



wwPDB EM Validation Summary Report ⓘ

Mar 20, 2026 – 07:45 AM UTC

PDB ID : 7LPD / pdb_00007lpd
EMDB ID : EMD-23478
Title : Cryo-EM structure of full-length TRPV1 with capsaicin at 48 degrees Celsius, in an intermediate state, class 2
Authors : Kwon, D.H.; Zhang, F.; Suo, Y.; Lee, S.-Y.
Deposited on : 2021-02-11
Resolution : 3.55 Å(reported)
Based on initial model : 5IRZ

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

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

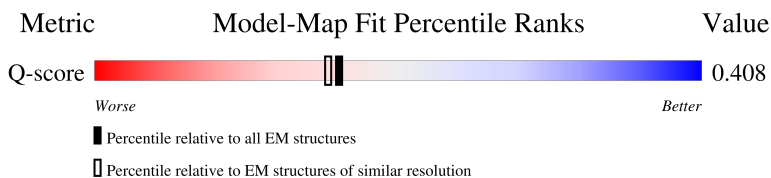
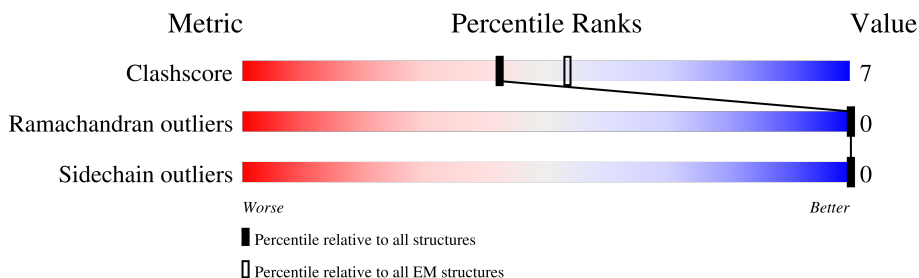
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.55 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	12819 (3.05 - 4.05)

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	868	19% 50% 7% 43%
1	B	868	19% 50% 7% 43%
1	C	868	19% 50% 7% 43%
1	D	868	19% 50% 8% 43%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15564 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 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	497	Total	C	N	O	S	0	0
			3733	2466	599	647	21		
1	B	497	Total	C	N	O	S	0	0
			3733	2466	599	647	21		
1	C	497	Total	C	N	O	S	0	0
			3733	2466	599	647	21		
1	D	497	Total	C	N	O	S	0	0
			3733	2466	599	647	21		

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	839	GLU	-	expression tag	UNP O35433
A	840	ASN	-	expression tag	UNP O35433
A	841	SER	-	expression tag	UNP O35433
A	842	LEU	-	expression tag	UNP O35433
A	843	GLU	-	expression tag	UNP O35433
A	844	VAL	-	expression tag	UNP O35433
A	845	LEU	-	expression tag	UNP O35433
A	846	PHE	-	expression tag	UNP O35433
A	847	GLN	-	expression tag	UNP O35433
A	848	GLY	-	expression tag	UNP O35433
A	849	PRO	-	expression tag	UNP O35433
A	850	ASP	-	expression tag	UNP O35433
A	851	TYR	-	expression tag	UNP O35433
A	852	LYS	-	expression tag	UNP O35433
A	853	ASP	-	expression tag	UNP O35433
A	854	ASP	-	expression tag	UNP O35433
A	855	ASP	-	expression tag	UNP O35433
A	856	ASP	-	expression tag	UNP O35433
A	857	LYS	-	expression tag	UNP O35433
A	858	ALA	-	expression tag	UNP O35433
A	859	HIS	-	expression tag	UNP O35433

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Chain	Residue	Modelled	Actual	Comment	Reference
A	860	HIS	-	expression tag	UNP O35433
A	861	HIS	-	expression tag	UNP O35433
A	862	HIS	-	expression tag	UNP O35433
A	863	HIS	-	expression tag	UNP O35433
A	864	HIS	-	expression tag	UNP O35433
A	865	HIS	-	expression tag	UNP O35433
A	866	HIS	-	expression tag	UNP O35433
A	867	HIS	-	expression tag	UNP O35433
A	868	HIS	-	expression tag	UNP O35433
B	839	GLU	-	expression tag	UNP O35433
B	840	ASN	-	expression tag	UNP O35433
B	841	SER	-	expression tag	UNP O35433
B	842	LEU	-	expression tag	UNP O35433
B	843	GLU	-	expression tag	UNP O35433
B	844	VAL	-	expression tag	UNP O35433
B	845	LEU	-	expression tag	UNP O35433
B	846	PHE	-	expression tag	UNP O35433
B	847	GLN	-	expression tag	UNP O35433
B	848	GLY	-	expression tag	UNP O35433
B	849	PRO	-	expression tag	UNP O35433
B	850	ASP	-	expression tag	UNP O35433
B	851	TYR	-	expression tag	UNP O35433
B	852	LYS	-	expression tag	UNP O35433
B	853	ASP	-	expression tag	UNP O35433
B	854	ASP	-	expression tag	UNP O35433
B	855	ASP	-	expression tag	UNP O35433
B	856	ASP	-	expression tag	UNP O35433
B	857	LYS	-	expression tag	UNP O35433
B	858	ALA	-	expression tag	UNP O35433
B	859	HIS	-	expression tag	UNP O35433
B	860	HIS	-	expression tag	UNP O35433
B	861	HIS	-	expression tag	UNP O35433
B	862	HIS	-	expression tag	UNP O35433
B	863	HIS	-	expression tag	UNP O35433
B	864	HIS	-	expression tag	UNP O35433
B	865	HIS	-	expression tag	UNP O35433
B	866	HIS	-	expression tag	UNP O35433
B	867	HIS	-	expression tag	UNP O35433
B	868	HIS	-	expression tag	UNP O35433
C	839	GLU	-	expression tag	UNP O35433
C	840	ASN	-	expression tag	UNP O35433
C	841	SER	-	expression tag	UNP O35433

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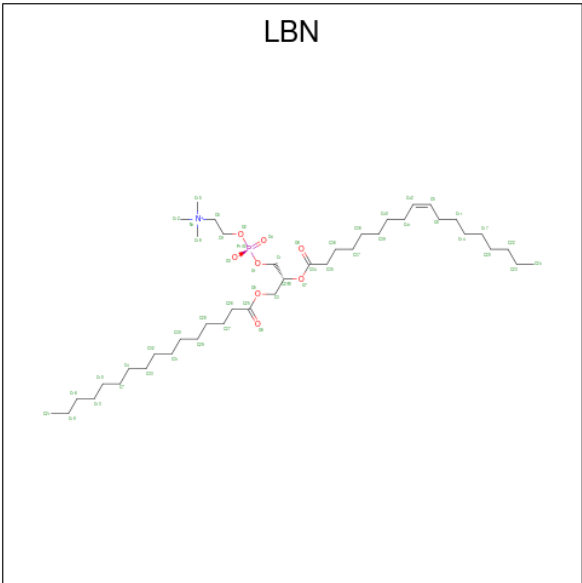
Chain	Residue	Modelled	Actual	Comment	Reference
C	842	LEU	-	expression tag	UNP O35433
C	843	GLU	-	expression tag	UNP O35433
C	844	VAL	-	expression tag	UNP O35433
C	845	LEU	-	expression tag	UNP O35433
C	846	PHE	-	expression tag	UNP O35433
C	847	GLN	-	expression tag	UNP O35433
C	848	GLY	-	expression tag	UNP O35433
C	849	PRO	-	expression tag	UNP O35433
C	850	ASP	-	expression tag	UNP O35433
C	851	TYR	-	expression tag	UNP O35433
C	852	LYS	-	expression tag	UNP O35433
C	853	ASP	-	expression tag	UNP O35433
C	854	ASP	-	expression tag	UNP O35433
C	855	ASP	-	expression tag	UNP O35433
C	856	ASP	-	expression tag	UNP O35433
C	857	LYS	-	expression tag	UNP O35433
C	858	ALA	-	expression tag	UNP O35433
C	859	HIS	-	expression tag	UNP O35433
C	860	HIS	-	expression tag	UNP O35433
C	861	HIS	-	expression tag	UNP O35433
C	862	HIS	-	expression tag	UNP O35433
C	863	HIS	-	expression tag	UNP O35433
C	864	HIS	-	expression tag	UNP O35433
C	865	HIS	-	expression tag	UNP O35433
C	866	HIS	-	expression tag	UNP O35433
C	867	HIS	-	expression tag	UNP O35433
C	868	HIS	-	expression tag	UNP O35433
D	839	GLU	-	expression tag	UNP O35433
D	840	ASN	-	expression tag	UNP O35433
D	841	SER	-	expression tag	UNP O35433
D	842	LEU	-	expression tag	UNP O35433
D	843	GLU	-	expression tag	UNP O35433
D	844	VAL	-	expression tag	UNP O35433
D	845	LEU	-	expression tag	UNP O35433
D	846	PHE	-	expression tag	UNP O35433
D	847	GLN	-	expression tag	UNP O35433
D	848	GLY	-	expression tag	UNP O35433
D	849	PRO	-	expression tag	UNP O35433
D	850	ASP	-	expression tag	UNP O35433
D	851	TYR	-	expression tag	UNP O35433
D	852	LYS	-	expression tag	UNP O35433
D	853	ASP	-	expression tag	UNP O35433

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Chain	Residue	Modelled	Actual	Comment	Reference
D	854	ASP	-	expression tag	UNP O35433
D	855	ASP	-	expression tag	UNP O35433
D	856	ASP	-	expression tag	UNP O35433
D	857	LYS	-	expression tag	UNP O35433
D	858	ALA	-	expression tag	UNP O35433
D	859	HIS	-	expression tag	UNP O35433
D	860	HIS	-	expression tag	UNP O35433
D	861	HIS	-	expression tag	UNP O35433
D	862	HIS	-	expression tag	UNP O35433
D	863	HIS	-	expression tag	UNP O35433
D	864	HIS	-	expression tag	UNP O35433
D	865	HIS	-	expression tag	UNP O35433
D	866	HIS	-	expression tag	UNP O35433
D	867	HIS	-	expression tag	UNP O35433
D	868	HIS	-	expression tag	UNP O35433

- Molecule 2 is 1-palmitoyl-2-oleoyl-sn-glycero-3-phosphocholine (CCD ID: LBN) (formula: C₄₂H₈₂NO₈P).



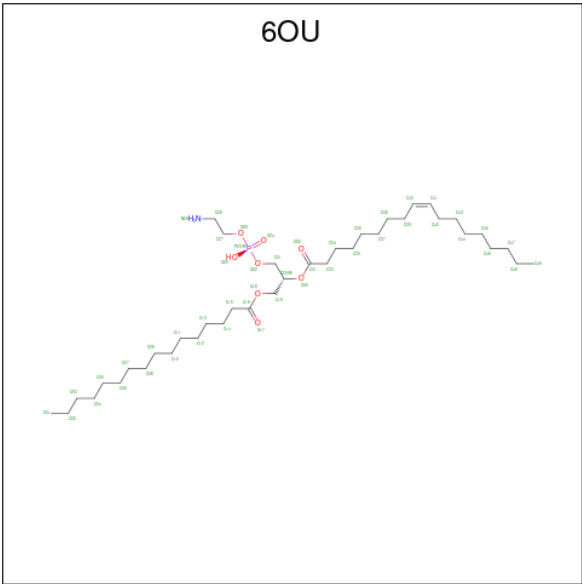
Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	N	O	P	0
			33	23	1	8	1	
2	A	1	Total	C	O			0
			17	12	5			
2	B	1	Total	C	N	O	P	0
			33	23	1	8	1	

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Mol	Chain	Residues	Atoms					AltConf		
2	B	1	Total	C	O				0	
			17	12	5					
2	C	1	Total	C	N	O	P			0
			33	23	1	8	1			
2	C	1	Total	C	O				0	
			17	12	5					
2	D	1	Total	C	N	O	P			0
			33	23	1	8	1			
2	D	1	Total	C	O				0	
			17	12	5					

- Molecule 3 is [(2 {R})-1-[2-azanylethoxy(oxidanyl)phosphoryl]oxy-3-hexadecanoyloxy-prop an-2-yl] ({Z})-octadec-9-enoate (CCD ID: 6OU) (formula: C₃₉H₇₆NO₈P).



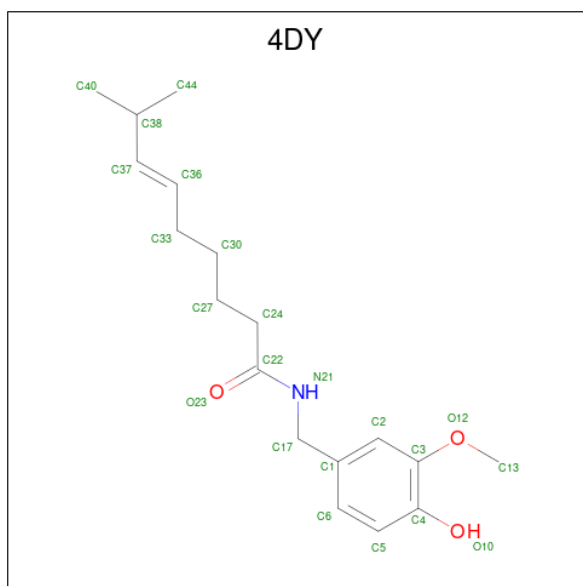
Mol	Chain	Residues	Atoms				AltConf
3	A	1	Total	C	O	P	0
			31	22	8	1	
3	A	1	Total	C	O	P	0
			24	15	8	1	
3	A	1	Total	C	O	P	0
			31	22	8	1	
3	A	1	Total	C	O	P	0
			24	15	8	1	
3	B	1	Total	C	O	P	0
			31	22	8	1	
3	B	1	Total	C	O	P	0
			24	15	8	1	

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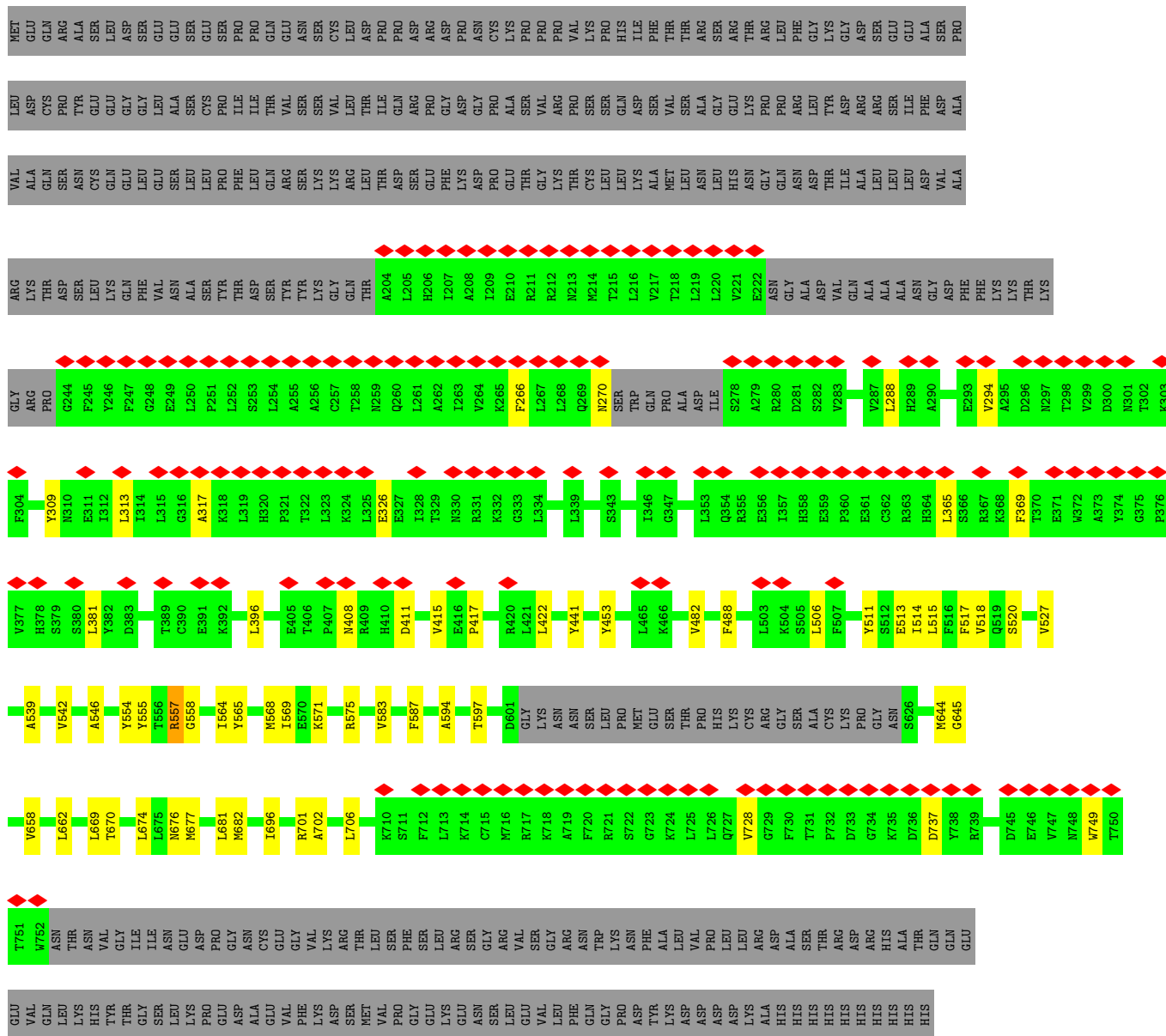
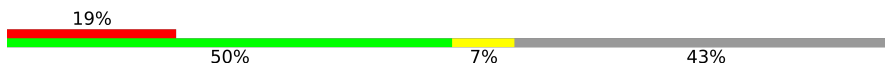
Mol	Chain	Residues	Atoms				AltConf
3	B	1	Total	C	O	P	0
			31	22	8	1	
3	C	1	Total	C	O	P	0
			31	22	8	1	
3	C	1	Total	C	O	P	0
			24	15	8	1	
3	C	1	Total	C	O	P	0
			31	22	8	1	
3	D	1	Total	C	O	P	0
			31	22	8	1	
3	D	1	Total	C	O	P	0
			31	22	8	1	

- Molecule 4 is (6E)-N-(4-hydroxy-3-methoxybenzyl)-8-methylnon-6-enamide (CCD ID: 4DY) (formula: $C_{18}H_{27}NO_3$).

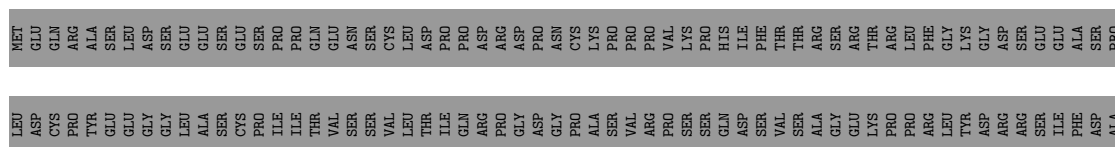
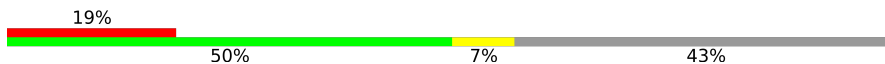


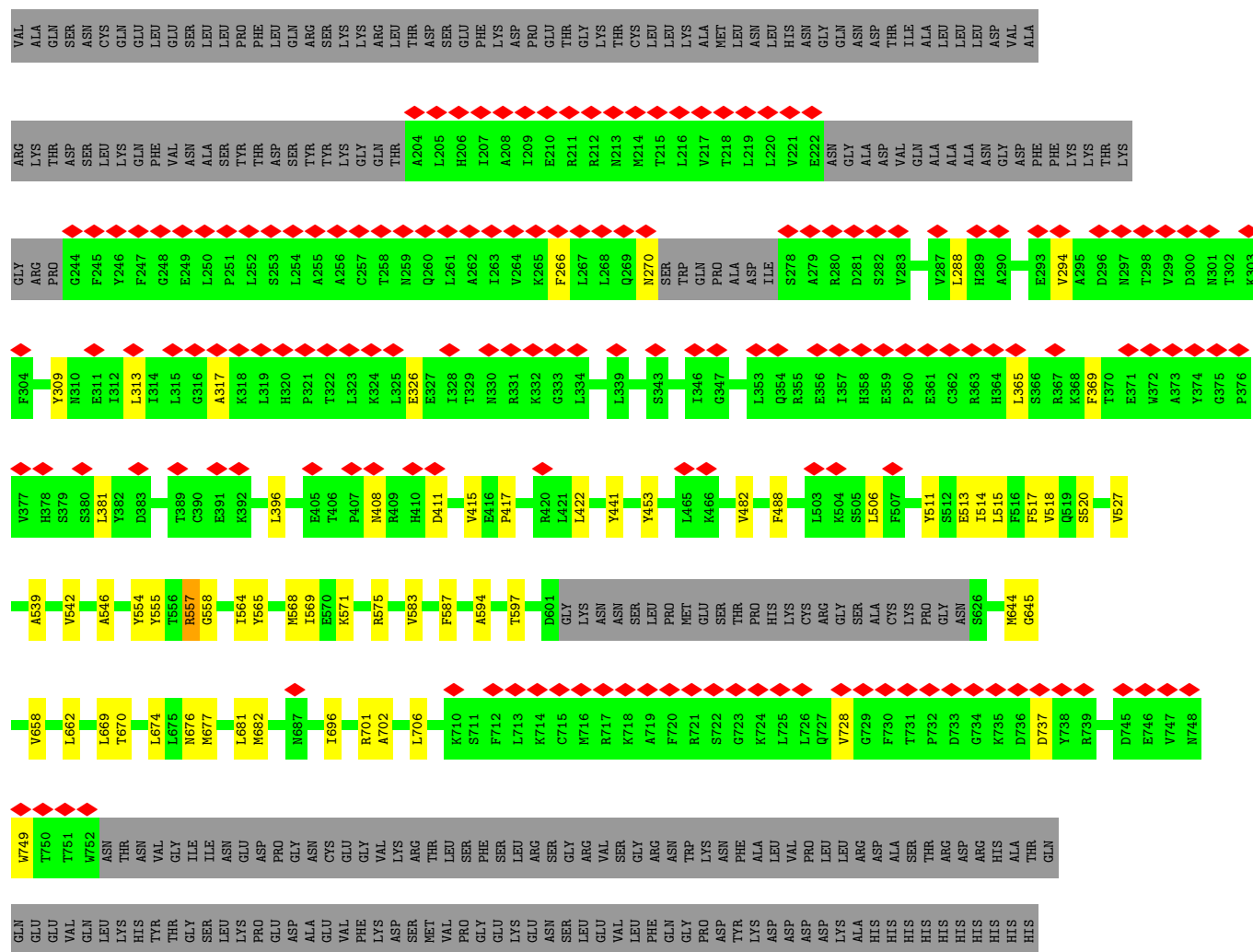
Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	N	O	0
			22	18	1	3	
4	B	1	Total	C	N	O	0
			22	18	1	3	
4	C	1	Total	C	N	O	0
			22	18	1	3	
4	D	1	Total	C	N	O	0
			22	18	1	3	

Chain B:



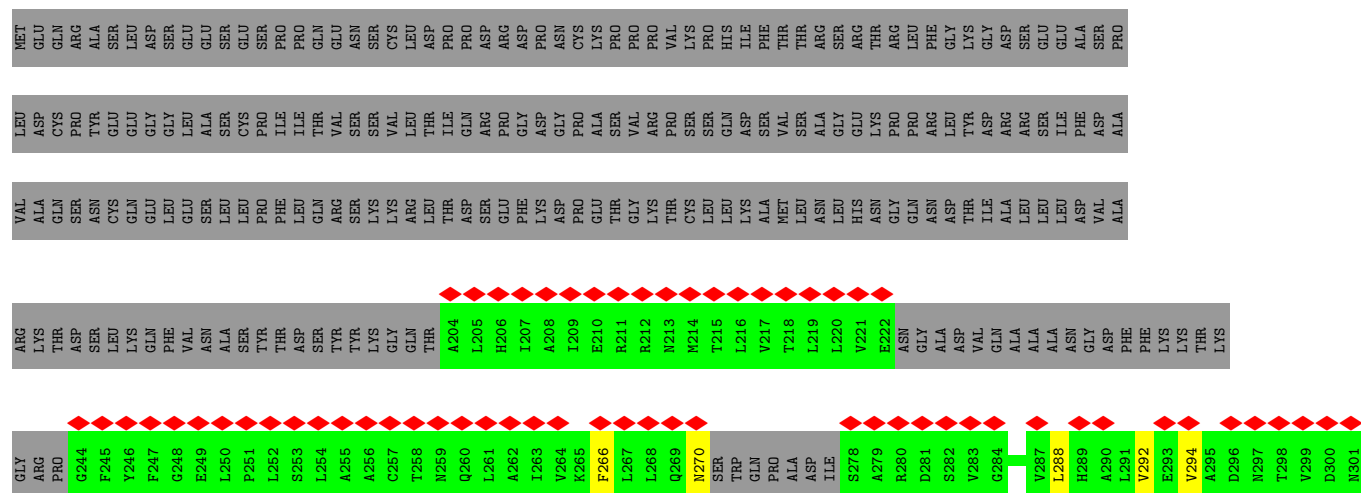
Chain C:





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

Chain D:



T302	K303	F304	Y309	N310	E311	I312	L313	I314	L315	G316	A317	K318	L319	H320	P321	T322	L323	K324	L325	E326	E327	I328	T329	N330	R331	K332	G333	L334	A338	L339	S343	I346	G347	V348	L353	Q354	R355	E356	I357	H358	E359	P360	E361	C362	R363	H364	L365	S366	R367	K368	F369	T370	E371	W372			
A373	Y374	G375	P376	V377	H378	S379	S380	L381	Y382	D383	T389	C390	E391	K392	L396	E405	T406	P407	N408	R409	H410	D411	V415	E416	P417	R420	L421	L422	Y441	Y453	L465	K466	V482	F488	L503	K504	S505	L506	F507	Y511	S512	E513	I514	L515	F516	F517	V518										
Q519	S520	V527	A539	V542	A546	Y554	Y555	T556	R557	G558	I564	Y565	M568	I569	E570	K571	R575	V583	F587	A594	T597	D601	GLY	LYS	ASN	ASN	SER	LEU	PRO	MET	GLU	SER	THR	PRO	HIS	LYS	CYS	ARG	GLY	SER	ALA	CYS	LYS	PRO	GLY	ASN											
S626	M644	G645	V658	L662	L669	T670	L674	L675	N676	M677	L681	M682	I696	R701	A702	L706	F712	L713	K714	C715	M716	R717	K718	A719	F720	R721	S722	G723	K724	L725	L726	Q727	V728	G729	F730	T731	P732	D733	G734	K735	D736	D737	Y738	R739	D745	E746	V747	N748									
W749	T750	T751	W752	ASN	THR	ASN	ASN	VAL	GLY	THR	ILE	ASN	GLU	ASP	PRO	GLY	ASN	ALA	GLU	CYS	VAL	PHE	GLY	LYS	VAL	ASP	SER	ARG	THR	LEU	SER	GLY	ARG	VAL	VAL	SER	ASN	ASN	PHE	ALA	LEU	VAL	PRO	ASP	ASP	ASP	ASP	LEU	LEU	ALA	HIS	HIS	HIS	HIS	HIS	HIS	GLN
GLN	GLU	GLU	VAL	GLN	LEU	LYS	HIS	TYR	THR	GLY	SER	LEU	LYS	PRO	GLU	ASP	ALA	GLU	CYS	VAL	PHE	GLY	LYS	ASP	SER	MET	VAL	PRO	GLY	PHE	GLN	GLY	PRO	ASP	TYR	LYS	ASP	ASP	ASP	ASP	ASP	LYS	ALA	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS				

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	18991	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.329	Depositor
Minimum map value	-1.503	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.074	Depositor
Recommended contour level	0.4	Depositor
Map size (Å)	271.104, 271.104, 271.104	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.059, 1.059, 1.059	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: 6OU, 4DY, LBN

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	0.34	1/3822 (0.0%)	0.52	5/5206 (0.1%)
1	B	0.34	1/3822 (0.0%)	0.52	5/5206 (0.1%)
1	C	0.34	1/3822 (0.0%)	0.52	5/5206 (0.1%)
1	D	0.34	1/3822 (0.0%)	0.52	5/5206 (0.1%)
All	All	0.34	4/15288 (0.0%)	0.52	20/20824 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	677	MET	C-O	-5.64	1.17	1.24
1	A	677	MET	C-O	-5.58	1.17	1.24
1	B	677	MET	C-O	-5.54	1.17	1.24
1	D	677	MET	C-O	-5.52	1.17	1.24

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	677	MET	N-CA-C	-8.70	100.99	111.69
1	B	677	MET	N-CA-C	-8.69	101.00	111.69
1	A	677	MET	N-CA-C	-8.69	101.00	111.69
1	C	677	MET	N-CA-C	-8.66	101.03	111.69
1	C	557	ARG	CA-C-N	6.75	127.42	120.00

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	3733	0	3512	64	0
1	B	3733	0	3512	66	0
1	C	3733	0	3512	65	0
1	D	3733	0	3512	65	0
2	A	50	0	0	3	0
2	B	50	0	0	3	0
2	C	50	0	0	3	0
2	D	50	0	0	3	0
3	A	110	0	0	0	0
3	B	86	0	0	0	0
3	C	86	0	0	0	0
3	D	62	0	0	0	0
4	A	22	0	0	0	0
4	B	22	0	0	0	0
4	C	22	0	0	0	0
4	D	22	0	0	0	0
All	All	15564	0	14048	198	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 7.

The worst 5 of 198 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:C:554:TYR:O	1:C:557:ARG:HG2	1.91	0.71
1:B:554:TYR:O	1:B:557:ARG:HG2	1.91	0.70
1:D:554:TYR:O	1:D:557:ARG:HG2	1.91	0.70
1:A:554:TYR:O	1:A:557:ARG:HG2	1.91	0.69
1:C:583:VAL:HG21	1:D:565:TYR:CD2	2.30	0.67

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	489/868 (56%)	476 (97%)	13 (3%)	0	100	100
1	B	489/868 (56%)	476 (97%)	13 (3%)	0	100	100
1	C	489/868 (56%)	476 (97%)	13 (3%)	0	100	100
1	D	489/868 (56%)	476 (97%)	13 (3%)	0	100	100
All	All	1956/3472 (56%)	1904 (97%)	52 (3%)	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	355/770 (46%)	355 (100%)	0	100	100
1	B	355/770 (46%)	355 (100%)	0	100	100
1	C	355/770 (46%)	355 (100%)	0	100	100
1	D	355/770 (46%)	355 (100%)	0	100	100
All	All	1420/3080 (46%)	1420 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	364	HIS
1	B	364	HIS
1	C	364	HIS
1	D	358	HIS
1	D	364	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

24 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	6OU	D	902	-	30,30,48	1.84	5 (16%)	33,35,53	1.02	2 (6%)
4	4DY	A	904	-	22,22,22	1.41	3 (13%)	27,27,27	1.09	1 (3%)
4	4DY	D	903	-	22,22,22	1.41	3 (13%)	27,27,27	1.09	1 (3%)
2	LBN	C	906	-	16,16,51	1.12	2 (12%)	18,18,59	1.00	0
4	4DY	C	904	-	22,22,22	1.41	3 (13%)	27,27,27	1.09	1 (3%)
3	6OU	B	902	-	30,30,48	1.84	5 (16%)	33,35,53	1.02	2 (6%)
2	LBN	D	901	-	32,32,51	1.18	2 (6%)	38,40,59	1.07	2 (5%)
2	LBN	A	906	-	16,16,51	1.13	2 (12%)	18,18,59	0.99	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	6OU	C	905	-	30,30,48	1.70	5 (16%)	33,35,53	0.93	2 (6%)
2	LBN	C	901	-	32,32,51	1.18	2 (6%)	38,40,59	1.07	2 (5%)
3	6OU	A	902	-	30,30,48	1.84	6 (20%)	33,35,53	1.02	2 (6%)
3	6OU	A	905	-	30,30,48	1.69	5 (16%)	33,35,53	0.93	2 (6%)
3	6OU	B	903	-	23,23,48	1.99	4 (17%)	26,28,53	1.03	2 (7%)
3	6OU	A	907	-	23,23,48	1.99	3 (13%)	26,28,53	1.02	2 (7%)
3	6OU	C	902	-	30,30,48	1.84	5 (16%)	33,35,53	1.02	2 (6%)
3	6OU	D	904	-	30,30,48	1.69	5 (16%)	33,35,53	0.93	2 (6%)
3	6OU	C	903	-	23,23,48	1.99	4 (17%)	26,28,53	1.03	2 (7%)
3	6OU	B	905	-	30,30,48	1.69	5 (16%)	33,35,53	0.93	2 (6%)
4	4DY	B	904	-	22,22,22	1.41	3 (13%)	27,27,27	1.09	1 (3%)
2	LBN	B	906	-	16,16,51	1.13	2 (12%)	18,18,59	1.00	0
3	6OU	A	903	-	23,23,48	1.99	3 (13%)	26,28,53	1.02	2 (7%)
2	LBN	B	901	-	32,32,51	1.18	2 (6%)	38,40,59	1.07	2 (5%)
2	LBN	A	901	-	32,32,51	1.18	2 (6%)	38,40,59	1.07	2 (5%)
2	LBN	D	905	-	16,16,51	1.13	2 (12%)	18,18,59	0.99	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	6OU	D	902	-	-	10/32/32/52	-
4	4DY	A	904	-	-	1/16/16/16	0/1/1/1
4	4DY	D	903	-	-	1/16/16/16	0/1/1/1
2	LBN	C	906	-	-	12/17/17/55	-
4	4DY	C	904	-	-	1/16/16/16	0/1/1/1
3	6OU	B	902	-	-	10/32/32/52	-
2	LBN	D	901	-	-	15/36/36/55	-
2	LBN	A	906	-	-	12/17/17/55	-
3	6OU	C	905	-	-	17/32/32/52	-
2	LBN	C	901	-	-	15/36/36/55	-
3	6OU	A	902	-	-	10/32/32/52	-
3	6OU	A	905	-	-	17/32/32/52	-
3	6OU	B	903	-	-	14/25/25/52	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	6OU	A	907	-	-	14/25/25/52	-
3	6OU	C	902	-	-	10/32/32/52	-
3	6OU	D	904	-	-	17/32/32/52	-
3	6OU	C	903	-	-	14/25/25/52	-
3	6OU	B	905	-	-	17/32/32/52	-
4	4DY	B	904	-	-	1/16/16/16	0/1/1/1
2	LBN	B	906	-	-	12/17/17/55	-
3	6OU	A	903	-	-	14/25/25/52	-
2	LBN	B	901	-	-	15/36/36/55	-
2	LBN	A	901	-	-	15/36/36/55	-
2	LBN	D	905	-	-	12/17/17/55	-

The worst 5 of 83 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	902	6OU	P23-O22	5.39	1.77	1.60
3	D	902	6OU	P23-O22	5.39	1.77	1.60
3	B	902	6OU	P23-O22	5.39	1.77	1.60
3	C	902	6OU	P23-O22	5.38	1.77	1.60
3	C	903	6OU	P23-O22	5.32	1.77	1.60

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	LBN	O3-P1-O4	3.48	128.64	112.44
2	D	901	LBN	O3-P1-O4	3.48	128.63	112.44
2	C	901	LBN	O3-P1-O4	3.48	128.61	112.44
2	A	901	LBN	O3-P1-O4	3.47	128.60	112.44
3	B	905	6OU	O25-P23-O24	3.45	124.29	110.83

There are no chirality outliers.

5 of 276 torsion outliers are listed below:

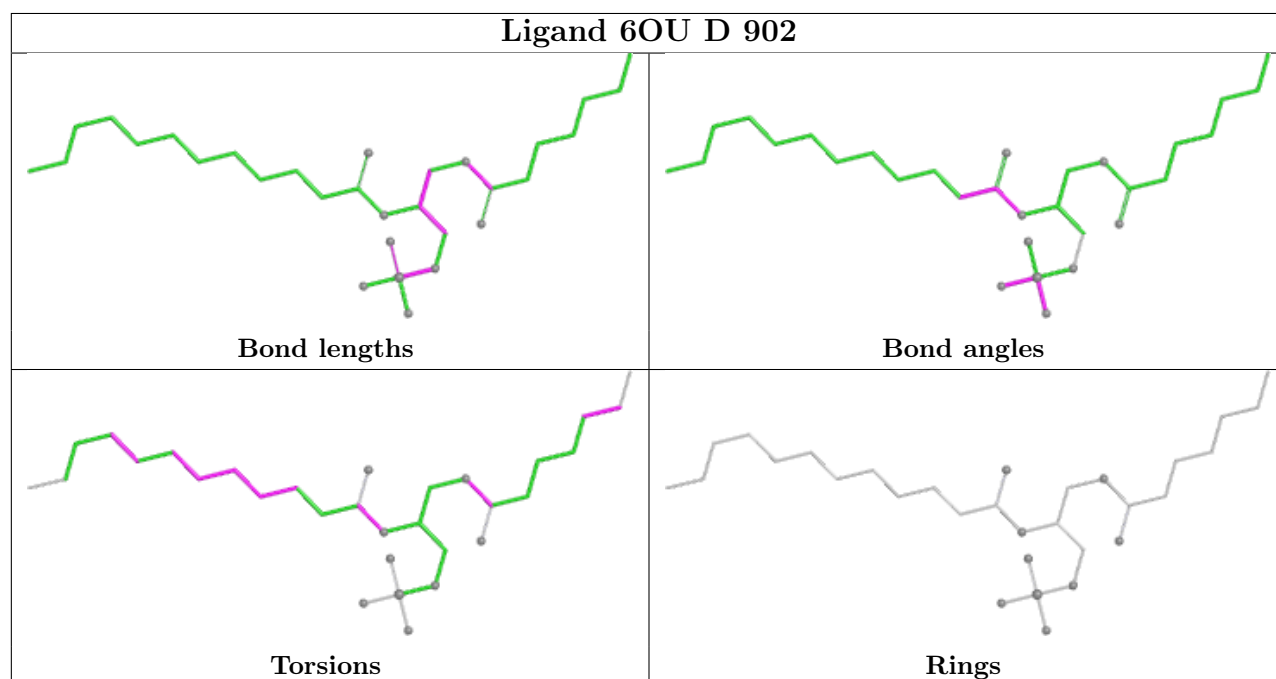
Mol	Chain	Res	Type	Atoms
2	A	901	LBN	C1-O1-P1-O3
2	A	901	LBN	N1-C6-C9-O2
2	A	906	LBN	O1-C1-C2-O7
2	B	901	LBN	C1-O1-P1-O3
2	B	901	LBN	N1-C6-C9-O2

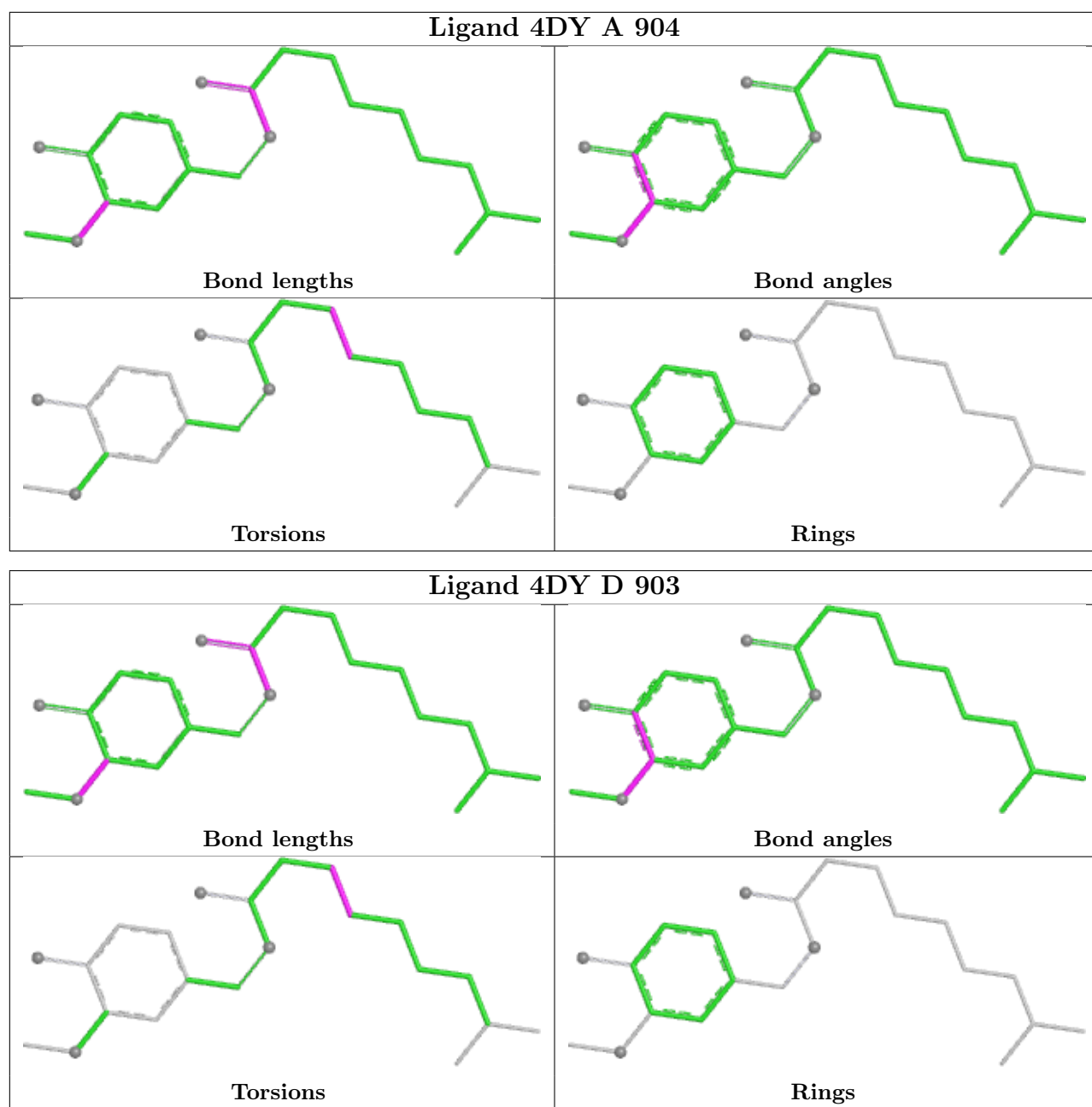
There are no ring outliers.

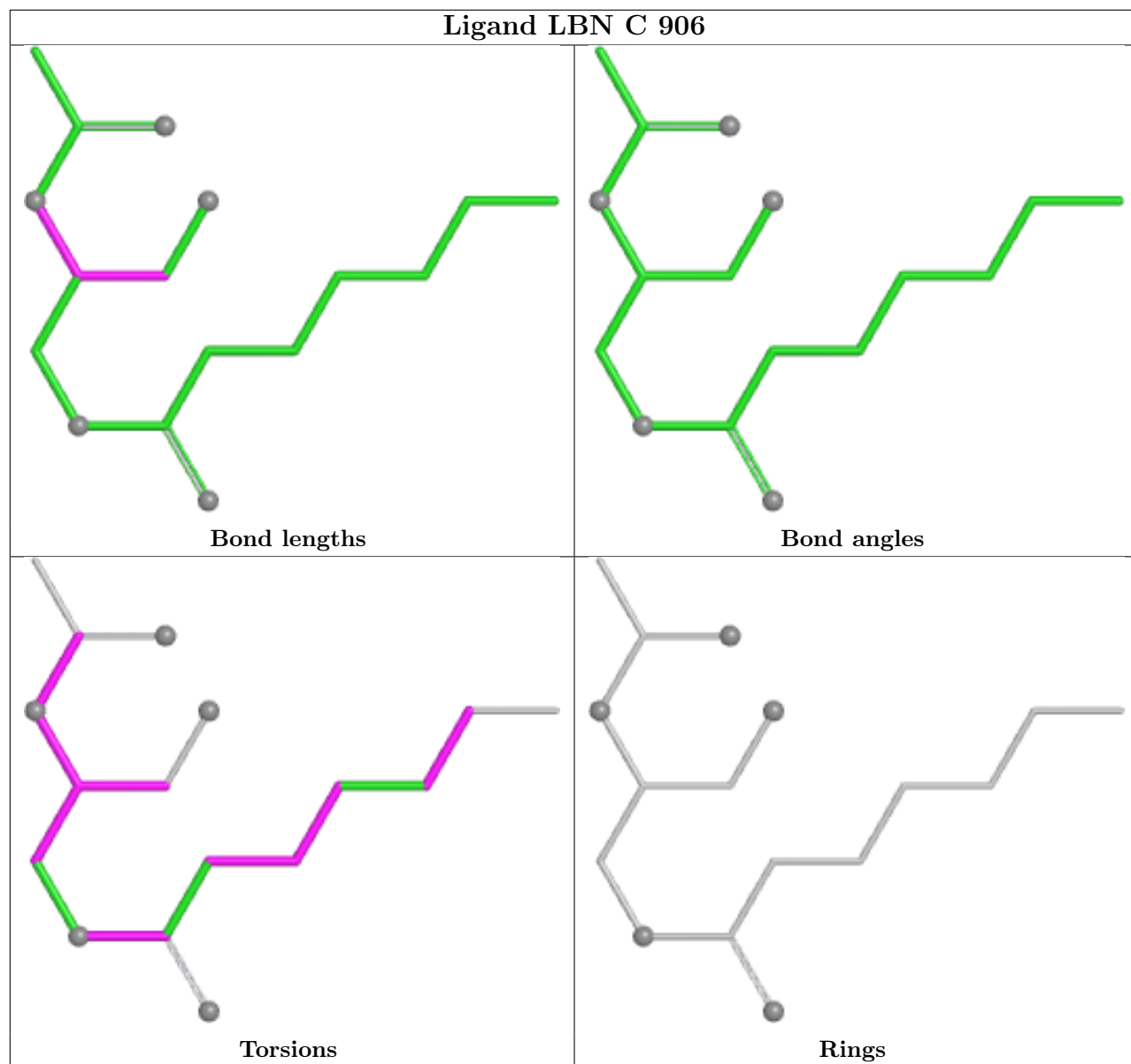
4 monomers are involved in 12 short contacts:

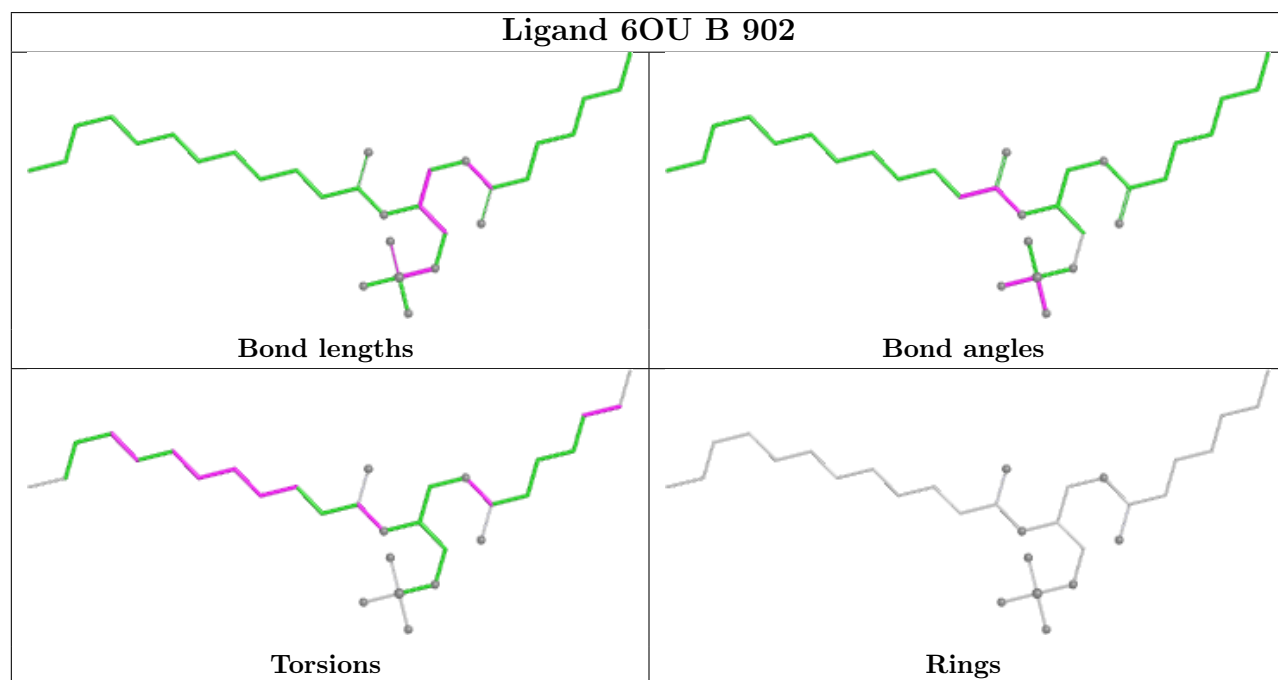
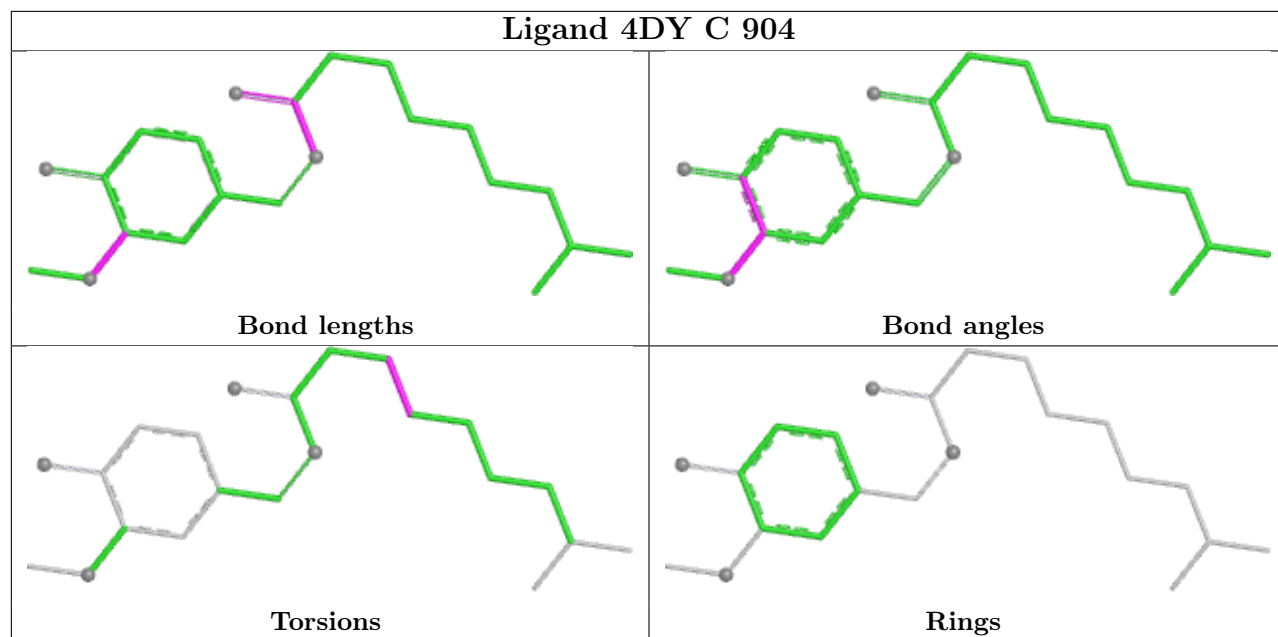
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	901	LBN	3	0
2	C	901	LBN	3	0
2	B	901	LBN	3	0
2	A	901	LBN	3	0

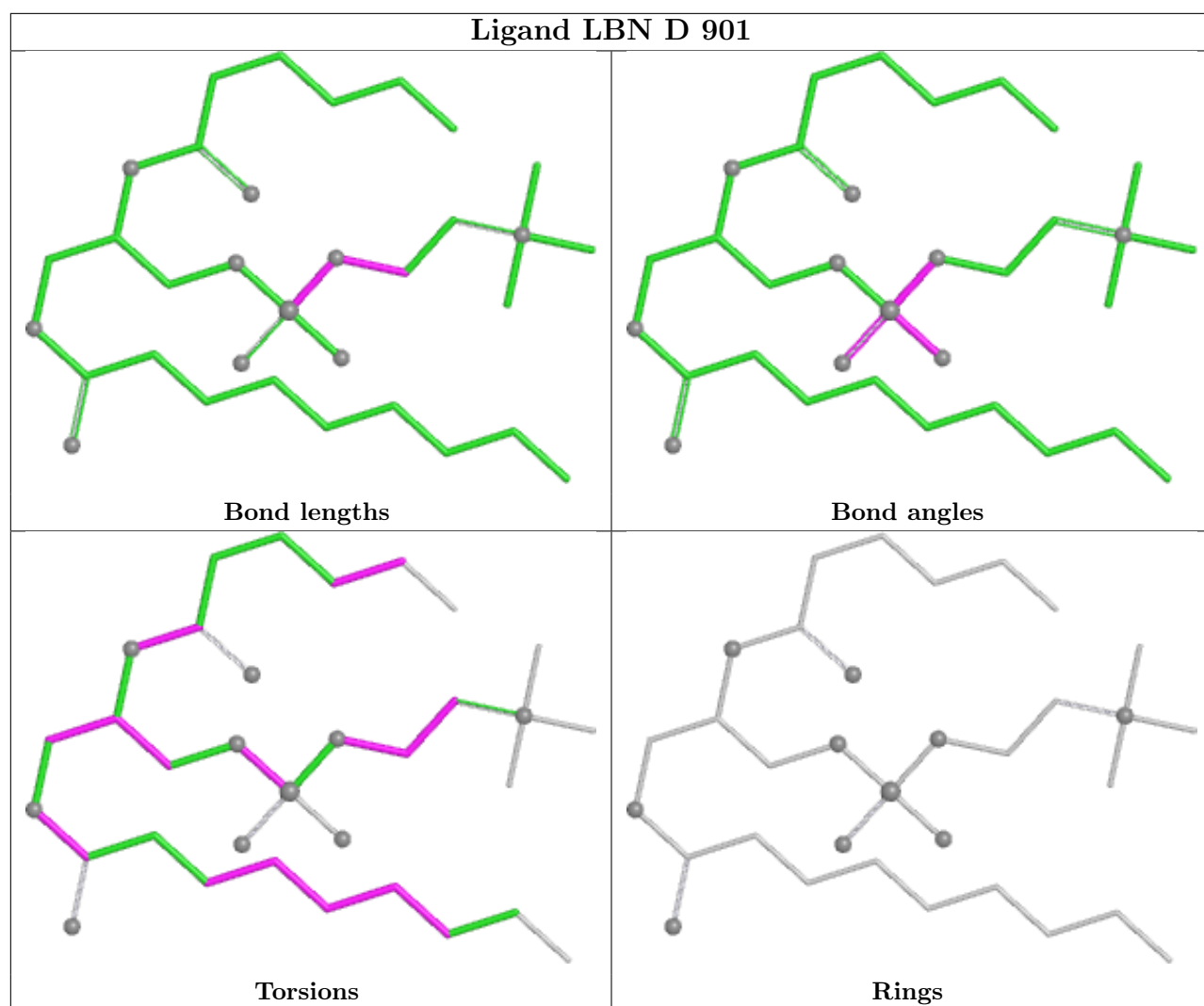
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

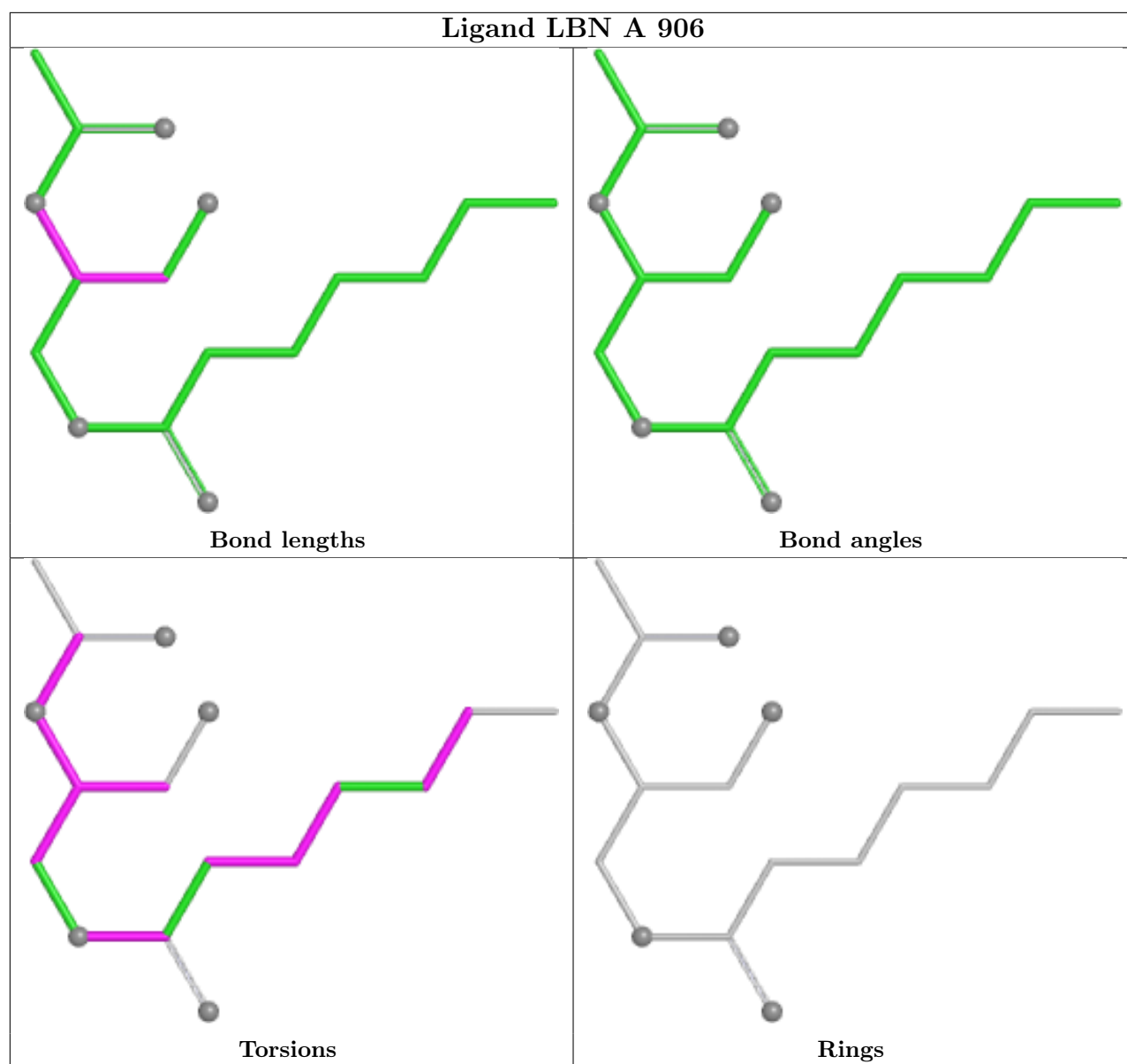


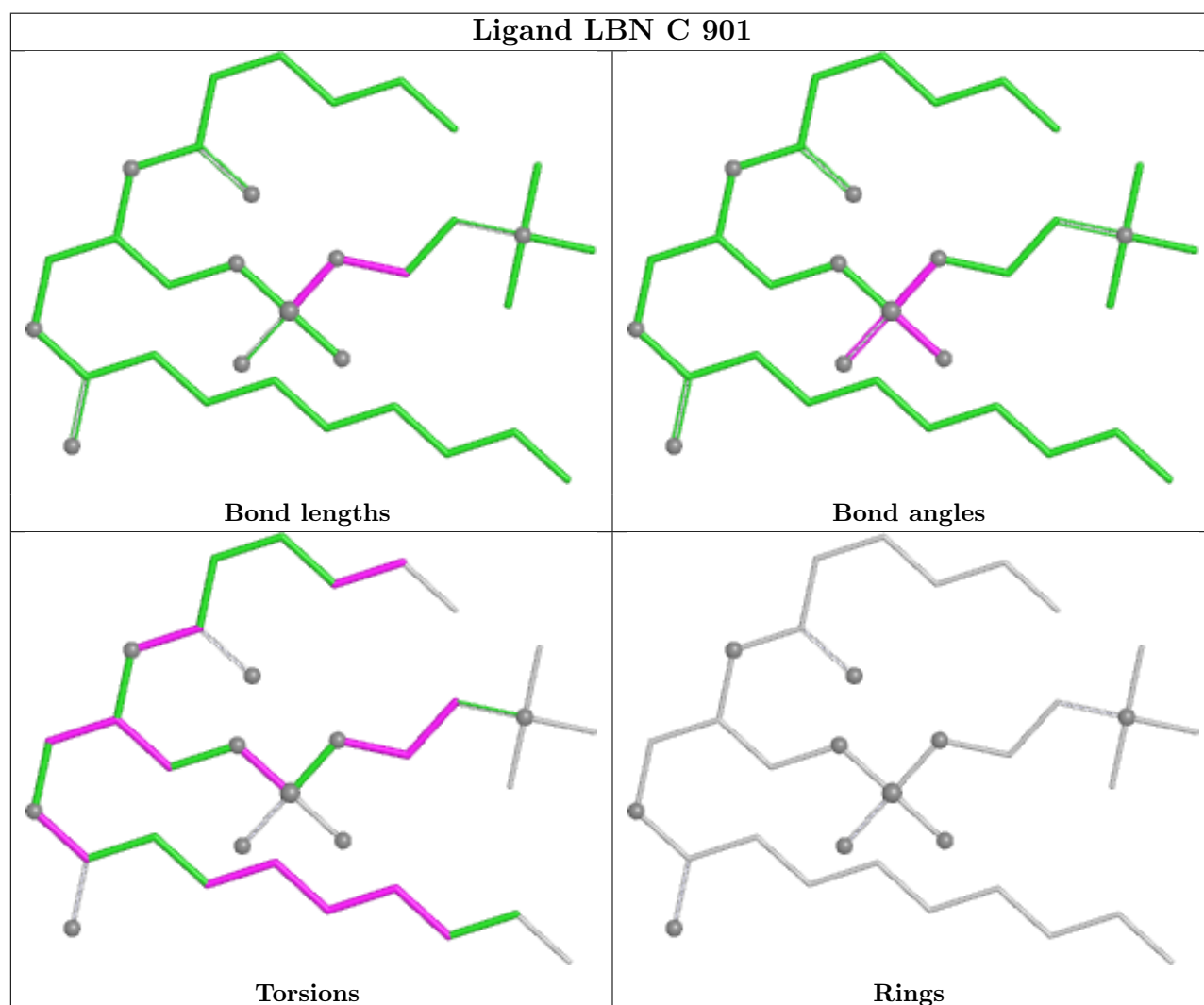
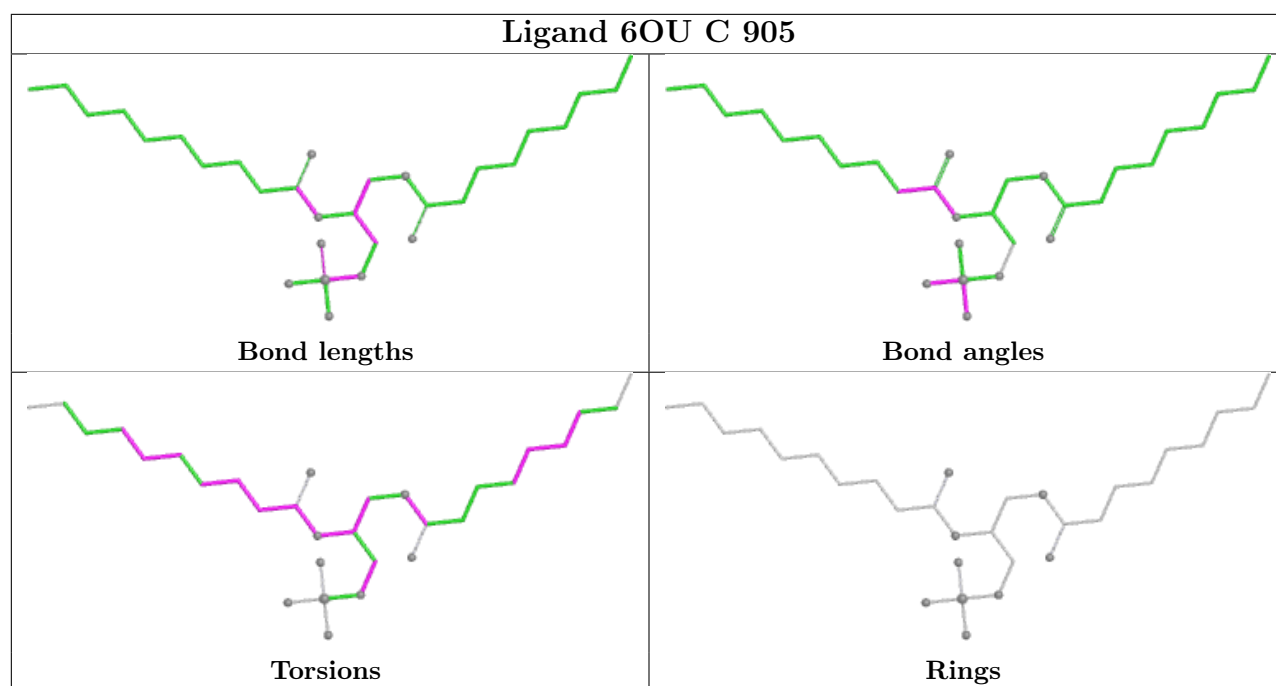


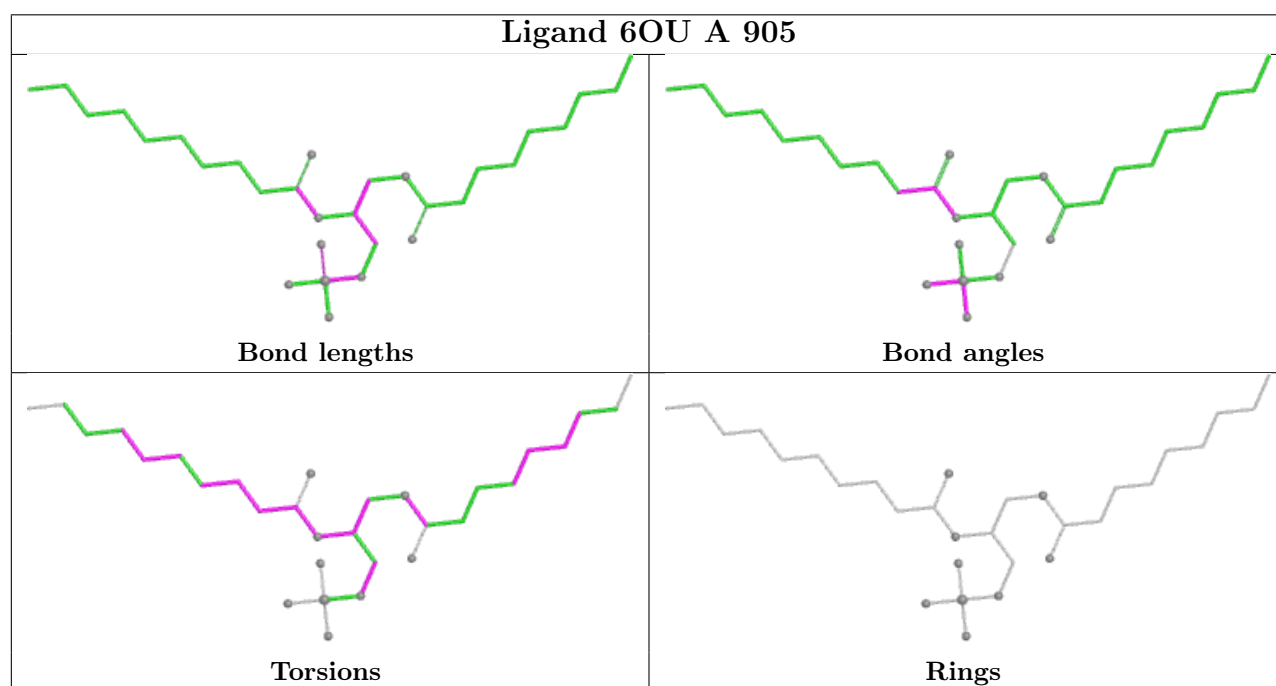
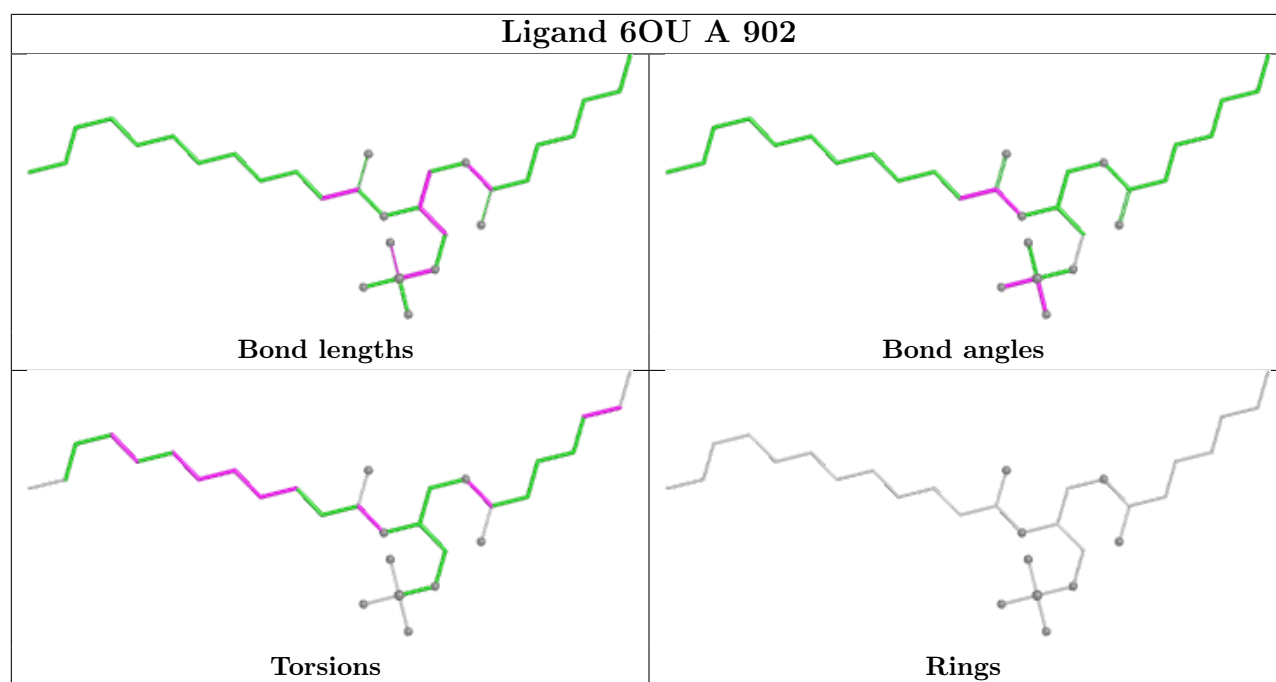


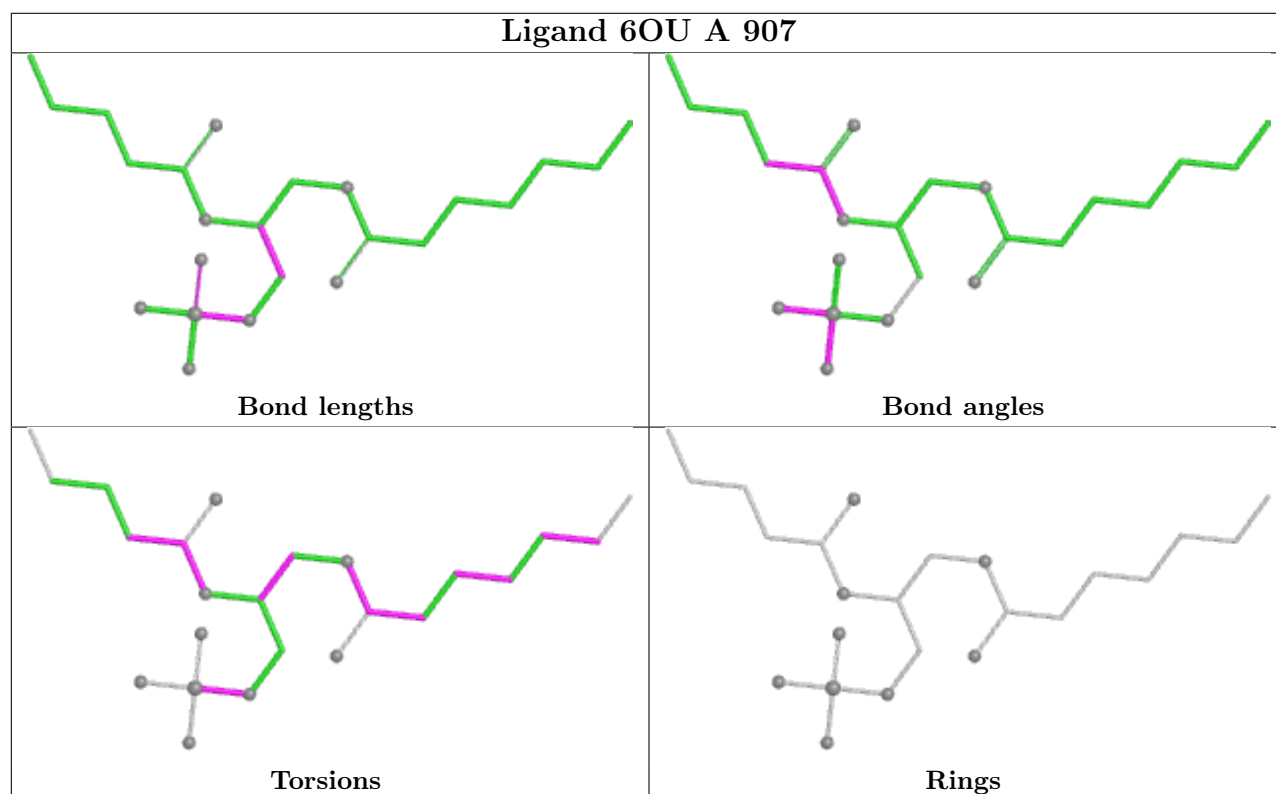
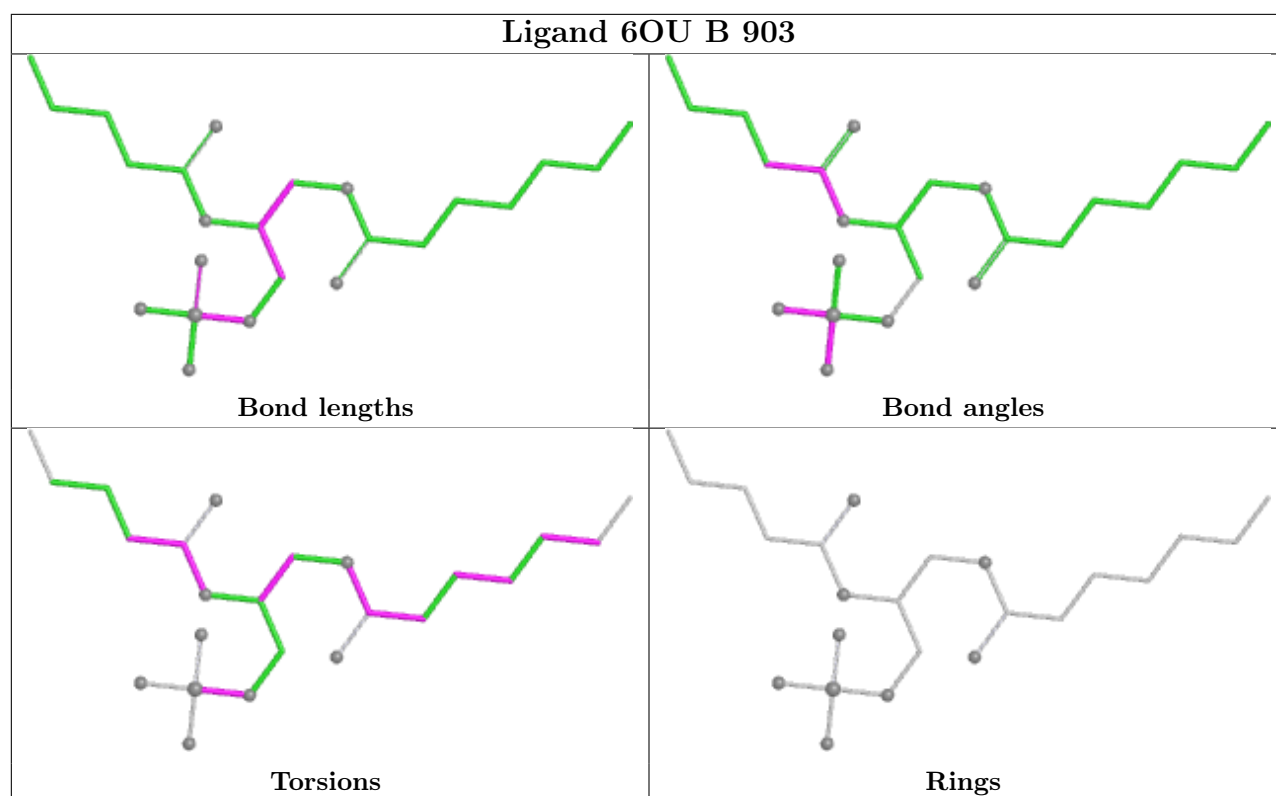


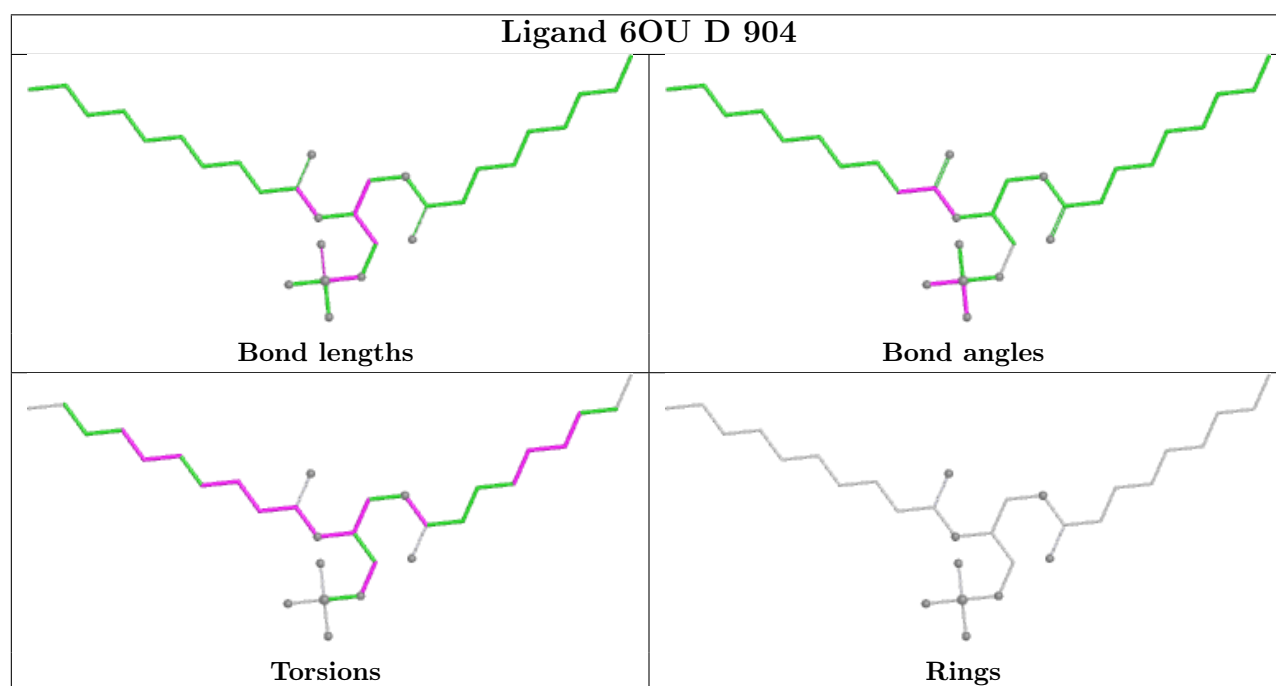
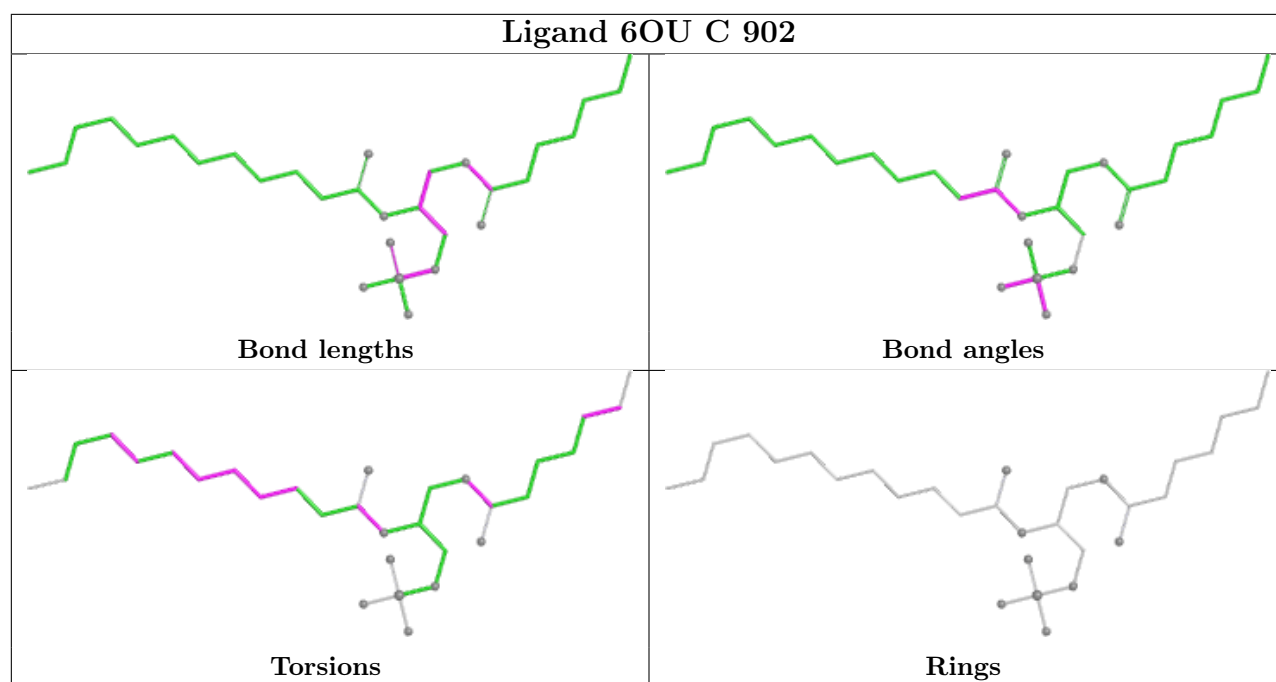


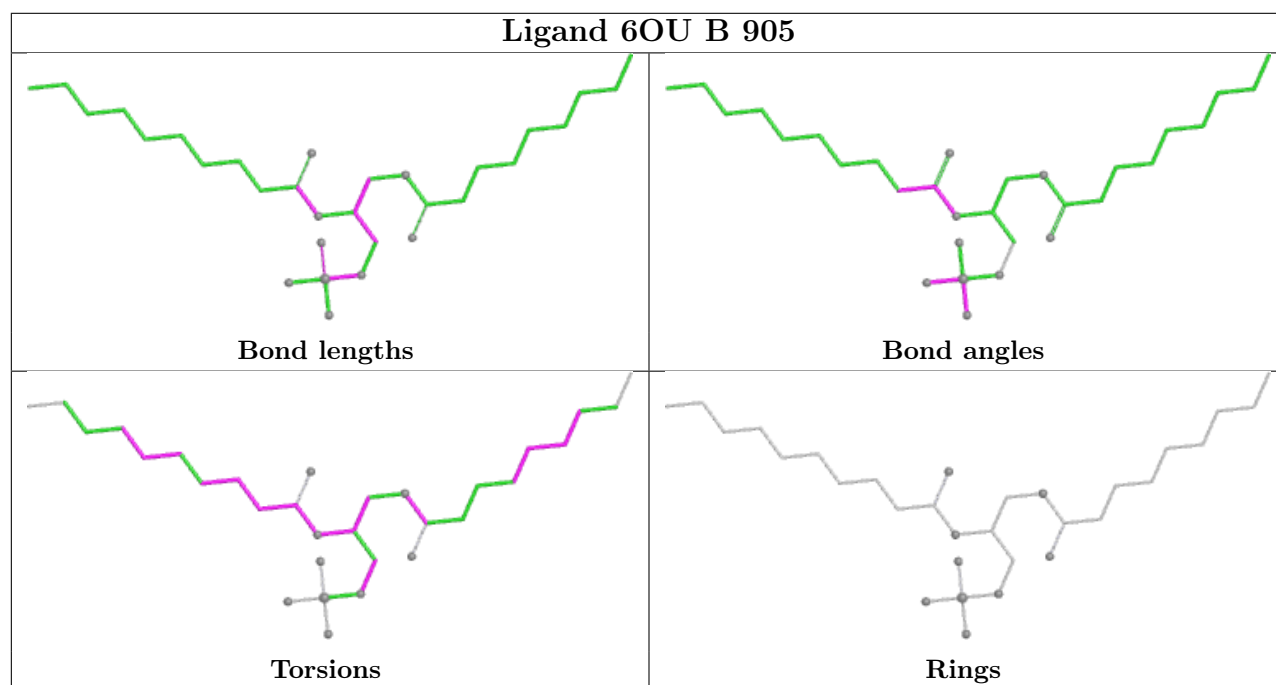
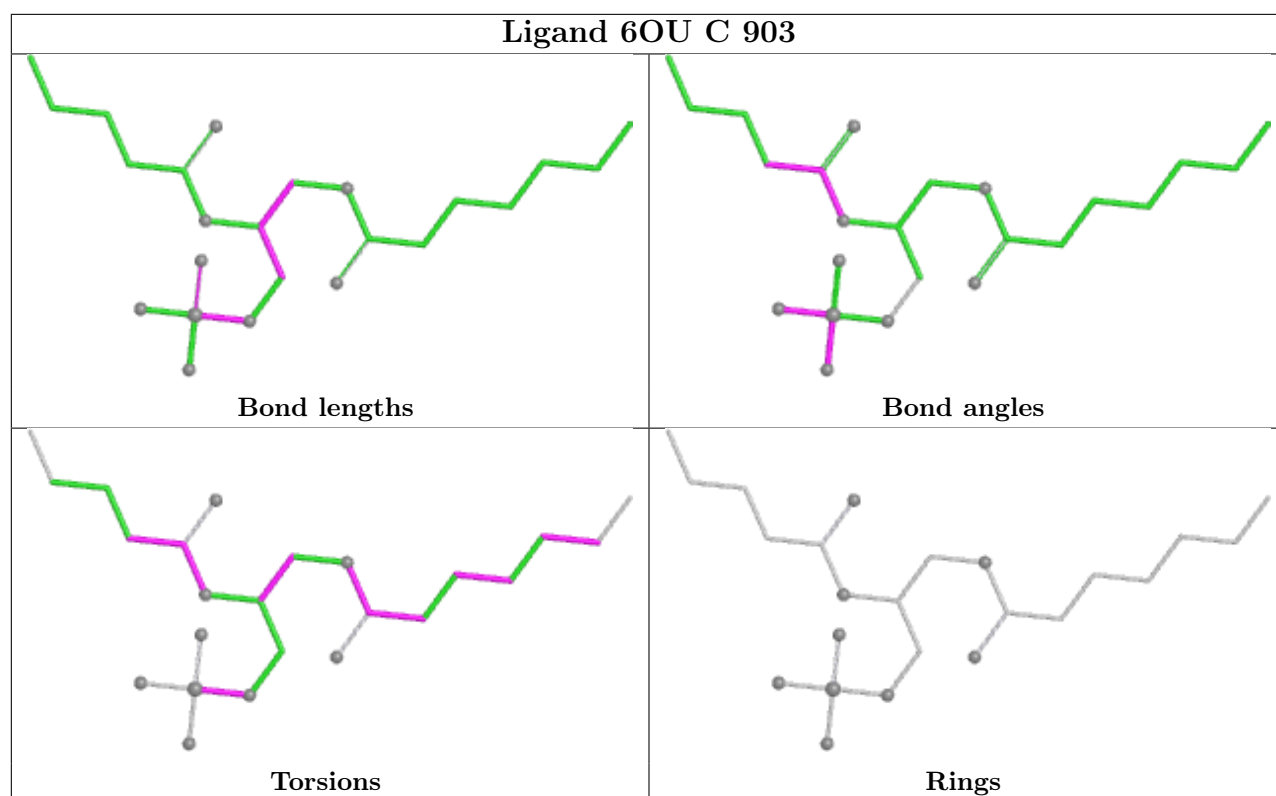


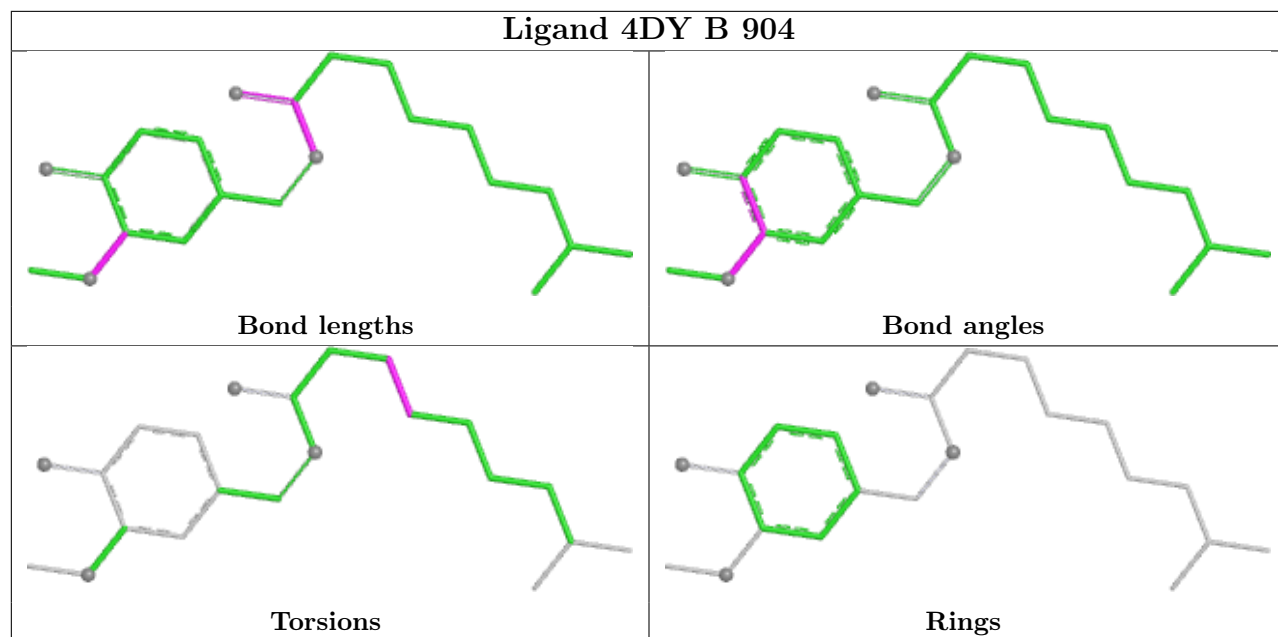


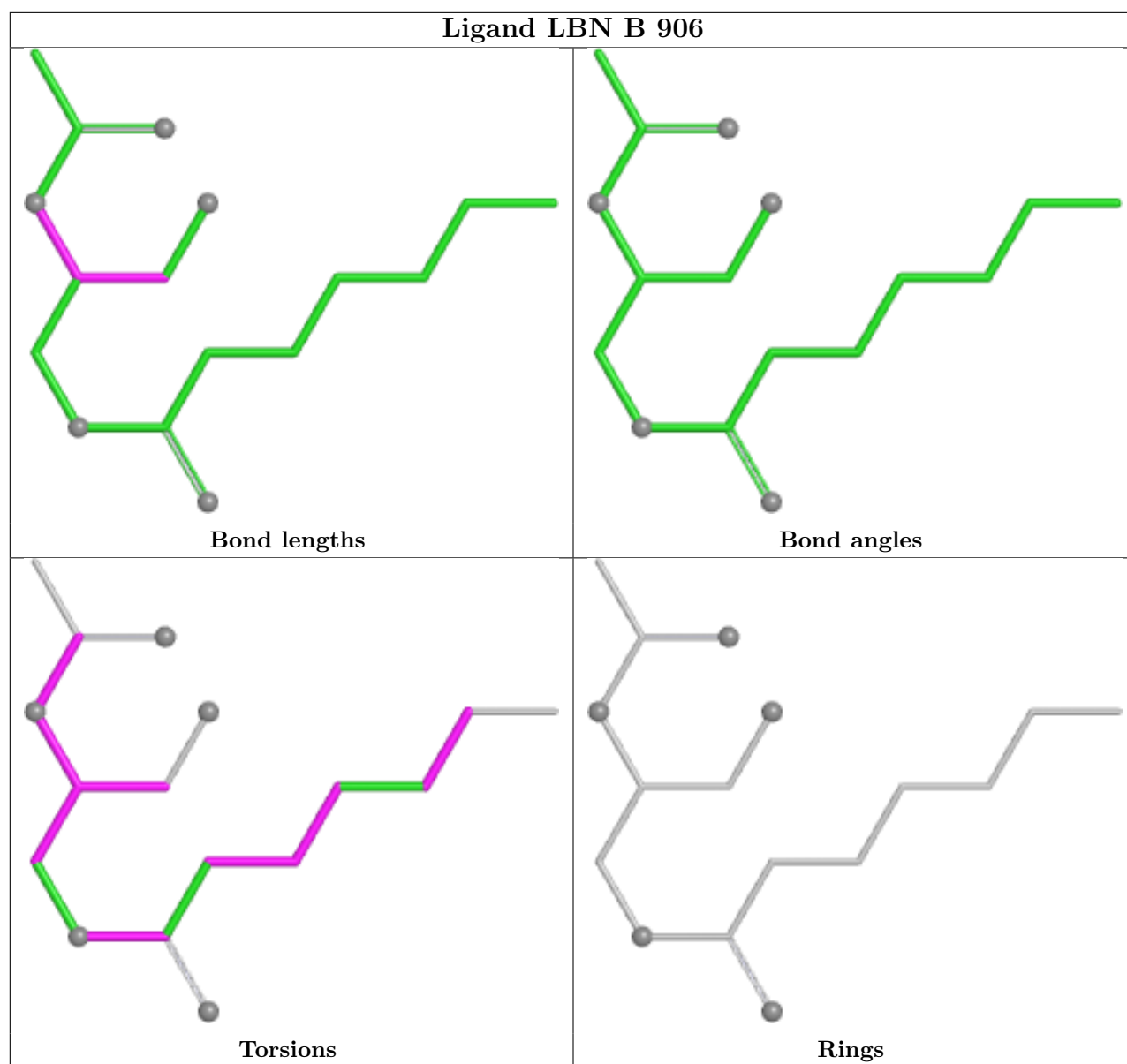


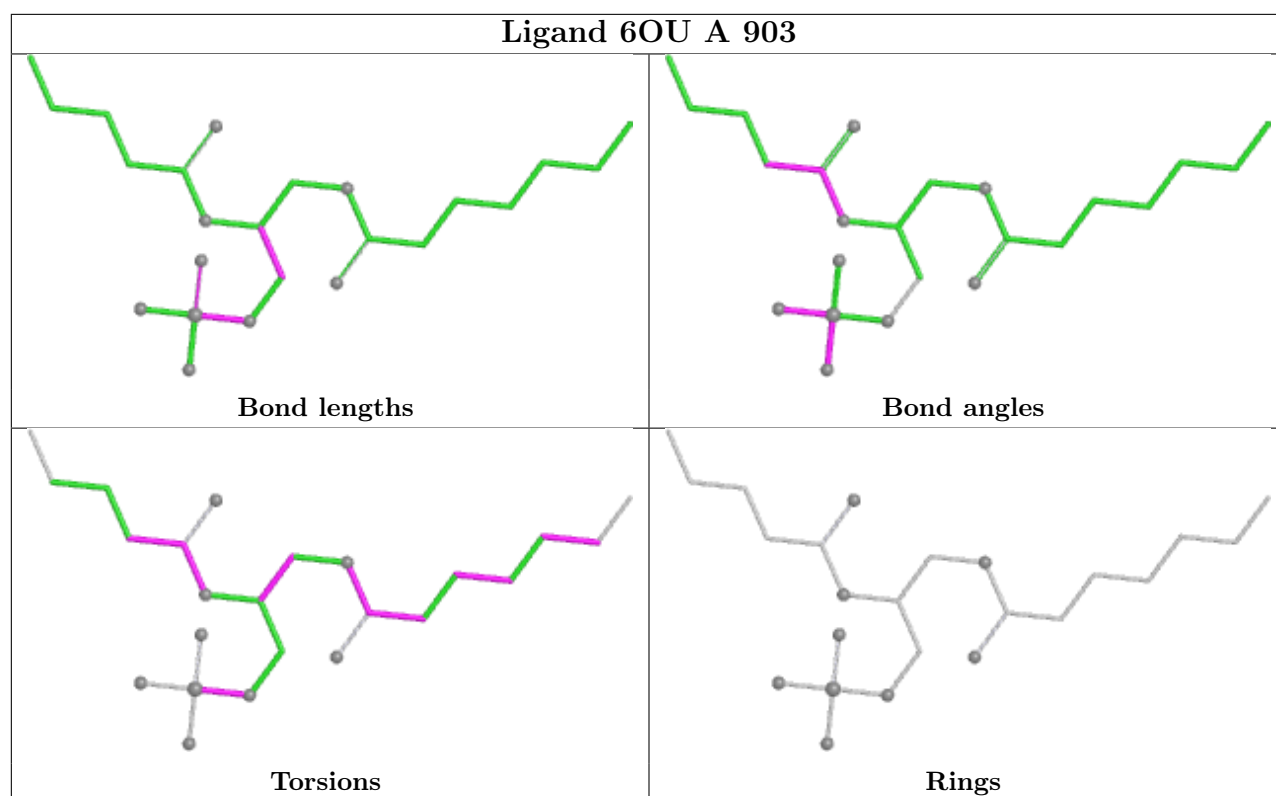


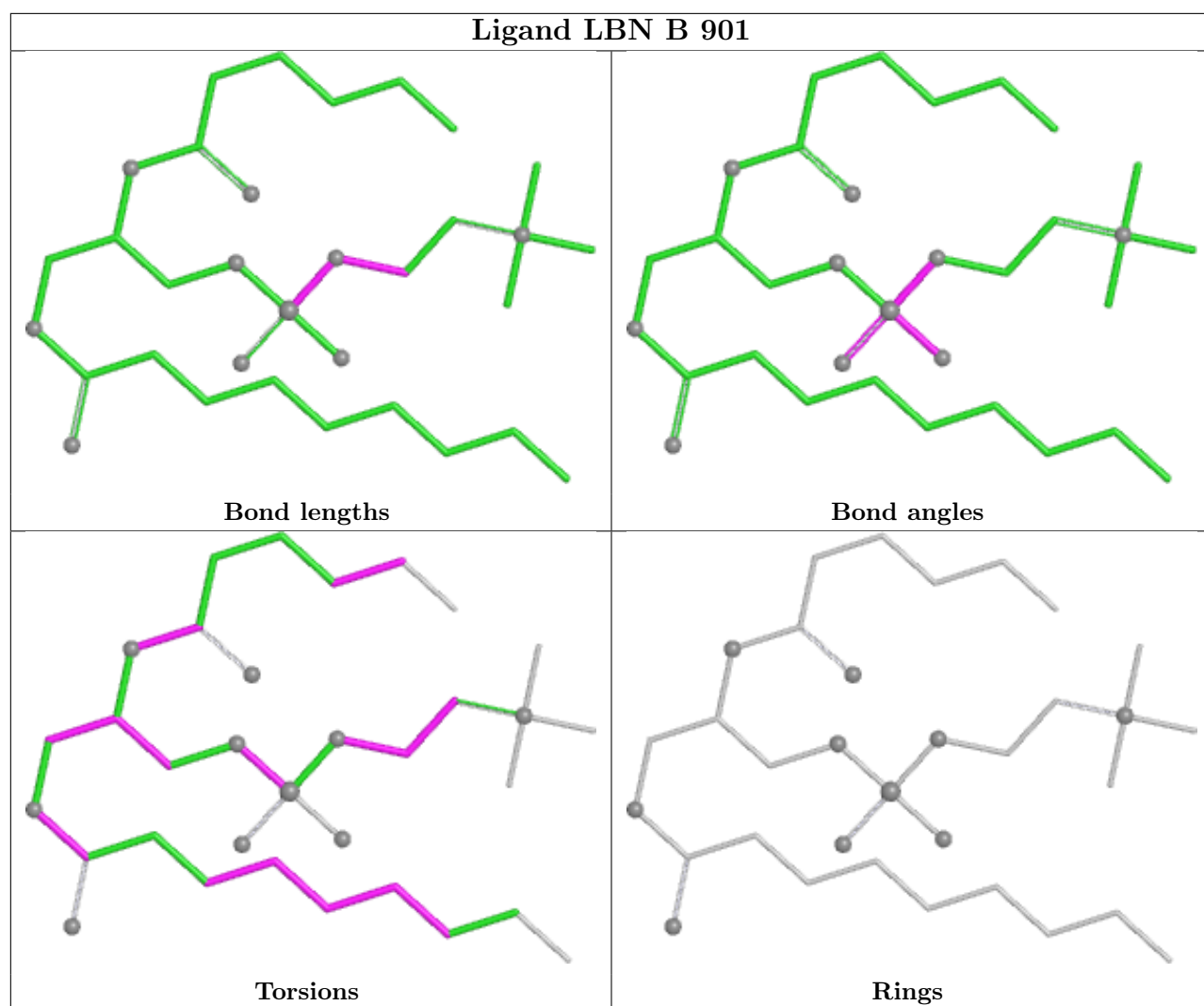


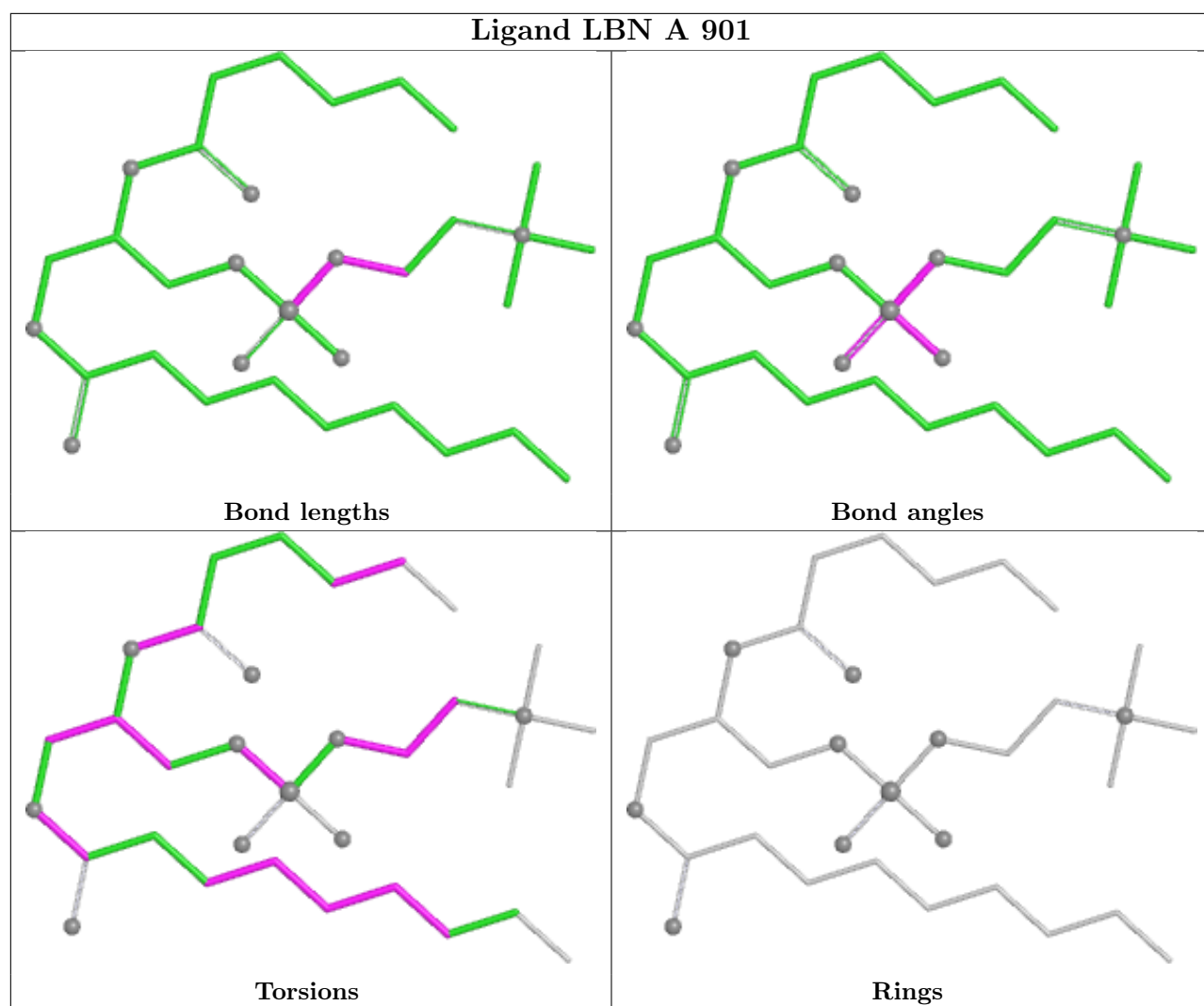


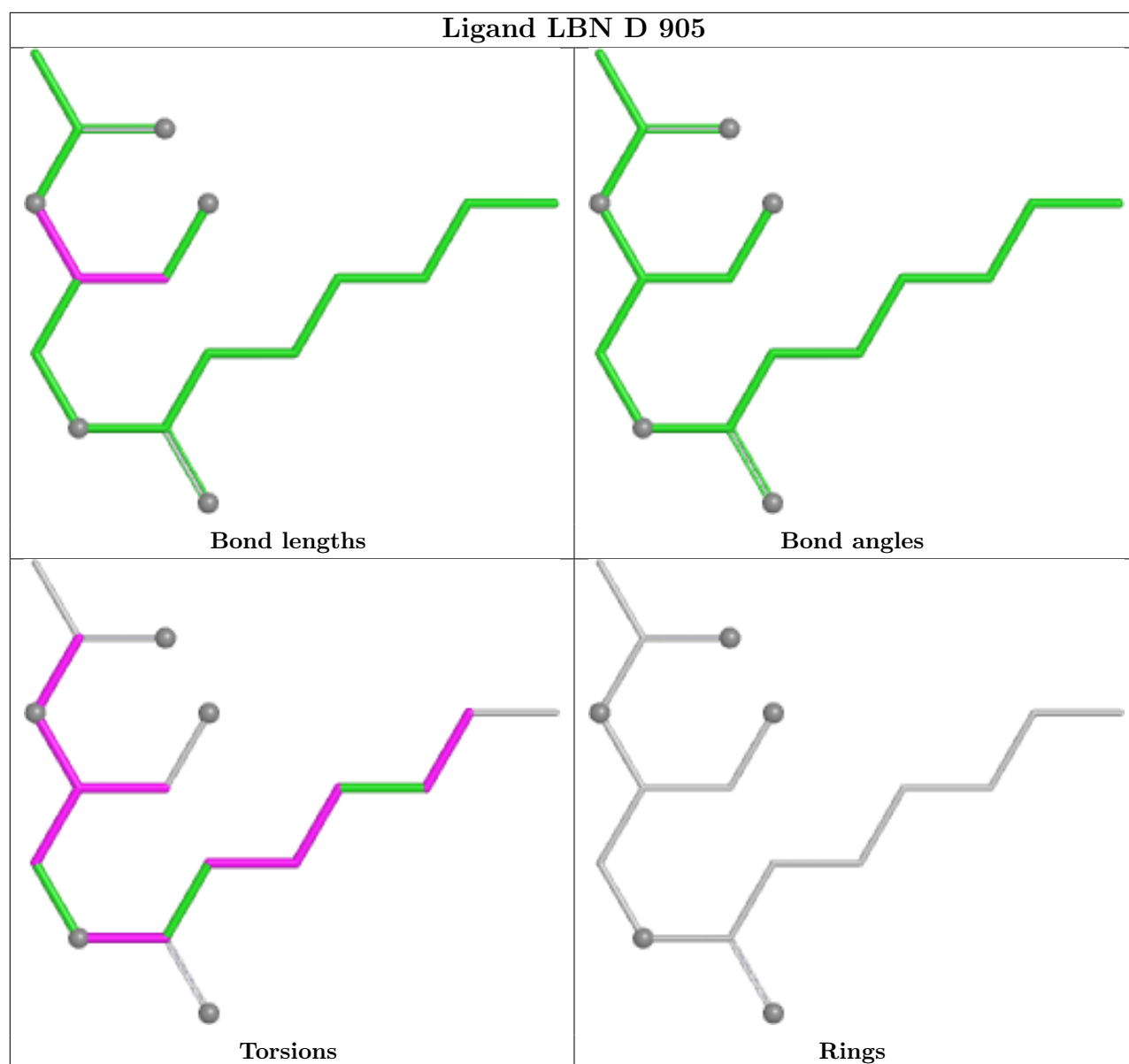












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

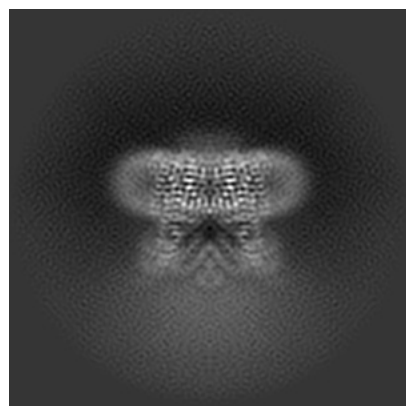
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23478. These allow visual inspection of the internal detail of the map and identification of artifacts.

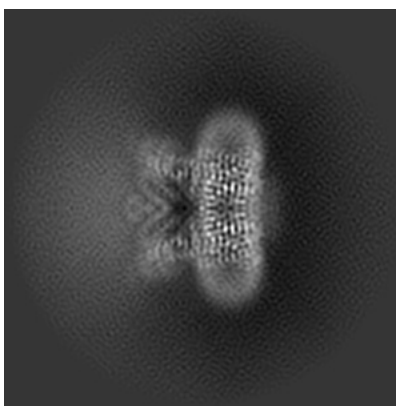
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

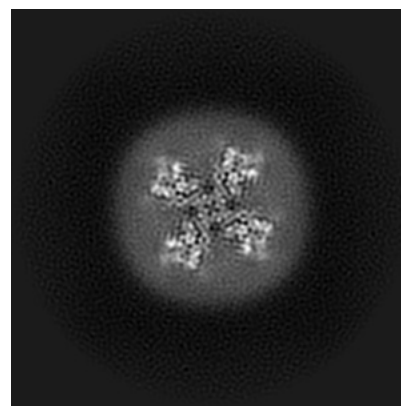
6.1.1 Primary map



X

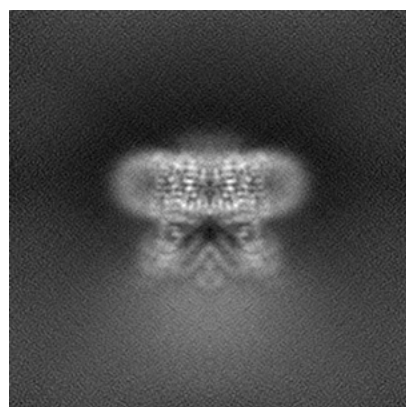


Y

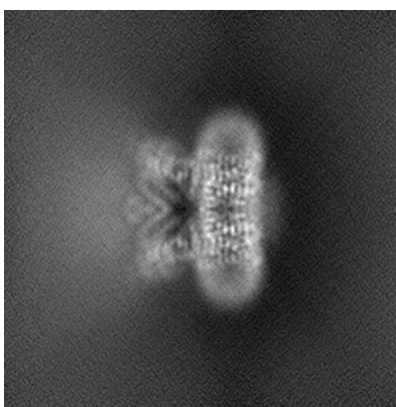


Z

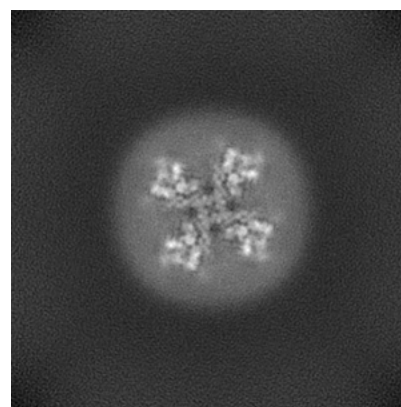
6.1.2 Raw map



X



Y

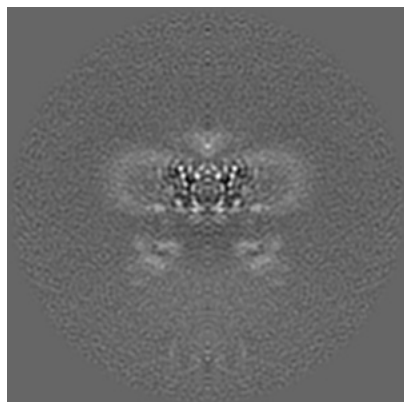


Z

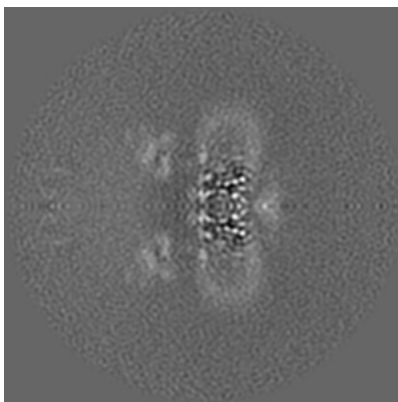
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

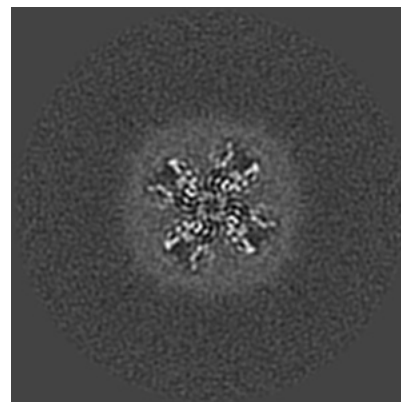
6.2.1 Primary map



X Index: 128

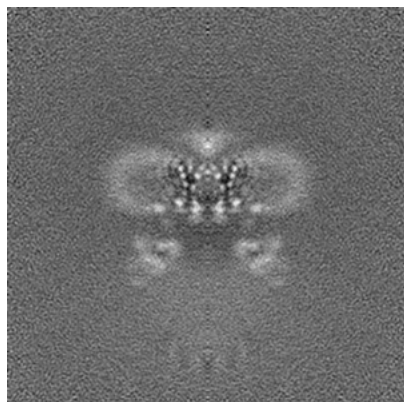


Y Index: 128

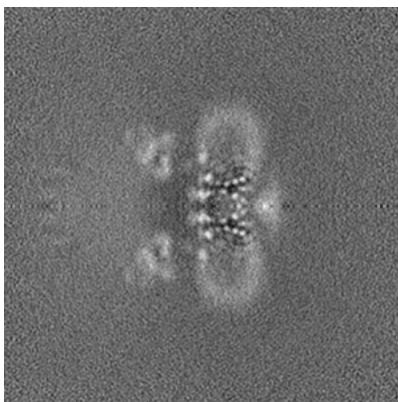


Z Index: 128

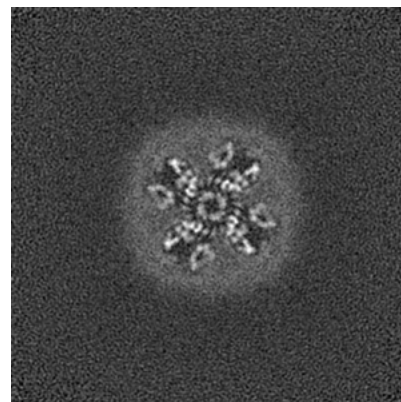
6.2.2 Raw map



X Index: 128



Y Index: 128

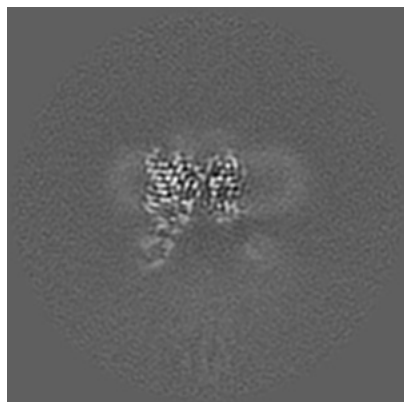


Z Index: 128

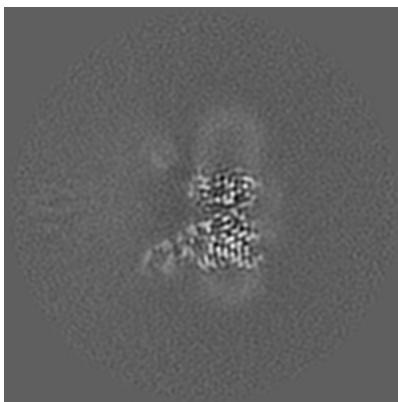
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

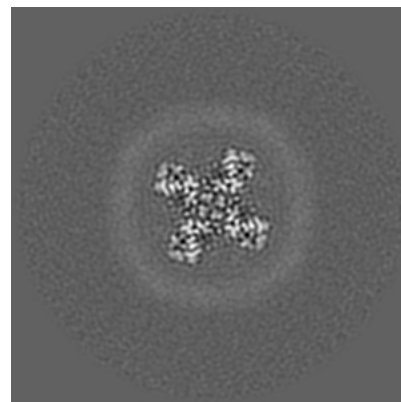
6.3.1 Primary map



X Index: 116

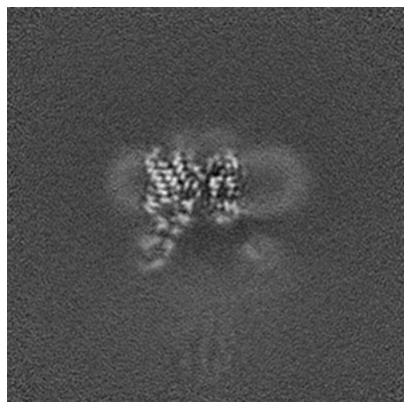


Y Index: 140

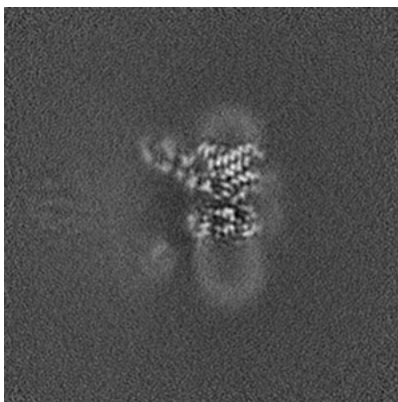


Z Index: 146

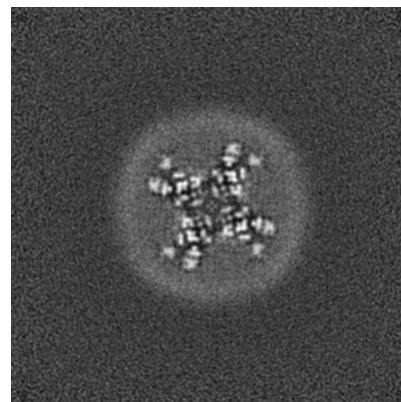
6.3.2 Raw map



X Index: 116



Y Index: 116

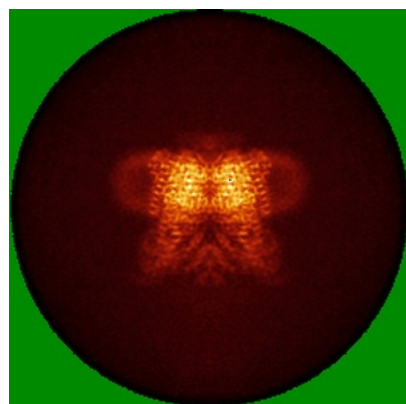


Z Index: 155

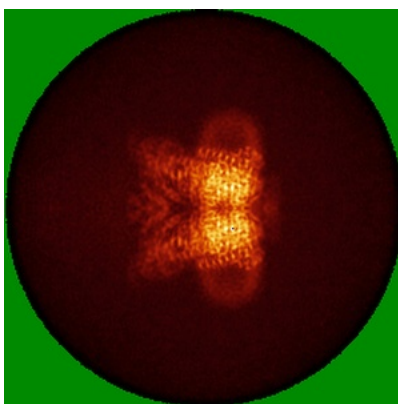
The images above show the largest variance slices of the map in three orthogonal directions.

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

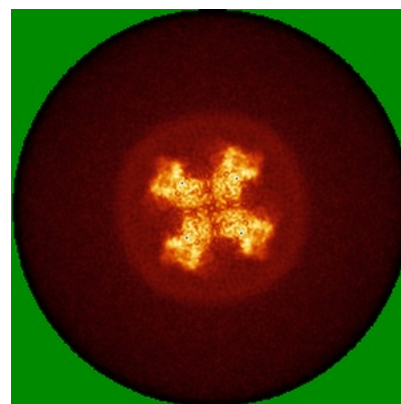
6.4.1 Primary map



X

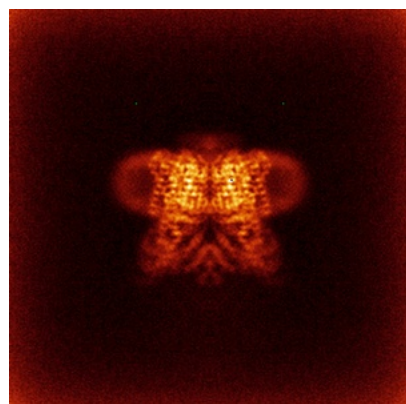


Y

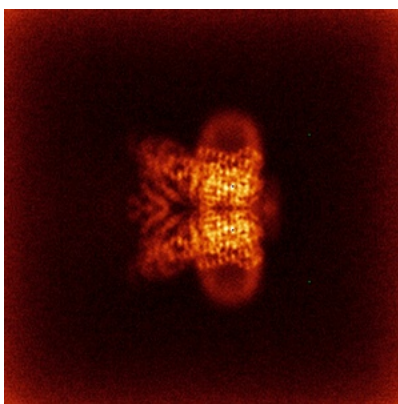


Z

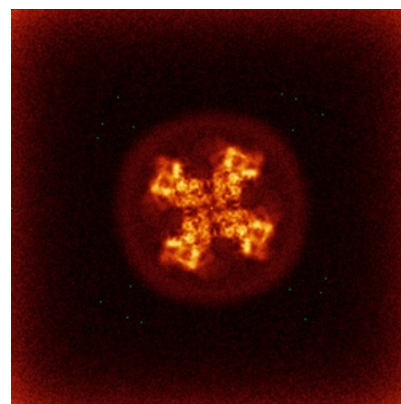
6.4.2 Raw map



X



Y

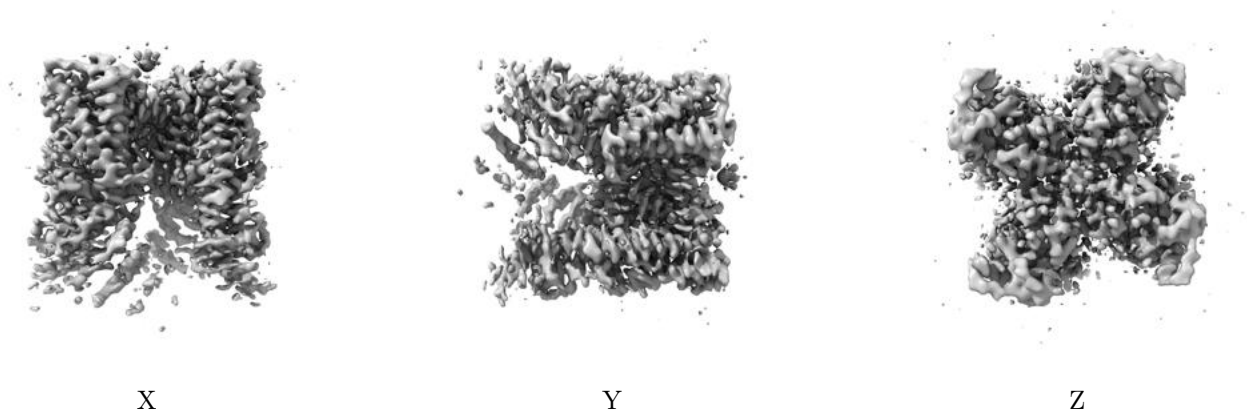


Z

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



The images above show the 3D surface view of the map at the recommended contour level 0.4. 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



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.

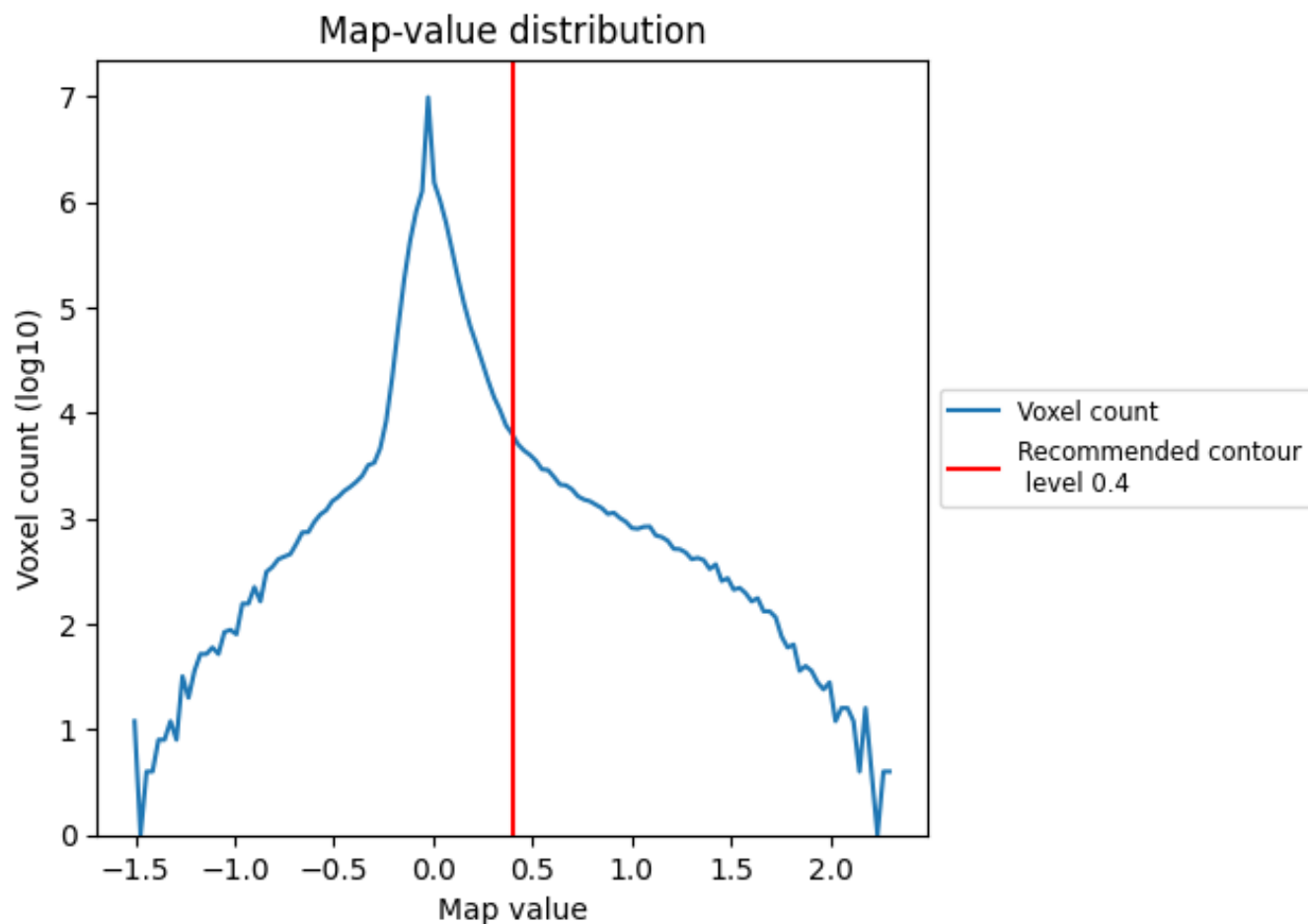
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

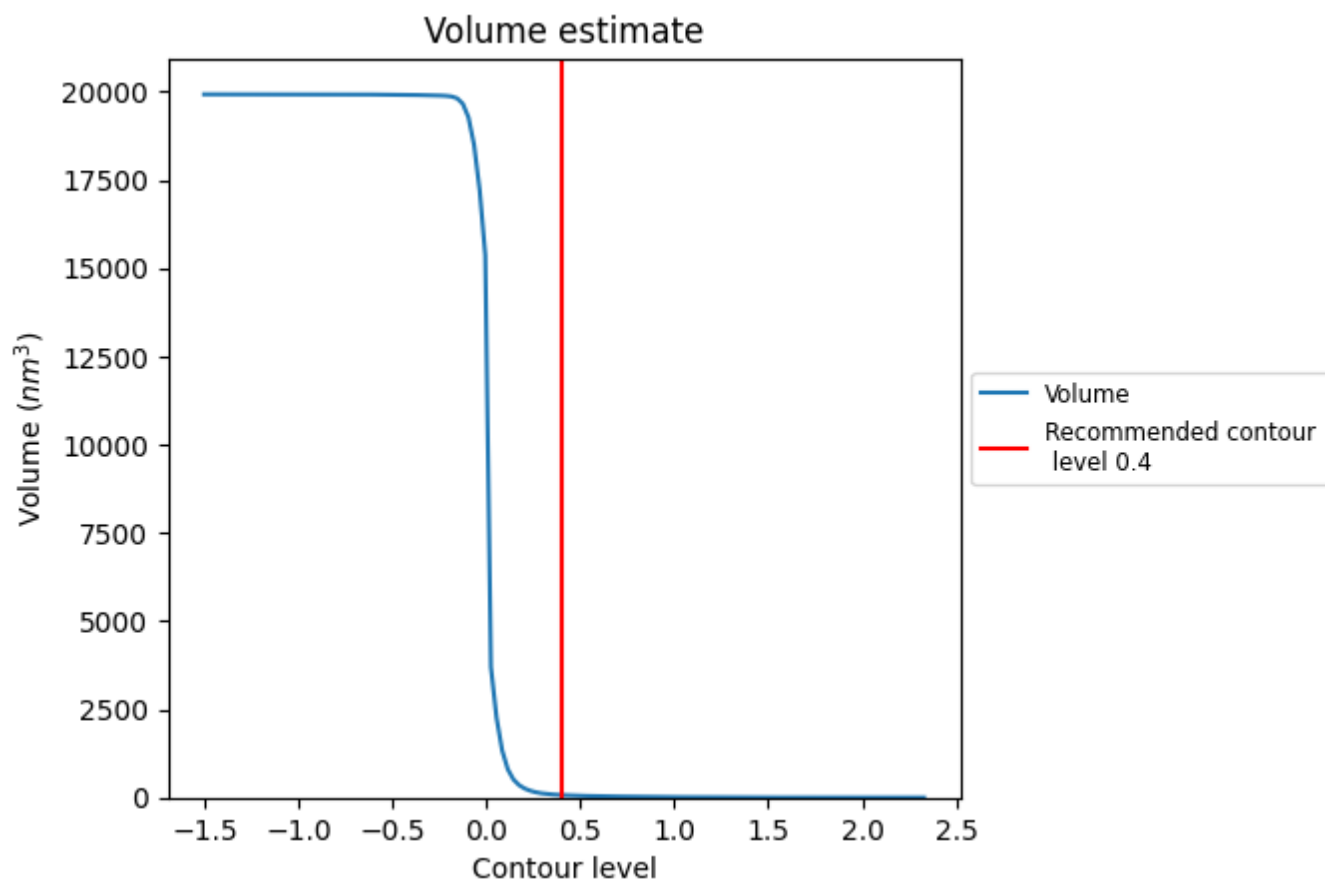
This section contains the results of statistical analysis of the map.

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.

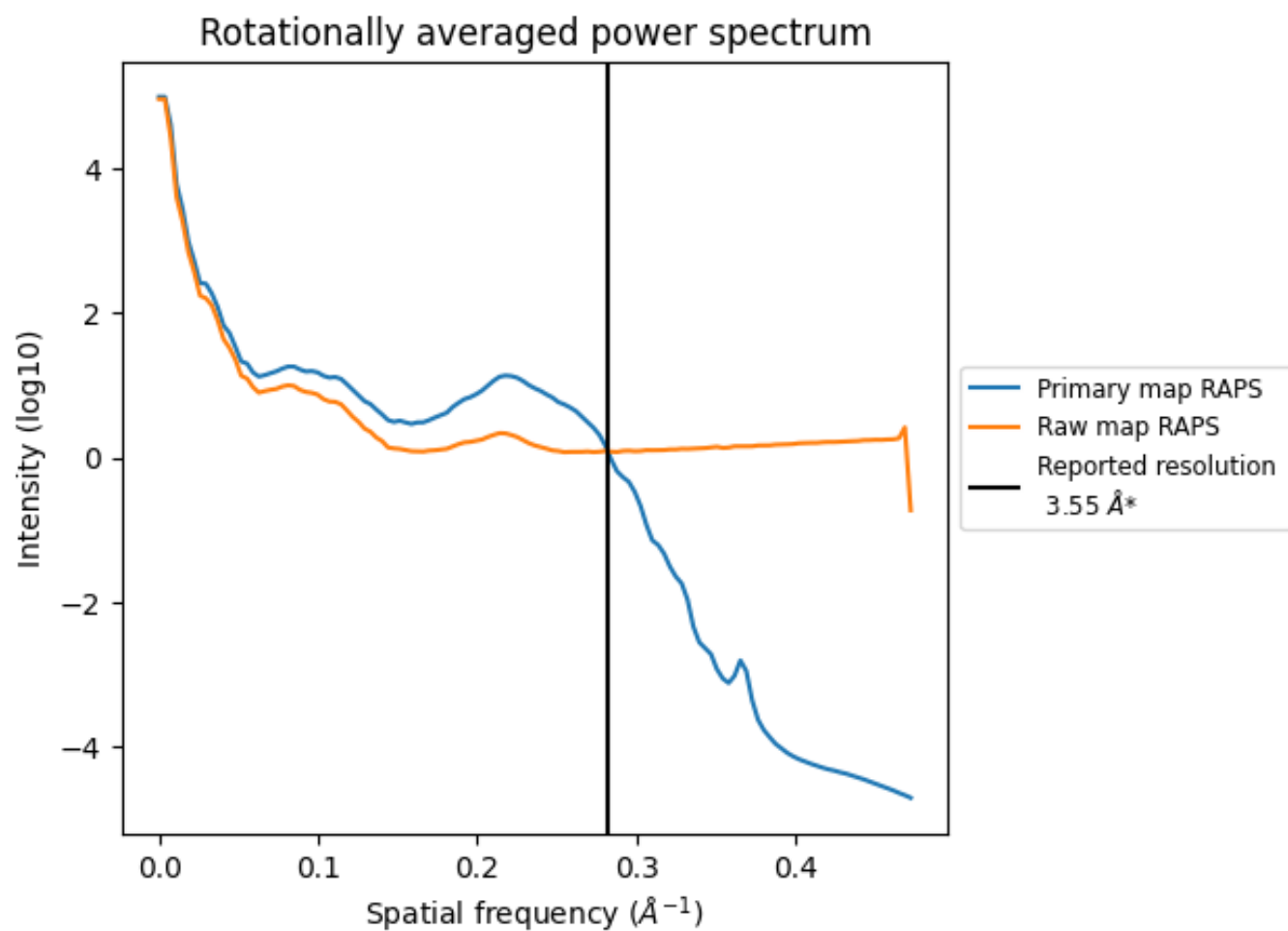
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 71 nm³; this corresponds to an approximate mass of 64 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 ⓘ

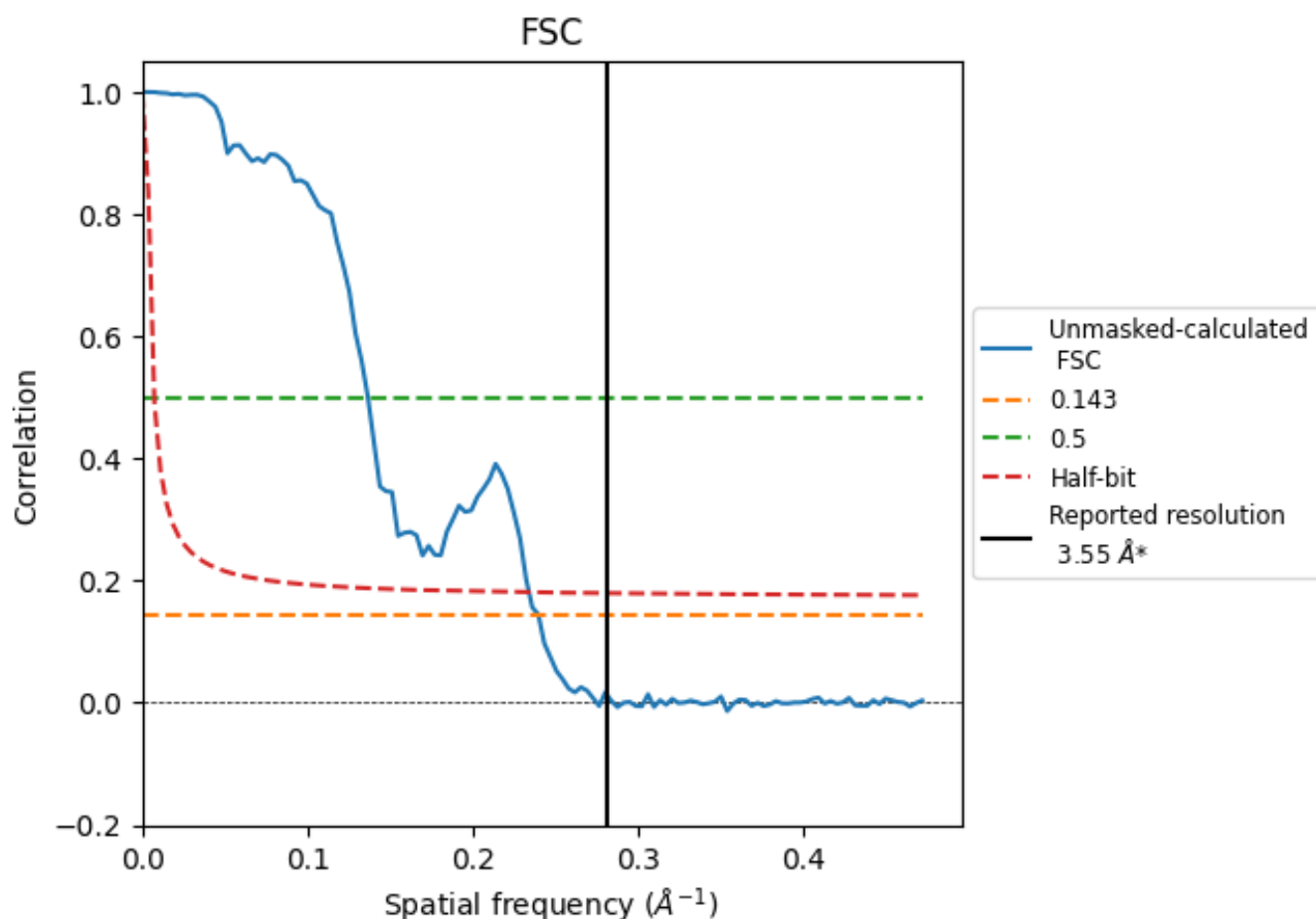


*Reported resolution corresponds to spatial frequency of 0.282 Å⁻¹

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 0.282 Å⁻¹

8.2 Resolution estimates [i](#)

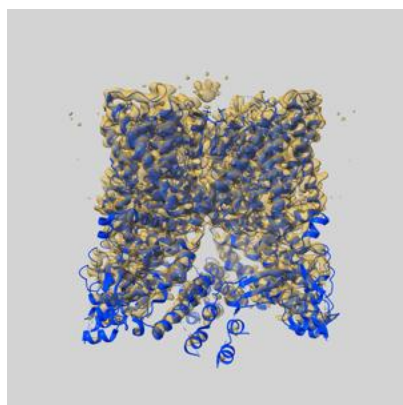
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.55	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.17	7.32	4.27

*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 4.17 differs from the reported value 3.55 by more than 10 %

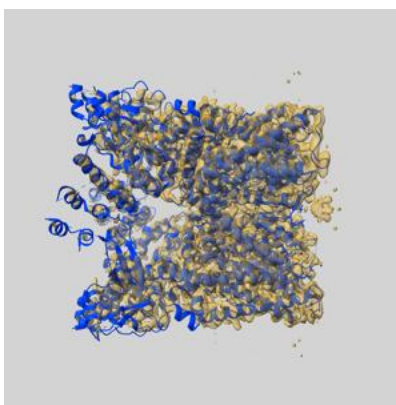
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23478 and PDB model 7LPD. Per-residue inclusion information can be found in section [3](#) on page [9](#).

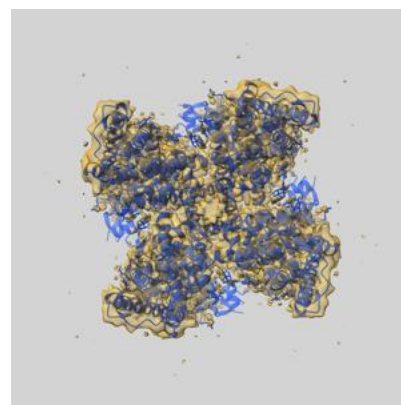
9.1 Map-model overlay [i](#)



X



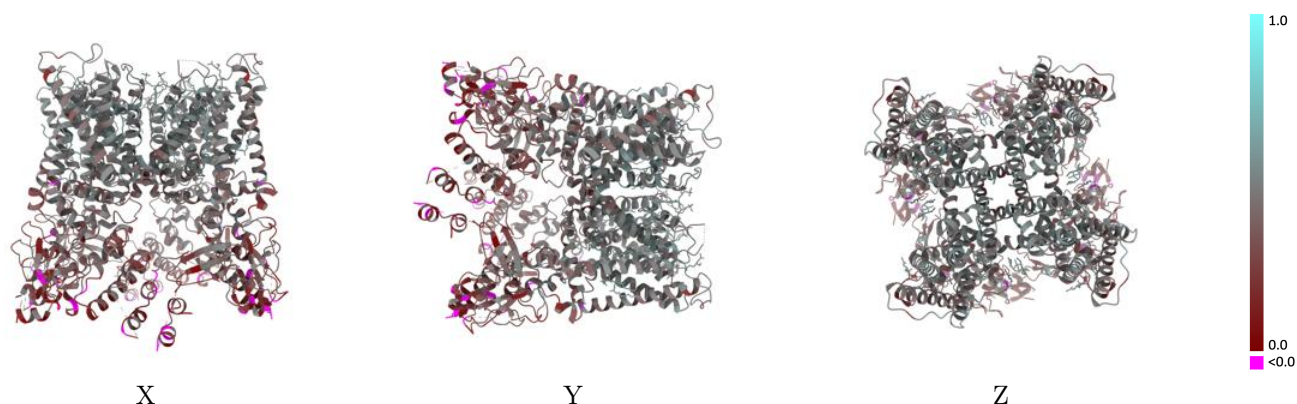
Y



Z

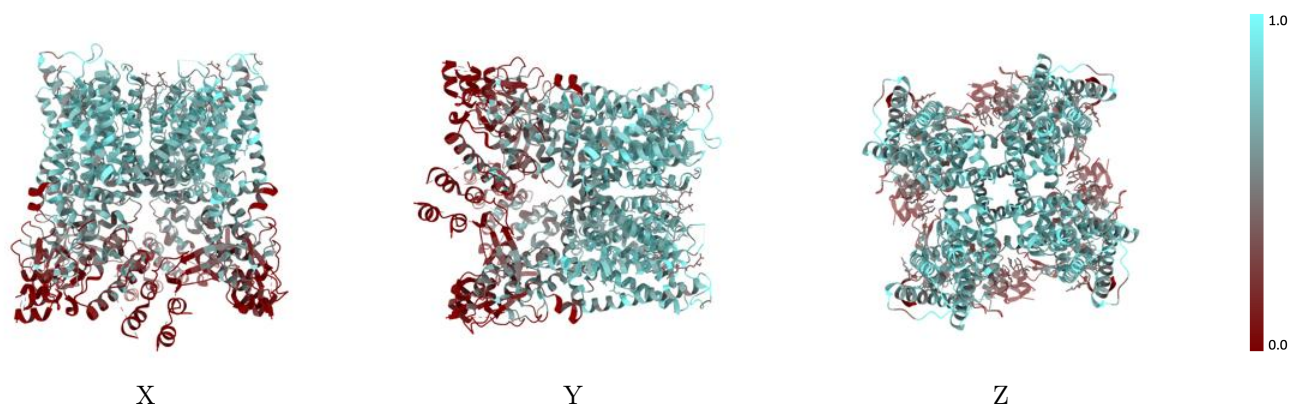
The images above show the 3D surface view of the map at the recommended contour level 0.4 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



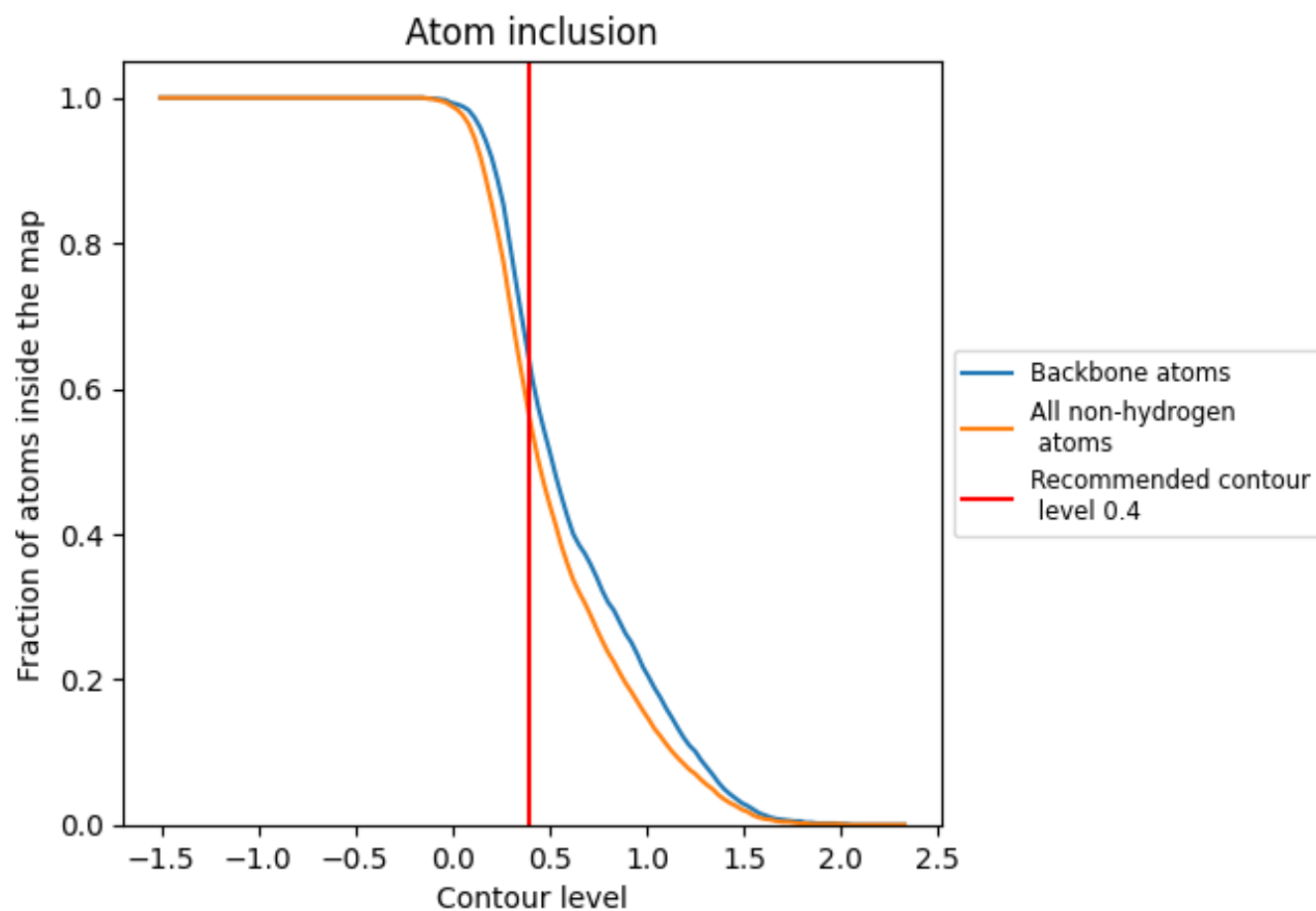
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.4).

9.4 Atom inclusion [i](#)



At the recommended contour level, 63% of all backbone atoms, 56% 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 (0.4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5560	0.4080
A	0.5540	0.4090
B	0.5560	0.4080
C	0.5570	0.4070
D	0.5550	0.4060

