



## Full wwPDB EM Validation Report ⓘ

May 26, 2024 – 04:36 AM EDT

PDB ID : 7RX2  
EMDB ID : EMD-24722  
Title : afTMEM16 in C22 lipid nanodiscs with MSP1E3 scaffold protein in the presence of Ca<sup>2+</sup>  
Authors : Falzone, M.E.; Accardi, A.  
Deposited on : 2021-08-21  
Resolution : 2.70 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev92  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
MolProbity : 4.02b-467  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36.2

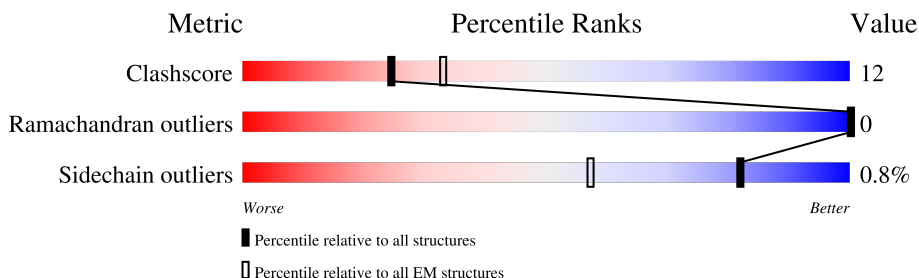
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 2.70 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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  $\geq 3$ , 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	735	 61% 22% 16%
1	B	735	 61% 22% 16%

2 Entry composition ⓘ

There are 4 unique types of molecules in this entry. The entry contains 10390 atoms, of which 12 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.

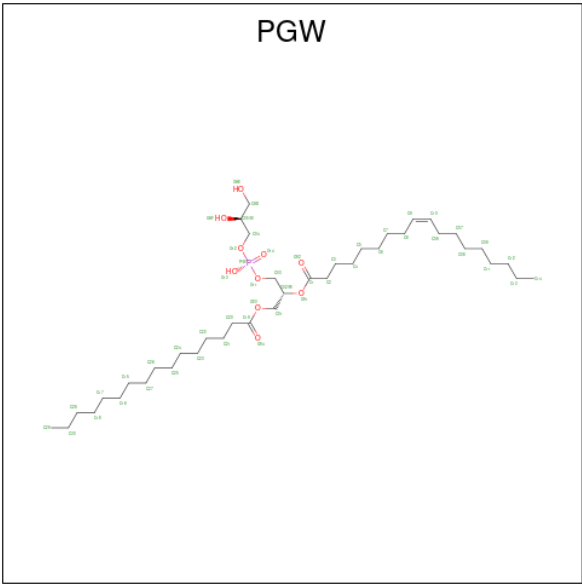
- Molecule 1 is a protein called afTMEM16 lipid scramblase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	615	Total	C	N	O	S	0	0
			4969	3251	832	867	19		
1	A	615	Total	C	N	O	S	0	0
			4969	3251	832	867	19		

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
2	B	2	Total	Ca	0
			2	2	
2	A	2	Total	Ca	0
			2	2	

- Molecule 3 is (1R)-2-{[(S)-{[(2S)-2,3-dihydroxypropyl]oxy}(hydroxy)phosphoryl]oxy}-1-[(hexadecanoyloxy)methyl]ethyl (9Z)-octadec-9-enoate (three-letter code: PGW) (formula: C<sub>40</sub>H<sub>77</sub>O<sub>10</sub>P).



Mol	Chain	Residues	Atoms				AltConf
3	B	1	Total	C	O		0
			26	21	5		
3	B	1	Total	C	O	P	0
			41	30	10	1	
3	B	1	Total	C	O	P	0
			26	17	8	1	
3	B	1	Total	C	O	P	0
			38	29	8	1	
3	B	1	Total	C	O	P	0
			30	21	8	1	
3	B	1	Total	C	O	P	0
			30	21	8	1	
3	B	1	Total	C	O	P	0
			24	15	8	1	
3	A	1	Total	C	O	P	0
			24	15	8	1	
3	A	1	Total	C	O		0
			26	21	5		
3	A	1	Total	C	O	P	0
			41	30	10	1	
3	A	1	Total	C	O	P	0
			26	17	8	1	
3	A	1	Total	C	O	P	0
			38	29	8	1	
3	A	1	Total	C	O	P	0
			30	21	8	1	
3	A	1	Total	C	O	P	0
			30	21	8	1	

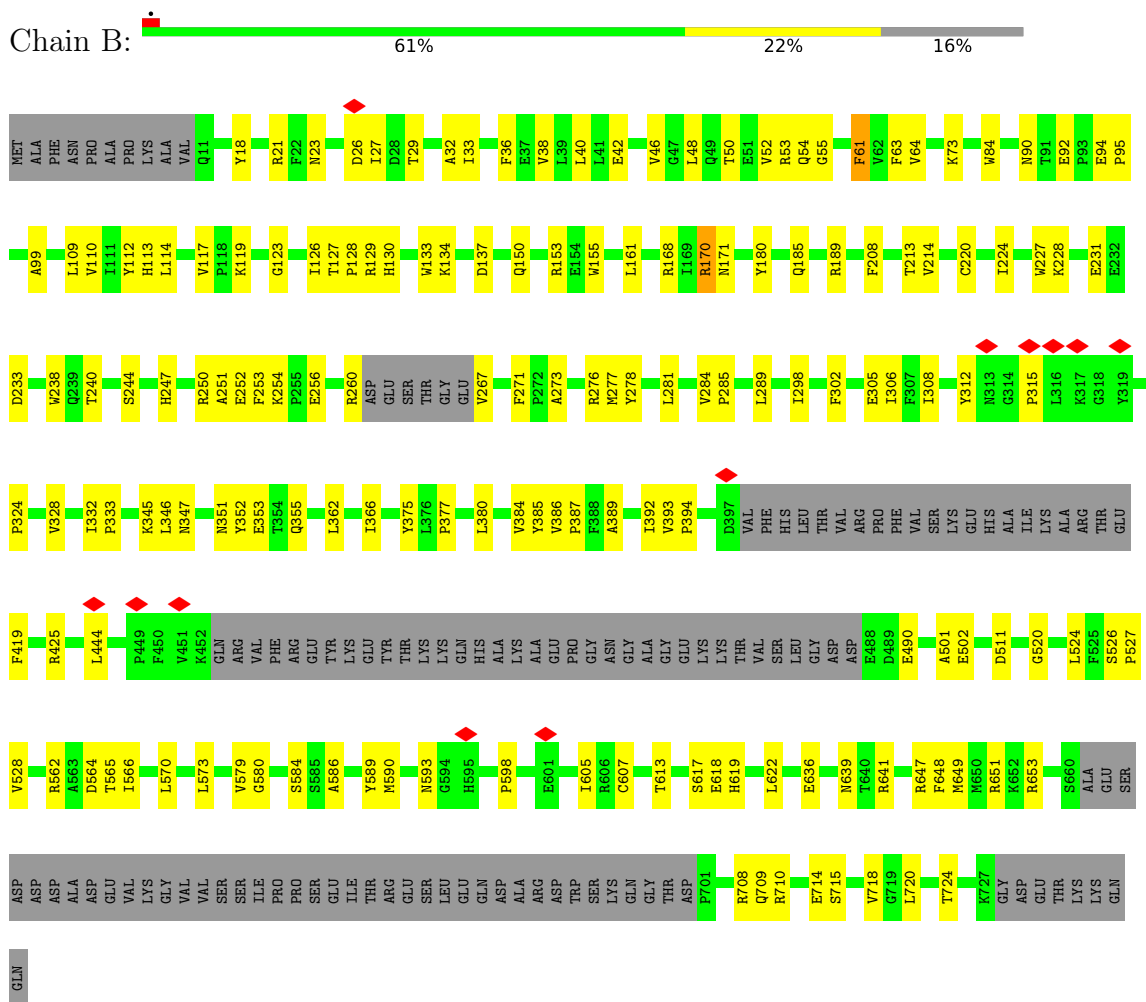
- Molecule 4 is water.

Mol	Chain	Residues	Atoms			AltConf
4	B	3	Total	H	O	0
			9	6	3	
4	A	3	Total	H	O	0
			9	6	3	

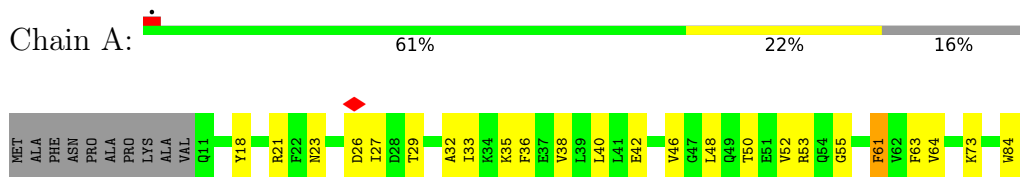
### 3 Residue-property plots

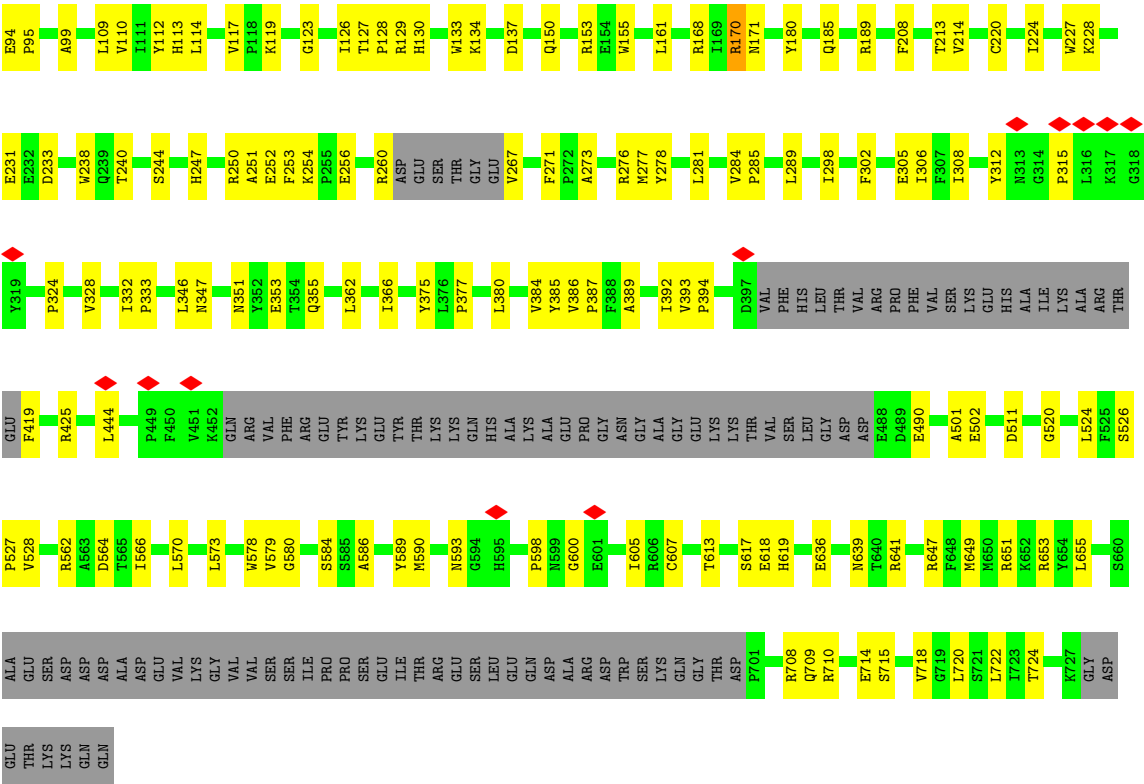
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: aTtMEM16 lipid scramblase



- Molecule 1: aTtMEM16 lipid scramblase





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	132332	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 ( $e^-/\text{\AA}^2$ )	42.7884	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.087	Depositor
Minimum map value	-0.046	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.013	Depositor
Map size ( $\text{\AA}$ )	271.3728, 271.3728, 271.3728	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.7067, 0.7067, 0.7067	Depositor

## 5 Model quality

### 5.1 Standard geometry

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

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	0.44	0/5101	0.47	0/6933
1	B	0.44	0/5101	0.47	0/6933
All	All	0.44	0/10202	0.47	0/13866

There are no bond length outliers.

There are no bond angle outliers.

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	4969	0	4947	122	0
1	B	4969	0	4947	125	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	215	0	239	17	0
3	B	215	0	239	15	0
4	A	3	6	0	0	0
4	B	3	6	0	0	0
All	All	10378	12	10372	244	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 12.



All (244) 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:A:129:ARG:HA	1:A:134:LYS:HA	1.40	1.03
1:B:129:ARG:HA	1:B:134:LYS:HA	1.40	0.98
1:B:189:ARG:HD3	3:B:805:PGW:H03A	1.47	0.96
1:A:189:ARG:HD3	3:A:1506:PGW:H03A	1.47	0.94
1:A:524:LEU:HD23	1:A:580:GLY:HA2	1.62	0.82
1:B:524:LEU:HD23	1:B:580:GLY:HA2	1.62	0.81
1:A:380:LEU:O	1:A:384:VAL:HG22	1.82	0.80
1:B:380:LEU:O	1:B:384:VAL:HG22	1.82	0.79
1:A:109:LEU:HD21	1:A:562:ARG:HD2	1.70	0.73
1:B:598:PRO:HG3	1:B:605:ILE:HD11	1.71	0.73
1:A:598:PRO:HG3	1:A:605:ILE:HD11	1.71	0.73
1:B:109:LEU:HD21	1:B:562:ARG:HD2	1.70	0.73
1:B:21:ARG:HA	1:B:61:PHE:HB3	1.72	0.72
1:A:21:ARG:HA	1:A:61:PHE:HB3	1.72	0.72
1:B:380:LEU:O	1:B:384:VAL:CG2	2.38	0.71
1:A:393:VAL:HG23	1:A:394:PRO:HD3	1.73	0.71
1:A:380:LEU:O	1:A:384:VAL:CG2	2.38	0.70
1:B:21:ARG:NH2	1:B:501:ALA:O	2.24	0.70
1:A:617:SER:HB2	3:A:1505:PGW:H16	1.72	0.70
1:B:617:SER:HB2	3:B:804:PGW:H16	1.72	0.70
1:A:21:ARG:NH2	1:A:501:ALA:O	2.24	0.70
1:B:393:VAL:HG23	1:B:394:PRO:HD3	1.73	0.70
1:A:90:ASN:HD21	1:A:647:ARG:HB2	1.58	0.69
3:B:806:PGW:C30	3:A:1501:PGW:C9	2.72	0.68
3:B:809:PGW:C9	3:A:1507:PGW:C30	2.71	0.68
1:B:619:HIS:NE2	1:A:618:GLU:OE2	2.27	0.68
1:B:90:ASN:HD21	1:B:647:ARG:HB2	1.58	0.67
1:B:490:GLU:OE1	1:A:708:ARG:NH2	2.27	0.65
1:A:21:ARG:HH21	1:A:501:ALA:HB1	1.64	0.63
1:B:21:ARG:HH21	1:B:501:ALA:HB1	1.64	0.61
1:B:586:ALA:O	1:B:590:MET:HG2	2.03	0.59
1:B:240:THR:O	1:B:564:ASP:HB3	2.03	0.59
1:A:586:ALA:O	1:A:590:MET:HG2	2.03	0.59
1:B:710:ARG:N	1:B:714:GLU:OE2	2.34	0.58
1:A:240:THR:O	1:A:564:ASP:HB3	2.02	0.58
1:B:618:GLU:OE2	1:A:619:HIS:NE2	2.34	0.58
1:B:651:ARG:NH2	1:A:84:TRP:O	2.33	0.57
1:A:110:VAL:O	1:A:114:LEU:HD13	2.04	0.57
1:B:110:VAL:O	1:B:114:LEU:HD13	2.05	0.57
1:A:129:ARG:HA	1:A:134:LYS:CA	2.25	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:710:ARG:N	1:A:714:GLU:OE2	2.34	0.56
1:A:284:VAL:HB	1:A:285:PRO:HD3	1.89	0.55
1:B:254:LYS:O	1:B:351:ASN:ND2	2.40	0.55
1:B:385:TYR:CE2	1:B:425:ARG:HD2	2.42	0.55
1:A:714:GLU:O	1:A:718:VAL:HG23	2.07	0.55
1:B:273:ALA:O	1:B:277:MET:HG3	2.07	0.55
1:B:284:VAL:HB	1:B:285:PRO:HD3	1.88	0.55
3:B:805:PGW:H3	3:B:805:PGW:H20A	1.89	0.55
1:A:273:ALA:O	1:A:277:MET:HG3	2.07	0.55
3:A:1506:PGW:H3	3:A:1506:PGW:H20A	1.89	0.55
1:B:619:HIS:HD2	3:A:1507:PGW:H5	1.72	0.55
1:A:254:LYS:O	1:A:351:ASN:ND2	2.40	0.54
1:B:129:ARG:HA	1:B:134:LYS:CA	2.25	0.54
1:A:385:TYR:CE2	1:A:425:ARG:HD2	2.42	0.54
1:A:48:LEU:HD13	1:A:64:VAL:HG11	1.89	0.54
1:B:48:LEU:HD13	1:B:64:VAL:HG11	1.89	0.54
1:B:589:TYR:HD1	3:B:804:PGW:H20A	1.73	0.54
1:A:109:LEU:CD2	1:A:562:ARG:HD2	2.38	0.53
1:A:208:PHE:CE1	1:A:528:VAL:HG22	2.43	0.53
1:B:346:LEU:HB2	3:B:808:PGW:H5	1.91	0.53
1:B:129:ARG:CA	1:B:134:LYS:HA	2.28	0.53
1:B:231:GLU:OE2	1:B:566:ILE:HG12	2.08	0.53
1:B:714:GLU:O	1:B:718:VAL:HG23	2.07	0.53
1:A:231:GLU:OE2	1:A:566:ILE:HG12	2.08	0.53
1:A:524:LEU:HD23	1:A:580:GLY:CA	2.38	0.53
1:B:84:TRP:O	1:A:651:ARG:NH2	2.37	0.53
1:B:127:THR:HG21	1:B:251:ALA:HB2	1.90	0.53
1:A:589:TYR:HD1	3:A:1505:PGW:H20A	1.73	0.53
1:B:224:ILE:HD12	1:B:573:LEU:HB3	1.91	0.53
1:A:46:VAL:HG21	1:A:123:GLY:HA2	1.91	0.53
1:A:253:PHE:CZ	1:A:353:GLU:HG2	2.44	0.53
1:B:208:PHE:CE1	1:B:528:VAL:HG22	2.43	0.53
1:B:393:VAL:HG21	1:B:419:PHE:CG	2.44	0.53
1:B:253:PHE:CZ	1:B:353:GLU:HG2	2.44	0.52
1:A:127:THR:HG21	1:A:251:ALA:HB2	1.90	0.52
1:A:393:VAL:HG21	1:A:419:PHE:CG	2.44	0.52
1:A:214:VAL:HG22	3:A:1505:PGW:H15A	1.91	0.52
1:B:46:VAL:HG21	1:B:123:GLY:HA2	1.91	0.52
1:A:636:GLU:OE1	1:A:641:ARG:HG3	2.10	0.52
1:A:129:ARG:CA	1:A:134:LYS:HA	2.28	0.52
1:A:224:ILE:HD12	1:A:573:LEU:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:VAL:HG22	3:B:804:PGW:H15A	1.91	0.52
1:B:636:GLU:OE1	1:B:641:ARG:HG3	2.10	0.52
1:A:250:ARG:HB3	1:A:355:GLN:HG3	1.92	0.51
1:B:54:GLN:OE1	1:A:722:LEU:HD22	2.11	0.51
1:A:308:ILE:HA	1:A:312:TYR:HB3	1.92	0.51
1:A:346:LEU:HB2	3:A:1509:PGW:H5	1.91	0.51
1:B:389:ALA:HA	1:B:392:ILE:HB	1.93	0.51
1:B:524:LEU:HD23	1:B:580:GLY:CA	2.38	0.51
1:A:389:ALA:HA	1:A:392:ILE:HB	1.93	0.51
1:B:649:MET:O	1:B:653:ARG:HG3	2.11	0.51
1:A:23:ASN:ND2	1:A:137:ASP:HB2	2.27	0.50
1:A:393:VAL:CG2	1:A:394:PRO:HD3	2.39	0.50
1:B:622:LEU:HD13	1:A:578:TRP:CH2	2.47	0.50
1:B:23:ASN:ND2	1:B:137:ASP:HB2	2.27	0.50
1:A:393:VAL:HG11	1:A:419:PHE:CE1	2.46	0.50
1:A:48:LEU:HD13	1:A:64:VAL:CG1	2.42	0.50
1:B:393:VAL:HG21	1:B:419:PHE:CD1	2.47	0.50
1:B:393:VAL:HG11	1:B:419:PHE:CE1	2.46	0.50
1:A:393:VAL:HG21	1:A:419:PHE:CD1	2.47	0.50
1:A:649:MET:O	1:A:653:ARG:HG3	2.11	0.50
1:B:113:HIS:O	1:B:117:VAL:HG22	2.12	0.49
1:A:180:TYR:CZ	1:A:511:ASP:HB3	2.48	0.49
1:B:48:LEU:HD13	1:B:64:VAL:CG1	2.42	0.49
1:B:109:LEU:CD2	1:B:562:ARG:HD2	2.38	0.49
1:B:250:ARG:HB3	1:B:355:GLN:HG3	1.92	0.49
1:A:332:ILE:HB	1:A:333:PRO:HD3	1.94	0.49
1:B:220:CYS:O	1:B:224:ILE:HG12	2.12	0.49
1:A:128:PRO:O	1:A:129:ARG:HG2	2.12	0.49
1:B:180:TYR:CZ	1:B:511:ASP:HB3	2.48	0.49
1:B:308:ILE:HA	1:B:312:TYR:HB3	1.92	0.49
1:B:524:LEU:HD21	1:B:579:VAL:HG12	1.93	0.49
1:B:128:PRO:O	1:B:129:ARG:HG2	2.12	0.49
1:B:332:ILE:HB	1:B:333:PRO:HD3	1.94	0.49
1:A:220:CYS:O	1:A:224:ILE:HG12	2.12	0.49
1:B:227:TRP:CE2	1:B:231:GLU:HG3	2.49	0.48
1:B:393:VAL:CG2	1:B:394:PRO:HD3	2.39	0.48
1:B:27:ILE:HG22	1:B:32:ALA:HB2	1.96	0.48
1:B:52:VAL:O	1:B:53:ARG:HD3	2.14	0.48
1:B:362:LEU:O	1:B:366:ILE:HG13	2.13	0.48
1:A:113:HIS:O	1:A:117:VAL:HG22	2.12	0.48
1:A:524:LEU:HD21	1:A:579:VAL:HG12	1.93	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:PHE:HE2	1:B:276:ARG:HB2	1.78	0.48
1:A:52:VAL:O	1:A:53:ARG:HD3	2.14	0.48
1:B:170:ARG:HD2	1:B:238:TRP:CD2	2.49	0.48
1:A:27:ILE:HG22	1:A:32:ALA:HB2	1.96	0.48
1:A:170:ARG:HD2	1:A:238:TRP:CD2	2.49	0.48
1:A:227:TRP:CE2	1:A:231:GLU:HG3	2.49	0.48
1:A:324:PRO:O	1:A:328:VAL:HG13	2.13	0.48
1:B:305:GLU:OE1	1:B:425:ARG:NH1	2.40	0.48
1:B:260:ARG:HA	1:B:267:VAL:HA	1.96	0.48
1:A:36:PHE:CE2	1:A:40:LEU:HD11	2.49	0.48
1:A:271:PHE:HE2	1:A:276:ARG:HB2	1.77	0.48
1:B:233:ASP:OD1	1:B:639:ASN:HB2	2.14	0.47
1:B:324:PRO:O	1:B:328:VAL:HG13	2.13	0.47
1:B:29:THR:O	1:B:33:ILE:HG22	2.15	0.47
1:A:362:LEU:O	1:A:366:ILE:HG13	2.13	0.47
1:B:84:TRP:HZ3	1:A:655:LEU:HB2	1.79	0.47
1:B:380:LEU:HA	1:B:384:VAL:HG22	1.97	0.47
1:B:90:ASN:ND2	1:B:647:ARG:HB2	2.27	0.47
1:B:94:GLU:HB3	1:B:95:PRO:HD2	1.97	0.47
1:B:150:GLN:NE2	1:B:168:ARG:HH12	2.13	0.47
1:B:347:ASN:ND2	1:B:362:LEU:HB2	2.30	0.47
1:A:29:THR:O	1:A:33:ILE:HG22	2.15	0.47
1:A:233:ASP:OD1	1:A:639:ASN:HB2	2.14	0.47
1:A:347:ASN:ND2	1:A:362:LEU:HB2	2.30	0.47
1:B:36:PHE:CE2	1:B:40:LEU:HD11	2.49	0.47
1:A:386:VAL:HB	1:A:387:PRO:HD3	1.97	0.47
1:A:73:LYS:HG2	1:A:99:ALA:O	2.15	0.46
1:A:260:ARG:HA	1:A:267:VAL:HA	1.96	0.46
1:A:380:LEU:HA	1:A:384:VAL:HG22	1.96	0.46
1:A:150:GLN:NE2	1:A:168:ARG:HH12	2.13	0.46
1:A:305:GLU:OE1	1:A:425:ARG:NH1	2.40	0.46
1:A:607:CYS:HB2	3:A:1504:PGW:H02	1.97	0.46
1:B:298:ILE:HD13	1:B:377:PRO:HD3	1.98	0.46
1:A:42:GLU:CD	1:A:119:LYS:HE2	2.36	0.46
1:A:94:GLU:HB3	1:A:95:PRO:HD2	1.97	0.46
1:B:73:LYS:HG2	1:B:99:ALA:O	2.16	0.46
1:A:90:ASN:ND2	1:A:647:ARG:HB2	2.27	0.46
1:A:709:GLN:OE1	1:A:715:SER:HA	2.16	0.46
1:B:42:GLU:CD	1:B:119:LYS:HE2	2.36	0.46
1:B:130:HIS:CD2	1:B:252:GLU:HG3	2.51	0.46
1:B:386:VAL:HB	1:B:387:PRO:HD3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:HIS:CD2	1:A:252:GLU:HG3	2.51	0.45
1:A:256:GLU:HG2	1:A:271:PHE:HD1	1.81	0.45
1:B:607:CYS:HB2	3:B:803:PGW:H02	1.97	0.45
1:A:298:ILE:HD13	1:A:377:PRO:HD3	1.98	0.45
1:B:380:LEU:O	1:B:384:VAL:HG23	2.17	0.45
1:B:590:MET:HA	3:B:804:PGW:HADA	1.99	0.45
1:B:613:THR:HB	3:B:804:PGW:H4A	1.99	0.45
1:A:256:GLU:HG2	1:A:271:PHE:CD1	2.52	0.45
1:B:170:ARG:HD2	1:B:238:TRP:CE2	2.52	0.45
3:B:806:PGW:H15	3:B:806:PGW:H18	1.62	0.45
1:B:709:GLN:OE1	1:B:715:SER:HA	2.16	0.45
1:A:613:THR:HB	3:A:1505:PGW:H4A	1.99	0.45
1:B:244:SER:HA	1:B:247:HIS:CE1	2.52	0.44
1:A:170:ARG:HD2	1:A:238:TRP:CE2	2.52	0.44
1:B:720:LEU:O	1:B:724:THR:HG23	2.17	0.44
1:A:302:PHE:O	1:A:306:ILE:HG12	2.17	0.44
1:B:256:GLU:HG2	1:B:271:PHE:CD1	2.52	0.44
1:B:302:PHE:O	1:B:306:ILE:HG12	2.17	0.44
1:A:18:TYR:OH	1:A:112:TYR:HB2	2.17	0.44
1:A:214:VAL:CG2	3:A:1505:PGW:H15A	2.48	0.44
1:A:244:SER:HA	1:A:247:HIS:CE1	2.52	0.44
1:A:312:TYR:OH	1:A:315:PRO:O	2.20	0.44
1:B:18:TYR:OH	1:B:112:TYR:HB2	2.17	0.44
1:A:38:VAL:O	1:A:42:GLU:HG2	2.18	0.44
1:A:590:MET:HA	3:A:1505:PGW:HADA	1.99	0.44
1:B:38:VAL:O	1:B:42:GLU:HG2	2.18	0.44
1:B:256:GLU:HG2	1:B:271:PHE:HD1	1.81	0.44
1:B:708:ARG:NH2	1:A:490:GLU:OE1	2.46	0.44
1:B:170:ARG:HG3	1:B:171:ASN:N	2.32	0.44
1:B:55:GLY:HA2	1:B:61:PHE:CE2	2.53	0.43
1:A:170:ARG:HG3	1:A:171:ASN:N	2.32	0.43
1:A:224:ILE:CD1	1:A:573:LEU:HB3	2.48	0.43
1:B:312:TYR:OH	1:B:315:PRO:O	2.20	0.43
1:A:289:LEU:HD13	3:A:1509:PGW:C10	2.49	0.43
1:A:55:GLY:HA2	1:A:61:PHE:CE2	2.53	0.43
1:A:720:LEU:O	1:A:724:THR:HG23	2.18	0.43
1:B:289:LEU:HD13	3:B:808:PGW:C10	2.49	0.43
1:A:278:TYR:O	1:A:281:LEU:HB2	2.19	0.43
1:B:214:VAL:CG2	3:B:804:PGW:H15A	2.48	0.43
1:A:127:THR:O	1:A:133:TRP:HB2	2.19	0.43
1:B:50:THR:HA	1:B:63:PHE:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:278:TYR:O	1:B:281:LEU:HB2	2.19	0.43
1:B:127:THR:O	1:B:133:TRP:HB2	2.19	0.43
1:B:161:LEU:HB2	1:B:185:GLN:OE1	2.19	0.43
1:B:224:ILE:CD1	1:B:573:LEU:HB3	2.48	0.43
1:B:648:PHE:CE1	1:A:85:LEU:HB3	2.54	0.43
1:B:26:ASP:O	1:B:27:ILE:HD13	2.19	0.42
1:B:213:THR:HG23	1:B:584:SER:HB3	2.01	0.42
1:A:213:THR:HG23	1:A:584:SER:HB3	2.01	0.42
1:A:26:ASP:O	1:A:27:ILE:HD13	2.19	0.42
1:A:380:LEU:O	1:A:384:VAL:HG23	2.16	0.42
1:A:598:PRO:O	1:A:600:GLY:N	2.50	0.42
1:A:50:THR:HA	1:A:63:PHE:O	2.19	0.42
1:A:502:GLU:HA	1:A:502:GLU:OE2	2.19	0.42
1:B:502:GLU:OE2	1:B:502:GLU:HA	2.19	0.42
1:A:161:LEU:HB2	1:A:185:GLN:OE1	2.19	0.42
1:B:150:GLN:HG3	1:B:153:ARG:NH2	2.35	0.42
1:B:155:TRP:CG	1:B:161:LEU:HD11	2.55	0.42
1:A:150:GLN:HG3	1:A:153:ARG:NH2	2.35	0.42
1:A:393:VAL:HG11	1:A:419:PHE:CZ	2.55	0.42
1:B:526:SER:N	1:B:527:PRO:HD2	2.35	0.42
1:A:155:TRP:CG	1:A:161:LEU:HD11	2.55	0.42
3:A:1507:PGW:H18	3:A:1507:PGW:H15	1.62	0.42
1:A:526:SER:N	1:A:527:PRO:HD2	2.35	0.42
1:B:375:TYR:OH	1:B:520:GLY:HA3	2.20	0.41
3:A:1501:PGW:H5	3:A:1501:PGW:H2	1.78	0.41
1:A:126:ILE:HA	1:A:133:TRP:CD1	2.56	0.41
1:A:228:LYS:HG2	1:A:570:LEU:HD21	2.02	0.41
1:B:393:VAL:HG11	1:B:419:PHE:CZ	2.55	0.41
1:B:345:LYS:HD3	1:B:345:LYS:HA	1.90	0.41
1:B:126:ILE:HA	1:B:133:TRP:CD1	2.56	0.41
1:B:92:GLU:OE1	1:B:653:ARG:NH1	2.54	0.41
1:B:278:TYR:HB3	3:B:807:PGW:H2A	2.03	0.41
1:A:92:GLU:OE1	1:A:653:ARG:NH1	2.54	0.41
1:A:375:TYR:OH	1:A:520:GLY:HA3	2.20	0.41
1:B:228:LYS:HG2	1:B:570:LEU:HD21	2.02	0.41
1:B:305:GLU:HB2	1:B:324:PRO:HG2	2.03	0.41
1:A:35:LYS:HA	1:A:38:VAL:HG22	2.03	0.40
1:B:250:ARG:HD3	1:B:352:TYR:O	2.21	0.40
1:B:231:GLU:OE1	1:B:565:THR:OG1	2.37	0.40
1:A:278:TYR:HB3	3:A:1508:PGW:H2A	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	605/735 (82%)	571 (94%)	34 (6%)	0	100	100
1	B	605/735 (82%)	571 (94%)	34 (6%)	0	100	100
All	All	1210/1470 (82%)	1142 (94%)	68 (6%)	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	529/646 (82%)	525 (99%)	4 (1%)	81	93
1	B	529/646 (82%)	525 (99%)	4 (1%)	81	93
All	All	1058/1292 (82%)	1050 (99%)	8 (1%)	82	93

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

Mol	Chain	Res	Type
1	B	61	PHE
1	B	170	ARG
1	B	444	LEU
1	B	593	ASN
1	A	61	PHE
1	A	170	ARG
1	A	444	LEU
1	A	593	ASN



Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	148	ASN
1	B	150	GLN
1	B	283	GLN
1	B	370	ASN
1	B	438	GLN
1	B	499	ASN
1	A	148	ASN
1	A	150	GLN
1	A	283	GLN
1	A	370	ASN
1	A	438	GLN
1	A	499	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 4 are monoatomic - leaving 14 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
3	PGW	B	806	-	37,37,50	1.09	5 (13%)	41,42,56	1.23	3 (7%)
3	PGW	B	803	-	25,25,50	1.18	3 (12%)	27,27,56	1.38	2 (7%)
3	PGW	A	1504	-	25,25,50	1.18	3 (12%)	27,27,56	1.38	2 (7%)
3	PGW	B	804	-	40,40,50	0.29	0	43,46,56	0.34	0
3	PGW	A	1508	-	29,29,50	1.21	4 (13%)	33,34,56	1.07	2 (6%)
3	PGW	A	1506	-	25,25,50	1.30	4 (16%)	29,30,56	1.44	3 (10%)
3	PGW	A	1501	-	23,23,50	0.49	0	27,28,56	0.64	1 (3%)
3	PGW	B	805	-	25,25,50	1.30	4 (16%)	29,30,56	1.44	3 (10%)
3	PGW	B	807	-	29,29,50	1.21	4 (13%)	33,34,56	1.07	2 (6%)
3	PGW	A	1507	-	37,37,50	1.09	5 (13%)	41,42,56	1.23	3 (7%)
3	PGW	A	1509	-	28,28,50	1.24	5 (17%)	30,32,56	1.28	2 (6%)
3	PGW	A	1505	-	40,40,50	0.29	0	43,46,56	0.34	0
3	PGW	B	808	-	28,28,50	1.24	5 (17%)	30,32,56	1.28	2 (6%)
3	PGW	B	809	-	23,23,50	0.48	0	27,28,56	0.64	1 (3%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PGW	B	806	-	-	18/39/39/55	-
3	PGW	B	803	-	-	9/27/27/55	-
3	PGW	A	1504	-	-	9/27/27/55	-
3	PGW	B	804	-	-	22/45/45/55	-
3	PGW	A	1508	-	-	8/31/31/55	-
3	PGW	A	1506	-	-	8/27/27/55	-
3	PGW	A	1501	-	-	8/25/25/55	-
3	PGW	B	805	-	-	8/27/27/55	-
3	PGW	B	807	-	-	8/31/31/55	-
3	PGW	A	1507	-	-	18/39/39/55	-
3	PGW	A	1509	-	-	15/28/28/55	-
3	PGW	A	1505	-	-	22/45/45/55	-
3	PGW	B	808	-	-	15/28/28/55	-
3	PGW	B	809	-	-	8/25/25/55	-

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1508	PGW	O01-C1	2.95	1.42	1.34
3	B	807	PGW	O01-C1	2.94	1.42	1.34
3	A	1504	PGW	O01-C1	2.89	1.42	1.34
3	B	803	PGW	O01-C1	2.87	1.42	1.34
3	B	805	PGW	O01-C1	2.86	1.42	1.34
3	A	1506	PGW	O01-C1	2.85	1.42	1.34
3	B	808	PGW	O01-C1	2.70	1.41	1.34
3	A	1509	PGW	O01-C1	2.69	1.41	1.34
3	A	1507	PGW	O01-C1	2.66	1.41	1.34
3	B	806	PGW	O01-C1	2.65	1.41	1.34
3	A	1507	PGW	O03-C01	-2.47	1.39	1.45
3	A	1506	PGW	O03-C19	2.45	1.40	1.33
3	B	805	PGW	O03-C19	2.45	1.40	1.33
3	B	806	PGW	O03-C01	-2.43	1.39	1.45
3	B	803	PGW	O03-C19	2.43	1.40	1.33
3	A	1504	PGW	O03-C19	2.42	1.40	1.33
3	B	805	PGW	O03-C01	-2.42	1.39	1.45
3	A	1506	PGW	O03-C01	-2.42	1.39	1.45
3	A	1509	PGW	O03-C19	2.39	1.40	1.33
3	A	1508	PGW	P-O12	2.39	1.64	1.54
3	B	807	PGW	P-O12	2.37	1.64	1.54
3	B	808	PGW	O03-C19	2.37	1.40	1.33
3	A	1509	PGW	O03-C01	-2.36	1.39	1.45
3	B	808	PGW	O03-C01	-2.34	1.39	1.45
3	B	805	PGW	P-O12	2.34	1.63	1.54
3	A	1506	PGW	P-O12	2.34	1.63	1.54
3	A	1508	PGW	O03-C19	2.31	1.40	1.33
3	B	807	PGW	O03-C19	2.30	1.40	1.33
3	A	1509	PGW	P-O12	2.29	1.63	1.54
3	B	808	PGW	P-O12	2.29	1.63	1.54
3	A	1507	PGW	P-O12	2.27	1.63	1.54
3	B	806	PGW	O03-C19	2.26	1.39	1.33
3	B	806	PGW	P-O12	2.26	1.63	1.54
3	A	1507	PGW	O03-C19	2.25	1.39	1.33
3	B	807	PGW	O03-C01	-2.24	1.40	1.45
3	A	1508	PGW	O03-C01	-2.23	1.40	1.45
3	B	806	PGW	O01-C02	-2.23	1.41	1.46
3	A	1507	PGW	O01-C02	-2.20	1.41	1.46
3	B	808	PGW	O01-C02	-2.16	1.41	1.46
3	A	1509	PGW	O01-C02	-2.15	1.41	1.46
3	A	1504	PGW	O03-C01	-2.15	1.40	1.45
3	B	803	PGW	O03-C01	-2.15	1.40	1.45

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	805	PGW	O01-C1-C2	4.93	122.12	111.50
3	A	1506	PGW	O01-C1-C2	4.93	122.12	111.50
3	A	1504	PGW	O01-C1-C2	4.27	120.70	111.50
3	B	803	PGW	O01-C1-C2	4.27	120.70	111.50
3	B	806	PGW	O01-C1-C2	3.88	119.87	111.50
3	B	808	PGW	O01-C1-C2	3.88	119.87	111.50
3	A	1509	PGW	O01-C1-C2	3.88	119.87	111.50
3	A	1507	PGW	O01-C1-C2	3.88	119.86	111.50
3	B	807	PGW	O01-C1-C2	3.26	118.53	111.50
3	A	1508	PGW	O01-C1-C2	3.26	118.52	111.50
3	B	803	PGW	O03-C19-C20	2.76	120.58	111.91
3	A	1504	PGW	O03-C19-C20	2.76	120.57	111.91
3	A	1507	PGW	O03-C19-C20	2.71	120.40	111.91
3	B	806	PGW	O03-C19-C20	2.70	120.37	111.91
3	B	805	PGW	O03-C19-C20	2.60	120.07	111.91
3	A	1506	PGW	O03-C19-C20	2.59	120.04	111.91
3	B	808	PGW	O03-C19-C20	2.55	119.90	111.91
3	A	1509	PGW	O03-C19-C20	2.54	119.89	111.91
3	B	807	PGW	O03-C19-C20	2.48	119.70	111.91
3	A	1508	PGW	O03-C19-C20	2.48	119.68	111.91
3	A	1501	PGW	O13-P-O14	2.33	119.78	110.68
3	B	809	PGW	O13-P-O14	2.32	119.76	110.68
3	A	1507	PGW	C02-O01-C1	-2.23	112.30	117.79
3	B	806	PGW	C02-O01-C1	-2.22	112.33	117.79
3	B	805	PGW	O01-C1-O02	-2.18	118.43	123.70
3	A	1506	PGW	O01-C1-O02	-2.17	118.45	123.70

There are no chirality outliers.

All (176) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	803	PGW	C01-C02-C03-O11
3	B	803	PGW	O01-C02-C03-O11
3	B	804	PGW	C03-O11-P-O13
3	B	804	PGW	C03-O11-P-O14
3	B	805	PGW	C03-O11-P-O12
3	B	805	PGW	C03-O11-P-O13
3	B	805	PGW	C03-O11-P-O14
3	B	805	PGW	O03-C01-C02-O01
3	B	806	PGW	C03-O11-P-O12
3	B	806	PGW	C03-O11-P-O13
3	B	806	PGW	C03-O11-P-O14
3	B	807	PGW	C10-C06-C07-C08

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Mol	Chain	Res	Type	Atoms
3	B	808	PGW	C2-C1-O01-C02
3	B	808	PGW	O02-C1-O01-C02
3	A	1504	PGW	C01-C02-C03-O11
3	A	1504	PGW	O01-C02-C03-O11
3	A	1505	PGW	C03-O11-P-O13
3	A	1505	PGW	C03-O11-P-O14
3	A	1506	PGW	C03-O11-P-O12
3	A	1506	PGW	C03-O11-P-O13
3	A	1506	PGW	C03-O11-P-O14
3	A	1506	PGW	O03-C01-C02-O01
3	A	1507	PGW	C03-O11-P-O12
3	A	1507	PGW	C03-O11-P-O13
3	A	1507	PGW	C03-O11-P-O14
3	A	1508	PGW	C10-C06-C07-C08
3	A	1509	PGW	C2-C1-O01-C02
3	A	1509	PGW	O02-C1-O01-C02
3	B	809	PGW	C2-C3-C4-C5
3	A	1501	PGW	C2-C3-C4-C5
3	B	807	PGW	C2-C3-C4-C5
3	A	1508	PGW	C2-C3-C4-C5
3	B	806	PGW	C1-C2-C3-C4
3	B	806	PGW	C19-C20-C21-C22
3	A	1507	PGW	C1-C2-C3-C4
3	A	1507	PGW	C19-C20-C21-C22
3	B	806	PGW	C20-C19-O03-C01
3	A	1507	PGW	C20-C19-O03-C01
3	B	804	PGW	C20-C19-O03-C01
3	A	1505	PGW	C20-C19-O03-C01
3	B	804	PGW	C1-C2-C3-C4
3	B	809	PGW	C1-C2-C3-C4
3	A	1501	PGW	C1-C2-C3-C4
3	A	1505	PGW	C1-C2-C3-C4
3	B	806	PGW	O04-C19-O03-C01
3	A	1507	PGW	O04-C19-O03-C01
3	B	804	PGW	C03-O11-P-O12
3	A	1505	PGW	C03-O11-P-O12
3	B	806	PGW	C20-C21-C22-C23
3	A	1505	PGW	C16-C15-C27-C26
3	A	1507	PGW	C20-C21-C22-C23
3	B	803	PGW	C2-C1-O01-C02
3	A	1504	PGW	C2-C1-O01-C02
3	B	804	PGW	C16-C15-C27-C26

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Mol	Chain	Res	Type	Atoms
3	B	804	PGW	O04-C19-O03-C01
3	A	1505	PGW	O04-C19-O03-C01
3	B	803	PGW	O02-C1-O01-C02
3	A	1504	PGW	O02-C1-O01-C02
3	B	804	PGW	C21-C22-C23-C24
3	A	1505	PGW	C21-C22-C23-C24
3	B	804	PGW	C25-C26-C27-C15
3	A	1505	PGW	C25-C26-C27-C15
3	B	803	PGW	C6-C7-C8-C9
3	A	1504	PGW	C6-C7-C8-C9
3	B	803	PGW	C22-C23-C24-C25
3	B	804	PGW	C24-C25-C26-C27
3	A	1504	PGW	C22-C23-C24-C25
3	A	1505	PGW	C24-C25-C26-C27
3	B	803	PGW	C19-C20-C21-C22
3	A	1504	PGW	C19-C20-C21-C22
3	B	804	PGW	C15-C16-C17-C18
3	A	1505	PGW	C15-C16-C17-C18
3	B	806	PGW	C24-C25-C26-C27
3	A	1507	PGW	C24-C25-C26-C27
3	B	806	PGW	C21-C22-C23-C24
3	A	1507	PGW	C21-C22-C23-C24
3	B	806	PGW	C15-C16-C17-C18
3	A	1507	PGW	C15-C16-C17-C18
3	B	808	PGW	C6-C7-C8-C9
3	A	1509	PGW	C6-C7-C8-C9
3	B	804	PGW	C17-C18-C28-C30
3	A	1505	PGW	C17-C18-C28-C30
3	B	808	PGW	C7-C8-C9-C10
3	A	1509	PGW	C7-C8-C9-C10
3	A	1505	PGW	C5-C6-C7-C8
3	B	804	PGW	C5-C6-C7-C8
3	B	804	PGW	C4-C5-C6-C7
3	A	1505	PGW	C4-C5-C6-C7
3	B	804	PGW	O01-C02-C03-O11
3	A	1505	PGW	O01-C02-C03-O11
3	B	808	PGW	C01-C02-C03-O11
3	A	1509	PGW	C01-C02-C03-O11
3	B	806	PGW	C22-C23-C24-C25
3	A	1507	PGW	C22-C23-C24-C25
3	B	808	PGW	C20-C19-O03-C01
3	A	1509	PGW	C20-C19-O03-C01

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Mol	Chain	Res	Type	Atoms
3	B	805	PGW	C02-C03-O11-P
3	A	1506	PGW	C02-C03-O11-P
3	B	804	PGW	C19-C20-C21-C22
3	A	1505	PGW	C19-C20-C21-C22
3	B	808	PGW	O01-C02-C03-O11
3	A	1509	PGW	O01-C02-C03-O11
3	B	808	PGW	C4-C5-C6-C7
3	A	1509	PGW	C4-C5-C6-C7
3	B	805	PGW	C5-C6-C7-C8
3	A	1506	PGW	C5-C6-C7-C8
3	B	807	PGW	C6-C7-C8-C9
3	A	1508	PGW	C6-C7-C8-C9
3	B	807	PGW	C01-C02-C03-O11
3	A	1508	PGW	C01-C02-C03-O11
3	B	805	PGW	O03-C01-C02-C03
3	A	1506	PGW	O03-C01-C02-C03
3	B	807	PGW	O01-C02-C03-O11
3	A	1508	PGW	O01-C02-C03-O11
3	B	808	PGW	O03-C01-C02-O01
3	A	1509	PGW	O03-C01-C02-O01
3	B	808	PGW	O04-C19-O03-C01
3	A	1509	PGW	O04-C19-O03-C01
3	B	804	PGW	C04-C05-CAD-OAE
3	A	1505	PGW	C04-C05-CAD-OAE
3	B	804	PGW	C23-C24-C25-C26
3	A	1505	PGW	C23-C24-C25-C26
3	B	804	PGW	C01-C02-C03-O11
3	A	1505	PGW	C01-C02-C03-O11
3	B	809	PGW	O04-C19-C20-C21
3	A	1501	PGW	O04-C19-C20-C21
3	B	809	PGW	C6-C7-C8-C9
3	A	1501	PGW	C6-C7-C8-C9
3	B	808	PGW	O03-C19-C20-C21
3	A	1509	PGW	O03-C19-C20-C21
3	B	804	PGW	C2-C3-C4-C5
3	A	1508	PGW	C4-C5-C6-C7
3	A	1505	PGW	C2-C3-C4-C5
3	B	807	PGW	C4-C5-C6-C7
3	B	804	PGW	C04-O12-P-O11
3	A	1505	PGW	C04-O12-P-O11
3	B	808	PGW	O03-C01-C02-C03
3	A	1509	PGW	O03-C01-C02-C03

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Mol	Chain	Res	Type	Atoms
3	B	805	PGW	C1-C2-C3-C4
3	A	1506	PGW	C1-C2-C3-C4
3	B	808	PGW	C19-C20-C21-C22
3	A	1509	PGW	C19-C20-C21-C22
3	B	804	PGW	C03-C02-O01-C1
3	B	807	PGW	C01-C02-O01-C1
3	A	1505	PGW	C03-C02-O01-C1
3	A	1508	PGW	C01-C02-O01-C1
3	B	809	PGW	O03-C19-C20-C21
3	A	1501	PGW	O03-C19-C20-C21
3	A	1507	PGW	C27-C15-C16-C17
3	B	806	PGW	C27-C15-C16-C17
3	B	809	PGW	O03-C01-C02-O01
3	A	1501	PGW	O03-C01-C02-O01
3	B	809	PGW	O01-C1-C2-C3
3	A	1501	PGW	O01-C1-C2-C3
3	B	807	PGW	C07-C06-C10-C9
3	A	1508	PGW	C07-C06-C10-C9
3	B	803	PGW	O03-C19-C20-C21
3	A	1504	PGW	O03-C19-C20-C21
3	B	808	PGW	C2-C3-C4-C5
3	A	1509	PGW	C2-C3-C4-C5
3	B	806	PGW	O01-C1-C2-C3
3	A	1507	PGW	O01-C1-C2-C3
3	B	806	PGW	C7-C8-C9-C10
3	A	1507	PGW	C7-C8-C9-C10
3	B	806	PGW	O03-C19-C20-C21
3	A	1507	PGW	O03-C19-C20-C21
3	B	803	PGW	O04-C19-C20-C21
3	A	1504	PGW	O04-C19-C20-C21
3	B	806	PGW	O02-C1-C2-C3
3	A	1507	PGW	O02-C1-C2-C3
3	B	806	PGW	O04-C19-C20-C21
3	A	1507	PGW	O04-C19-C20-C21
3	B	809	PGW	C5-C6-C7-C8
3	A	1501	PGW	C5-C6-C7-C8
3	B	808	PGW	O01-C1-C2-C3
3	A	1509	PGW	O01-C1-C2-C3

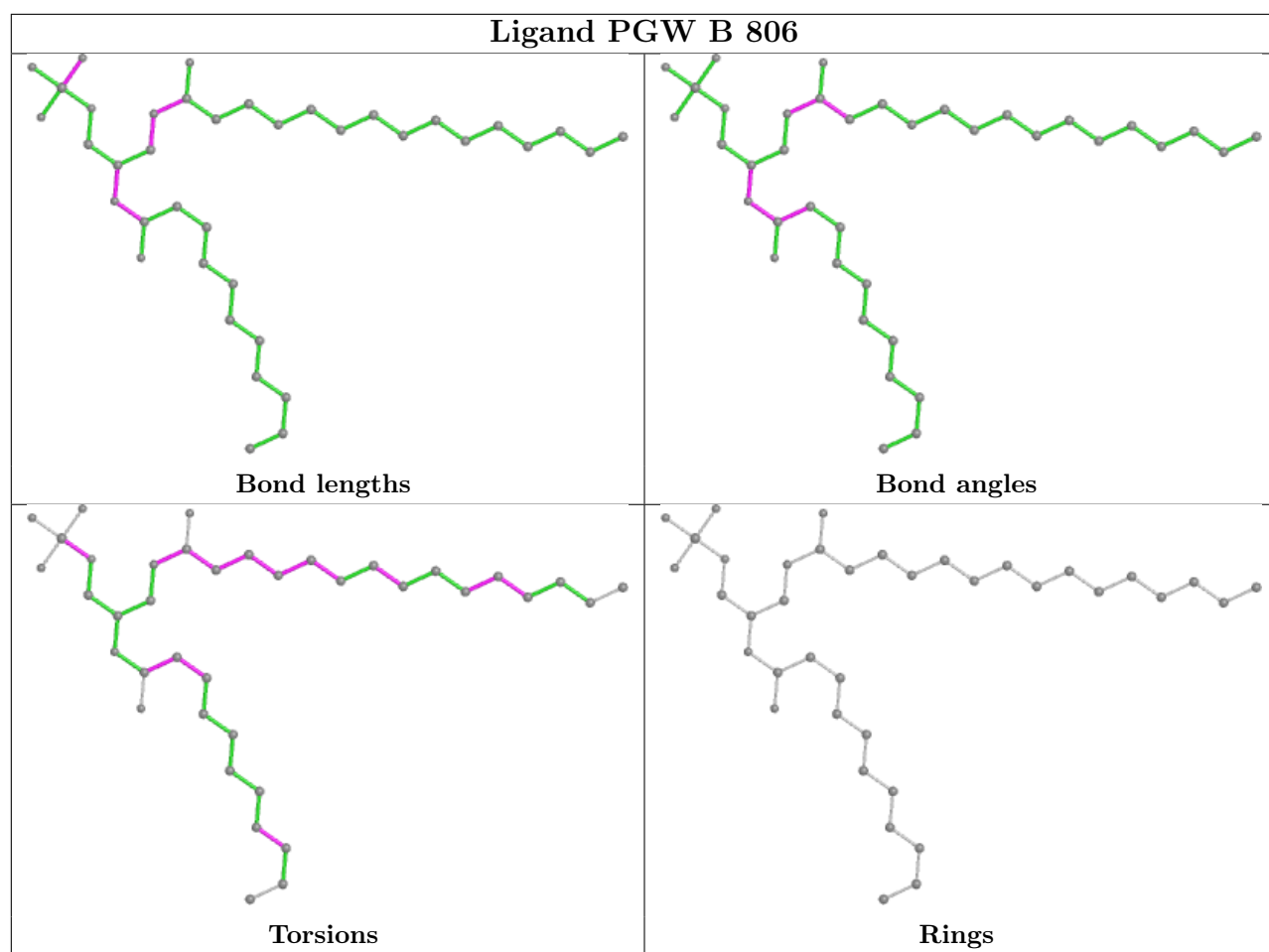
There are no ring outliers.

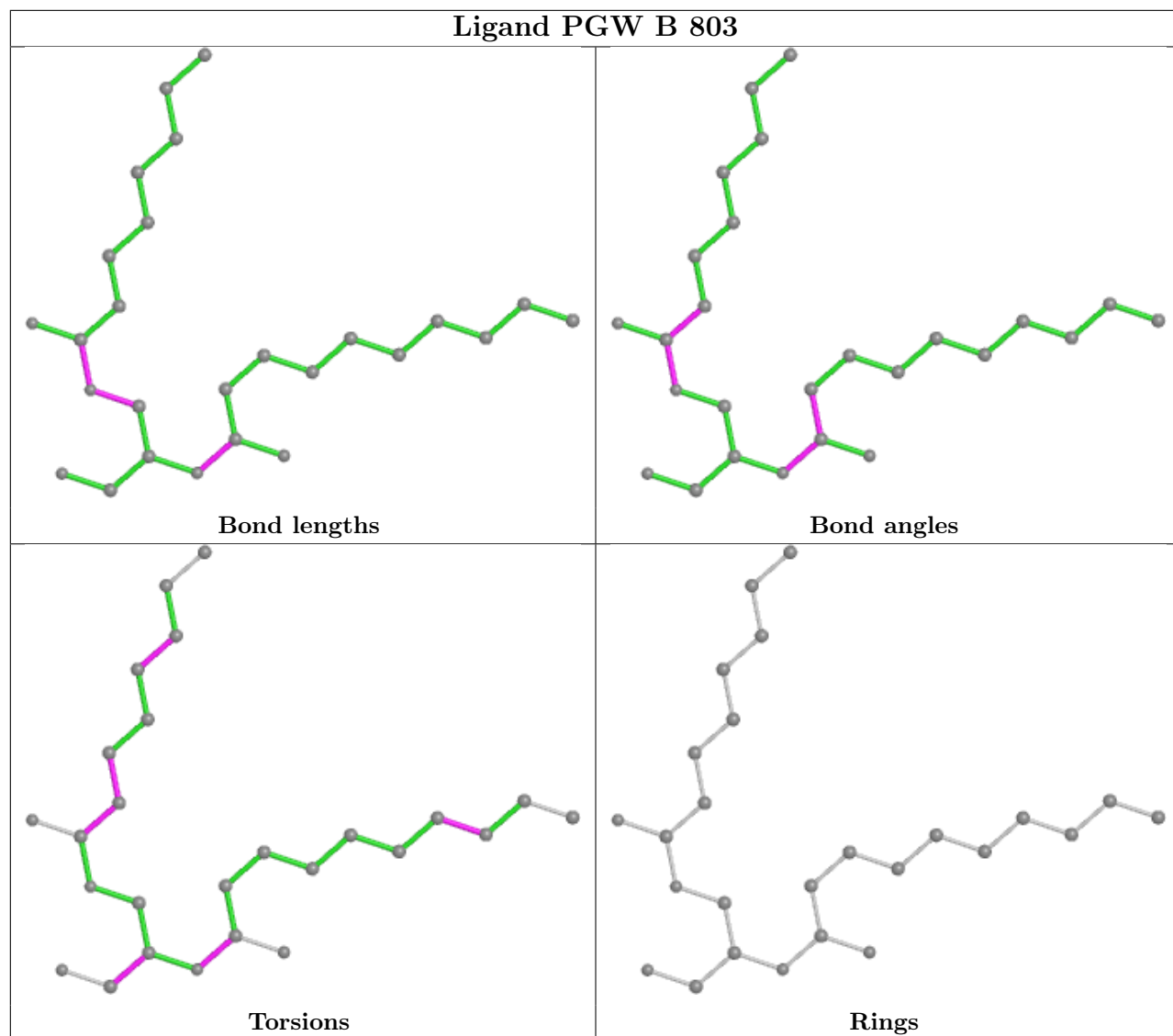
14 monomers are involved in 30 short contacts:

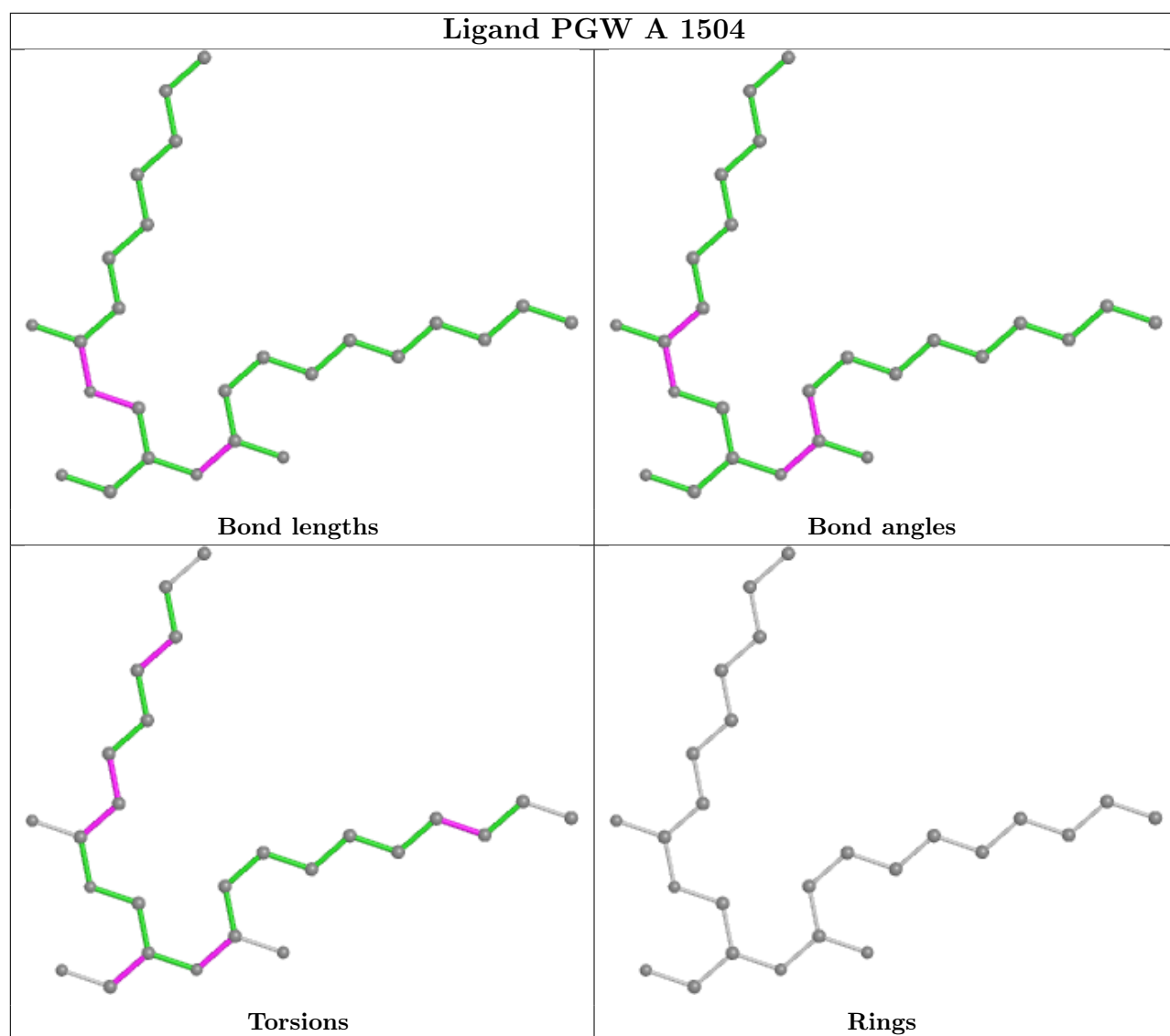
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	806	PGW	2	0
3	B	803	PGW	1	0
3	A	1504	PGW	1	0
3	B	804	PGW	6	0
3	A	1508	PGW	1	0
3	A	1506	PGW	2	0
3	A	1501	PGW	2	0
3	B	805	PGW	2	0
3	B	807	PGW	1	0
3	A	1507	PGW	3	0
3	A	1509	PGW	2	0
3	A	1505	PGW	6	0
3	B	808	PGW	2	0
3	B	809	PGW	1	0

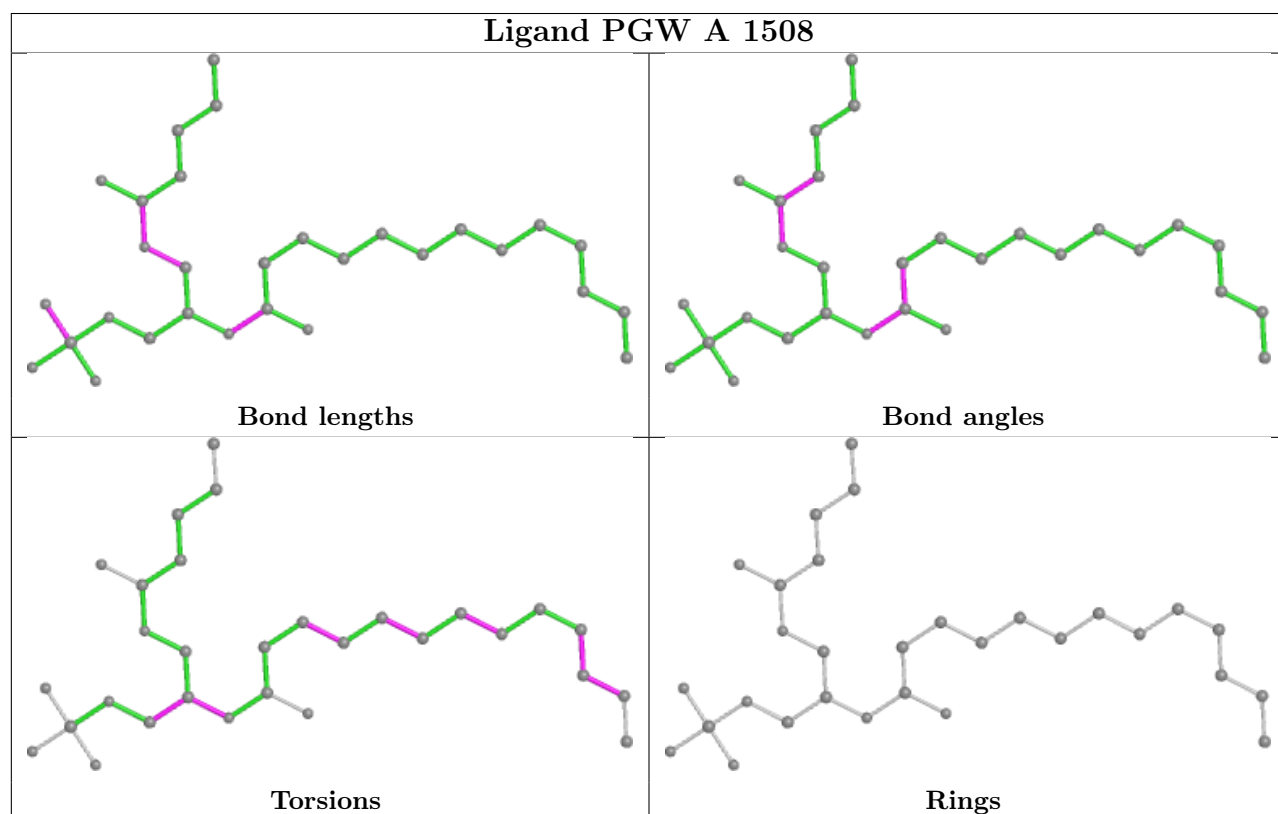
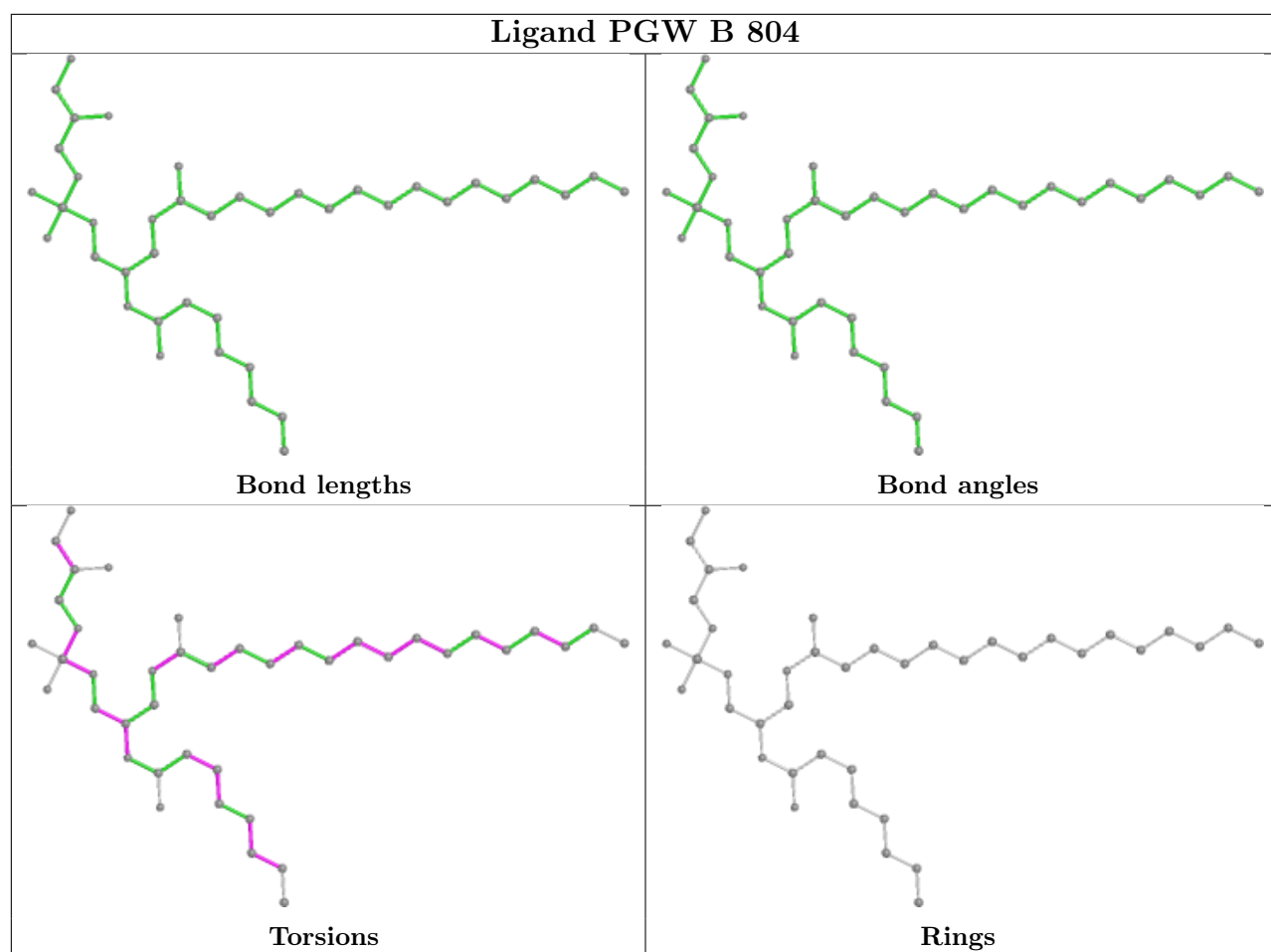
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



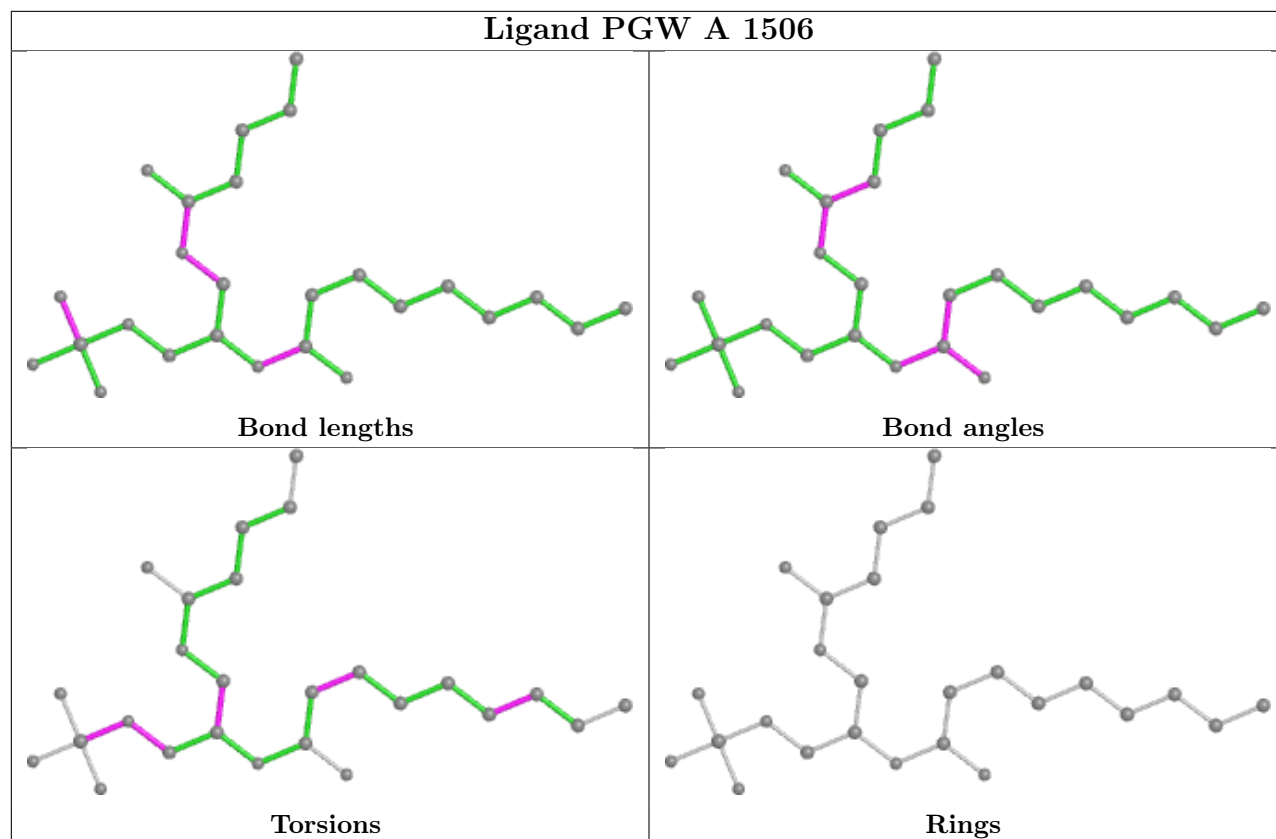




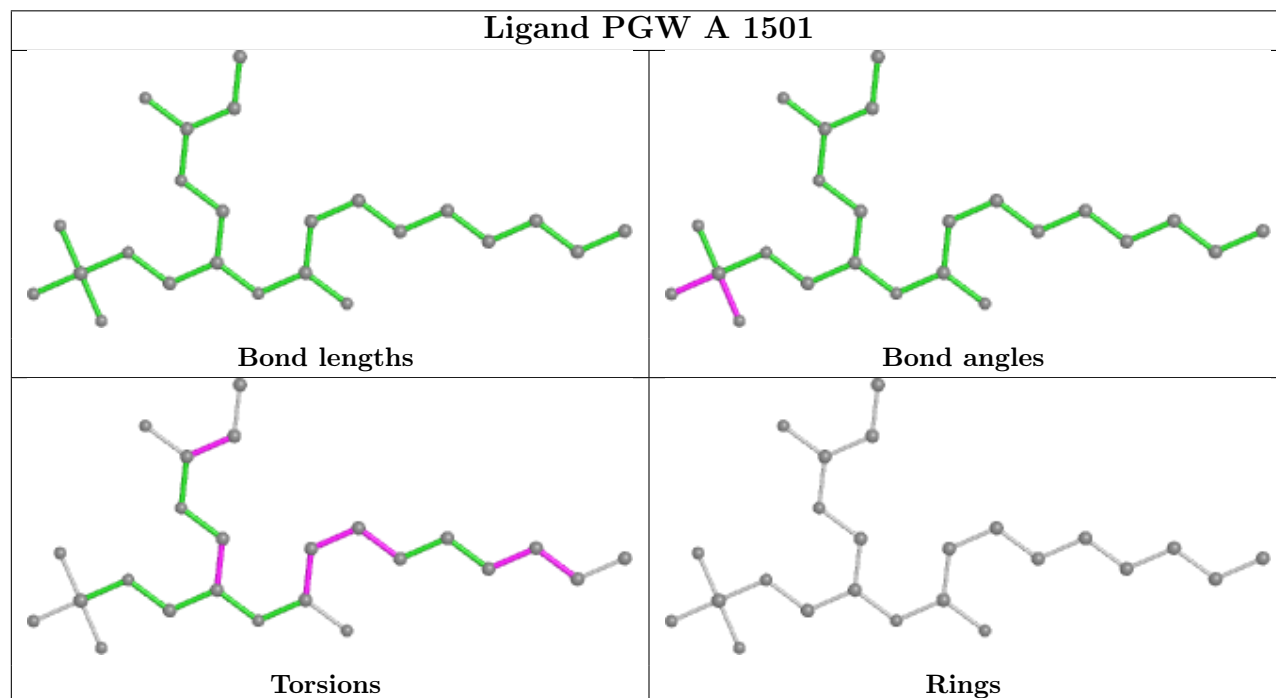


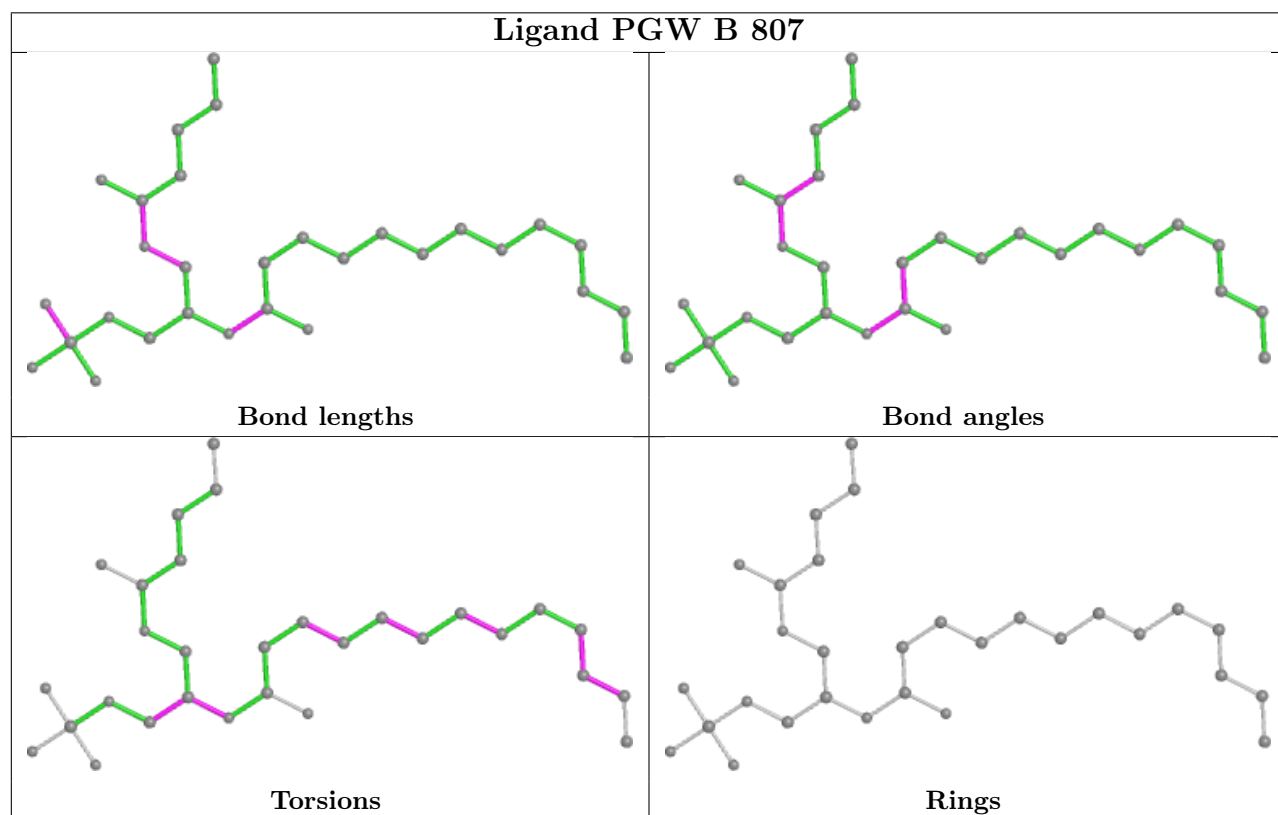
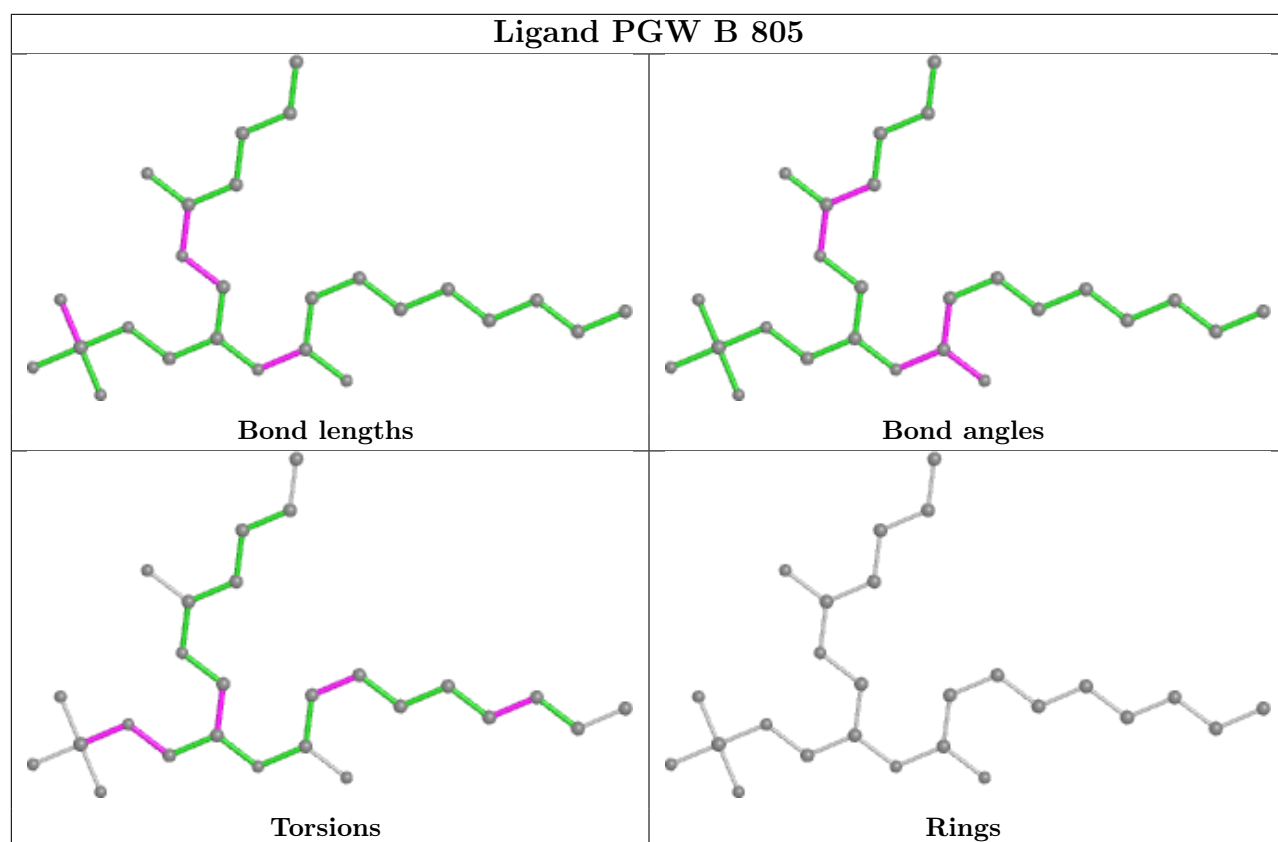


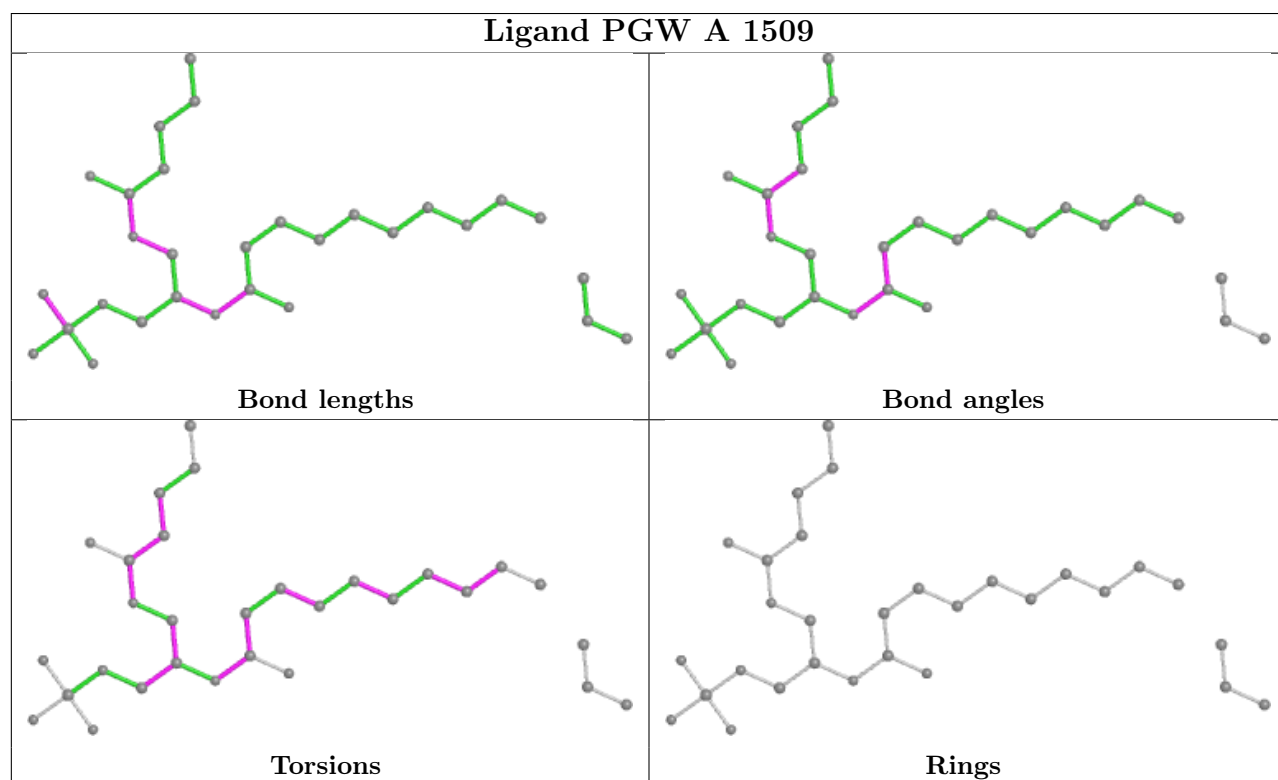
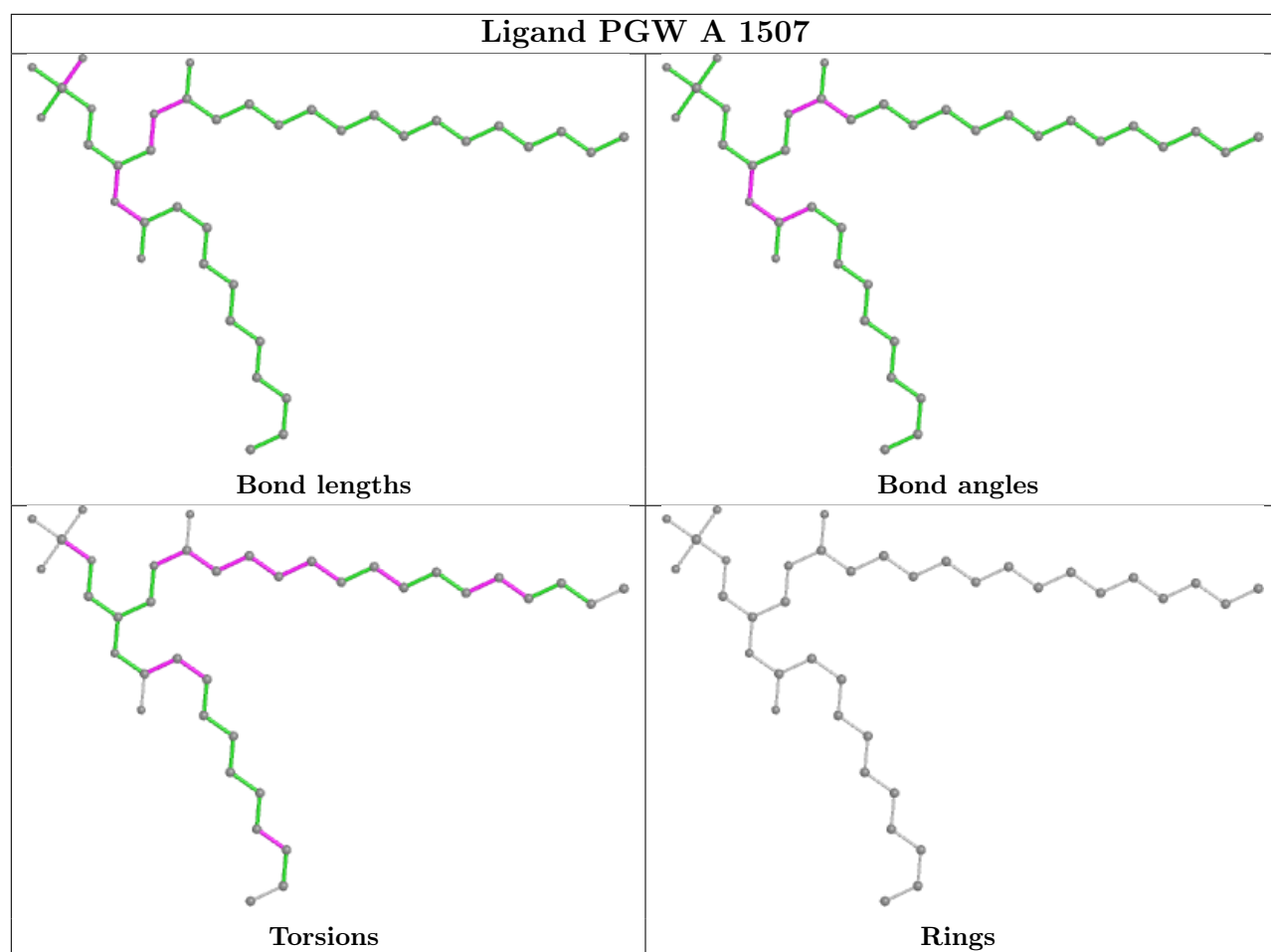
## Ligand PGW A 1506

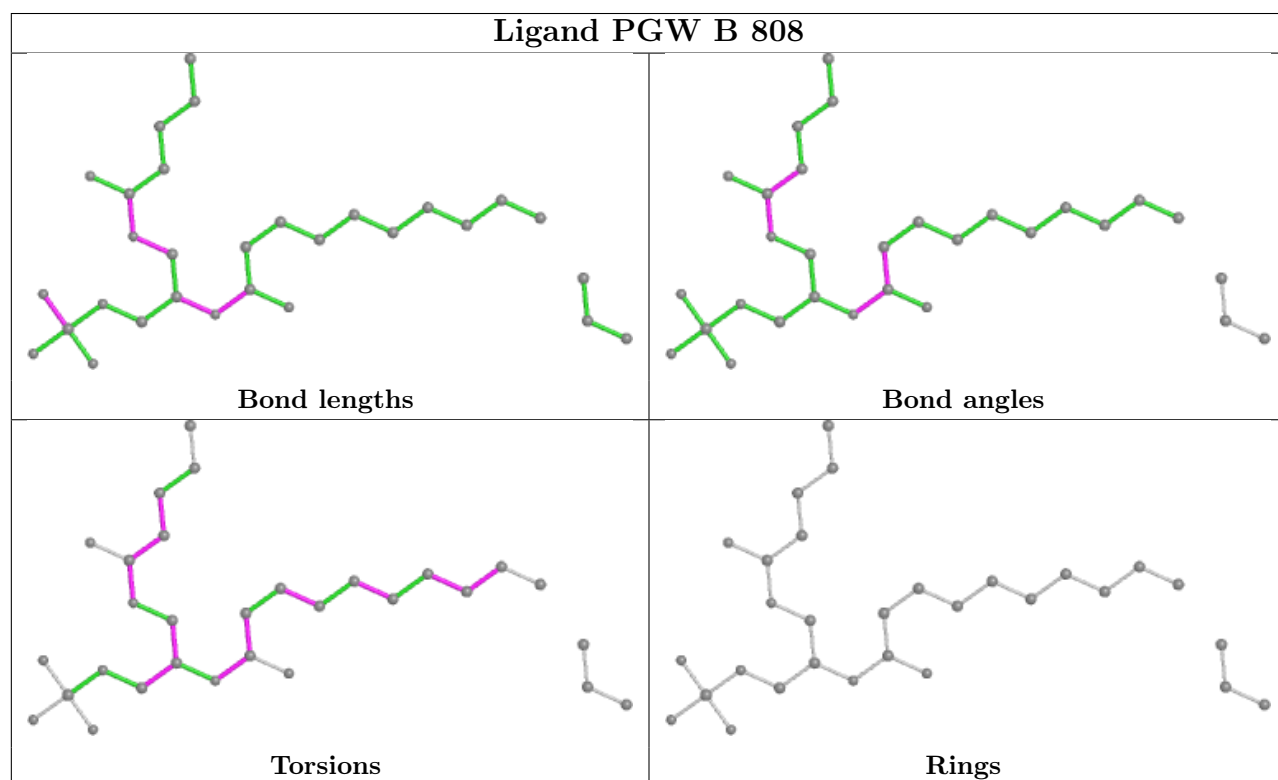
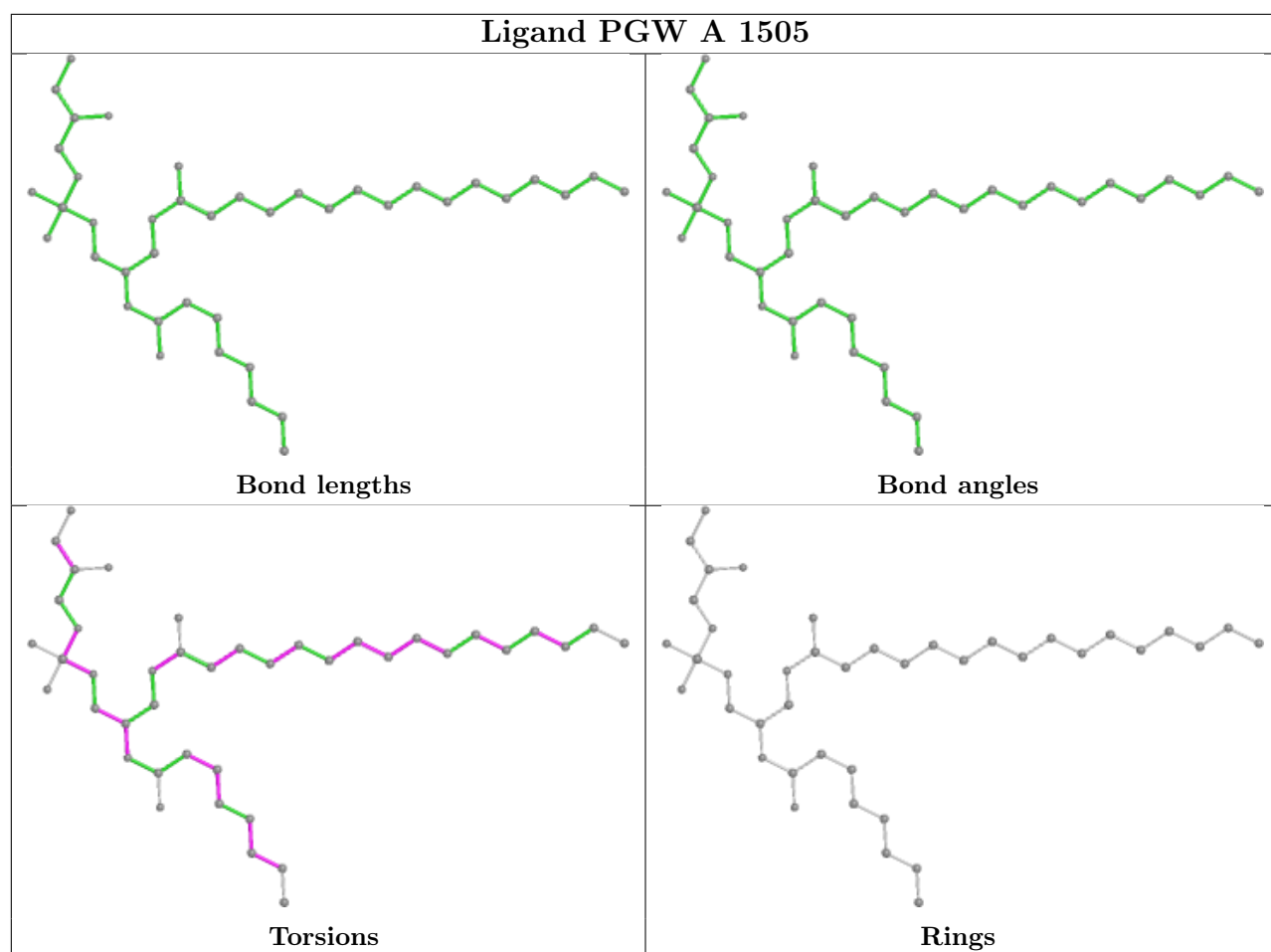


## Ligand PGW A 1501

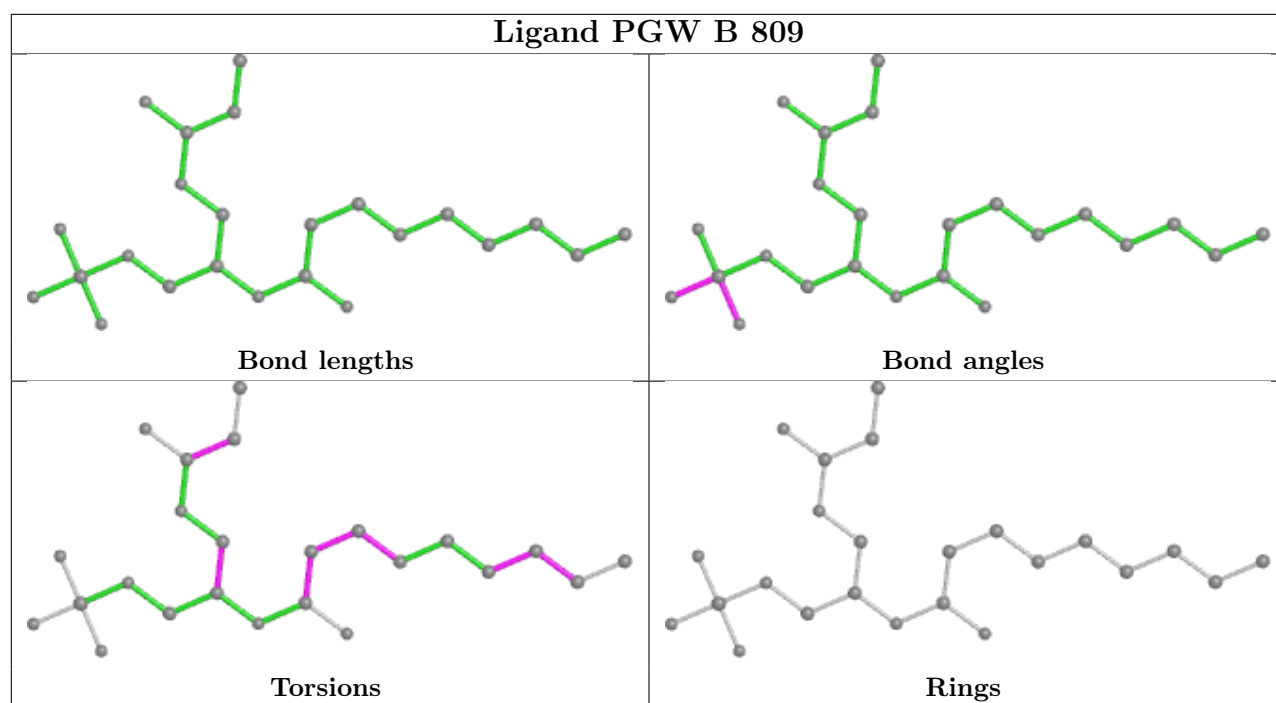












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

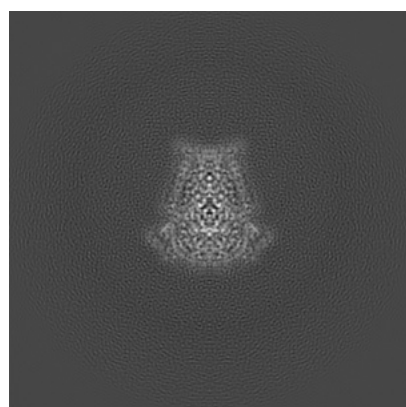
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24722. These allow visual inspection of the internal detail of the map and identification of artifacts.

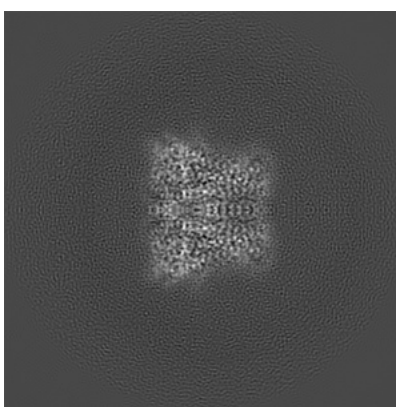
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

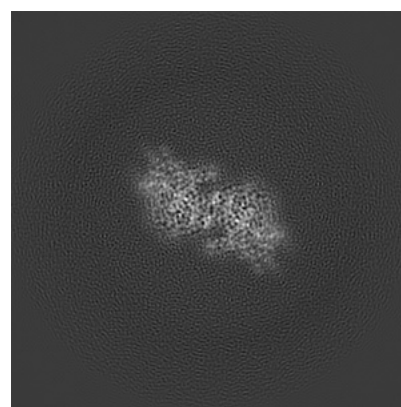
#### 6.1.1 Primary map



X



Y

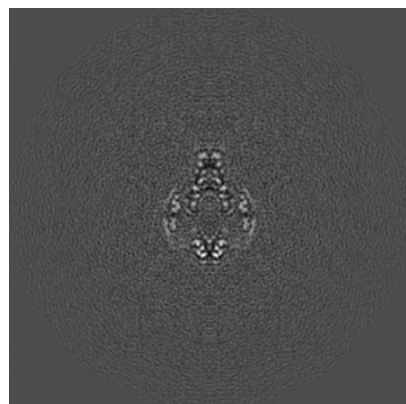


Z

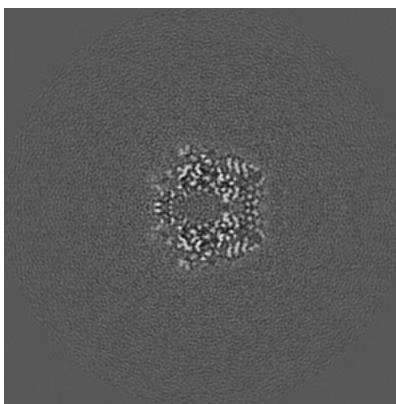
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

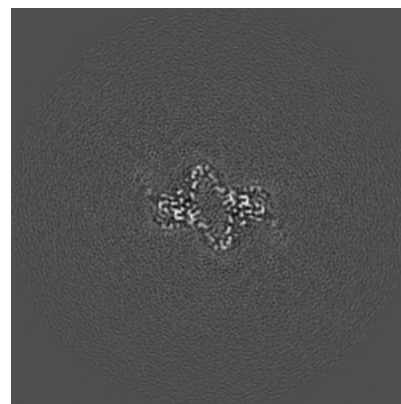
#### 6.2.1 Primary map



X Index: 192



Y Index: 192

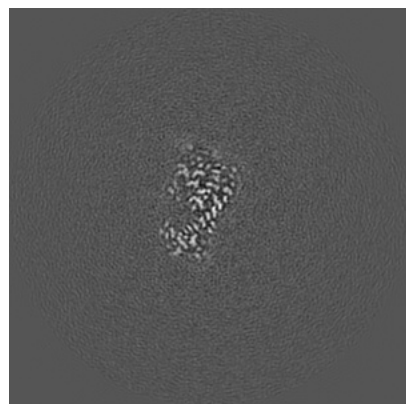


Z Index: 192

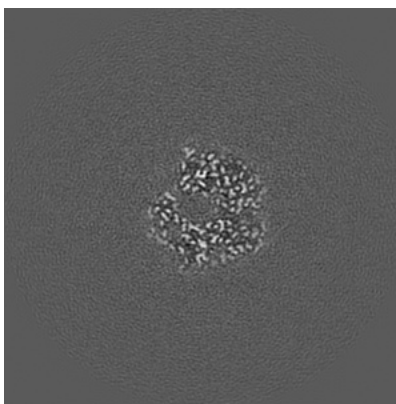
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

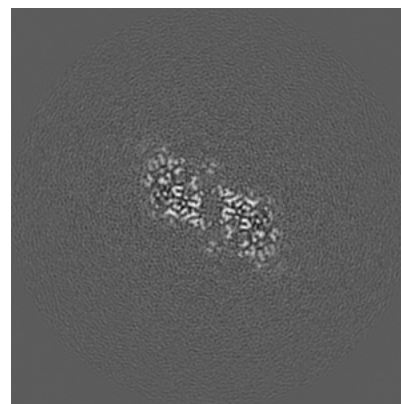
### 6.3.1 Primary map



X Index: 222



Y Index: 196

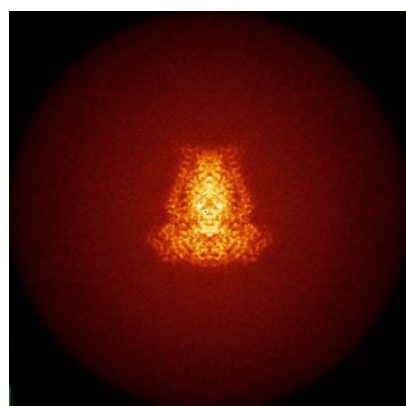


Z Index: 172

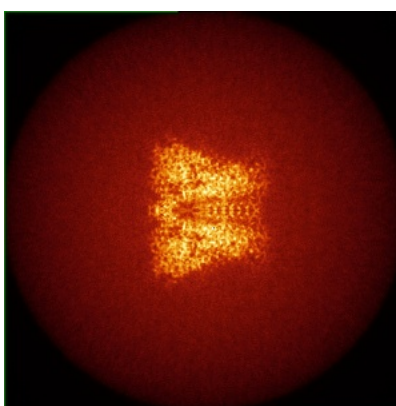
The images above show the largest variance slices of the map in three orthogonal directions.

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

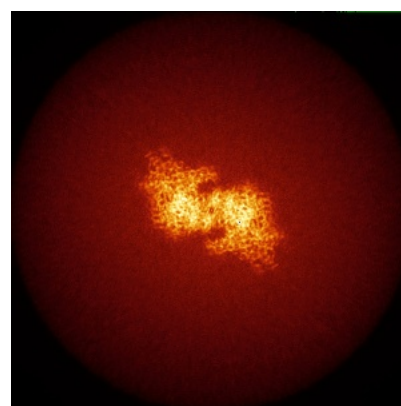
### 6.4.1 Primary map



X



Y

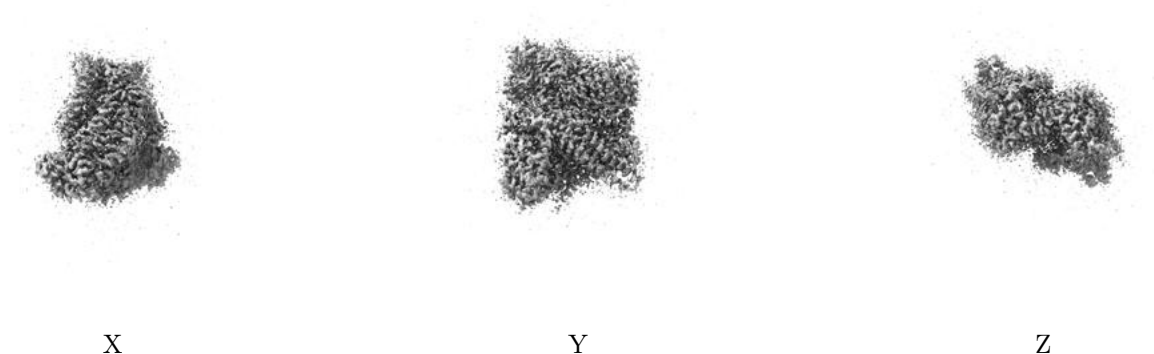


Z

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

## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.013. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

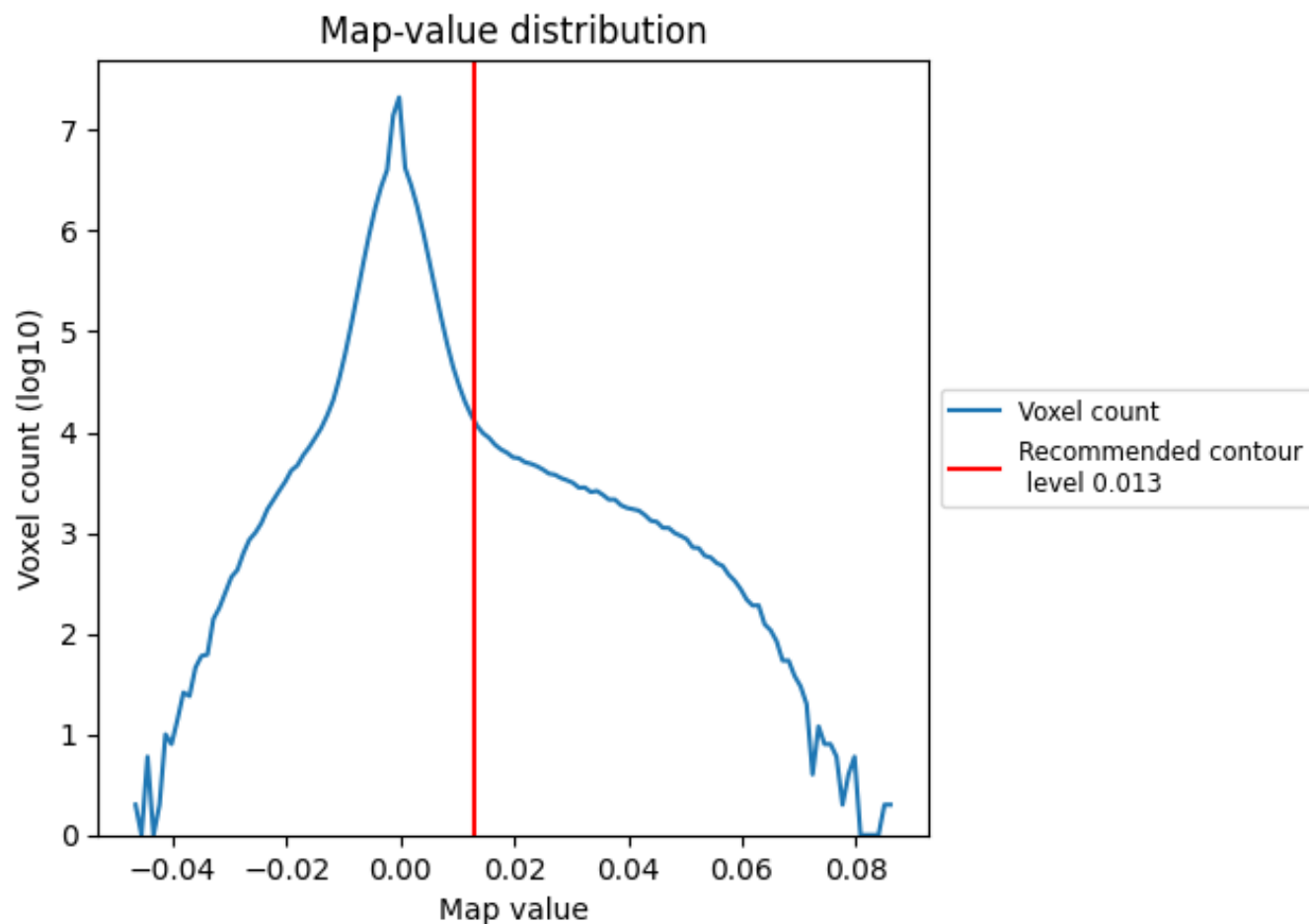
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

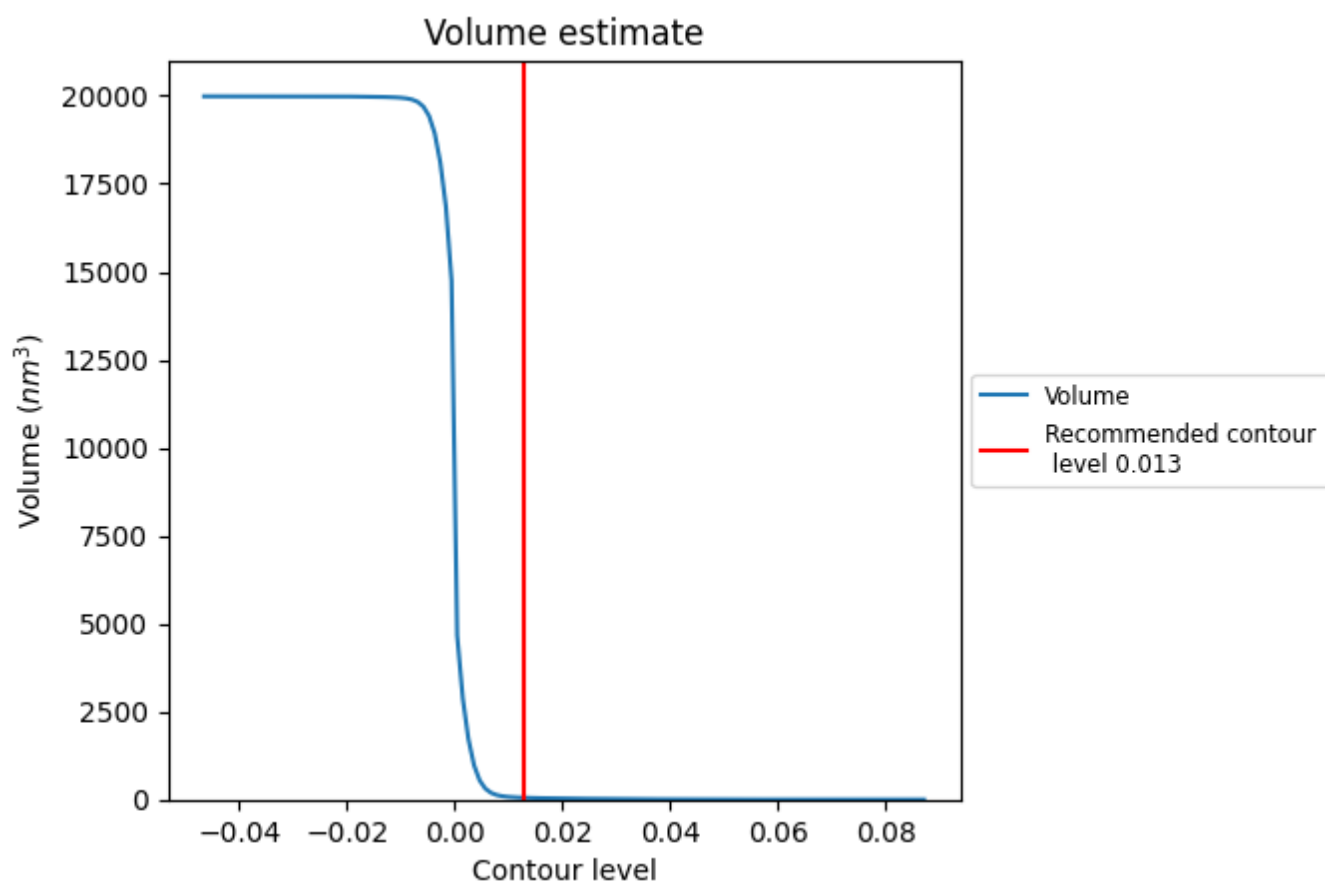
This section contains the results of statistical analysis of the map.

### 7.1 Map-value distribution [i](#)



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

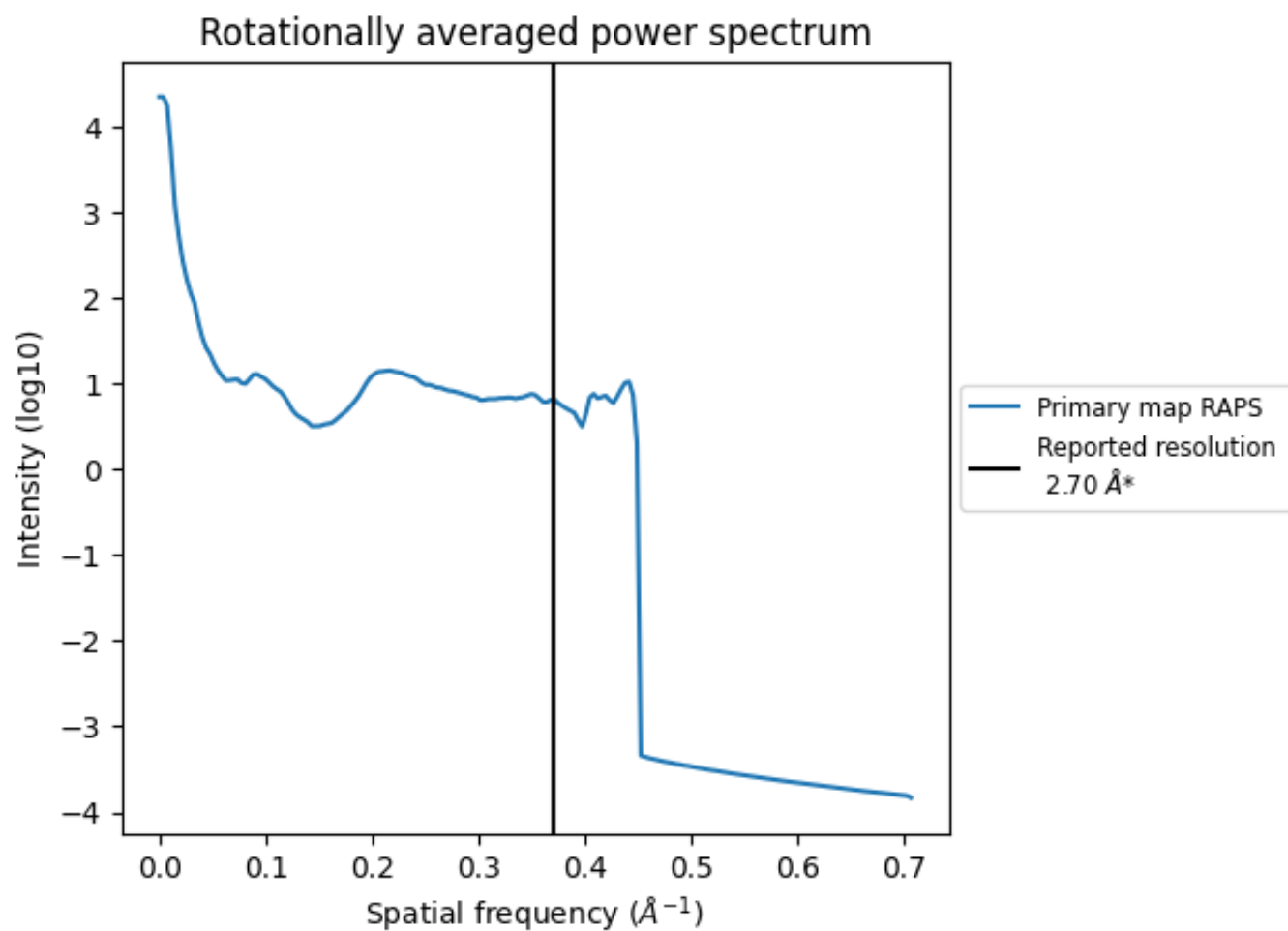
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 51 nm<sup>3</sup>; this corresponds to an approximate mass of 46 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 ⓘ

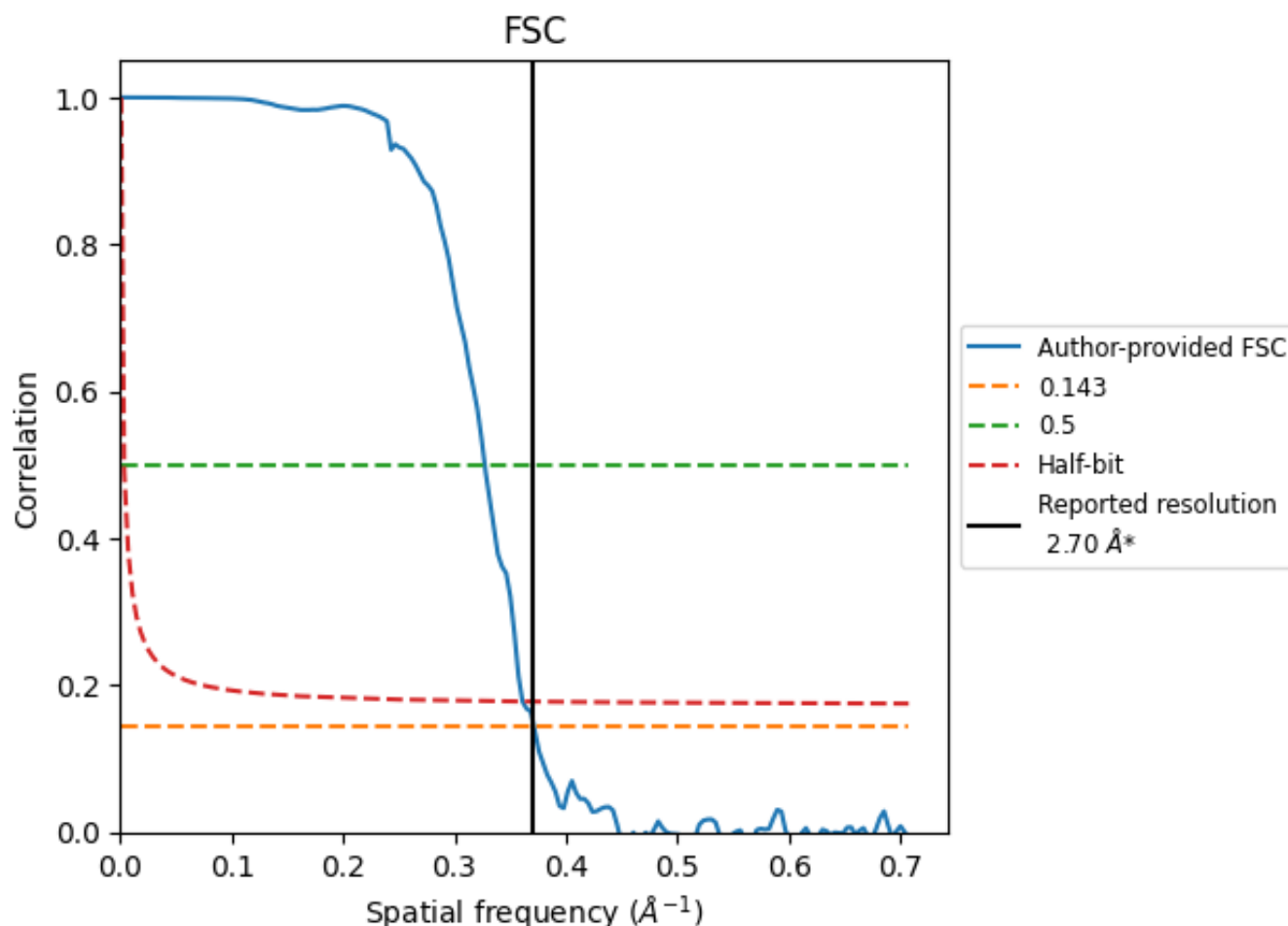


\*Reported resolution corresponds to spatial frequency of 0.370 Å<sup>-1</sup>

## 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 0.370 Å<sup>-1</sup>



## 8.2 Resolution estimates [i](#)

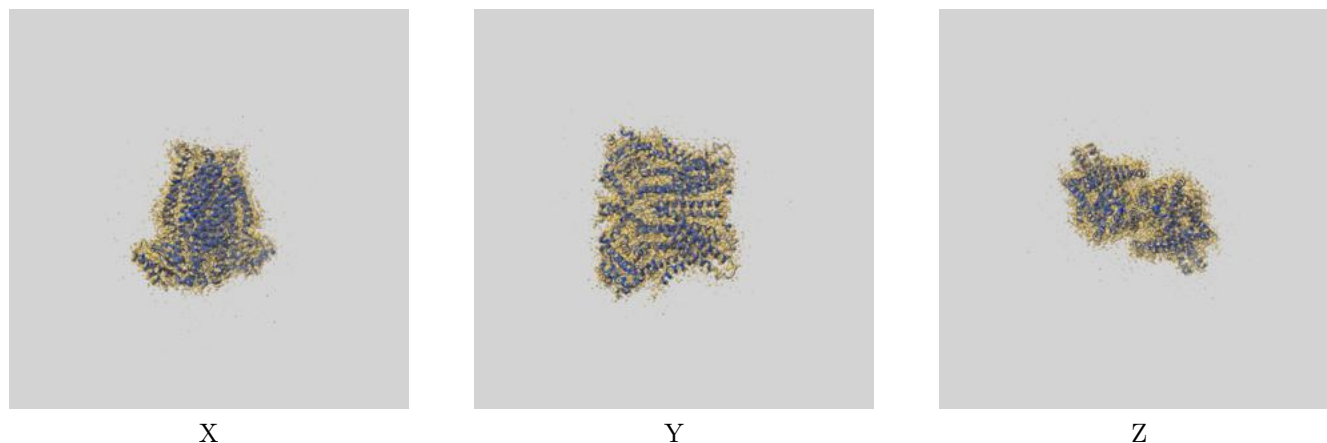
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.70	-	-
Author-provided FSC curve	2.69	3.06	2.77
Unmasked-calculated*	-	-	-

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

## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-24722 and PDB model 7RX2. Per-residue inclusion information can be found in section [3](#) on page [5](#).

### 9.1 Map-model overlay [i](#)



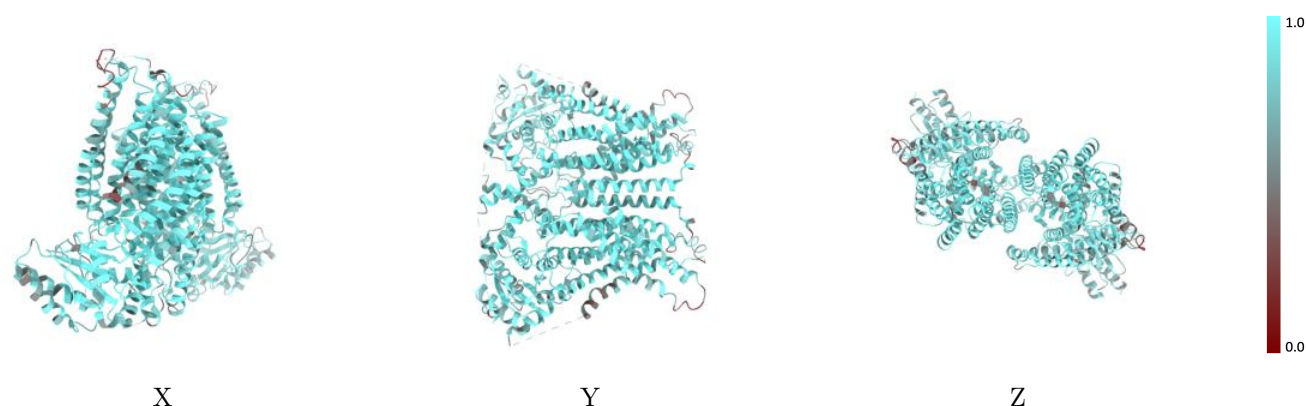
The images above show the 3D surface view of the map at the recommended contour level 0.013 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



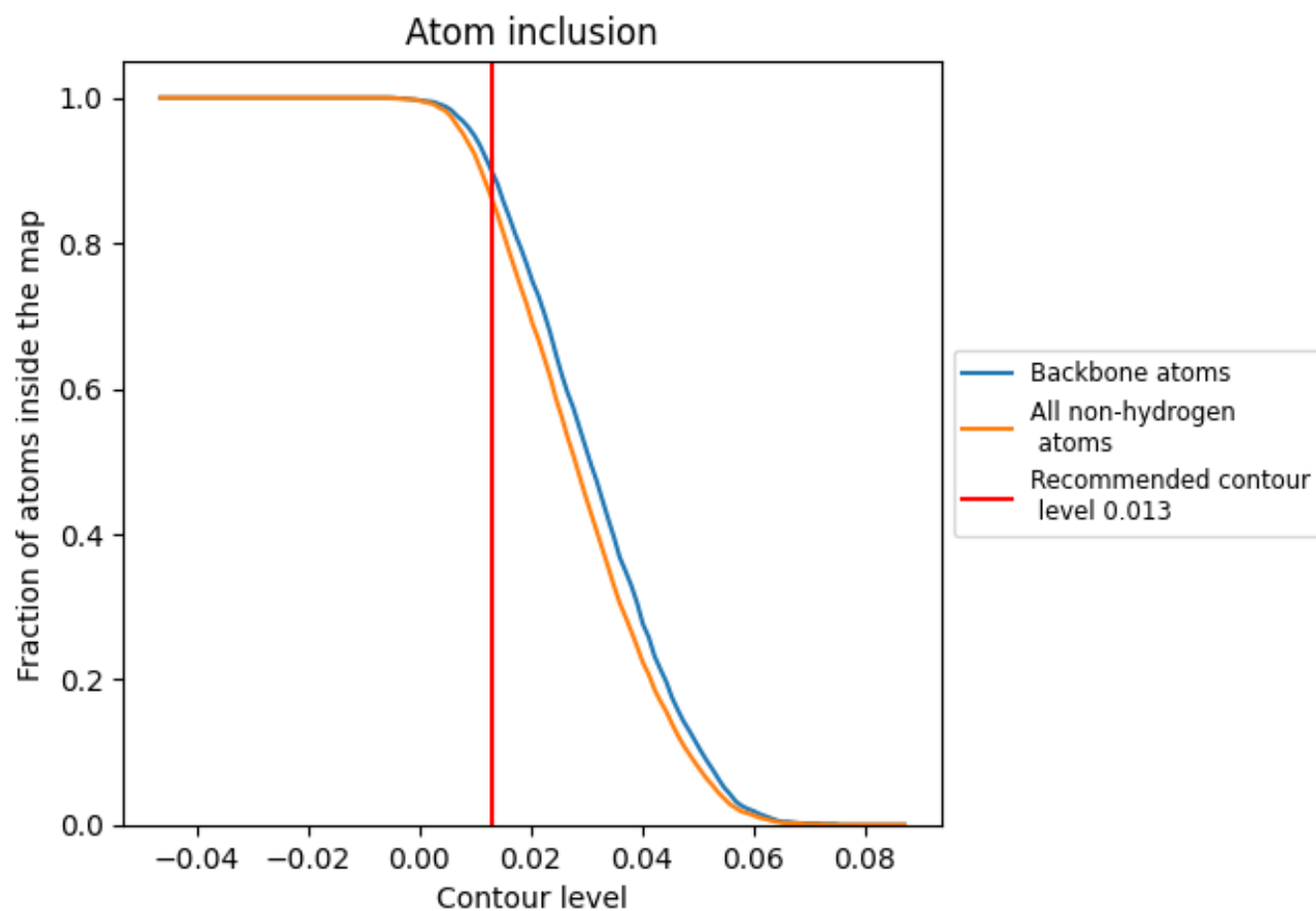
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.013).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 86% 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 (0.013) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.8620	<div><div></div></div> 0.6320
A	<div><div></div></div> 0.8630	<div><div></div></div> 0.6320
B	<div><div></div></div> 0.8620	<div><div></div></div> 0.6320

