



Full wwPDB EM Validation Report ⓘ

Mar 27, 2026 – 03:16 PM UTC

PDB ID : 7RXH / pdb_00007rxh
EMDB ID : EMD-24731
Title : afTMEM16 in C18 lipid nanodiscs with MSP1E3 scaffold protein in the presence of Ca²⁺, monomer with extra lipids
Authors : Falzone, M.E.; Accardi, A.
Deposited on : 2021-08-23
Resolution : 2.30 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

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

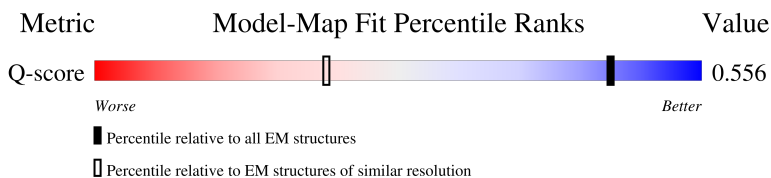
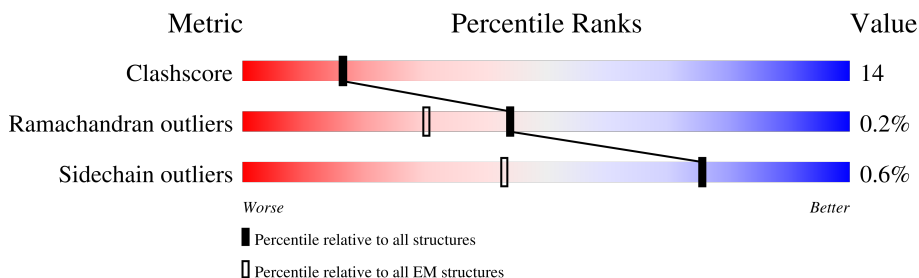
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.30 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	4254 (1.80 - 2.80)

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	735	 63% 22% 15%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5631 atoms, of which 48 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.

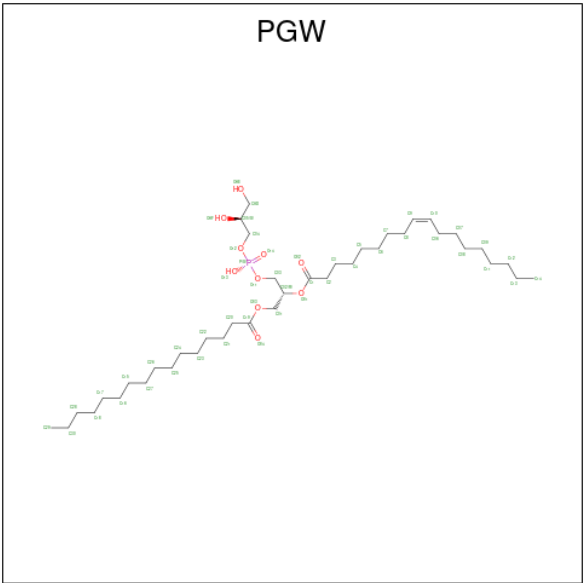
- Molecule 1 is a protein called afTMEM16 lipid scramblase.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	626	5062	3303	849	892	18	0	0

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
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 (CCD ID: PGW) (formula: C₄₀H₇₇O₁₀P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	O	P	
3	A	1	45	34	10	1	0
3	A	1	25	16	8	1	0

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Mol	Chain	Residues	Atoms				AltConf
3	A	1	Total	C	O	P	0
			34	25	8	1	
3	A	1	Total	C	O		0
			38	33	5		
3	A	1	Total	C	O	P	0
			31	22	8	1	
3	A	1	Total	C	O	P	0
			25	15	9	1	
3	A	1	Total	C	O	P	0
			32	23	8	1	
3	A	1	Total	C	O	P	0
			27	18	8	1	
3	A	1	Total	C	O	P	0
			28	19	8	1	
3	A	1	Total	C	O	P	0
			29	20	8	1	
3	A	1	Total	C	O	P	0
			44	33	10	1	
3	A	1	Total	C	O	P	0
			23	14	8	1	
3	A	1	Total	C	O	P	0
			28	19	8	1	
3	A	1	Total	C	O	P	0
			28	19	8	1	
3	A	1	Total	C	O	P	0
			22	13	8	1	
3	A	1	Total	C	O	P	0
			36	27	8	1	

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			AltConf
4	A	24	Total	H	O	0
			72	48	24	

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	1058829	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 ($e^-/\text{\AA}^2$)	42.8227	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	0.040	Depositor
Minimum map value	-0.015	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.0049	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: CA, PGW

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	0.22	0/5197	0.42	0/7070

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	5062	0	5018	135	0
2	A	2	0	0	0	0
3	A	495	0	558	52	0
4	A	24	48	0	4	0
All	All	5583	48	5576	157	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 14.

All (157) 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)
3:A:812:PGW:H23A	3:A:818:PGW:H2A	1.28	1.10
3:A:812:PGW:H23A	3:A:818:PGW:C2	1.84	1.06
1:A:604:THR:HG22	3:A:803:PGW:HAD	1.56	0.86
1:A:338:VAL:HG13	3:A:814:PGW:H2	1.59	0.85
1:A:349:TYR:OH	3:A:818:PGW:H03A	1.78	0.84
3:A:814:PGW:H20	3:A:814:PGW:H3A	1.57	0.83
1:A:109:LEU:HD21	1:A:562:ARG:HD2	1.60	0.83
3:A:814:PGW:H20	3:A:814:PGW:C3	2.12	0.79
1:A:304:ILE:HG22	1:A:324:PRO:HG3	1.65	0.78
1:A:393:VAL:HG23	1:A:394:PRO:HD3	1.65	0.77
1:A:452:LYS:HE2	1:A:456:PHE:CE2	2.20	0.77
1:A:389:ALA:HA	1:A:392:ILE:HB	1.70	0.74
3:A:812:PGW:H23A	3:A:818:PGW:H2	1.69	0.72
1:A:598:PRO:HG3	1:A:605:ILE:HD11	1.72	0.71
1:A:313:ASN:H	1:A:415:ALA:HB1	1.57	0.70
1:A:17:ASP:OD1	1:A:107:ARG:NH1	2.25	0.69
1:A:209:SER:HA	3:A:808:PGW:H04	1.75	0.69
1:A:452:LYS:HE2	1:A:456:PHE:HE2	1.56	0.69
1:A:184:LEU:HD23	1:A:544:LEU:HD23	1.75	0.68
1:A:291:ALA:HA	1:A:372:ILE:HD13	1.77	0.66
1:A:496:ARG:NH1	1:A:500:GLU:OE2	2.29	0.66
3:A:805:PGW:H02	3:A:806:PGW:H01A	1.78	0.66
3:A:814:PGW:H4A	3:A:814:PGW:H22	1.78	0.66
1:A:318:GLY:HA2	1:A:321:VAL:HG22	1.79	0.65
1:A:21:ARG:NH2	1:A:501:ALA:O	2.29	0.65
3:A:812:PGW:C23	3:A:818:PGW:H2A	2.19	0.64
1:A:189:ARG:HD3	3:A:810:PGW:H03A	1.79	0.63
1:A:294:LEU:HD11	1:A:369:VAL:HG13	1.80	0.63
1:A:228:LYS:HG2	1:A:570:LEU:HD21	1.79	0.63
1:A:452:LYS:CE	1:A:456:PHE:CE2	2.83	0.62
1:A:539:ASN:HB2	3:A:807:PGW:H09	1.81	0.62
1:A:310:GLU:OE2	1:A:425:ARG:NH2	2.31	0.62
1:A:52:VAL:CG1	1:A:60:LEU:HD23	2.30	0.62
1:A:109:LEU:CD2	1:A:562:ARG:HD2	2.28	0.62
1:A:313:ASN:N	1:A:415:ALA:O	2.34	0.60
1:A:338:VAL:CG1	3:A:814:PGW:H5	2.33	0.59
1:A:610:LEU:HD21	3:A:806:PGW:H25A	1.84	0.59
1:A:583:THR:HG22	3:A:806:PGW:H18A	1.84	0.58
1:A:119:LYS:HA	1:A:123:GLY:O	2.03	0.58
1:A:284:VAL:HB	1:A:285:PRO:HD3	1.86	0.58
1:A:29:THR:O	1:A:33:ILE:HG12	2.04	0.58
1:A:273:ALA:O	1:A:277:MET:HG3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:VAL:CG1	3:A:814:PGW:H2	2.33	0.56
1:A:305:GLU:OE1	1:A:425:ARG:NH1	2.36	0.55
1:A:253:PHE:CZ	1:A:353:GLU:HG2	2.41	0.55
1:A:151:CYS:HB3	1:A:155:TRP:CZ2	2.42	0.54
1:A:593:ASN:HD21	1:A:597:GLY:HA2	1.73	0.53
3:A:815:PGW:O02	3:A:815:PGW:H03	2.08	0.52
1:A:323:ILE:HB	1:A:324:PRO:HD3	1.92	0.52
1:A:386:VAL:HG12	1:A:598:PRO:HG2	1.91	0.52
1:A:607:CYS:O	1:A:611:LEU:HD23	2.10	0.52
1:A:188:PHE:CE1	1:A:541:TRP:HB2	2.44	0.51
1:A:494:LEU:O	1:A:498:ARG:HG3	2.10	0.51
1:A:21:ARG:NH1	1:A:137:ASP:OD2	2.42	0.51
3:A:812:PGW:O14	3:A:812:PGW:H01A	2.10	0.51
1:A:54:GLN:NE2	1:A:56:ASP:O	2.28	0.51
1:A:620:LEU:HG	3:A:803:PGW:H28	1.92	0.51
1:A:21:ARG:HD2	1:A:137:ASP:HB3	1.93	0.50
1:A:448:VAL:CG2	1:A:449:PRO:HD3	2.41	0.50
1:A:316:LEU:O	1:A:320:LEU:HG	2.11	0.50
1:A:180:TYR:CZ	1:A:511:ASP:HB3	2.46	0.50
1:A:607:CYS:SG	3:A:805:PGW:H2A	2.52	0.50
1:A:94:GLU:HB3	1:A:95:PRO:HD2	1.93	0.49
1:A:317:LYS:O	1:A:321:VAL:HG13	2.12	0.49
3:A:809:PGW:H24A	3:A:810:PGW:O04	2.13	0.49
1:A:161:LEU:N	1:A:185:GLN:OE1	2.43	0.49
1:A:227:TRP:CE2	1:A:231:GLU:HG3	2.48	0.49
1:A:24:TYR:CE2	1:A:60:LEU:HD12	2.47	0.49
1:A:52:VAL:HG11	1:A:60:LEU:HD23	1.93	0.49
1:A:444:LEU:HA	1:A:448:VAL:HG22	1.93	0.49
1:A:211:ILE:HD13	3:A:808:PGW:H6A	1.95	0.49
1:A:313:ASN:N	1:A:415:ALA:HB1	2.27	0.49
1:A:250:ARG:HD3	4:A:2701:HOH:O	2.12	0.49
1:A:393:VAL:CG2	1:A:394:PRO:HD3	2.40	0.48
1:A:613:THR:CG2	3:A:803:PGW:H22	2.43	0.48
1:A:516:CYS:SG	1:A:569:TRP:HB3	2.54	0.48
1:A:386:VAL:HB	1:A:387:PRO:HD3	1.95	0.48
1:A:104:GLU:HG3	1:A:107:ARG:NH2	2.29	0.48
1:A:208:PHE:O	3:A:808:PGW:HADA	2.14	0.48
1:A:151:CYS:HB3	1:A:155:TRP:CH2	2.48	0.48
1:A:375:TYR:OH	1:A:520:GLY:HA3	2.14	0.48
3:A:818:PGW:H25A	3:A:818:PGW:H15A	1.46	0.47
1:A:190:PHE:O	1:A:194:PRO:HD2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:ILE:HG22	1:A:228:LYS:HE2	1.97	0.46
1:A:276:ARG:HH11	1:A:351:ASN:HB3	1.79	0.46
1:A:432:TYR:CE2	3:A:816:PGW:H23A	2.50	0.46
1:A:453:GLN:HE22	1:A:554:GLU:HA	1.81	0.46
1:A:320:LEU:HD22	1:A:323:ILE:HD11	1.97	0.46
1:A:611:LEU:HD21	3:A:805:PGW:H6A	1.98	0.45
3:A:814:PGW:H20	3:A:814:PGW:H3	1.95	0.45
1:A:641:ARG:O	1:A:645:ILE:HG12	2.16	0.45
3:A:812:PGW:C24	3:A:818:PGW:H4A	2.47	0.45
1:A:619:HIS:HB3	3:A:813:PGW:H06A	1.98	0.45
1:A:338:VAL:HG11	3:A:814:PGW:H5	1.98	0.45
1:A:320:LEU:HD13	1:A:403:VAL:HB	2.00	0.44
1:A:338:VAL:HG11	3:A:814:PGW:C5	2.47	0.44
1:A:254:LYS:O	1:A:351:ASN:ND2	2.48	0.44
1:A:289:LEU:HD22	3:A:814:PGW:C23	2.46	0.44
1:A:302:PHE:O	1:A:306:ILE:HG12	2.18	0.44
1:A:212:TYR:OH	1:A:526:SER:HB2	2.17	0.44
1:A:294:LEU:CD1	1:A:369:VAL:HG13	2.47	0.44
1:A:425:ARG:HD2	4:A:2722:HOH:O	2.17	0.44
1:A:613:THR:HG23	3:A:803:PGW:H22	2.00	0.44
1:A:197:PHE:HD1	3:A:809:PGW:H30A	1.81	0.44
1:A:240:THR:O	1:A:243:VAL:HB	2.18	0.44
1:A:438:GLN:HE22	1:A:518:GLN:HE21	1.66	0.44
1:A:593:ASN:ND2	1:A:597:GLY:HA2	2.32	0.44
1:A:604:THR:CG2	3:A:803:PGW:HAD	2.37	0.44
1:A:324:PRO:O	1:A:328:VAL:HG23	2.18	0.44
1:A:424:ASP:HB3	1:A:428:LYS:HE2	1.99	0.43
1:A:194:PRO:HD3	1:A:219:TRP:CD1	2.53	0.43
1:A:150:GLN:OE1	1:A:153:ARG:NH1	2.41	0.43
1:A:231:GLU:OE2	1:A:566:ILE:HG12	2.18	0.43
1:A:92:GLU:OE2	1:A:653:ARG:HD2	2.19	0.43
1:A:209:SER:CA	3:A:808:PGW:H04	2.45	0.43
1:A:318:GLY:HA2	1:A:321:VAL:CG2	2.48	0.43
1:A:625:ARG:NH2	4:A:2704:HOH:O	2.29	0.43
3:A:818:PGW:H06A	3:A:818:PGW:H8	1.74	0.43
1:A:237:ARG:HA	1:A:642:ARG:HH12	1.84	0.43
1:A:318:GLY:CA	1:A:321:VAL:HG22	2.49	0.42
1:A:23:ASN:HD22	1:A:59:SER:HB3	1.85	0.42
1:A:83:ASP:OD2	1:A:562:ARG:NH2	2.52	0.42
1:A:113:HIS:O	1:A:117:VAL:HG22	2.19	0.42
1:A:210:ILE:HG12	1:A:588:VAL:HG11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:ASN:ND2	1:A:362:LEU:HB2	2.34	0.42
1:A:451:VAL:O	1:A:455:VAL:HG23	2.20	0.42
3:A:805:PGW:O04	3:A:805:PGW:H03	2.19	0.42
1:A:285:PRO:HD3	3:A:818:PGW:H06A	2.02	0.42
1:A:349:TYR:HH	3:A:818:PGW:H03A	1.79	0.42
3:A:805:PGW:H01A	3:A:806:PGW:H02	2.01	0.42
1:A:291:ALA:HA	1:A:372:ILE:CD1	2.48	0.42
1:A:33:ILE:O	1:A:37:GLU:HG3	2.18	0.42
1:A:619:HIS:CD2	3:A:813:PGW:H6	2.55	0.42
1:A:619:HIS:CB	3:A:813:PGW:H06A	2.49	0.42
3:A:809:PGW:C2	3:A:809:PGW:H01	2.49	0.42
1:A:144:ASP:OD2	1:A:147:THR:OG1	2.31	0.41
1:A:282:LEU:C	1:A:285:PRO:HD2	2.45	0.41
1:A:314:GLY:N	1:A:415:ALA:HB1	2.35	0.41
1:A:494:LEU:HD23	1:A:494:LEU:HA	1.86	0.41
3:A:816:PGW:H20A	3:A:816:PGW:H23	1.71	0.41
1:A:54:GLN:O	1:A:498:ARG:NH1	2.48	0.41
1:A:209:SER:OG	1:A:211:ILE:HG22	2.21	0.41
3:A:806:PGW:H26A	3:A:806:PGW:H16	1.83	0.41
1:A:341:THR:O	1:A:345:LYS:HD3	2.20	0.41
1:A:518:GLN:HG2	1:A:536:PHE:CE2	2.56	0.41
3:A:805:PGW:C01	3:A:806:PGW:H02	2.51	0.41
1:A:261:ASP:OD1	1:A:261:ASP:N	2.51	0.41
1:A:332:ILE:HD13	1:A:332:ILE:HA	1.88	0.41
1:A:370:ASN:HD21	1:A:513:ARG:HE	1.69	0.40
1:A:20:ILE:HB	1:A:62:VAL:HB	2.03	0.40
1:A:214:VAL:HG22	3:A:803:PGW:H15	2.03	0.40
1:A:444:LEU:HA	1:A:448:VAL:CG2	2.51	0.40
1:A:448:VAL:HG23	1:A:449:PRO:HD3	2.03	0.40
1:A:23:ASN:ND2	1:A:59:SER:HB3	2.37	0.40
1:A:207:SER:OG	1:A:595:HIS:NE2	2.51	0.40
1:A:425:ARG:HD3	4:A:2717:HOH:O	2.22	0.40
3:A:811:PGW:H20A	3:A:811:PGW:H01A	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	622/735 (85%)	604 (97%)	17 (3%)	1 (0%)	43	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	420	SER

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	540/646 (84%)	537 (99%)	3 (1%)	78	89

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	403	VAL
1	A	417	THR
1	A	421	ILE

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

Mol	Chain	Res	Type
1	A	23	ASN
1	A	148	ASN

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Mol	Chain	Res	Type
1	A	171	ASN
1	A	230	GLN
1	A	283	GLN
1	A	370	ASN
1	A	400	HIS
1	A	453	GLN
1	A	518	GLN
1	A	539	ASN
1	A	540	ASN
1	A	593	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 ⓘ

Of 18 ligands modelled in this entry, 2 are monoatomic - leaving 16 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	A	816	-	27,27,50	0.47	0	30,32,56	0.60	1 (3%)
3	PGW	A	807	-	30,30,50	1.23	5 (16%)	33,35,56	1.23	2 (6%)
3	PGW	A	808	-	24,24,50	1.07	2 (8%)	26,29,56	1.04	1 (3%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PGW	A	818	-	35,35,50	1.14	5 (14%)	38,40,56	1.19	2 (5%)
3	PGW	A	810	-	26,26,50	1.31	4 (15%)	29,31,56	1.32	2 (6%)
3	PGW	A	804	-	24,24,50	1.33	5 (20%)	27,29,56	1.35	3 (11%)
3	PGW	A	809	-	31,31,50	1.18	4 (12%)	34,36,56	1.23	2 (5%)
3	PGW	A	803	-	44,44,50	0.99	4 (9%)	47,50,56	1.13	2 (4%)
3	PGW	A	806	-	37,37,50	0.99	4 (10%)	39,39,56	1.21	2 (5%)
3	PGW	A	811	-	27,27,50	1.26	5 (18%)	30,32,56	1.27	2 (6%)
3	PGW	A	814	-	22,22,50	1.39	5 (22%)	25,27,56	1.31	2 (8%)
3	PGW	A	812	-	28,28,50	1.26	5 (17%)	31,33,56	1.27	2 (6%)
3	PGW	A	813	-	43,43,50	1.02	4 (9%)	46,49,56	1.11	2 (4%)
3	PGW	A	805	-	33,33,50	1.16	5 (15%)	36,38,56	1.21	2 (5%)
3	PGW	A	817	-	21,21,50	1.42	5 (23%)	24,26,56	1.34	2 (8%)
3	PGW	A	815	-	27,27,50	0.47	0	30,32,56	0.59	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	A	816	-	-	12/29/29/55	-
3	PGW	A	807	-	-	10/32/32/55	-
3	PGW	A	808	-	-	10/28/28/55	-
3	PGW	A	818	-	-	18/37/37/55	-
3	PGW	A	810	-	-	15/28/28/55	-
3	PGW	A	804	-	-	11/26/26/55	-
3	PGW	A	809	-	-	18/33/33/55	-
3	PGW	A	803	-	-	21/49/49/55	-
3	PGW	A	806	-	-	14/39/39/55	-
3	PGW	A	811	-	-	16/29/29/55	-
3	PGW	A	814	-	-	10/24/24/55	-
3	PGW	A	812	-	-	17/30/30/55	-
3	PGW	A	813	-	-	21/48/48/55	-
3	PGW	A	805	-	-	16/35/35/55	-
3	PGW	A	817	-	-	11/23/23/55	-
3	PGW	A	815	-	-	17/29/29/55	-

All (62) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	810	PGW	O01-C1	3.12	1.43	1.34
3	A	807	PGW	O01-C1	2.96	1.42	1.34
3	A	809	PGW	O01-C1	2.96	1.42	1.34
3	A	817	PGW	O01-C1	2.92	1.42	1.34
3	A	814	PGW	O01-C1	2.89	1.42	1.34
3	A	818	PGW	O01-C1	2.88	1.42	1.34
3	A	806	PGW	O01-C1	2.88	1.42	1.34
3	A	805	PGW	O01-C1	2.84	1.42	1.34
3	A	811	PGW	O01-C1	2.81	1.42	1.34
3	A	804	PGW	O01-C1	2.80	1.42	1.34
3	A	803	PGW	O01-C1	2.79	1.42	1.34
3	A	808	PGW	O01-C1	2.78	1.42	1.34
3	A	812	PGW	O01-C1	2.74	1.42	1.34
3	A	813	PGW	O01-C1	2.69	1.41	1.34
3	A	813	PGW	O03-C19	2.58	1.40	1.33
3	A	818	PGW	O03-C19	2.47	1.40	1.33
3	A	807	PGW	P-O12	2.47	1.64	1.54
3	A	809	PGW	O03-C19	2.46	1.40	1.33
3	A	810	PGW	O03-C19	2.46	1.40	1.33
3	A	814	PGW	O03-C19	2.45	1.40	1.33
3	A	804	PGW	P-O12	2.44	1.63	1.54
3	A	817	PGW	O03-C19	2.44	1.40	1.33
3	A	812	PGW	P-O12	2.44	1.63	1.54
3	A	814	PGW	P-O12	2.44	1.63	1.54
3	A	805	PGW	P-O12	2.44	1.63	1.54
3	A	804	PGW	O03-C01	-2.43	1.39	1.45
3	A	807	PGW	O03-C19	2.43	1.40	1.33
3	A	817	PGW	P-O12	2.43	1.63	1.54
3	A	818	PGW	P-O12	2.42	1.63	1.54
3	A	810	PGW	P-O12	2.41	1.63	1.54
3	A	811	PGW	P-O12	2.41	1.63	1.54
3	A	809	PGW	P-O12	2.41	1.63	1.54
3	A	809	PGW	O03-C01	-2.38	1.39	1.45
3	A	805	PGW	O03-C19	2.37	1.40	1.33
3	A	803	PGW	O03-C19	2.37	1.40	1.33
3	A	805	PGW	O03-C01	-2.37	1.39	1.45
3	A	804	PGW	O03-C19	2.36	1.40	1.33
3	A	806	PGW	O03-C01	-2.36	1.39	1.45
3	A	812	PGW	O03-C19	2.36	1.40	1.33
3	A	811	PGW	O03-C01	-2.34	1.39	1.45
3	A	811	PGW	O03-C19	2.33	1.40	1.33
3	A	807	PGW	O03-C01	-2.32	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	812	PGW	O03-C01	-2.32	1.40	1.45
3	A	814	PGW	O03-C01	-2.32	1.40	1.45
3	A	818	PGW	O03-C01	-2.31	1.40	1.45
3	A	803	PGW	O03-C01	-2.31	1.40	1.45
3	A	817	PGW	O03-C01	-2.30	1.40	1.45
3	A	810	PGW	O03-C01	-2.30	1.40	1.45
3	A	806	PGW	O03-C19	2.29	1.40	1.33
3	A	813	PGW	O03-C01	-2.26	1.40	1.45
3	A	813	PGW	O01-C02	-2.22	1.41	1.46
3	A	812	PGW	O01-C02	-2.19	1.41	1.46
3	A	811	PGW	O01-C02	-2.14	1.41	1.46
3	A	804	PGW	O01-C02	-2.09	1.41	1.46
3	A	814	PGW	O01-C02	-2.07	1.41	1.46
3	A	805	PGW	O01-C02	-2.07	1.41	1.46
3	A	803	PGW	O01-C02	-2.07	1.41	1.46
3	A	818	PGW	O01-C02	-2.07	1.41	1.46
3	A	806	PGW	O01-C02	-2.06	1.41	1.46
3	A	817	PGW	O01-C02	-2.06	1.41	1.46
3	A	807	PGW	O01-C02	-2.06	1.41	1.46
3	A	808	PGW	O01-C02	-2.04	1.41	1.46

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	810	PGW	O01-C1-C2	4.16	120.49	111.48
3	A	804	PGW	O01-C1-C2	4.12	120.39	111.48
3	A	806	PGW	O01-C1-C2	4.11	120.37	111.48
3	A	805	PGW	O01-C1-C2	4.10	120.35	111.48
3	A	813	PGW	O01-C1-C2	4.05	120.25	111.48
3	A	807	PGW	O01-C1-C2	4.04	120.22	111.48
3	A	812	PGW	O01-C1-C2	4.04	120.22	111.48
3	A	803	PGW	O01-C1-C2	4.02	120.19	111.48
3	A	809	PGW	O01-C1-C2	3.98	120.09	111.48
3	A	818	PGW	O01-C1-C2	3.98	120.09	111.48
3	A	811	PGW	O01-C1-C2	3.97	120.07	111.48
3	A	817	PGW	O01-C1-C2	3.95	120.03	111.48
3	A	814	PGW	O01-C1-C2	3.87	119.85	111.48
3	A	808	PGW	O01-C1-C2	3.79	119.68	111.48
3	A	810	PGW	O03-C19-C20	2.94	120.79	111.83
3	A	807	PGW	O03-C19-C20	2.84	120.48	111.83
3	A	813	PGW	O03-C19-C20	2.82	120.44	111.83
3	A	803	PGW	O03-C19-C20	2.74	120.19	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	806	PGW	O03-C19-C20	2.72	120.12	111.83
3	A	811	PGW	O03-C19-C20	2.67	119.97	111.83
3	A	814	PGW	O03-C19-C20	2.67	119.97	111.83
3	A	804	PGW	O03-C19-C20	2.65	119.92	111.83
3	A	817	PGW	O03-C19-C20	2.63	119.87	111.83
3	A	805	PGW	O03-C19-C20	2.63	119.86	111.83
3	A	809	PGW	O03-C19-C20	2.62	119.84	111.83
3	A	818	PGW	O03-C19-C20	2.62	119.82	111.83
3	A	812	PGW	O03-C19-C20	2.55	119.60	111.83
3	A	816	PGW	O13-P-O14	2.35	119.98	110.83
3	A	815	PGW	O13-P-O14	2.34	119.94	110.83
3	A	804	PGW	C02-O01-C1	-2.18	112.58	117.80

There are no chirality outliers.

All (237) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	803	PGW	C04-C05-CAD-OAE
3	A	803	PGW	O03-C01-C02-O01
3	A	804	PGW	C03-O11-P-O12
3	A	804	PGW	C03-O11-P-O14
3	A	805	PGW	C03-O11-P-O12
3	A	805	PGW	C03-O11-P-O13
3	A	805	PGW	C03-O11-P-O14
3	A	805	PGW	C2-C1-O01-C02
3	A	806	PGW	C01-C02-C03-O11
3	A	806	PGW	O01-C02-C03-O11
3	A	809	PGW	C2-C1-O01-C02
3	A	810	PGW	C03-O11-P-O13
3	A	810	PGW	C03-O11-P-O14
3	A	810	PGW	C2-C1-O01-C02
3	A	811	PGW	C03-O11-P-O12
3	A	811	PGW	C03-O11-P-O13
3	A	811	PGW	C03-O11-P-O14
3	A	811	PGW	O04-C19-O03-C01
3	A	811	PGW	C20-C19-O03-C01
3	A	813	PGW	OAF-C05-CAD-OAE
3	A	813	PGW	C04-C05-CAD-OAE
3	A	813	PGW	O12-C04-C05-CAD
3	A	815	PGW	C03-O11-P-O12
3	A	815	PGW	C03-O11-P-O13
3	A	815	PGW	C03-O11-P-O14

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Mol	Chain	Res	Type	Atoms
3	A	815	PGW	C2-C1-O01-C02
3	A	816	PGW	C2-C1-O01-C02
3	A	817	PGW	C03-O11-P-O12
3	A	817	PGW	C03-O11-P-O13
3	A	817	PGW	C03-O11-P-O14
3	A	818	PGW	C03-O11-P-O12
3	A	818	PGW	C03-O11-P-O13
3	A	818	PGW	C10-C06-C07-C08
3	A	812	PGW	O04-C19-O03-C01
3	A	809	PGW	O02-C1-O01-C02
3	A	810	PGW	O02-C1-O01-C02
3	A	815	PGW	O02-C1-O01-C02
3	A	816	PGW	O02-C1-O01-C02
3	A	809	PGW	C20-C19-O03-C01
3	A	812	PGW	C20-C19-O03-C01
3	A	814	PGW	O04-C19-O03-C01
3	A	805	PGW	O02-C1-O01-C02
3	A	814	PGW	C20-C19-O03-C01
3	A	809	PGW	O04-C19-O03-C01
3	A	803	PGW	C2-C1-O01-C02
3	A	816	PGW	C3-C4-C5-C6
3	A	813	PGW	O12-C04-C05-OAF
3	A	813	PGW	C19-C20-C21-C22
3	A	803	PGW	C17-C18-C28-C30
3	A	818	PGW	C25-C26-C27-C15
3	A	804	PGW	C1-C2-C3-C4
3	A	816	PGW	C1-C2-C3-C4
3	A	816	PGW	C19-C20-C21-C22
3	A	806	PGW	C19-C20-C21-C22
3	A	808	PGW	C1-C2-C3-C4
3	A	815	PGW	C1-C2-C3-C4
3	A	817	PGW	C19-C20-C21-C22
3	A	818	PGW	C1-C2-C3-C4
3	A	806	PGW	C16-C15-C27-C26
3	A	803	PGW	O02-C1-O01-C02
3	A	813	PGW	C10-C06-C07-C08
3	A	811	PGW	C1-C2-C3-C4
3	A	809	PGW	C19-C20-C21-C22
3	A	811	PGW	C24-C25-C26-C27
3	A	806	PGW	C2-C1-O01-C02
3	A	806	PGW	O02-C1-O01-C02
3	A	814	PGW	C2-C1-O01-C02

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Mol	Chain	Res	Type	Atoms
3	A	804	PGW	C21-C22-C23-C24
3	A	805	PGW	C23-C24-C25-C26
3	A	810	PGW	C2-C3-C4-C5
3	A	818	PGW	C2-C3-C4-C5
3	A	805	PGW	C2-C3-C4-C5
3	A	813	PGW	C24-C25-C26-C27
3	A	817	PGW	C2-C1-O01-C02
3	A	805	PGW	C16-C15-C27-C26
3	A	813	PGW	C16-C15-C27-C26
3	A	816	PGW	C20-C21-C22-C23
3	A	811	PGW	C22-C23-C24-C25
3	A	813	PGW	C22-C23-C24-C25
3	A	810	PGW	C20-C21-C22-C23
3	A	813	PGW	C5-C6-C7-C8
3	A	813	PGW	C25-C26-C27-C15
3	A	811	PGW	C23-C24-C25-C26
3	A	816	PGW	C4-C5-C6-C7
3	A	814	PGW	O02-C1-O01-C02
3	A	817	PGW	O02-C1-O01-C02
3	A	803	PGW	C24-C25-C26-C27
3	A	810	PGW	C1-C2-C3-C4
3	A	818	PGW	C24-C25-C26-C27
3	A	809	PGW	C21-C22-C23-C24
3	A	803	PGW	C23-C24-C25-C26
3	A	805	PGW	C21-C22-C23-C24
3	A	806	PGW	C22-C23-C24-C25
3	A	808	PGW	C2-C1-O01-C02
3	A	812	PGW	C2-C1-O01-C02
3	A	815	PGW	C2-C3-C4-C5
3	A	817	PGW	C21-C22-C23-C24
3	A	816	PGW	C21-C22-C23-C24
3	A	810	PGW	C19-C20-C21-C22
3	A	812	PGW	O02-C1-O01-C02
3	A	812	PGW	C1-C2-C3-C4
3	A	811	PGW	O03-C01-C02-O01
3	A	808	PGW	C3-C4-C5-C6
3	A	812	PGW	C7-C8-C9-C10
3	A	807	PGW	C02-C03-O11-P
3	A	811	PGW	C21-C22-C23-C24
3	A	803	PGW	C19-C20-C21-C22
3	A	803	PGW	OAF-C05-CAD-OAE
3	A	815	PGW	C4-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
3	A	803	PGW	C6-C7-C8-C9
3	A	806	PGW	C6-C7-C8-C9
3	A	815	PGW	C5-C6-C7-C8
3	A	804	PGW	C01-C02-C03-O11
3	A	805	PGW	C01-C02-C03-O11
3	A	807	PGW	C5-C6-C7-C8
3	A	818	PGW	C19-C20-C21-C22
3	A	809	PGW	C20-C21-C22-C23
3	A	808	PGW	O02-C1-O01-C02
3	A	809	PGW	O03-C01-C02-C03
3	A	806	PGW	C10-C06-C07-C08
3	A	818	PGW	C23-C24-C25-C26
3	A	813	PGW	C23-C24-C25-C26
3	A	808	PGW	C4-C5-C6-C7
3	A	818	PGW	C03-O11-P-O14
3	A	803	PGW	C16-C15-C27-C26
3	A	818	PGW	C22-C23-C24-C25
3	A	817	PGW	C20-C21-C22-C23
3	A	818	PGW	C5-C6-C7-C8
3	A	804	PGW	C3-C4-C5-C6
3	A	804	PGW	C20-C21-C22-C23
3	A	804	PGW	C22-C23-C24-C25
3	A	814	PGW	C2-C3-C4-C5
3	A	803	PGW	C18-C28-C30-C29
3	A	806	PGW	C21-C22-C23-C24
3	A	809	PGW	C15-C16-C17-C18
3	A	809	PGW	C23-C24-C25-C26
3	A	809	PGW	C01-C02-C03-O11
3	A	814	PGW	C01-C02-C03-O11
3	A	810	PGW	C24-C25-C26-C27
3	A	803	PGW	O03-C01-C02-C03
3	A	810	PGW	O03-C01-C02-C03
3	A	811	PGW	O03-C01-C02-C03
3	A	803	PGW	C22-C23-C24-C25
3	A	805	PGW	O01-C02-C03-O11
3	A	809	PGW	O01-C02-C03-O11
3	A	810	PGW	O01-C02-C03-O11
3	A	815	PGW	O01-C02-C03-O11
3	A	816	PGW	O01-C02-C03-O11
3	A	811	PGW	C3-C4-C5-C6
3	A	810	PGW	C03-O11-P-O12
3	A	815	PGW	C19-C20-C21-C22

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Mol	Chain	Res	Type	Atoms
3	A	807	PGW	C19-C20-C21-C22
3	A	809	PGW	C03-C02-O01-C1
3	A	815	PGW	C03-C02-O01-C1
3	A	806	PGW	C15-C16-C17-C18
3	A	809	PGW	C16-C17-C18-C28
3	A	805	PGW	C6-C7-C8-C9
3	A	814	PGW	O01-C02-C03-O11
3	A	808	PGW	O03-C01-C02-C03
3	A	816	PGW	C20-C19-O03-C01
3	A	808	PGW	O03-C01-C02-O01
3	A	810	PGW	O03-C01-C02-O01
3	A	813	PGW	C21-C22-C23-C24
3	A	808	PGW	C2-C3-C4-C5
3	A	817	PGW	C20-C19-O03-C01
3	A	812	PGW	C19-C20-C21-C22
3	A	815	PGW	C22-C23-C24-C25
3	A	812	PGW	C6-C7-C8-C9
3	A	811	PGW	C01-C02-C03-O11
3	A	815	PGW	C01-C02-C03-O11
3	A	816	PGW	C01-C02-C03-O11
3	A	810	PGW	C22-C23-C24-C25
3	A	807	PGW	C06-C07-C08-C09
3	A	816	PGW	O04-C19-O03-C01
3	A	817	PGW	O04-C19-O03-C01
3	A	812	PGW	C2-C3-C4-C5
3	A	811	PGW	O01-C02-C03-O11
3	A	807	PGW	C10-C06-C07-C08
3	A	803	PGW	C20-C21-C22-C23
3	A	809	PGW	O03-C01-C02-O01
3	A	812	PGW	O03-C01-C02-O01
3	A	812	PGW	O03-C01-C02-C03
3	A	813	PGW	O03-C01-C02-C03
3	A	818	PGW	C20-C19-O03-C01
3	A	813	PGW	C1-C2-C3-C4
3	A	805	PGW	C22-C23-C24-C25
3	A	808	PGW	C04-O12-P-O14
3	A	813	PGW	C03-O11-P-O14
3	A	818	PGW	O04-C19-O03-C01
3	A	803	PGW	C15-C16-C17-C18
3	A	810	PGW	C01-C02-C03-O11
3	A	805	PGW	O03-C19-C20-C21
3	A	818	PGW	O03-C19-C20-C21

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Mol	Chain	Res	Type	Atoms
3	A	804	PGW	O01-C02-C03-O11
3	A	807	PGW	O01-C02-C03-O11
3	A	812	PGW	C20-C21-C22-C23
3	A	809	PGW	C02-C03-O11-P
3	A	813	PGW	O03-C01-C02-O01
3	A	805	PGW	C3-C4-C5-C6
3	A	814	PGW	O01-C1-C2-C3
3	A	818	PGW	C4-C5-C6-C7
3	A	804	PGW	C19-C20-C21-C22
3	A	803	PGW	C21-C22-C23-C24
3	A	806	PGW	C17-C18-C28-C30
3	A	807	PGW	C4-C5-C6-C7
3	A	803	PGW	C2-C3-C4-C5
3	A	815	PGW	C3-C4-C5-C6
3	A	803	PGW	O03-C19-C20-C21
3	A	808	PGW	C02-C01-O03-C19
3	A	807	PGW	C2-C3-C4-C5
3	A	815	PGW	C21-C22-C23-C24
3	A	813	PGW	C15-C16-C17-C18
3	A	814	PGW	C20-C21-C22-C23
3	A	813	PGW	O01-C02-C03-O11
3	A	815	PGW	O03-C19-C20-C21
3	A	817	PGW	O03-C01-C02-O01
3	A	812	PGW	C02-C03-O11-P
3	A	806	PGW	C4-C5-C6-C7
3	A	804	PGW	C03-O11-P-O13
3	A	806	PGW	C7-C8-C9-C10
3	A	803	PGW	C5-C6-C7-C8
3	A	809	PGW	C22-C23-C24-C25
3	A	813	PGW	C7-C8-C9-C10
3	A	807	PGW	C01-C02-C03-O11
3	A	818	PGW	O02-C1-O01-C02
3	A	812	PGW	O01-C1-C2-C3
3	A	805	PGW	C4-C5-C6-C7
3	A	814	PGW	O03-C01-C02-O01
3	A	812	PGW	C3-C4-C5-C6
3	A	811	PGW	O01-C1-C2-C3
3	A	812	PGW	O02-C1-C2-C3
3	A	812	PGW	O03-C19-C20-C21
3	A	813	PGW	C01-C02-C03-O11
3	A	807	PGW	C07-C06-C10-C9
3	A	803	PGW	C1-C2-C3-C4

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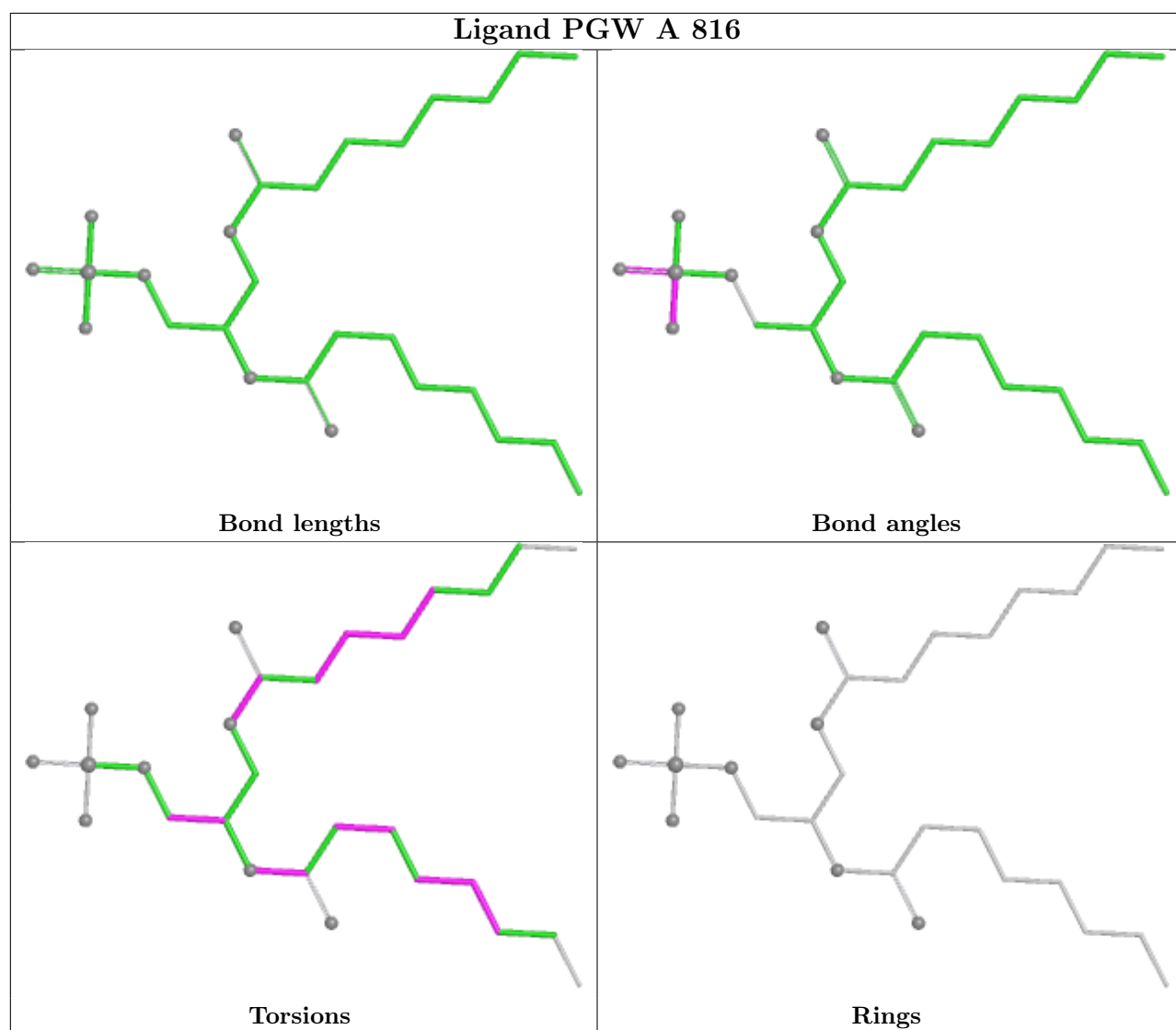
Mol	Chain	Res	Type	Atoms
3	A	818	PGW	C2-C1-O01-C02
3	A	809	PGW	O01-C1-C2-C3

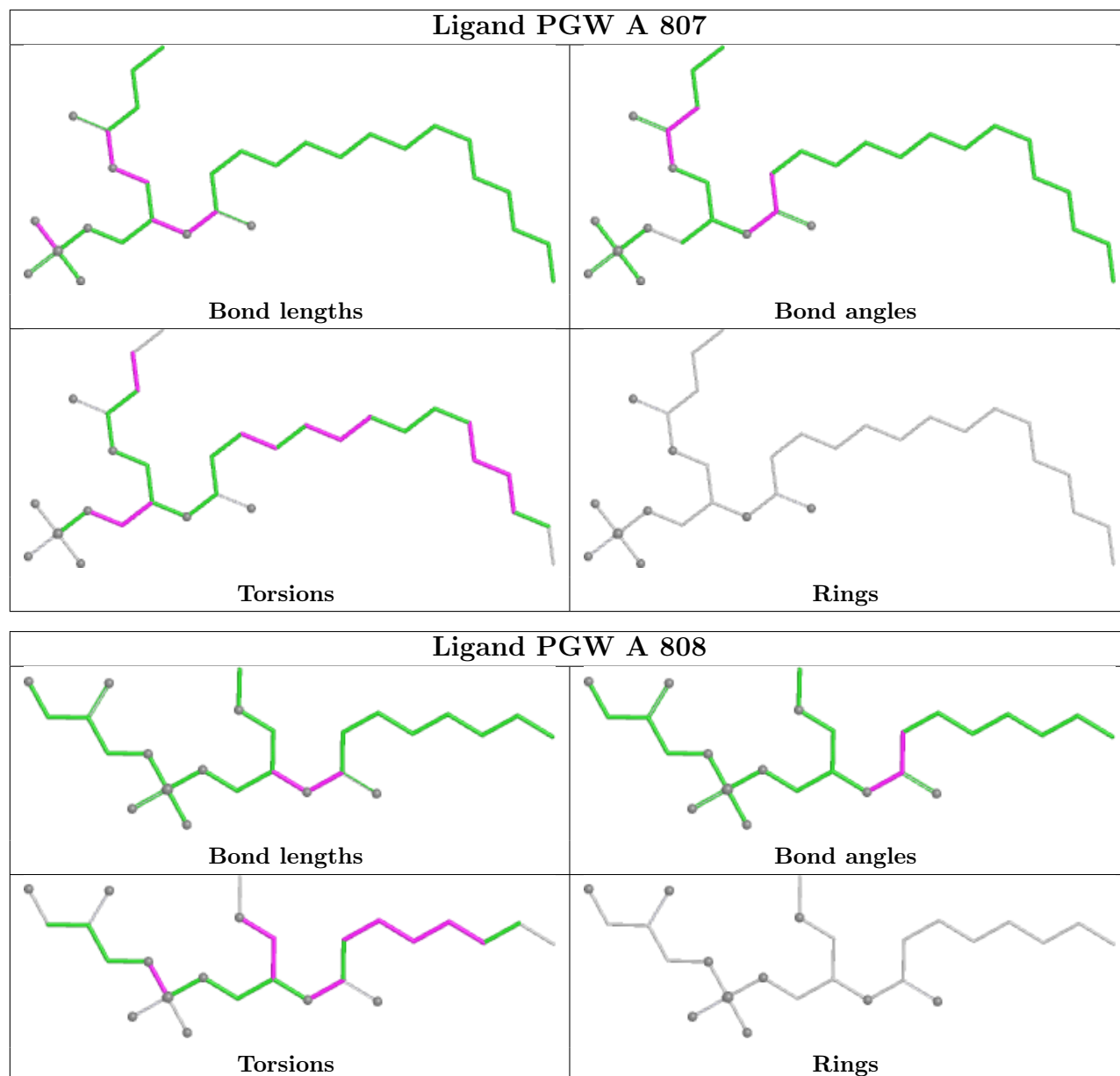
There are no ring outliers.

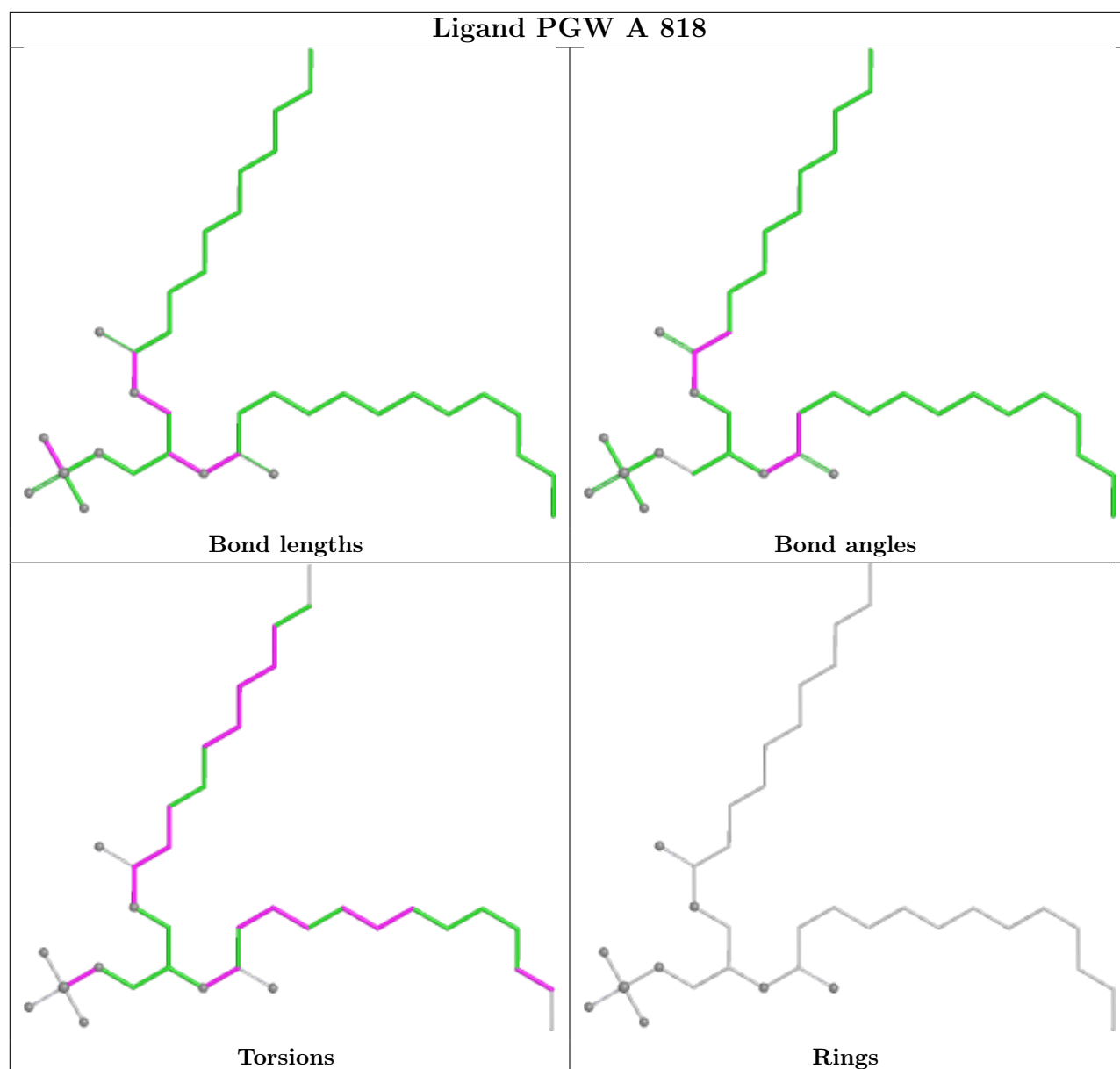
14 monomers are involved in 52 short contacts:

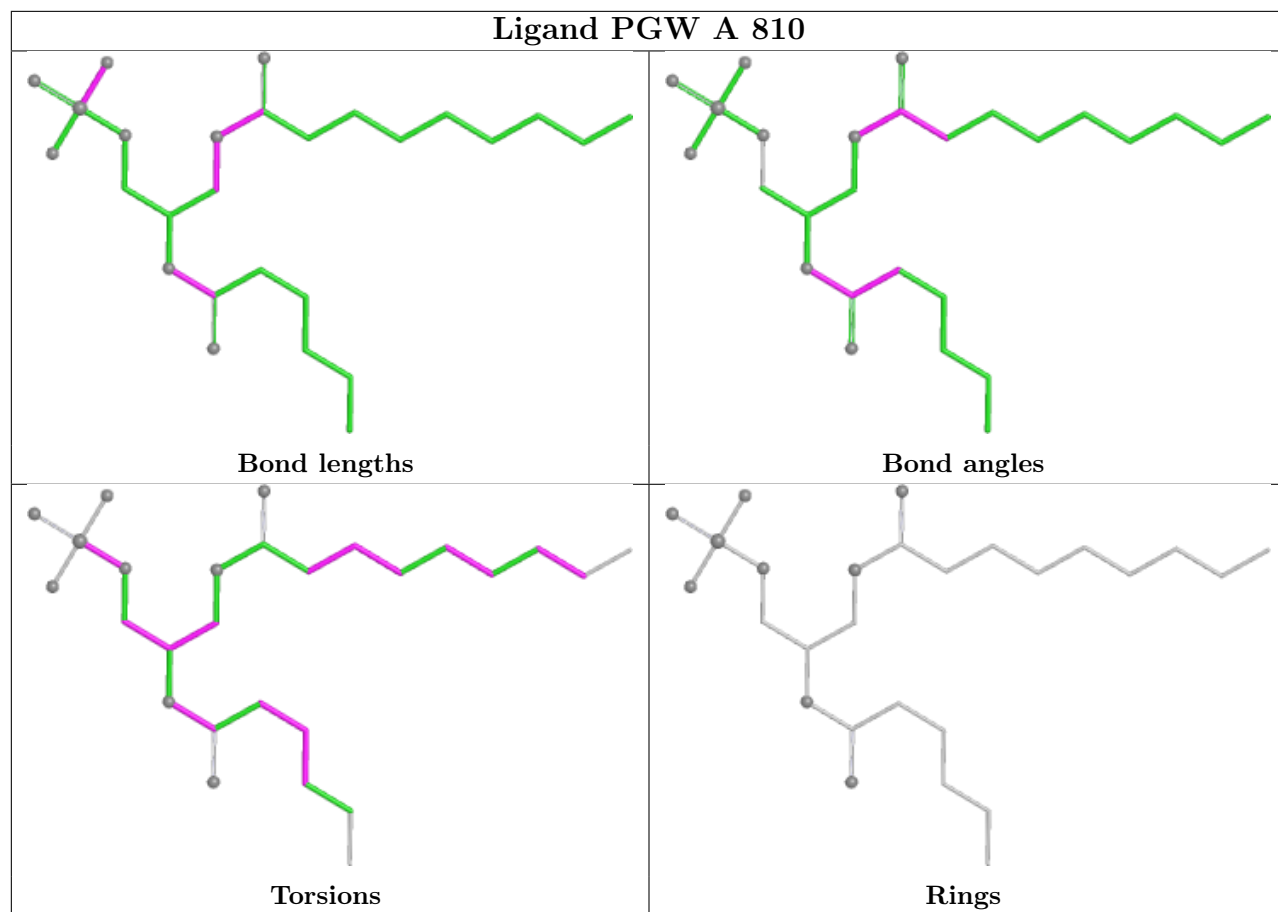
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	816	PGW	2	0
3	A	807	PGW	1	0
3	A	808	PGW	4	0
3	A	818	PGW	10	0
3	A	810	PGW	2	0
3	A	809	PGW	3	0
3	A	803	PGW	6	0
3	A	806	PGW	6	0
3	A	811	PGW	1	0
3	A	814	PGW	10	0
3	A	812	PGW	6	0
3	A	813	PGW	3	0
3	A	805	PGW	6	0
3	A	815	PGW	1	0

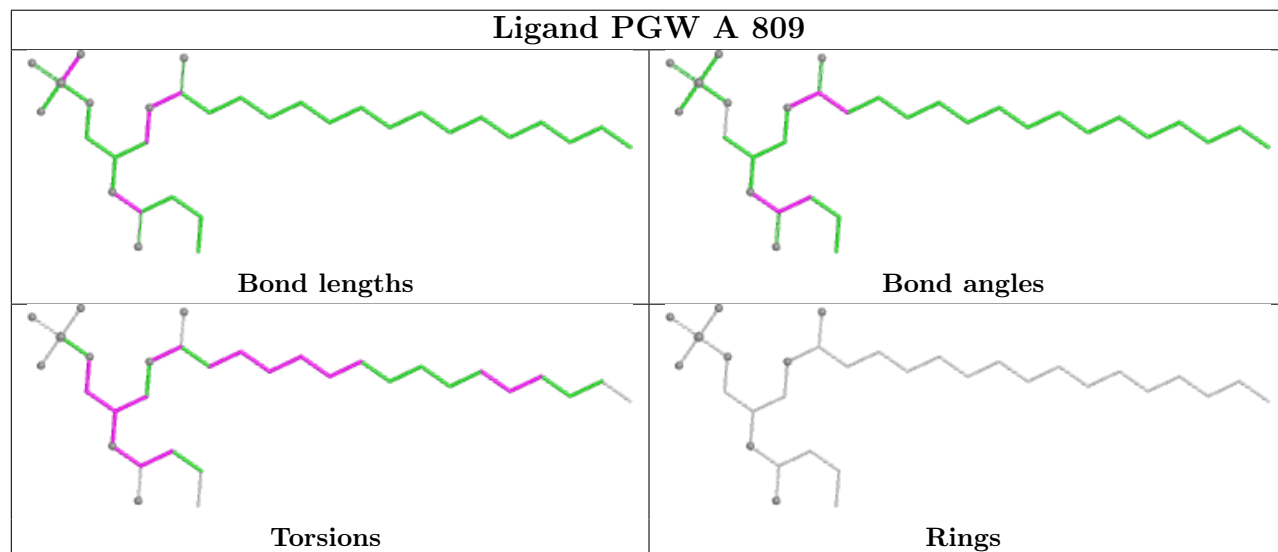
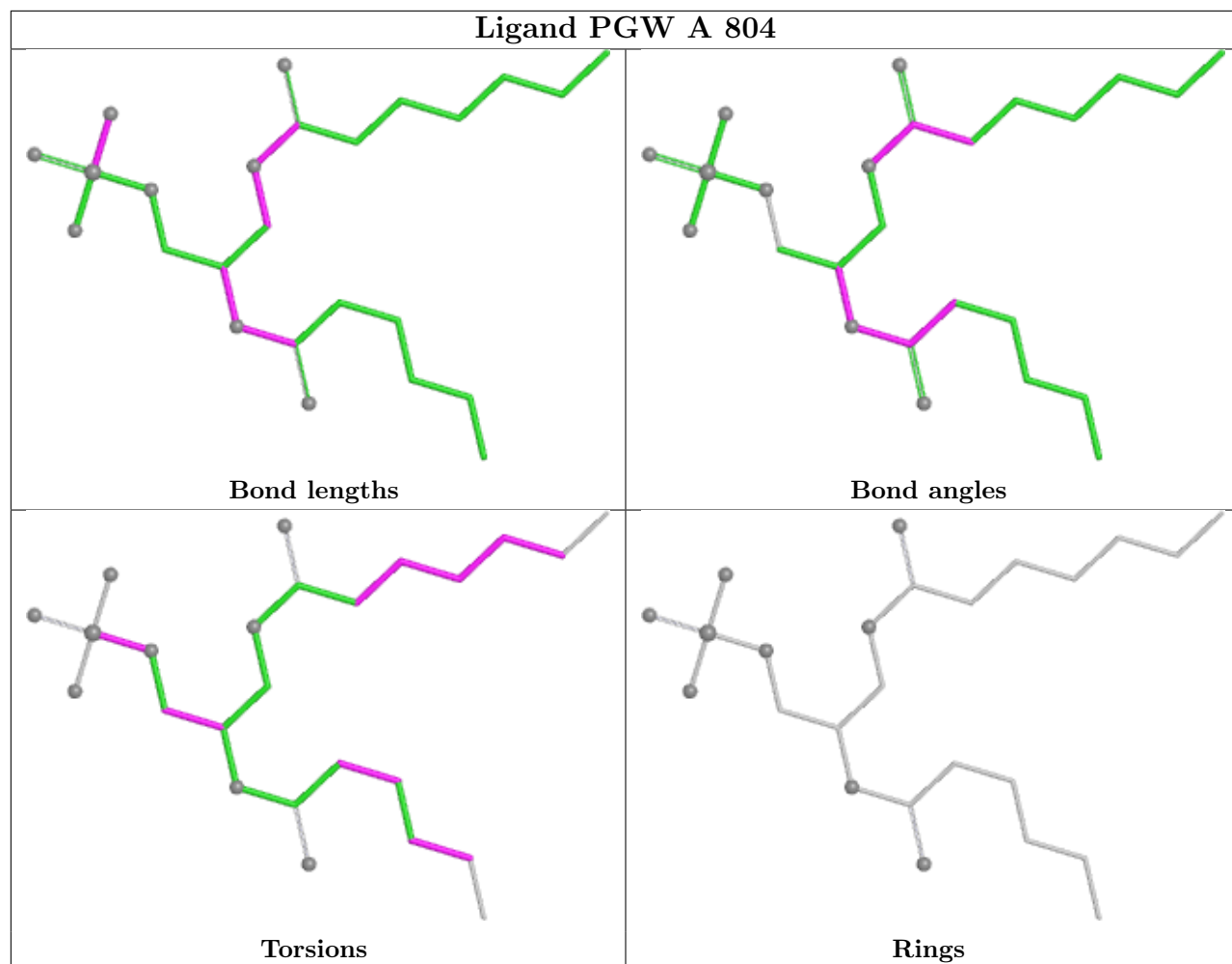
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

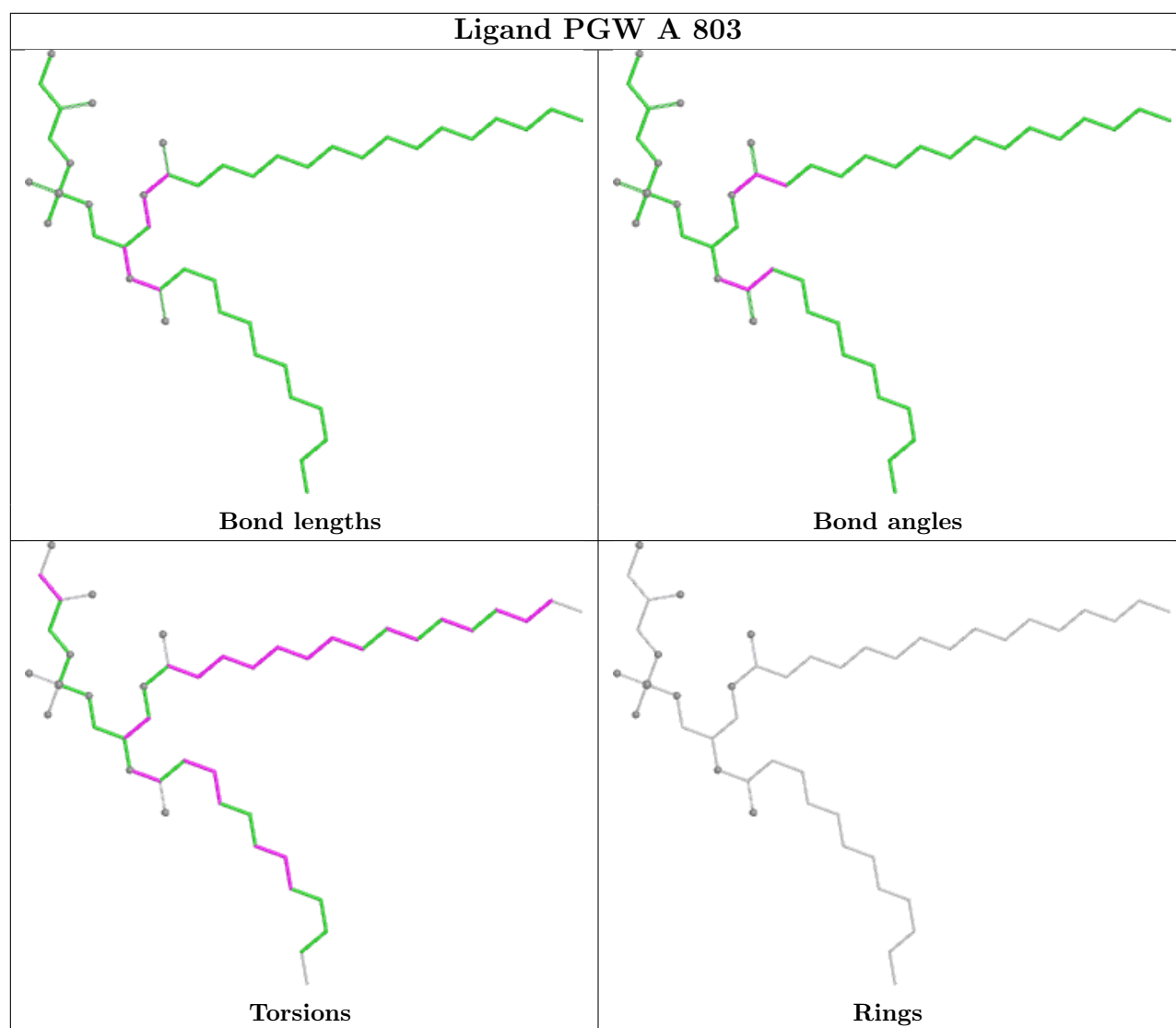


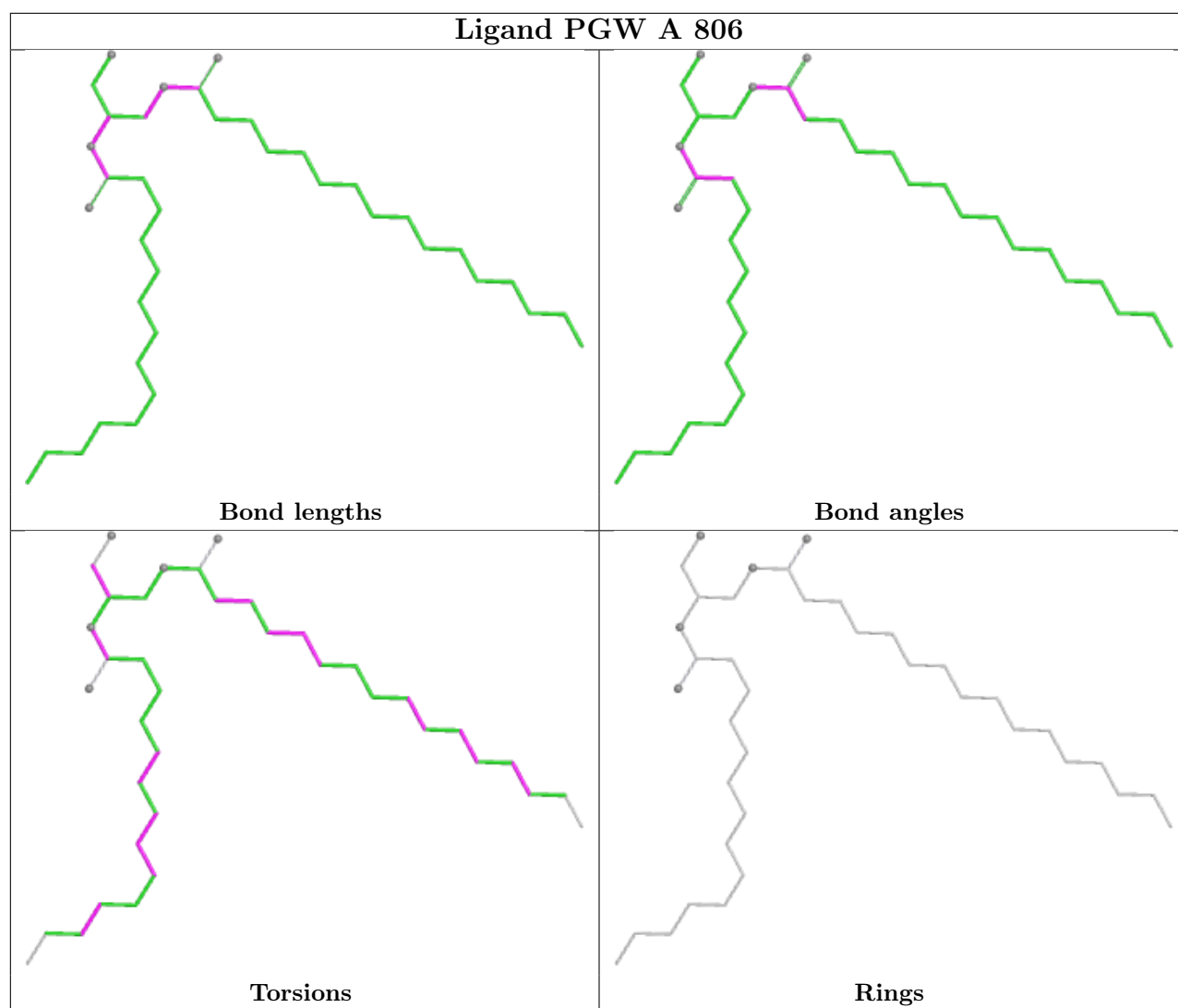


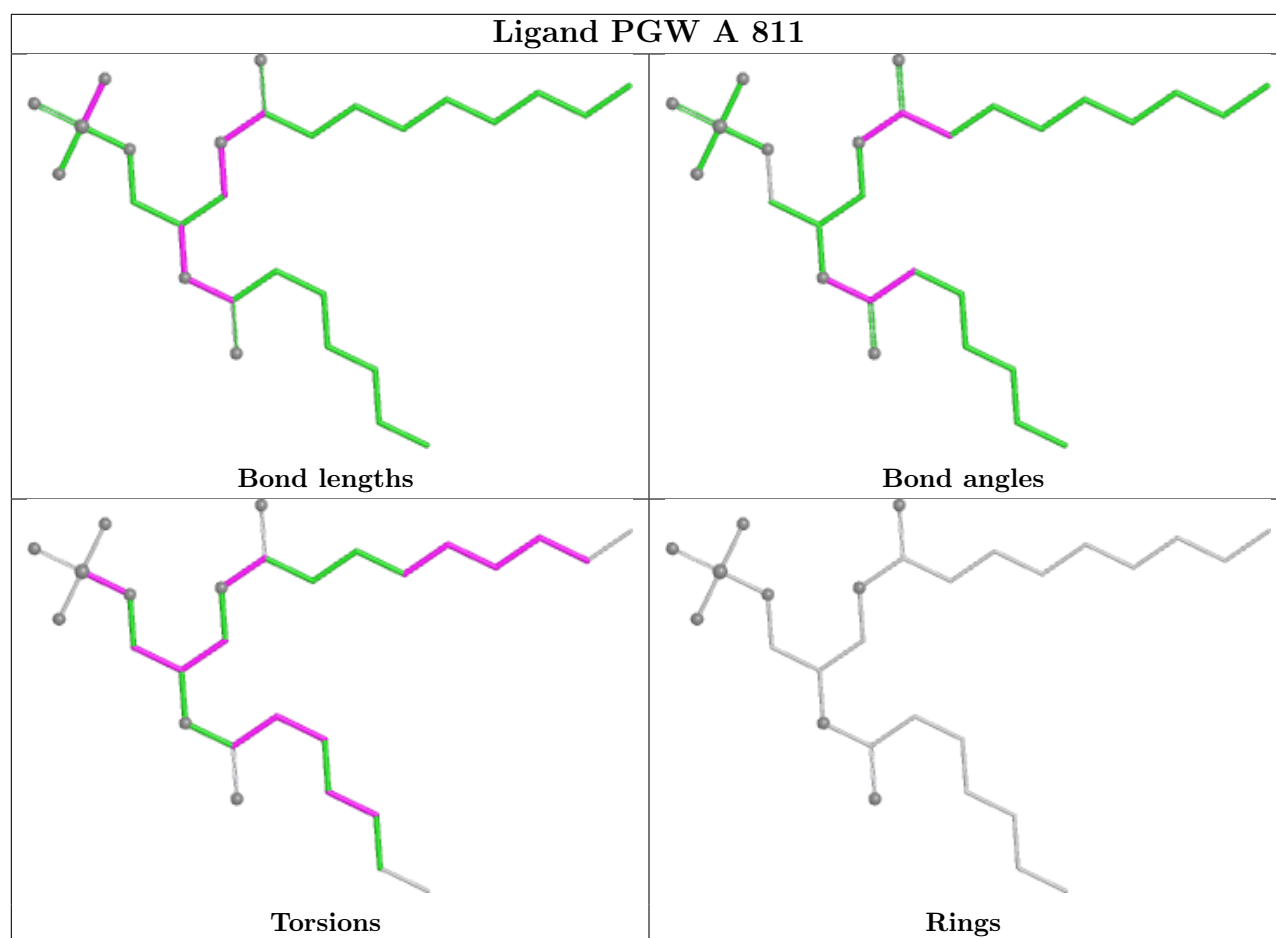


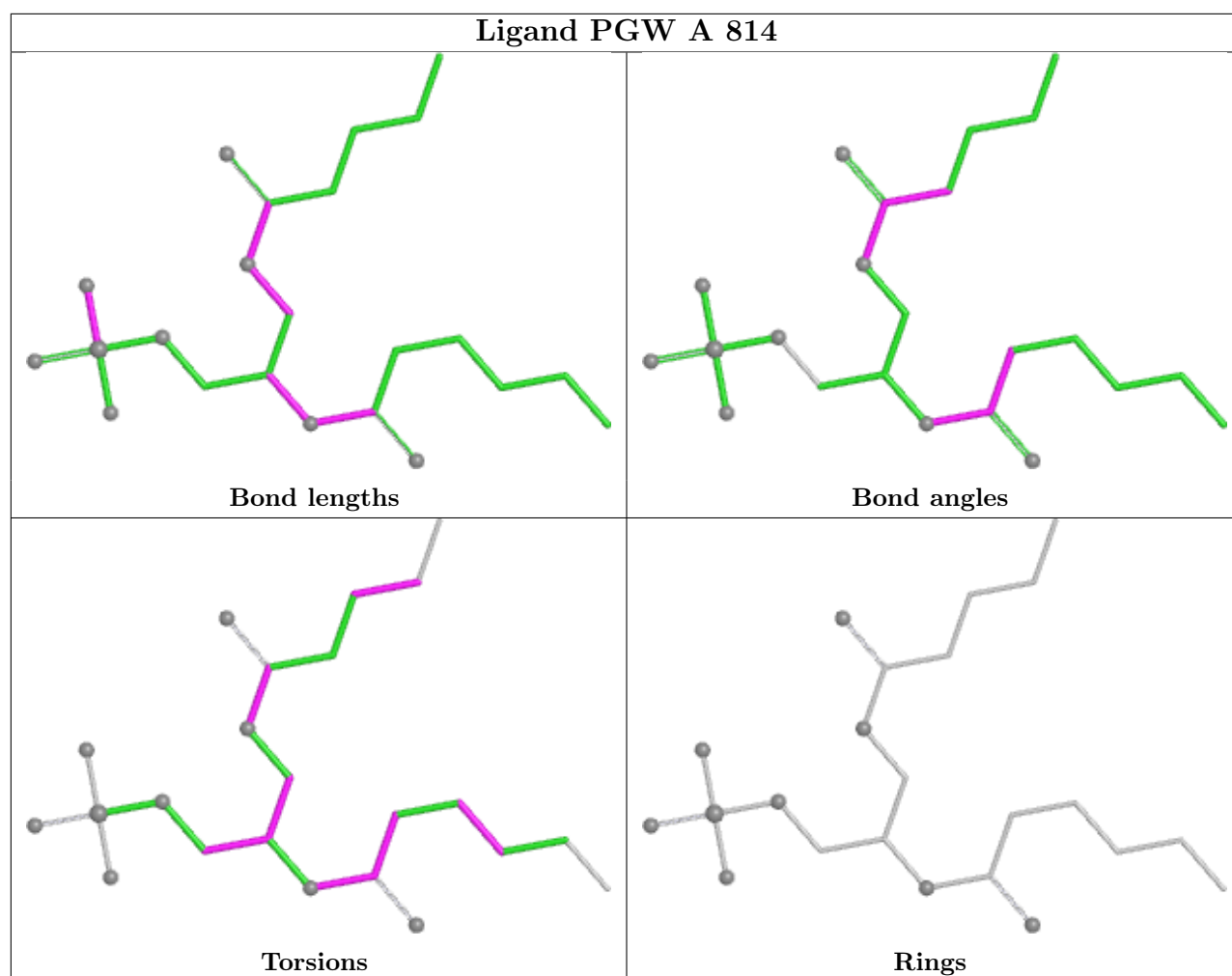


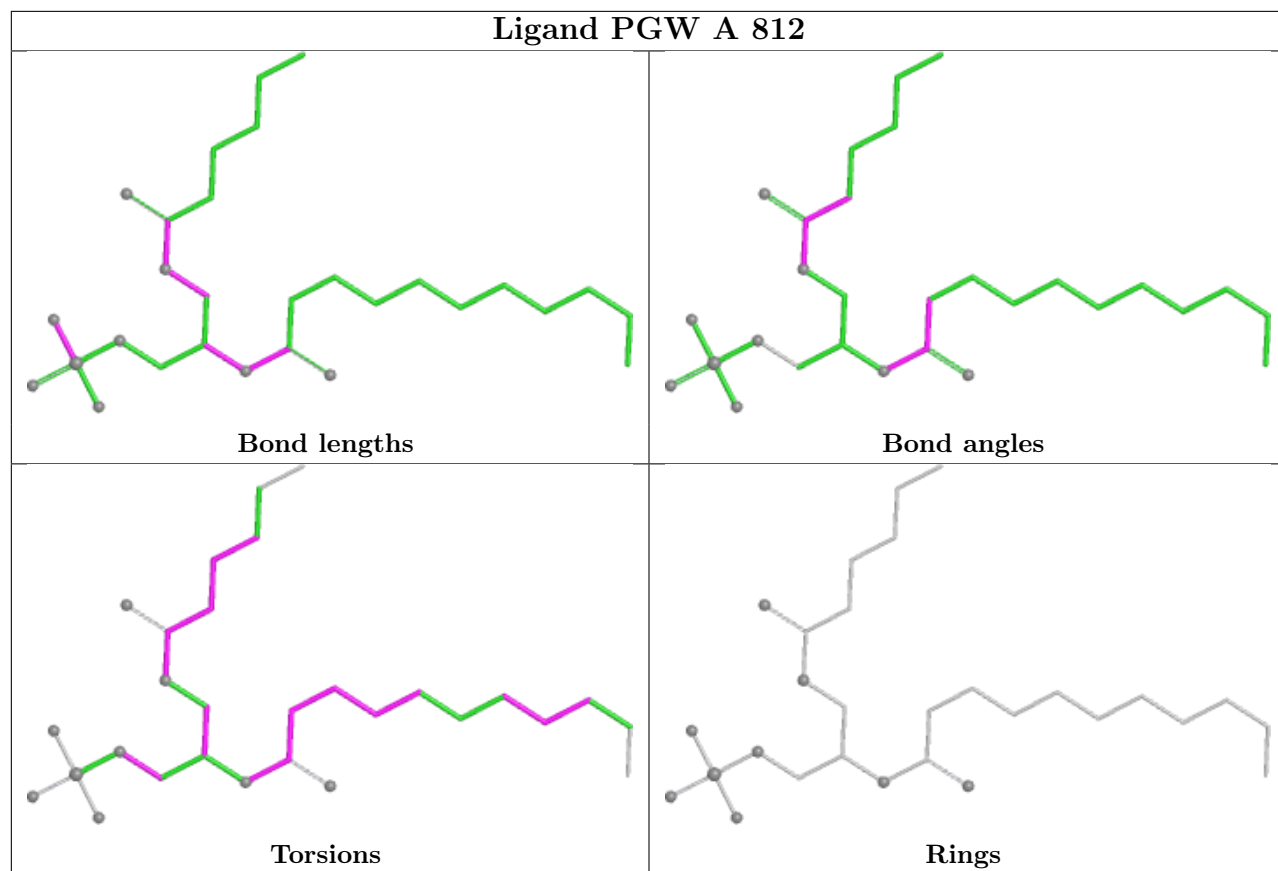


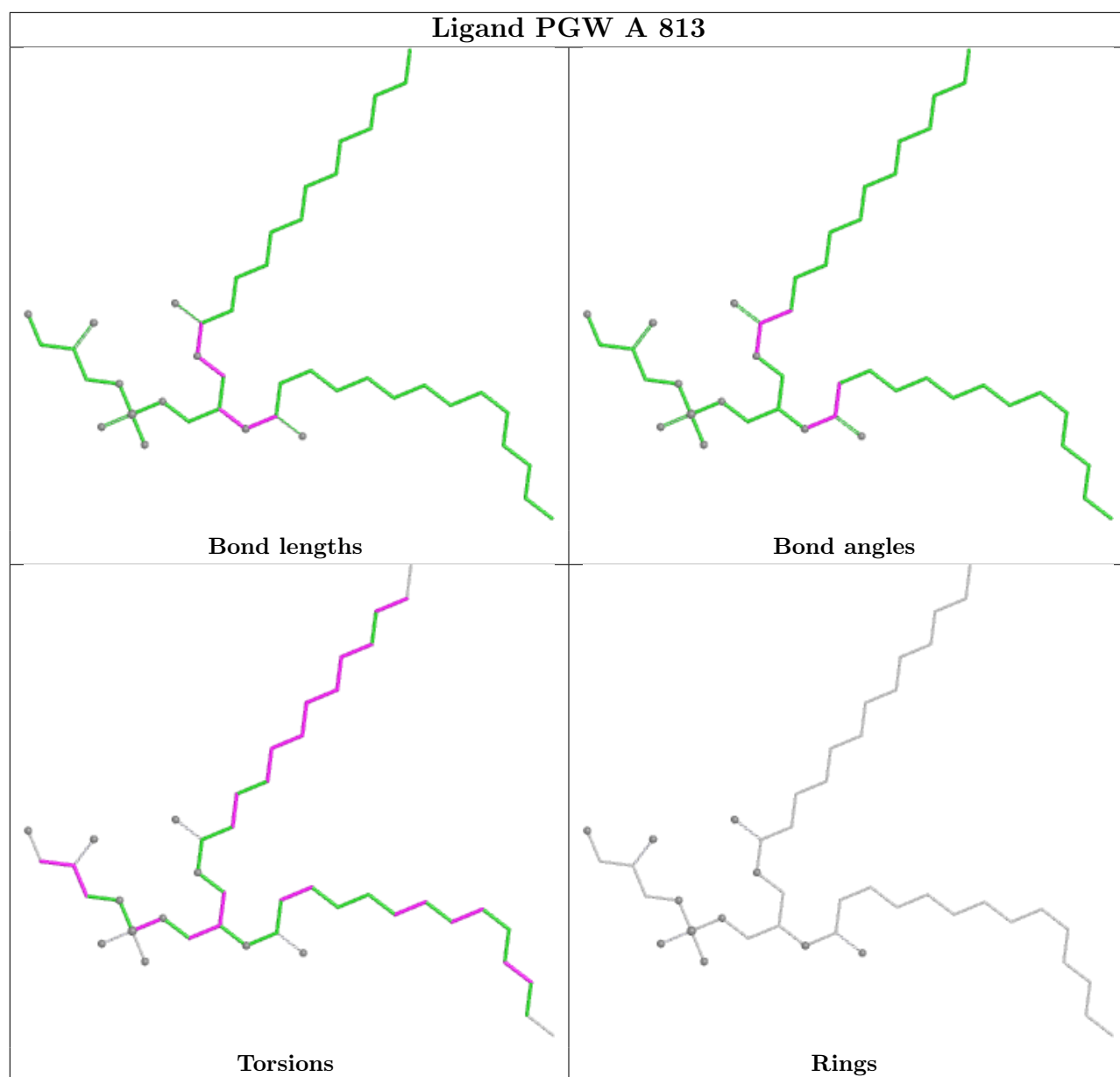


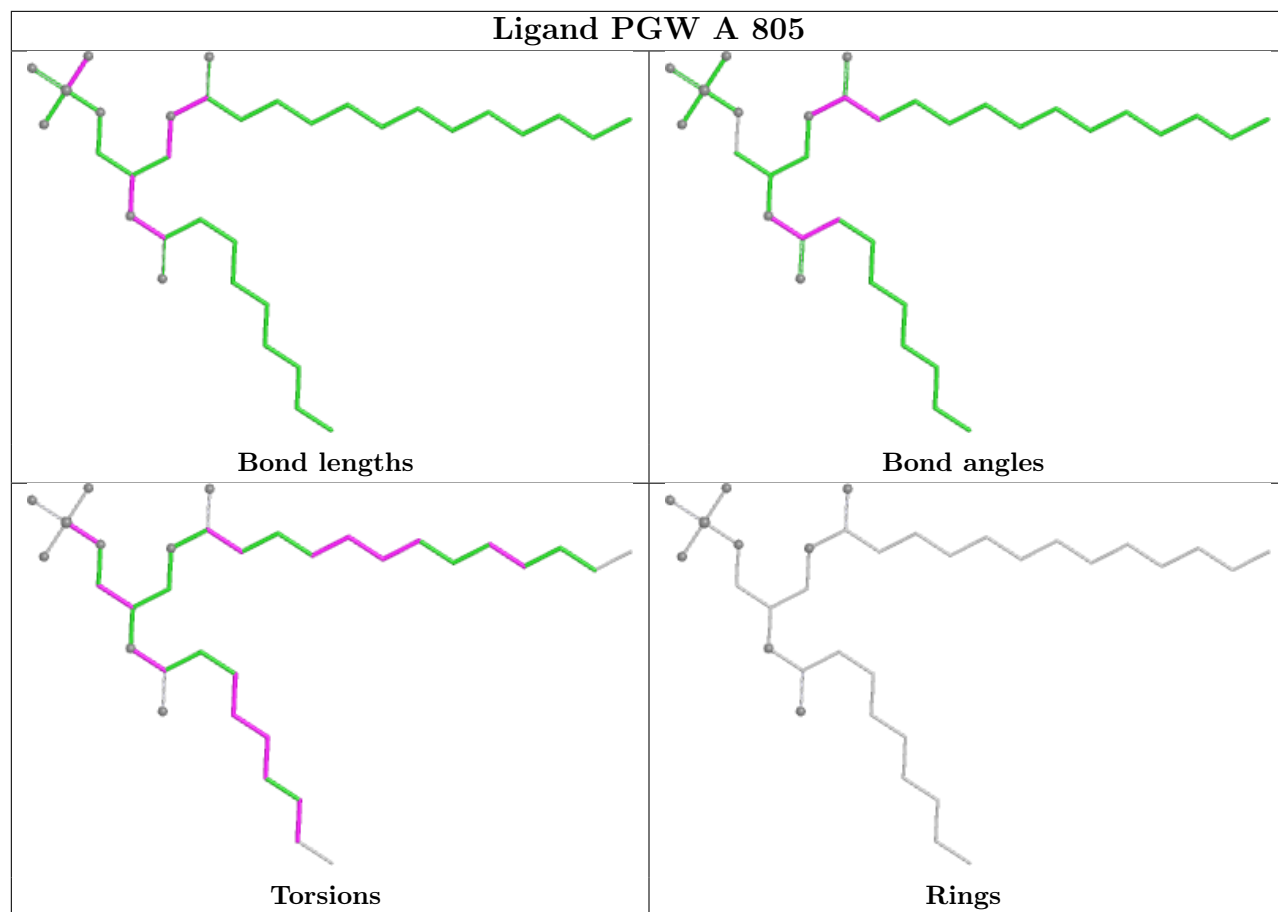


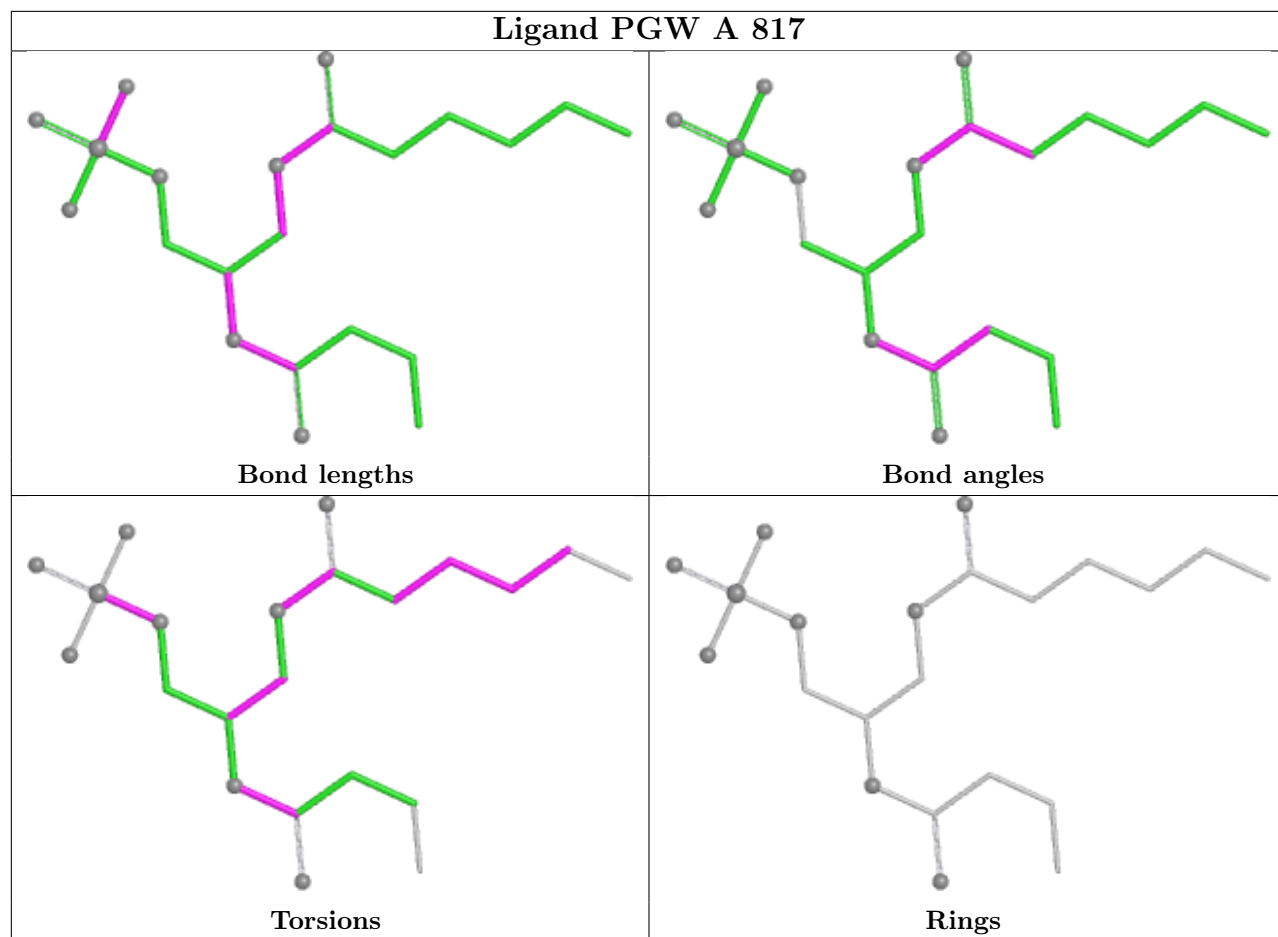


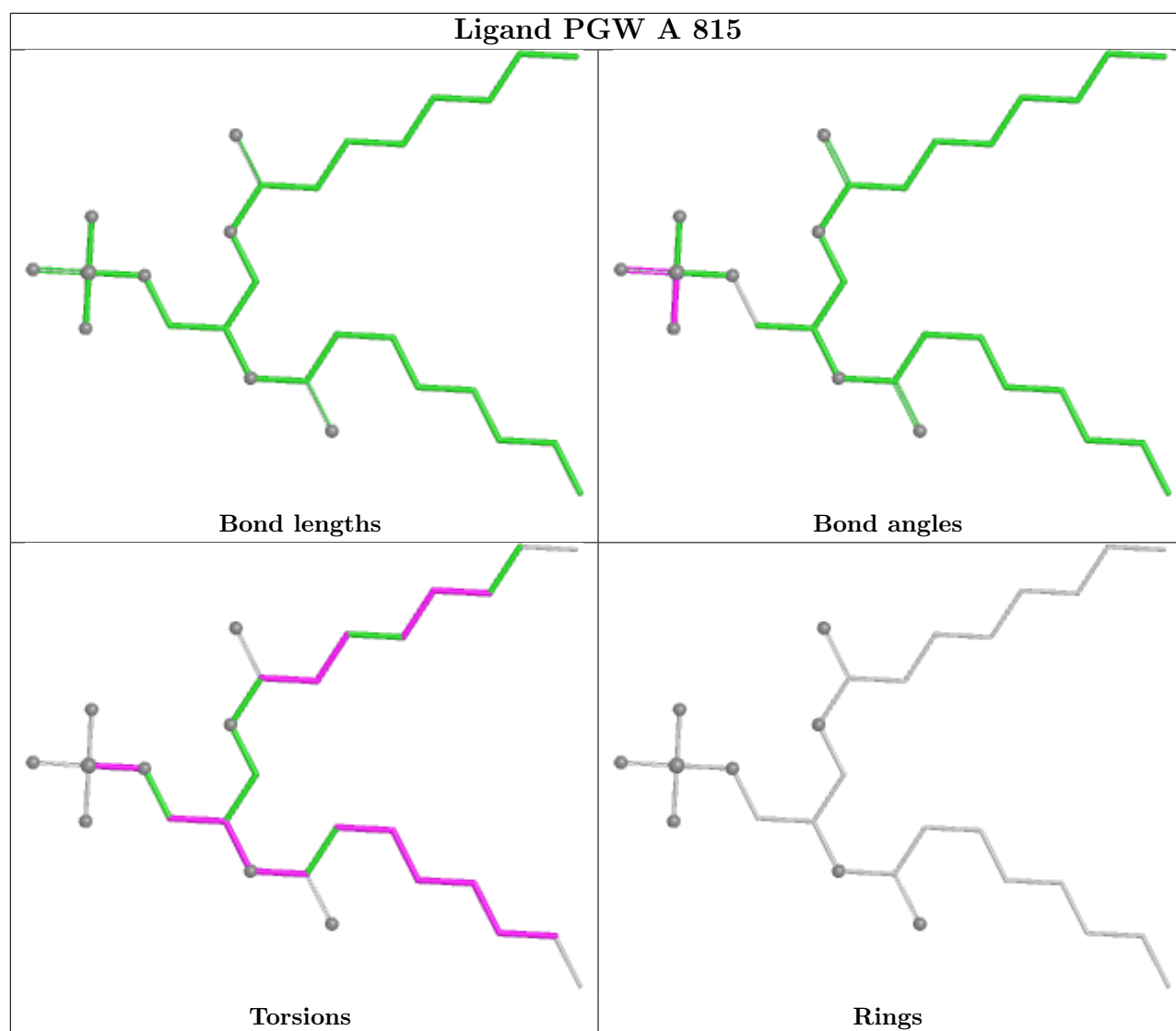












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

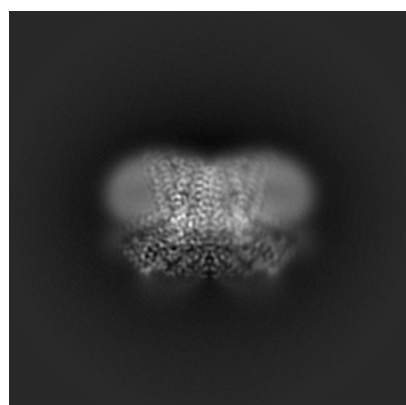
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24731. These allow visual inspection of the internal detail of the map and identification of artifacts.

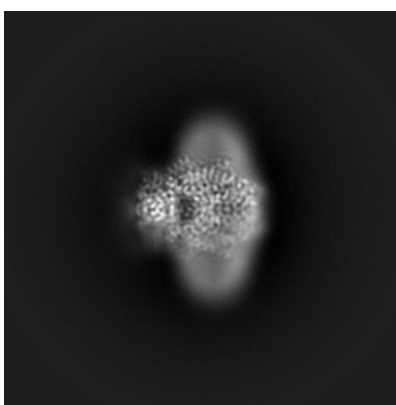
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

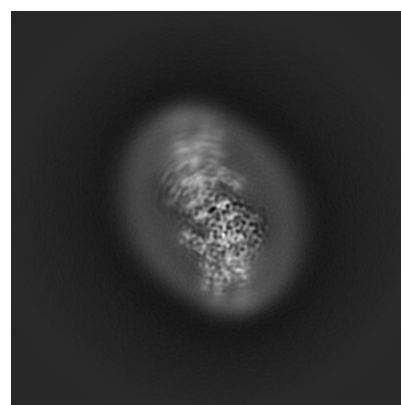
6.1.1 Primary map



X



Y

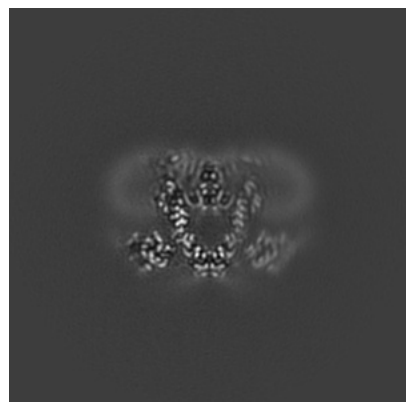


Z

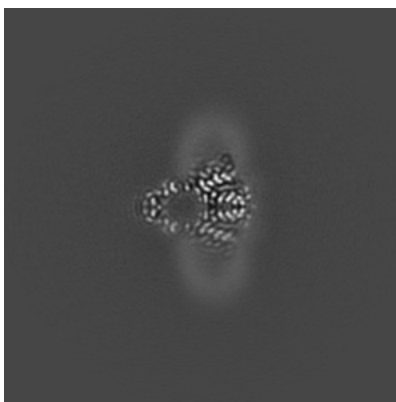
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

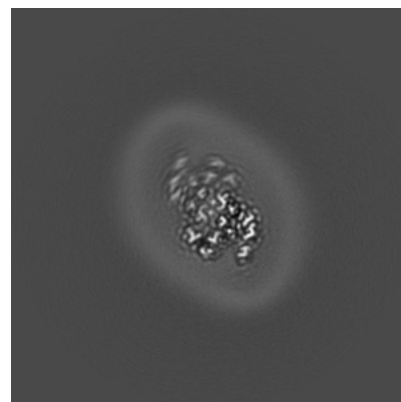
6.2.1 Primary map



X Index: 192



Y Index: 192

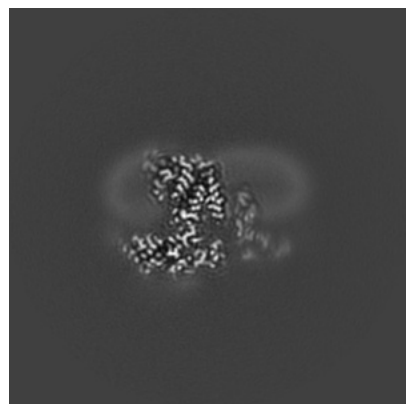


Z Index: 192

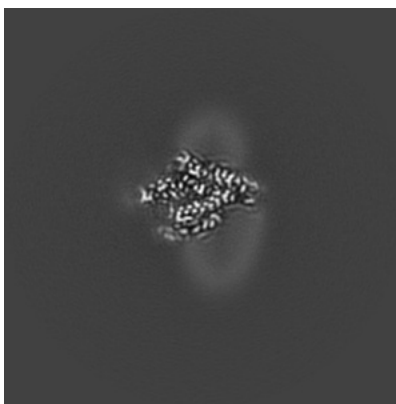
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

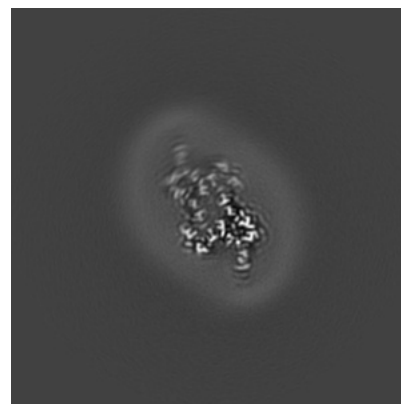
6.3.1 Primary map



X Index: 203



Y Index: 163

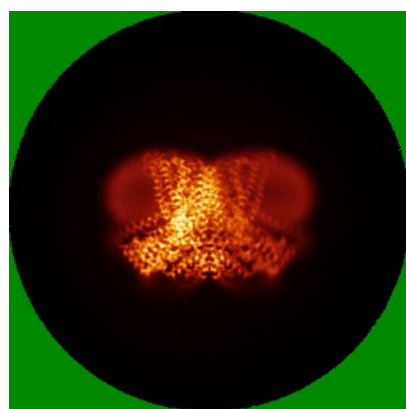


Z Index: 185

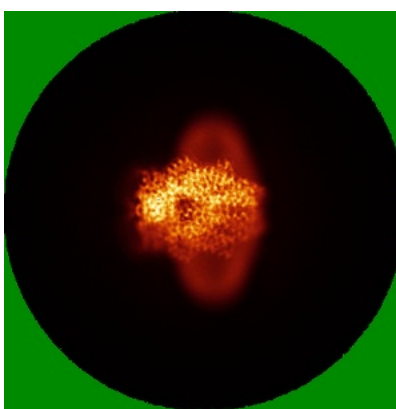
The images above show the largest variance slices of the map in three orthogonal directions.

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

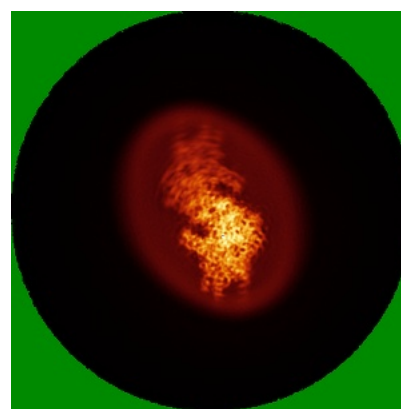
6.4.1 Primary map



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 0.0049. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

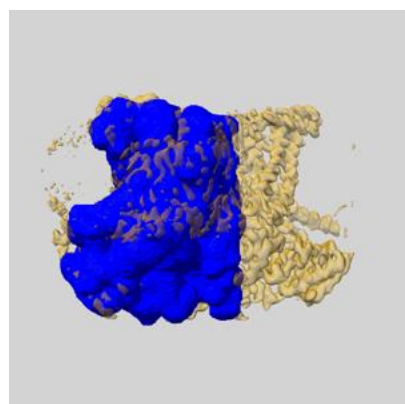
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

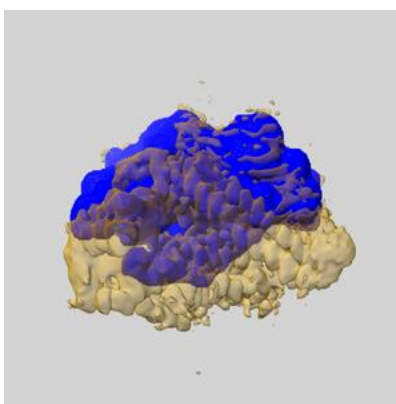
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

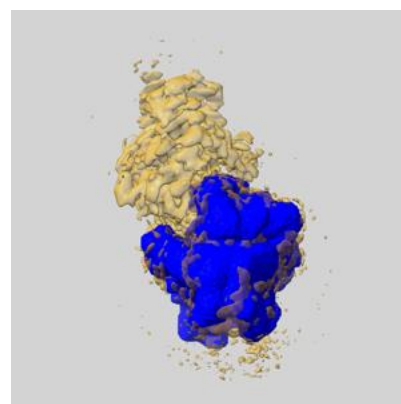
6.6.1 emd_24731_msk_1.map [i](#)



X



Y

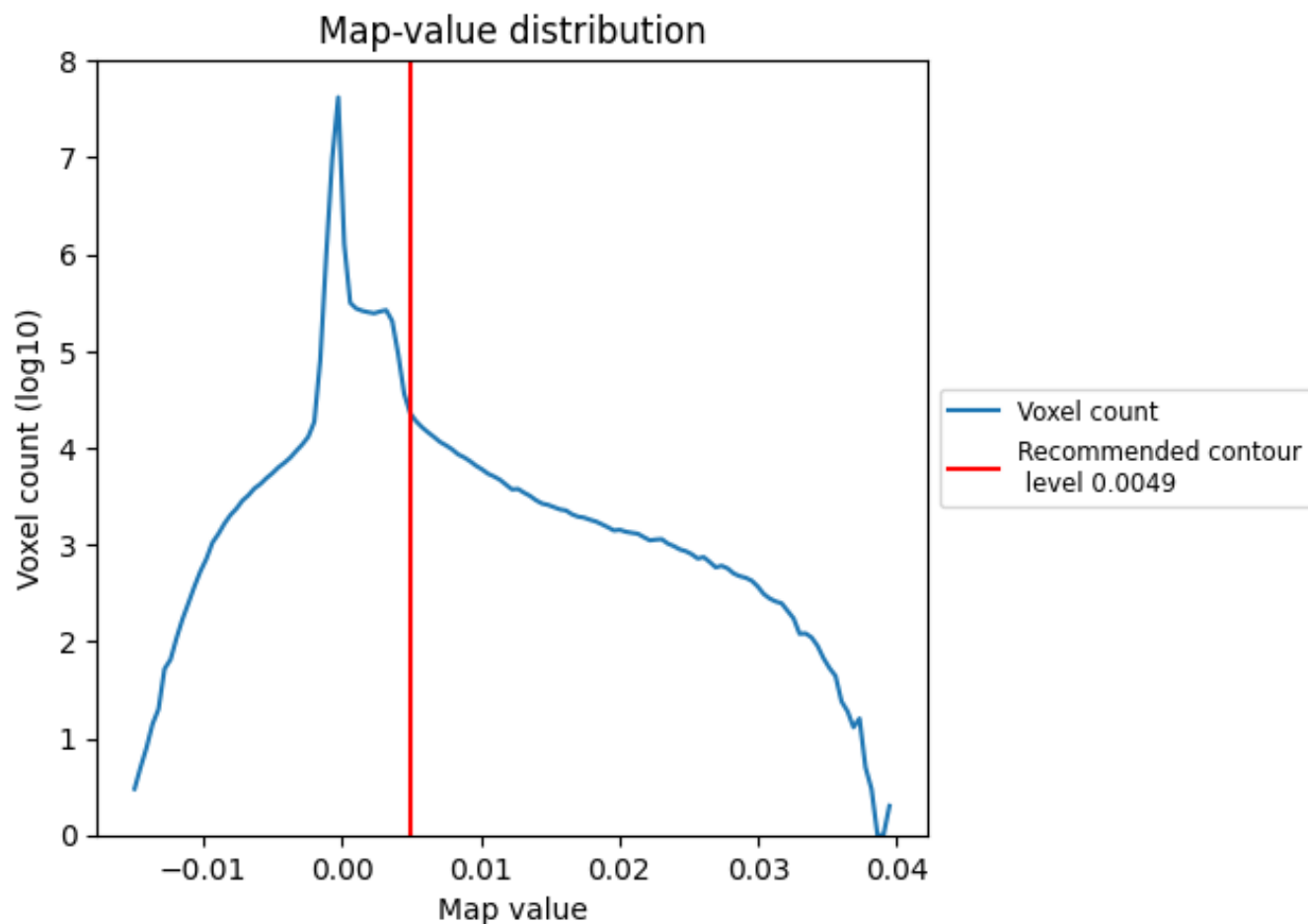


Z

7 Map analysis [i](#)

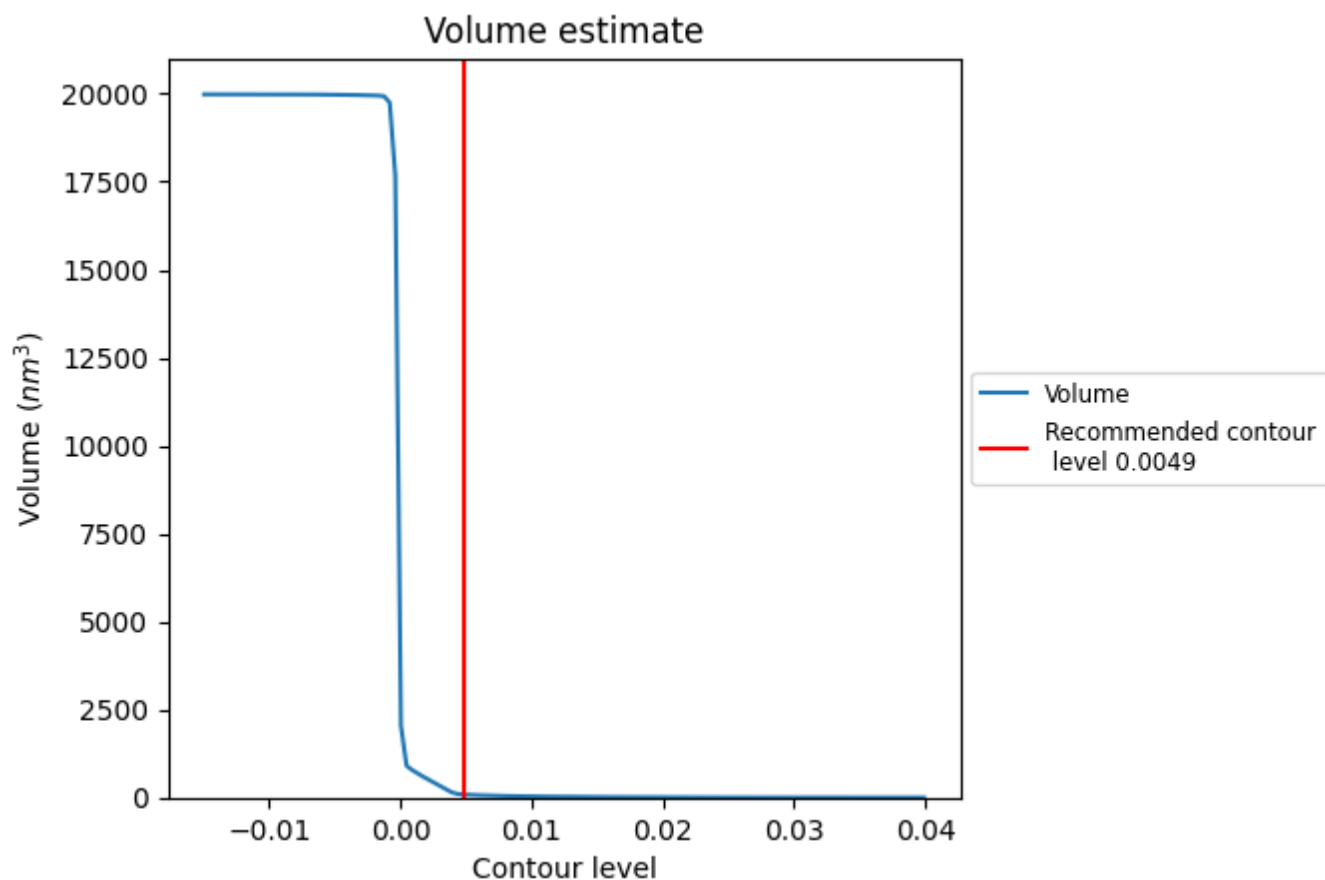
This section contains the results of statistical analysis of the map.

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.

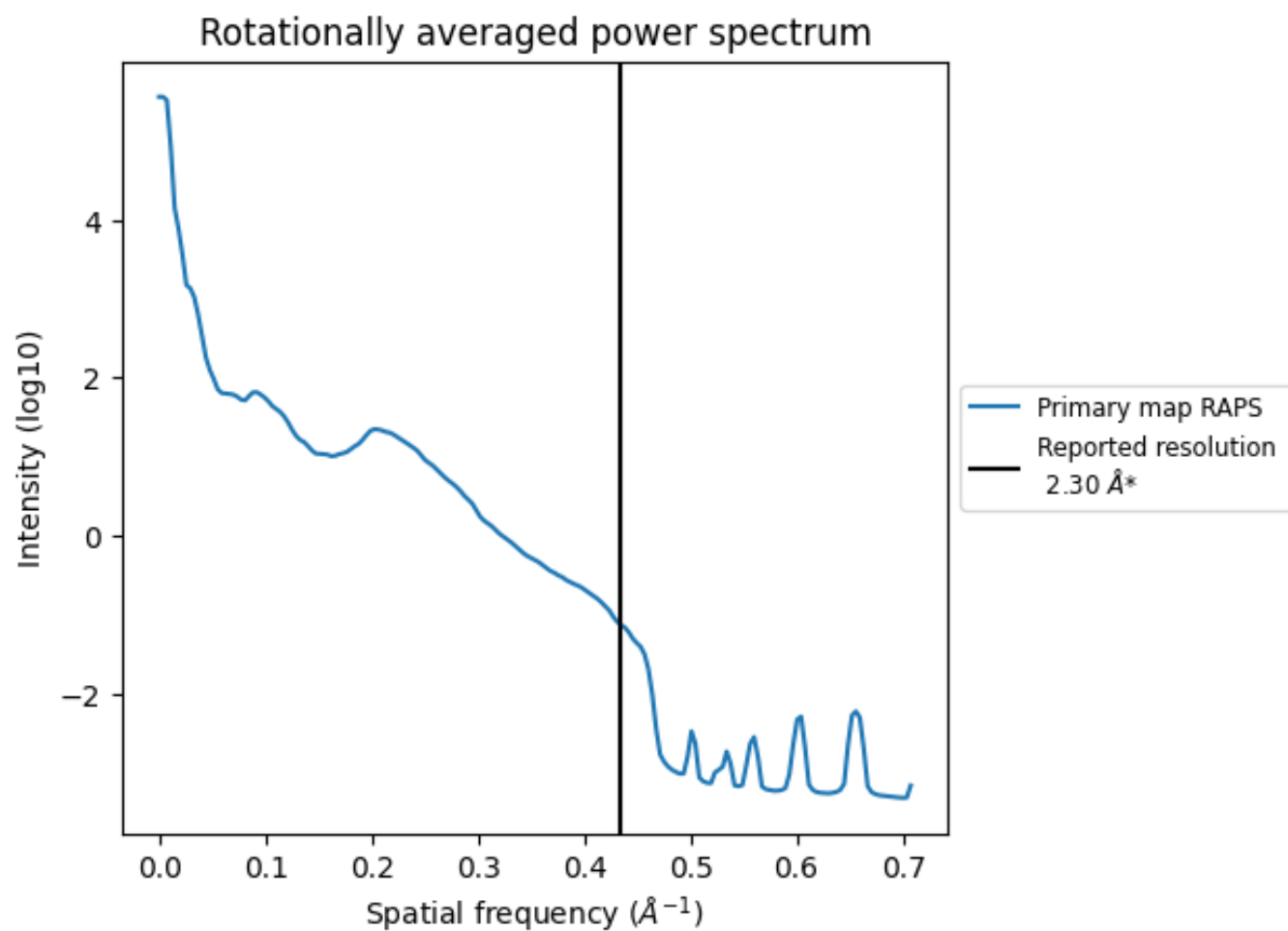
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 85 nm^3 ; this corresponds to an approximate mass of 77 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 ⓘ

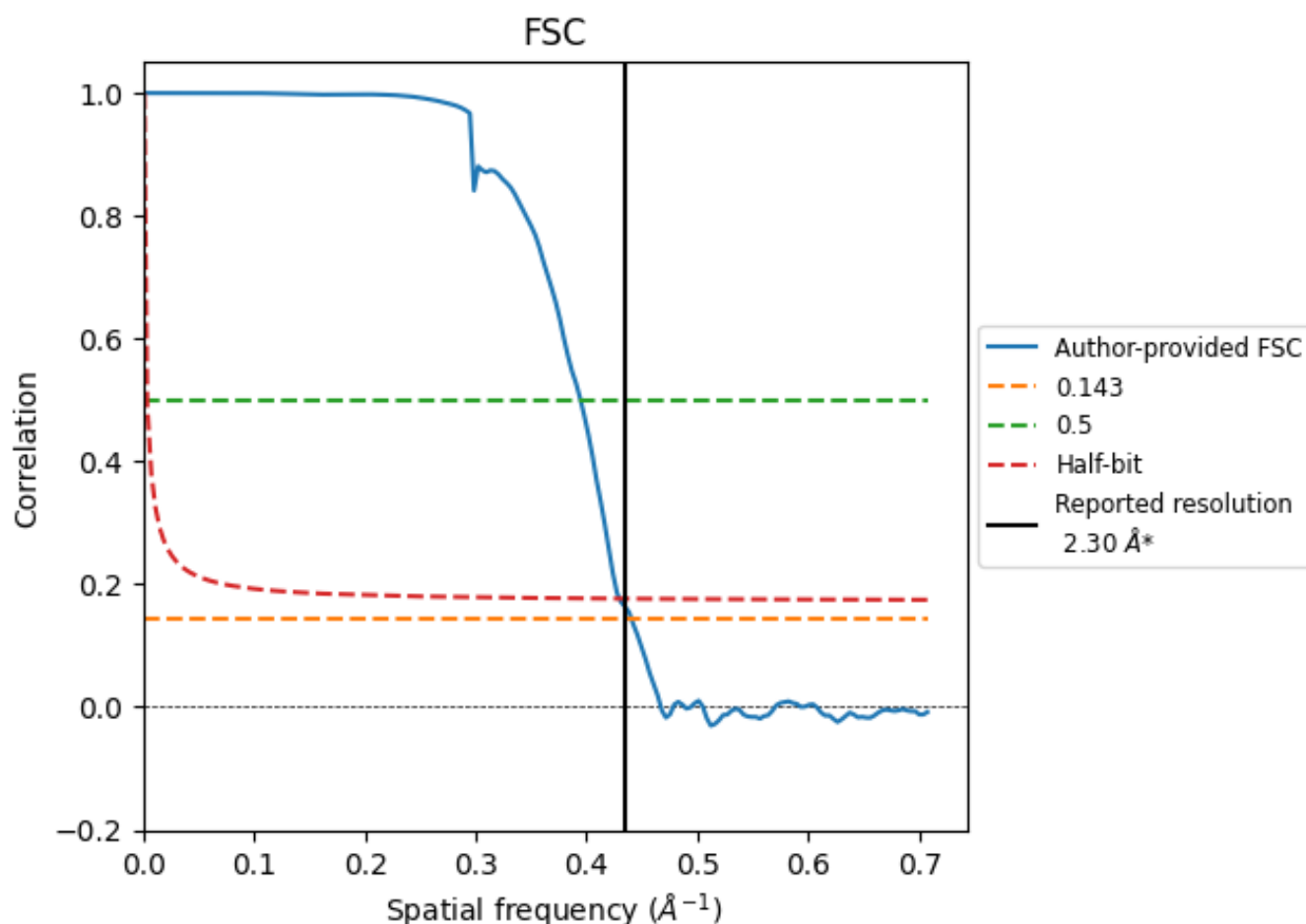


*Reported resolution corresponds to spatial frequency of 0.435 Å⁻¹

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 0.435 Å⁻¹

8.2 Resolution estimates [i](#)

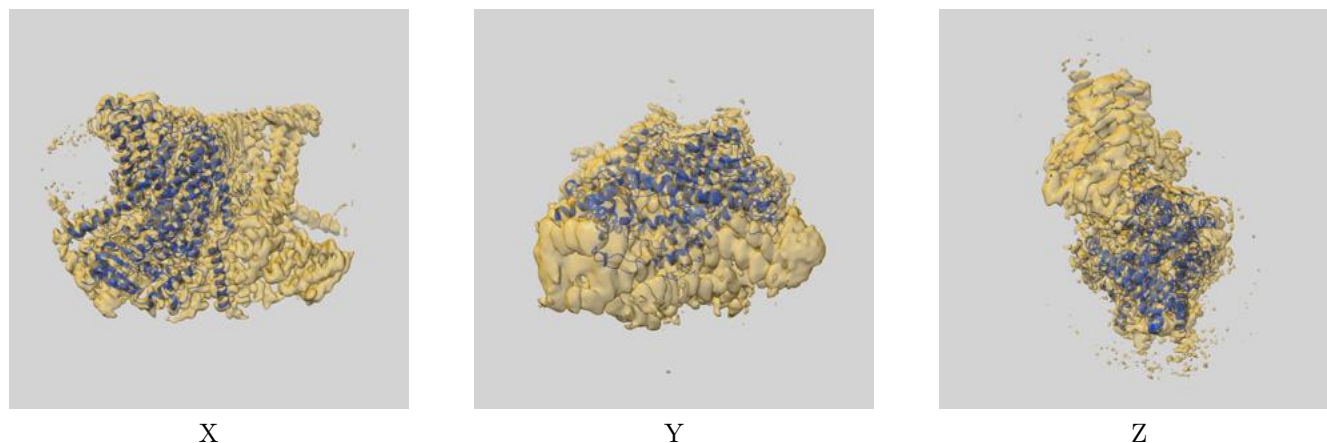
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.30	-	-
Author-provided FSC curve	2.27	2.53	2.33
Unmasked-calculated*	-	-	-

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

9 Map-model fit [i](#)

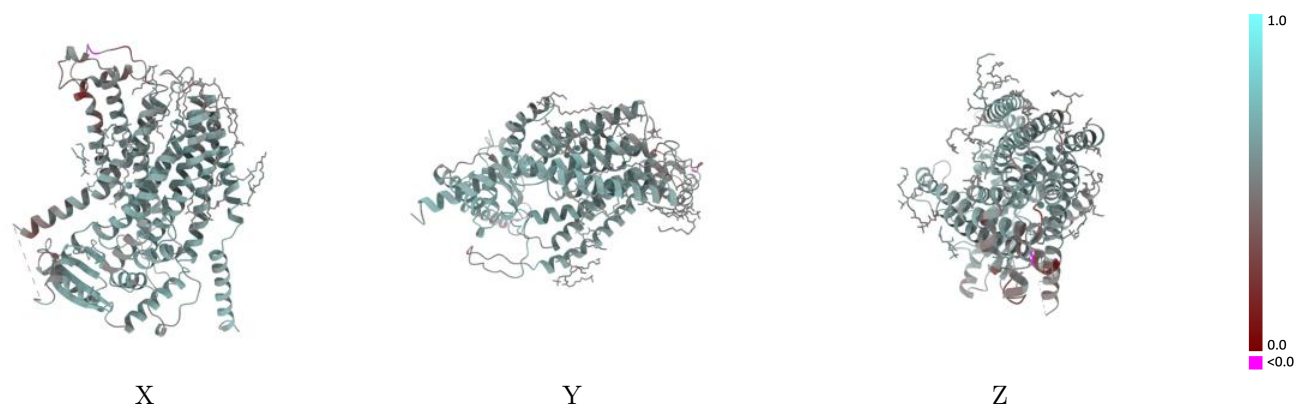
This section contains information regarding the fit between EMDB map EMD-24731 and PDB model 7RXH. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

9.1 Map-model overlay [i](#)



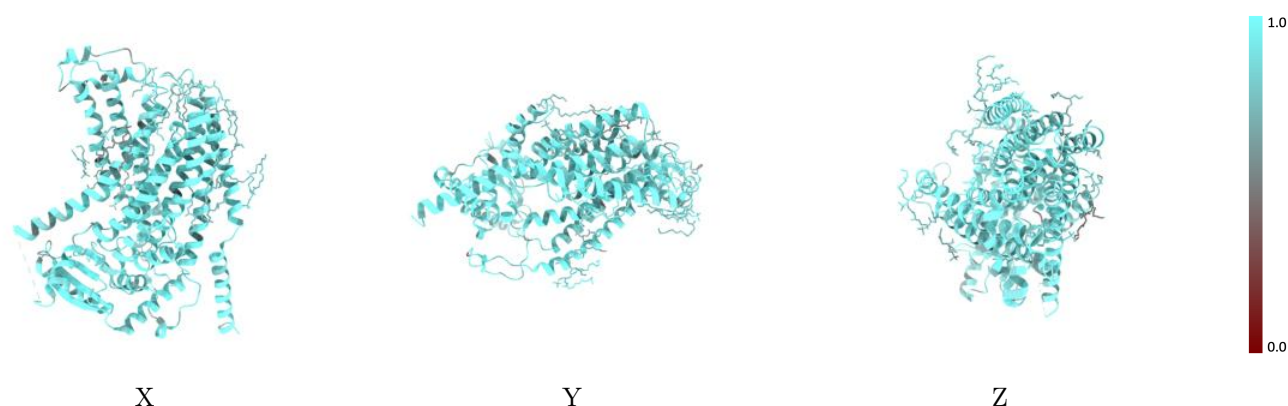
The images above show the 3D surface view of the map at the recommended contour level 0.0049 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



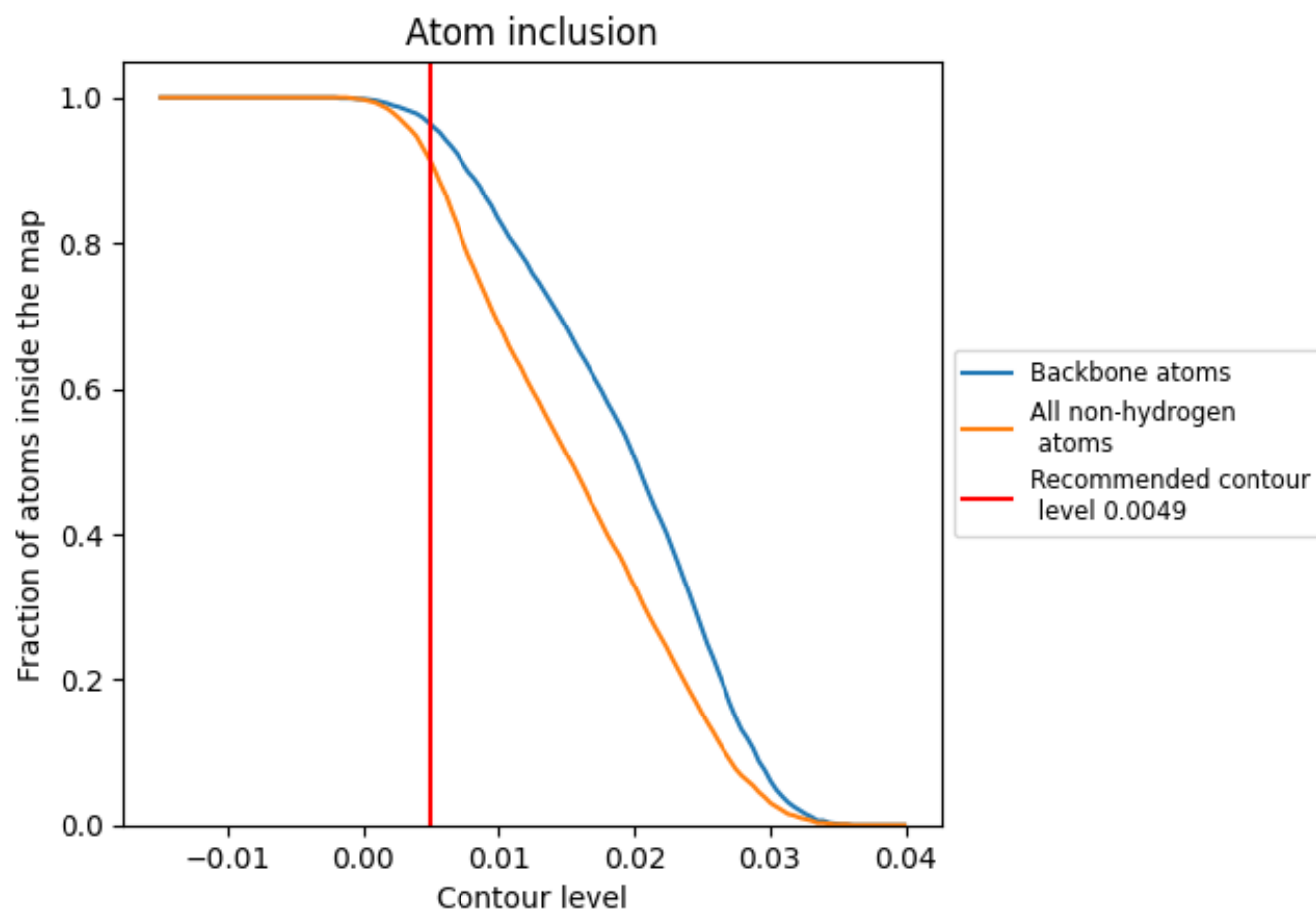
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0049).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% 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 (0.0049) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9140	0.5560
A	0.9140	0.5560

