



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 11:13 AM UTC

PDB ID : 8DEU / pdb_00008deu
EMDB ID : EMD-27399
Title : Cryo-electron microscopy structure of Neisseria gonorrhoeae multidrug efflux pump MtrD with CASP peptide complex
Authors : Lyu, M.; Yu, E.W.
Deposited on : 2022-06-21
Resolution : 2.95 Å(reported)

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

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<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
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The types of validation reports are described at

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

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

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

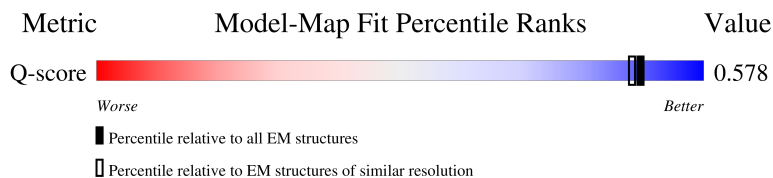
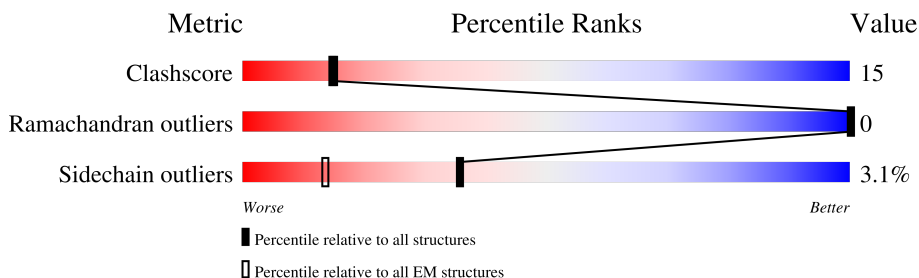
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.95 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	13114 (2.45 - 3.45)

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	1067	
1	B	1067	
1	C	1067	
2	D	9	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 29143 atoms, of which 5014 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 Efflux pump membrane transporter.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	1041	Total	C	H	N	O	S	0	0
			9455	5001	1658	1297	1454	45		
1	B	1040	Total	C	H	N	O	S	0	0
			9438	4995	1650	1295	1453	45		
1	C	1040	Total	C	H	N	O	S	0	0
			9442	4997	1651	1296	1453	45		

There are 75 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	738	VAL	ILE	conflict	UNP A0A6V7GUB3
A	752	SER	GLY	conflict	UNP A0A6V7GUB3
A	757	SER	ASN	conflict	UNP A0A6V7GUB3
A	774	GLY	ALA	conflict	UNP A0A6V7GUB3
A	775	ASP	SER	conflict	UNP A0A6V7GUB3
A	850	SER	GLY	conflict	UNP A0A6V7GUB3
A	871	ILE	LEU	conflict	UNP A0A6V7GUB3
A	872	ALA	ILE	conflict	UNP A0A6V7GUB3
A	875	ALA	GLY	conflict	UNP A0A6V7GUB3
A	878	ALA	VAL	conflict	UNP A0A6V7GUB3
A	879	VAL	ALA	conflict	UNP A0A6V7GUB3
A	899	LEU	ILE	conflict	UNP A0A6V7GUB3
A	902	MET	ILE	conflict	UNP A0A6V7GUB3
A	907	ALA	ILE	conflict	UNP A0A6V7GUB3
A	919	PHE	THR	conflict	UNP A0A6V7GUB3
A	921	GLY	LEU	conflict	UNP A0A6V7GUB3
A	922	LEU	MET	conflict	UNP A0A6V7GUB3
A	925	SER	GLY	conflict	UNP A0A6V7GUB3
A	926	VAL	ILE	conflict	UNP A0A6V7GUB3
A	928	SER	ALA	conflict	UNP A0A6V7GUB3
A	1052	ARG	LYS	conflict	UNP A0A6V7GUB3
A	1055	SER	ALA	conflict	UNP A0A6V7GUB3
A	1056	LYS	GLU	conflict	UNP A0A6V7GUB3
A	1065	LYS	SER	conflict	UNP A0A6V7GUB3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1067	TYR	HIS	conflict	UNP A0A6V7GUB3
B	738	VAL	ILE	conflict	UNP A0A6V7GUB3
B	752	SER	GLY	conflict	UNP A0A6V7GUB3
B	757	SER	ASN	conflict	UNP A0A6V7GUB3
B	774	GLY	ALA	conflict	UNP A0A6V7GUB3
B	775	ASP	SER	conflict	UNP A0A6V7GUB3
B	850	SER	GLY	conflict	UNP A0A6V7GUB3
B	871	ILE	LEU	conflict	UNP A0A6V7GUB3
B	872	ALA	ILE	conflict	UNP A0A6V7GUB3
B	875	ALA	GLY	conflict	UNP A0A6V7GUB3
B	878	ALA	VAL	conflict	UNP A0A6V7GUB3
B	879	VAL	ALA	conflict	UNP A0A6V7GUB3
B	899	LEU	ILE	conflict	UNP A0A6V7GUB3
B	902	MET	ILE	conflict	UNP A0A6V7GUB3
B	907	ALA	ILE	conflict	UNP A0A6V7GUB3
B	919	PHE	THR	conflict	UNP A0A6V7GUB3
B	921	GLY	LEU	conflict	UNP A0A6V7GUB3
B	922	LEU	MET	conflict	UNP A0A6V7GUB3
B	925	SER	GLY	conflict	UNP A0A6V7GUB3
B	926	VAL	ILE	conflict	UNP A0A6V7GUB3
B	928	SER	ALA	conflict	UNP A0A6V7GUB3
B	1052	ARG	LYS	conflict	UNP A0A6V7GUB3
B	1055	SER	ALA	conflict	UNP A0A6V7GUB3
B	1056	LYS	GLU	conflict	UNP A0A6V7GUB3
B	1065	LYS	SER	conflict	UNP A0A6V7GUB3
B	1067	TYR	HIS	conflict	UNP A0A6V7GUB3
C	738	VAL	ILE	conflict	UNP A0A6V7GUB3
C	752	SER	GLY	conflict	UNP A0A6V7GUB3
C	757	SER	ASN	conflict	UNP A0A6V7GUB3
C	774	GLY	ALA	conflict	UNP A0A6V7GUB3
C	775	ASP	SER	conflict	UNP A0A6V7GUB3
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C	872	ALA	ILE	conflict	UNP A0A6V7GUB3
C	875	ALA	GLY	conflict	UNP A0A6V7GUB3
C	878	ALA	VAL	conflict	UNP A0A6V7GUB3
C	879	VAL	ALA	conflict	UNP A0A6V7GUB3
C	899	LEU	ILE	conflict	UNP A0A6V7GUB3
C	902	MET	ILE	conflict	UNP A0A6V7GUB3
C	907	ALA	ILE	conflict	UNP A0A6V7GUB3
C	919	PHE	THR	conflict	UNP A0A6V7GUB3
C	921	GLY	LEU	conflict	UNP A0A6V7GUB3

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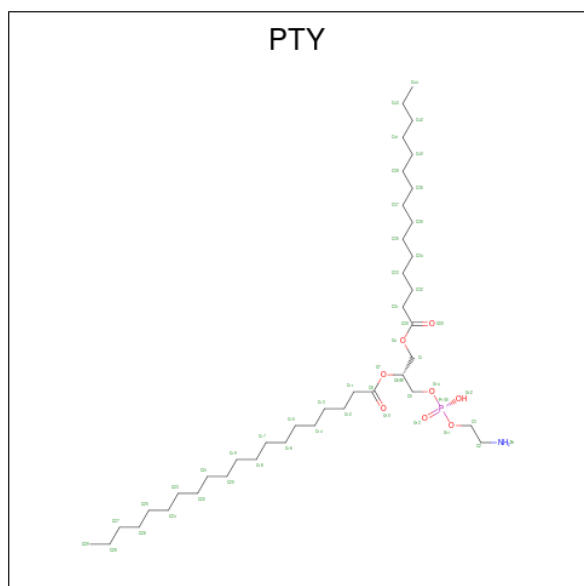
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Chain	Residue	Modelled	Actual	Comment	Reference
C	922	LEU	MET	conflict	UNP A0A6V7GUB3
C	925	SER	GLY	conflict	UNP A0A6V7GUB3
C	926	VAL	ILE	conflict	UNP A0A6V7GUB3
C	928	SER	ALA	conflict	UNP A0A6V7GUB3
C	1052	ARG	LYS	conflict	UNP A0A6V7GUB3
C	1055	SER	ALA	conflict	UNP A0A6V7GUB3
C	1056	LYS	GLU	conflict	UNP A0A6V7GUB3
C	1065	LYS	SER	conflict	UNP A0A6V7GUB3
C	1067	TYR	HIS	conflict	UNP A0A6V7GUB3

- Molecule 2 is a protein called CASP peptide.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	9	Total	C	N	O	S	
			64	36	11	15	2	

- Molecule 3 is PHOSPHATIDYLETHANOLAMINE (CCD ID: PTY) (formula: C₄₀H₈₀NO₈P).



Mol	Chain	Residues	Atoms			AltConf
3	A	1	Total	C		0
			10	10		
3	A	1	Total	C		0
			10	10		
3	A	1	Total	C	O	0
			26	22	4	

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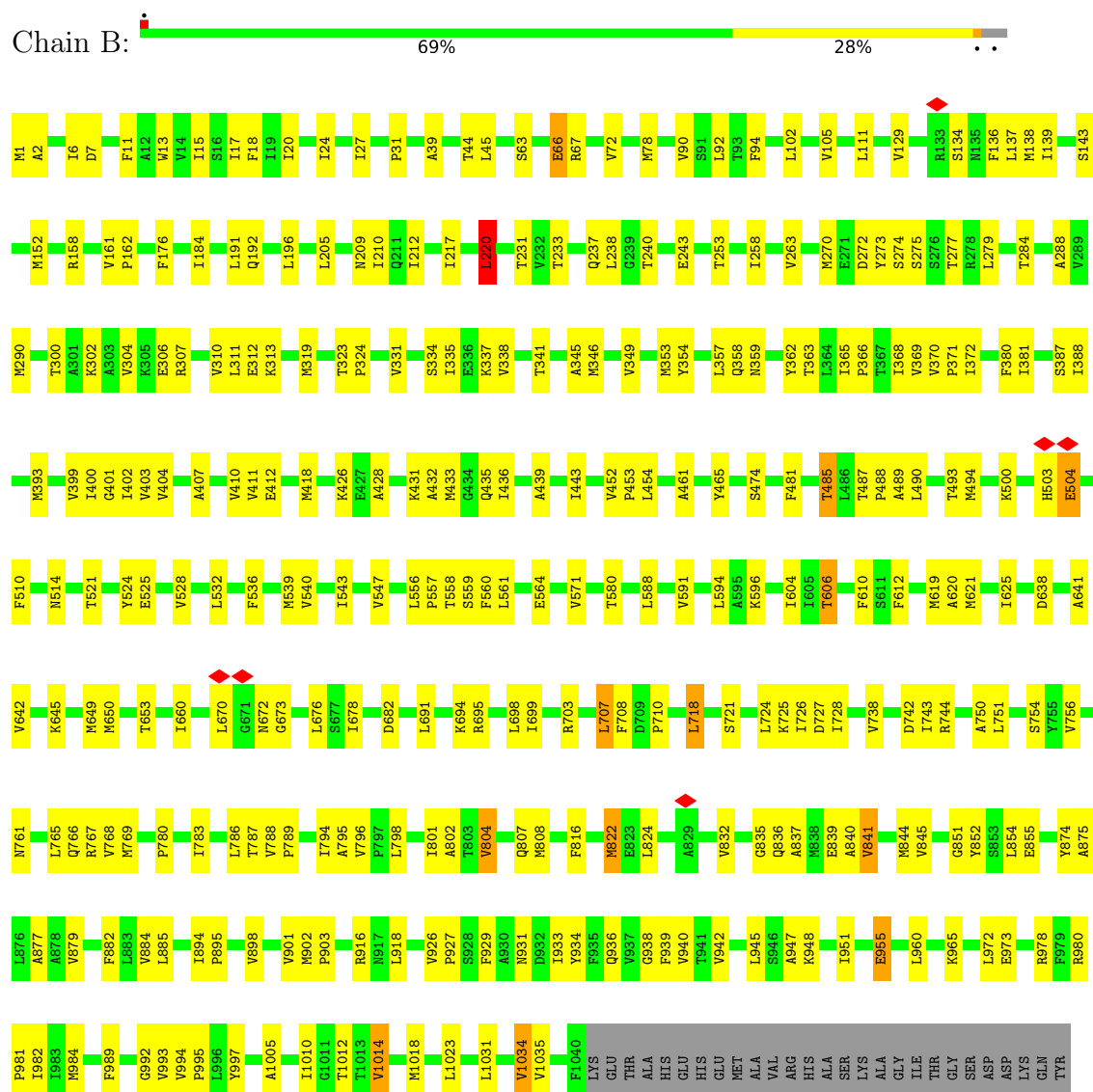
Mol	Chain	Residues	Atoms					AltConf
3	A	1	Total	C	O			0
			17	15	2			
3	A	1	Total	C	N	O	P	0
			30	20	1	8	1	
3	A	1	Total	C	O			0
			23	21	2			
3	A	1	Total	C	O			0
			31	27	4			
3	A	1	Total	C	O			0
			13	11	2			
3	A	1	Total	C	H	O	P	0
			94	30	55	8	1	
3	A	1	Total	C				0
			10	10				
3	A	1	Total	C				0
			28	28				
3	A	1	Total	C				0
			12	12				
3	B	1	Total	C	N	O	P	0
			37	27	1	8	1	
3	B	1	Total	C	O			0
			22	20	2			
3	B	1	Total	C	O			0
			38	35	3			
3	B	1	Total	C	N	O	P	0
			35	25	1	8	1	
3	B	1	Total	C	N	O	P	0
			44	34	1	8	1	
3	B	1	Total	C	O			0
			33	29	4			
3	C	1	Total	C	N	O	P	0
			45	35	1	8	1	
3	C	1	Total	C				0
			29	29				
3	C	1	Total	C	O			0
			30	26	4			
3	C	1	Total	C	N	O	P	0
			35	25	1	8	1	
3	C	1	Total	C				0
			26	26				
3	C	1	Total	C	O			0
			21	20	1			

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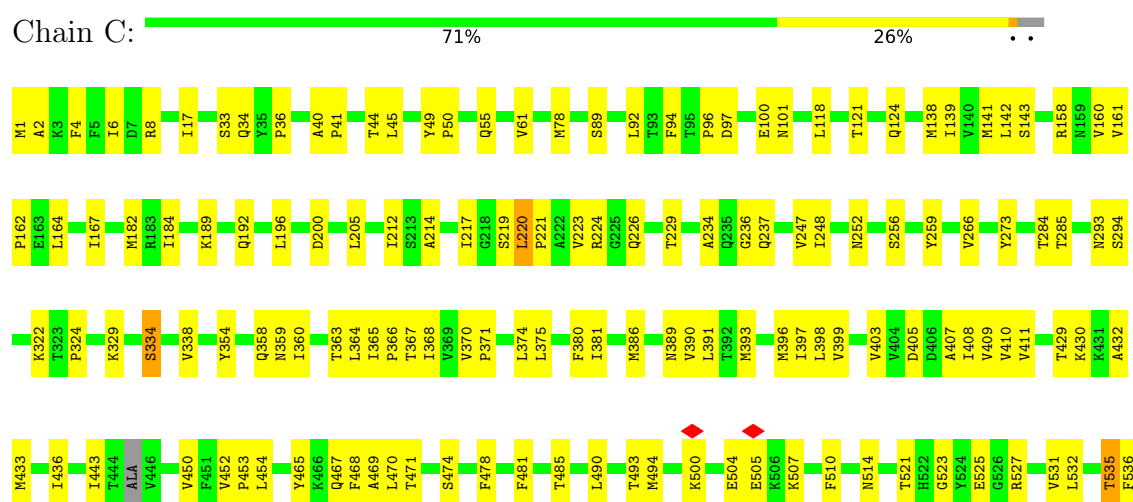
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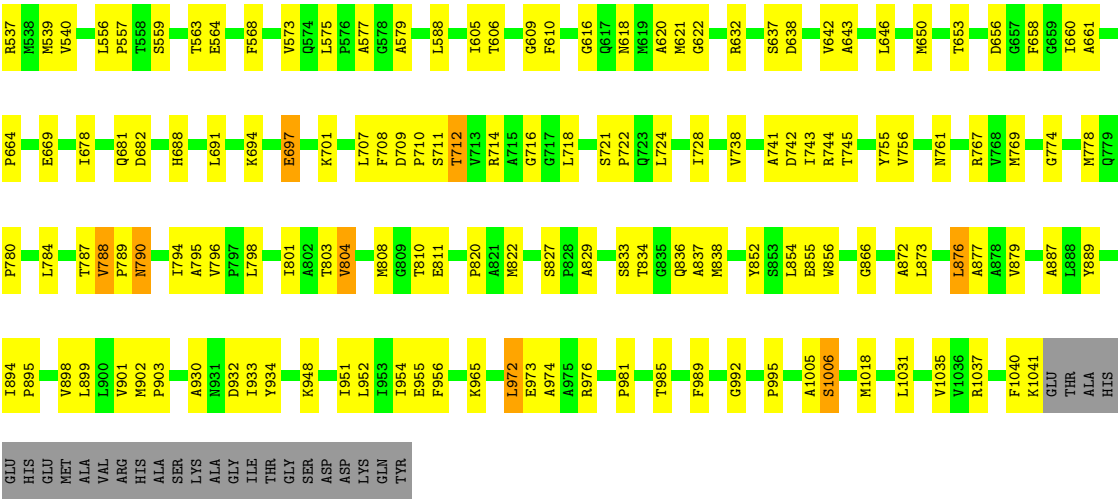
Mol	Chain	Residues	Atoms	AltConf
3	C	1	Total C 10 10	0
3	C	1	Total C O 15 13 2	0
3	C	1	Total C 20 20	0

• Molecule 1: Efflux pump membrane transporter



• Molecule 1: Efflux pump membrane transporter





● Molecule 2: CASP peptide



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	129244	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 ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	750	Depositor
Maximum defocus (nm)	1750	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.248	Depositor
Minimum map value	-1.339	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.071	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	324.0, 324.0, 324.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PTY

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	0.32	0/7939	0.54	11/10770 (0.1%)
1	B	0.27	0/7930	0.44	4/10759 (0.0%)
1	C	0.30	0/7931	0.54	13/10757 (0.1%)
2	D	0.50	0/64	1.08	0/85
All	All	0.30	0/23864	0.51	28/32371 (0.1%)

There are no bond length outliers.

The worst 5 of 28 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	756	VAL	N-CA-C	10.79	121.62	110.72
1	A	143	SER	N-CA-C	9.24	124.19	111.17
1	A	61	VAL	N-CA-C	9.15	119.43	111.56
1	A	756	VAL	N-CA-C	8.51	119.76	111.67
1	A	712	THR	N-CA-C	7.74	119.80	111.36

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	7797	1658	7973	261	0
1	B	7788	1650	7960	232	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	7791	1651	7964	220	0
2	D	64	0	53	12	0
3	A	249	55	365	34	0
3	B	209	0	300	28	0
3	C	231	0	352	21	0
All	All	24129	5014	24967	736	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 15.

The worst 5 of 736 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:556:LEU:HD12	1:B:557:PRO:HD2	1.31	1.13
1:B:290:MET:HE3	2:D:126:LYS:HE3	1.36	1.07
1:C:478:PHE:CD1	1:C:478:PHE:CZ	2.38	1.03
1:B:212:ILE:HG12	1:C:744:ARG:HG3	1.38	1.01
1:C:577:ALA:HB1	1:C:721:SER:HB3	1.45	0.96

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	1039/1067 (97%)	987 (95%)	52 (5%)	0	100	100
1	B	1038/1067 (97%)	997 (96%)	41 (4%)	0	100	100
1	C	1036/1067 (97%)	986 (95%)	50 (5%)	0	100	100
2	D	7/9 (78%)	4 (57%)	3 (43%)	0	100	100
All	All	3120/3210 (97%)	2974 (95%)	146 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	833/853 (98%)	806 (97%)	27 (3%)	34	59
1	B	832/853 (98%)	804 (97%)	28 (3%)	32	58
1	C	833/853 (98%)	811 (97%)	22 (3%)	40	65
2	D	8/8 (100%)	8 (100%)	0	100	100
All	All	2506/2567 (98%)	2429 (97%)	77 (3%)	36	59

5 of 77 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	142	LEU
1	C	803	THR
1	C	220	LEU
1	C	535	THR
1	C	972	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	104	GLN
1	C	235	GLN
1	C	963	GLN
1	C	165	GLN
1	C	413	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

27 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	PTY	C	1108	-	14,14,49	1.03	1 (7%)	14,14,54	1.21	0
3	PTY	B	1106	-	32,32,49	0.86	2 (6%)	33,33,54	1.18	2 (6%)
3	PTY	A	1108	-	12,12,49	1.05	2 (16%)	12,12,54	1.17	1 (8%)
3	PTY	A	1112	-	11,11,49	0.28	0	10,10,54	0.86	0
3	PTY	B	1101	-	36,36,49	1.02	4 (11%)	39,41,54	1.16	2 (5%)
3	PTY	A	1102	-	9,9,49	0.28	0	8,8,54	0.85	0
3	PTY	B	1104	-	34,34,49	1.05	4 (11%)	37,39,54	1.20	2 (5%)
3	PTY	C	1105	-	23,23,49	0.30	0	20,20,54	0.78	0
3	PTY	A	1107	-	29,29,49	0.91	3 (10%)	29,29,54	1.16	2 (6%)
3	PTY	A	1106	-	21,21,49	0.75	1 (4%)	20,20,54	1.00	0
3	PTY	C	1103	-	28,28,49	0.87	2 (7%)	28,28,54	1.23	1 (3%)
3	PTY	A	1104	-	16,16,49	0.83	1 (6%)	16,16,54	1.15	1 (6%)
3	PTY	A	1111	-	26,26,49	0.29	0	24,24,54	0.86	0
3	PTY	C	1102	-	27,27,49	0.33	0	25,25,54	0.86	0
3	PTY	C	1106	-	19,19,49	1.08	1 (5%)	17,17,54	0.92	0
3	PTY	A	1109	-	38,38,49	1.05	5 (13%)	41,43,54	1.03	2 (4%)
3	PTY	C	1104	-	34,34,49	1.06	2 (5%)	37,39,54	1.17	3 (8%)
3	PTY	B	1102	-	20,20,49	0.84	2 (10%)	19,19,54	1.05	1 (5%)
3	PTY	C	1101	-	44,44,49	0.93	4 (9%)	47,49,54	1.15	2 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PTY	B	1103	-	36,36,49	0.97	2 (5%)	35,35,54	0.99	0
3	PTY	A	1103	-	24,24,49	0.95	2 (8%)	24,24,54	1.21	2 (8%)
3	PTY	A	1110	-	9,9,49	0.29	0	8,8,54	0.83	0
3	PTY	B	1105	-	43,43,49	0.93	4 (9%)	46,48,54	1.16	2 (4%)
3	PTY	C	1109	-	19,19,49	0.28	0	18,18,54	0.91	0
3	PTY	C	1107	-	9,9,49	0.28	0	8,8,54	0.87	0
3	PTY	A	1105	-	29,29,49	1.13	4 (13%)	32,34,54	1.19	2 (6%)
3	PTY	A	1101	-	9,9,49	0.28	0	8,8,54	0.85	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
3	PTY	C	1108	-	-	8/12/12/53	-
3	PTY	B	1106	-	-	10/32/32/53	-
3	PTY	A	1108	-	-	0/11/11/53	-
3	PTY	A	1112	-	-	3/9/9/53	-
3	PTY	B	1101	-	-	20/40/40/53	-
3	PTY	A	1102	-	-	0/7/7/53	-
3	PTY	B	1104	-	-	21/38/38/53	-
3	PTY	C	1105	-	-	4/17/17/53	-
3	PTY	A	1107	-	-	8/26/26/53	-
3	PTY	A	1106	-	-	5/17/17/53	-
3	PTY	C	1103	-	-	5/24/24/53	-
3	PTY	A	1104	-	-	4/14/14/53	-
3	PTY	A	1111	-	-	3/22/22/53	-
3	PTY	C	1102	-	-	15/23/23/53	-
3	PTY	C	1106	-	-	5/15/15/53	-
3	PTY	A	1109	-	-	13/40/40/53	-
3	PTY	C	1104	-	-	17/38/38/53	-
3	PTY	B	1102	-	-	4/17/17/53	-
3	PTY	C	1101	-	-	14/48/48/53	-
3	PTY	B	1103	-	-	13/32/32/53	-
3	PTY	A	1103	-	-	6/20/20/53	-
3	PTY	A	1110	-	-	2/7/7/53	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PTY	B	1105	-	-	20/47/47/53	-
3	PTY	C	1109	-	-	6/17/17/53	-
3	PTY	C	1107	-	-	0/7/7/53	-
3	PTY	A	1105	-	-	13/33/33/53	-
3	PTY	A	1101	-	-	2/7/7/53	-

The worst 5 of 46 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1106	PTY	O7-C8	-4.57	1.18	1.42
3	B	1103	PTY	O4-C30	-4.56	1.18	1.42
3	C	1104	PTY	O4-C30	4.04	1.45	1.33
3	C	1104	PTY	O7-C8	3.89	1.45	1.34
3	C	1108	PTY	O7-C8	3.20	1.41	1.30

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1105	PTY	O7-C8-C11	4.22	120.61	111.48
3	B	1105	PTY	O7-C8-C11	4.09	120.33	111.48
3	C	1101	PTY	O7-C8-C11	3.86	119.83	111.48
3	C	1104	PTY	O7-C8-C11	3.81	119.73	111.48
3	B	1104	PTY	O7-C8-C11	3.78	119.65	111.48

There are no chirality outliers.

5 of 221 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1105	PTY	N1-C2-C3-O11
3	A	1105	PTY	C5-O14-P1-O13
3	B	1101	PTY	C3-O11-P1-O13
3	B	1101	PTY	C3-O11-P1-O14
3	B	1101	PTY	C5-O14-P1-O11

There are no ring outliers.

24 monomers are involved in 80 short contacts:

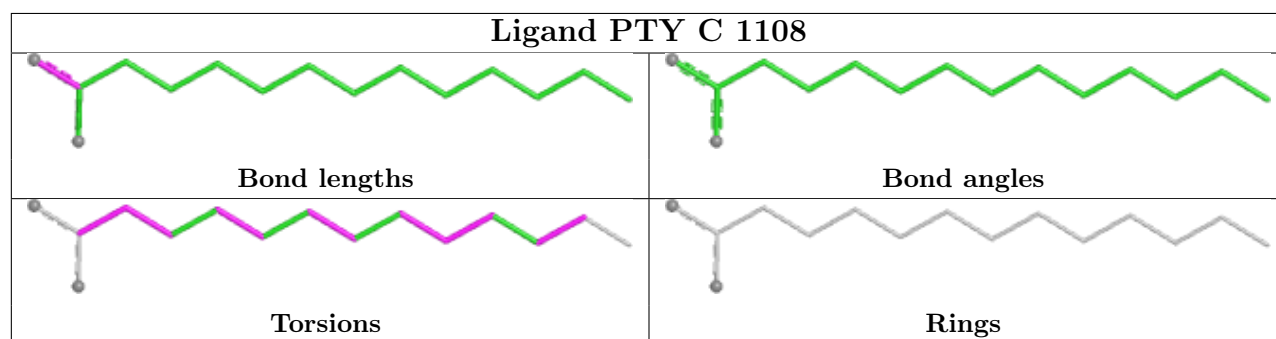
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1108	PTY	3	0
3	B	1106	PTY	7	0

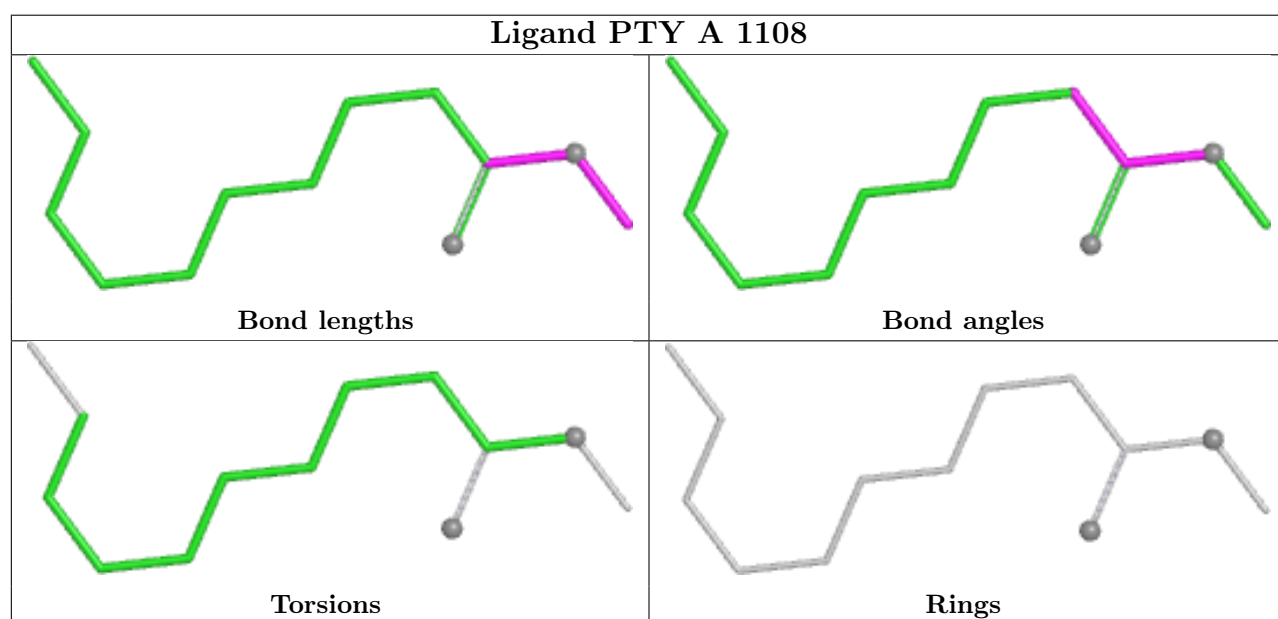
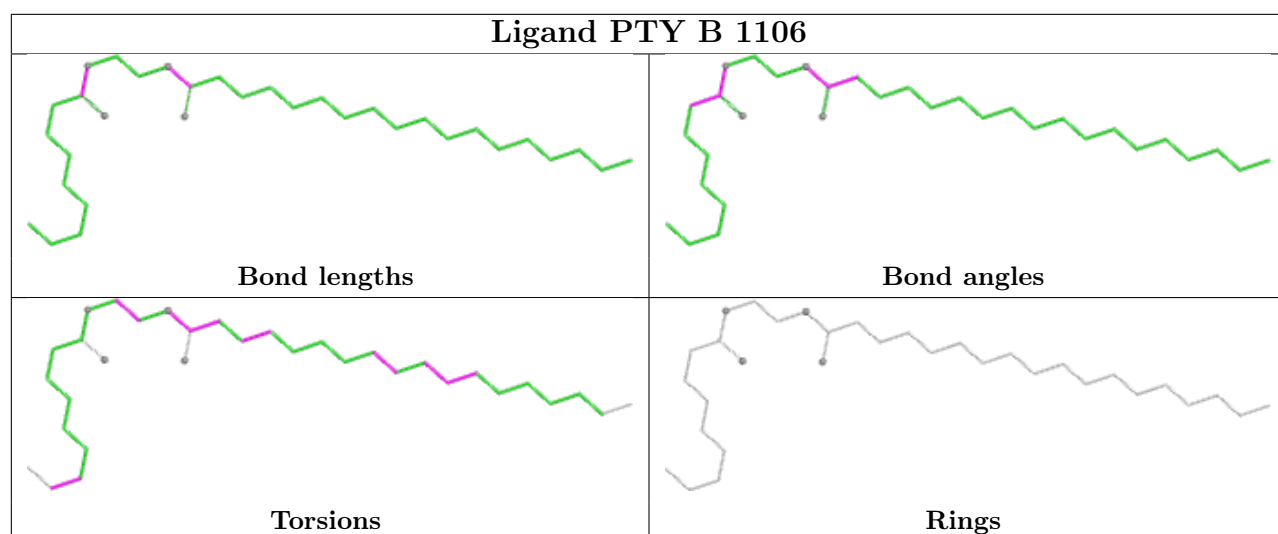
Continued on next page...

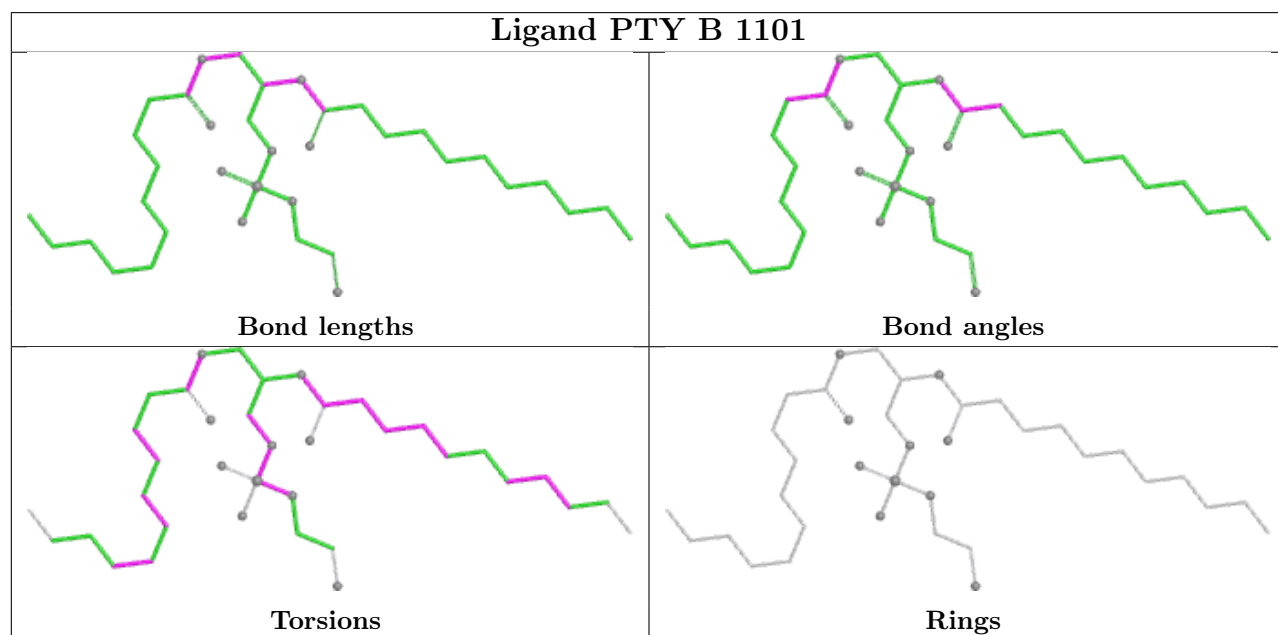
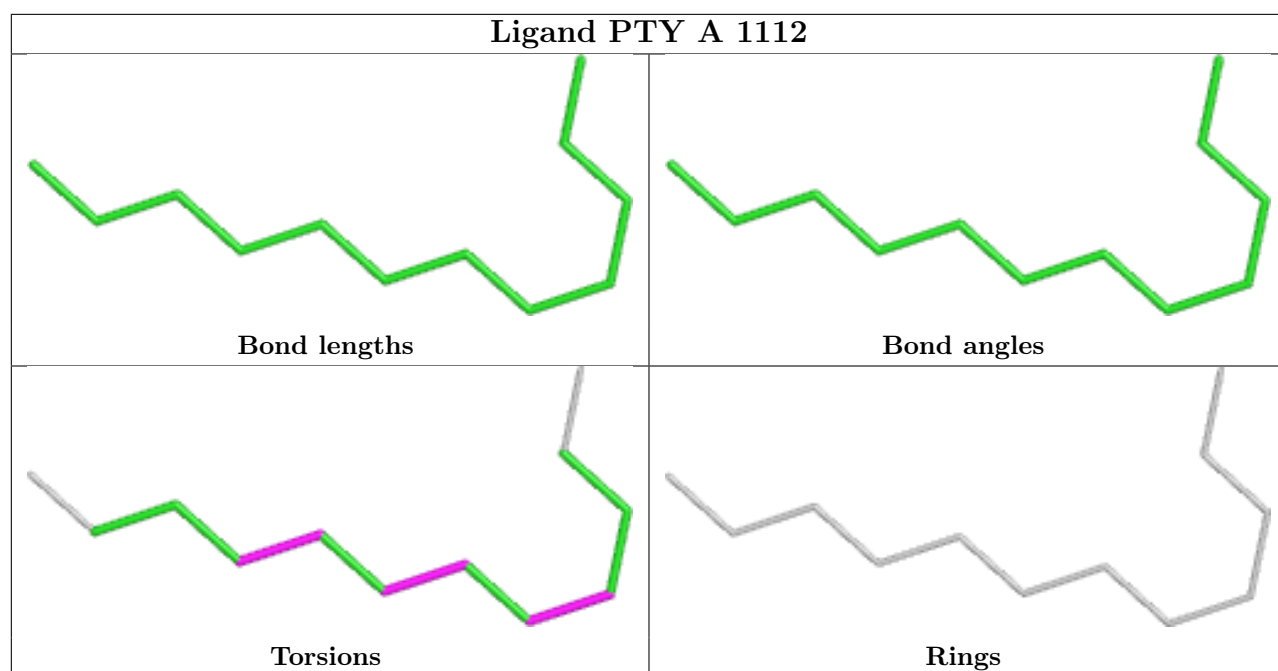
Continued from previous page...

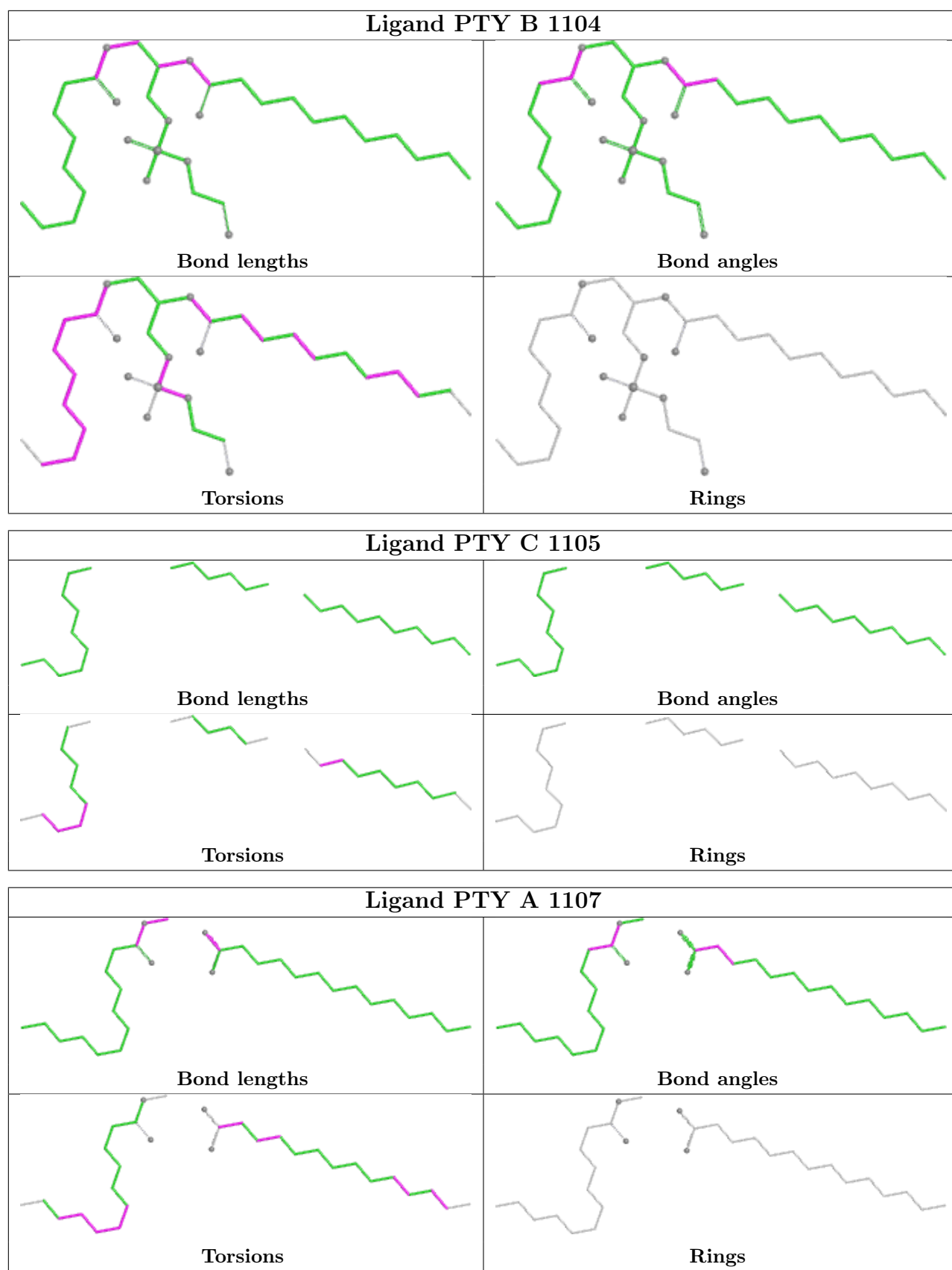
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1108	PTY	2	0
3	A	1112	PTY	1	0
3	B	1101	PTY	7	0
3	A	1102	PTY	2	0
3	B	1104	PTY	3	0
3	C	1105	PTY	3	0
3	A	1107	PTY	4	0
3	A	1106	PTY	5	0
3	C	1103	PTY	8	0
3	A	1104	PTY	8	0
3	A	1111	PTY	4	0
3	C	1102	PTY	1	0
3	C	1106	PTY	3	0
3	A	1109	PTY	2	0
3	C	1104	PTY	5	0
3	B	1102	PTY	1	0
3	C	1101	PTY	1	0
3	B	1103	PTY	13	0
3	A	1103	PTY	6	0
3	C	1109	PTY	3	0
3	A	1105	PTY	2	0
3	A	1101	PTY	4	0

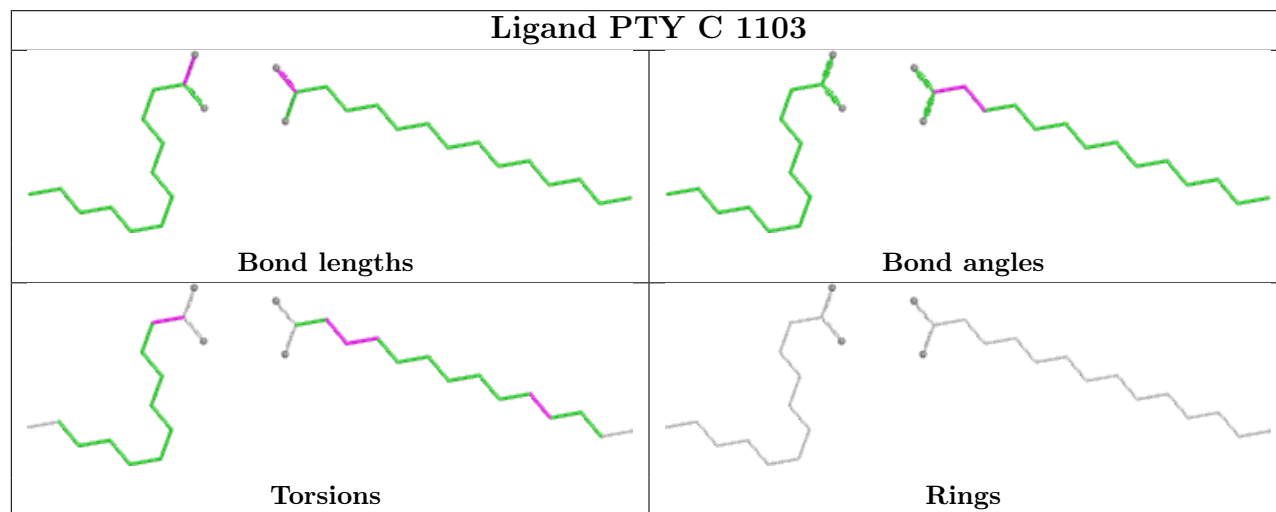
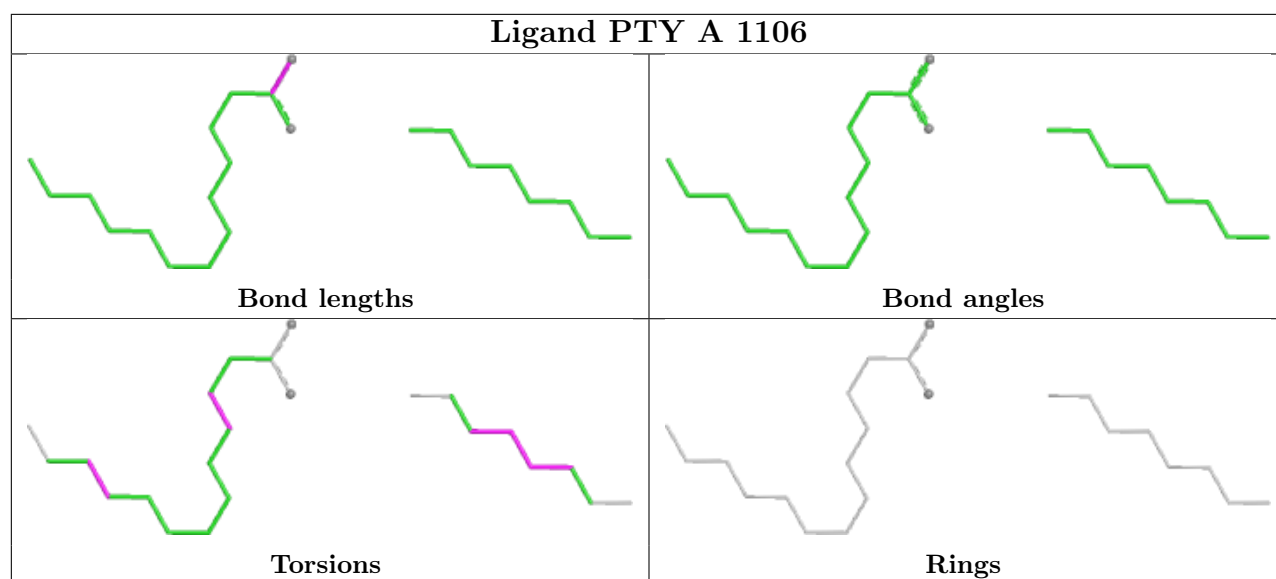
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

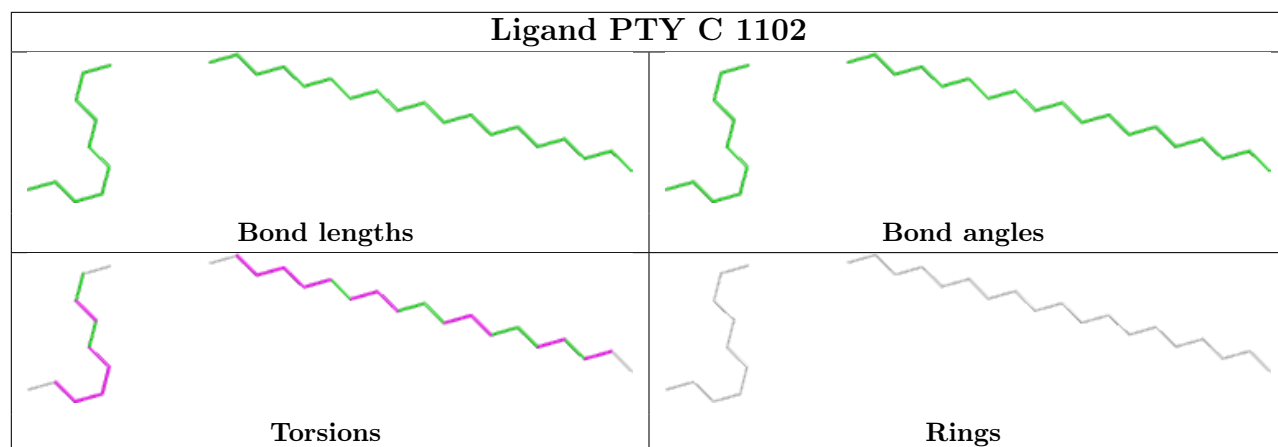
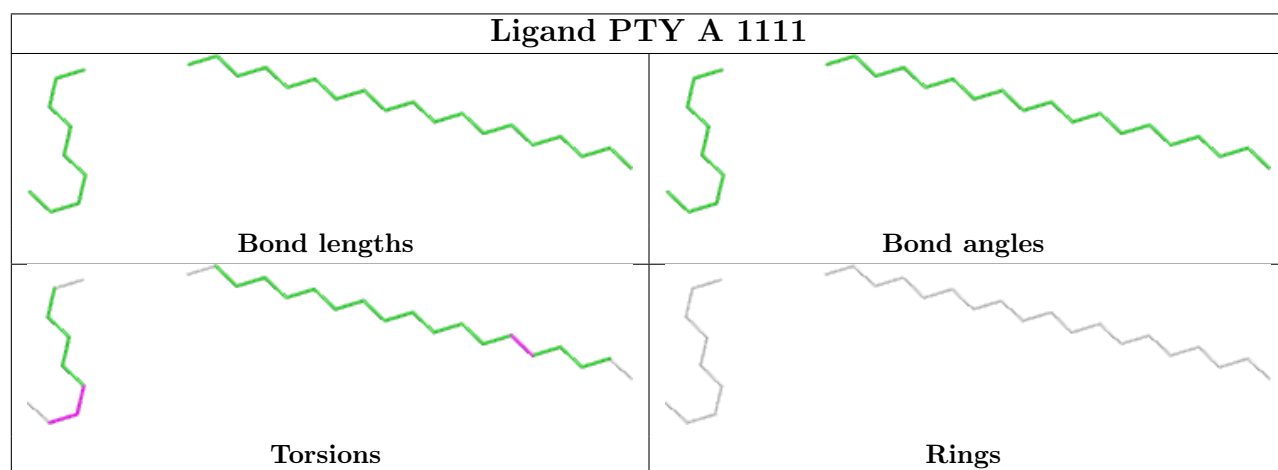
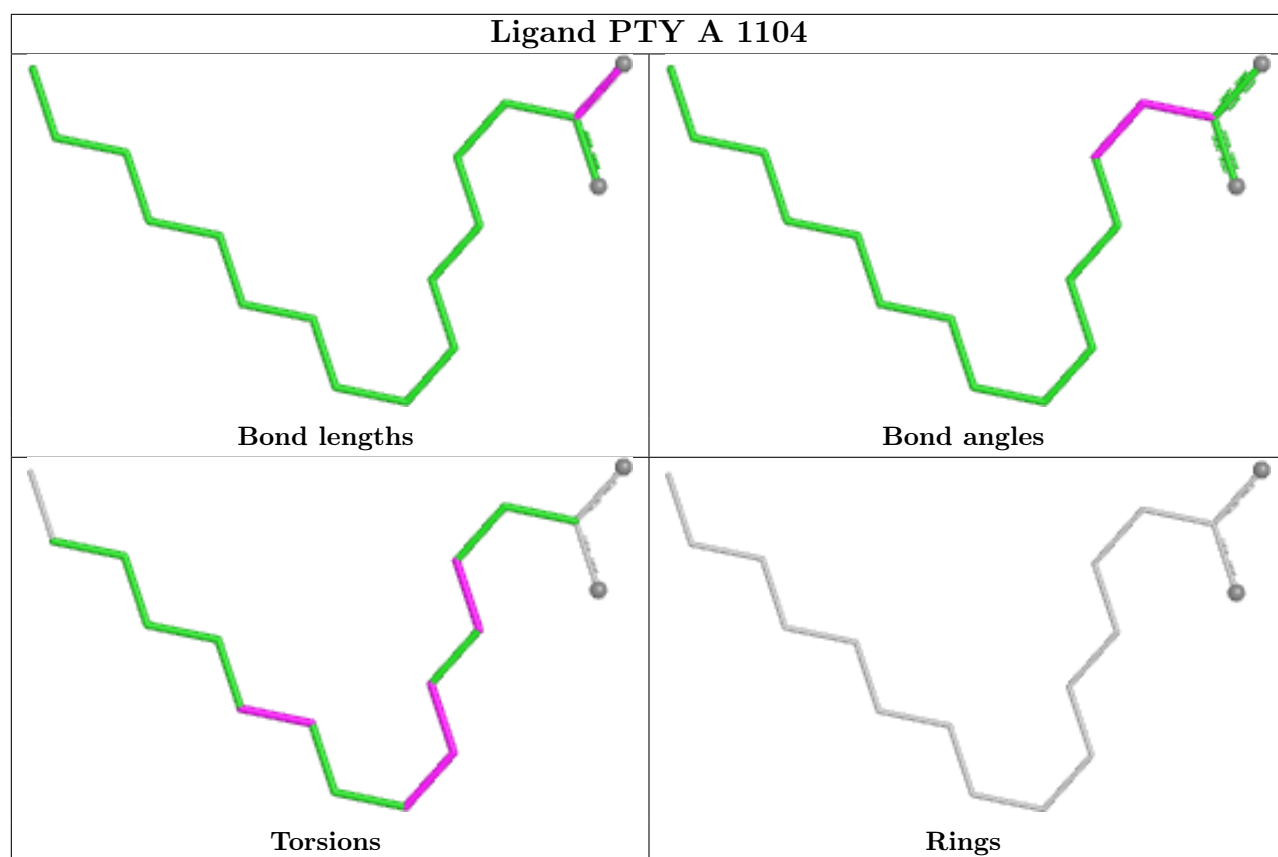


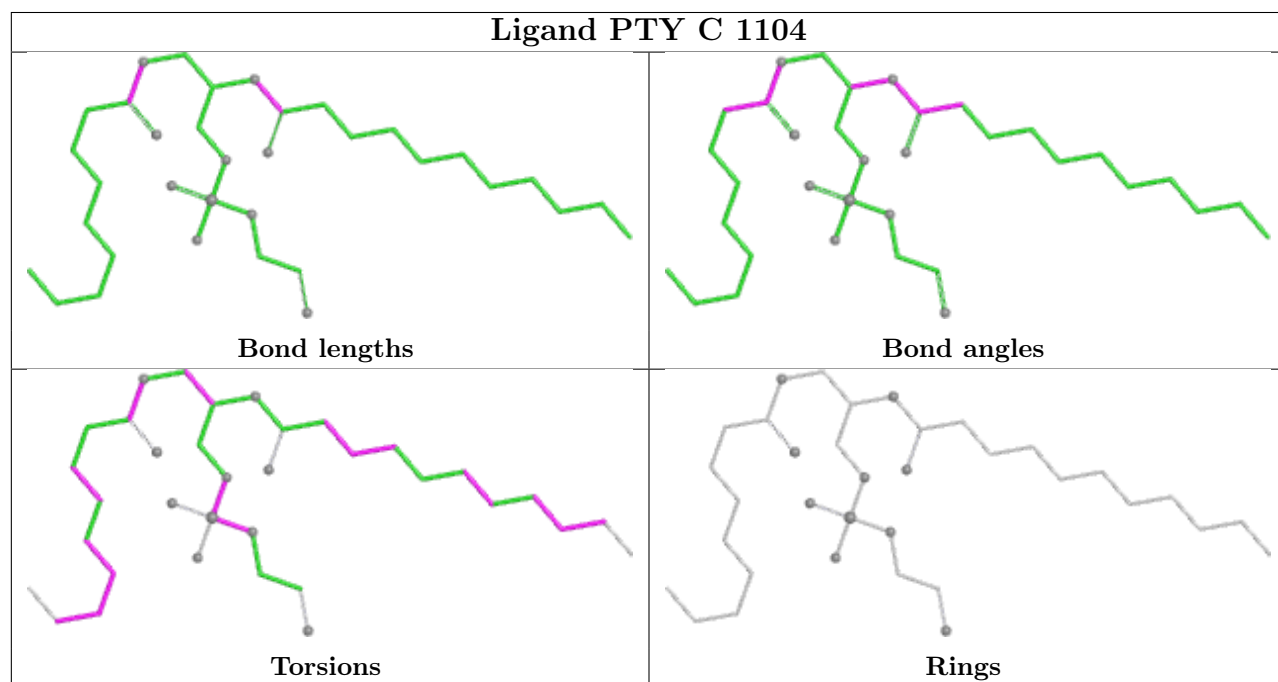
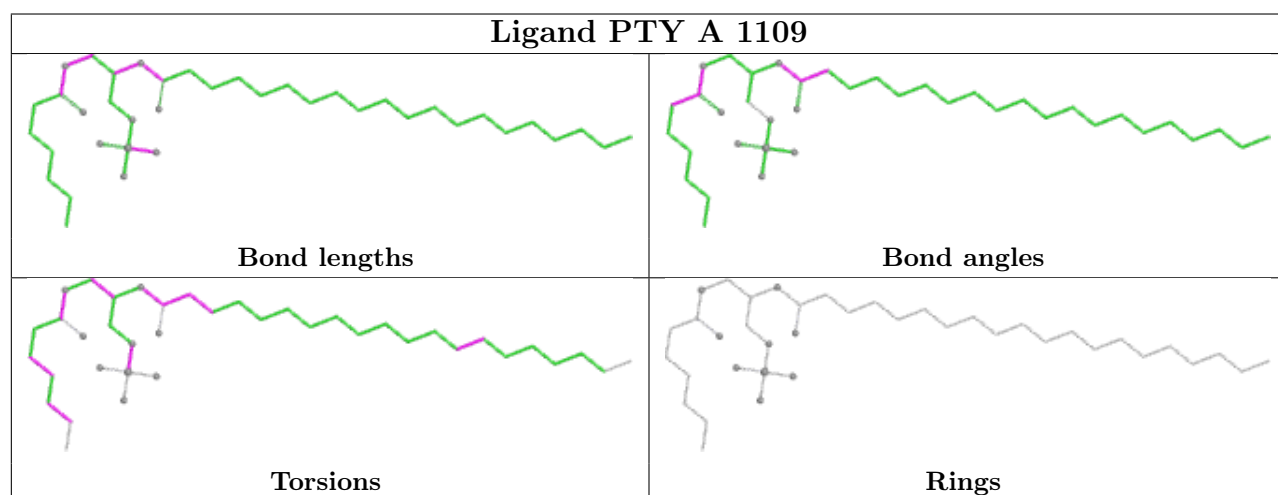
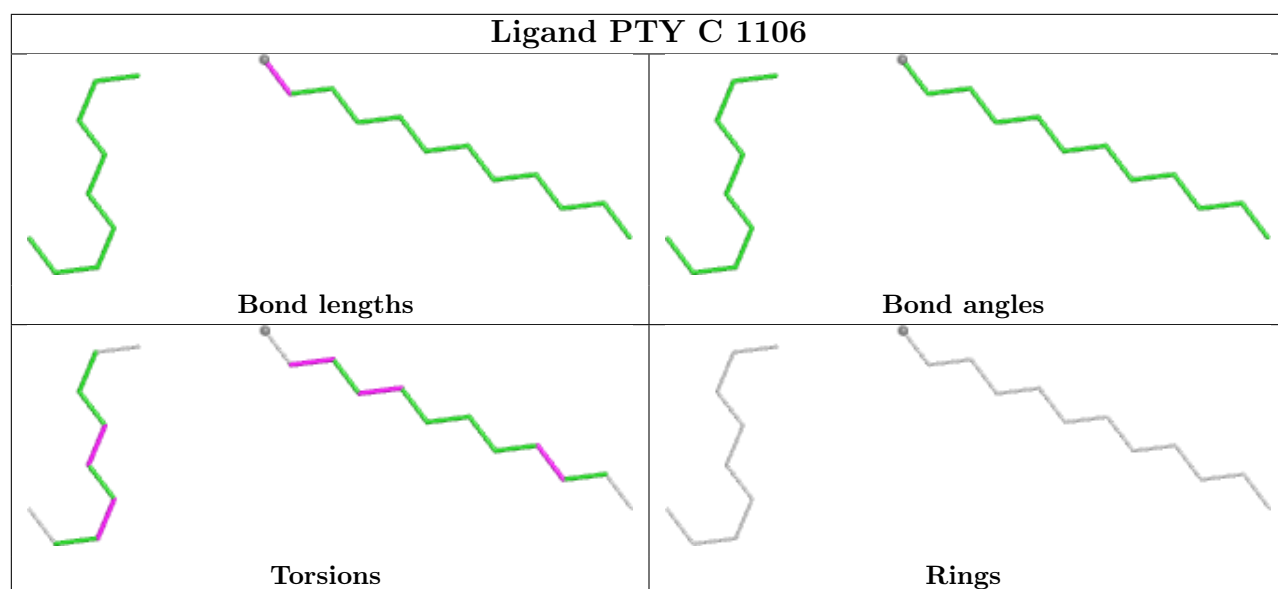


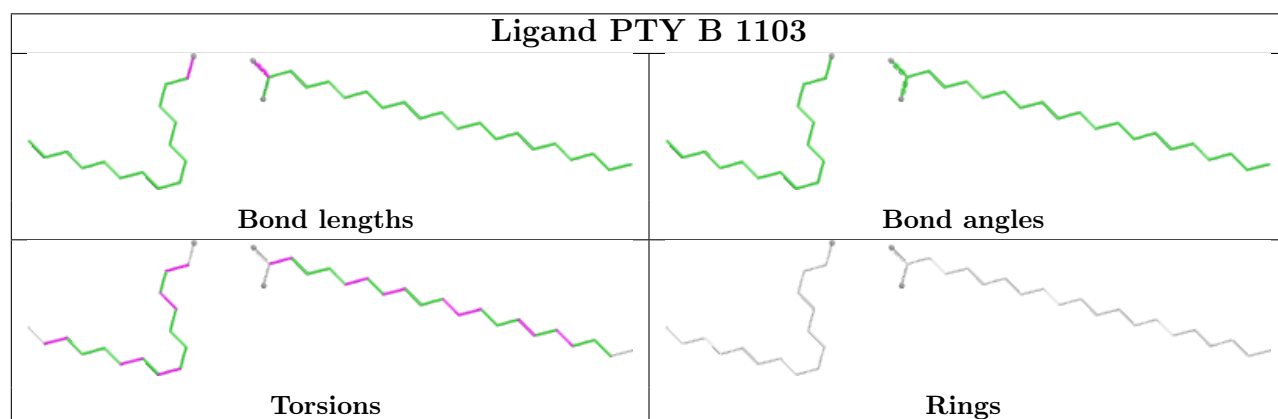
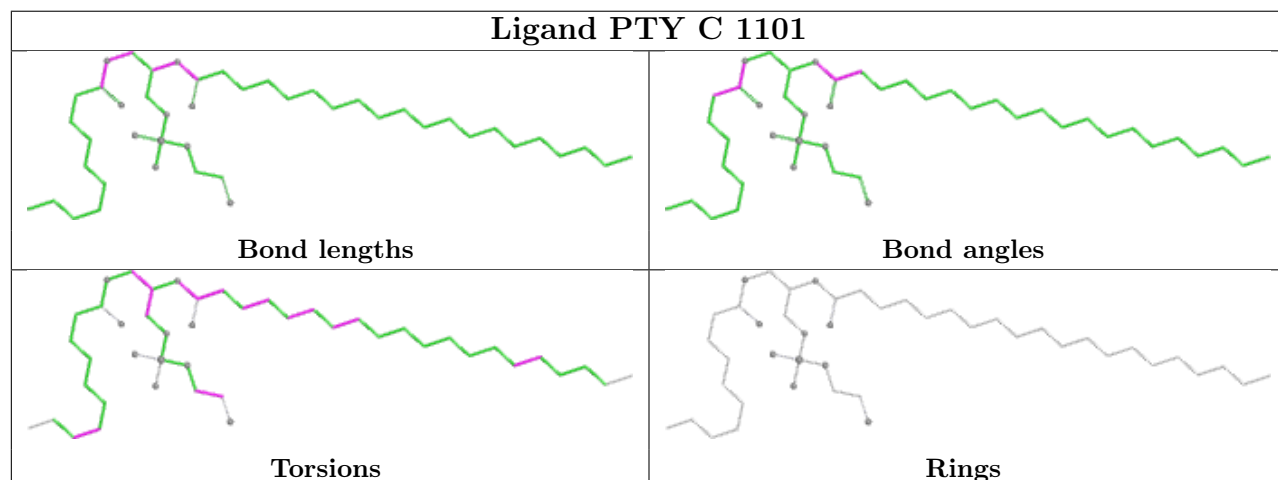
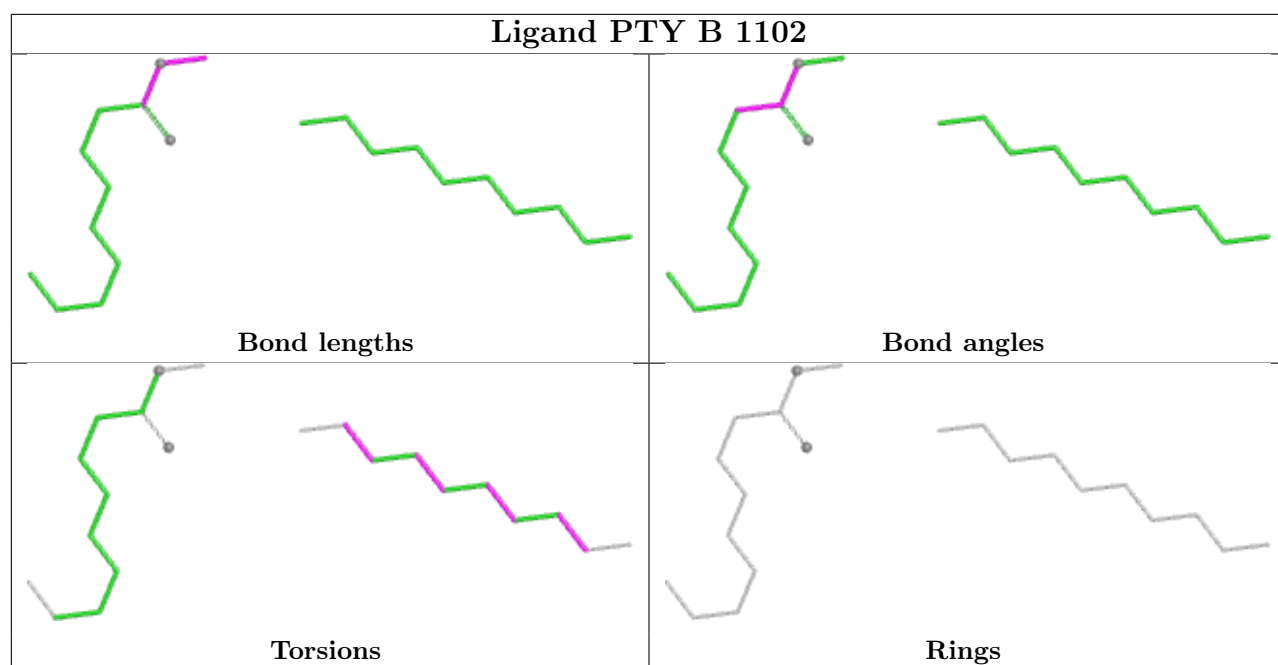


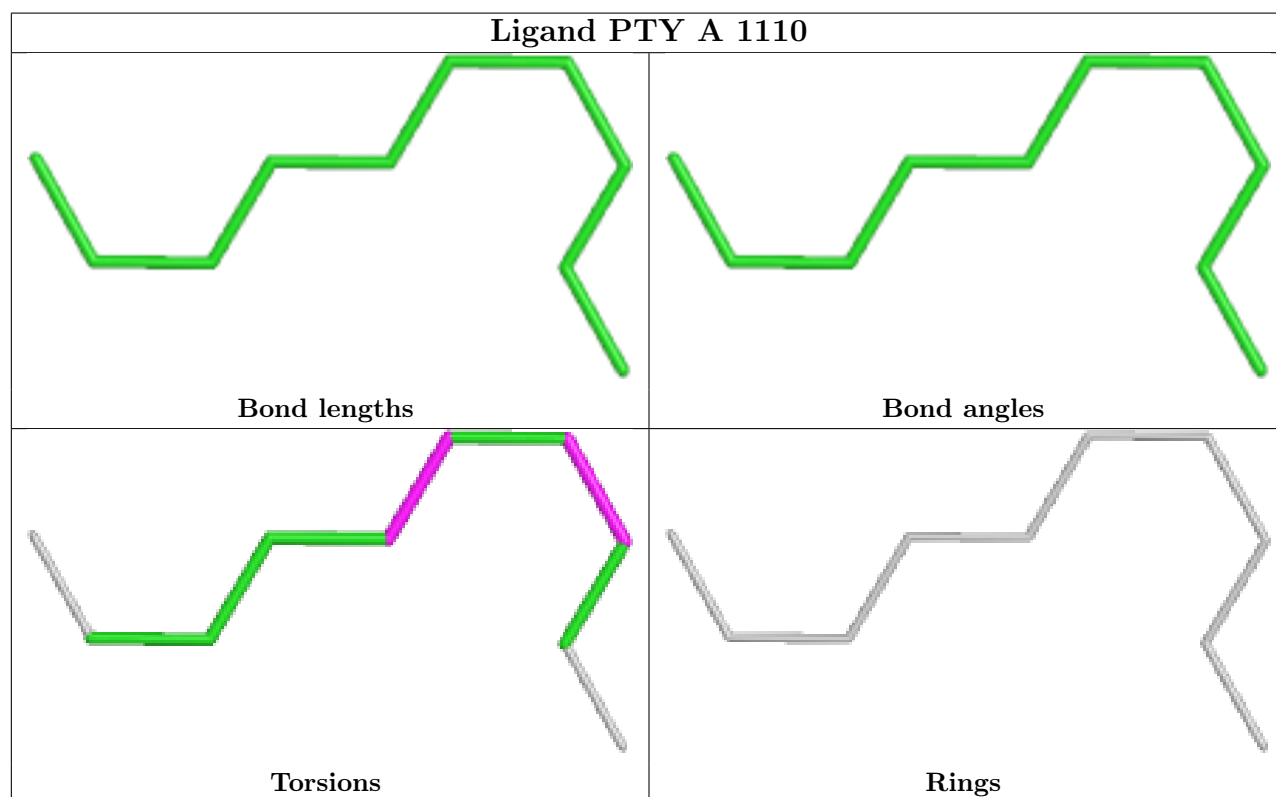
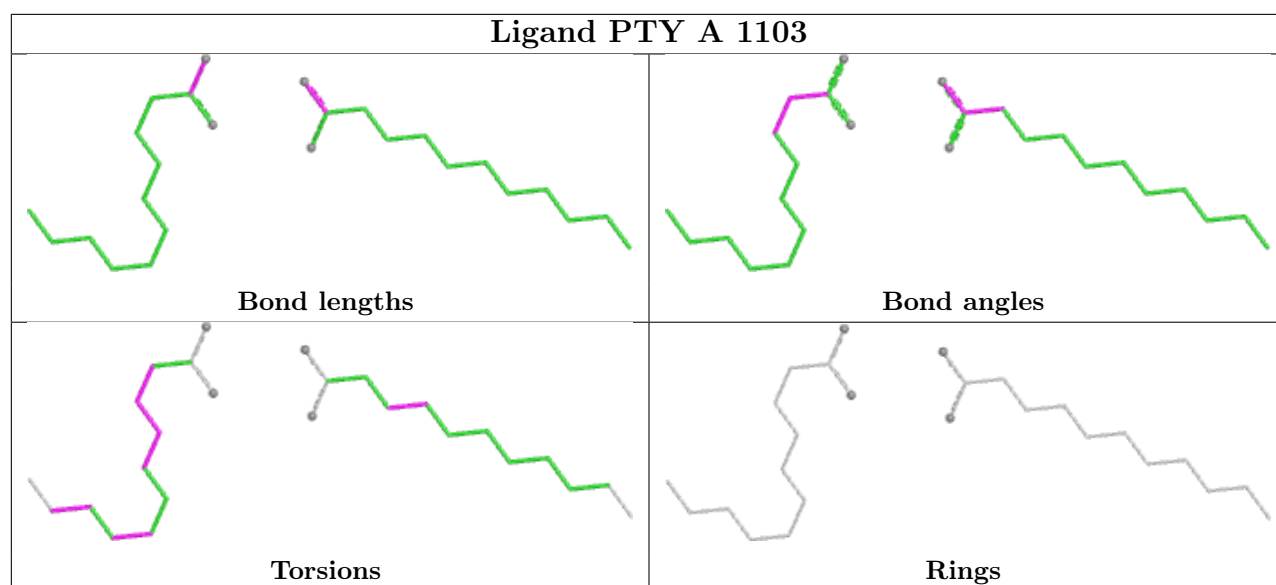


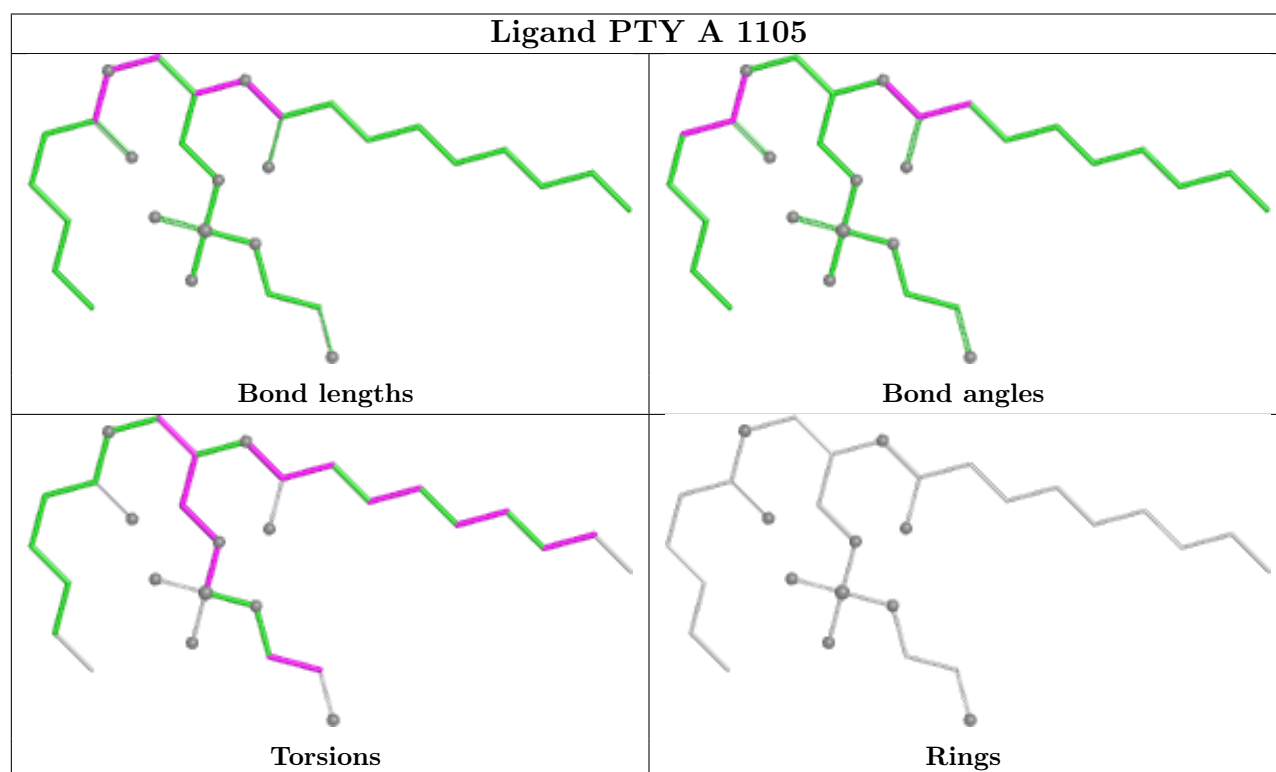
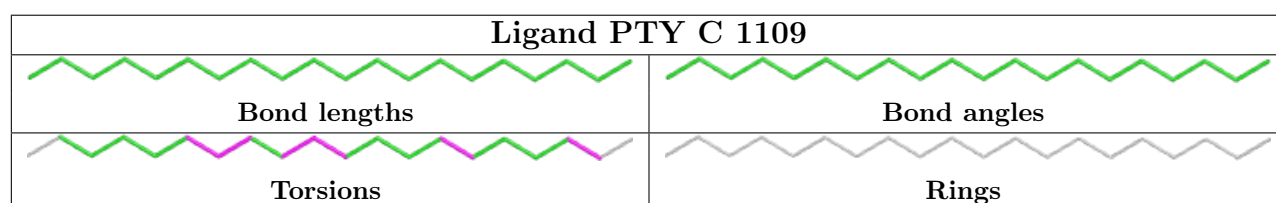
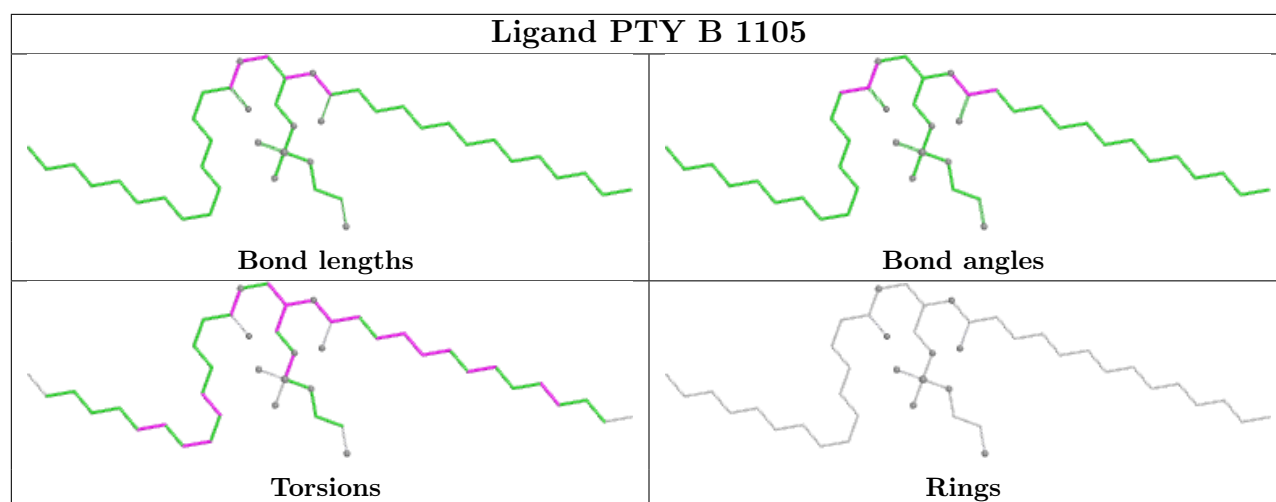


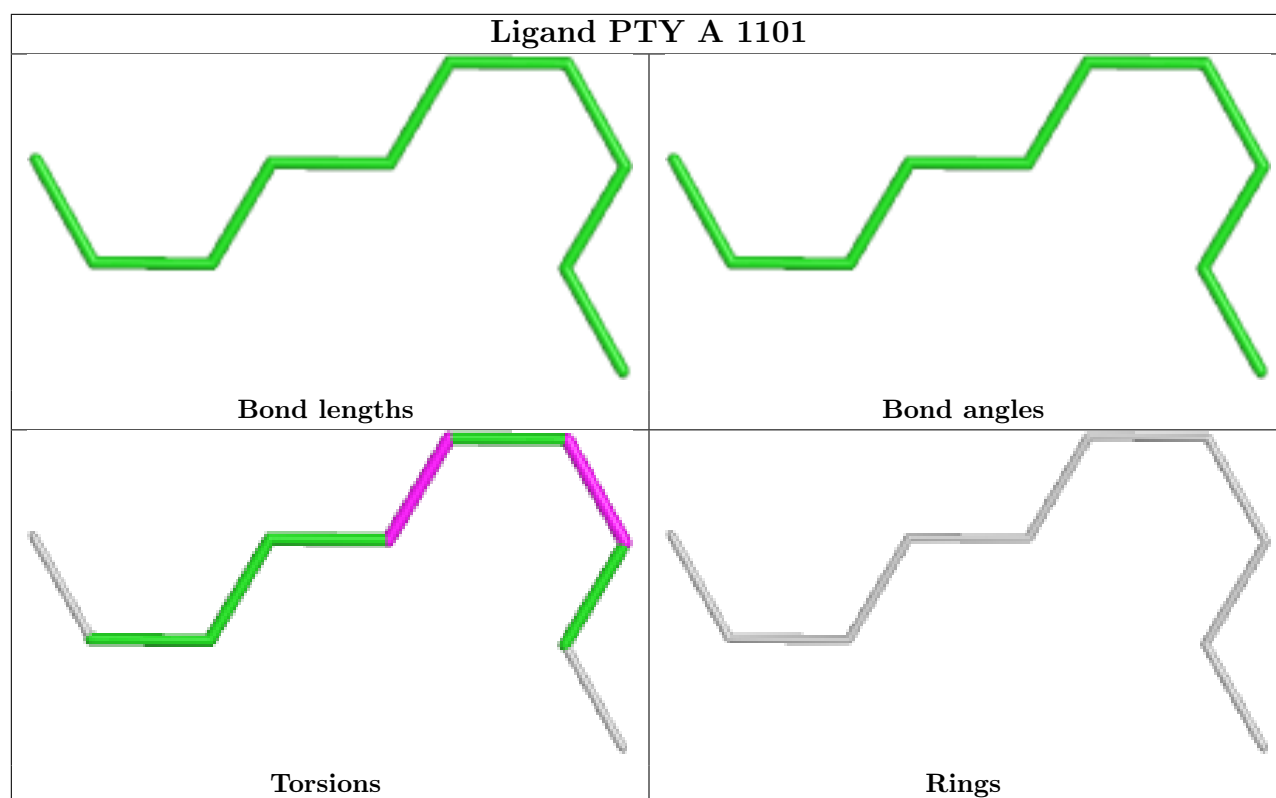












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

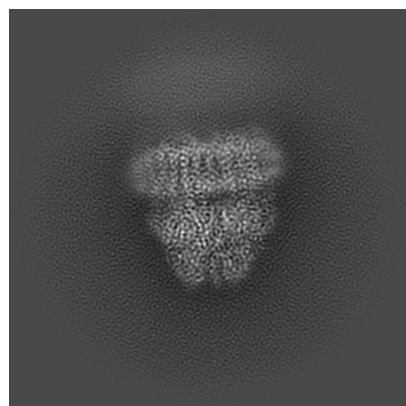
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-27399. These allow visual inspection of the internal detail of the map and identification of artifacts.

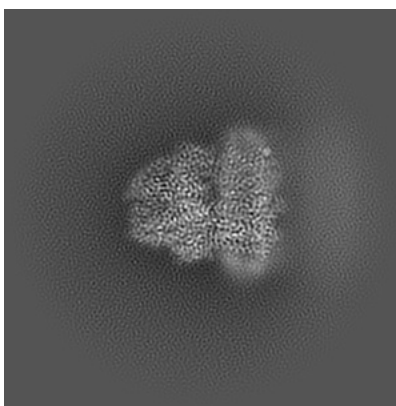
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

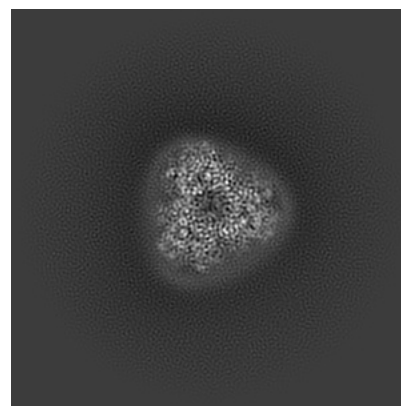
6.1.1 Primary map



X

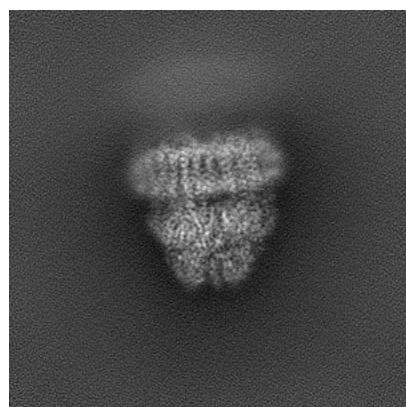


Y

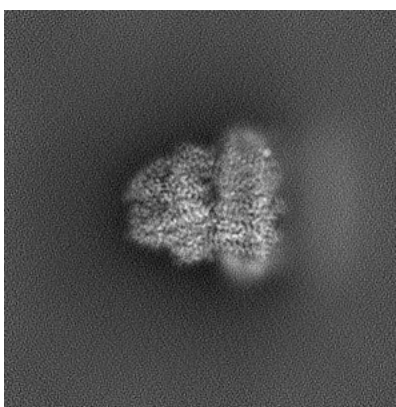


Z

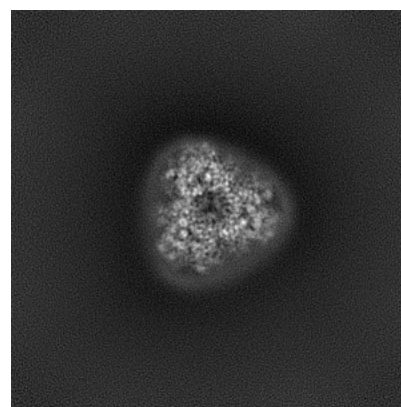
6.1.2 Raw map



X



Y

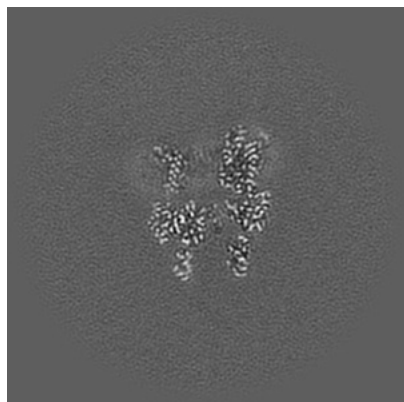


Z

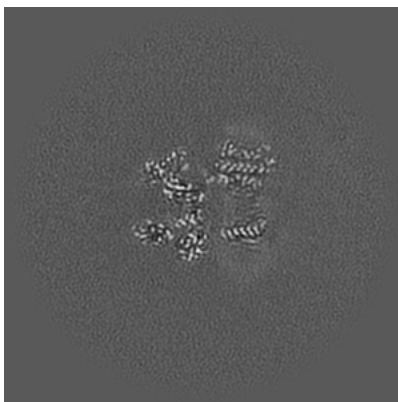
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

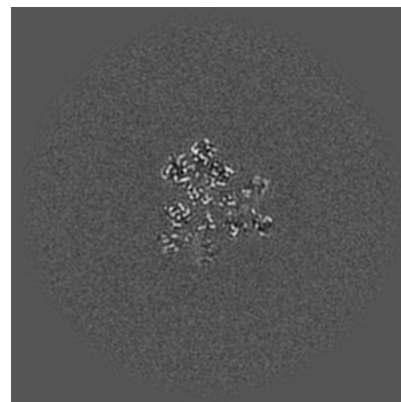
6.2.1 Primary map



X Index: 150

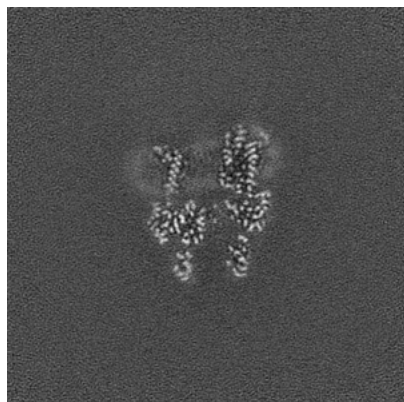


Y Index: 150

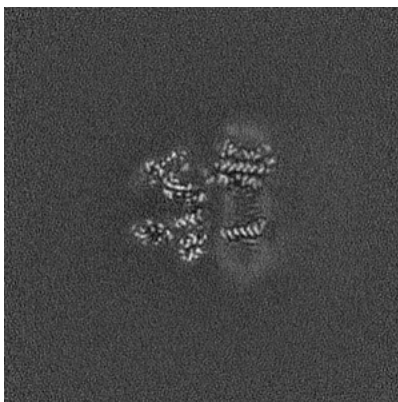


Z Index: 150

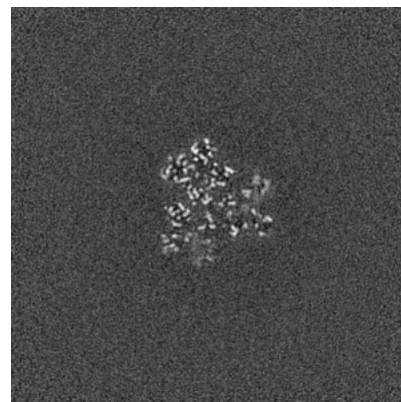
6.2.2 Raw map



X Index: 150



Y Index: 150

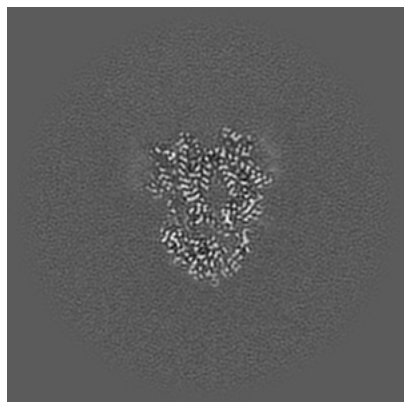


Z Index: 150

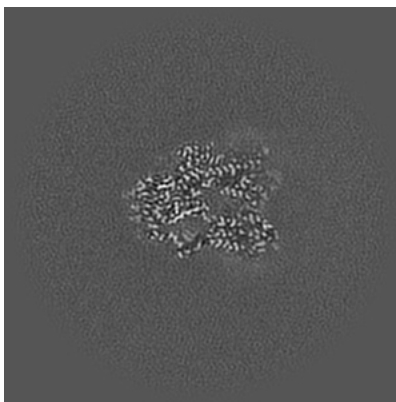
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

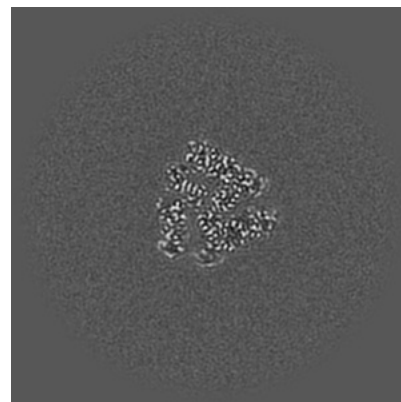
6.3.1 Primary map



X Index: 129

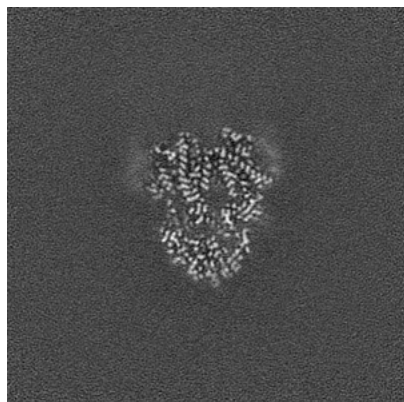


Y Index: 134

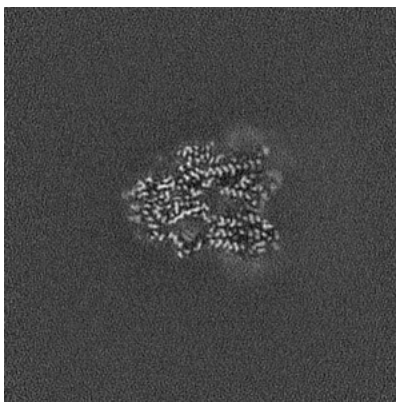


Z Index: 138

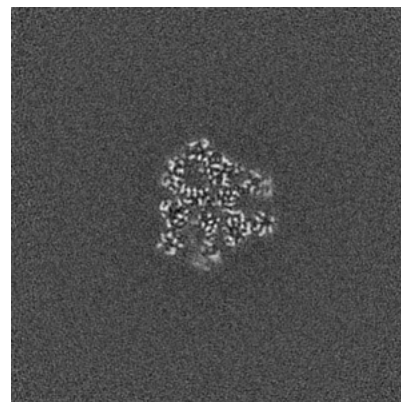
6.3.2 Raw map



X Index: 129



Y Index: 134

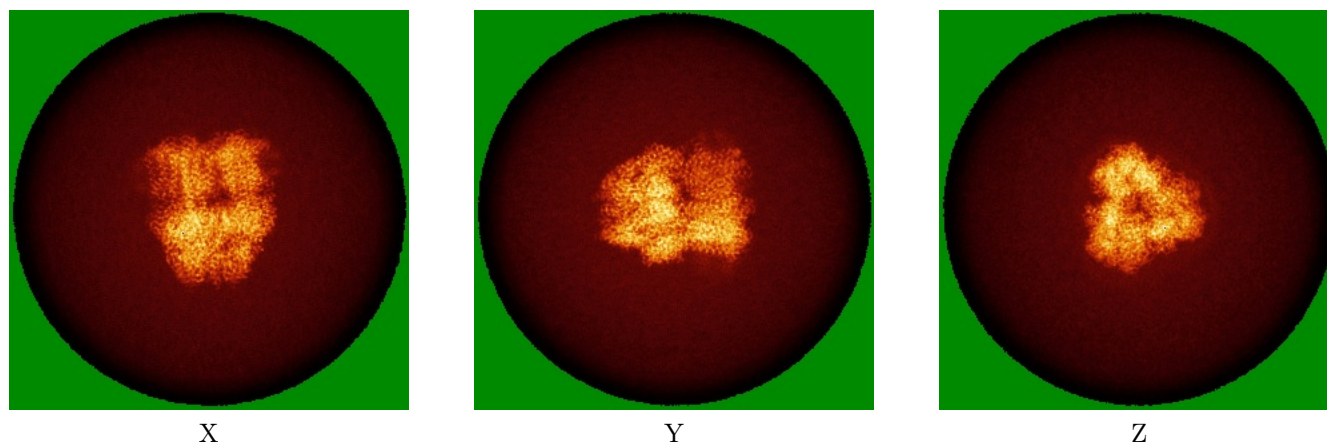


Z Index: 144

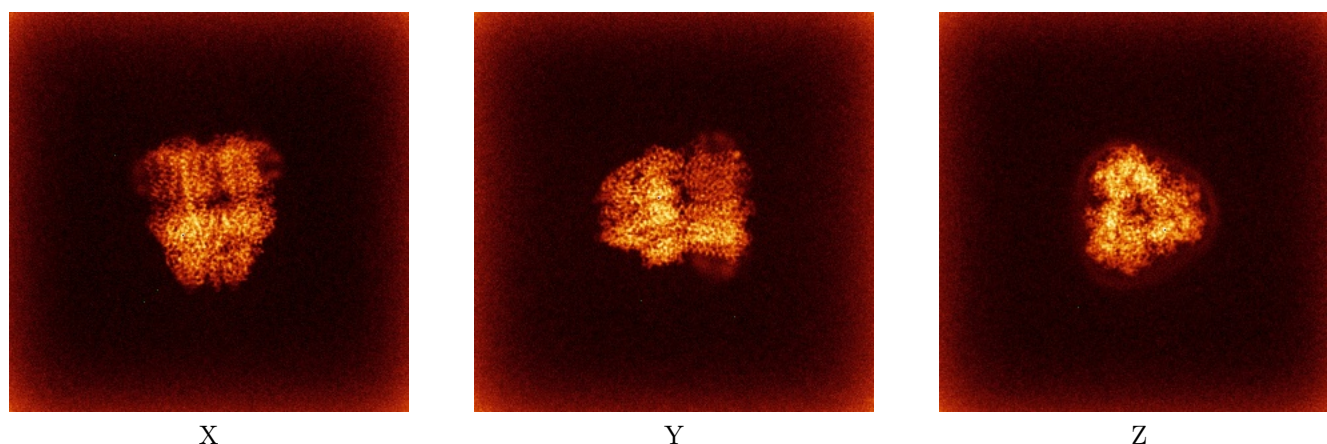
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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

6.5 Orthogonal surface views [i](#)

This section was not generated.

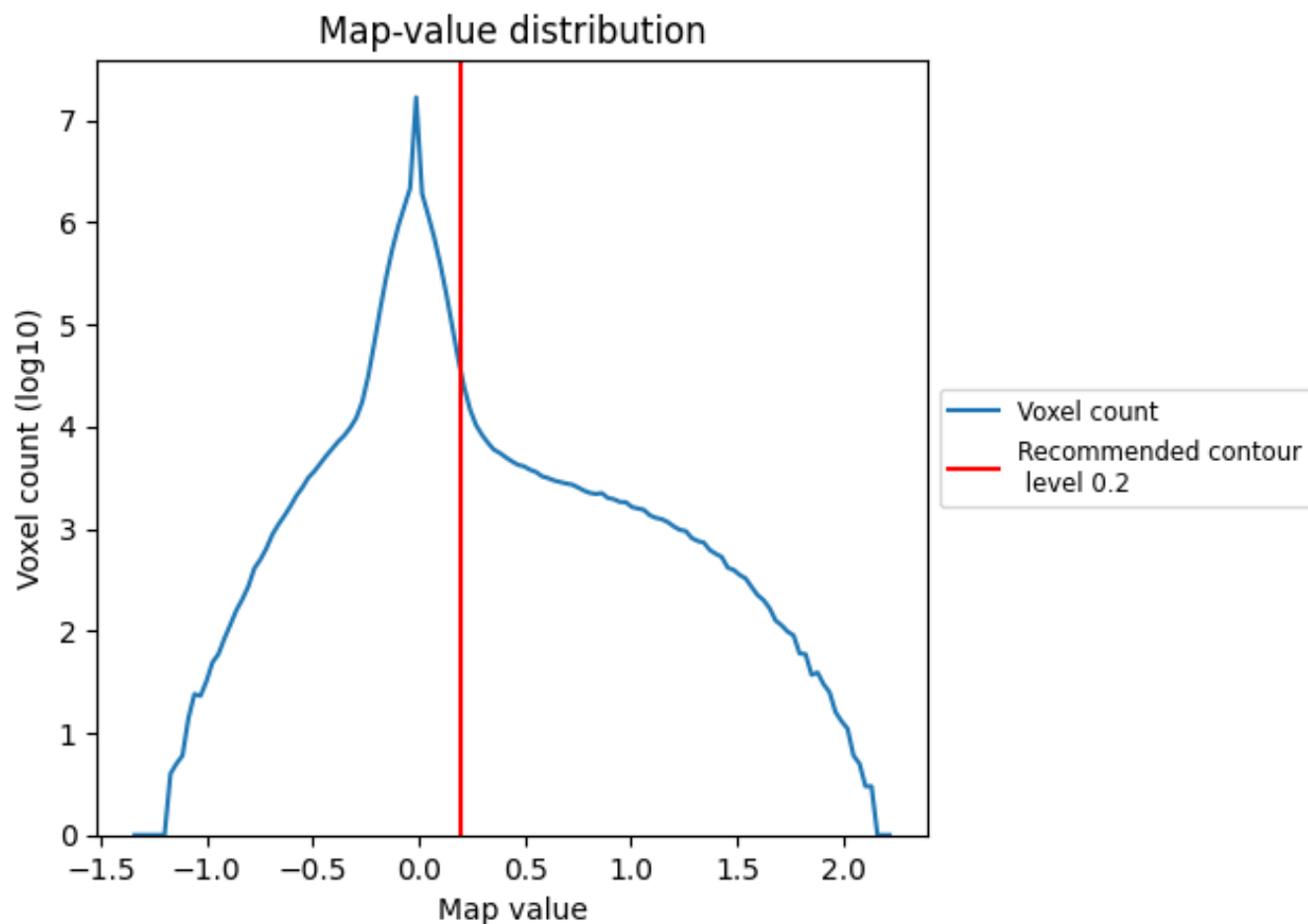
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

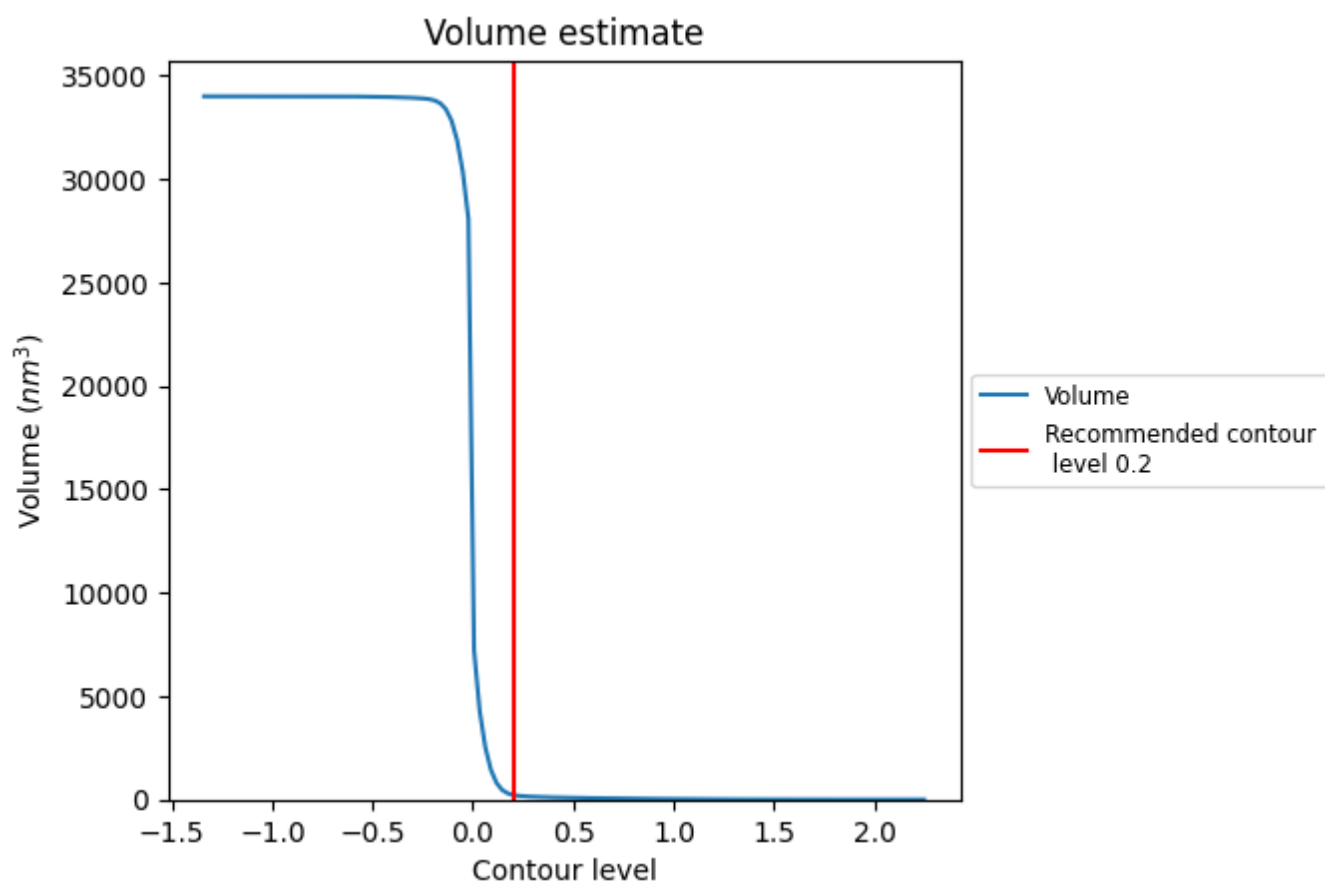
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

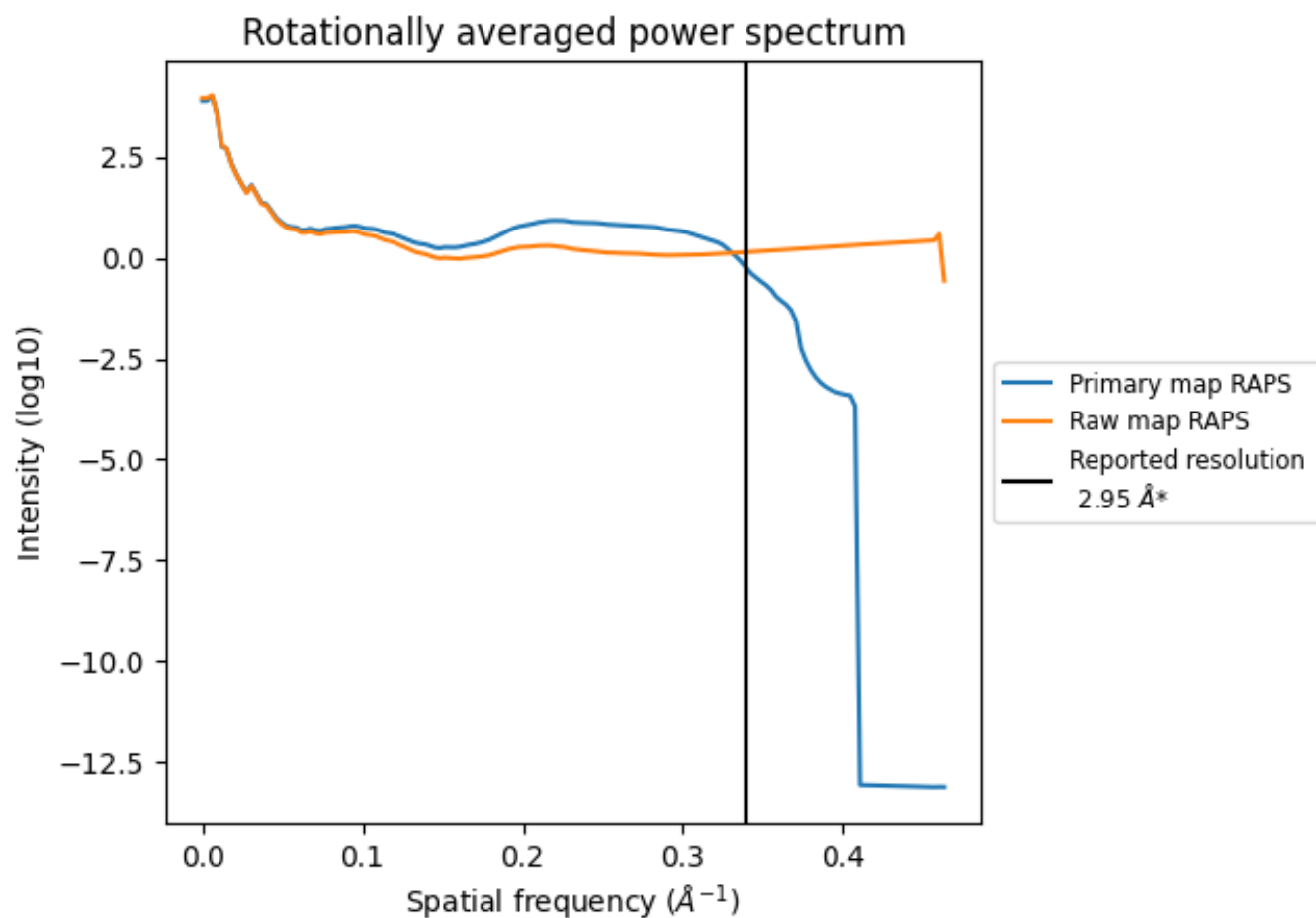
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 224 nm³; this corresponds to an approximate mass of 203 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 ⓘ

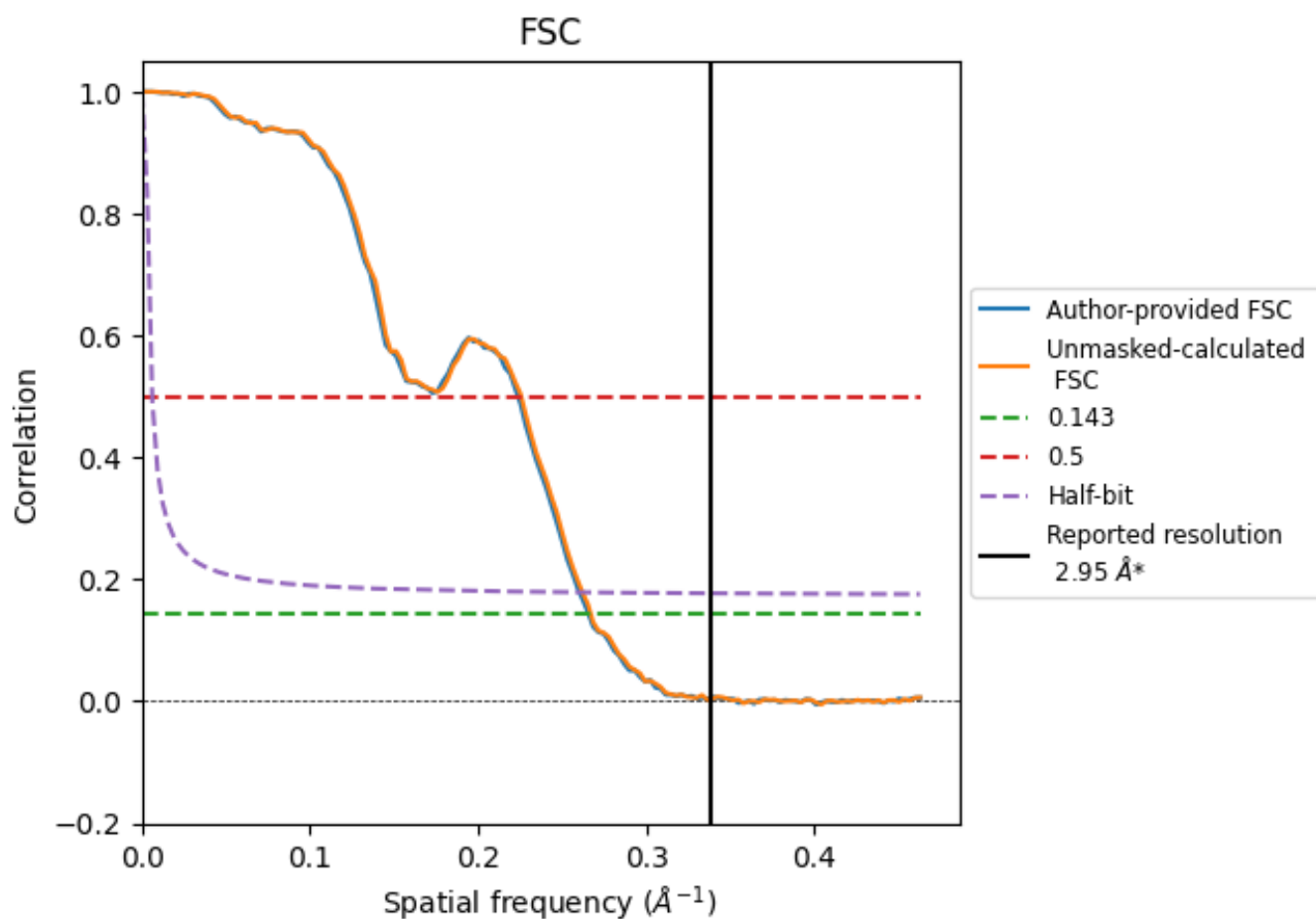


*Reported resolution corresponds to spatial frequency of 0.339 Å⁻¹

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 0.339 Å⁻¹

8.2 Resolution estimates

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.95	-	-
Author-provided FSC curve	3.76	4.47	3.84
Unmasked-calculated*	3.74	4.43	3.81

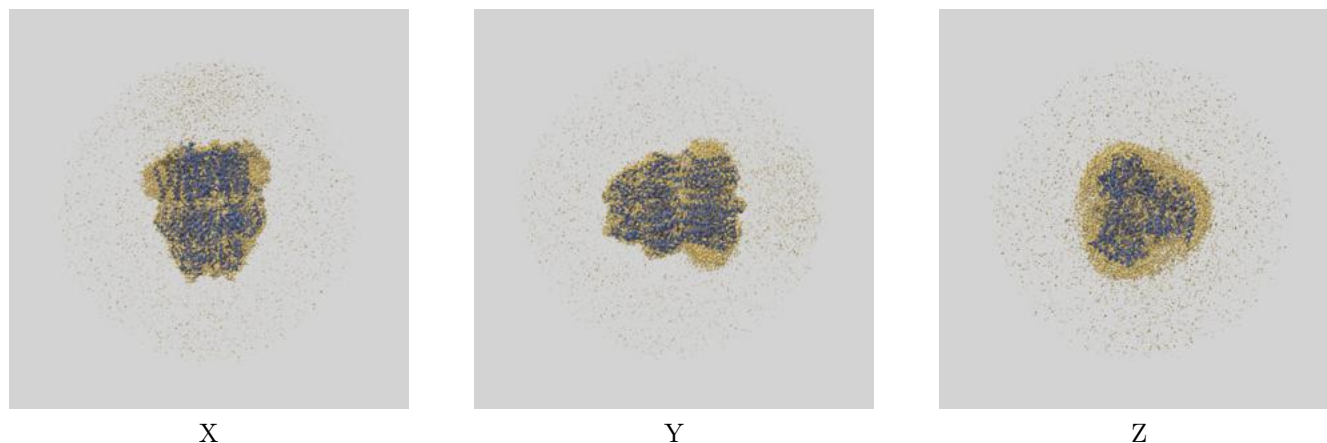
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from author-provided FSC intersecting FSC 0.143 CUT-OFF 3.76 differs from the reported value 2.95 by more than 10 %

The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.74 differs from the reported value 2.95 by more than 10 %

9 Map-model fit [i](#)

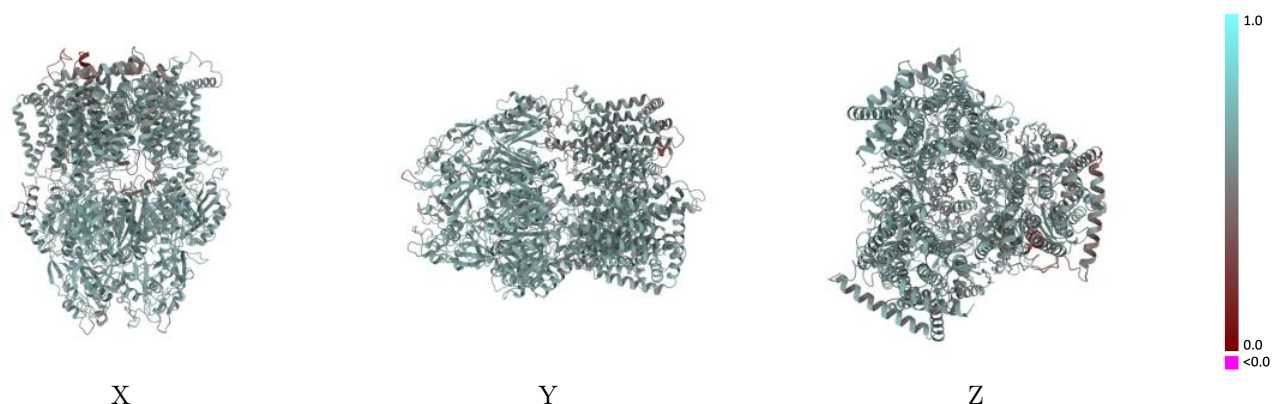
This section contains information regarding the fit between EMDB map EMD-27399 and PDB model 8DEU. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.2 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](#)

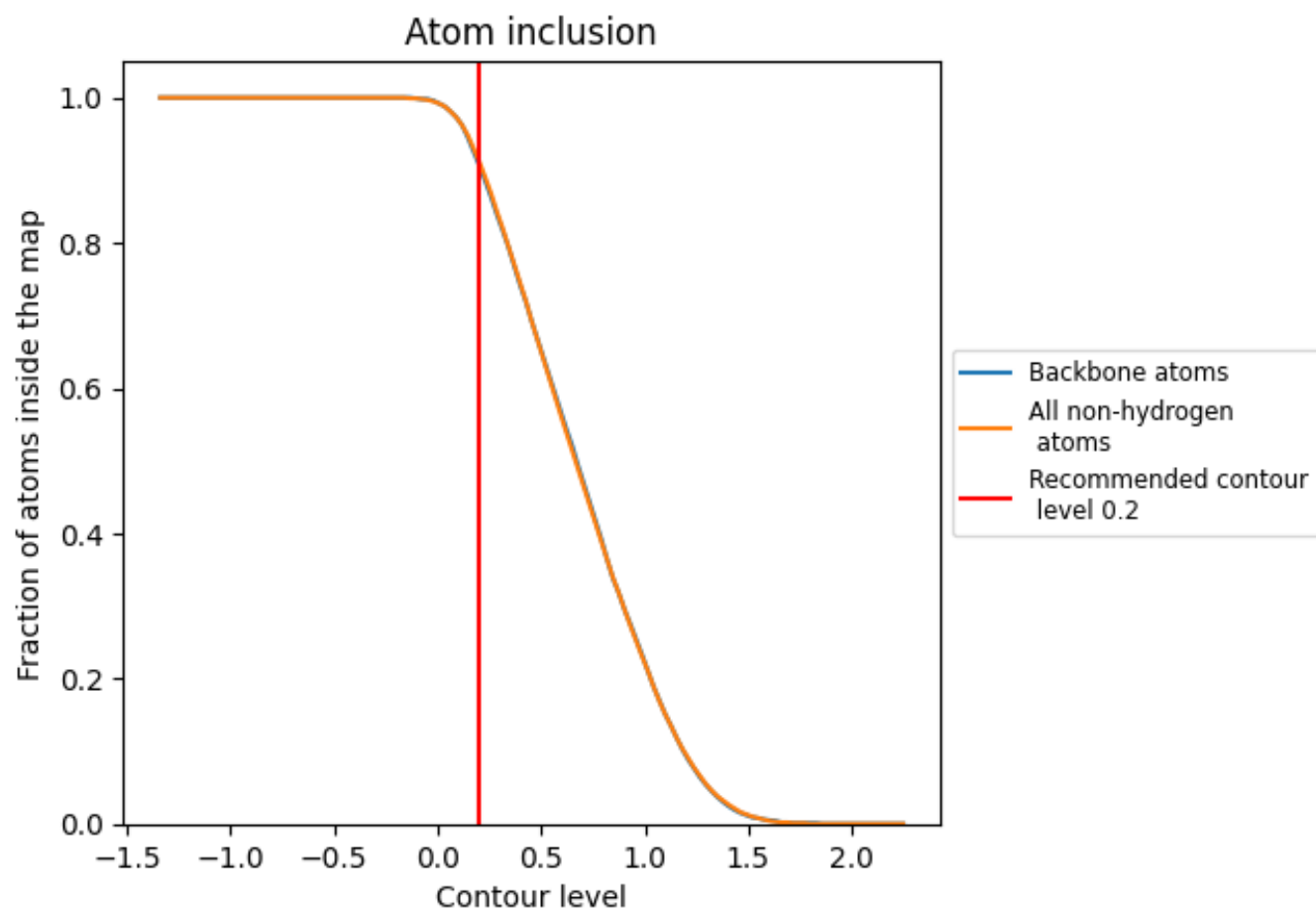


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 91% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9130	0.5780
A	0.8900	0.5640
B	0.9160	0.5820
C	0.9370	0.5910
D	0.5000	0.3540

