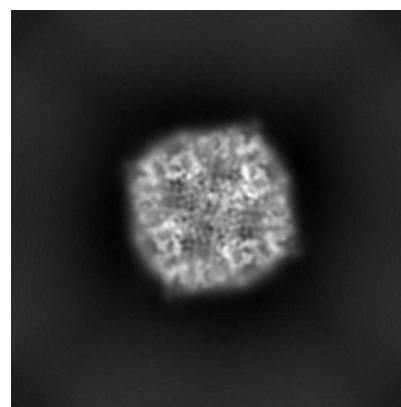
6.1.2 Raw map



X



Y

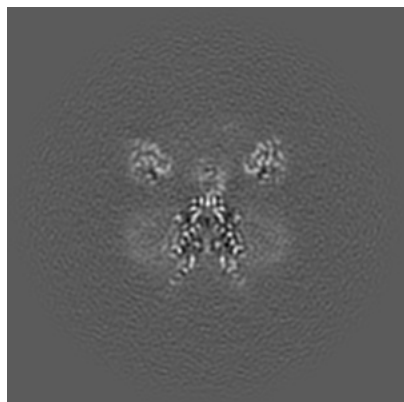


Z

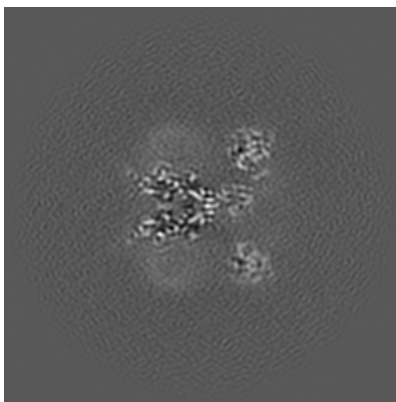
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

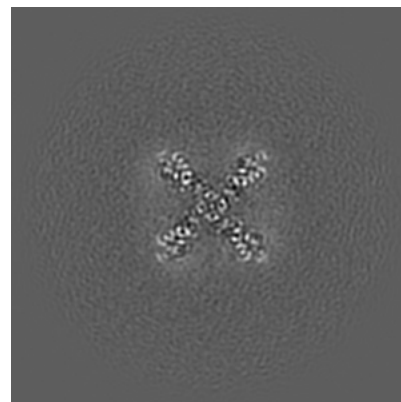
6.2.1 Primary map



X Index: 128

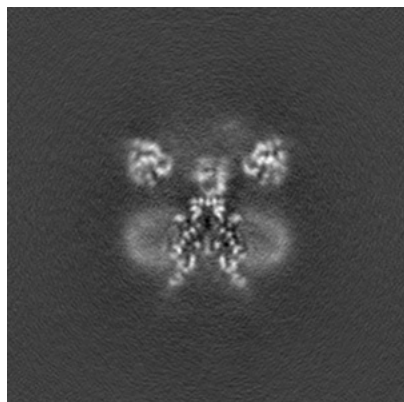


Y Index: 128

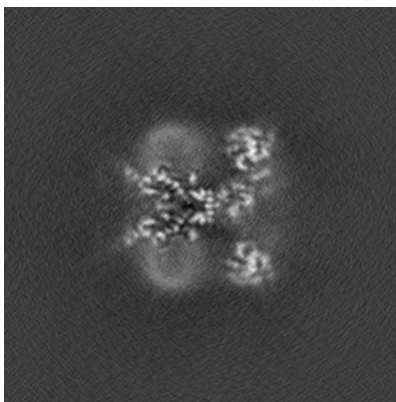


Z Index: 128

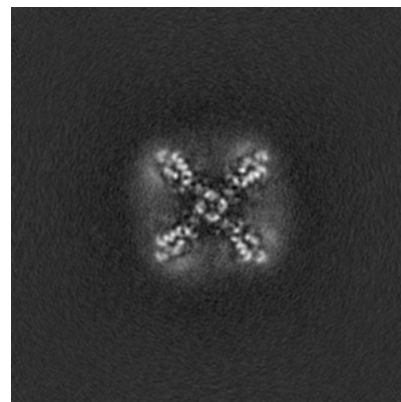
6.2.2 Raw map



X Index: 128



Y Index: 128

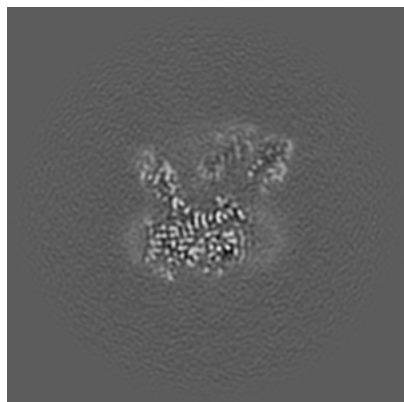


Z Index: 128

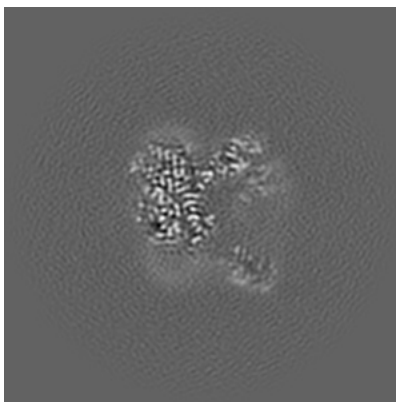
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

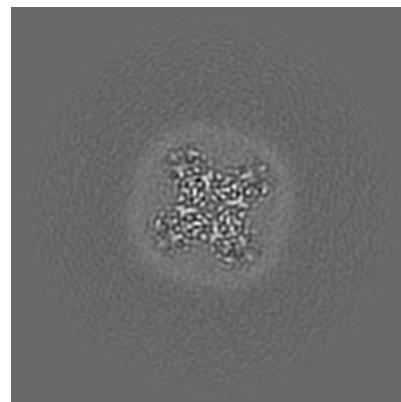
6.3.1 Primary map



X Index: 142

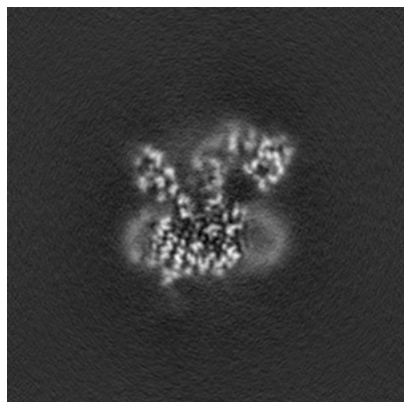


Y Index: 142

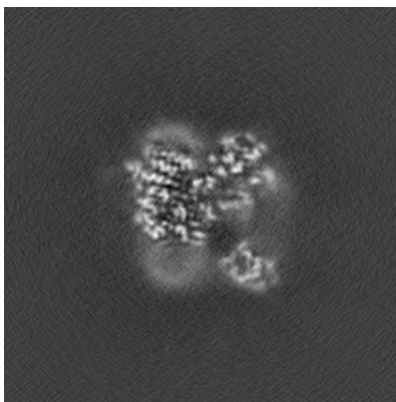


Z Index: 102

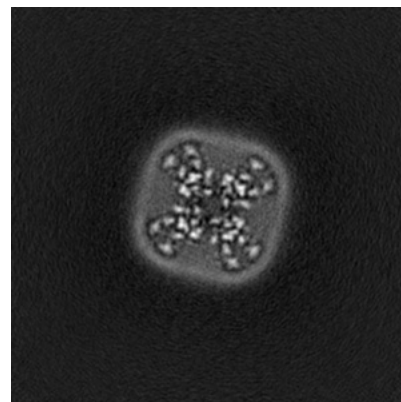
6.3.2 Raw map



X Index: 138



Y Index: 138

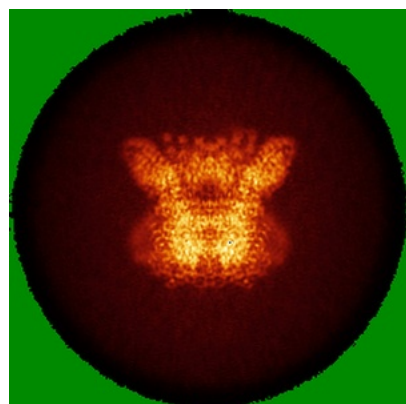


Z Index: 96

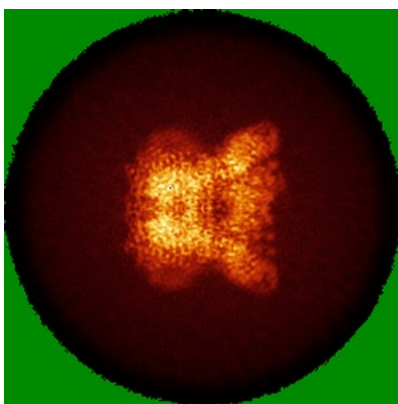
The images above show the largest variance slices of the map in three orthogonal directions.

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

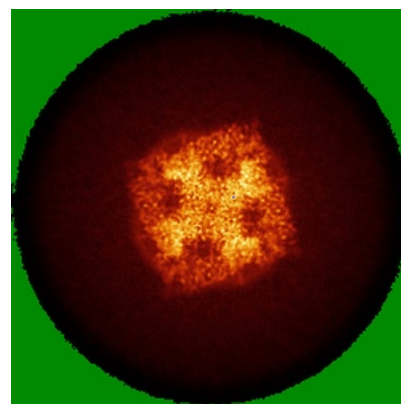
6.4.1 Primary map



X

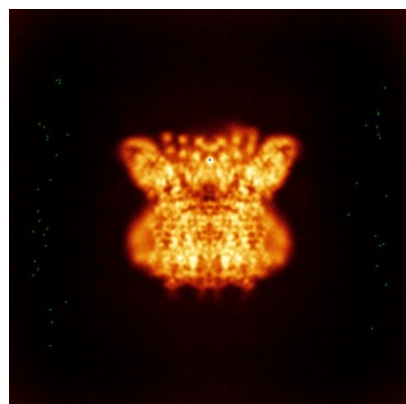


Y

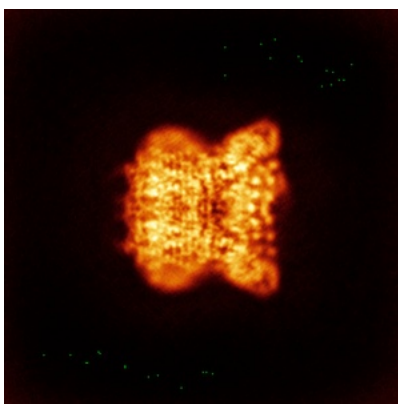


Z

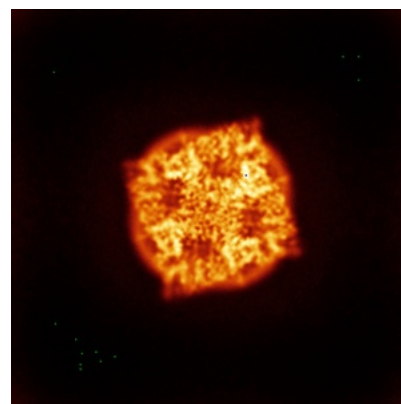
6.4.2 Raw map



X



Y

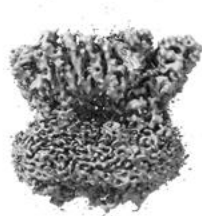


Z

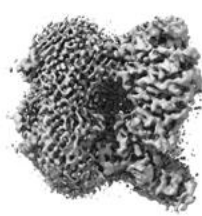
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

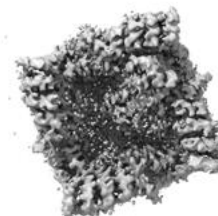
6.5.1 Primary map



X



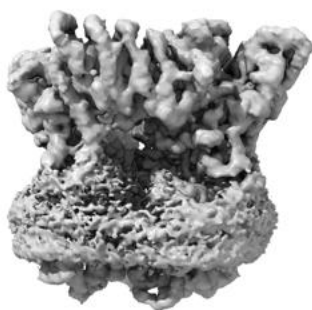
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.326. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

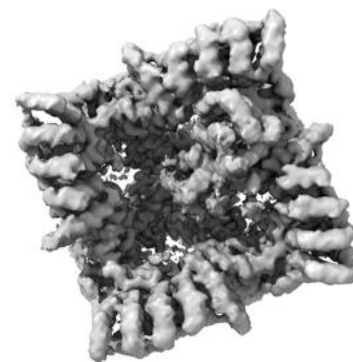
6.5.2 Raw map



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.

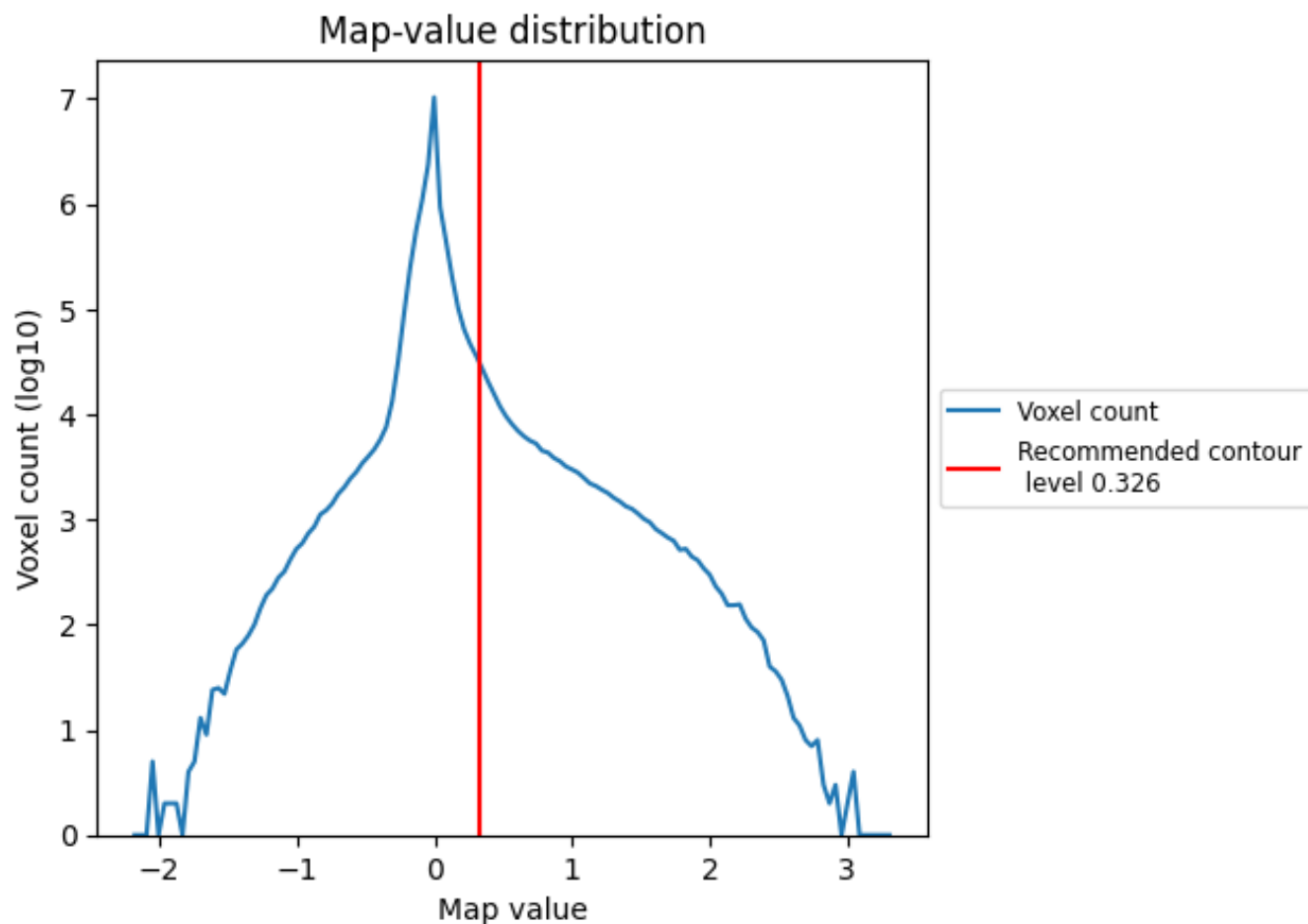
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

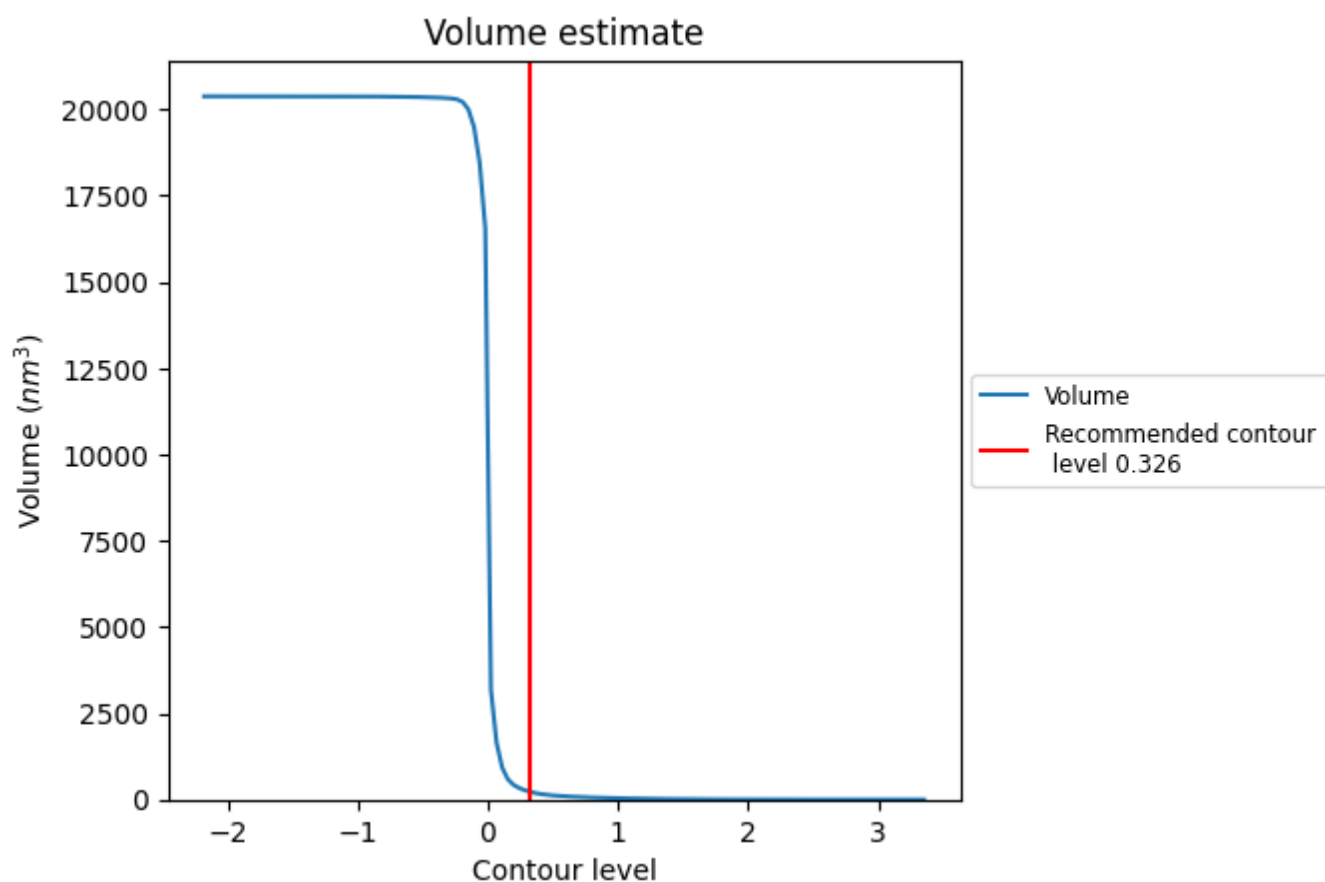
This section contains the results of statistical analysis of the map.

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.

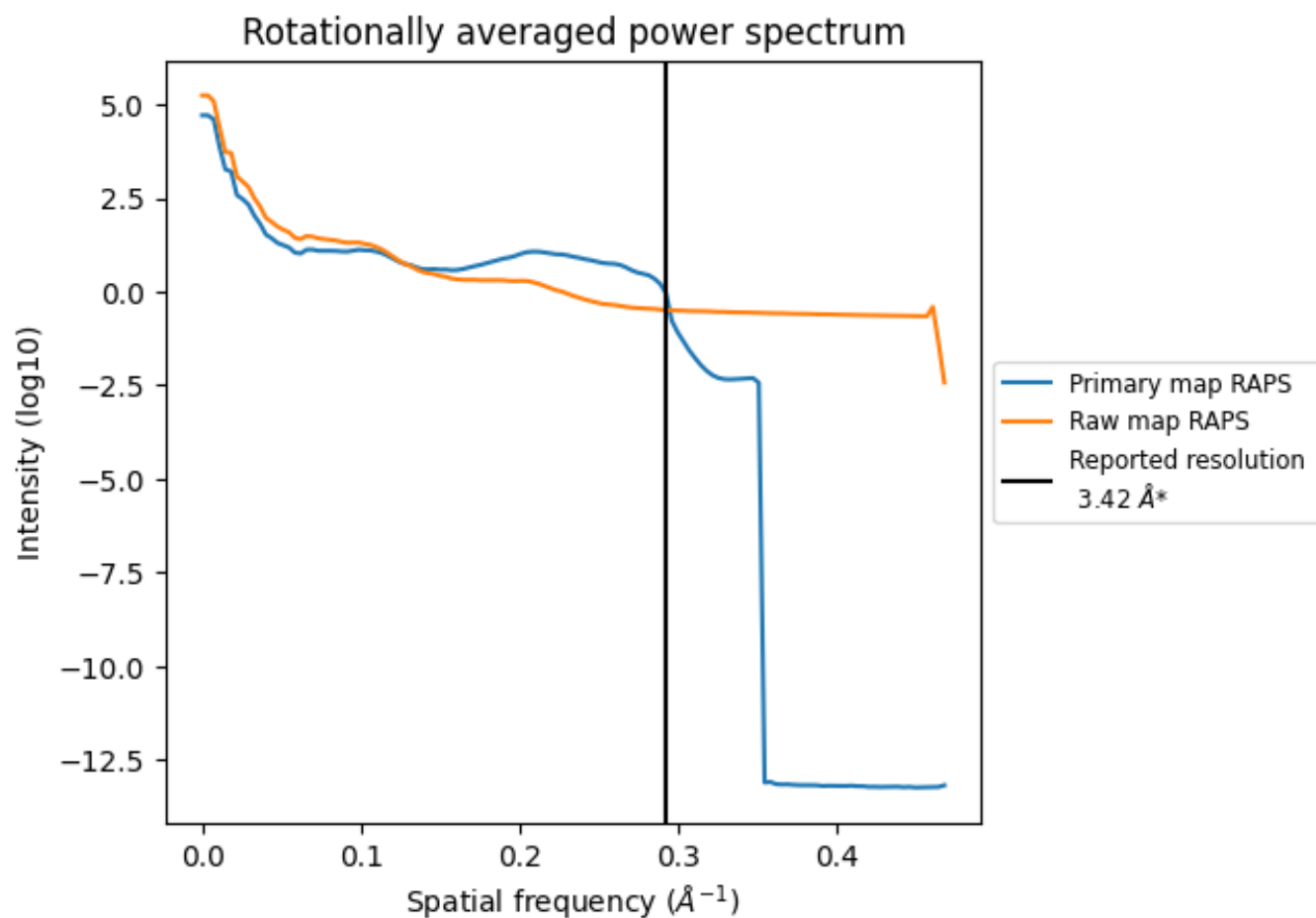
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 223 nm³; this corresponds to an approximate mass of 202 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 ⓘ

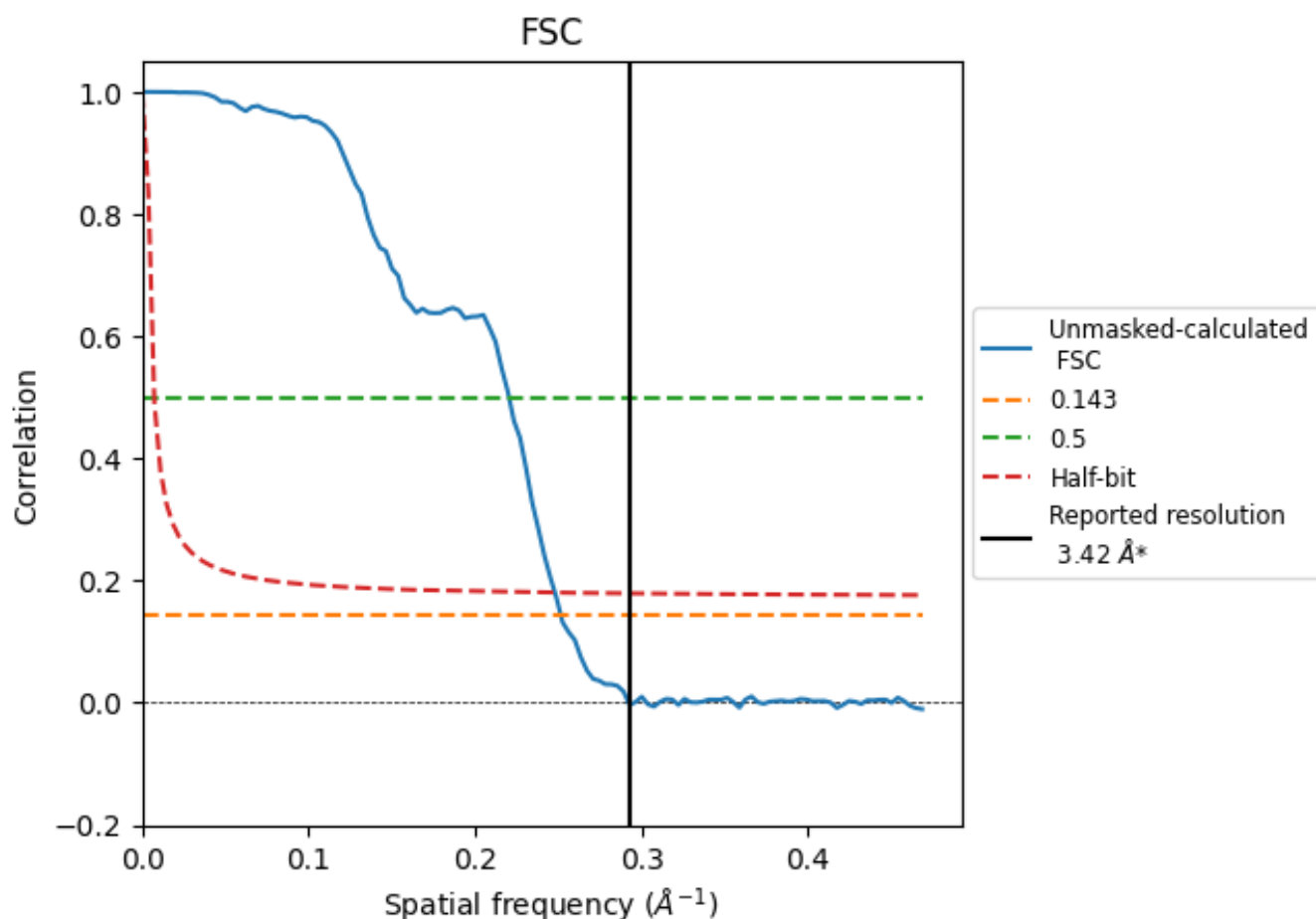


*Reported resolution corresponds to spatial frequency of 0.292 \AA^{-1}

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.292 Å⁻¹

8.2 Resolution estimates [i](#)

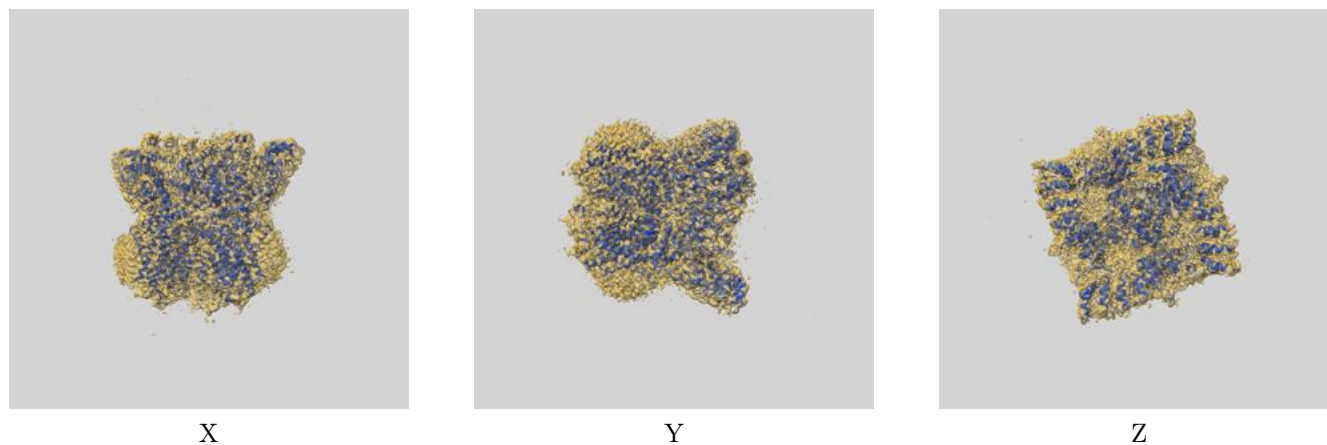
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.42	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.97	4.54	4.03

*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 3.97 differs from the reported value 3.42 by more than 10 %

9 Map-model fit [i](#)

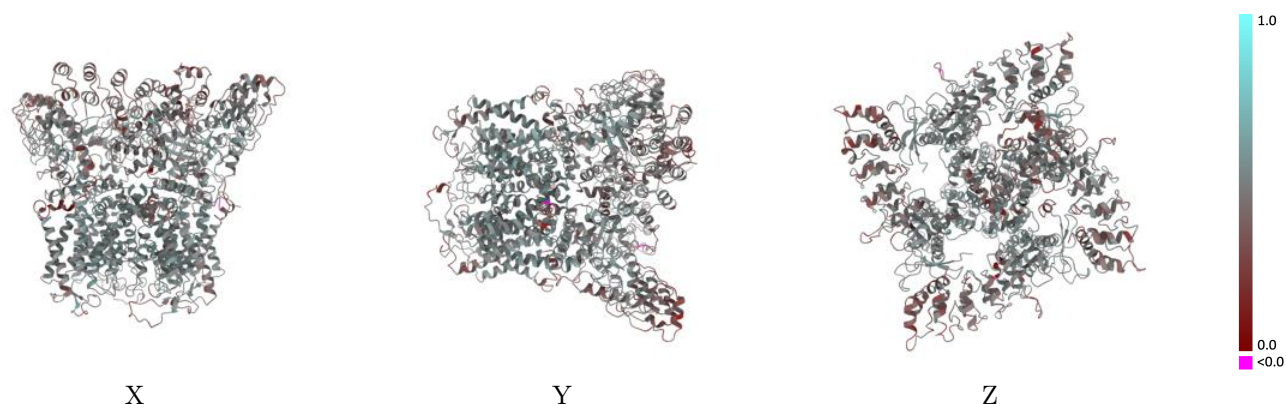
This section contains information regarding the fit between EMDB map EMD-45934 and PDB model 9CUI. Per-residue inclusion information can be found in section [3](#) on page [9](#).

9.1 Map-model overlay [i](#)



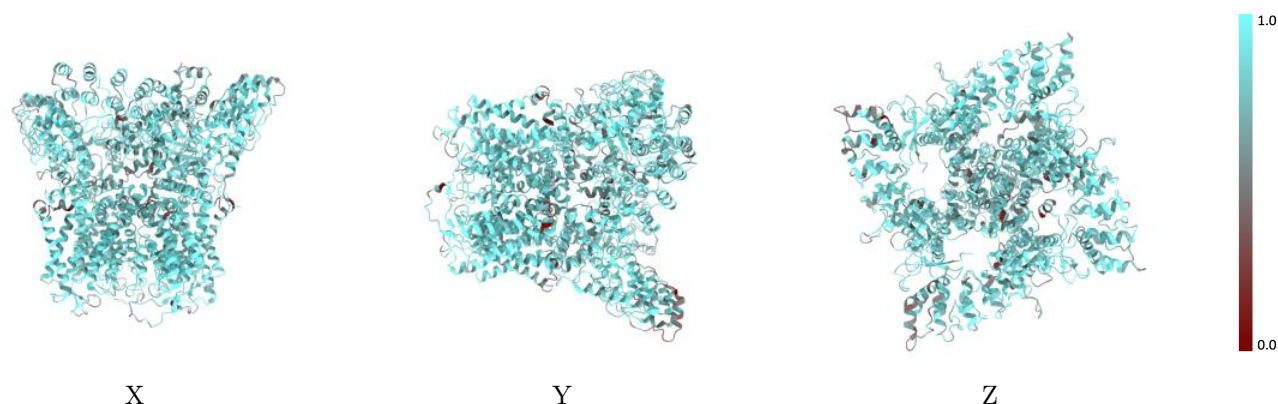
The images above show the 3D surface view of the map at the recommended contour level 0.326 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



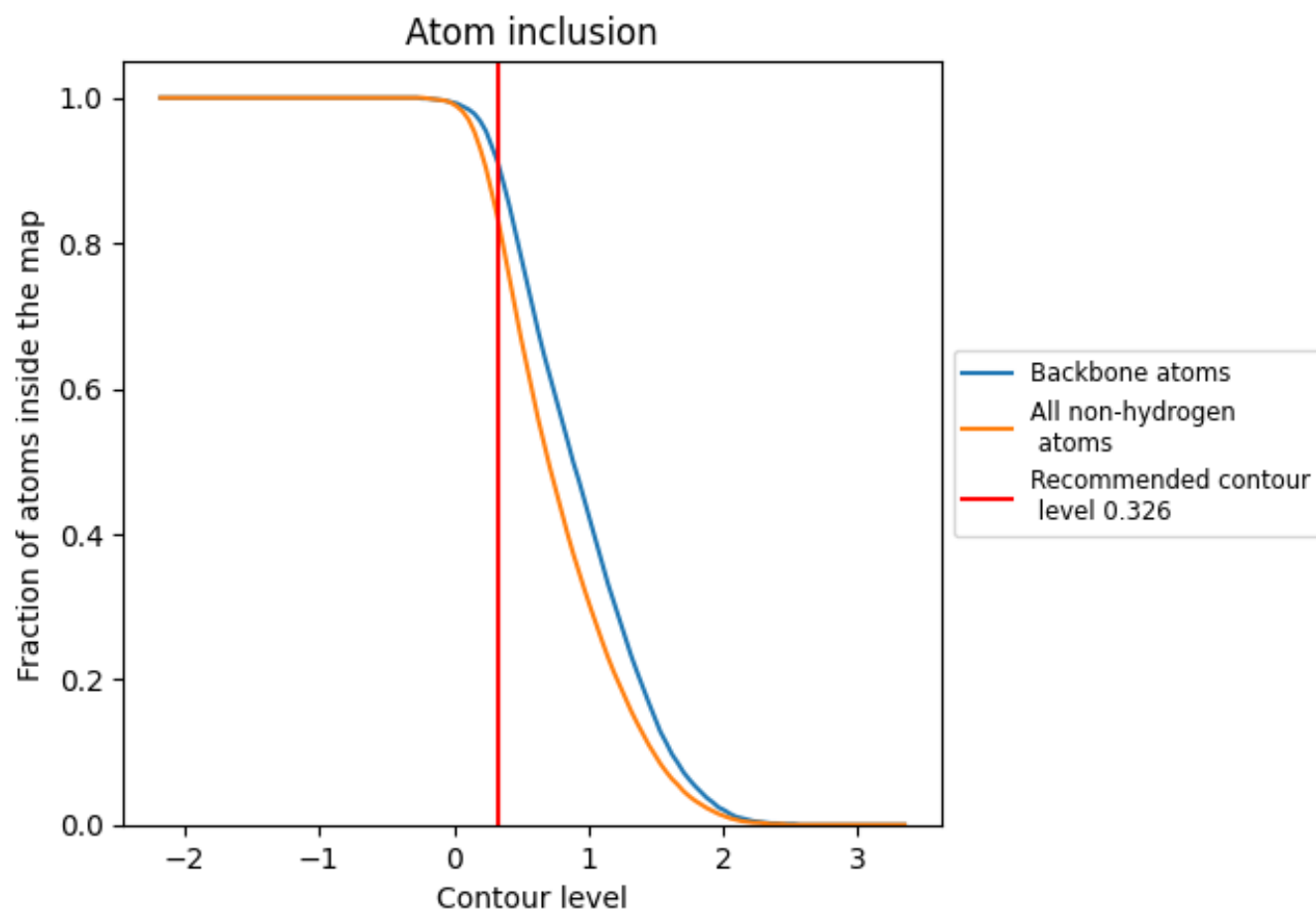
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.326).

9.4 Atom inclusion [i](#)



At the recommended contour level, 91% 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 (0.326) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8350	<div></div> 0.4680
A	<div></div> 0.8320	<div></div> 0.4680
B	<div></div> 0.8340	<div></div> 0.4690
C	<div></div> 0.8380	<div></div> 0.4680
D	<div></div> 0.8580	<div></div> 0.4820
E	<div></div> 0.7410	<div></div> 0.4070

