



Full wwPDB EM Validation Report ⓘ

Oct 13, 2025 – 10:16 AM JST

PDB ID : 8Z4O / pdb_00008z4o
EMDB ID : EMD-39768
Title : MERS-CoV post-fusion S ectodomain trimer
Authors : Wang, X.; Wang, Z.
Deposited on : 2024-04-17
Resolution : 3.79 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.dev129
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
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) : 2.46

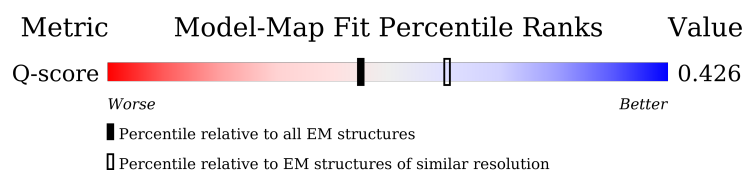
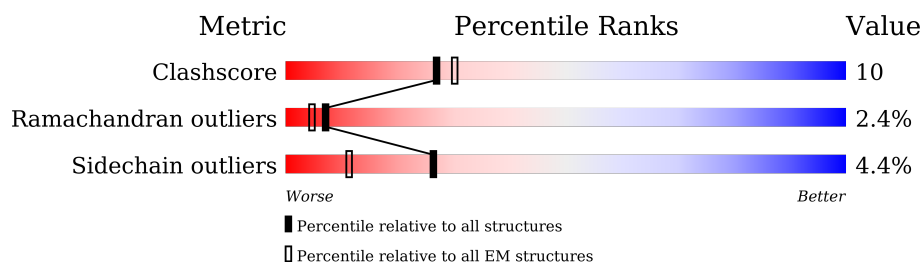
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.79 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	210492	15764	-
Ramachandran outliers	207382	16835	-
Sidechain outliers	206894	16415	-
Q-score	-	25397	10018 (3.29 - 4.29)

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	539	
1	B	539	
1	C	539	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 7515 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.

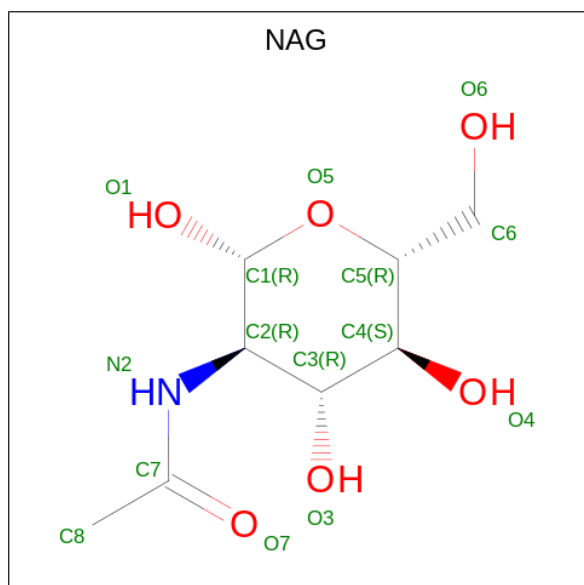
- Molecule 1 is a protein called Spike protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	358	Total	C	N	O	S	0	0
			2439	1536	405	487	11		
1	B	358	Total	C	N	O	S	0	0
			2437	1529	408	489	11		
1	C	357	Total	C	N	O	S	0	0
			2415	1516	404	484	11		

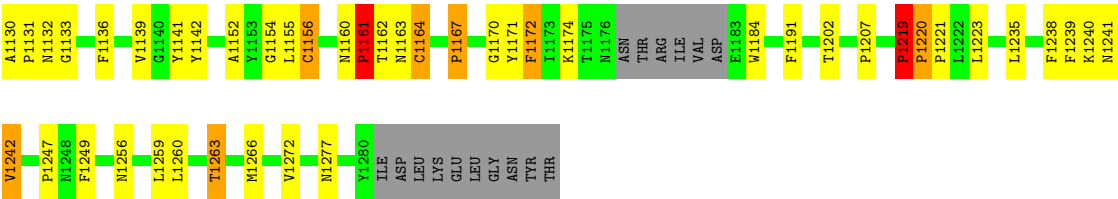
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1020	GLN	HIS	conflict	UNP K9N5Q8
B	1020	GLN	HIS	conflict	UNP K9N5Q8
C	1020	GLN	HIS	conflict	UNP K9N5Q8

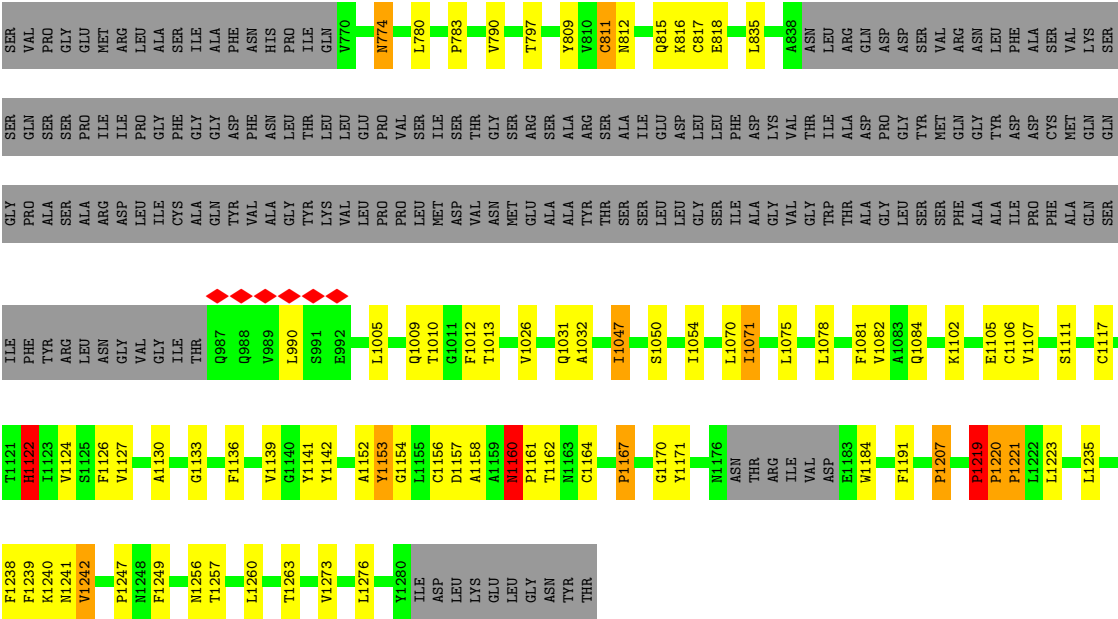
- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



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



● Molecule 1: Spike protein S2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	36570	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TECNAI F30	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.743	Depositor
Minimum map value	-1.494	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.043	Depositor
Recommended contour level	0.2	Depositor
Map size (\AA)	365.228, 365.228, 365.228	wwPDB
Map dimensions	340, 340, 340	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.0742, 1.0742, 1.0742	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	1.04	1/2465 (0.0%)	1.71	27/3322 (0.8%)
1	B	1.02	2/2463 (0.1%)	1.70	20/3318 (0.6%)
1	C	1.02	1/2441 (0.0%)	1.72	29/3291 (0.9%)
All	All	1.03	4/7369 (0.1%)	1.71	76/9931 (0.8%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1167	PRO	N-CD	-17.30	1.23	1.47
1	C	1167	PRO	N-CD	-14.59	1.27	1.47
1	B	1167	PRO	N-CD	-13.20	1.29	1.47
1	B	1161	PRO	N-CD	-7.05	1.37	1.47

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1167	PRO	CB-CA-C	-16.14	91.30	111.64
1	A	1167	PRO	CB-CA-C	-15.40	92.23	111.64
1	C	1167	PRO	CB-CA-C	-14.24	93.09	111.46
1	A	1167	PRO	N-CA-CB	-9.38	94.83	103.27
1	C	1161	PRO	N-CA-C	9.01	124.91	111.03
1	A	1167	PRO	CA-N-CD	8.78	124.29	112.00
1	C	1160	ASN	CB-CA-C	8.43	126.77	110.17
1	A	783	PRO	CB-CA-C	-8.39	101.07	111.64
1	C	1153	TYR	N-CA-C	-7.80	101.89	111.33
1	C	1142	TYR	CB-CA-C	-7.69	96.96	108.87
1	C	1167	PRO	N-CA-CB	-7.66	96.49	103.31
1	B	1167	PRO	N-CA-CB	-7.62	96.41	103.27
1	C	1263	THR	N-CA-C	-7.62	103.87	113.01
1	A	1263	THR	N-CA-C	-7.59	104.02	113.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1263	THR	N-CA-C	-7.30	104.25	113.01
1	B	1167	PRO	CA-N-CD	7.29	122.21	112.00
1	B	783	PRO	CB-CA-C	-7.22	102.15	111.46
1	C	1167	PRO	CA-N-CD	6.99	121.78	112.00
1	B	1142	TYR	CB-CA-C	-6.96	98.02	109.15
1	A	1136	PHE	CB-CA-C	-6.88	98.92	110.19
1	A	1153	TYR	N-CA-C	-6.83	103.06	111.33
1	A	1249	PHE	CB-CA-C	6.61	120.21	109.90
1	C	1249	PHE	CB-CA-C	6.42	119.92	109.90
1	C	1122	HIS	CB-CA-C	6.23	120.03	109.75
1	C	1136	PHE	CB-CA-C	-6.22	99.98	110.19
1	B	1241	ASN	N-CA-C	-6.17	104.56	111.28
1	B	1249	PHE	CB-CA-C	6.14	119.48	109.90
1	C	1071	ILE	N-CA-C	-6.09	104.40	110.62
1	C	1241	ASN	N-CA-C	-6.09	104.64	111.28
1	B	1172	PHE	CB-CA-C	-6.08	99.83	109.80
1	C	1220	PRO	N-CA-C	6.07	118.10	110.70
1	A	1226	SER	N-CA-C	6.06	117.96	111.36
1	A	1220	PRO	N-CA-C	6.05	118.08	110.70
1	B	815	GLN	N-CA-C	-6.00	105.81	113.01
1	A	1241	ASN	N-CA-C	-5.97	104.77	111.28
1	B	1220	PRO	N-CA-C	5.96	117.98	110.70
1	B	1031	GLN	N-CA-C	-5.94	104.81	111.28
1	A	1071	ILE	N-CA-C	-5.91	104.59	110.62
1	C	816	LYS	N-CA-C	-5.81	104.95	111.28
1	B	1084	GLN	N-CA-CB	5.79	118.63	110.12
1	C	1157	ASP	CB-CA-C	5.79	118.49	110.16
1	A	1084	GLN	N-CA-CB	5.75	118.57	110.12
1	B	1219	PRO	N-CA-CB	-5.74	97.51	103.08
1	B	1161	PRO	N-CA-C	5.65	124.11	112.47
1	A	824	TYR	CA-C-N	-5.64	117.06	121.86
1	A	824	TYR	C-N-CA	-5.64	117.06	121.86
1	C	1031	GLN	N-CA-C	-5.64	105.14	111.28
1	B	1136	PHE	CB-CA-C	-5.63	100.95	110.19
1	B	816	LYS	N-CA-C	-5.62	105.15	111.28
1	A	816	LYS	N-CA-C	-5.59	105.19	111.28
1	C	774	ASN	CA-CB-CG	-5.58	107.02	112.60
1	C	1219	PRO	N-CA-CB	-5.58	97.67	103.08
1	A	1136	PHE	CA-CB-CG	5.55	119.35	113.80
1	A	1207	PRO	CB-CA-C	-5.44	105.69	113.09
1	A	812	ASN	CB-CA-C	-5.43	104.45	109.83
1	A	1031	GLN	N-CA-C	-5.43	105.36	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1071	ILE	N-CA-C	-5.40	105.11	110.62
1	C	1154	GLY	CA-C-N	-5.39	115.36	122.85
1	C	1154	GLY	C-N-CA	-5.39	115.36	122.85
1	A	1219	PRO	N-CA-CB	-5.39	97.86	103.08
1	A	1171	TYR	CA-C-O	-5.38	115.63	121.87
1	C	1171	TYR	CA-C-O	-5.37	115.71	121.56
1	C	1122	HIS	CA-CB-CG	5.34	119.14	113.80
1	C	1207	PRO	CB-CA-C	-5.33	105.84	113.09
1	C	783	PRO	CB-CA-C	-5.33	102.77	111.56
1	B	1171	TYR	CA-C-O	-5.32	115.76	121.56
1	A	1047	ILE	N-CA-C	-5.19	105.32	110.62
1	A	825	GLY	N-CA-C	5.12	117.75	111.45
1	A	1116	PHE	CA-C-O	-5.12	115.85	121.58
1	C	1084	GLN	N-CA-CB	5.10	117.41	110.01
1	B	1128	VAL	CA-C-O	-5.09	115.45	120.90
1	C	1047	ILE	N-CA-C	-5.08	105.44	110.62
1	A	831	ILE	N-CA-C	-5.02	105.50	110.62
1	A	1026	VAL	N-CA-C	-5.02	105.65	110.72
1	C	1084	GLN	CB-CA-C	-5.01	103.01	110.88
1	C	815	GLN	N-CA-C	-5.01	107.00	113.01

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	2439	0	2041	59	0
1	B	2437	0	2021	50	0
1	C	2415	0	1986	50	0
2	A	70	0	65	0	0
2	B	70	0	65	0	0
2	C	84	0	78	0	0
All	All	7515	0	6256	133	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 10.

All (133) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1156:CYS:HA	1:C:1164:CYS:HB3	1.60	0.84
1:A:1050:SER:O	1:A:1054:ILE:HG12	1.83	0.78
1:C:1050:SER:O	1:C:1054:ILE:HG12	1.85	0.76
1:B:1160:ASN:HD22	1:B:1161:PRO:HD2	1.51	0.74
1:A:1172:PHE:HZ	1:A:1202:THR:HG22	1.53	0.73
1:B:1152:ALA:C	1:B:1154:GLY:H	1.96	0.72
1:A:1166:ALA:HB2	1:A:1205:VAL:CG1	2.20	0.72
1:B:1272:VAL:HG11	1:C:1010:THR:OG1	1.90	0.72
1:B:1240:LYS:HB2	1:B:1242:VAL:HG13	1.70	0.71
1:A:1166:ALA:HB2	1:A:1205:VAL:HG11	1.74	0.69
1:A:809:TYR:HE2	1:A:1078:LEU:HB2	1.57	0.68
1:C:1240:LYS:HB2	1:C:1242:VAL:HG13	1.74	0.68
1:A:1260:LEU:HD13	1:C:1026:VAL:HG11	1.77	0.65
1:A:809:TYR:CE2	1:A:1078:LEU:HB2	2.31	0.65
1:A:1033:LEU:HD21	1:C:1032:ALA:HB1	1.79	0.65
1:C:780:LEU:HD22	1:C:780:LEU:C	2.22	0.64
1:C:1256:ASN:O	1:C:1257:THR:HG23	1.97	0.64
1:A:1166:ALA:CB	1:A:1205:VAL:CG1	2.75	0.64
1:B:1156:CYS:N	1:B:1164:CYS:SG	2.70	0.63
1:B:1152:ALA:C	1:B:1154:GLY:N	2.53	0.63
1:A:1010:THR:O	1:A:1013:THR:OG1	2.16	0.62
1:A:1240:LYS:HB2	1:A:1242:VAL:HG13	1.82	0.62
1:A:1067:ILE:HD13	1:B:1235:LEU:HD11	1.82	0.61
1:A:1047:ILE:HG21	1:C:1047:ILE:HD12	1.82	0.61
1:C:1256:ASN:O	1:C:1257:THR:CG2	2.49	0.61
1:C:1167:PRO:HB3	1:C:1170:GLY:O	2.01	0.61
1:B:797:THR:HG22	1:B:1133:GLY:HA2	1.83	0.60
1:B:1026:VAL:HG11	1:C:1260:LEU:HD13	1.83	0.60
1:C:835:LEU:HD13	1:C:1082:VAL:HG13	1.83	0.60
1:A:1081:PHE:HB3	1:B:1223:LEU:HD13	1.84	0.60
1:C:780:LEU:HD22	1:C:780:LEU:O	2.02	0.60
1:B:1130:ALA:HB3	1:B:1133:GLY:O	2.02	0.60
1:C:809:TYR:CE2	1:C:1078:LEU:HB2	2.36	0.59
1:A:1223:LEU:HD13	1:C:1081:PHE:HB3	1.83	0.59
1:C:817:CYS:SG	1:C:818:GLU:N	2.76	0.59
1:A:817:CYS:HA	1:A:1071:ILE:HD11	1.84	0.58
1:A:1047:ILE:HD12	1:B:1047:ILE:HG21	1.85	0.58
1:B:1081:PHE:HB3	1:C:1223:LEU:HD13	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1166:ALA:CB	1:A:1205:VAL:HG13	2.33	0.57
1:A:1167:PRO:HB3	1:A:1170:GLY:O	2.03	0.57
1:A:1067:ILE:HD13	1:B:1235:LEU:CD1	2.33	0.57
1:A:998:ALA:HA	1:A:1001:PHE:HB3	1.87	0.57
1:B:809:TYR:CE1	1:B:1078:LEU:HB2	2.40	0.57
1:A:1130:ALA:HB3	1:A:1133:GLY:O	2.05	0.57
1:A:1166:ALA:CB	1:A:1205:VAL:HG11	2.34	0.57
1:A:1067:ILE:HD11	1:B:1235:LEU:HD21	1.86	0.56
1:A:811:CYS:SG	1:A:812:ASN:N	2.78	0.56
1:B:1160:ASN:ND2	1:B:1161:PRO:HD2	2.19	0.56
1:B:1152:ALA:HB2	1:B:1184:TRP:CH2	2.41	0.56
1:C:1256:ASN:C	1:C:1257:THR:HG23	2.31	0.56
1:C:817:CYS:HB2	1:C:1071:ILE:HD11	1.87	0.55
1:B:1155:LEU:HD22	1:B:1164:CYS:SG	2.46	0.55
1:C:1130:ALA:HB3	1:C:1133:GLY:O	2.06	0.55
1:B:1167:PRO:HB3	1:B:1170:GLY:O	2.08	0.54
1:B:1272:VAL:CG1	1:C:1010:THR:OG1	2.56	0.54
1:C:809:TYR:HE2	1:C:1078:LEU:HB2	1.74	0.53
1:B:1047:ILE:HD12	1:C:1047:ILE:HG21	1.89	0.53
1:C:1152:ALA:HB2	1:C:1184:TRP:CH2	2.45	0.52
1:C:1162:THR:O	1:C:1207:PRO:HB3	2.10	0.51
1:C:1235:LEU:HG	1:C:1239:PHE:HB2	1.91	0.51
1:A:797:THR:HG22	1:A:1133:GLY:HA2	1.91	0.51
1:C:1152:ALA:O	1:C:1153:TYR:C	2.54	0.50
1:C:797:THR:HG22	1:C:1133:GLY:HA2	1.93	0.50
1:B:1156:CYS:HB3	1:B:1164:CYS:HA	1.92	0.50
1:A:1152:ALA:HB2	1:A:1184:TRP:CH2	2.46	0.50
1:B:1005:LEU:O	1:B:1009:GLN:HG2	2.12	0.49
1:A:1166:ALA:HB3	1:A:1205:VAL:HG13	1.95	0.49
1:A:1026:VAL:HG11	1:B:1260:LEU:HD13	1.95	0.49
1:A:1263:THR:O	1:A:1266:MET:HB3	2.12	0.49
1:A:806:CYS:HB3	1:A:827:PHE:CE2	2.48	0.49
1:A:1238:PHE:HZ	1:C:1070:LEU:HB2	1.77	0.49
1:A:1033:LEU:HD21	1:C:1032:ALA:CB	2.41	0.48
1:C:1005:LEU:O	1:C:1009:GLN:HG2	2.14	0.48
1:B:996:LEU:O	1:B:1000:LYS:HG2	2.14	0.48
1:A:1162:THR:O	1:A:1207:PRO:HB3	2.14	0.47
1:A:1269:LEU:CD2	1:C:1012:PHE:CG	2.97	0.47
1:C:1107:VAL:HG11	1:C:1127:VAL:HG12	1.96	0.47
1:A:1043:THR:O	1:A:1047:ILE:HG12	2.14	0.47
1:A:810:VAL:HG11	1:A:1075:LEU:HD21	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1102:LYS:O	1:A:1106:CYS:HB2	2.13	0.47
1:C:1273:VAL:HA	1:C:1276:LEU:HD12	1.95	0.47
1:B:1235:LEU:HG	1:B:1239:PHE:HB2	1.96	0.46
1:A:987:GLN:HE21	1:C:990:LEU:HD22	1.79	0.46
1:C:1105:GLU:O	1:C:1111:SER:HB2	2.15	0.46
1:A:1070:LEU:HB2	1:B:1238:PHE:HZ	1.79	0.46
1:A:1235:LEU:HG	1:A:1239:PHE:HB2	1.96	0.46
1:C:1120:GLY:HA3	1:C:1141:TYR:O	2.16	0.46
1:C:1156:CYS:HA	1:C:1164:CYS:CB	2.39	0.46
1:B:835:LEU:HD13	1:B:1082:VAL:HG23	1.98	0.46
1:B:1160:ASN:O	1:B:1162:THR:N	2.48	0.45
1:C:1010:THR:O	1:C:1013:THR:OG1	2.26	0.45
1:B:1070:LEU:HB2	1:C:1238:PHE:HZ	1.81	0.45
1:A:1116:PHE:HD1	1:A:1116:PHE:HA	1.69	0.45
1:A:987:GLN:NE2	1:C:990:LEU:HD22	2.32	0.45
1:A:1005:LEU:O	1:A:1009:GLN:HG2	2.17	0.44
1:C:1102:LYS:O	1:C:1106:CYS:HB2	2.18	0.44
1:B:1075:LEU:HD23	1:B:1075:LEU:HA	1.74	0.44
1:C:1075:LEU:HA	1:C:1075:LEU:HD23	1.68	0.44
1:A:1166:ALA:HB2	1:A:1205:VAL:HG13	1.94	0.44
1:A:1269:LEU:HD23	1:C:1012:PHE:CG	2.53	0.44
1:A:821:LEU:HD13	1:A:1071:ILE:HG21	2.01	0.43
1:A:810:VAL:HG11	1:A:1075:LEU:CD2	2.48	0.43
1:B:790:VAL:HG22	1:B:1139:VAL:HG22	2.00	0.43
1:B:1152:ALA:O	1:B:1154:GLY:N	2.51	0.43
1:B:1162:THR:O	1:B:1207:PRO:HB3	2.17	0.43
1:A:1067:ILE:CD1	1:B:1235:LEU:HD21	2.49	0.43
1:A:1130:ALA:HB1	1:A:1131:PRO:CD	2.49	0.43
1:B:1105:GLU:O	1:B:1111:SER:HB2	2.18	0.43
1:A:1276:LEU:HA	1:A:1279:SER:HB3	2.01	0.42
1:A:1152:ALA:O	1:A:1153:TYR:C	2.62	0.42
1:B:1174:LYS:HB2	1:B:1184:TRP:CZ3	2.54	0.42
1:A:835:LEU:HD12	1:A:835:LEU:HA	1.75	0.42
1:B:1259:LEU:HD12	1:B:1259:LEU:HA	1.88	0.42
1:C:1122:HIS:CD2	1:C:1122:HIS:C	2.97	0.42
1:A:810:VAL:HG21	1:A:1075:LEU:HD21	2.02	0.42
1:C:1256:ASN:C	1:C:1257:THR:CG2	2.93	0.42
1:A:821:LEU:HD12	1:A:821:LEU:HA	1.78	0.42
1:B:1263:THR:O	1:B:1266:MET:HB3	2.20	0.42
1:A:810:VAL:HG21	1:A:1075:LEU:CD2	2.49	0.42
1:A:1105:GLU:O	1:A:1111:SER:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:790:VAL:HG22	1:C:1139:VAL:HG22	2.02	0.41
1:B:1120:GLY:HA3	1:B:1141:TYR:O	2.21	0.41
1:B:835:LEU:HD12	1:B:835:LEU:HA	1.80	0.41
1:B:1079:ASN:HA	1:B:1082:VAL:HG12	2.03	0.41
1:B:1130:ALA:HB1	1:B:1131:PRO:HD2	2.03	0.41
1:B:1240:LYS:HB2	1:B:1242:VAL:CG1	2.46	0.41
1:A:1078:LEU:CD1	1:A:1082:VAL:HG23	2.51	0.40
1:B:1070:LEU:HD12	1:B:1070:LEU:HA	1.90	0.40
1:B:1070:LEU:HD22	1:C:1238:PHE:CE2	2.57	0.40
1:B:1005:LEU:HA	1:B:1005:LEU:HD12	1.76	0.40
1:A:1105:GLU:CD	1:B:1113:ARG:HH21	2.28	0.40
1:B:1172:PHE:HZ	1:B:1202:THR:HG22	1.87	0.40
1:C:811:CYS:SG	1:C:812:ASN:N	2.93	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	352/539 (65%)	329 (94%)	16 (4%)	7 (2%)	6	32
1	B	352/539 (65%)	327 (93%)	17 (5%)	8 (2%)	5	30
1	C	351/539 (65%)	325 (93%)	16 (5%)	10 (3%)	4	27
All	All	1055/1617 (65%)	981 (93%)	49 (5%)	25 (2%)	7	29

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1219	PRO
1	A	1247	PRO
1	B	1117	CYS
1	B	1161	PRO

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Mol	Chain	Res	Type
1	B	1219	PRO
1	B	1247	PRO
1	C	1117	CYS
1	C	1124	VAL
1	C	1219	PRO
1	C	1247	PRO
1	A	1117	CYS
1	A	1124	VAL
1	C	1158	ALA
1	B	1124	VAL
1	A	1220	PRO
1	C	1220	PRO
1	A	1221	PRO
1	A	1242	VAL
1	B	811	CYS
1	B	1220	PRO
1	B	1242	VAL
1	C	811	CYS
1	C	1221	PRO
1	C	1242	VAL
1	C	1160	ASN

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	216/458 (47%)	208 (96%)	8 (4%)	29	53
1	B	217/458 (47%)	204 (94%)	13 (6%)	16	41
1	C	208/458 (45%)	201 (97%)	7 (3%)	32	55
All	All	641/1374 (47%)	613 (96%)	28 (4%)	26	48

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	778	PHE

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Mol	Chain	Res	Type
1	A	1116	PHE
1	A	1122	HIS
1	A	1126	PHE
1	A	1163	ASN
1	A	1191	PHE
1	A	1219	PRO
1	A	1221	PRO
1	B	774	ASN
1	B	824	TYR
1	B	1122	HIS
1	B	1126	PHE
1	B	1132	ASN
1	B	1156	CYS
1	B	1163	ASN
1	B	1164	CYS
1	B	1191	PHE
1	B	1219	PRO
1	B	1221	PRO
1	B	1256	ASN
1	B	1277	ASN
1	C	774	ASN
1	C	1122	HIS
1	C	1126	PHE
1	C	1160	ASN
1	C	1191	PHE
1	C	1219	PRO
1	C	1221	PRO

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	1002	ASN
1	A	1029	ASN
1	A	1122	HIS
1	A	1217	ASN
1	B	774	ASN
1	B	1023	GLN
1	B	1029	ASN
1	B	1122	HIS
1	B	1160	ASN
1	C	1023	GLN
1	C	1029	ASN

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Mol	Chain	Res	Type
1	C	1217	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

16 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	$\# Z > 2$
2	NAG	C	1302	1	14,14,15	0.36	0	17,19,21	0.97	1 (5%)
2	NAG	B	1302	1	14,14,15	0.34	0	17,19,21	0.67	0
2	NAG	C	1301	1	14,14,15	0.23	0	17,19,21	1.12	1 (5%)
2	NAG	C	1303	1	14,14,15	0.37	0	17,19,21	0.63	0
2	NAG	A	1303	1	14,14,15	0.39	0	17,19,21	0.79	0
2	NAG	A	1302	1	14,14,15	0.34	0	17,19,21	1.07	2 (11%)
2	NAG	B	1303	1	14,14,15	0.39	0	17,19,21	0.64	0
2	NAG	B	1304	1	14,14,15	0.31	0	17,19,21	0.58	0
2	NAG	A	1301	1	14,14,15	0.33	0	17,19,21	1.06	1 (5%)
2	NAG	A	1304	1	14,14,15	0.28	0	17,19,21	0.64	0
2	NAG	B	1301	1	14,14,15	0.37	0	17,19,21	0.77	0
2	NAG	B	1305	1	14,14,15	0.30	0	17,19,21	0.65	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	1304	1	14,14,15	0.35	0	17,19,21	0.49	0
2	NAG	A	1305	1	14,14,15	0.33	0	17,19,21	0.74	0
2	NAG	C	1305	1	14,14,15	0.42	0	17,19,21	1.20	1 (5%)
2	NAG	C	1306	1	14,14,15	0.34	0	17,19,21	0.84	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
2	NAG	C	1302	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1302	1	-	3/6/23/26	0/1/1/1
2	NAG	C	1301	1	-	3/6/23/26	0/1/1/1
2	NAG	C	1303	1	-	2/6/23/26	0/1/1/1
2	NAG	A	1303	1	-	2/6/23/26	0/1/1/1
2	NAG	A	1302	1	-	4/6/23/26	0/1/1/1
2	NAG	B	1303	1	-	2/6/23/26	0/1/1/1
2	NAG	B	1304	1	-	3/6/23/26	0/1/1/1
2	NAG	A	1301	1	-	3/6/23/26	0/1/1/1
2	NAG	A	1304	1	-	3/6/23/26	0/1/1/1
2	NAG	B	1301	1	-	2/6/23/26	0/1/1/1
2	NAG	B	1305	1	-	2/6/23/26	0/1/1/1
2	NAG	C	1304	1	-	2/6/23/26	0/1/1/1
2	NAG	A	1305	1	-	2/6/23/26	0/1/1/1
2	NAG	C	1305	1	-	4/6/23/26	0/1/1/1
2	NAG	C	1306	1	-	2/6/23/26	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(°)
2	C	1305	NAG	C2-N2-C7	3.19	127.45	122.90
2	A	1302	NAG	C1-O5-C5	3.05	116.33	112.19
2	C	1301	NAG	C2-N2-C7	2.94	127.09	122.90
2	A	1301	NAG	C2-N2-C7	2.76	126.84	122.90
2	C	1302	NAG	C4-C3-C2	-2.50	107.35	111.02
2	A	1302	NAG	C2-N2-C7	2.24	126.10	122.90

There are no chirality outliers.

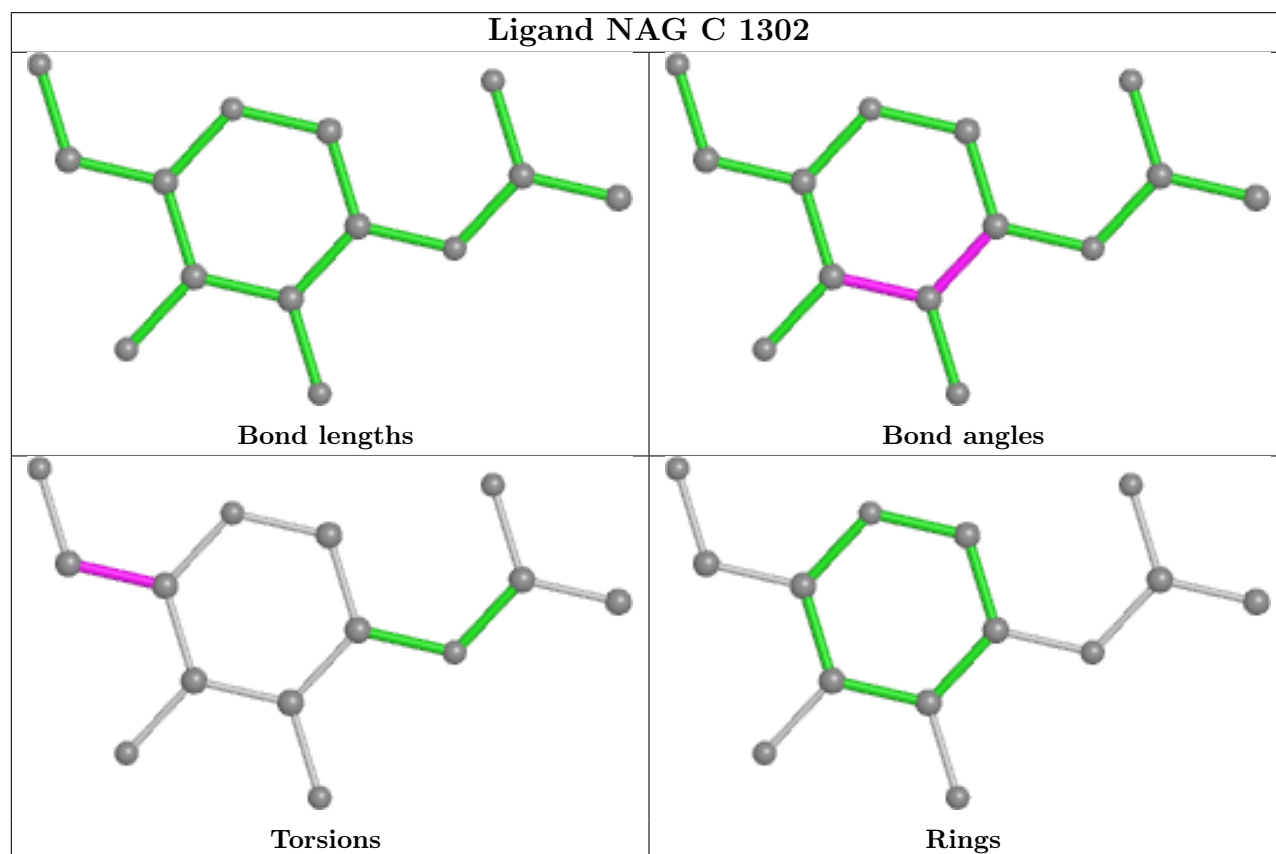
All (40) torsion outliers are listed below:

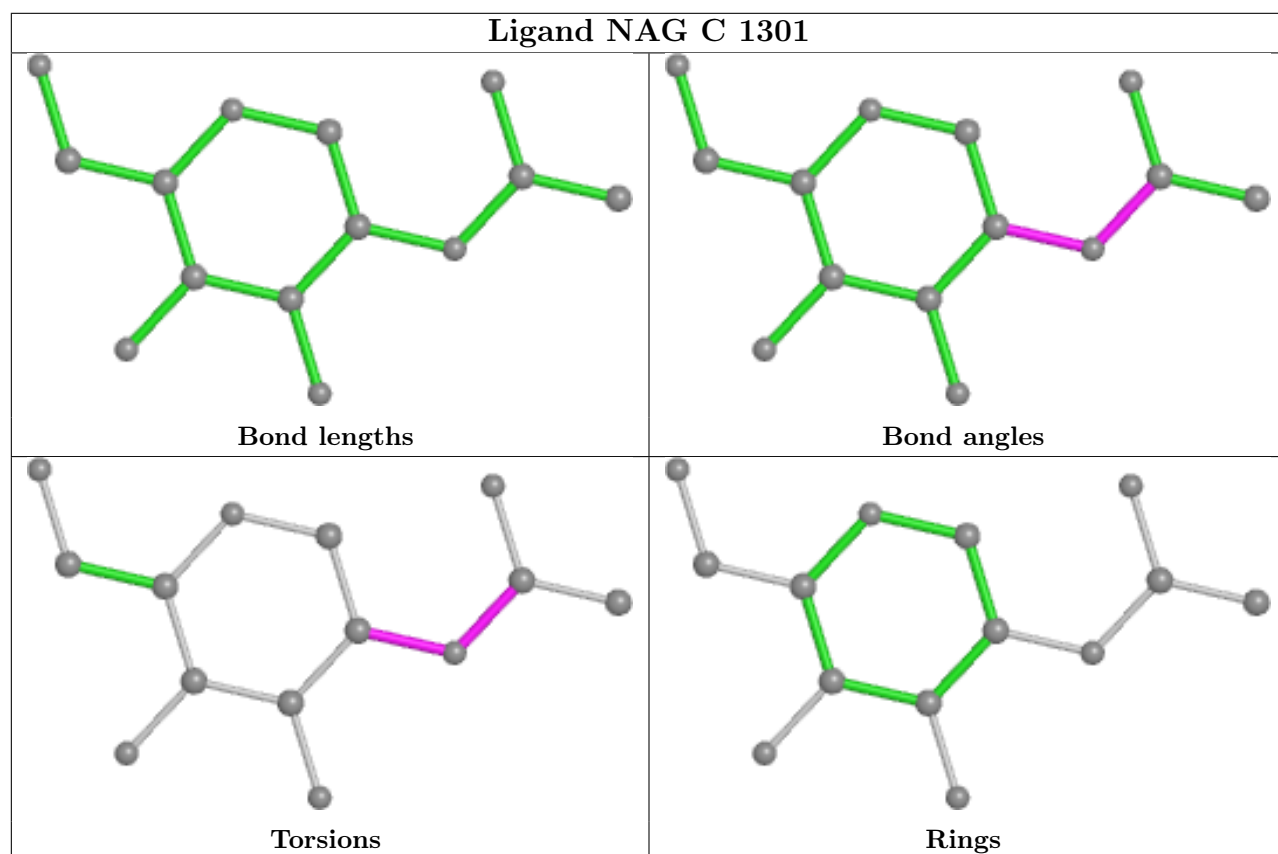
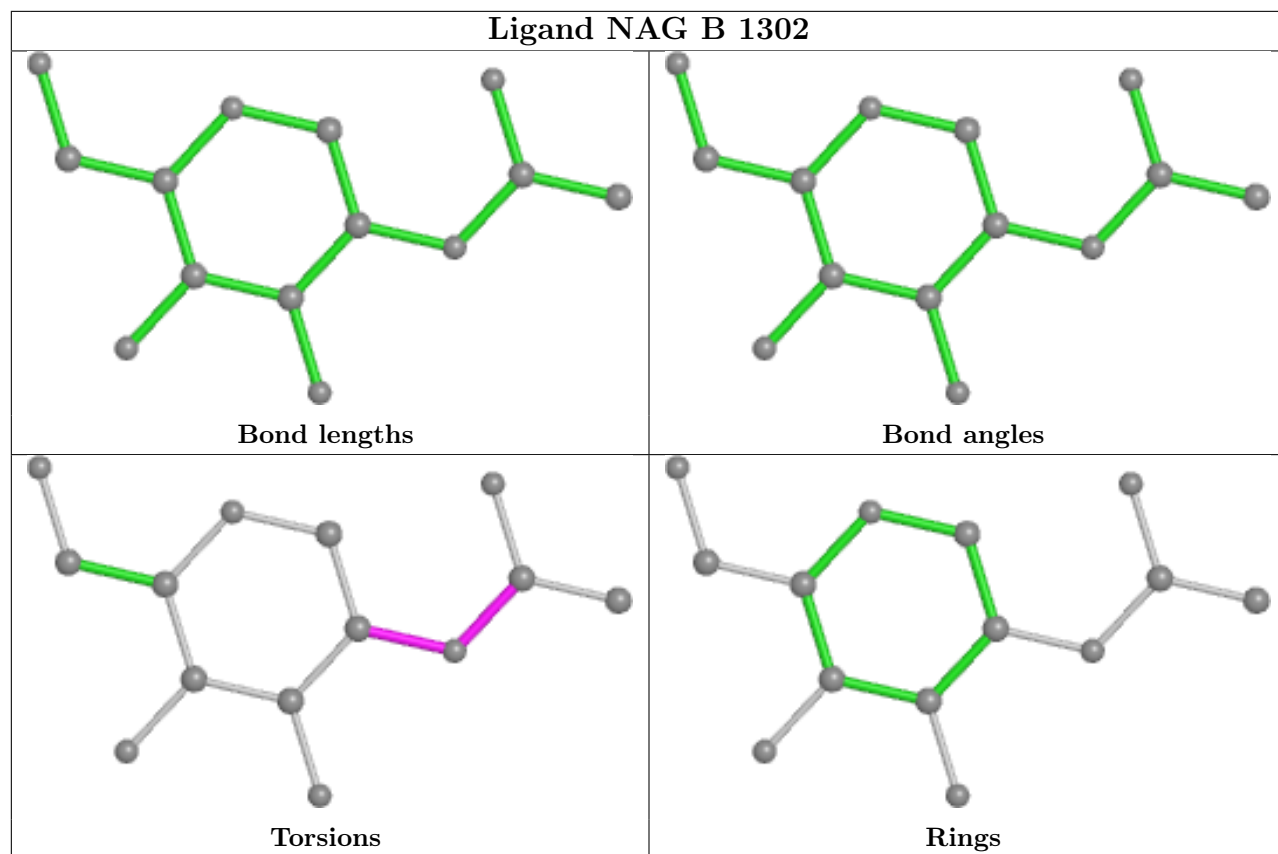
Mol	Chain	Res	Type	Atoms
2	A	1301	NAG	C8-C7-N2-C2
2	A	1301	NAG	O7-C7-N2-C2
2	A	1302	NAG	C1-C2-N2-C7
2	A	1302	NAG	C8-C7-N2-C2
2	A	1302	NAG	O7-C7-N2-C2
2	A	1303	NAG	C8-C7-N2-C2
2	A	1303	NAG	O7-C7-N2-C2
2	A	1305	NAG	C8-C7-N2-C2
2	A	1305	NAG	O7-C7-N2-C2
2	B	1302	NAG	O7-C7-N2-C2
2	B	1303	NAG	C8-C7-N2-C2
2	B	1303	NAG	O7-C7-N2-C2
2	C	1301	NAG	C8-C7-N2-C2
2	C	1301	NAG	O7-C7-N2-C2
2	C	1303	NAG	C8-C7-N2-C2
2	C	1303	NAG	O7-C7-N2-C2
2	C	1305	NAG	O7-C7-N2-C2
2	C	1306	NAG	O7-C7-N2-C2
2	B	1302	NAG	C8-C7-N2-C2
2	C	1305	NAG	C8-C7-N2-C2
2	C	1306	NAG	C8-C7-N2-C2
2	A	1301	NAG	C1-C2-N2-C7
2	B	1304	NAG	C8-C7-N2-C2
2	B	1305	NAG	C8-C7-N2-C2
2	B	1305	NAG	O7-C7-N2-C2
2	B	1301	NAG	C8-C7-N2-C2
2	B	1304	NAG	O7-C7-N2-C2
2	C	1304	NAG	C8-C7-N2-C2
2	B	1301	NAG	O7-C7-N2-C2
2	C	1304	NAG	O7-C7-N2-C2
2	C	1301	NAG	C1-C2-N2-C7
2	A	1304	NAG	C8-C7-N2-C2
2	C	1305	NAG	C1-C2-N2-C7
2	A	1302	NAG	O5-C5-C6-O6
2	A	1304	NAG	O5-C5-C6-O6
2	B	1304	NAG	O5-C5-C6-O6
2	B	1302	NAG	C3-C2-N2-C7
2	C	1305	NAG	C3-C2-N2-C7
2	A	1304	NAG	O7-C7-N2-C2
2	C	1302	NAG	C4-C5-C6-O6

There are no ring outliers.

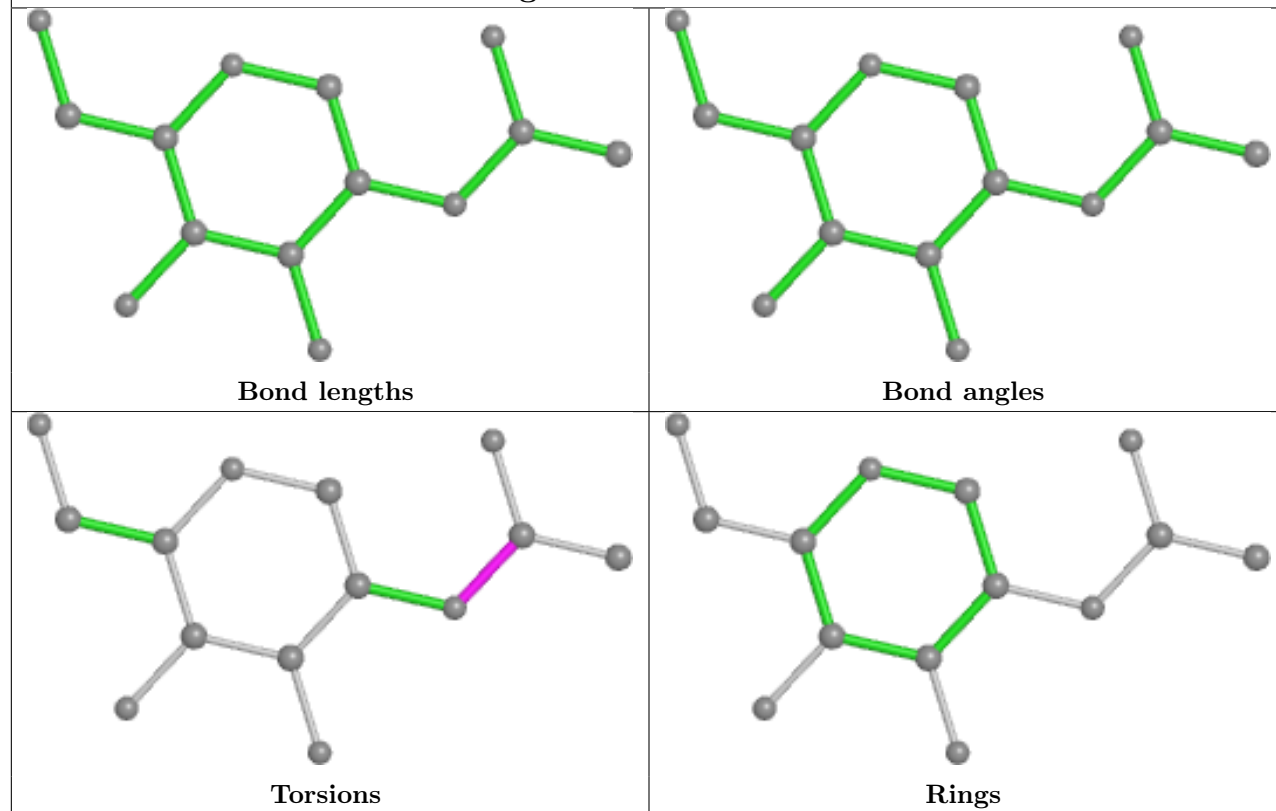
No monomer is involved in short contacts.

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

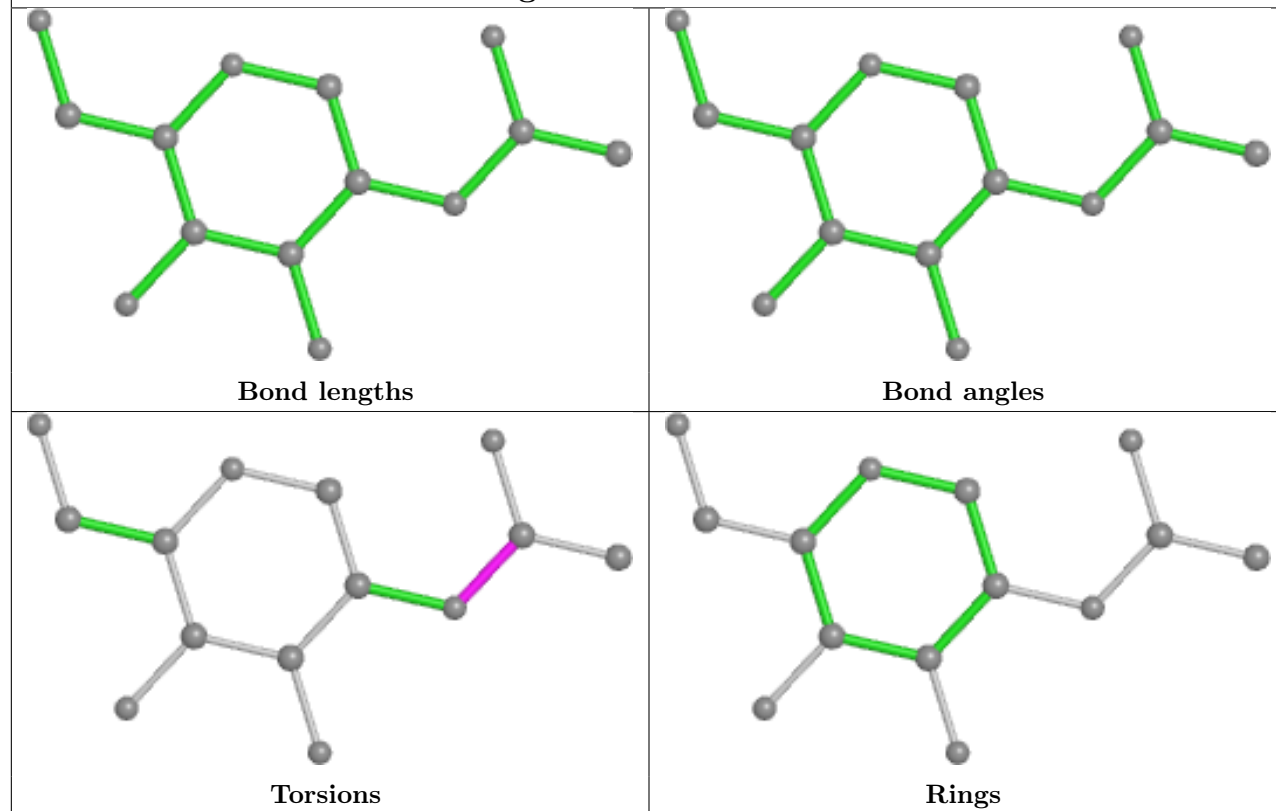


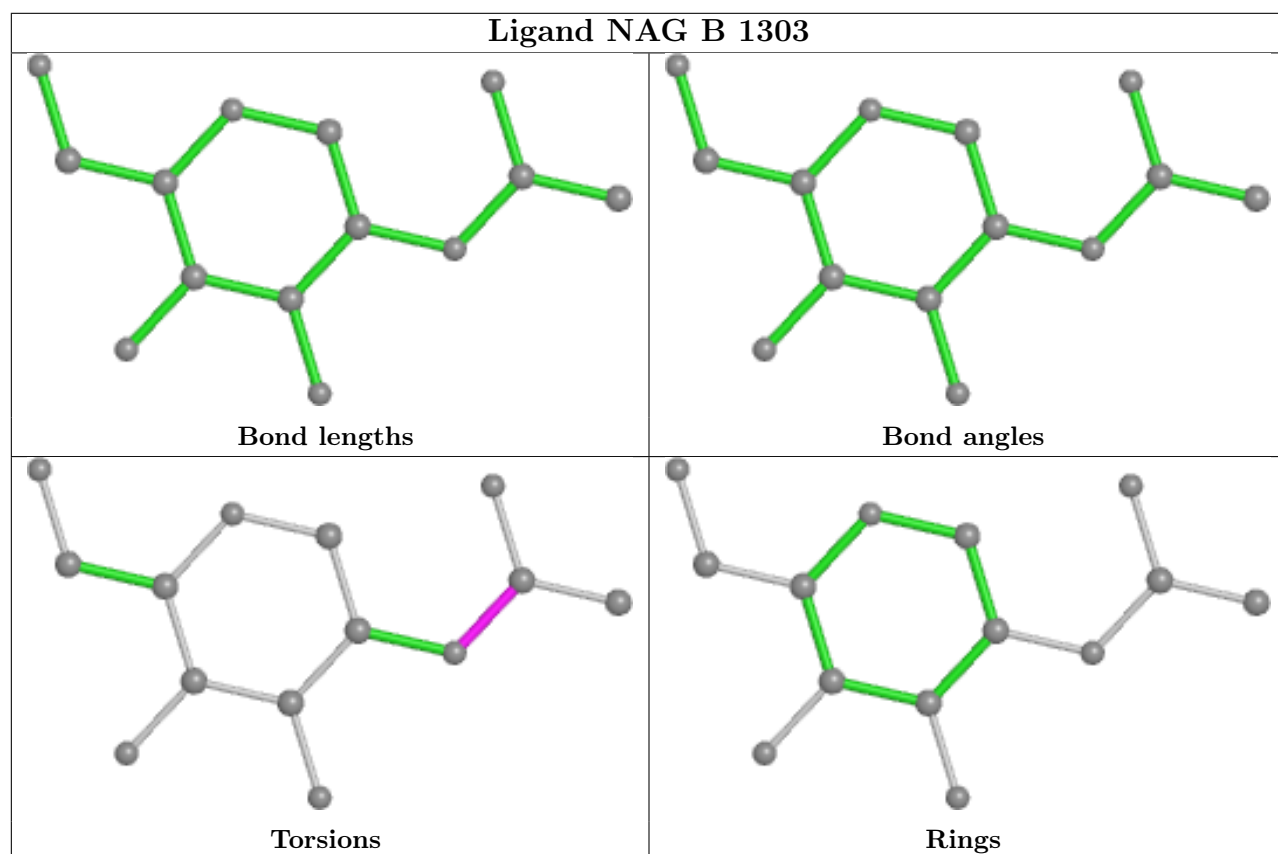
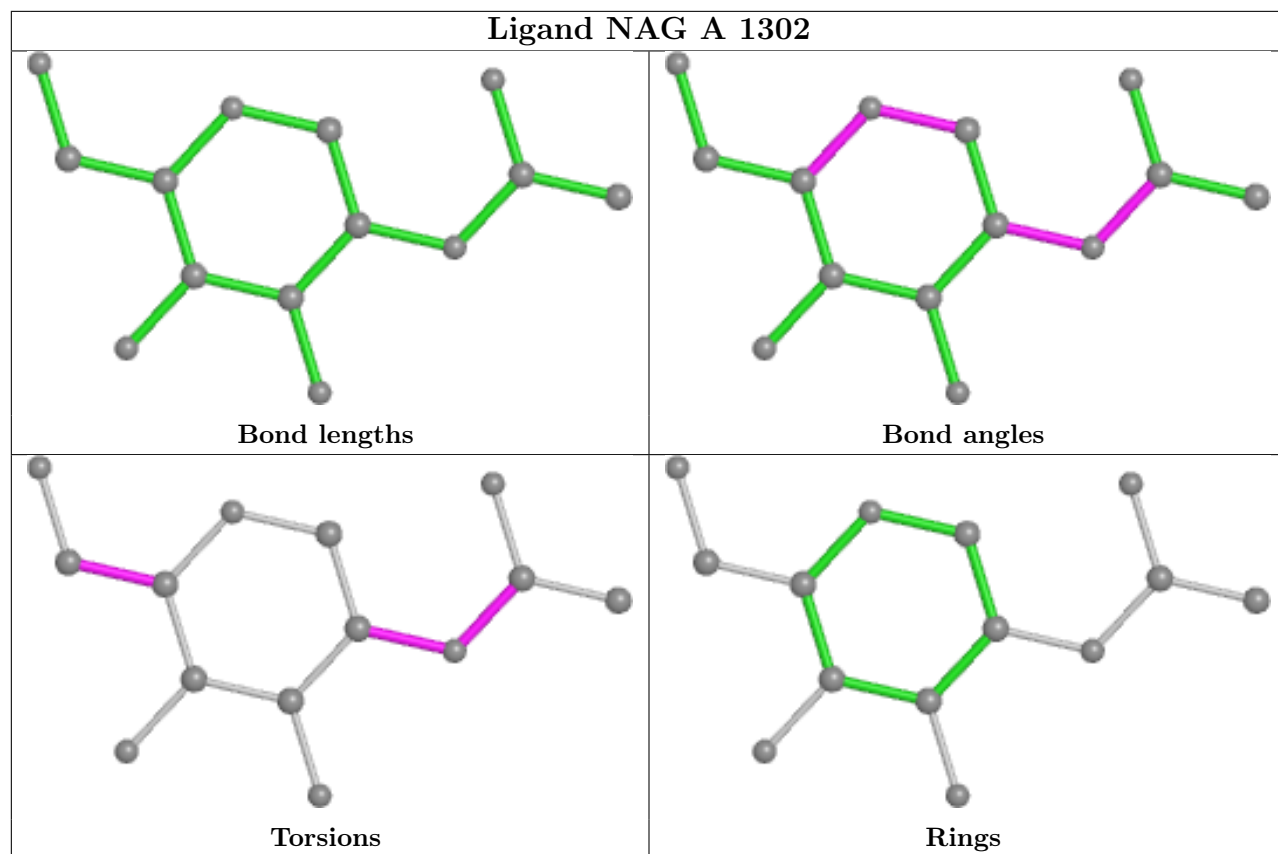


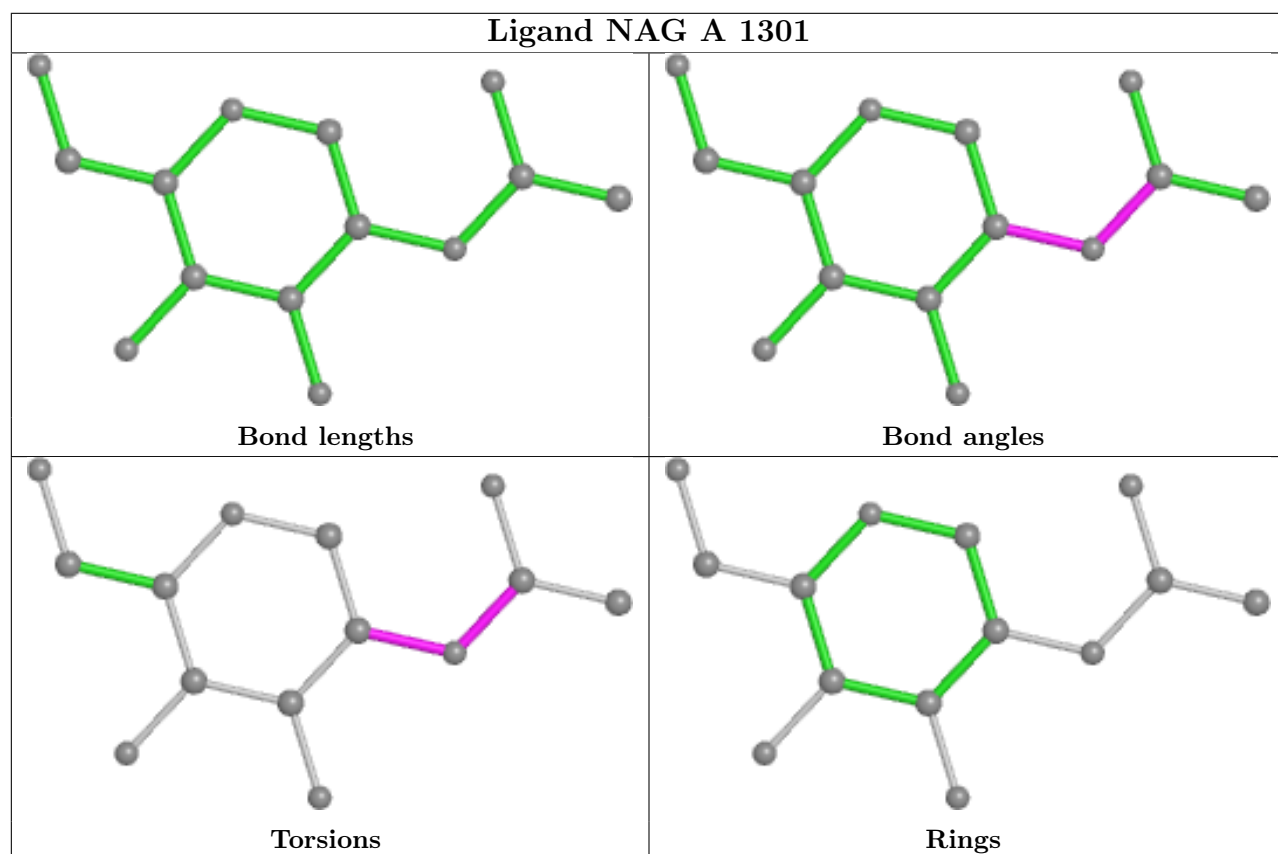
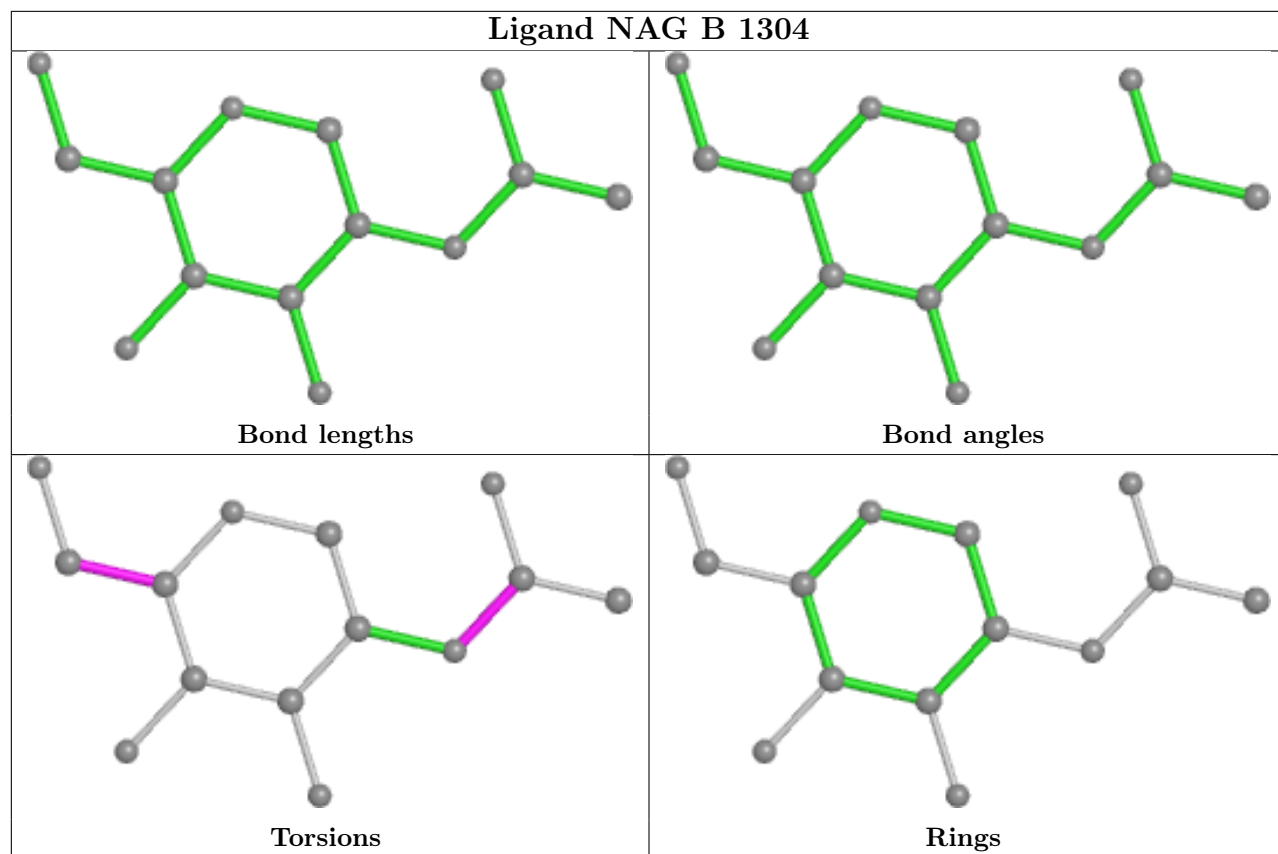
Ligand NAG C 1303

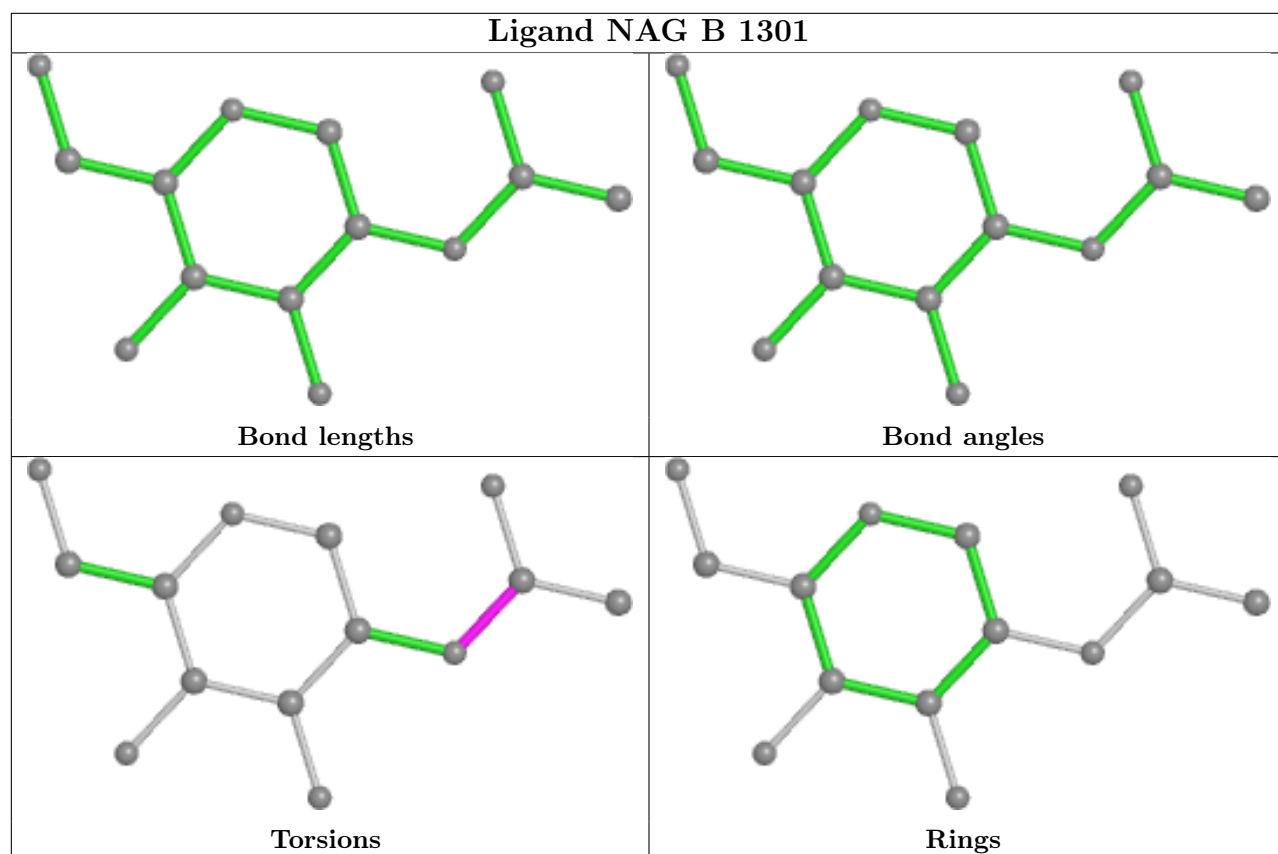
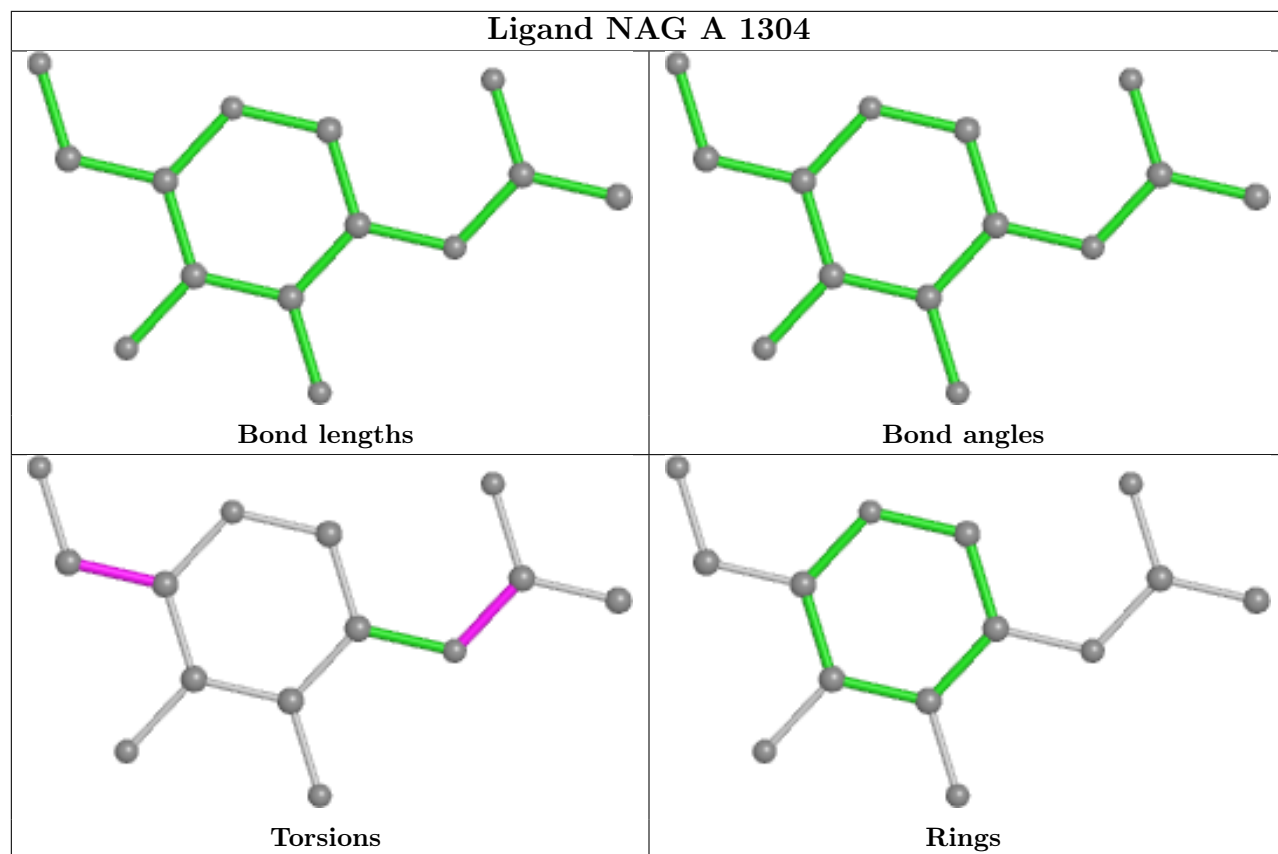


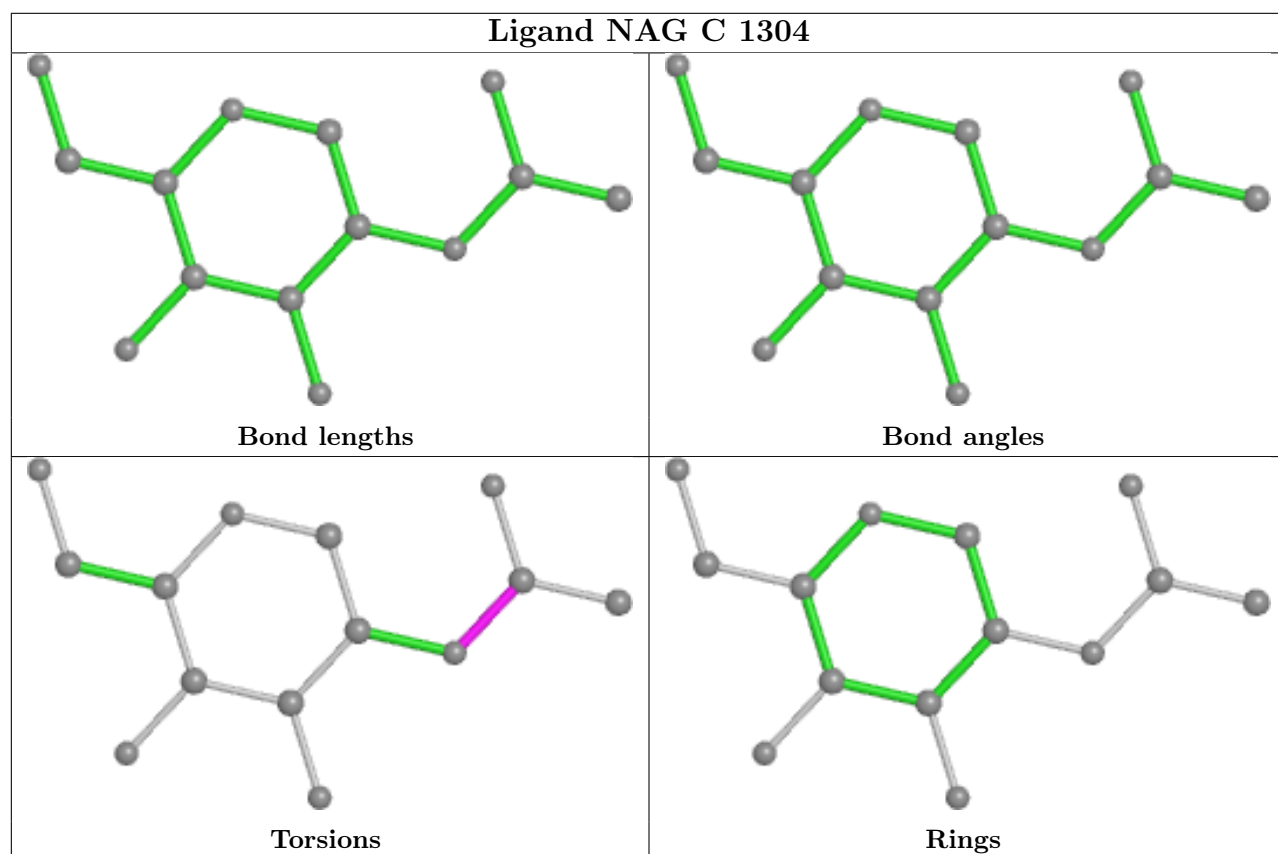
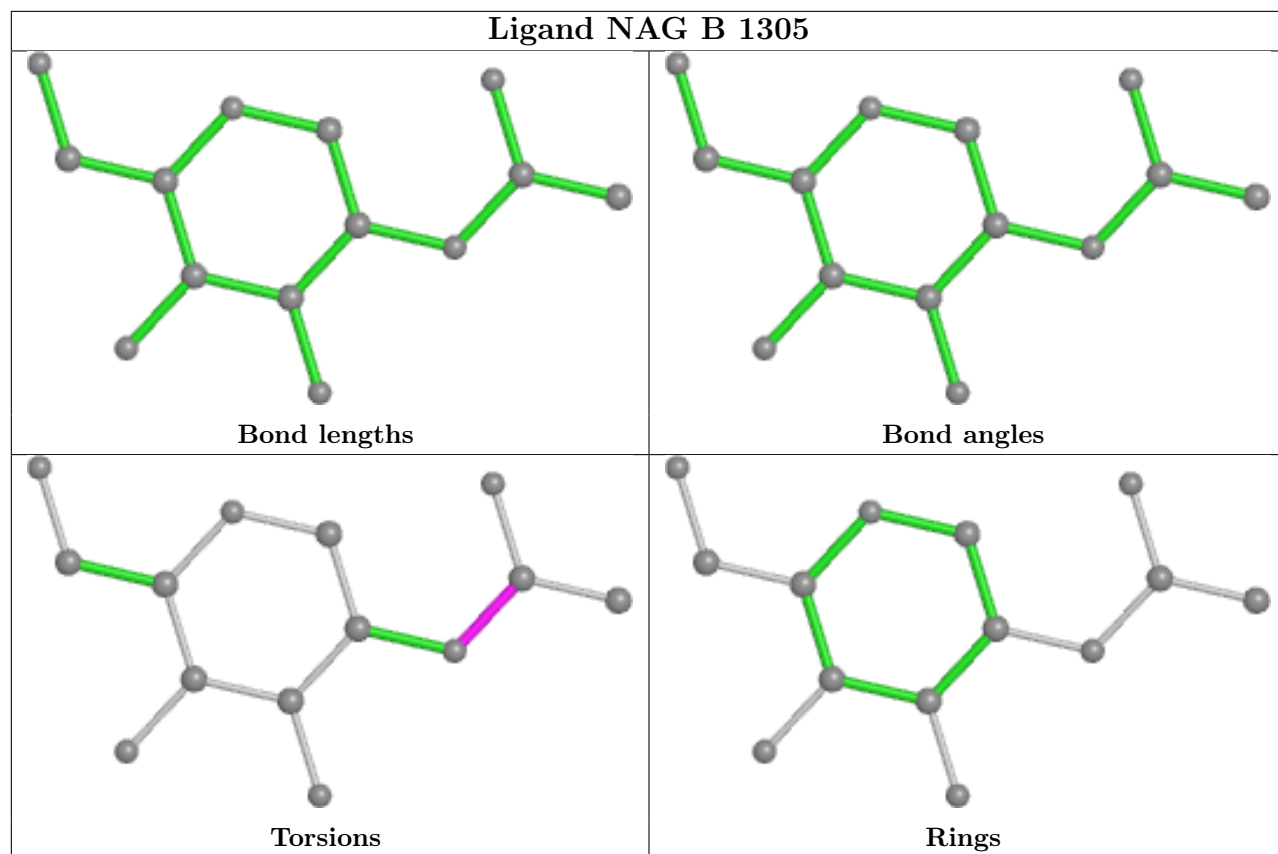
Ligand NAG A 1303

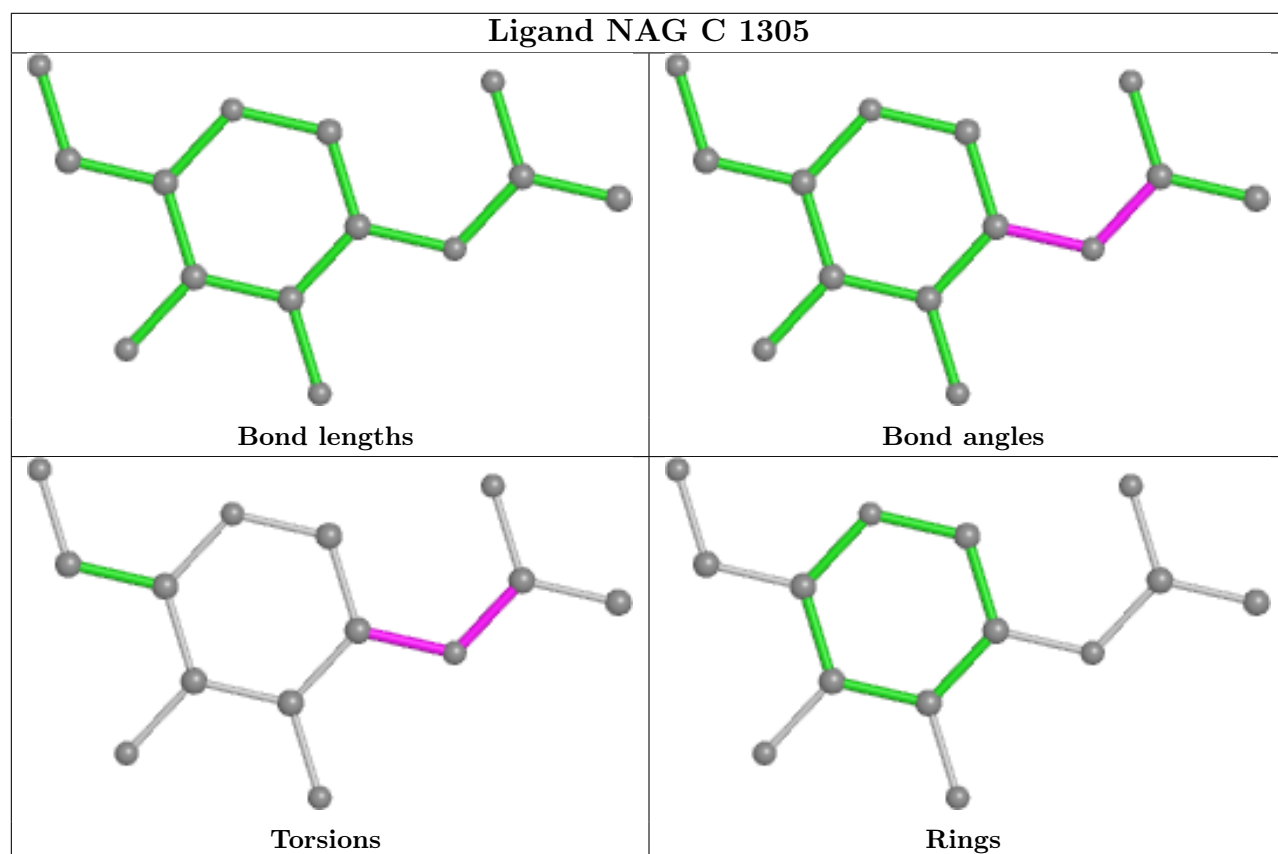
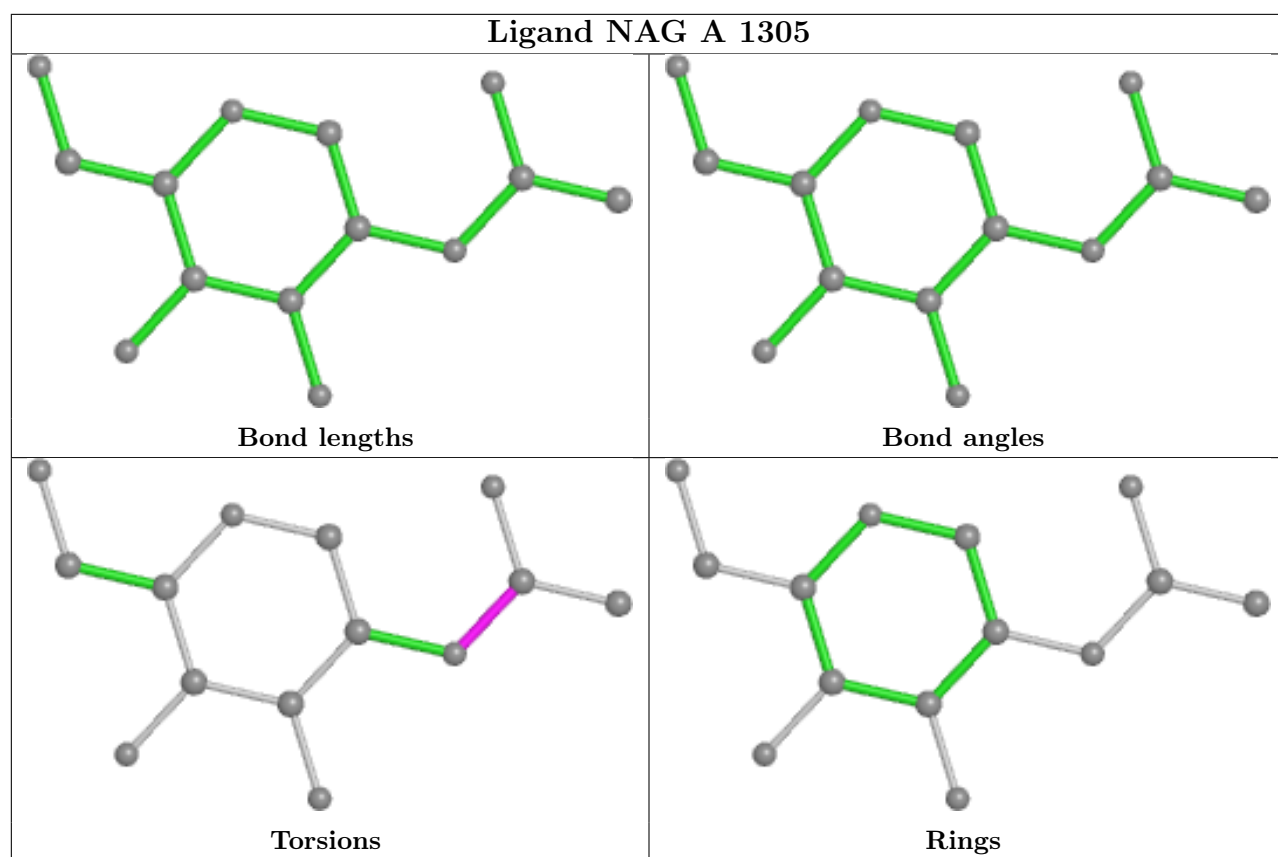


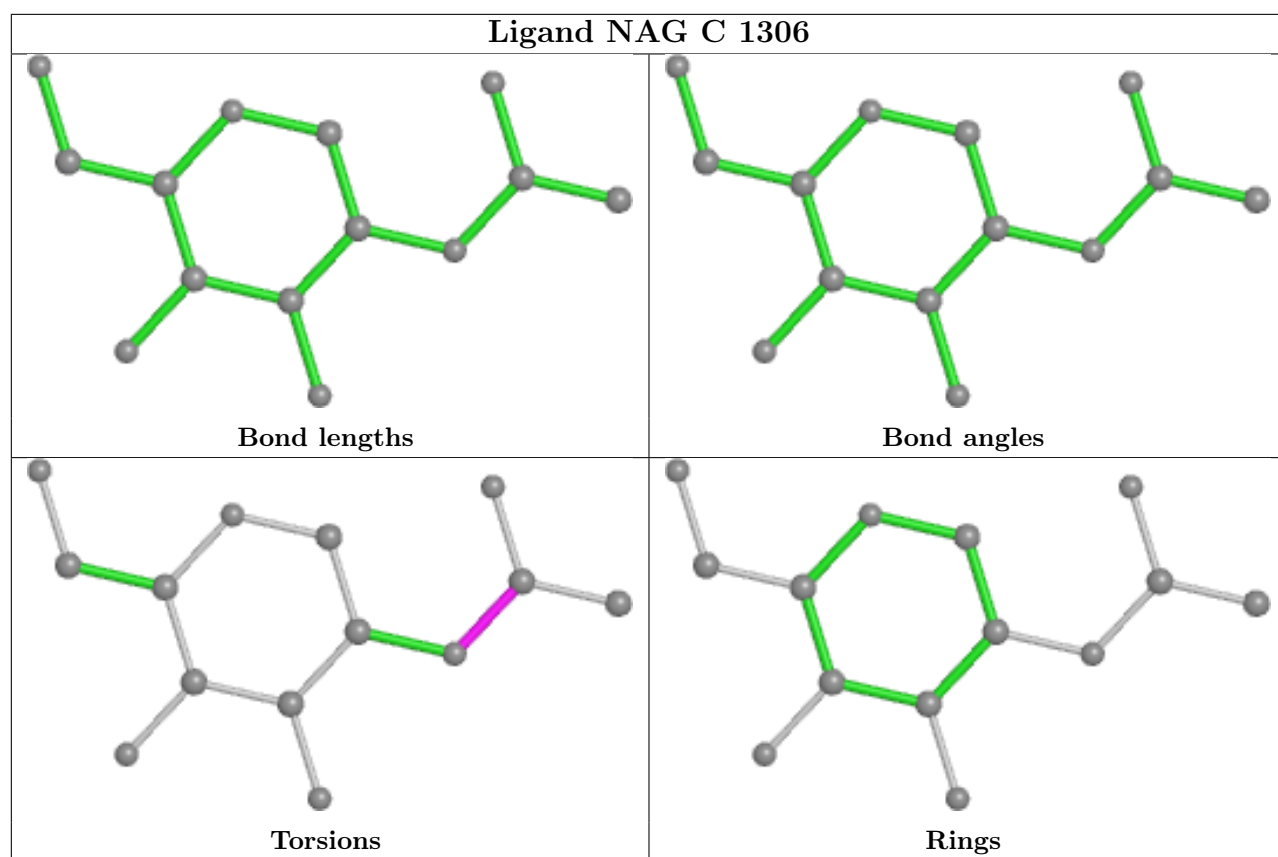












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

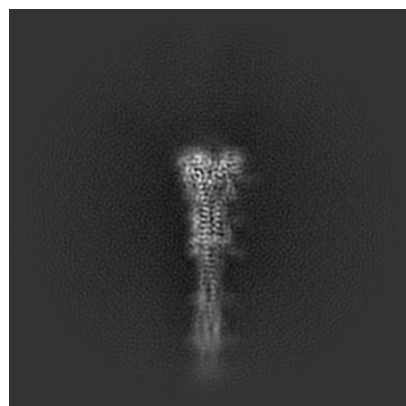
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-39768. These allow visual inspection of the internal detail of the map and identification of artifacts.

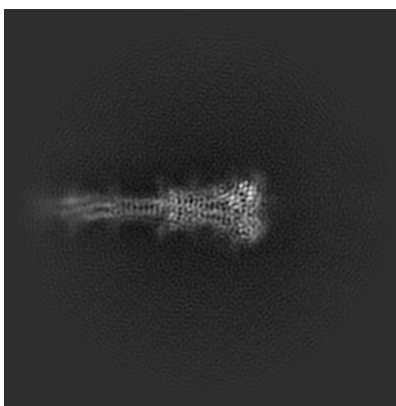
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

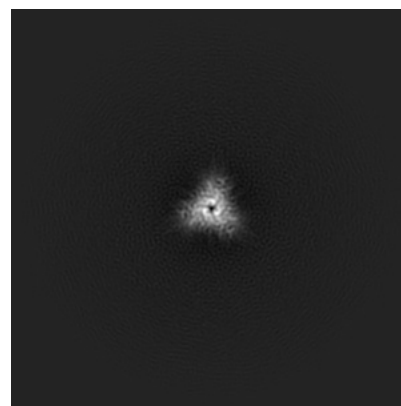
6.1.1 Primary map



X

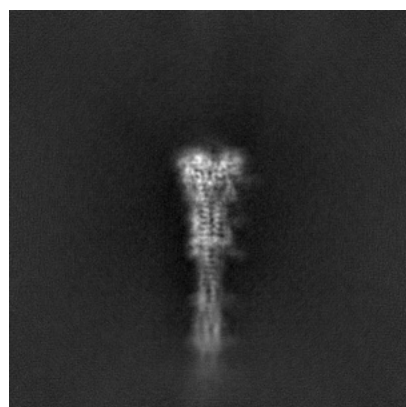


Y

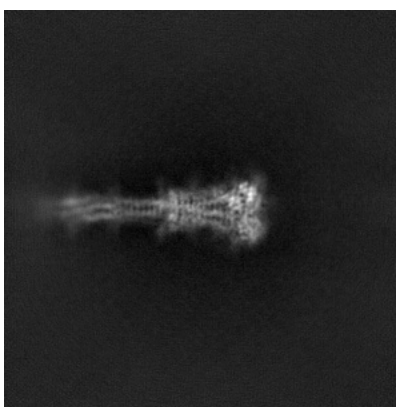


Z

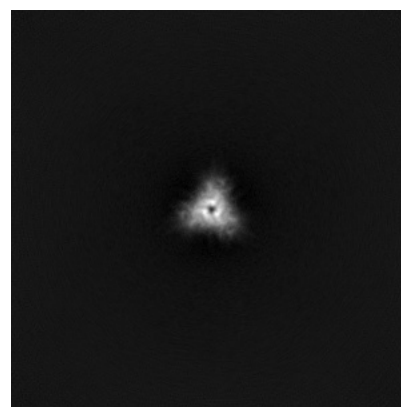
6.1.2 Raw map



X



Y



Z

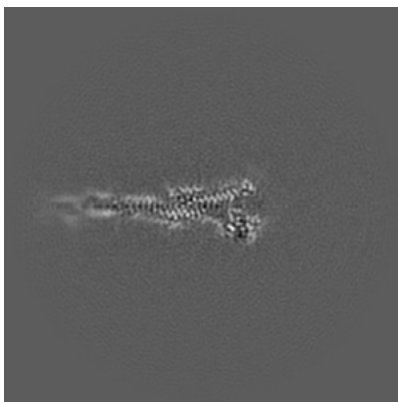
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

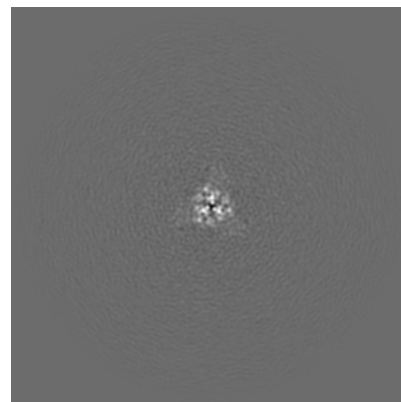
6.2.1 Primary map



X Index: 170

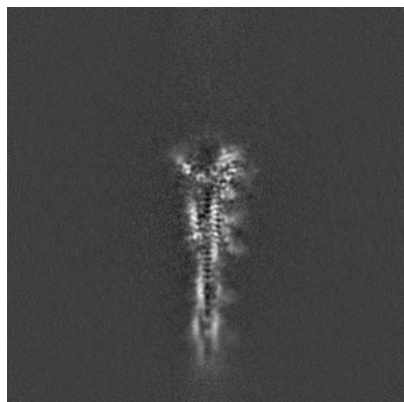


Y Index: 170

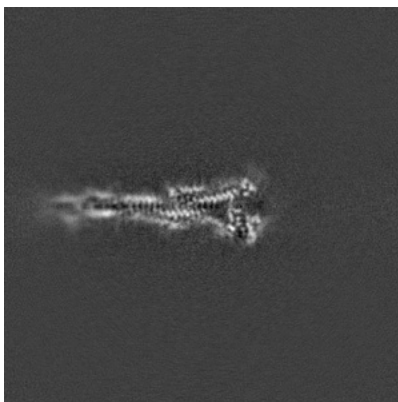


Z Index: 170

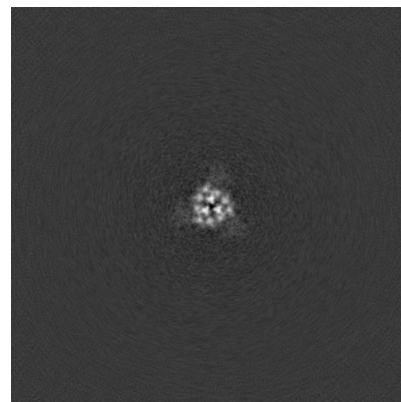
6.2.2 Raw map



X Index: 170



Y Index: 170

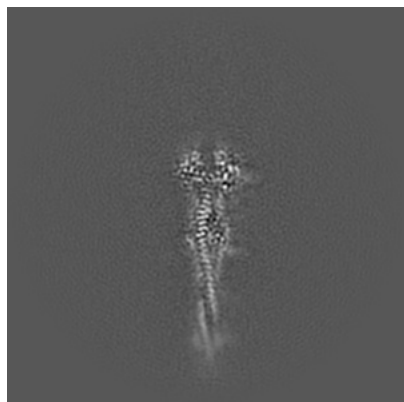


Z Index: 170

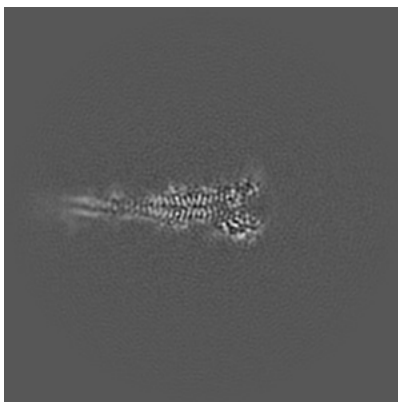
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

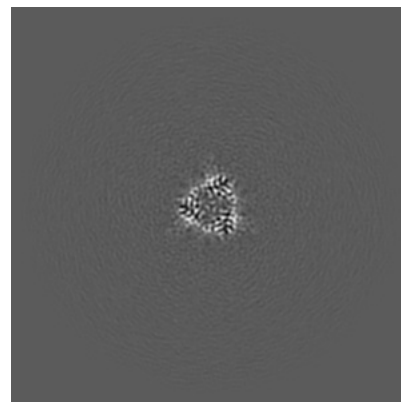
6.3.1 Primary map



X Index: 175

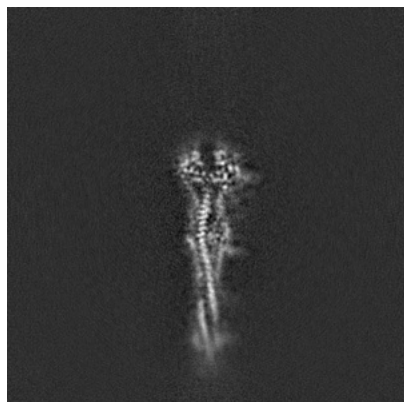


Y Index: 167

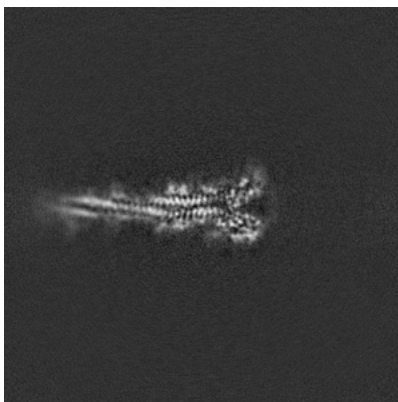


Z Index: 201

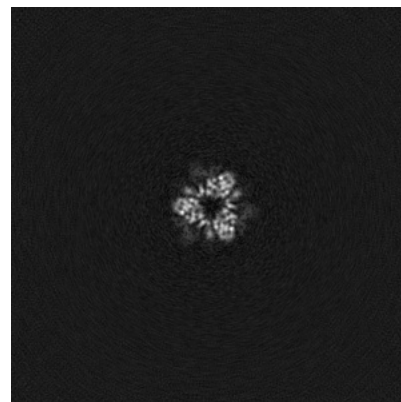
6.3.2 Raw map



X Index: 175



Y Index: 166

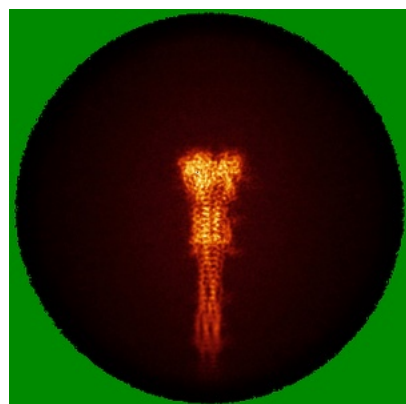


Z Index: 209

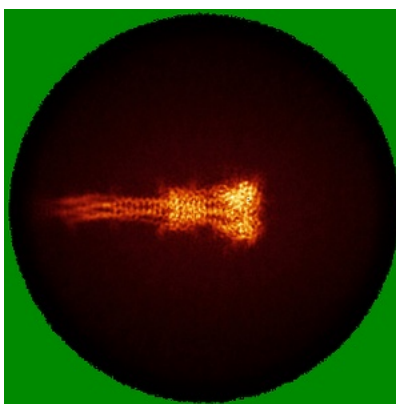
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) ⓘ

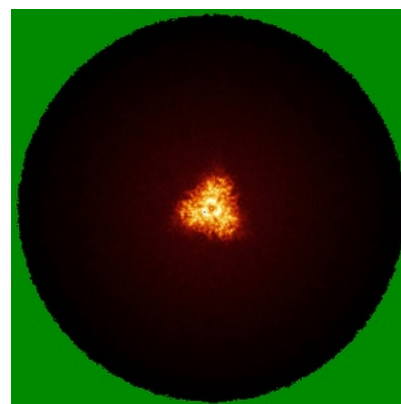
6.4.1 Primary map



X

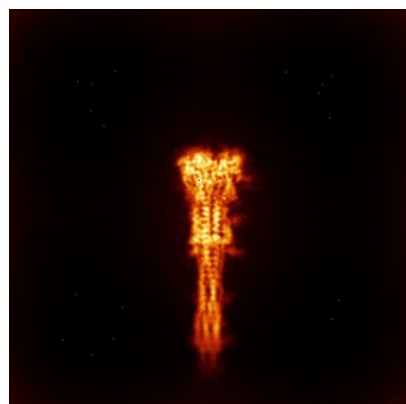


Y

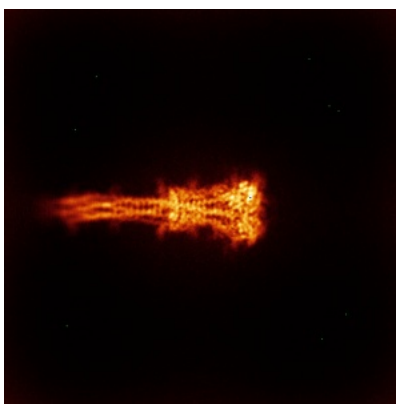


Z

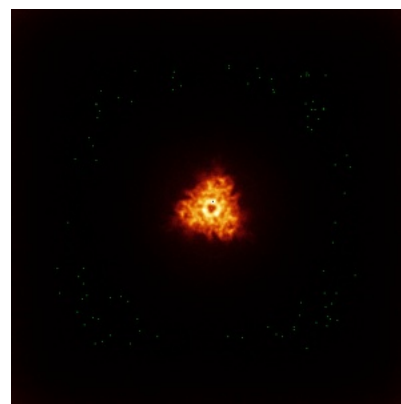
6.4.2 Raw map



X



Y

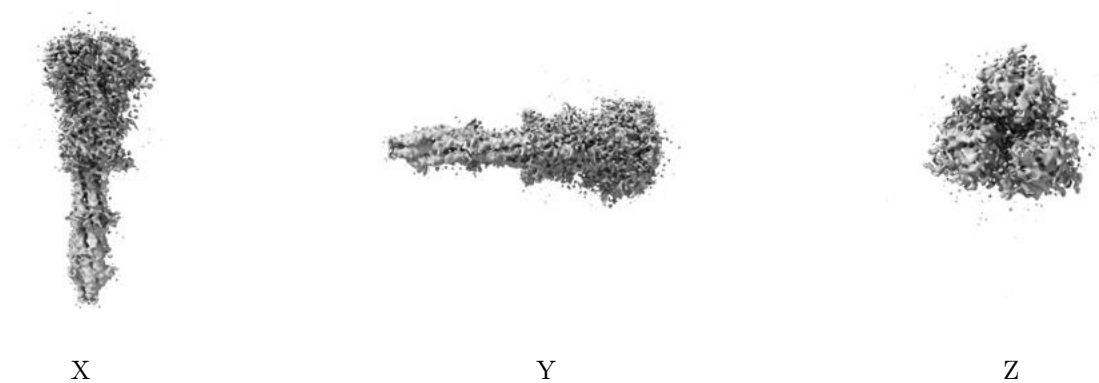


Z

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

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

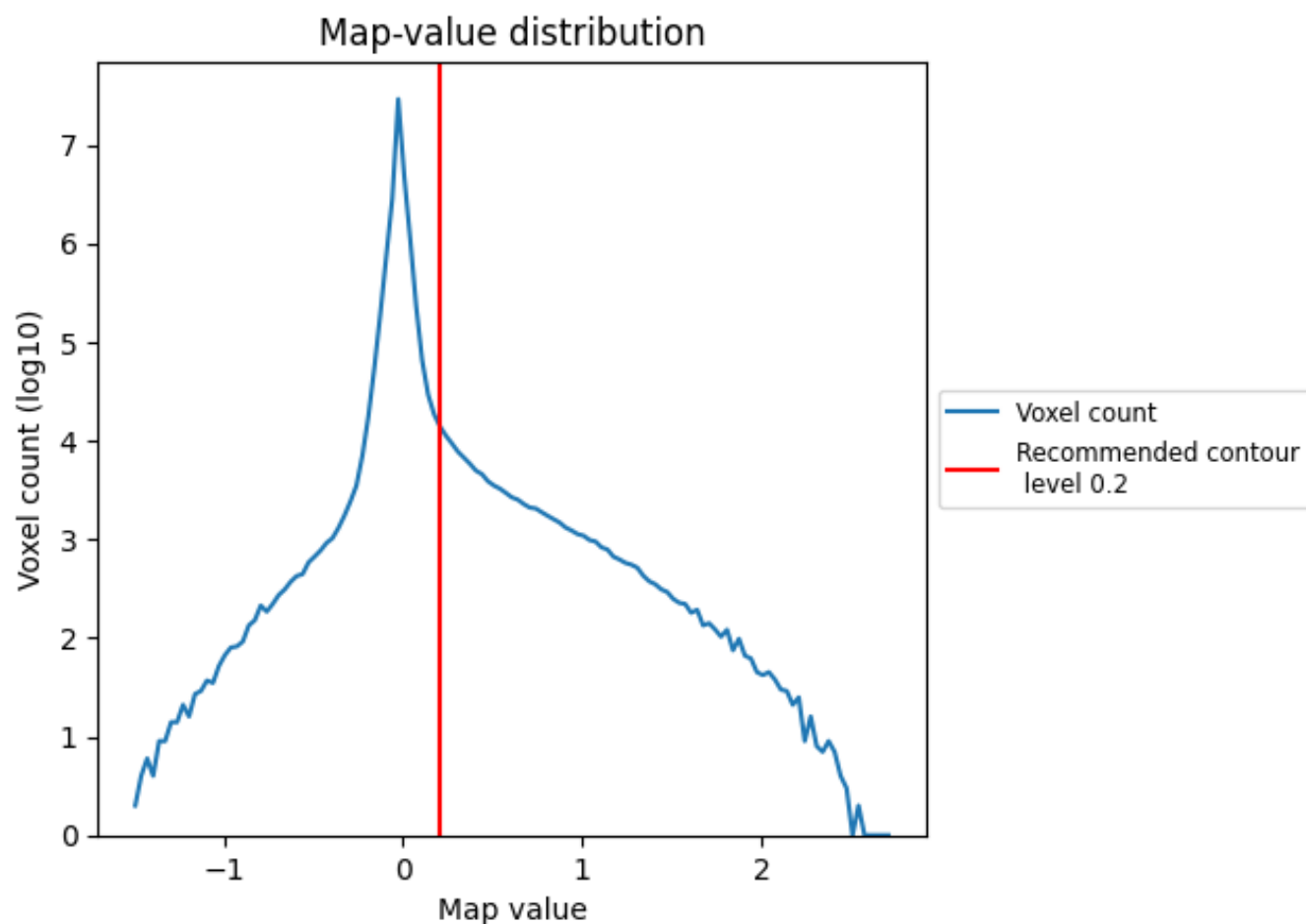
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

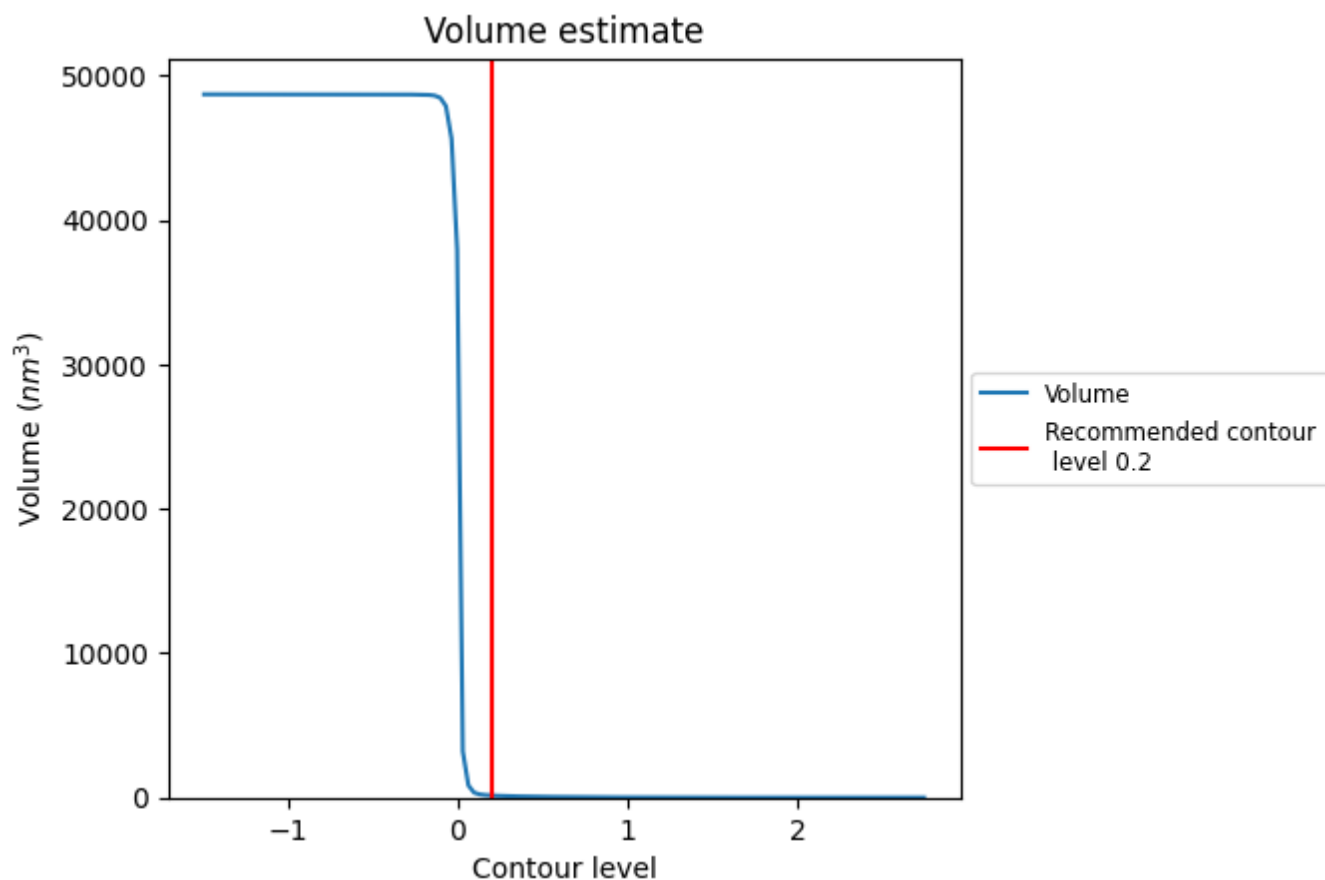
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

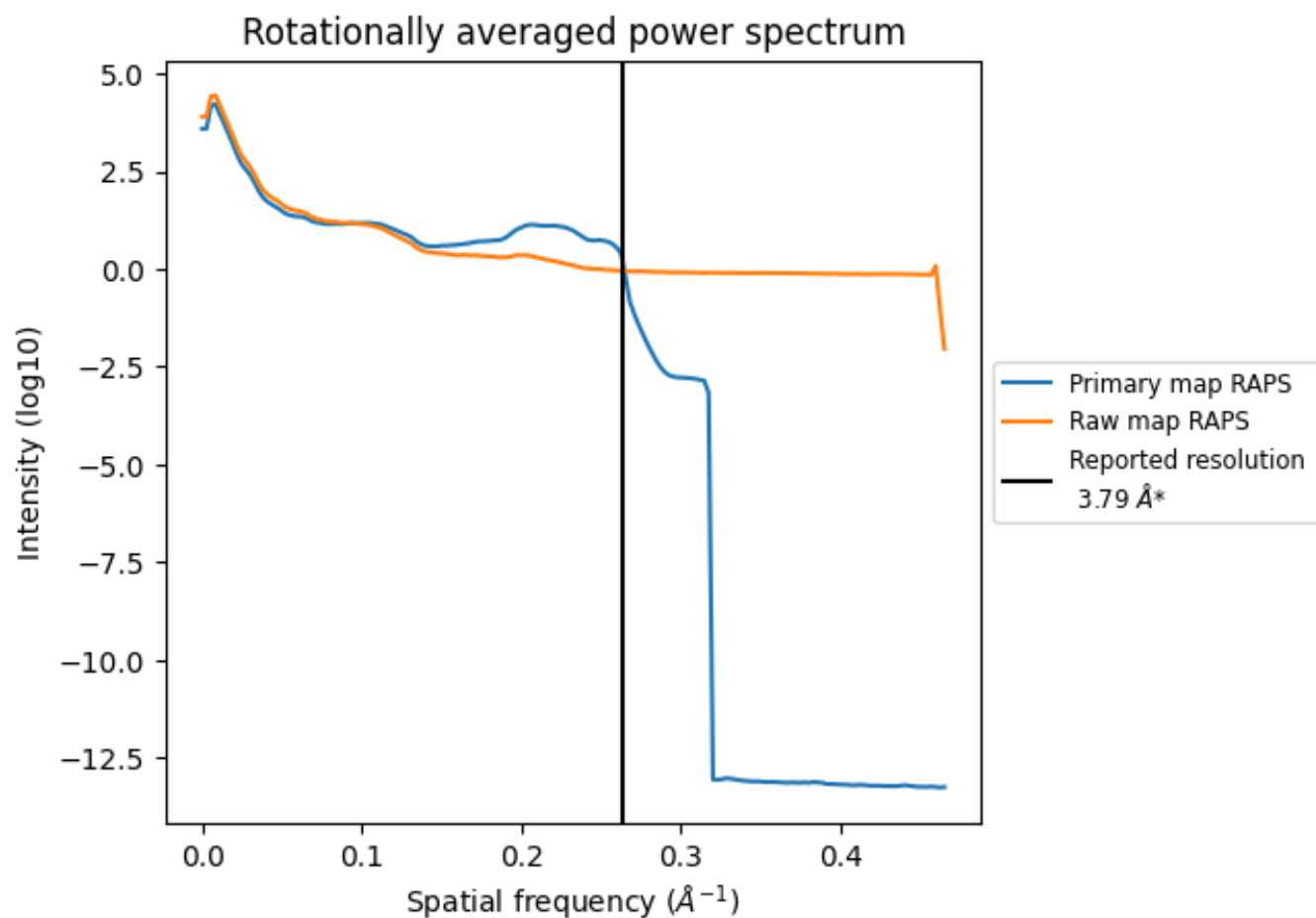
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 143 nm³; this corresponds to an approximate mass of 129 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 ⓘ

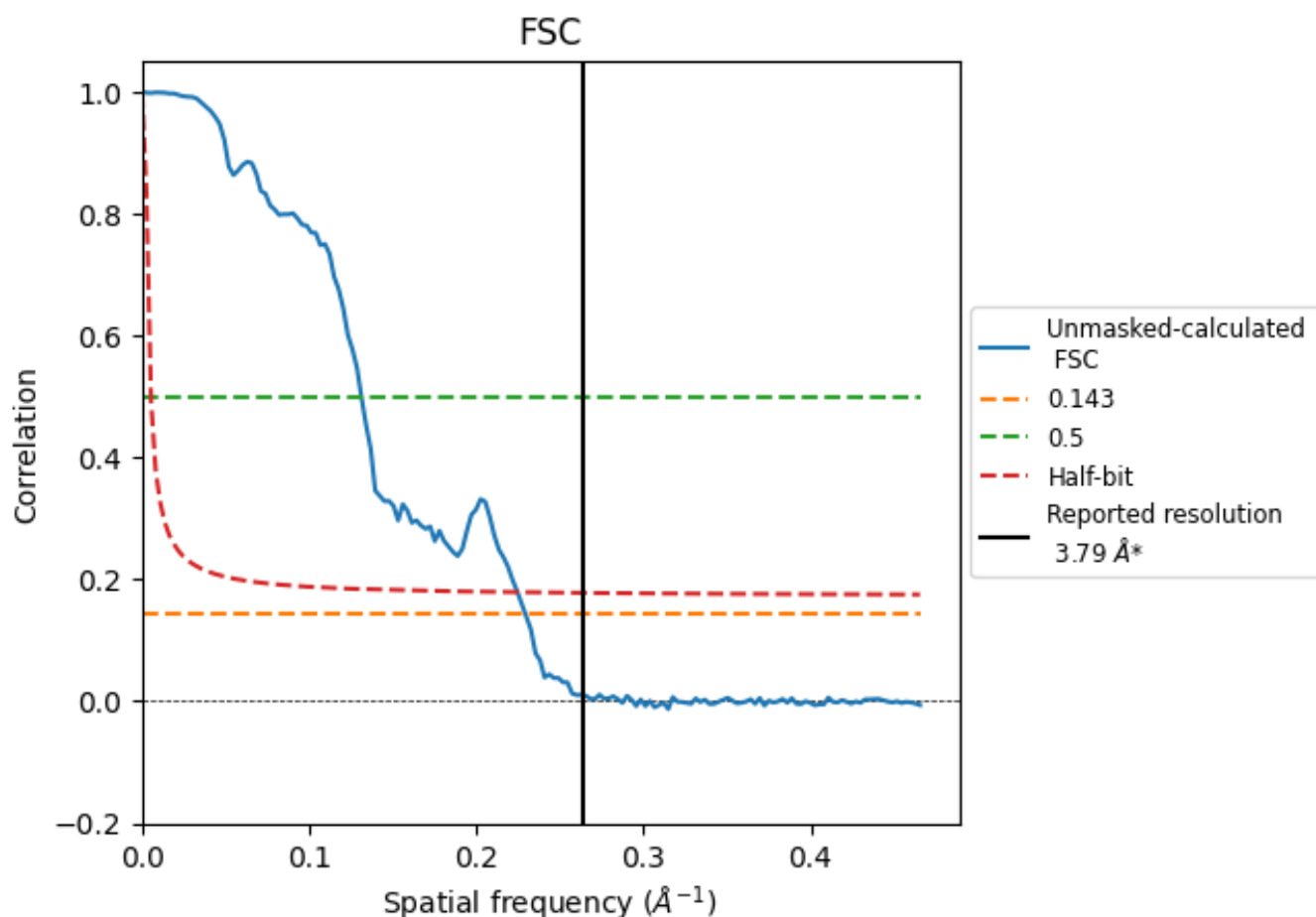


*Reported resolution corresponds to spatial frequency of 0.264 Å⁻¹

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 0.264 \AA^{-1}

8.2 Resolution estimates [i](#)

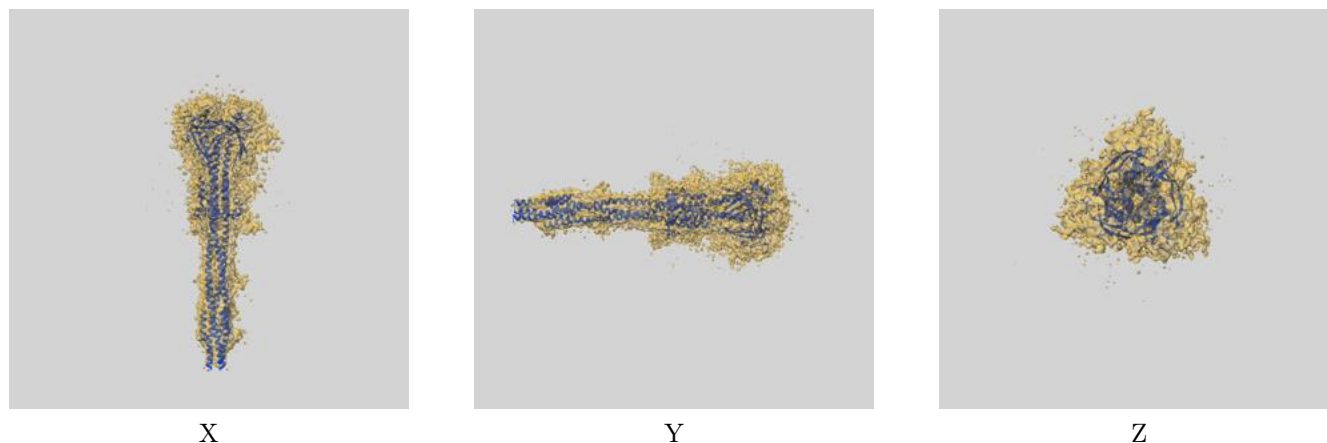
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.79	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.36	7.62	4.45

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.36 differs from the reported value 3.79 by more than 10 %

9 Map-model fit [i](#)

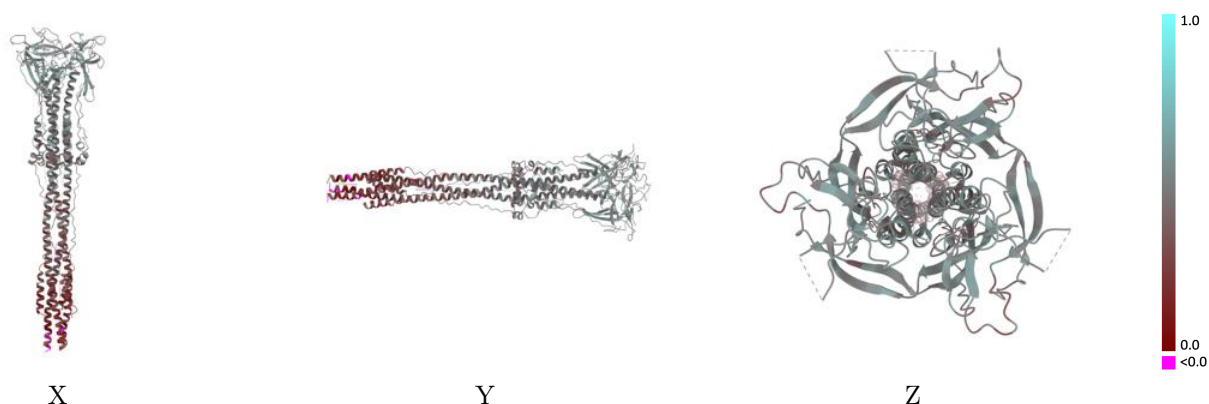
This section contains information regarding the fit between EMDB map EMD-39768 and PDB model 8Z4O. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



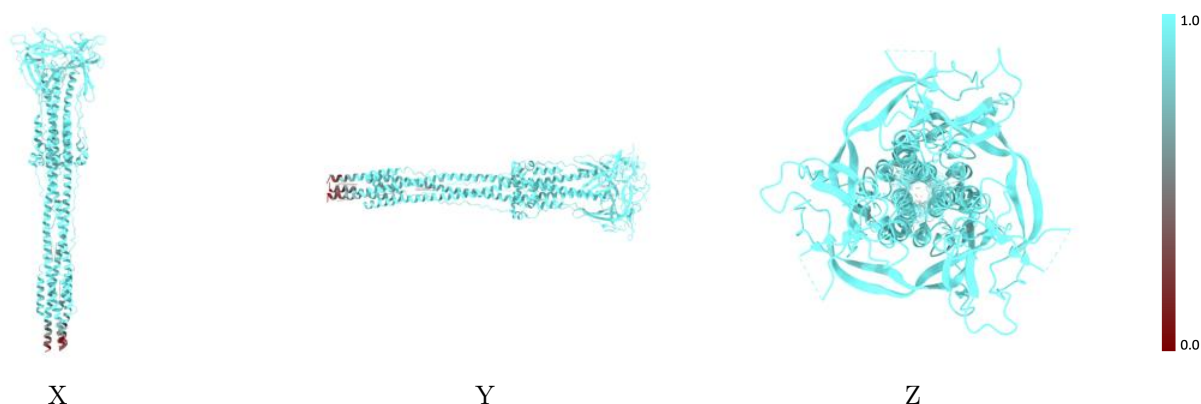
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



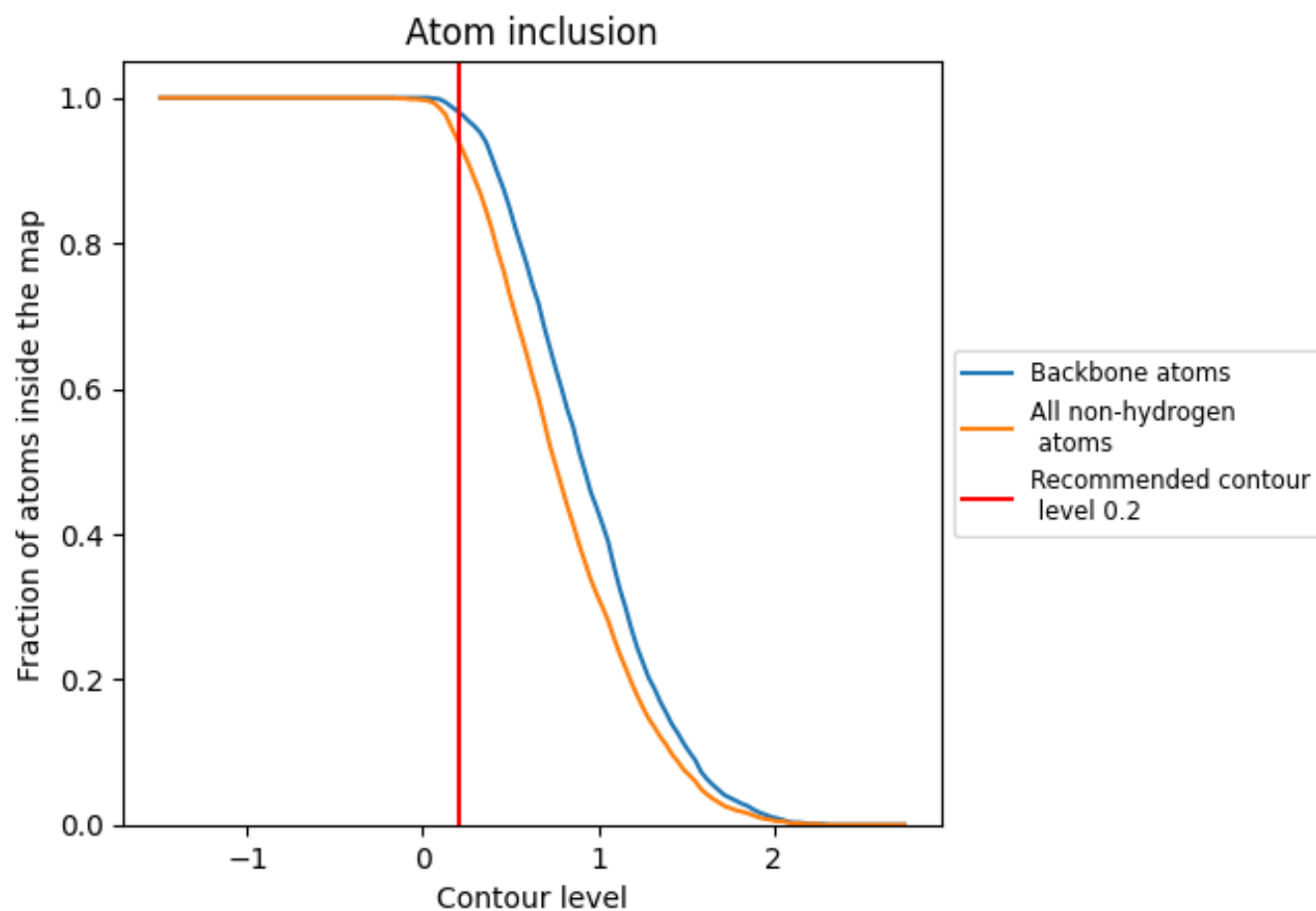
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



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 (0.2).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 94% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.9410	<div><div></div></div> 0.4260
A	<div><div></div></div> 0.9330	<div><div></div></div> 0.4250
B	<div><div></div></div> 0.9460	<div><div></div></div> 0.4240
C	<div><div></div></div> 0.9440	<div><div></div></div> 0.4300

