



Full wwPDB EM Validation Report ⓘ

Jan 1, 2025 – 08:07 AM EST

PDB ID : 8S9B
EMDB ID : EMD-40238
Title : Cryo-EM structure of Nav1.7 with LCM
Authors : Fan, X.; Huang, J.; Yan, N.
Deposited on : 2023-03-27
Resolution : 2.90 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.dev113
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.40

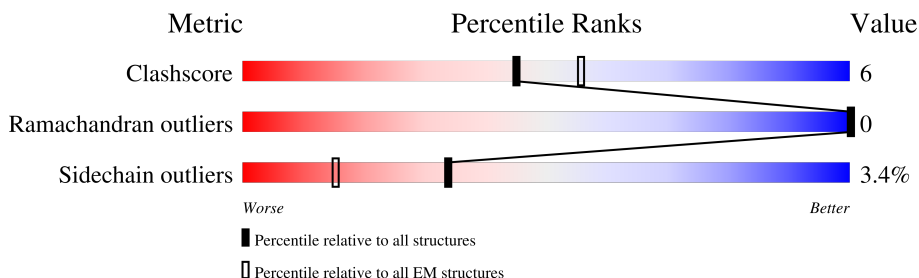
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.90 Å.

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



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1988	
2	B	218	
3	C	215	
4	D	2	
4	E	2	
4	F	2	

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 13409 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 Sodium channel protein type 9 subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1273	Total	C	N	O	S	0	0
			10262	6802	1613	1770	77		

- Molecule 2 is a protein called Sodium channel subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	173	Total	C	N	O	S	0	0
			1416	902	232	272	10		

- Molecule 3 is a protein called Sodium channel subunit beta-2.

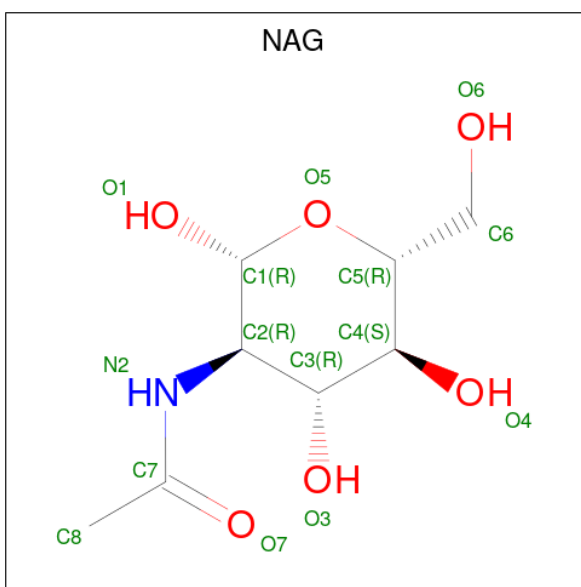
Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	119	Total	C	N	O	S	0	0
			971	609	172	180	10		

- Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



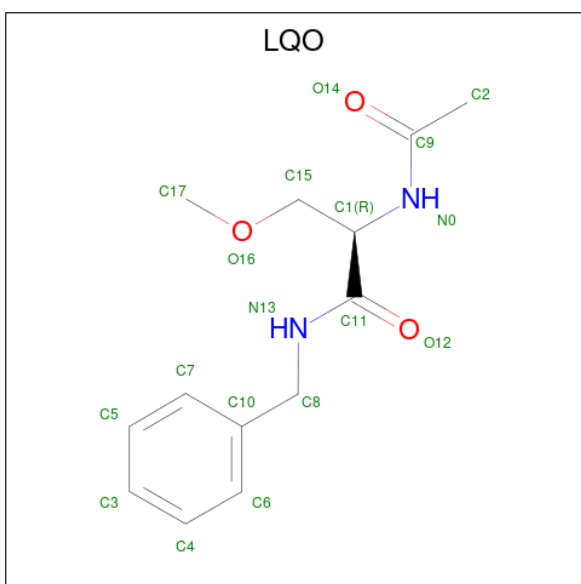
Mol	Chain	Residues	Atoms				AltConf	Trace
4	D	2	Total	C	N	O	0	0
			28	16	2	10		
4	E	2	Total	C	N	O	0	0
			28	16	2	10		
4	F	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



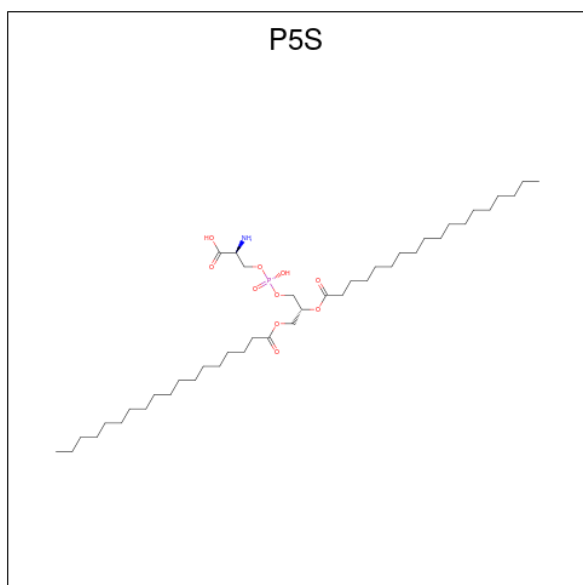
Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	

- Molecule 6 is lacosamide (three-letter code: LQO) (formula: $C_{13}H_{18}N_2O_3$) (labeled as "Ligand of Interest" by depositor).



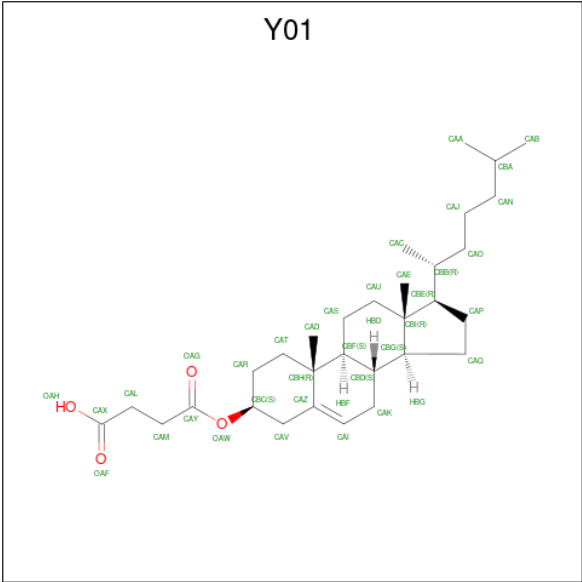
Mol	Chain	Residues	Atoms				AltConf
6	A	1	Total	C	N	O	0
			18	13	2	3	
6	A	1	Total	C	N	O	0
			18	13	2	3	

- Molecule 7 is O-[(R)-{[(2R)-2,3-bis(octadecanoyloxy)propyl]oxy} (hydroxy)phosphoryl]-L-serine (three-letter code: P5S) (formula: C₄₂H₈₂NO₁₀P).



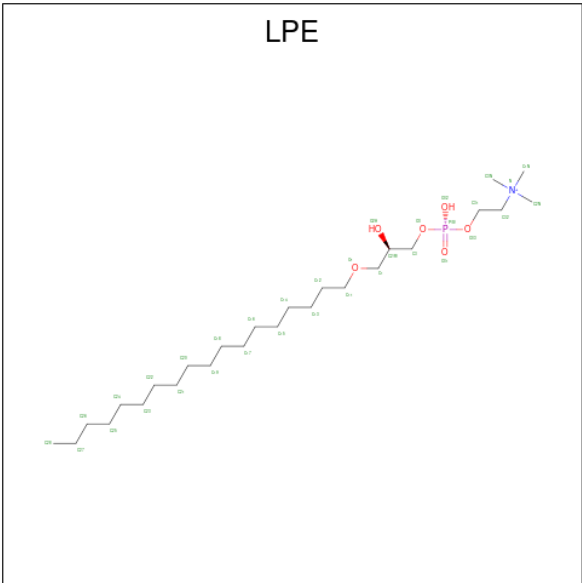
Mol	Chain	Residues	Atoms					AltConf
7	A	1	Total	C	N	O	P	0
			35	24	1	9	1	
7	A	1	Total	C	N	O	P	0
			41	31	1	8	1	

- Molecule 8 is CHOLESTEROL HEMISUCCINATE (three-letter code: Y01) (formula: C₃₁H₅₀O₄).



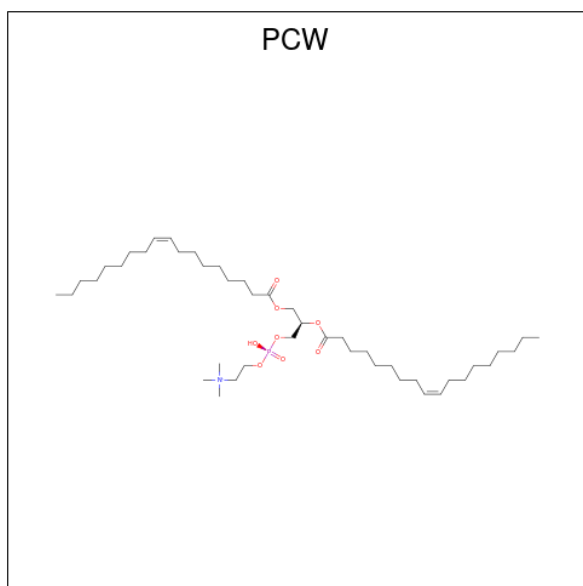
Mol	Chain	Residues	Atoms			AltConf
8	A	1	Total	C	O	0
			35	31	4	
8	A	1	Total	C	O	0
			35	31	4	
8	A	1	Total	C	O	0
			35	31	4	
8	A	1	Total	C	O	0
			35	31	4	

- Molecule 9 is 1-O-OCTADECYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: LPE) (formula: C₂₆H₅₇NO₆P).



Mol	Chain	Residues	Atoms					AltConf
9	A	1	Total	C	N	O	P	0
			25	17	1	6	1	
9	A	1	Total	C	N	O	P	0
			28	20	1	6	1	
9	A	1	Total	C	N	O	P	0
			25	17	1	6	1	
9	A	1	Total	C	N	O	P	0
			25	17	1	6	1	
9	A	1	Total	C	N	O	P	0
			25	17	1	6	1	
9	A	1	Total	C	N	O	P	0
			25	17	1	6	1	
9	A	1	Total	C	N	O	P	0
			25	17	1	6	1	

- Molecule 10 is 1,2-DIOLEOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: PCW) (formula: $C_{44}H_{85}NO_8P$).



Mol	Chain	Residues	Atoms					AltConf
10	A	1	Total	C	N	O	P	0
			47	37	1	8	1	
10	A	1	Total	C	N	O	P	0
			44	34	1	8	1	
10	A	1	Total	C	N	O	P	0
			44	34	1	8	1	

- Molecule 11 is water.

Mol	Chain	Residues	Atoms		AltConf
11	A	21	Total 21	O 21	0
11	B	9	Total 9	O 9	0




- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  50% 50%

MAG1
MAG2

- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  100%

MAG1
MAG2

- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  50% 50%

MAG1
MAG2

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	300001	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)	1200	Depositor
Maximum defocus (nm)	1400	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	6.825	Depositor
Minimum map value	-3.611	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.127	Depositor
Recommended contour level	0.7	Depositor
Map size (\AA)	356.47998, 356.47998, 356.47998	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.114, 1.114, 1.114	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, PCW, LQO, P5S, Y01, LPE

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	0.35	0/10511	0.50	0/14239
2	B	0.36	0/1442	0.50	0/1949
3	C	0.33	0/993	0.58	0/1343
All	All	0.35	0/12946	0.50	0/17531

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	217	ARG	Sidechain
1	A	835	ARG	Sidechain

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	10262	0	10499	117	0
2	B	1416	0	1380	26	0
3	C	971	0	930	18	0
4	D	28	0	25	0	0
4	E	28	0	25	2	0
4	F	28	0	25	0	0
5	A	14	0	13	0	0
5	B	42	0	39	2	0
6	A	36	0	0	0	0
7	A	76	0	96	1	0
8	A	140	0	196	5	0
9	A	203	0	283	0	0
10	A	135	0	184	4	0
11	A	21	0	0	1	0
11	B	9	0	0	2	0
All	All	13409	0	13695	164	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 6.

All (164) 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:1324:VAL:HG21	1:A:1455:VAL:HG21	1.36	1.06
2:B:65:GLU:N	2:B:65:GLU:OE1	2.07	0.88
1:A:737:ILE:HG22	1:A:797:ALA:HB2	1.58	0.86
1:A:78:LEU:H	1:A:89:ILE:HD12	1.43	0.83
1:A:1324:VAL:CG2	1:A:1455:VAL:HG21	2.11	0.81
4:E:2:NAG:O3	4:E:2:NAG:O7	2.04	0.76
1:A:1324:VAL:HG21	1:A:1455:VAL:CG2	2.16	0.74
1:A:1180:ASN:O	1:A:1184:THR:HG23	1.87	0.74
1:A:741:VAL:O	1:A:741:VAL:HG12	1.86	0.74
1:A:802:GLU:N	1:A:802:GLU:OE1	2.21	0.74
2:B:46:ARG:NE	2:B:48:GLU:OE2	2.22	0.73
2:B:71:LEU:CB	2:B:80:LEU:HD23	2.20	0.72
1:A:281:GLU:OE1	1:A:281:GLU:N	2.20	0.70
1:A:1280:TYR:HB3	1:A:1283:LEU:CD1	2.24	0.67
1:A:1574:PHE:HD1	1:A:1580:ILE:HG12	1.58	0.67
2:B:71:LEU:HB3	2:B:80:LEU:HD23	1.77	0.67
3:C:100:GLU:OE1	3:C:100:GLU:N	2.28	0.67
2:B:78:LEU:HD21	2:B:80:LEU:HD21	1.78	0.66
1:A:71:VAL:HG11	1:A:93:LYS:HG3	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1625:LYS:O	1:A:1631:ARG:NH1	2.28	0.65
8:A:2007:Y01:HAA1	8:A:2007:Y01:HBB	1.78	0.65
2:B:114:ASN:ND2	5:B:303:NAG:O7	2.29	0.64
1:A:1751:VAL:HG23	1:A:1752:VAL:HG23	1.80	0.64
1:A:78:LEU:H	1:A:89:ILE:CD1	2.09	0.64
1:A:1458:ASP:OD1	1:A:1462:GLN:HG2	2.01	0.61
1:A:1644:LEU:HD11	1:A:1759:ILE:HD11	1.81	0.60
1:A:401:VAL:HG11	1:A:960:LEU:CD2	2.32	0.59
2:B:71:LEU:HB2	2:B:80:LEU:HD23	1.84	0.59
2:B:81:GLU:HB2	11:B:401:HOH:O	2.02	0.59
1:A:1505:GLN:HA	1:A:1508:ILE:HD12	1.85	0.59
3:C:135:ARG:NE	3:C:135:ARG:HA	2.19	0.58
1:A:53:LEU:O	1:A:99:ARG:NH2	2.37	0.57
3:C:72:CYS:SG	3:C:75:CYS:N	2.77	0.57
1:A:741:VAL:O	1:A:741:VAL:CG1	2.51	0.57
2:B:69:LYS:NZ	11:B:401:HOH:O	2.32	0.57
3:C:70:GLN:NE2	3:C:72:CYS:O	2.38	0.57
1:A:210:VAL:HA	1:A:213:LEU:HD12	1.85	0.57
1:A:348:SER:HB3	10:A:2010:PCW:H352	1.87	0.56
1:A:413:ILE:CG2	1:A:417:LYS:HE3	2.34	0.56
1:A:1280:TYR:HB3	1:A:1283:LEU:HD12	1.87	0.56
1:A:89:ILE:HD11	1:A:99:ARG:HH12	1.72	0.55
1:A:1299:ARG:NH2	11:A:2101:HOH:O	2.40	0.55
1:A:737:ILE:CG2	1:A:797:ALA:HB2	2.31	0.55
1:A:1580:ILE:HD13	10:A:2019:PCW:H122	1.89	0.55
1:A:78:LEU:N	1:A:89:ILE:HD12	2.18	0.54
1:A:179:GLY:O	1:A:185:ARG:NH1	2.40	0.54
1:A:1504:ILE:HG13	1:A:1505:GLN:H	1.71	0.54
3:C:61:LYS:N	3:C:61:LYS:HD3	2.22	0.54
3:C:51:THR:HG22	3:C:109:ASP:OD1	2.08	0.54
1:A:1742:SER:O	1:A:1746:ILE:HD12	2.08	0.54
1:A:1750:VAL:HG22	1:A:1754:MET:HE2	1.89	0.53
2:B:78:LEU:HD23	2:B:80:LEU:HG	1.90	0.53
3:C:135:ARG:HA	3:C:135:ARG:HE	1.74	0.53
1:A:418:GLN:HA	1:A:418:GLN:OE1	2.09	0.53
3:C:123:GLY:O	3:C:143:HIS:ND1	2.40	0.53
1:A:1574:PHE:HB3	1:A:1580:ILE:HD11	1.91	0.53
1:A:1453:ILE:O	1:A:1457:ILE:HG12	2.09	0.52
2:B:78:LEU:CD2	2:B:80:LEU:HG	2.41	0.51
1:A:346:THR:HG22	1:A:347:PHE:N	2.26	0.51
1:A:1191:HIS:CD2	1:A:1193:TRP:HB2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:837:PHE:CE1	8:A:2007:Y01:HAA3	2.46	0.50
4:E:2:NAG:HO3	4:E:2:NAG:C7	2.19	0.50
1:A:1508:ILE:HG23	1:A:1564:LEU:HD22	1.94	0.50
1:A:24:GLU:HG2	1:A:86:LYS:HZ2	1.76	0.50
1:A:413:ILE:HG23	1:A:417:LYS:HE3	1.92	0.50
3:C:68:THR:HB	3:C:77:GLU:HB3	1.94	0.50
1:A:1260:TRP:CD2	8:A:2005:Y01:HAE3	2.47	0.49
1:A:1352:ASN:O	1:A:1356:GLY:N	2.38	0.49
2:B:64:THR:HG22	2:B:65:GLU:H	1.76	0.49
1:A:1184:THR:HG21	2:B:182:TYR:HB2	1.94	0.49
1:A:1223:LYS:O	1:A:1227:GLU:HG2	2.11	0.49
1:A:1282:ASP:N	1:A:1282:ASP:OD1	2.46	0.49
2:B:78:LEU:HD21	2:B:80:LEU:CD2	2.42	0.49
1:A:791:MET:SD	1:A:816:VAL:HG11	2.53	0.49
1:A:1499:ARG:HB3	1:A:1499:ARG:NH1	2.28	0.49
2:B:64:THR:HG22	2:B:65:GLU:N	2.28	0.48
1:A:146:ASN:N	1:A:146:ASN:OD1	2.47	0.48
1:A:1584:VAL:O	1:A:1588:ILE:HG23	2.13	0.48
1:A:810:ILE:HD12	1:A:811:PHE:N	2.27	0.48
1:A:1377:SER:OG	1:A:1378:GLN:N	2.44	0.48
1:A:223:LYS:O	1:A:227:VAL:HG22	2.14	0.48
1:A:395:ASN:ND2	1:A:1755:TYR:CD1	2.81	0.48
1:A:1220:LYS:O	1:A:1224:ILE:HG12	2.13	0.48
1:A:1574:PHE:CD1	1:A:1580:ILE:HG12	2.42	0.48
2:B:27:GLU:N	2:B:27:GLU:OE1	2.47	0.48
2:B:42:SER:O	2:B:42:SER:OG	2.31	0.48
1:A:1385:LEU:O	1:A:1388:ASN:ND2	2.47	0.47
1:A:1442:ILE:O	1:A:1446:PHE:HB3	2.14	0.47
3:C:107:LYS:O	3:C:107:LYS:HG2	2.13	0.47
1:A:16:THR:HG22	1:A:17:LYS:N	2.29	0.47
8:A:2020:Y01:HAC2	8:A:2020:Y01:HAN2	1.96	0.47
2:B:116:SER:HB3	2:B:146:VAL:HG12	1.97	0.47
1:A:735:LYS:HG3	1:A:736:CYS:N	2.30	0.47
1:A:810:ILE:O	1:A:814:LEU:HD12	2.14	0.47
1:A:1408:TRP:CE3	1:A:1412:MET:CE	2.97	0.47
1:A:248:VAL:HG21	1:A:400:VAL:HG21	1.97	0.46
1:A:401:VAL:HG11	1:A:960:LEU:HD23	1.98	0.46
2:B:168:VAL:O	2:B:172:ILE:HG22	2.15	0.46
3:C:105:PRO:C	3:C:107:LYS:N	2.67	0.46
1:A:1177:ILE:N	1:A:1177:ILE:HD12	2.32	0.45
1:A:1280:TYR:CD1	1:A:1280:TYR:N	2.83	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:182:TYR:C	2:B:182:TYR:CD1	2.89	0.45
1:A:86:LYS:HB3	1:A:102:ALA:HB3	1.98	0.45
1:A:811:PHE:HA	1:A:814:LEU:HD12	1.98	0.45
3:C:58:VAL:HG21	3:C:105:PRO:O	2.17	0.45
2:B:55:THR:HB	2:B:104:LEU:HD22	1.98	0.45
1:A:901:ASP:OD1	1:A:902:ASP:N	2.49	0.45
1:A:1558:PHE:O	1:A:1561:GLU:HG3	2.17	0.45
1:A:12:PHE:O	1:A:70:MET:HE3	2.17	0.44
3:C:68:THR:OG1	3:C:126:ASN:HB3	2.17	0.44
1:A:1543:MET:HA	1:A:1546:VAL:HG22	1.99	0.44
2:B:71:LEU:HB3	2:B:80:LEU:CD2	2.45	0.44
1:A:1486:LYS:HA	1:A:1639:MET:HE2	1.99	0.44
1:A:1228:TYR:O	1:A:1232:ILE:HG13	2.18	0.44
1:A:1494:GLN:N	1:A:1494:GLN:OE1	2.50	0.44
1:A:23:ILE:HG21	1:A:86:LYS:HA	1.99	0.44
1:A:221:ALA:O	1:A:224:THR:OG1	2.31	0.44
1:A:1577:GLY:HA2	10:A:2012:PCW:H121	1.99	0.44
1:A:33:LYS:O	1:A:34:GLU:C	2.57	0.44
3:C:96:GLN:OE1	3:C:98:ARG:NE	2.49	0.43
1:A:1360:PRO:O	1:A:1363:GLN:O	2.36	0.43
1:A:211:SER:O	1:A:215:THR:OG1	2.32	0.43
1:A:179:GLY:O	1:A:185:ARG:NH2	2.51	0.43
2:B:96:ARG:NH2	5:B:301:NAG:O7	2.51	0.43
1:A:1512:VAL:HG13	1:A:1565:LYS:HG2	2.01	0.43
3:C:110:VAL:HG13	3:C:110:VAL:O	2.18	0.43
1:A:320:ASP:OD1	1:A:320:ASP:N	2.47	0.43
2:B:34:MET:O	2:B:111:VAL:HG23	2.19	0.43
3:C:108:TYR:CD1	3:C:108:TYR:N	2.86	0.43
1:A:1348:TYR:CE1	1:A:1384:ASN:HB2	2.54	0.43
1:A:1432:TYR:HB2	8:A:2007:Y01:HAT2	2.01	0.43
1:A:54:GLU:HG2	1:A:55:ALA:N	2.34	0.42
1:A:264:LEU:O	1:A:268:MET:HB2	2.20	0.42
1:A:1673:ASP:N	1:A:1673:ASP:OD1	2.53	0.42
1:A:1512:VAL:CG1	1:A:1565:LYS:HG2	2.50	0.42
1:A:1683:THR:HG22	1:A:1684:PHE:N	2.34	0.42
1:A:935:CYS:C	1:A:944:CYS:SG	2.98	0.42
1:A:1676:ASN:O	1:A:1680:ASN:ND2	2.52	0.42
1:A:1494:GLN:NE2	10:A:2012:PCW:H81	2.34	0.42
3:C:67:TRP:CE2	3:C:127:CYS:HB3	2.55	0.41
1:A:183:PHE:O	1:A:189:ASN:ND2	2.42	0.41
1:A:743:ASP:OD1	1:A:994:ARG:NH2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:932:MET:HG2	1:A:936:MET:HE2	2.02	0.41
1:A:1694:ILE:HG21	1:A:1703:LEU:HD12	2.02	0.41
3:C:126:ASN:HD22	3:C:127:CYS:N	2.18	0.41
1:A:1317:ALA:HA	1:A:1459:ASN:OD1	2.20	0.41
1:A:147:ASN:O	1:A:147:ASN:CG	2.58	0.41
1:A:251:LEU:HB2	1:A:1630:ILE:HG22	2.01	0.41
1:A:1246:ILE:HD11	7:A:2004:P5S:H55A	2.01	0.41
1:A:1402:VAL:HA	1:A:1408:TRP:HB3	2.03	0.41
1:A:26:ARG:HD2	1:A:83:ALA:O	2.21	0.41
1:A:1552:VAL:O	1:A:1556:ILE:HG23	2.20	0.41
1:A:89:ILE:HD11	1:A:99:ARG:NH1	2.35	0.41
1:A:1445:SER:O	1:A:1445:SER:OG	2.37	0.41
1:A:810:ILE:HD12	1:A:810:ILE:C	2.41	0.41
1:A:1655:MET:HG2	1:A:1743:TYR:OH	2.21	0.41
1:A:928:TRP:HZ3	1:A:932:MET:HE1	1.86	0.41
1:A:979:ILE:HG23	1:A:980:GLU:O	2.21	0.41
1:A:90:VAL:HG12	1:A:98:PHE:HB2	2.03	0.40
1:A:765:HIS:HB2	1:A:767:PRO:HD2	2.03	0.40
1:A:1659:ALA:O	1:A:1663:MET:HG3	2.22	0.40
1:A:1707:ILE:O	1:A:1707:ILE:HG22	2.21	0.40
2:B:29:GLU:HA	2:B:145:GLU:O	2.21	0.40
2:B:84:GLU:H	2:B:84:GLU:CD	2.25	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	1261/1988 (63%)	1199 (95%)	62 (5%)	0	100	100
2	B	171/218 (78%)	164 (96%)	7 (4%)	0	100	100
3	C	117/215 (54%)	100 (86%)	17 (14%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1549/2421 (64%)	1463 (94%)	86 (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	1136/1778 (64%)	1099 (97%)	37 (3%)	33	68
2	B	157/190 (83%)	154 (98%)	3 (2%)	52	81
3	C	111/193 (58%)	103 (93%)	8 (7%)	12	35
All	All	1404/2161 (65%)	1356 (97%)	48 (3%)	34	67

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	MET
1	A	86	LYS
1	A	119	SER
1	A	130	MET
1	A	176	PHE
1	A	193	PHE
1	A	198	PHE
1	A	217	ARG
1	A	730	TRP
1	A	738	TYR
1	A	766	HIS
1	A	770	GLU
1	A	772	PHE
1	A	798	MET
1	A	814	LEU
1	A	819	SER
1	A	853	ASN
1	A	934	ASP

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Mol	Chain	Res	Type
1	A	950	MET
1	A	952	MET
1	A	1014	PHE
1	A	1180	ASN
1	A	1193	TRP
1	A	1293	ARG
1	A	1318	ILE
1	A	1408	TRP
1	A	1450	ASN
1	A	1459	ASN
1	A	1486	LYS
1	A	1504	ILE
1	A	1511	LEU
1	A	1520	SER
1	A	1528	ASN
1	A	1603	PHE
1	A	1661	PHE
1	A	1711	LYS
1	A	1761	GLU
2	B	93	ASN
2	B	95	SER
2	B	148	ASP
3	C	30	MET
3	C	44	SER
3	C	61	LYS
3	C	64	SER
3	C	94	ARG
3	C	97	ASP
3	C	126	ASN
3	C	135	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	HIS
1	A	408	GLN
1	A	780	ASN
1	A	1461	ASN
3	C	126	ASN
3	C	139	HIS

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 ⓘ

6 monosaccharides 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	$\# Z > 2$
4	NAG	D	1	1,4	14,14,15	0.34	0	17,19,21	0.62	0
4	NAG	D	2	4	14,14,15	0.28	0	17,19,21	0.99	2 (11%)
4	NAG	E	1	1,4	14,14,15	0.40	0	17,19,21	1.67	3 (17%)
4	NAG	E	2	4	14,14,15	0.28	0	17,19,21	0.72	0
4	NAG	F	1	2,4	14,14,15	0.32	0	17,19,21	0.68	0
4	NAG	F	2	4	14,14,15	0.35	0	17,19,21	0.91	1 (5%)

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
4	NAG	D	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1
4	NAG	E	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	E	2	4	-	1/6/23/26	0/1/1/1
4	NAG	F	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	1	NAG	C4-C3-C2	-4.27	104.75	111.02
4	E	1	NAG	C1-O5-C5	3.06	116.28	112.19
4	E	1	NAG	O4-C4-C3	2.65	116.62	110.38
4	F	2	NAG	C1-O5-C5	2.31	115.29	112.19
4	D	2	NAG	C2-N2-C7	2.18	125.83	122.90
4	D	2	NAG	C1-C2-N2	-2.00	107.28	110.43

There are no chirality outliers.

All (3) torsion outliers are listed below:

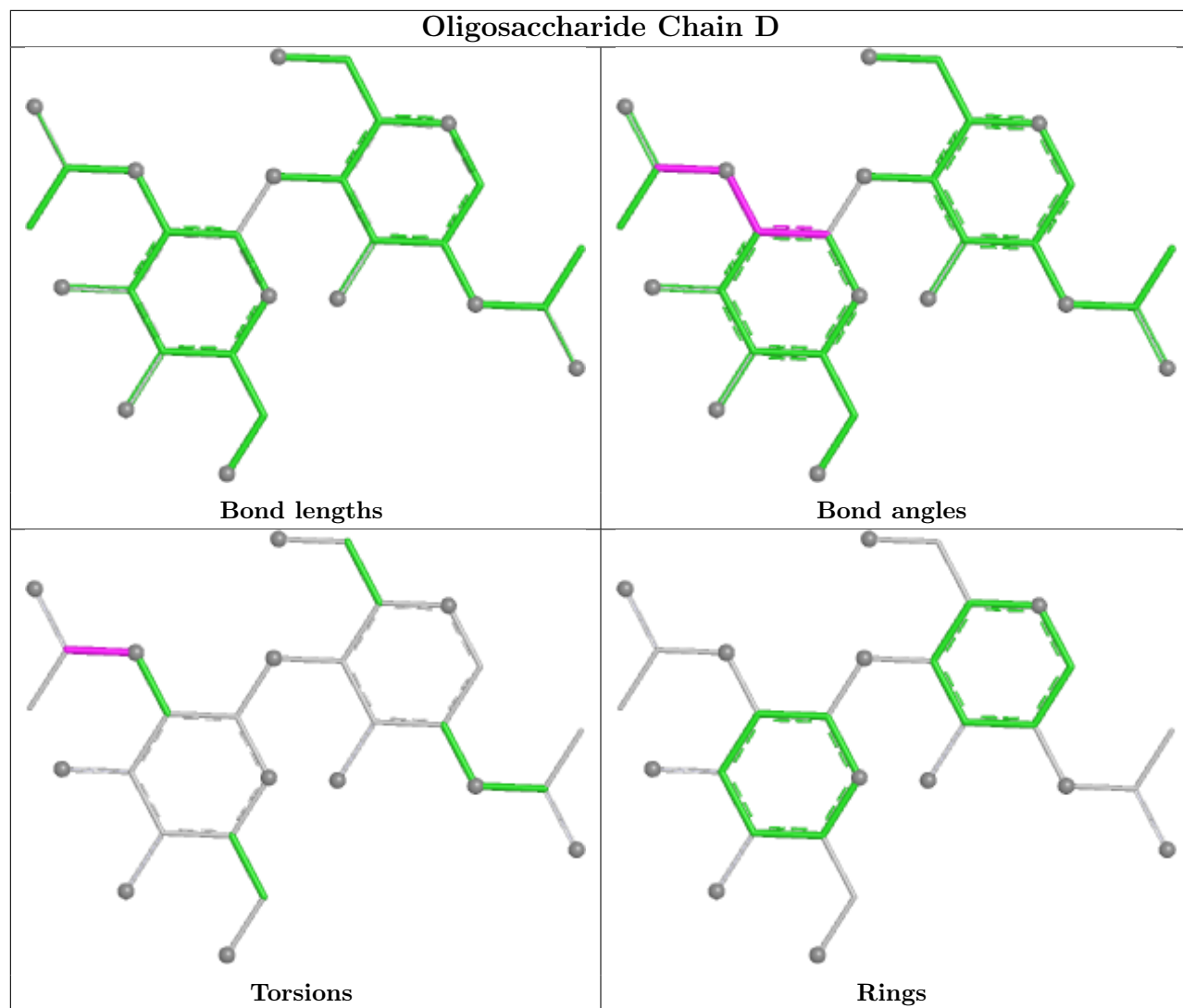
Mol	Chain	Res	Type	Atoms
4	D	2	NAG	C8-C7-N2-C2
4	D	2	NAG	O7-C7-N2-C2
4	E	2	NAG	C3-C2-N2-C7

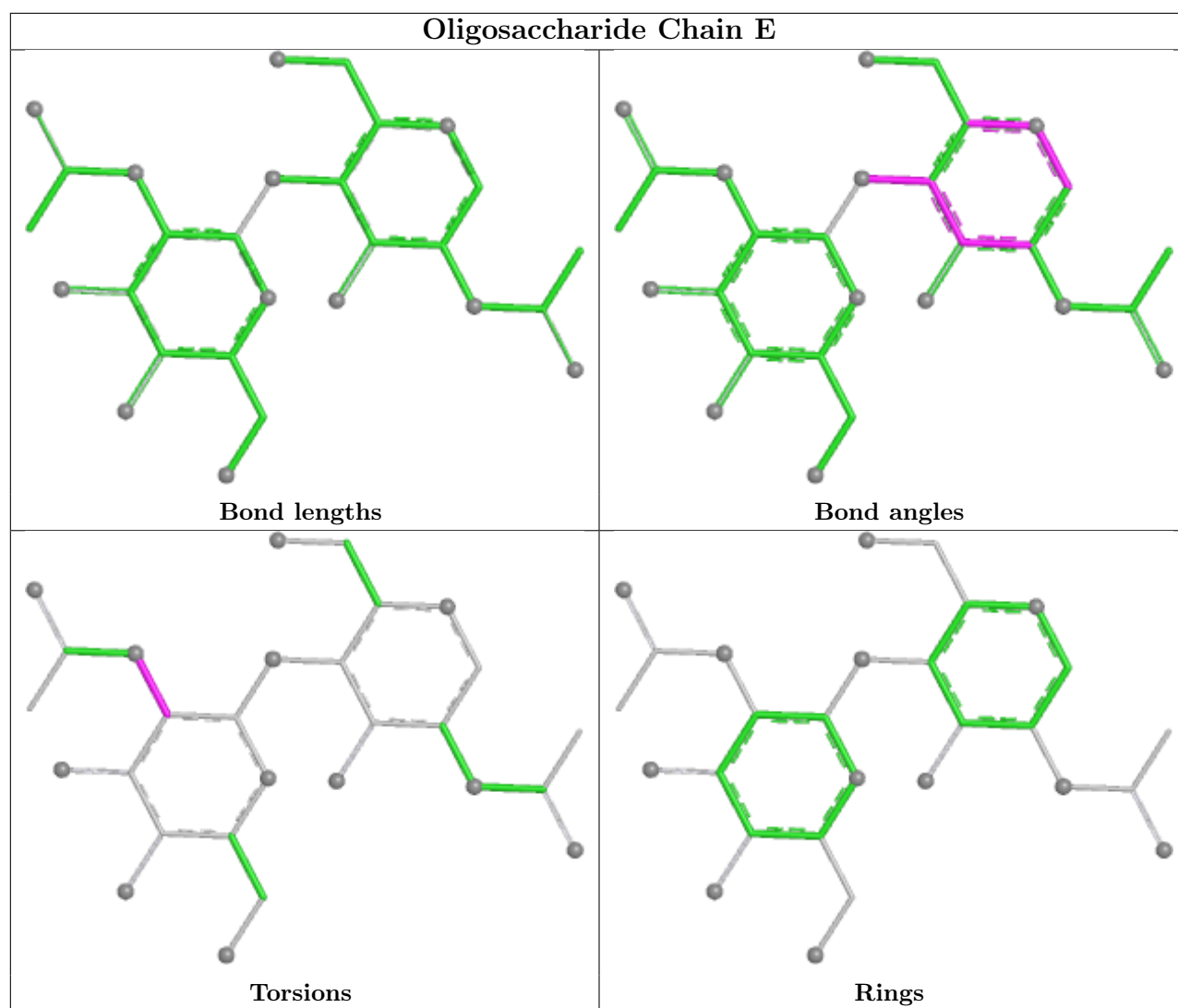
There are no ring outliers.

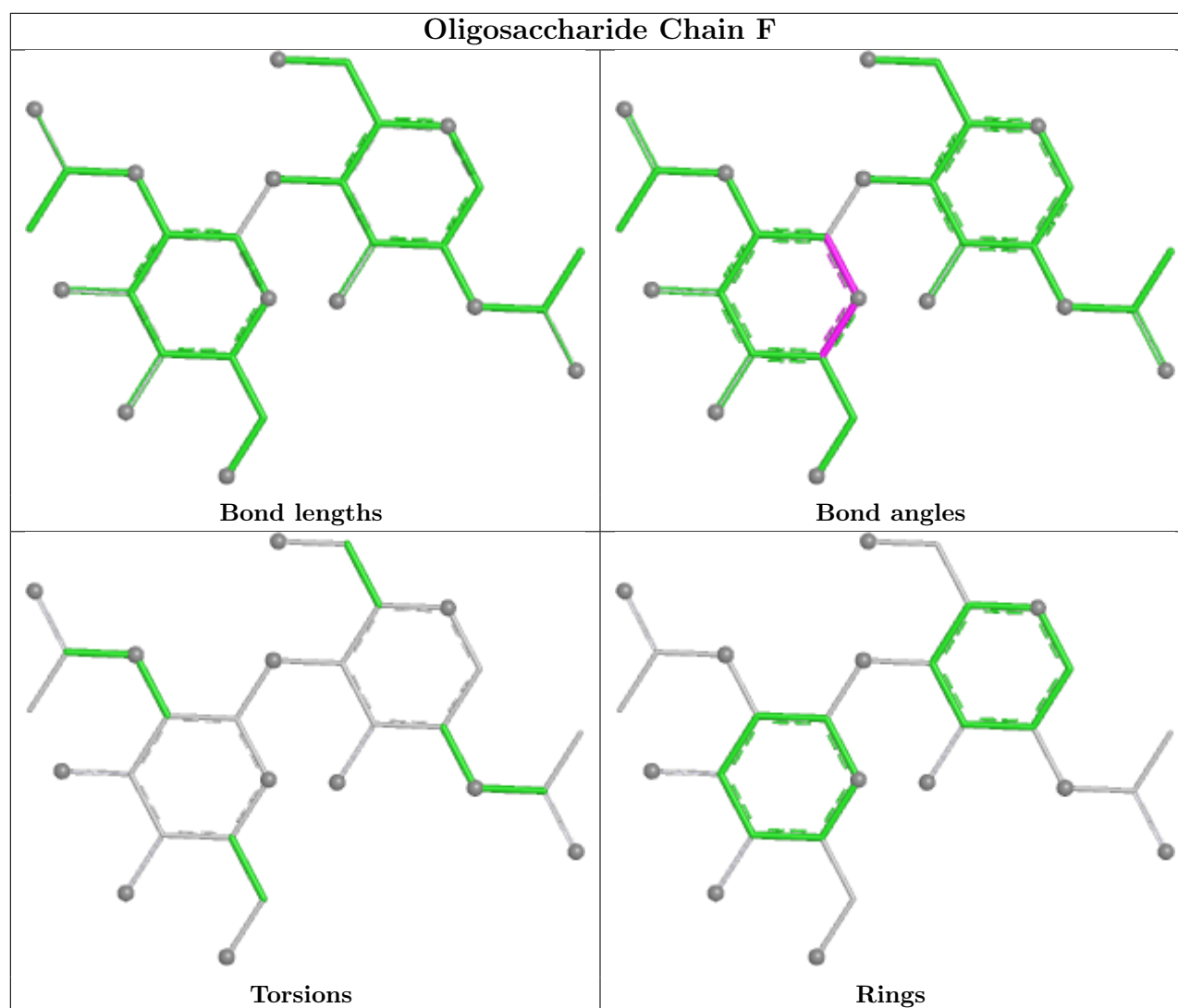
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	2	NAG	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.







5.6 Ligand geometry [i](#)

23 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	$\# Z > 2$
8	Y01	A	2006	-	38,38,38	0.48	0	57,57,57	0.54	0
9	LPE	A	2014	-	24,24,33	0.28	0	28,30,39	0.37	0
6	LQO	A	2003	-	18,18,18	3.67	9 (50%)	22,22,22	0.86	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	LPE	A	2013	-	24,24,33	0.45	0	28,30,39	0.59	0
8	Y01	A	2005	-	38,38,38	0.54	0	57,57,57	0.93	3 (5%)
9	LPE	A	2015	-	24,24,33	0.27	0	28,30,39	0.33	0
6	LQO	A	2002	-	18,18,18	3.61	10 (55%)	22,22,22	1.95	4 (18%)
8	Y01	A	2020	-	38,38,38	0.56	0	57,57,57	0.82	1 (1%)
5	NAG	B	303	2	14,14,15	0.33	0	17,19,21	0.56	0
8	Y01	A	2007	-	38,38,38	0.49	0	57,57,57	0.95	4 (7%)
9	LPE	A	2018	-	24,24,33	0.28	0	28,30,39	0.39	0
7	P5S	A	2011	-	40,40,53	0.52	0	43,45,60	0.37	0
9	LPE	A	2016	-	24,24,33	0.28	0	28,30,39	0.36	0
9	LPE	A	2009	-	27,27,33	0.27	0	31,33,39	0.36	0
5	NAG	A	2001	1	14,14,15	0.35	0	17,19,21	0.61	0
10	PCW	A	2012	-	43,43,53	0.47	0	49,51,61	0.51	0
7	P5S	A	2004	-	33,34,53	0.47	0	34,40,60	0.56	0
5	NAG	B	301	2	14,14,15	0.31	0	17,19,21	0.62	0
9	LPE	A	2017	-	24,24,33	0.28	0	28,30,39	0.37	0
10	PCW	A	2019	-	43,43,53	0.32	0	49,51,61	0.40	0
10	PCW	A	2010	-	46,46,53	0.32	0	52,54,61	0.36	0
9	LPE	A	2008	-	24,24,33	0.32	0	25,27,39	0.41	0
5	NAG	B	302	2	14,14,15	0.31	0	17,19,21	0.74	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
8	Y01	A	2006	-	-	6/19/77/77	0/4/4/4
9	LPE	A	2014	-	-	4/25/25/34	-
6	LQO	A	2003	-	-	2/16/16/16	0/1/1/1
9	LPE	A	2013	-	-	12/25/25/34	-
8	Y01	A	2005	-	-	8/19/77/77	0/4/4/4
9	LPE	A	2015	-	-	3/25/25/34	-
6	LQO	A	2002	-	-	2/16/16/16	0/1/1/1
8	Y01	A	2020	-	-	9/19/77/77	0/4/4/4
5	NAG	B	303	2	-	1/6/23/26	0/1/1/1
8	Y01	A	2007	-	-	4/19/77/77	0/4/4/4
9	LPE	A	2018	-	-	4/25/25/34	-
7	P5S	A	2011	-	-	20/44/44/59	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	LPE	A	2016	-	-	5/25/25/34	-
9	LPE	A	2009	-	-	7/28/28/34	-
5	NAG	A	2001	1	-	0/6/23/26	0/1/1/1
10	PCW	A	2012	-	-	25/47/47/57	-
7	P5S	A	2004	-	-	10/39/39/59	-
5	NAG	B	301	2	-	0/6/23/26	0/1/1/1
9	LPE	A	2017	-	-	1/25/25/34	-
10	PCW	A	2019	-	-	8/47/47/57	-
10	PCW	A	2010	-	-	11/50/50/57	-
9	LPE	A	2008	-	-	4/25/25/34	-
5	NAG	B	302	2	-	0/6/23/26	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	2003	LQO	C4-C6	7.92	1.52	1.38
6	A	2002	LQO	C4-C6	6.93	1.50	1.38
6	A	2003	LQO	C7-C10	6.86	1.52	1.38
6	A	2003	LQO	C3-C5	6.25	1.52	1.38
6	A	2002	LQO	C3-C5	6.14	1.51	1.38
6	A	2002	LQO	C5-C7	-5.69	1.29	1.38
6	A	2002	LQO	C7-C10	5.39	1.49	1.38
6	A	2003	LQO	C11-N13	5.06	1.45	1.33
6	A	2002	LQO	C11-N13	4.94	1.45	1.33
6	A	2002	LQO	C6-C10	-4.42	1.30	1.38
6	A	2003	LQO	C5-C7	-4.36	1.31	1.38
6	A	2002	LQO	C3-C4	-3.92	1.29	1.38
6	A	2003	LQO	C6-C10	-3.85	1.31	1.38
6	A	2003	LQO	C9-N0	3.47	1.45	1.34
6	A	2002	LQO	O14-C9	-3.17	1.16	1.23
6	A	2003	LQO	C3-C4	-3.10	1.31	1.38
6	A	2002	LQO	O12-C11	-3.07	1.17	1.23
6	A	2003	LQO	O12-C11	-2.51	1.18	1.23
6	A	2002	LQO	C9-N0	2.51	1.42	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	2002	LQO	C15-C1-N0	-4.93	100.42	111.31
6	A	2002	LQO	C10-C8-N13	-4.25	104.12	113.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	2002	LQO	C1-C11-N13	3.33	123.68	116.54
8	A	2005	Y01	CAR-CBC-CAV	-2.98	106.84	110.97
8	A	2020	Y01	CAP-CBE-CBB	2.90	116.57	112.18
8	A	2007	Y01	CBF-CBD-CBG	-2.47	105.86	109.09
8	A	2005	Y01	CBC-CAV-CAZ	-2.40	107.91	111.45
8	A	2007	Y01	CAS-CBF-CBD	-2.26	108.62	111.78
8	A	2007	Y01	CBH-CBF-CBD	2.20	115.92	112.71
6	A	2002	LQO	C7-C10-C6	2.13	121.40	118.23
8	A	2005	Y01	CAQ-CBG-CBI	2.06	106.26	103.84
8	A	2007	Y01	CAK-CBD-CBF	2.01	112.04	109.72

There are no chirality outliers.

All (146) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	2003	LQO	C11-C1-C15-O16
6	A	2003	LQO	N0-C1-C15-O16
7	A	2004	P5S	O-C-CA-CB
7	A	2004	P5S	OXT-C-CA-CB
9	A	2008	LPE	C32-C31-O33-P
9	A	2009	LPE	C2-C3-O3-P
9	A	2013	LPE	C3-O3-P-O31
9	A	2013	LPE	C3-O3-P-O32
9	A	2013	LPE	C3-O3-P-O33
9	A	2013	LPE	C31-O33-P-O3
9	A	2013	LPE	C31-O33-P-O32
9	A	2013	LPE	O33-C31-C32-N
9	A	2014	LPE	C32-C31-O33-P
9	A	2015	LPE	C31-O33-P-O31
9	A	2016	LPE	C31-O33-P-O31
9	A	2016	LPE	C32-C31-O33-P
9	A	2018	LPE	C31-O33-P-O31
9	A	2018	LPE	C32-C31-O33-P
10	A	2012	PCW	O4P-C4-C5-N
10	A	2012	PCW	C32-C31-O2-C2
10	A	2012	PCW	C1-O3P-P-O1P
10	A	2012	PCW	C1-O3P-P-O2P
10	A	2012	PCW	C1-O3P-P-O4P
10	A	2012	PCW	C4-O4P-P-O2P
10	A	2019	PCW	C12-C11-O3-C3
10	A	2019	PCW	O11-C11-O3-C3
10	A	2012	PCW	O11-C11-O3-C3

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Mol	Chain	Res	Type	Atoms
10	A	2012	PCW	O31-C31-O2-C2
10	A	2012	PCW	C12-C11-O3-C3
8	A	2020	Y01	CAJ-CAO-CBB-CAC
9	A	2013	LPE	C1-C2-C3-O3
8	A	2020	Y01	CAJ-CAO-CBB-CBE
8	A	2005	Y01	CAJ-CAO-CBB-CBE
8	A	2006	Y01	CAO-CBB-CBE-CBI
8	A	2020	Y01	CAO-CBB-CBE-CBI
7	A	2011	P5S	C2-C1-O19-C17
8	A	2005	Y01	CAJ-CAO-CBB-CAC
8	A	2005	Y01	CAR-CBC-OAW-CAY
7	A	2011	P5S	C17-C20-C21-C22
9	A	2013	LPE	O2H-C2-C3-O3
10	A	2012	PCW	C31-C32-C33-C34
9	A	2018	LPE	C2-C1-O1-C11
8	A	2005	Y01	CAV-CBC-OAW-CAY
8	A	2006	Y01	CAC-CBB-CBE-CBI
7	A	2011	P5S	C39-C38-O37-C2
8	A	2020	Y01	CAM-CAY-OAW-CBC
8	A	2006	Y01	CAC-CBB-CBE-CAP
8	A	2020	Y01	OAG-CAY-OAW-CBC
8	A	2020	Y01	CAC-CBB-CBE-CBI
6	A	2002	LQO	N0-C1-C15-O16
9	A	2013	LPE	C2-C3-O3-P
7	A	2004	P5S	OXT-C-CA-N
8	A	2020	Y01	CAC-CBB-CBE-CAP
8	A	2005	Y01	CAJ-CAN-CBA-CAB
8	A	2007	Y01	CAJ-CAN-CBA-CAA
9	A	2013	LPE	C11-C12-C13-C14
10	A	2012	PCW	C13-C14-C15-C16
8	A	2007	Y01	CAJ-CAN-CBA-CAB
10	A	2012	PCW	C32-C33-C34-C35
9	A	2008	LPE	C11-C12-C13-C14
7	A	2011	P5S	O47-C38-O37-C2
7	A	2011	P5S	C25-C26-C27-C28
8	A	2005	Y01	CAJ-CAN-CBA-CAA
7	A	2004	P5S	C39-C38-O37-C2
7	A	2011	P5S	C28-C29-C30-C31
7	A	2004	P5S	C40-C41-C42-C43
6	A	2002	LQO	C11-C1-C15-O16
7	A	2004	P5S	O47-C38-O37-C2
7	A	2011	P5S	O19-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
10	A	2012	PCW	C16-C17-C18-C19
10	A	2012	PCW	C21-C22-C23-C24
7	A	2011	P5S	C39-C40-C41-C42
10	A	2012	PCW	C35-C36-C37-C38
9	A	2013	LPE	C13-C14-C15-C16
7	A	2011	P5S	C20-C17-O19-C1
7	A	2011	P5S	C43-C44-C45-C46
7	A	2011	P5S	C32-C33-C34-C35
10	A	2019	PCW	C17-C18-C19-C20
10	A	2010	PCW	O2-C31-C32-C33
7	A	2011	P5S	C1-C2-C3-O16
10	A	2012	PCW	C24-C25-C26-C27
7	A	2004	P5S	C45-C46-C48-C49
7	A	2004	P5S	O-C-CA-N
7	A	2011	P5S	O37-C2-C3-O16
10	A	2012	PCW	C36-C37-C38-C39
10	A	2012	PCW	C23-C24-C25-C26
9	A	2009	LPE	C1-C2-C3-O3
7	A	2011	P5S	O18-C17-O19-C1
9	A	2009	LPE	O2H-C2-C3-O3
9	A	2009	LPE	C32-C31-O33-P
9	A	2013	LPE	C32-C31-O33-P
10	A	2010	PCW	C5-C4-O4P-P
10	A	2012	PCW	C5-C4-O4P-P
10	A	2019	PCW	C5-C4-O4P-P
10	A	2012	PCW	C18-C19-C20-C21
10	A	2010	PCW	O2-C2-C3-O3
7	A	2004	P5S	N-CA-CB-OG
9	A	2009	LPE	O33-C31-C32-N
9	A	2014	LPE	O33-C31-C32-N
9	A	2015	LPE	O33-C31-C32-N
9	A	2016	LPE	O33-C31-C32-N
9	A	2017	LPE	O33-C31-C32-N
9	A	2018	LPE	O33-C31-C32-N
10	A	2010	PCW	O4P-C4-C5-N
10	A	2019	PCW	O4P-C4-C5-N
8	A	2006	Y01	CAR-CBC-OAW-CAY
8	A	2020	Y01	CAO-CBB-CBE-CAP
10	A	2010	PCW	C37-C38-C39-C40
7	A	2011	P5S	O19-C1-C2-O37
10	A	2010	PCW	C14-C15-C16-C17
9	A	2008	LPE	C31-O33-P-O31

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Mol	Chain	Res	Type	Atoms
9	A	2014	LPE	C31-O33-P-O31
10	A	2010	PCW	C4-O4P-P-O2P
10	A	2019	PCW	C33-C34-C35-C36
9	A	2015	LPE	C2-C3-O3-P
8	A	2006	Y01	CAO-CBB-CBE-CAP
7	A	2011	P5S	C27-C28-C29-C30
9	A	2016	LPE	O1-C1-C2-O2H
9	A	2009	LPE	O1-C11-C12-C13
10	A	2012	PCW	C15-C16-C17-C18
8	A	2006	Y01	CAV-CBC-OAW-CAY
10	A	2012	PCW	C19-C20-C21-C22
9	A	2008	LPE	O1-C1-C2-O2H
7	A	2004	P5S	C52-C53-C54-C55
10	A	2012	PCW	C17-C18-C19-C20
5	B	303	NAG	C1-C2-N2-C7
10	A	2010	PCW	O31-C31-C32-C33
9	A	2009	LPE	C2-C1-O1-C11
9	A	2016	LPE	C2-C1-O1-C11
10	A	2010	PCW	C19-C20-C21-C22
10	A	2019	PCW	C19-C20-C21-C22
10	A	2019	PCW	C32-C33-C34-C35
10	A	2010	PCW	C1-C2-C3-O3
8	A	2005	Y01	CAM-CAY-OAW-CBC
10	A	2010	PCW	C34-C35-C36-C37
8	A	2005	Y01	OAG-CAY-OAW-CBC
7	A	2011	P5S	O37-C38-C39-C40
10	A	2012	PCW	O3-C11-C12-C13
8	A	2007	Y01	CAN-CAJ-CAO-CBB
8	A	2020	Y01	CAJ-CAN-CBA-CAB
9	A	2014	LPE	C2-C1-O1-C11
7	A	2011	P5S	O47-C38-C39-C40
7	A	2011	P5S	C21-C22-C23-C24
10	A	2012	PCW	O11-C11-C12-C13
7	A	2011	P5S	C31-C32-C33-C34
8	A	2007	Y01	CAL-CAM-CAY-OAW

There are no ring outliers.

9 monomers are involved in 12 short contacts:

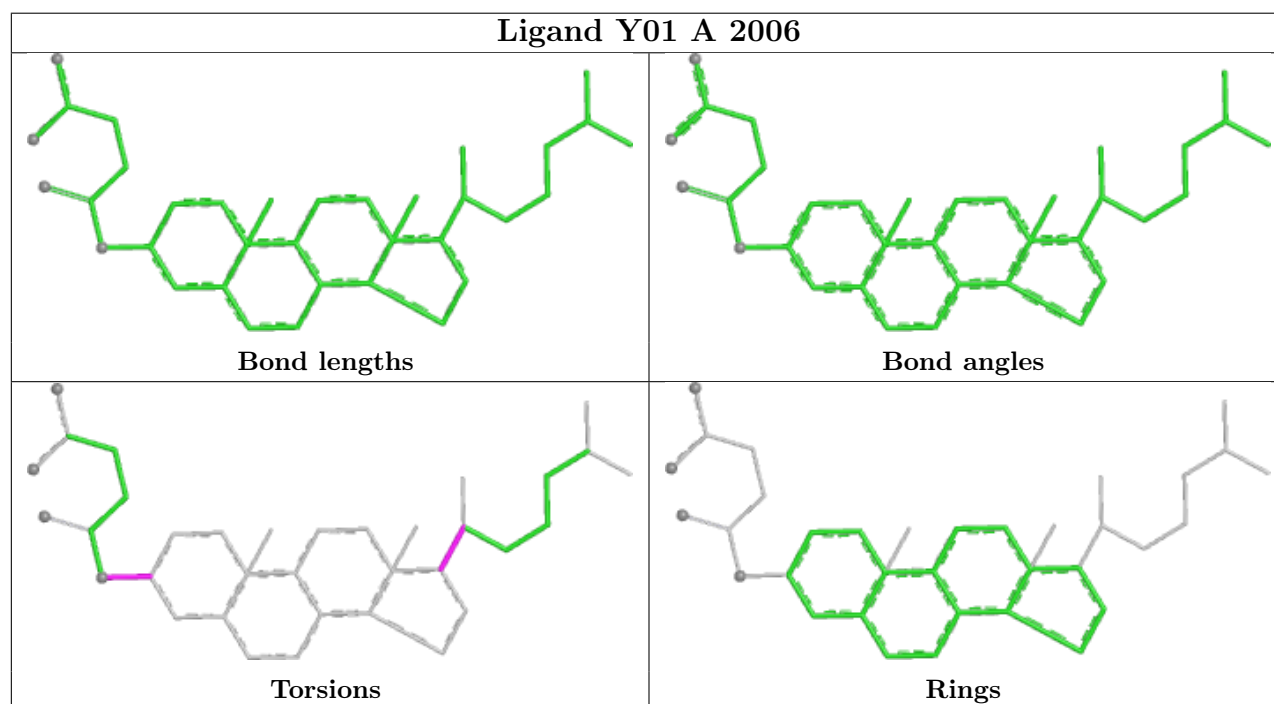
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	2005	Y01	1	0
8	A	2020	Y01	1	0

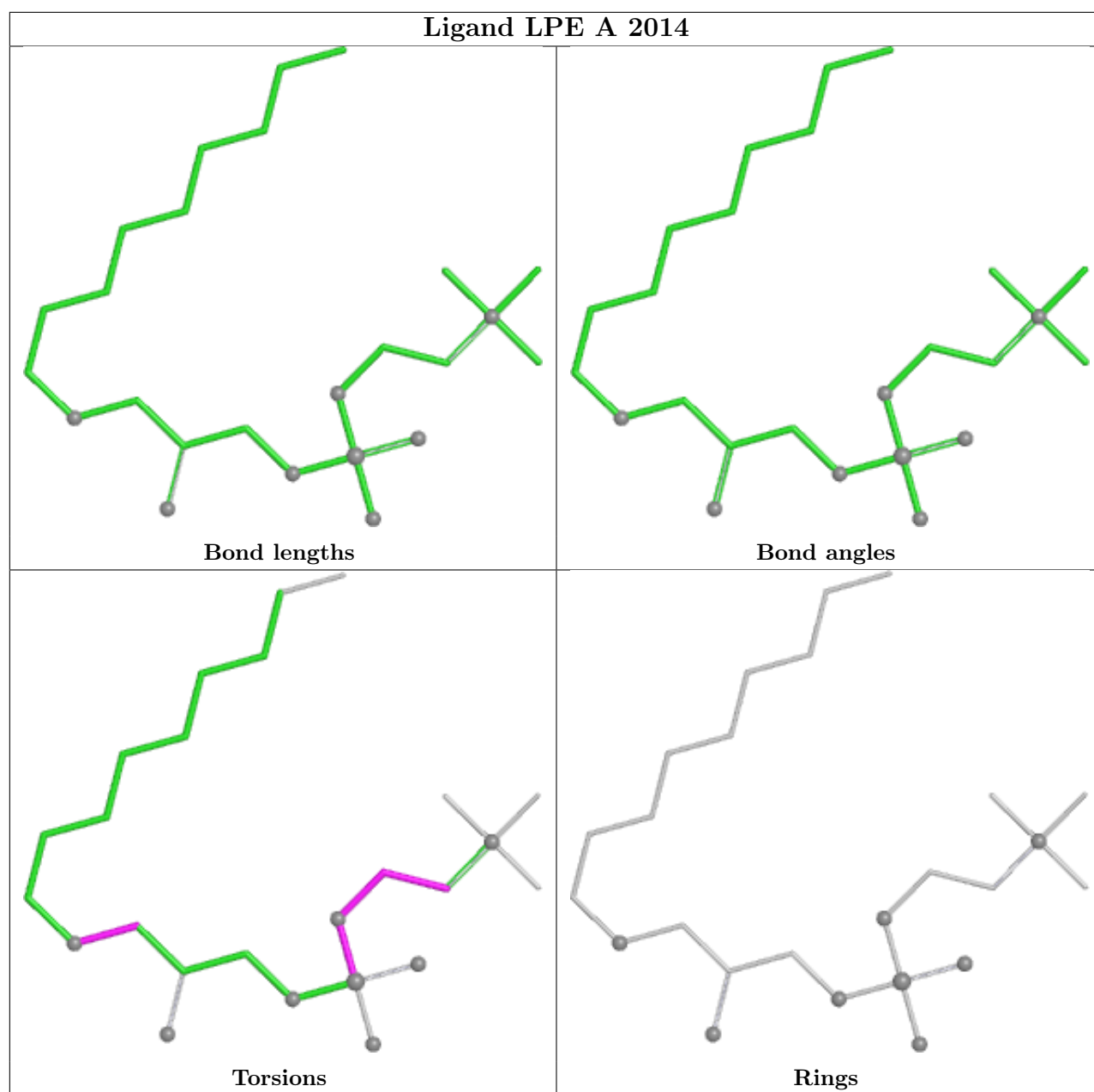
Continued on next page...

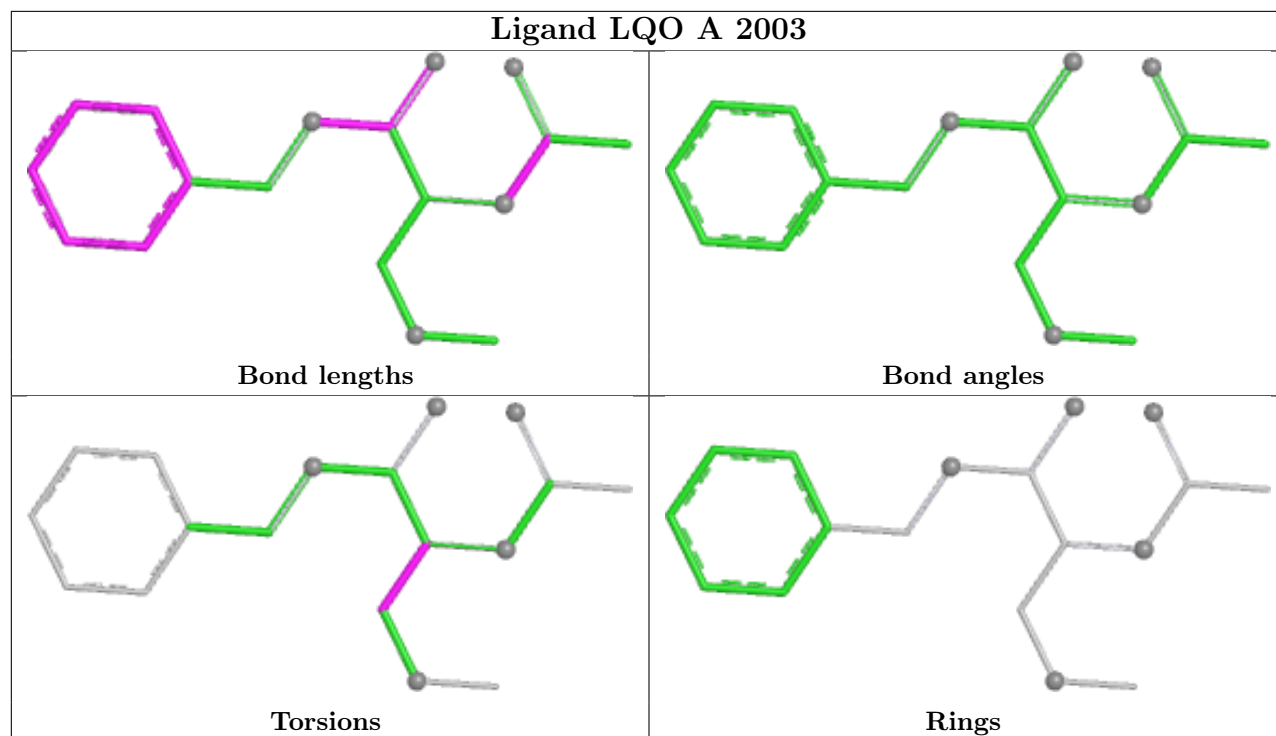
Continued from previous page...

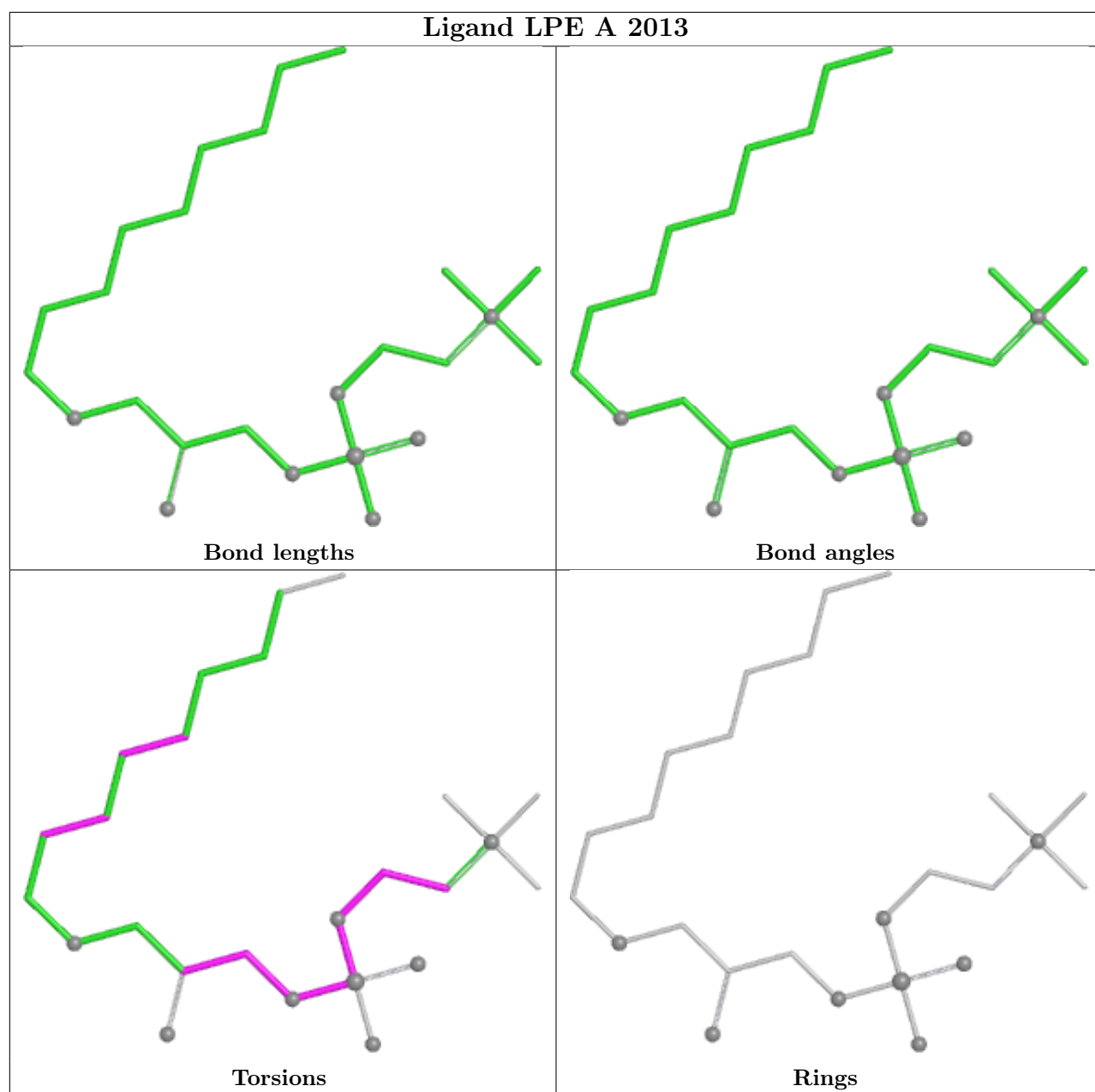
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	303	NAG	1	0
8	A	2007	Y01	3	0
10	A	2012	PCW	2	0
7	A	2004	P5S	1	0
5	B	301	NAG	1	0
10	A	2019	PCW	1	0
10	A	2010	PCW	1	0

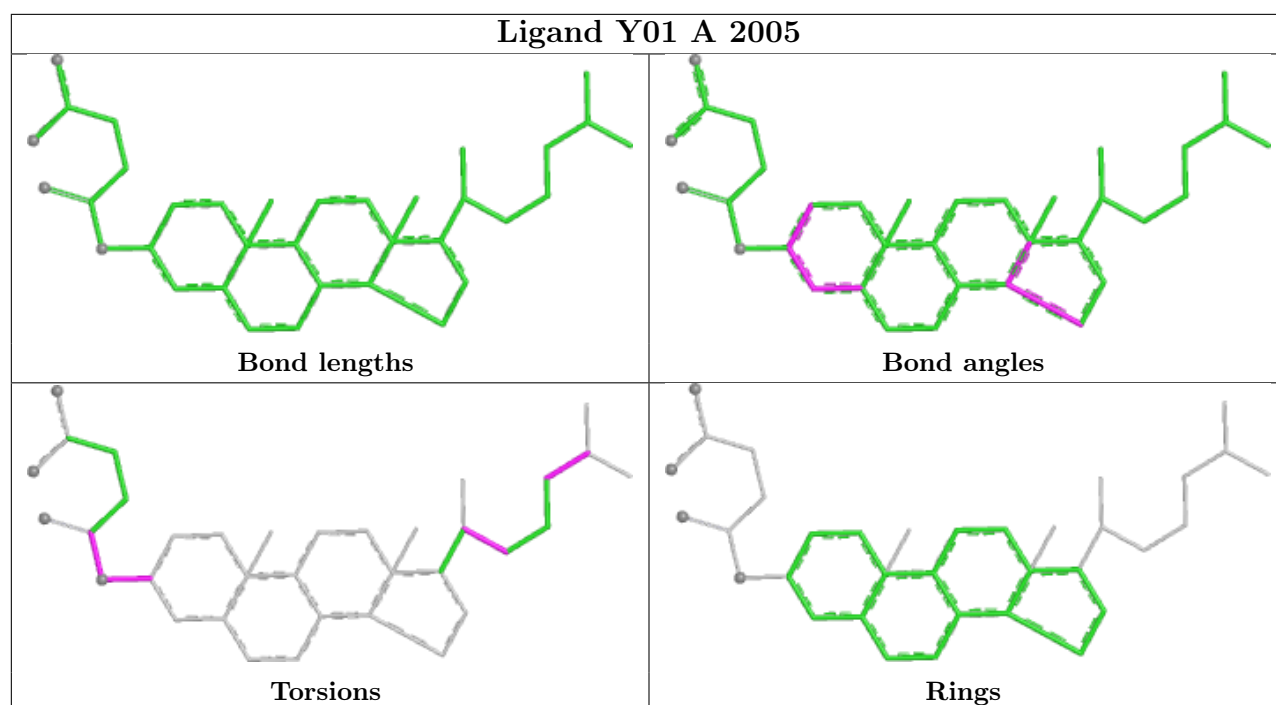
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

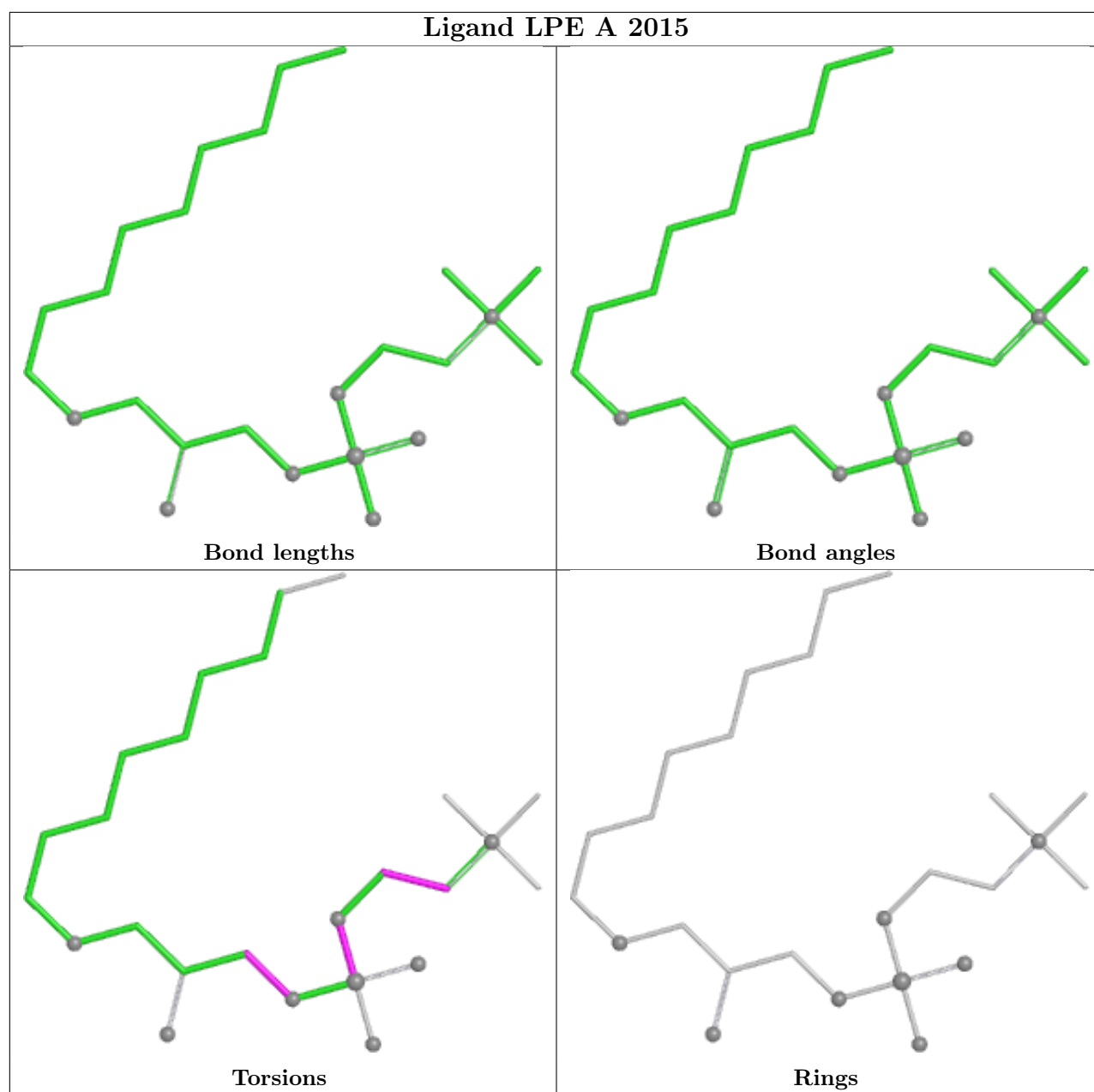


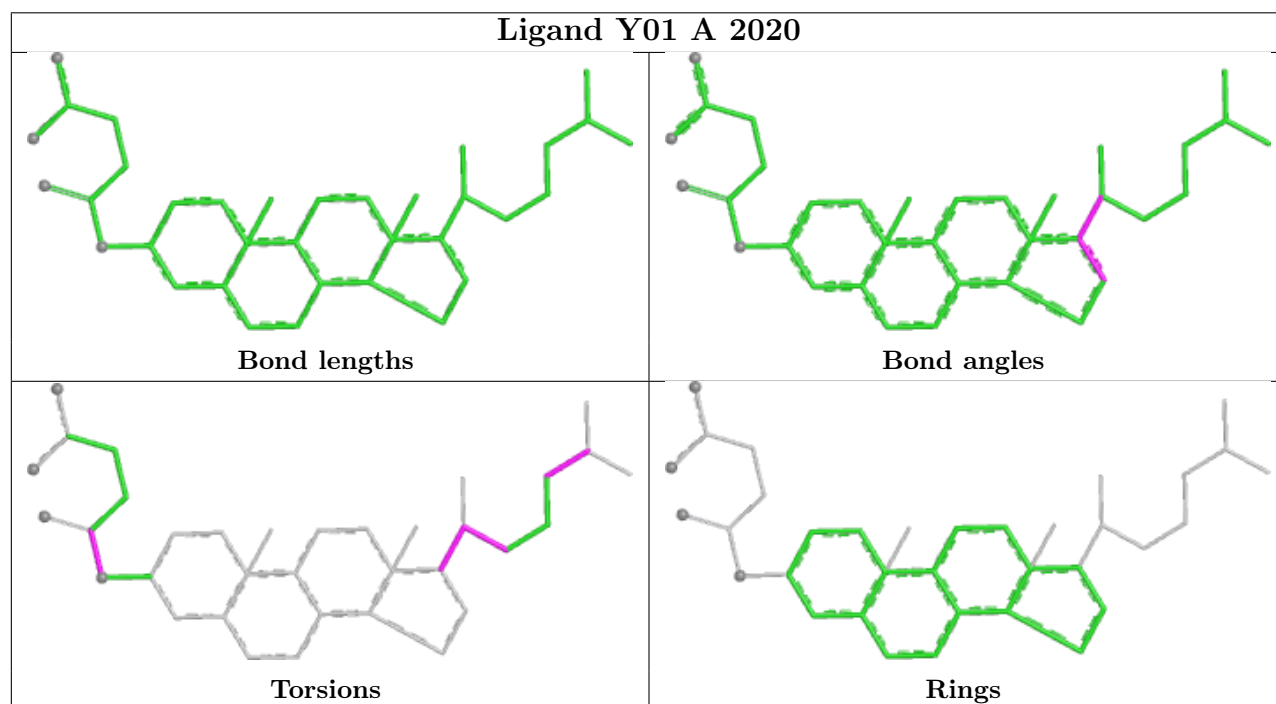
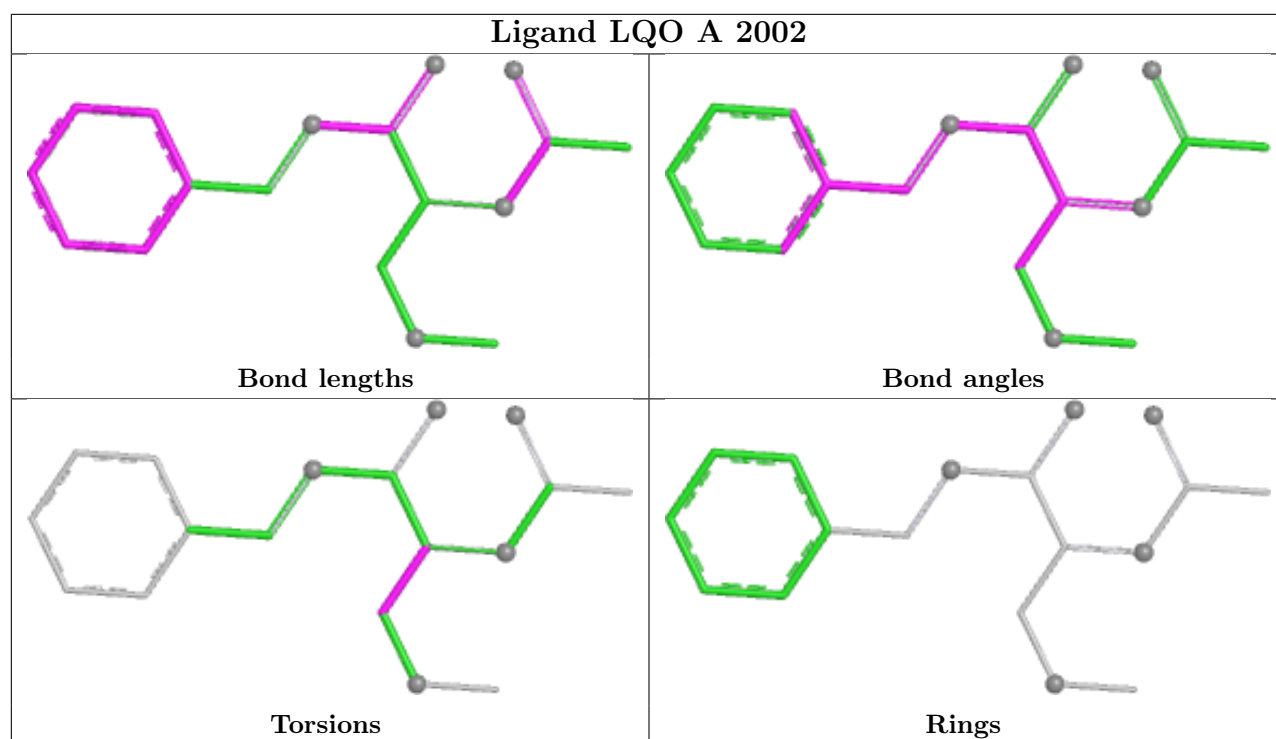


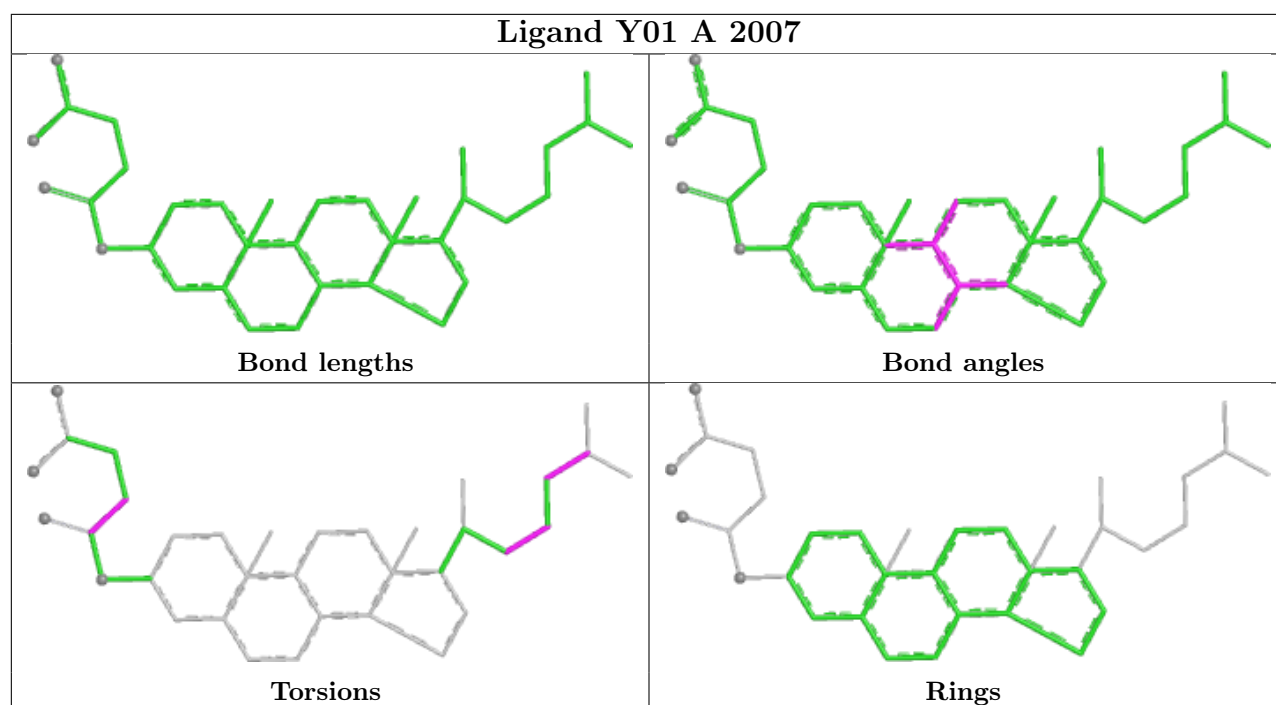


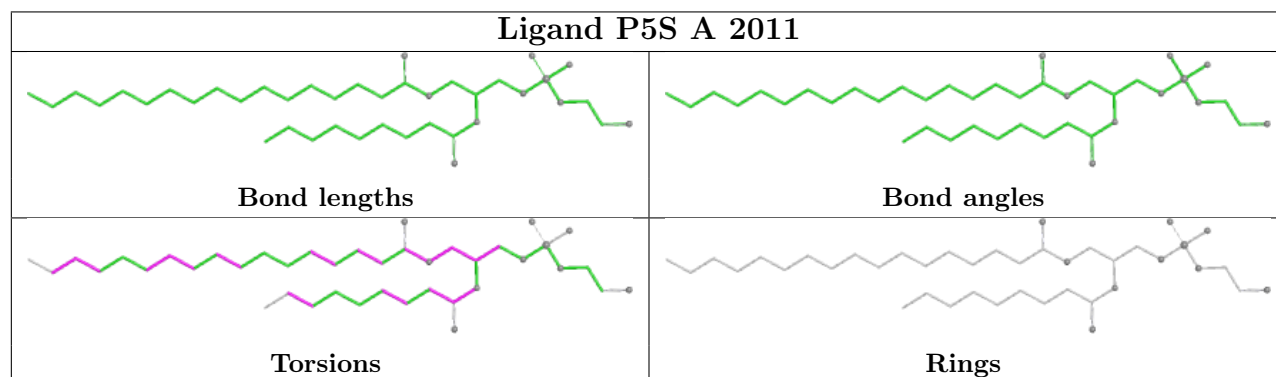
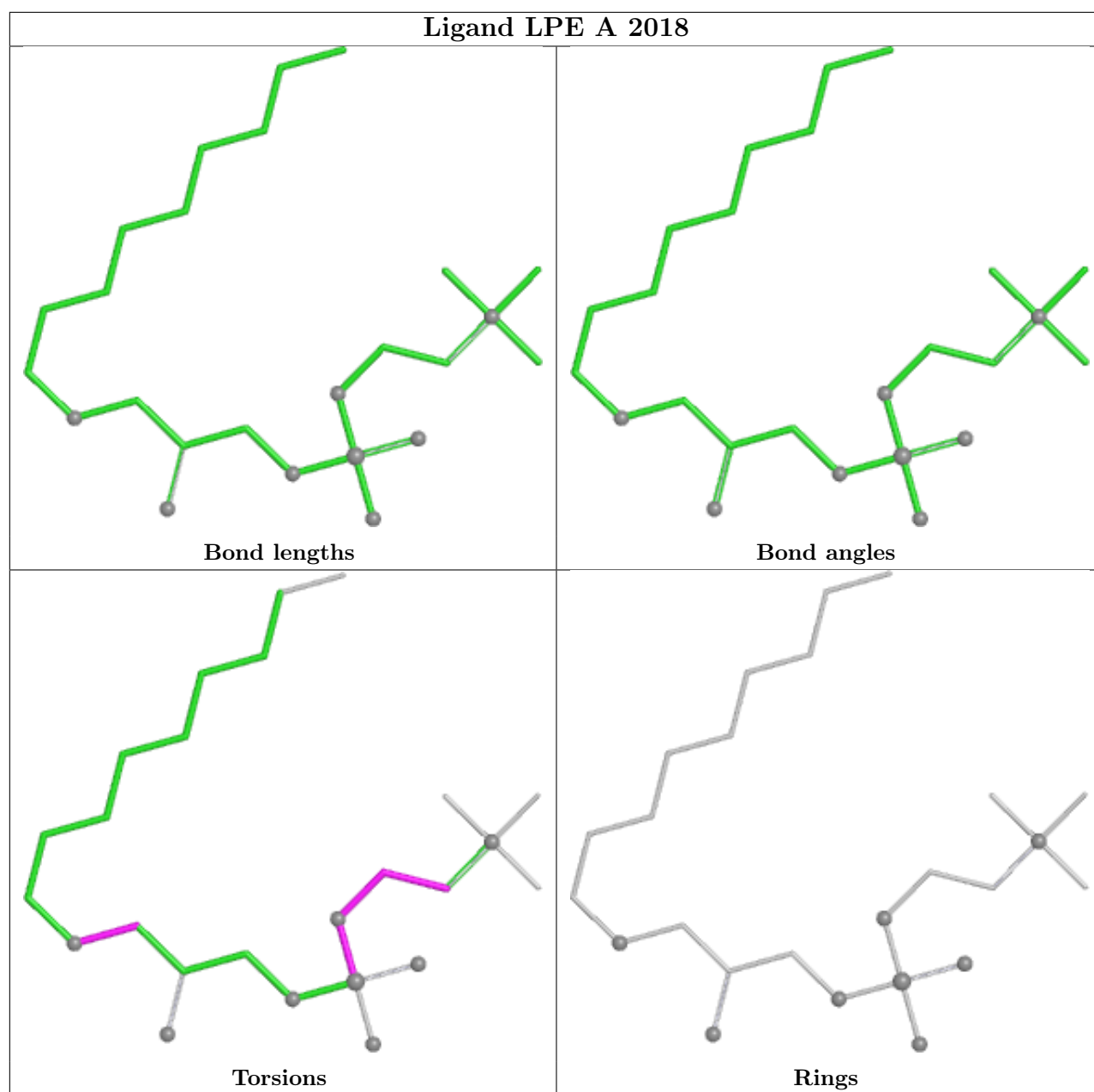


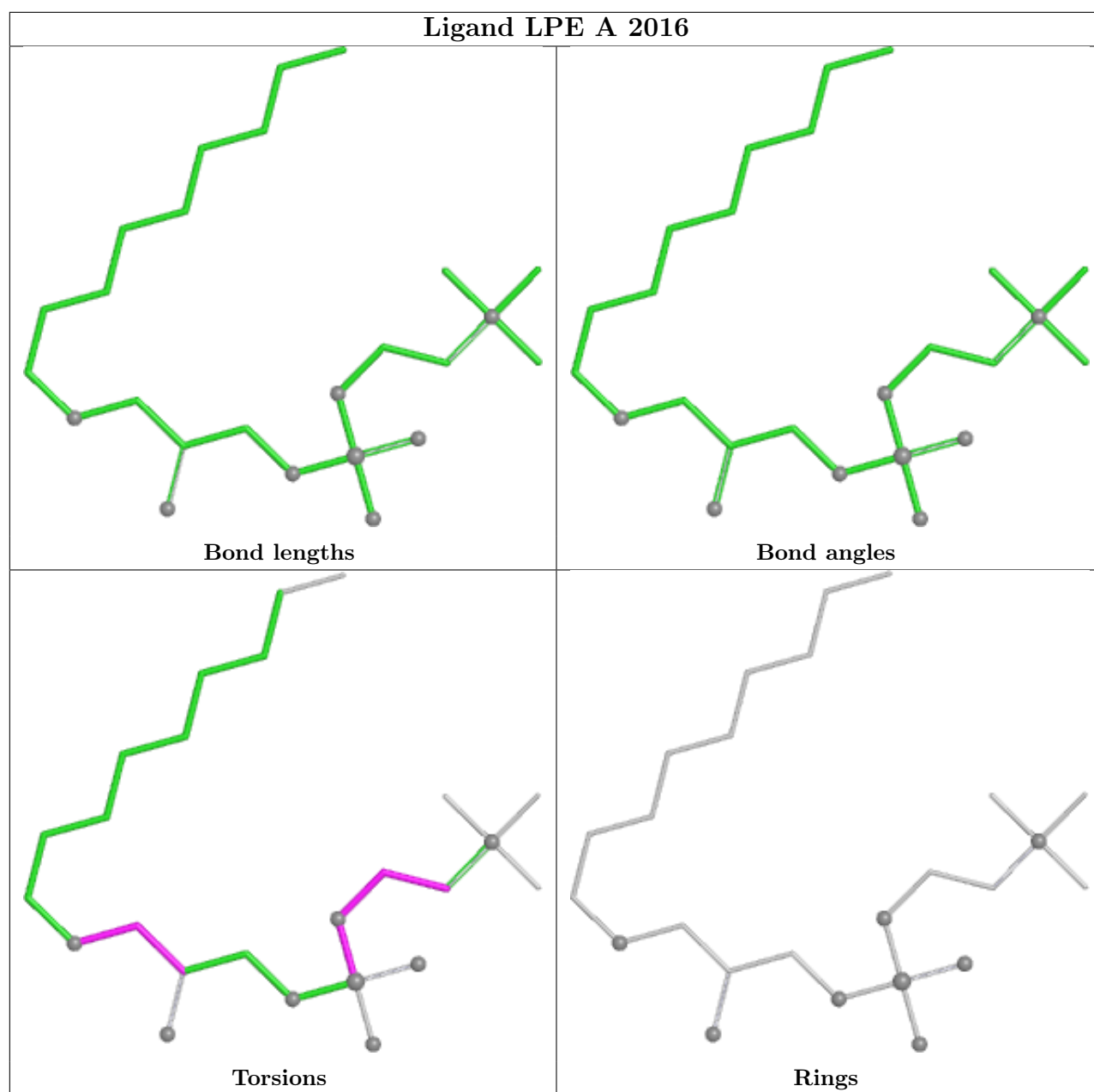


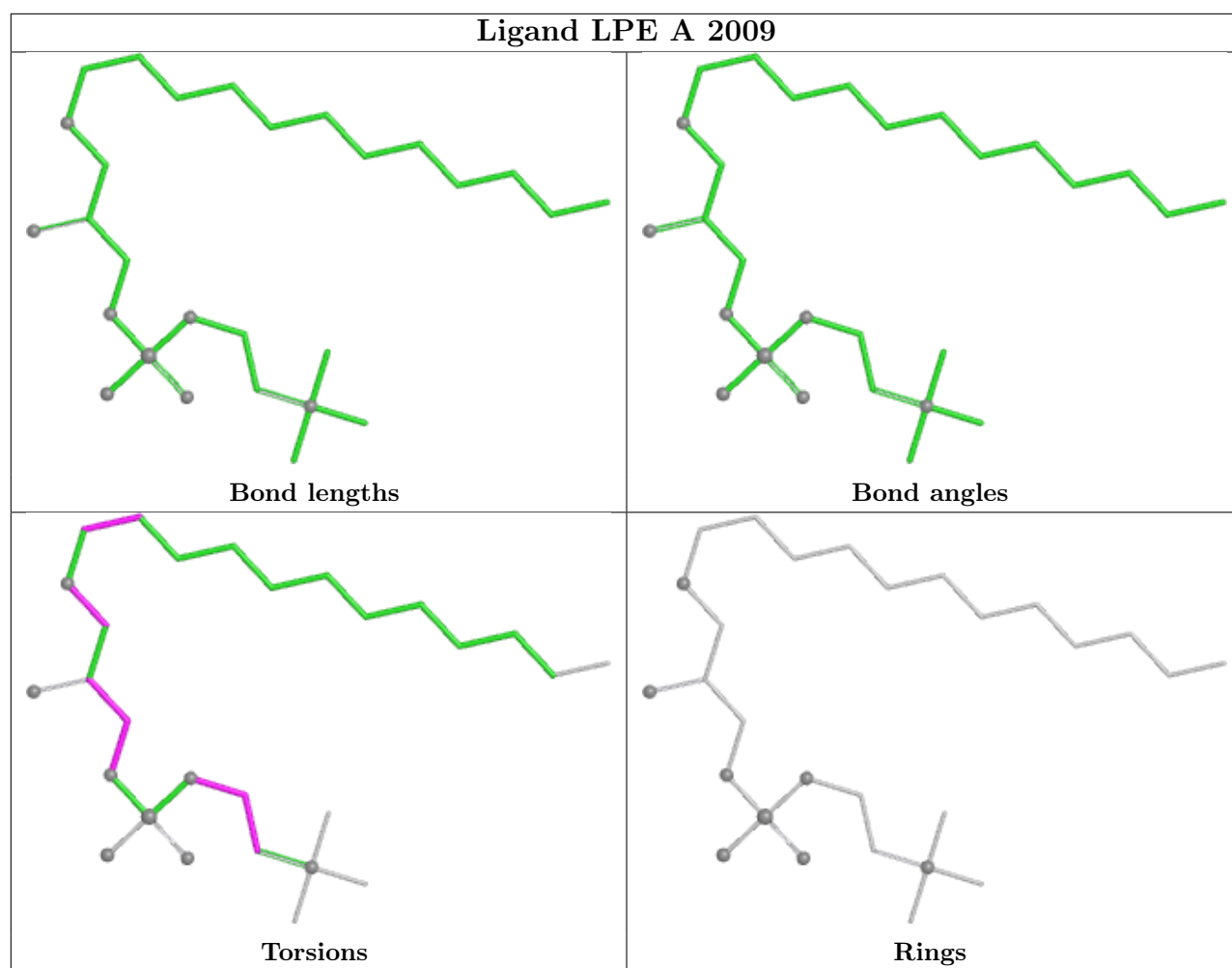


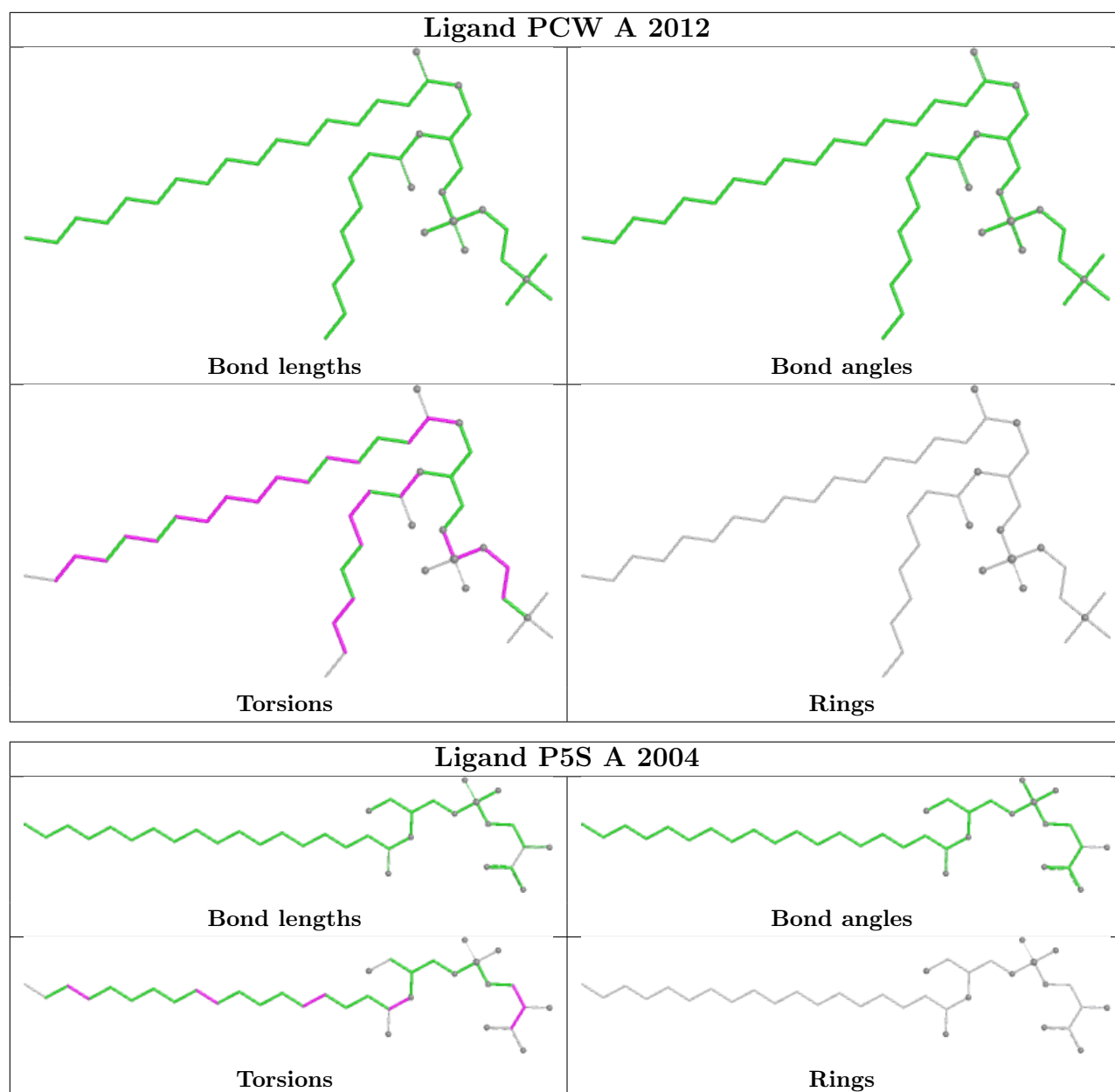


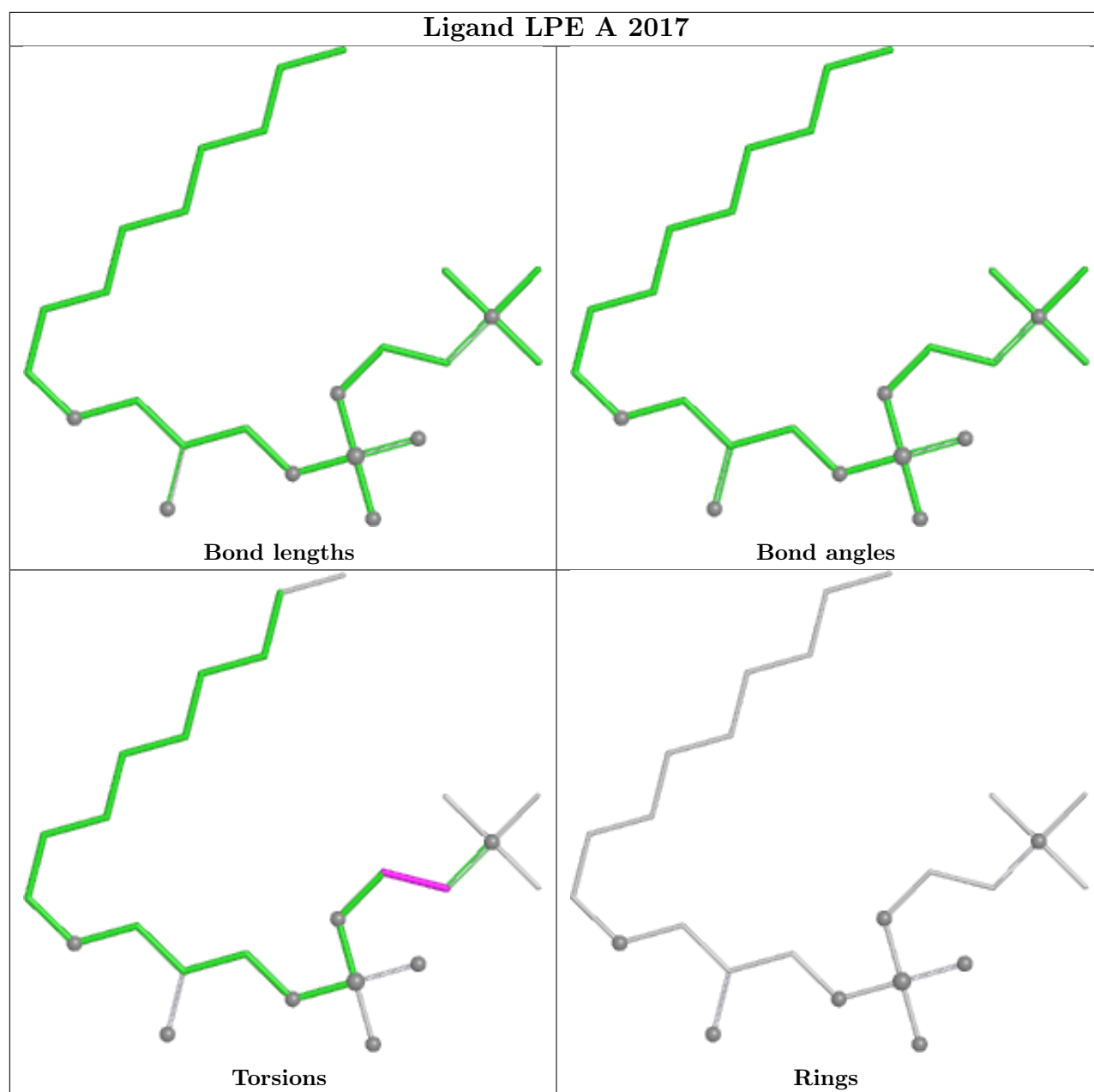


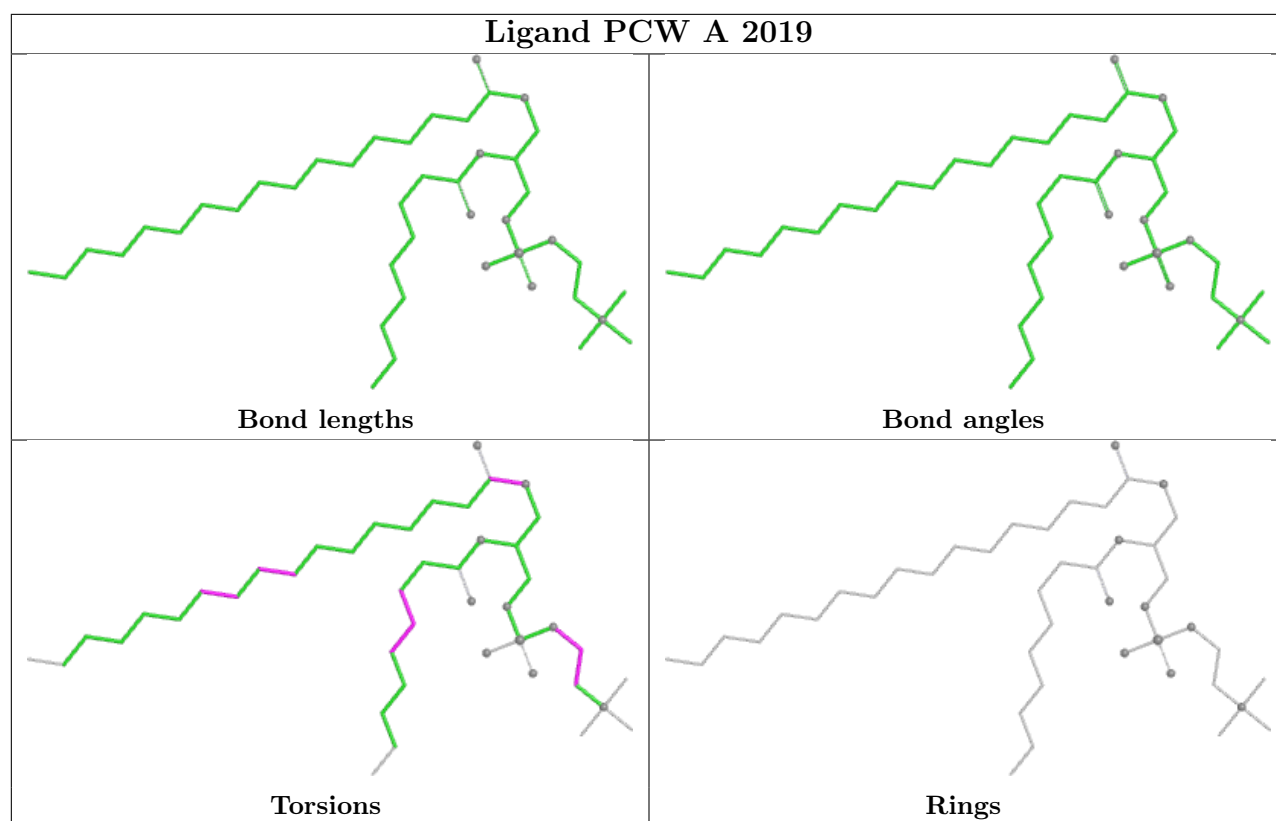


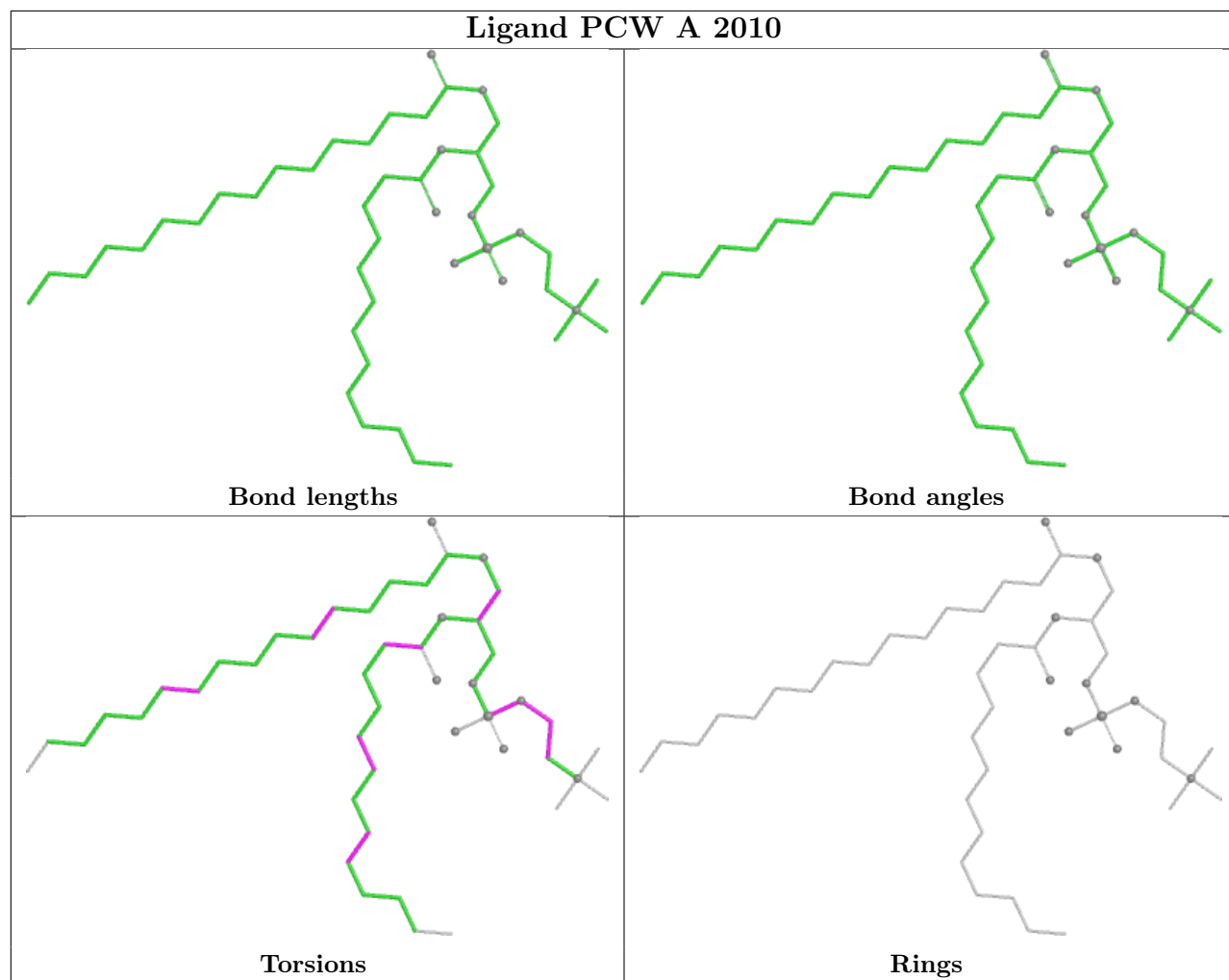


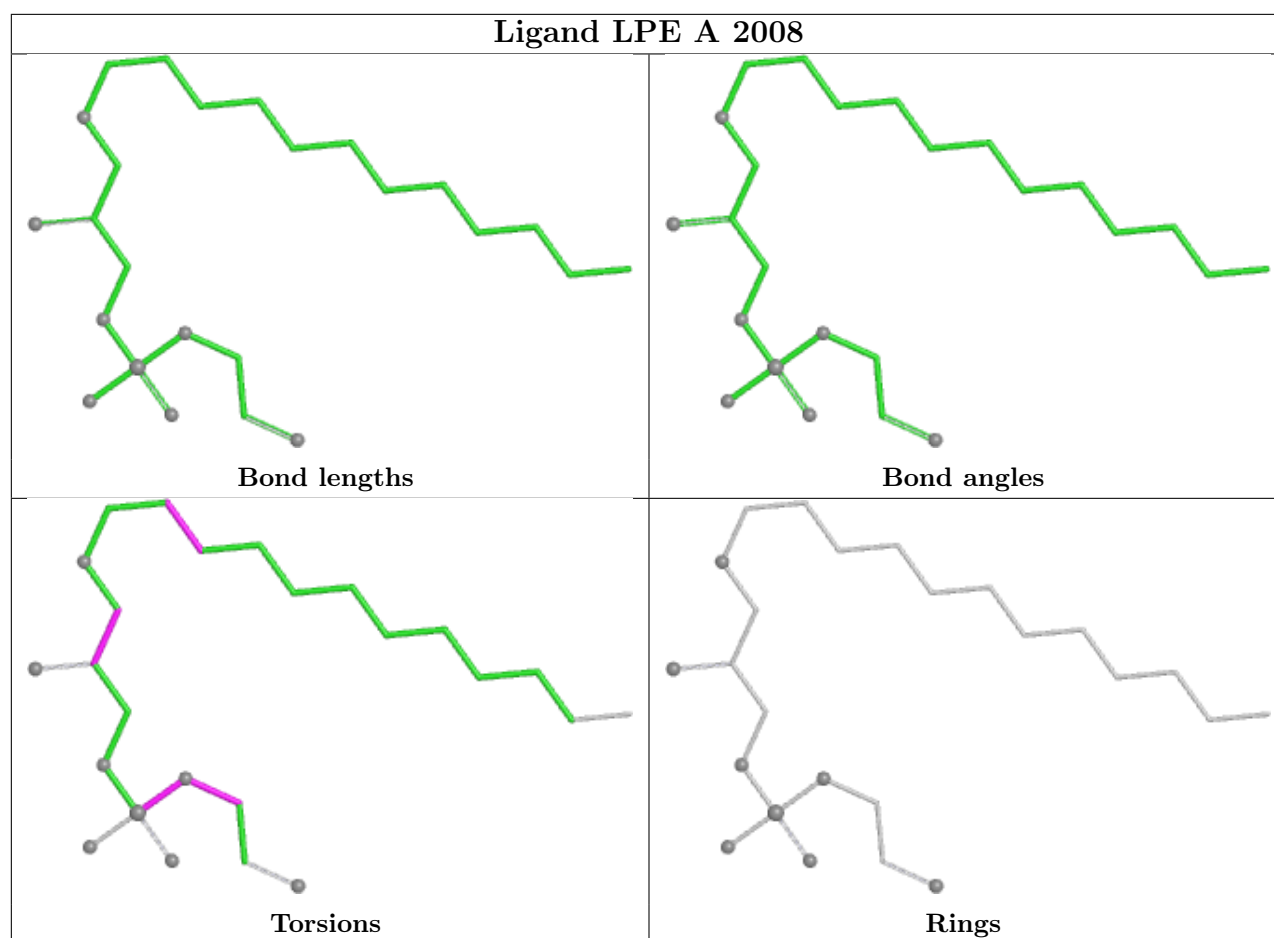












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

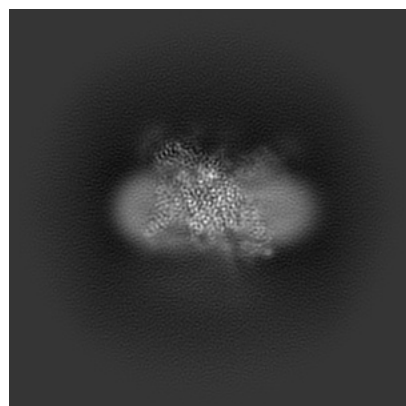
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40238. These allow visual inspection of the internal detail of the map and identification of artifacts.

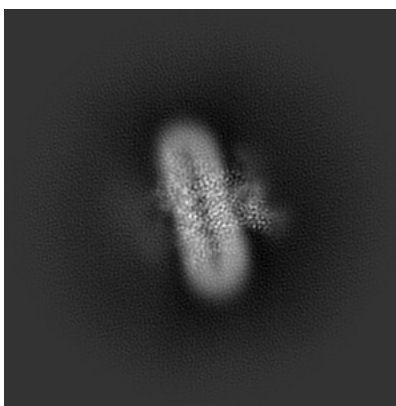
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

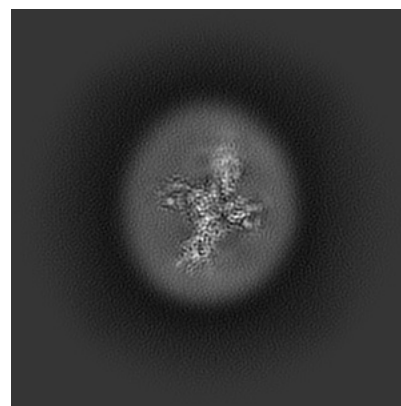
6.1.1 Primary map



X

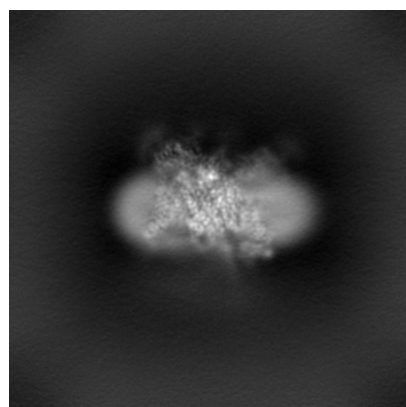


Y

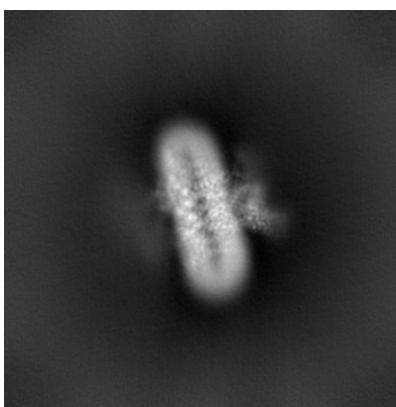


Z

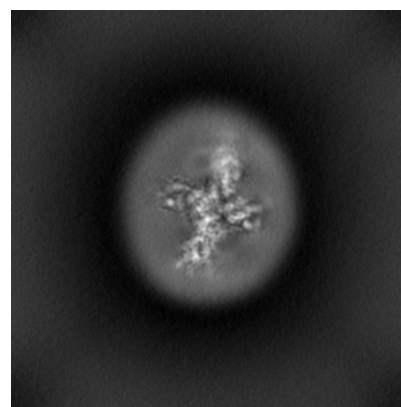
6.1.2 Raw map



X



Y

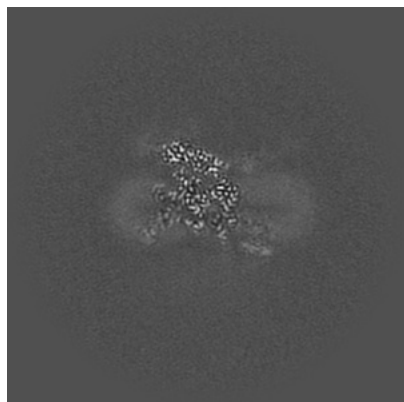


Z

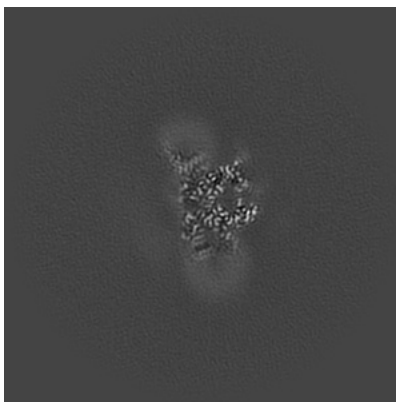
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

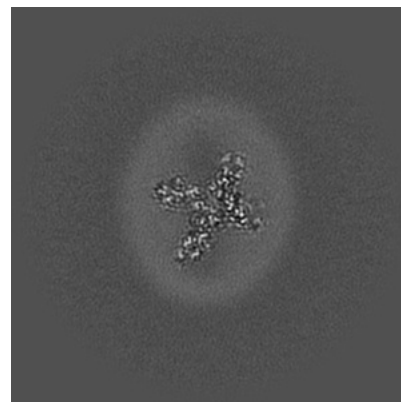
6.2.1 Primary map



X Index: 160

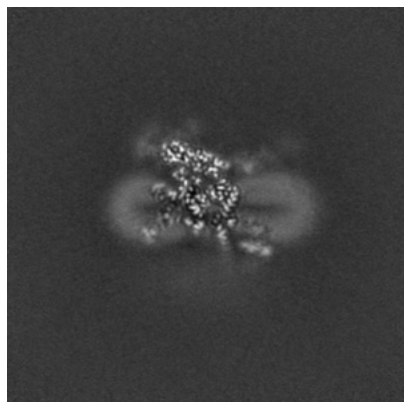


Y Index: 160

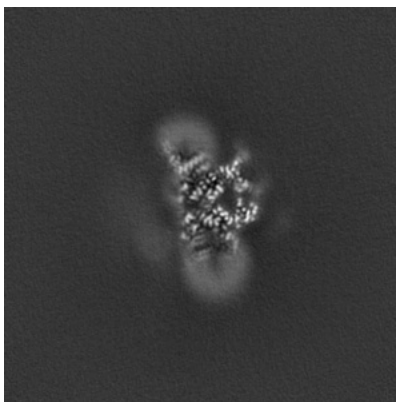


Z Index: 160

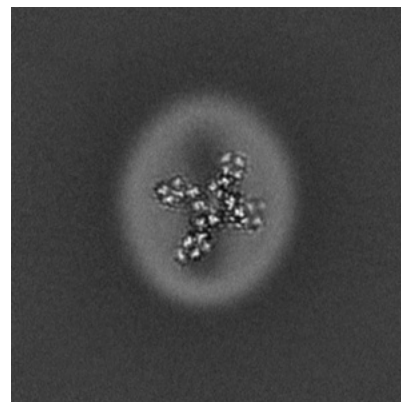
6.2.2 Raw map



X Index: 160



Y Index: 160

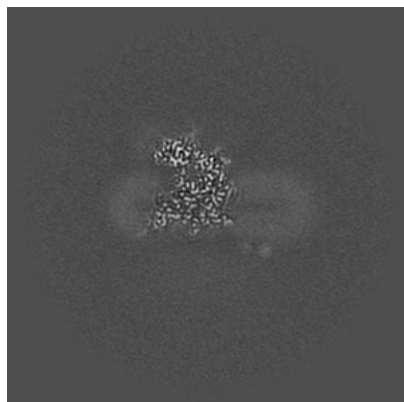


Z Index: 160

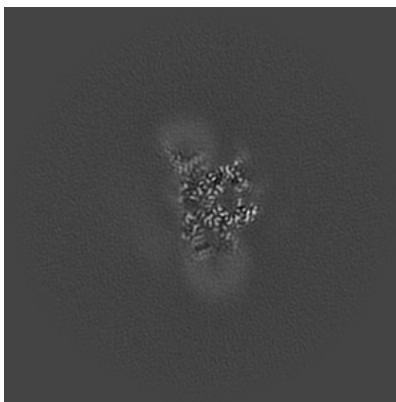
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

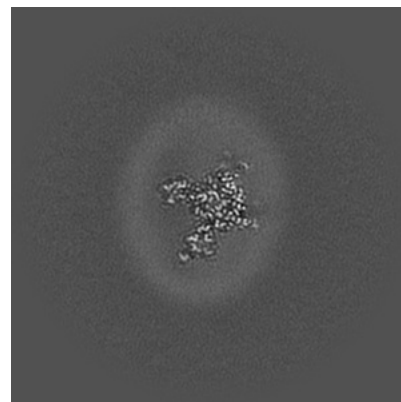
6.3.1 Primary map



X Index: 153

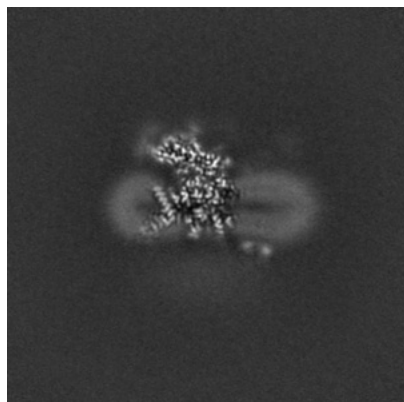


Y Index: 160

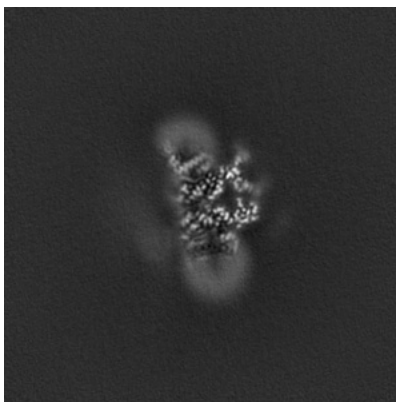


Z Index: 167

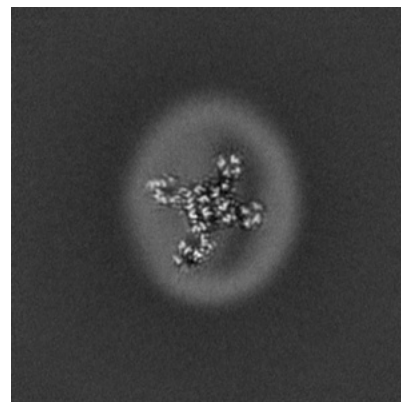
6.3.2 Raw map



X Index: 156



Y Index: 161

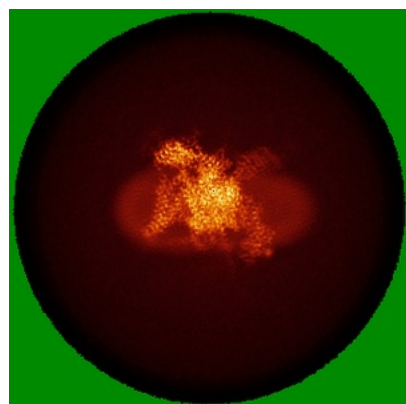


Z Index: 152

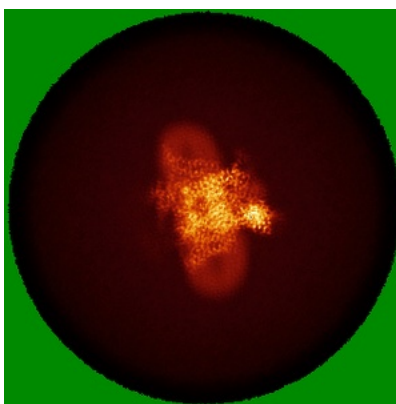
The images above show the largest variance slices of the map in three orthogonal directions.

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

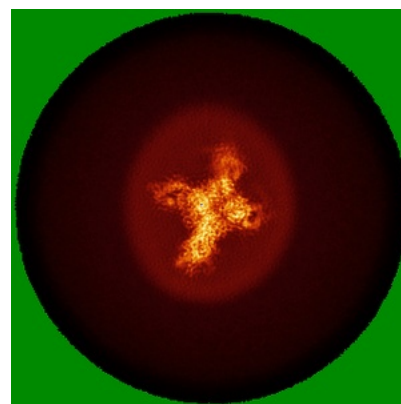
6.4.1 Primary map



X

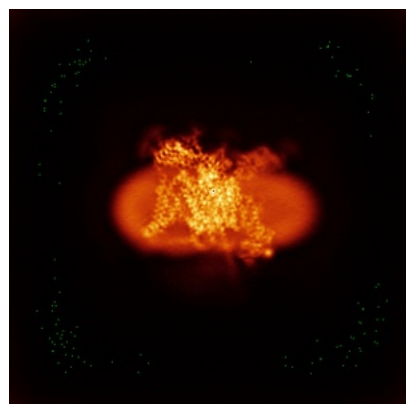


Y

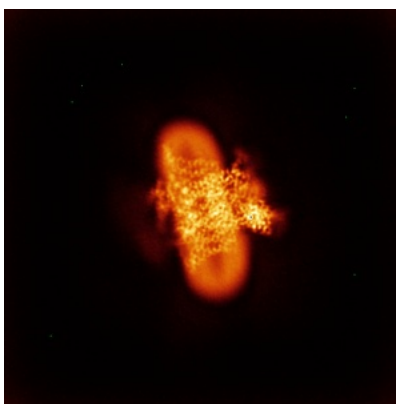


Z

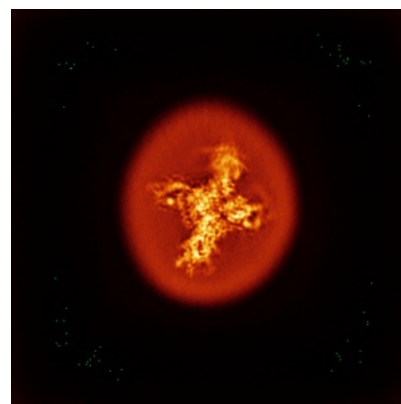
6.4.2 Raw map



X



Y

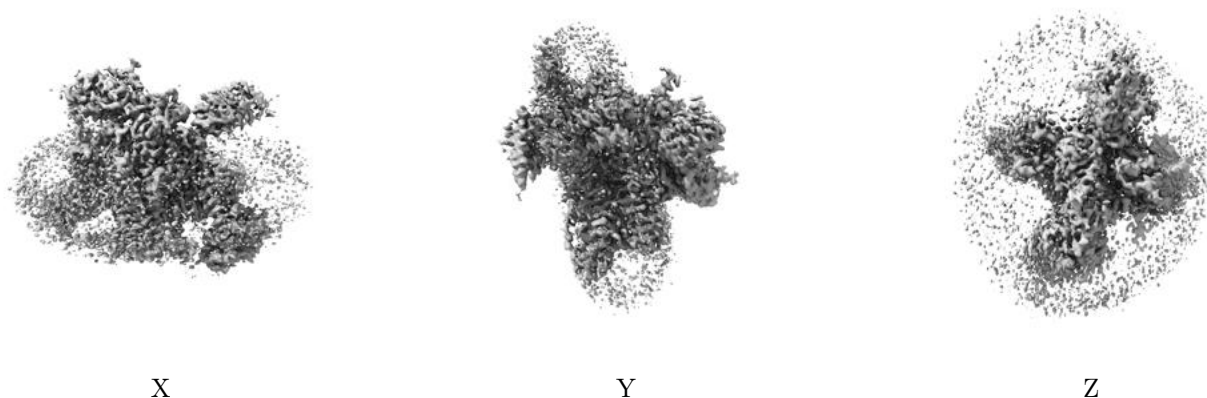


Z

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 0.7. 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.

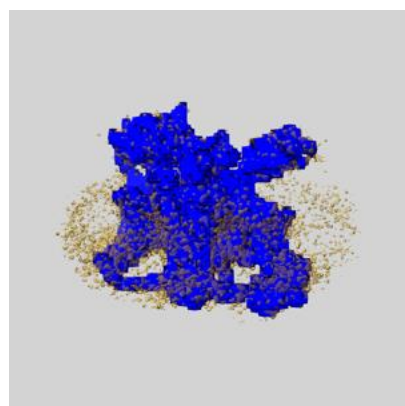
6.6 Mask visualisation [i](#)

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

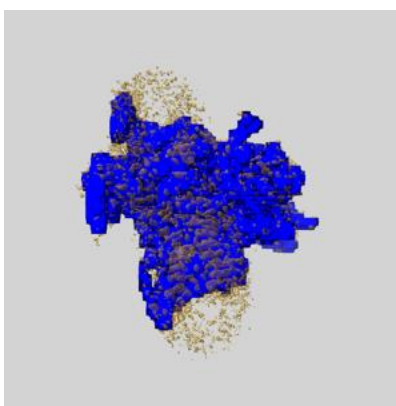
A mask typically either:

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

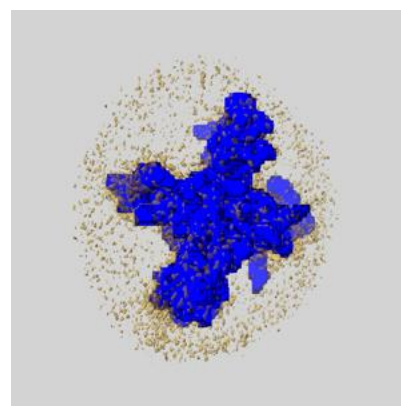
6.6.1 emd_40238_msk_1.map [i](#)



X



Y

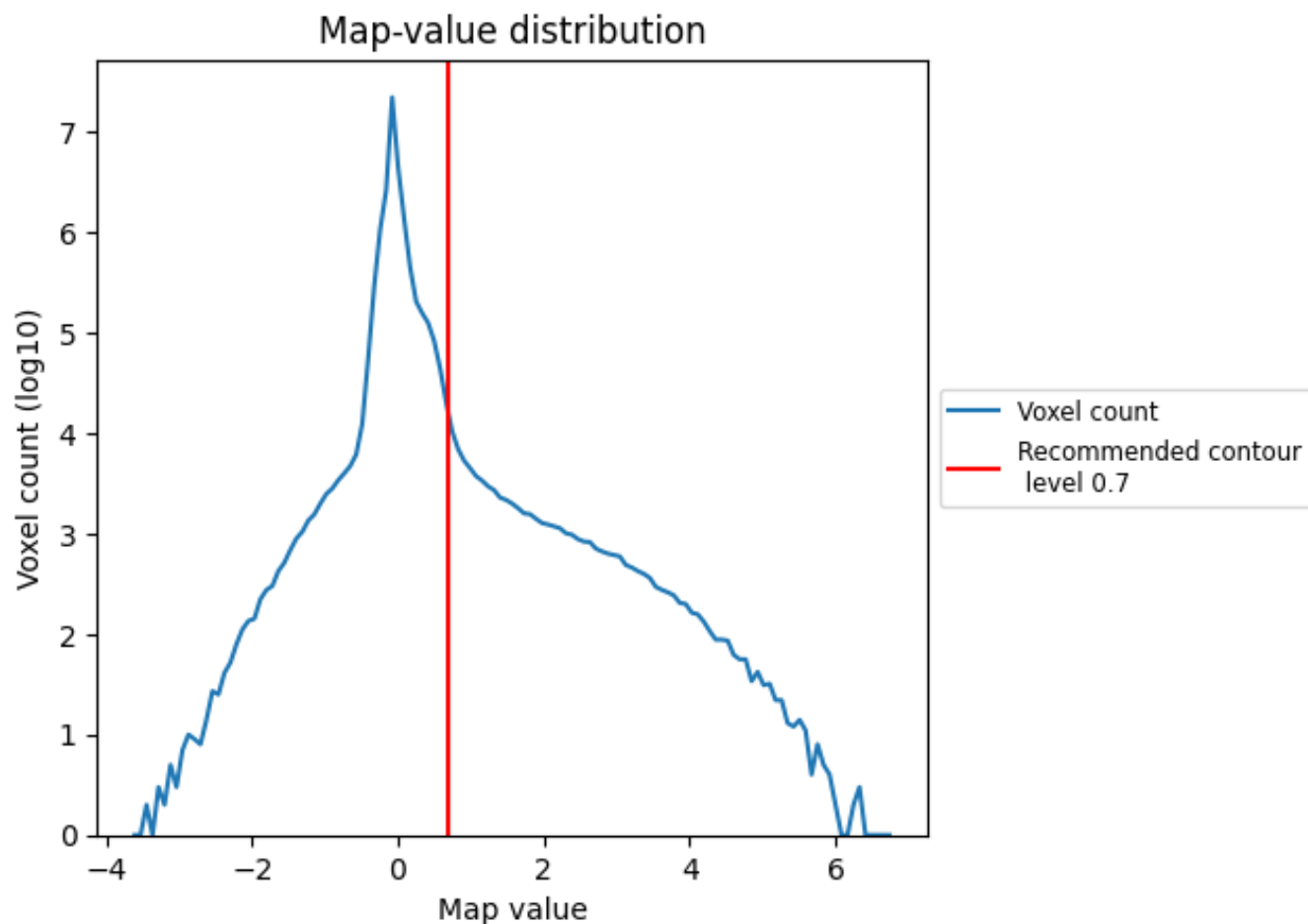


Z

7 Map analysis [i](#)

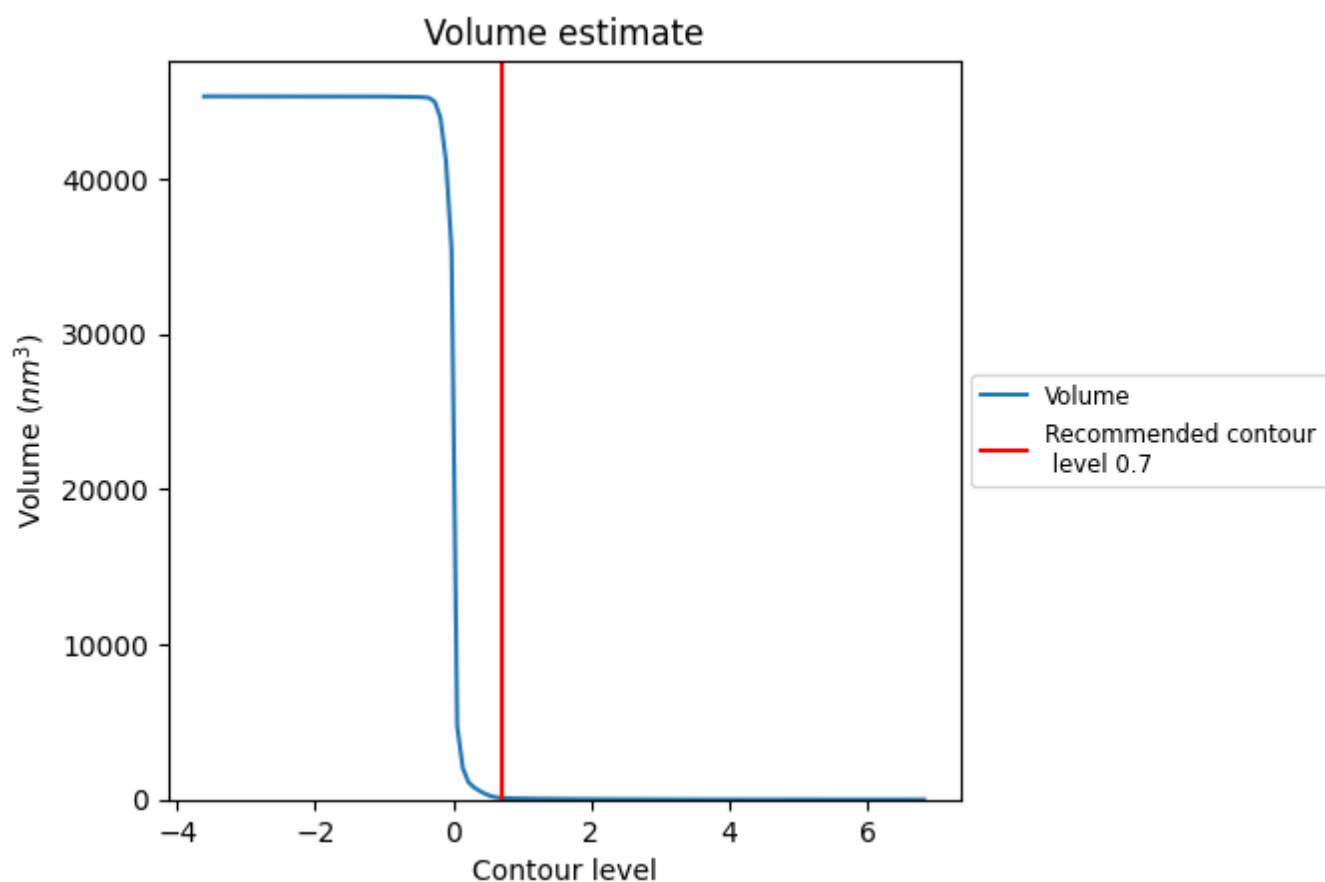
This section contains the results of statistical analysis of the map.

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.

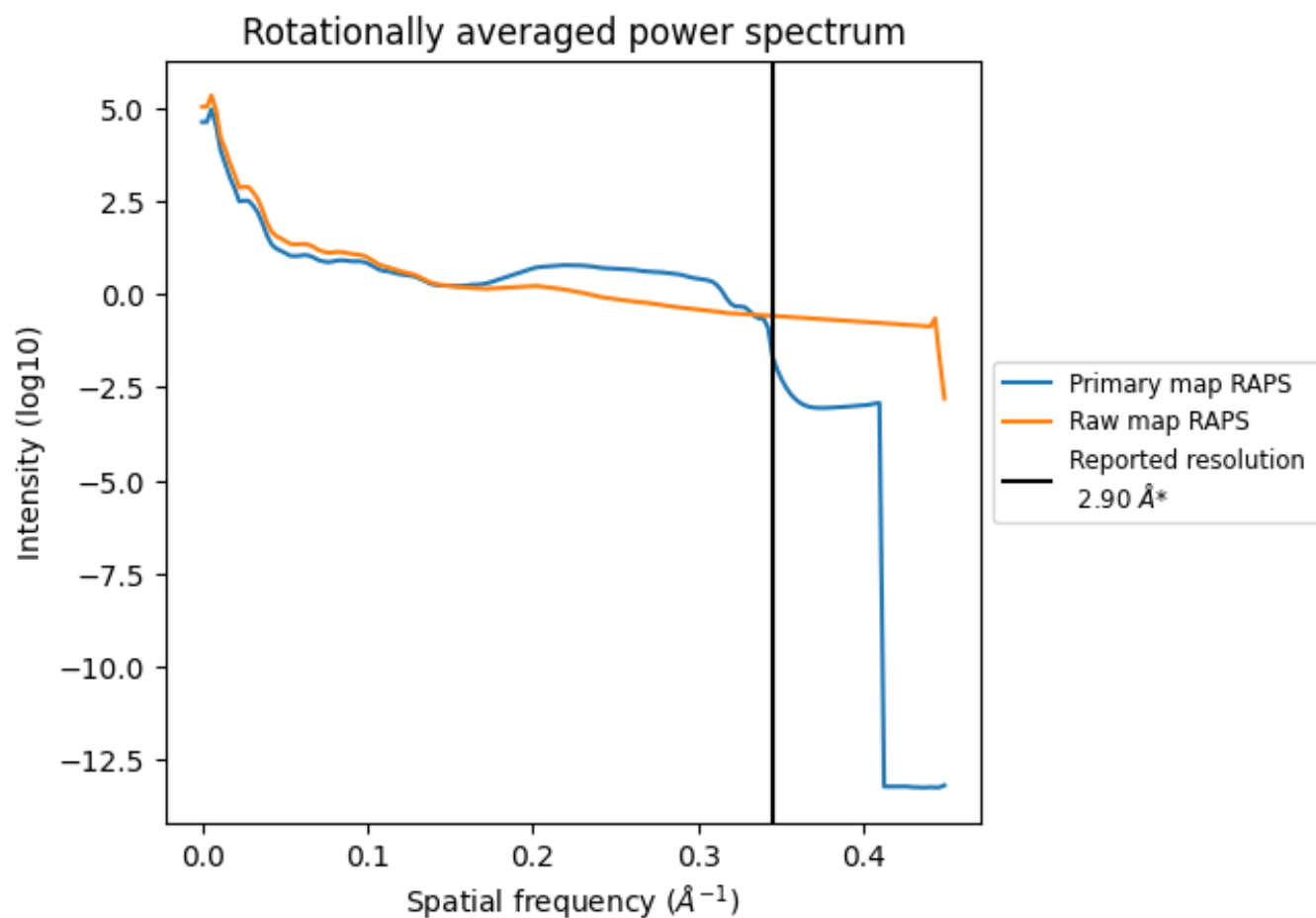
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 111 nm³; this corresponds to an approximate mass of 100 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 ⓘ

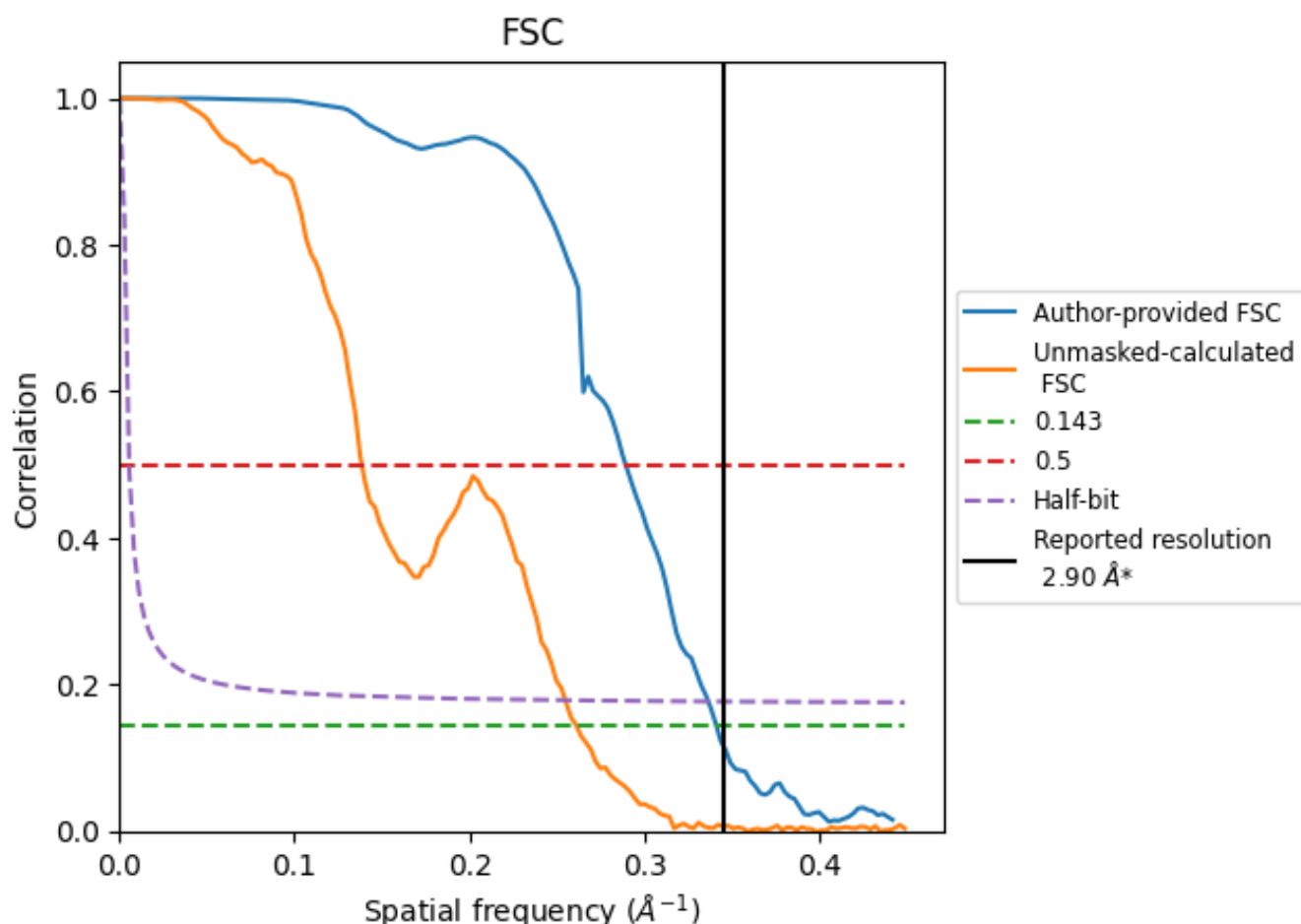


*Reported resolution corresponds to spatial frequency of 0.345 \AA^{-1}

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.345 \AA^{-1}

8.2 Resolution estimates [i](#)

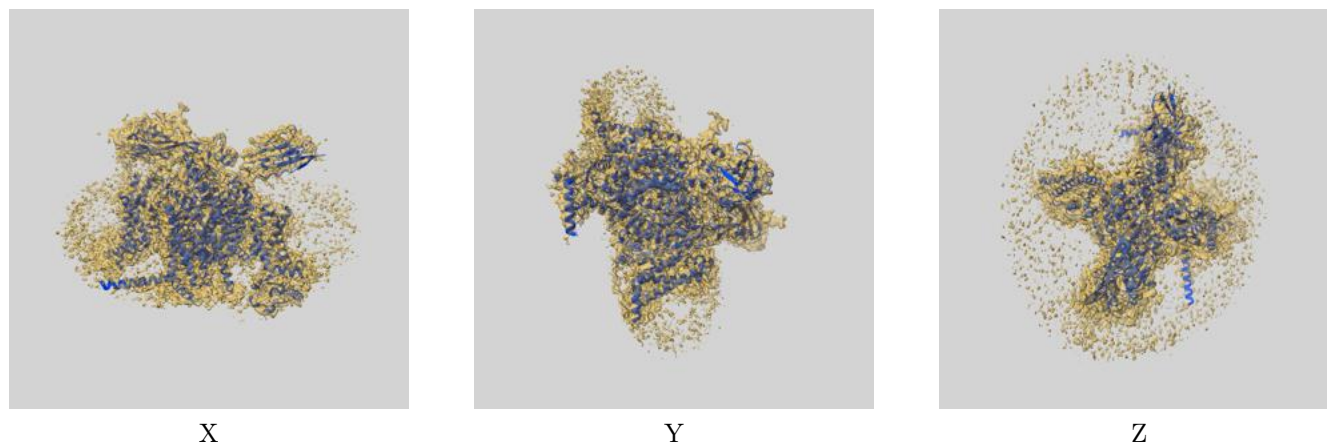
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.90	-	-
Author-provided FSC curve	2.93	3.46	2.97
Unmasked-calculated*	3.83	7.23	3.92

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.83 differs from the reported value 2.9 by more than 10 %

9 Map-model fit [i](#)

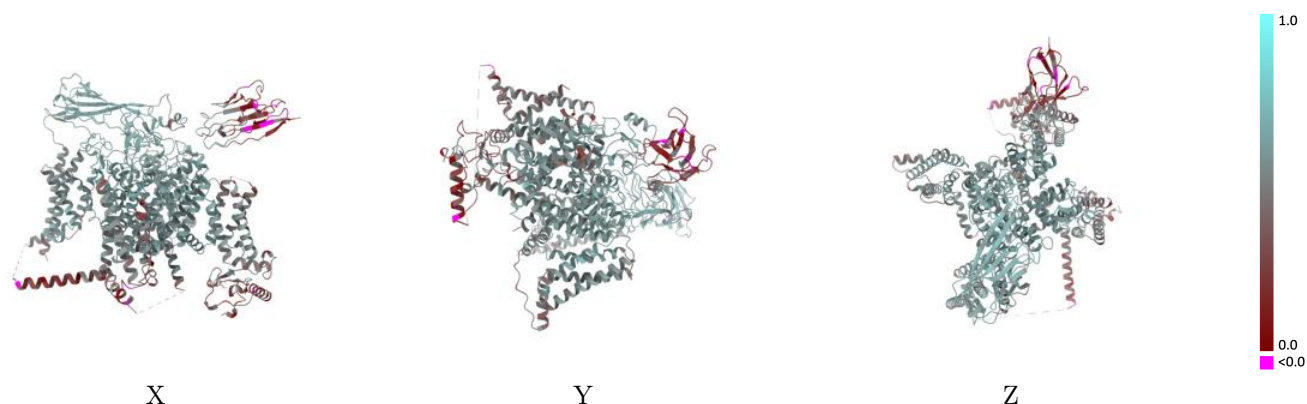
This section contains information regarding the fit between EMDB map EMD-40238 and PDB model 8S9B. Per-residue inclusion information can be found in [section 3](#) on [page 9](#).

9.1 Map-model overlay [i](#)



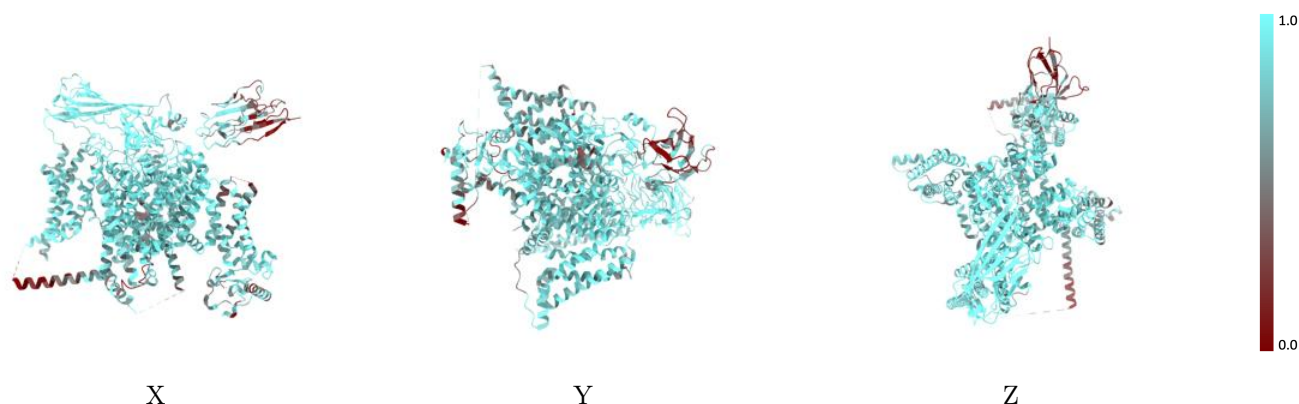
The images above show the 3D surface view of the map at the recommended contour level 0.7 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



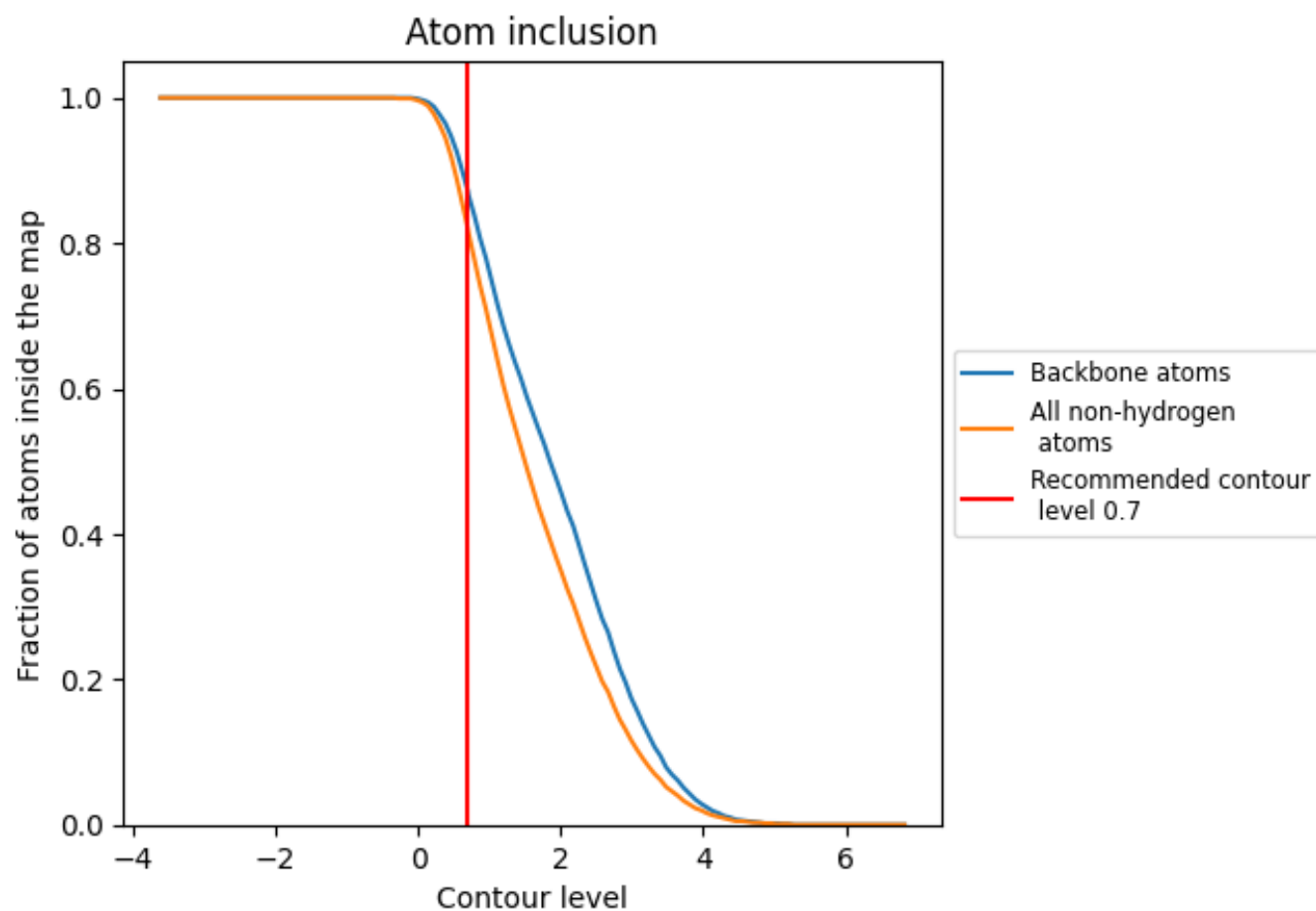
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.7).

9.4 Atom inclusion [i](#)



At the recommended contour level, 87% of all backbone atoms, 82% 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 (0.7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div>0.8200</div>	<div><div></div>0.5130</div>
A	<div><div></div>0.8340</div>	<div><div></div>0.5240</div>
B	<div><div></div>0.9190</div>	<div><div></div>0.5730</div>
C	<div><div></div>0.5140</div>	<div><div></div>0.3090</div>
D	<div><div></div>0.7500</div>	<div><div></div>0.4410</div>
E	<div><div></div>0.8570</div>	<div><div></div>0.4920</div>
F	<div><div></div>0.9290</div>	<div><div></div>0.5570</div>

1.0

0.0

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