



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 05:41 PM UTC

PDB ID : 8FU9 / pdb_00008fu9
EMDB ID : EMD-29456
Title : Structure of Covid Spike variant deltaN25 with one erect RBD
Authors : Yu, X.; Juraszek, J.; Rutten, L.; Bakkers, M.J.G.; Blokland, S.; Van den Broek, N.J.F.; Verwilligen, A.Y.W.; Abeywickrema, P.; Vingerhoets, J.; Neefs, J.; Bakhsh, S.A.M.; Roychoudhury, P.; Greninger, A.; Sharma, S.; Langedijk, J.P.M.
Deposited on : 2023-01-16
Resolution : 3.52 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

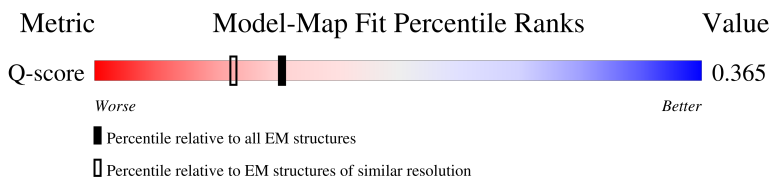
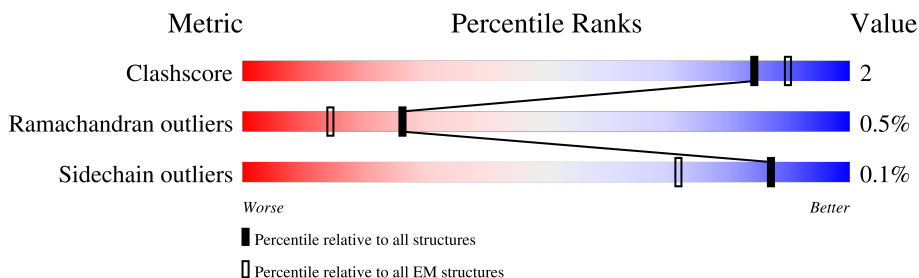
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.52 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13017 (3.02 - 4.02)

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	1188	
1	B	1188	
1	C	1188	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 25313 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 glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1058	Total	C	N	O	S	0	0
			8265	5264	1383	1580	38		
1	B	1057	Total	C	N	O	S	0	0
			8262	5265	1383	1576	38		
1	C	1056	Total	C	N	O	S	0	0
			8254	5258	1381	1577	38		

There are 93 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	TRP	deletion	UNP P0DTC2
A	?	-	PHE	deletion	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	ALA	deletion	UNP P0DTC2
A	?	-	ILE	deletion	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	?	-	SER	deletion	UNP P0DTC2
A	?	-	GLY	deletion	UNP P0DTC2
A	?	-	THR	deletion	UNP P0DTC2
A	?	-	ASN	deletion	UNP P0DTC2
A	?	-	GLY	deletion	UNP P0DTC2
A	?	-	THR	deletion	UNP P0DTC2
A	?	-	ARG	deletion	UNP P0DTC2
A	?	-	SER	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	?	-	LEU	deletion	UNP P0DTC2
A	?	-	THR	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	?	-	GLY	deletion	UNP P0DTC2
A	233	ASN	ASP	engineered mutation	UNP P0DTC2
A	432	GLN	LEU	engineered mutation	UNP P0DTC2
A	451	GLN	GLU	conflict	UNP P0DTC2
A	470	SER	PHE	engineered mutation	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	594	ASN	ASP	engineered mutation	UNP P0DTC2
A	662	SER	ARG	conflict	UNP P0DTC2
A	665	GLY	ARG	conflict	UNP P0DTC2
A	839	ASN	THR	engineered mutation	UNP P0DTC2
A	872	PRO	ALA	engineered mutation	UNP P0DTC2
A	922	PRO	ALA	engineered mutation	UNP P0DTC2
A	967	PRO	VAL	engineered mutation	UNP P0DTC2
B	?	-	TRP	deletion	UNP P0DTC2
B	?	-	PHE	deletion	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	ALA	deletion	UNP P0DTC2
B	?	-	ILE	deletion	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	?	-	SER	deletion	UNP P0DTC2
B	?	-	GLY	deletion	UNP P0DTC2
B	?	-	THR	deletion	UNP P0DTC2
B	?	-	ASN	deletion	UNP P0DTC2
B	?	-	GLY	deletion	UNP P0DTC2
B	?	-	THR	deletion	UNP P0DTC2
B	?	-	ARG	deletion	UNP P0DTC2
B	?	-	SER	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	?	-	LEU	deletion	UNP P0DTC2
B	?	-	THR	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	?	-	GLY	deletion	UNP P0DTC2
B	233	ASN	ASP	engineered mutation	UNP P0DTC2
B	432	GLN	LEU	engineered mutation	UNP P0DTC2
B	451	GLN	GLU	conflict	UNP P0DTC2
B	470	SER	PHE	engineered mutation	UNP P0DTC2
B	594	ASN	ASP	engineered mutation	UNP P0DTC2
B	662	SER	ARG	conflict	UNP P0DTC2
B	665	GLY	ARG	conflict	UNP P0DTC2
B	839	ASN	THR	engineered mutation	UNP P0DTC2
B	872	PRO	ALA	engineered mutation	UNP P0DTC2
B	922	PRO	ALA	engineered mutation	UNP P0DTC2
B	967	PRO	VAL	engineered mutation	UNP P0DTC2
C	?	-	TRP	deletion	UNP P0DTC2
C	?	-	PHE	deletion	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	ALA	deletion	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	ILE	deletion	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	?	-	SER	deletion	UNP P0DTC2
C	?	-	GLY	deletion	UNP P0DTC2
C	?	-	THR	deletion	UNP P0DTC2
C	?	-	ASN	deletion	UNP P0DTC2
C	?	-	GLY	deletion	UNP P0DTC2
C	?	-	THR	deletion	UNP P0DTC2
C	?	-	ARG	deletion	UNP P0DTC2
C	?	-	SER	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	?	-	LEU	deletion	UNP P0DTC2
C	?	-	THR	deletion	UNP P0DTC2
C	?	-	PRO	deletion	UNP P0DTC2
C	?	-	GLY	deletion	UNP P0DTC2
C	233	ASN	ASP	engineered mutation	UNP P0DTC2
C	432	GLN	LEU	engineered mutation	UNP P0DTC2
C	451	GLN	GLU	conflict	UNP P0DTC2
C	470	SER	PHE	engineered mutation	UNP P0DTC2
C	594	ASN	ASP	engineered mutation	UNP P0DTC2
C	662	SER	ARG	conflict	UNP P0DTC2
C	665	GLY	ARG	conflict	UNP P0DTC2
C	839	ASN	THR	engineered mutation	UNP P0DTC2
C	872	PRO	ALA	engineered mutation	UNP P0DTC2
C	922	PRO	ALA	engineered mutation	UNP P0DTC2
C	967	PRO	VAL	engineered mutation	UNP P0DTC2

- 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	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	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	

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Mol	Chain	Residues	Atoms				AltConf
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	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	
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	

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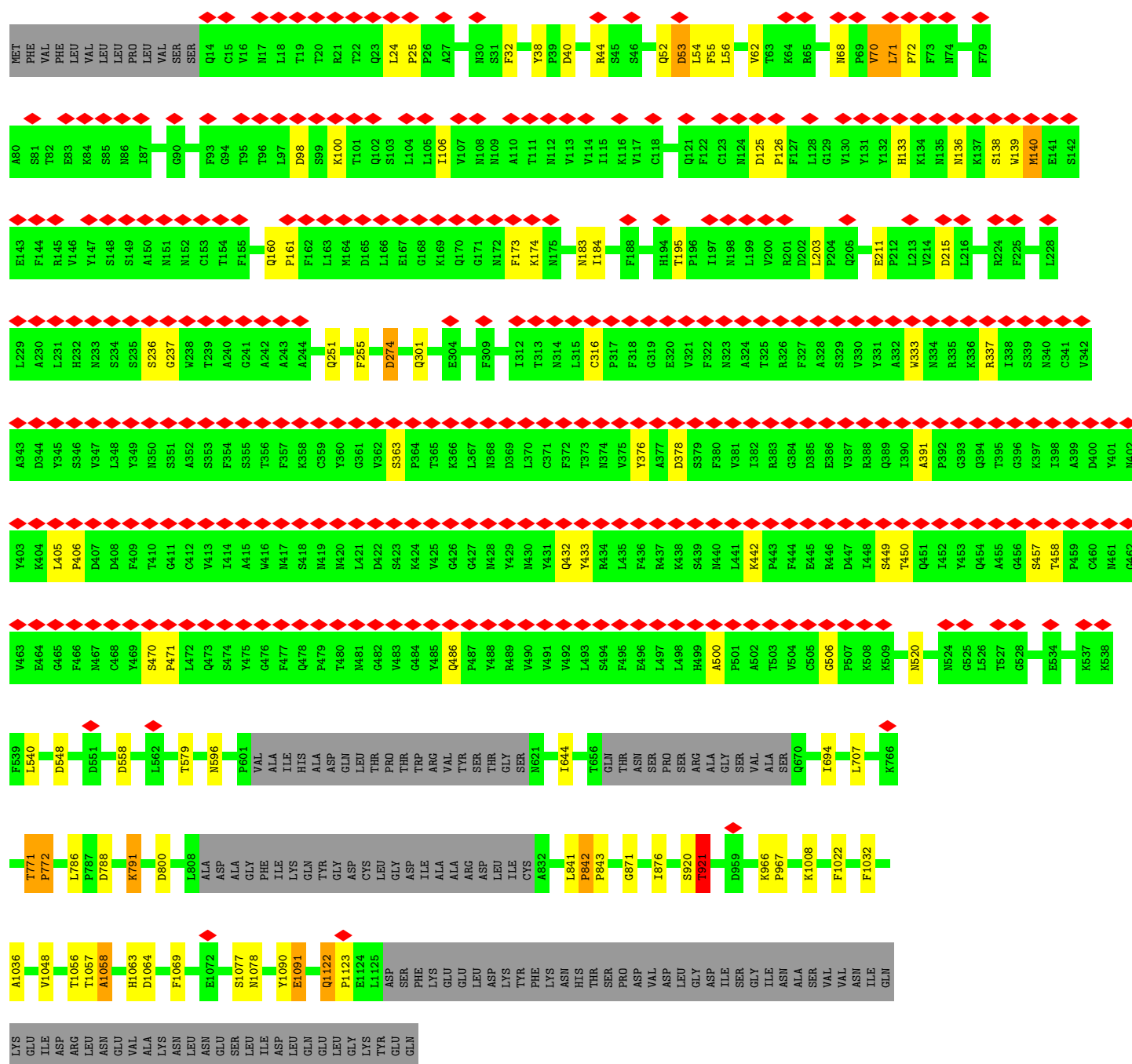
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Mol	Chain	Residues	Atoms				AltConf
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	

GLY
LYS
TYR
GLU
GLN

• Molecule 1: Spike glycoprotein

Chain B: 



• Molecule 1: Spike glycoprotein

Chain C: 



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	207525	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 ($e^-/\text{\AA}^2$)	52	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.064	Depositor
Minimum map value	-0.033	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.007	Depositor
Map size ($\text{\AA}$)	366.168, 366.168, 366.168	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.8322, 0.8322, 0.8322	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	0.63	1/8455 (0.0%)	1.03	89/11507 (0.8%)
1	B	0.61	0/8452	0.97	91/11503 (0.8%)
1	C	0.62	0/8444	0.98	90/11492 (0.8%)
All	All	0.62	1/25351 (0.0%)	0.99	270/34502 (0.8%)

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	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	689	ASN	CB-CG	8.40	1.73	1.52

All (270) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	689	ASN	CB-CG-OD1	-23.03	74.74	120.80
1	A	689	ASN	CB-CG-ND2	-20.77	85.25	116.40
1	A	689	ASN	OD1-CG-ND2	9.21	131.81	122.60
1	A	24	LEU	CA-C-N	9.13	126.23	119.66
1	A	24	LEU	C-N-CA	9.13	126.23	119.66
1	A	211	GLU	CA-C-N	8.45	128.39	119.85
1	A	211	GLU	C-N-CA	8.45	128.39	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	24	LEU	CA-C-N	8.24	125.68	119.66
1	B	24	LEU	C-N-CA	8.24	125.68	119.66
1	B	841	LEU	CA-C-N	8.03	125.44	119.66
1	B	841	LEU	C-N-CA	8.03	125.44	119.66
1	B	506	GLY	CA-C-N	7.73	127.66	119.85
1	B	506	GLY	C-N-CA	7.73	127.66	119.85
1	A	38	TYR	CA-C-N	7.69	127.70	119.78
1	A	38	TYR	C-N-CA	7.69	127.70	119.78
1	A	195	THR	CA-C-N	7.68	127.44	119.76
1	A	195	THR	C-N-CA	7.68	127.44	119.76
1	B	195	THR	CA-C-N	7.67	127.68	119.78
1	B	195	THR	C-N-CA	7.67	127.68	119.78
1	C	38	TYR	CA-C-N	7.58	127.48	120.21
1	C	38	TYR	C-N-CA	7.58	127.48	120.21
1	C	478	GLN	CA-C-N	7.53	127.23	119.56
1	C	478	GLN	C-N-CA	7.53	127.23	119.56
1	C	842	PRO	CA-C-N	7.49	127.44	120.03
1	C	842	PRO	C-N-CA	7.49	127.44	120.03
1	B	38	TYR	CA-C-N	7.45	127.46	119.78
1	B	38	TYR	C-N-CA	7.45	127.46	119.78
1	A	772	PRO	CA-C-N	7.43	127.06	119.56
1	A	772	PRO	C-N-CA	7.43	127.06	119.56
1	C	195	THR	CA-C-N	7.41	127.37	120.03
1	C	195	THR	C-N-CA	7.41	127.37	120.03
1	A	842	PRO	CA-C-N	7.38	127.31	119.85
1	A	842	PRO	C-N-CA	7.38	127.31	119.85
1	C	25	PRO	CA-C-N	7.37	127.53	120.31
1	C	25	PRO	C-N-CA	7.37	127.53	120.31
1	C	211	GLU	CA-C-N	7.32	127.08	119.76
1	C	211	GLU	C-N-CA	7.32	127.08	119.76
1	B	68	ASN	CA-C-N	7.29	127.73	119.93
1	B	68	ASN	C-N-CA	7.29	127.73	119.93
1	B	842	PRO	CA-C-N	7.21	127.21	119.78
1	B	842	PRO	C-N-CA	7.21	127.21	119.78
1	B	772	PRO	CA-C-N	7.19	126.82	119.56
1	B	772	PRO	C-N-CA	7.19	126.82	119.56
1	A	1119	ASP	CA-C-N	7.18	127.34	119.87
1	A	1119	ASP	C-N-CA	7.18	127.34	119.87
1	B	921	THR	CA-C-N	7.16	127.20	119.90
1	B	921	THR	C-N-CA	7.16	127.20	119.90
1	A	921	THR	CA-C-N	7.12	126.80	119.82
1	A	921	THR	C-N-CA	7.12	126.80	119.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	301	GLN	CA-C-N	7.08	127.05	119.76
1	C	301	GLN	C-N-CA	7.08	127.05	119.76
1	B	56	LEU	CA-C-N	7.03	127.00	119.76
1	B	56	LEU	C-N-CA	7.03	127.00	119.76
1	A	68	ASN	CA-C-N	7.02	126.94	119.85
1	A	68	ASN	C-N-CA	7.02	126.94	119.85
1	C	772	PRO	CA-C-N	7.01	126.64	119.56
1	C	772	PRO	C-N-CA	7.01	126.64	119.56
1	A	71	LEU	CA-C-N	6.98	126.69	119.64
1	A	71	LEU	C-N-CA	6.98	126.69	119.64
1	A	1058	ALA	CA-C-N	6.91	126.54	119.56
1	A	1058	ALA	C-N-CA	6.91	126.54	119.56
1	C	203	LEU	CA-C-N	6.88	126.80	119.85
1	C	203	LEU	C-N-CA	6.88	126.80	119.85
1	B	788	ASP	CA-C-N	6.87	126.50	119.56
1	B	788	ASP	C-N-CA	6.87	126.50	119.56
1	C	500	ALA	CA-C-N	6.84	126.88	119.90
1	C	500	ALA	C-N-CA	6.84	126.88	119.90
1	C	1058	ALA	CA-C-N	6.84	126.47	119.56
1	C	1058	ALA	C-N-CA	6.84	126.47	119.56
1	C	70	VAL	CA-C-N	6.80	133.95	121.70
1	C	70	VAL	C-N-CA	6.80	133.95	121.70
1	C	558	ASP	CA-C-N	6.79	126.49	119.56
1	C	558	ASP	C-N-CA	6.79	126.49	119.56
1	A	558	ASP	CA-C-N	6.76	126.46	119.56
1	A	558	ASP	C-N-CA	6.76	126.46	119.56
1	A	791	LYS	CA-C-N	6.72	126.35	119.56
1	A	791	LYS	C-N-CA	6.72	126.35	119.56
1	A	363	SER	CA-C-N	6.67	126.37	119.56
1	A	363	SER	C-N-CA	6.67	126.37	119.56
1	B	70	VAL	CA-C-N	6.64	133.65	121.70
1	B	70	VAL	C-N-CA	6.64	133.65	121.70
1	B	25	PRO	CA-C-N	6.62	126.60	119.78
1	B	25	PRO	C-N-CA	6.62	126.60	119.78
1	A	70	VAL	CA-C-N	6.60	133.58	121.70
1	A	70	VAL	C-N-CA	6.60	133.58	121.70
1	B	274	ASP	CA-C-N	6.57	127.09	119.47
1	B	274	ASP	C-N-CA	6.57	127.09	119.47
1	B	791	LYS	CA-C-N	6.52	126.15	119.56
1	B	791	LYS	C-N-CA	6.52	126.15	119.56
1	A	203	LEU	CA-C-N	6.50	126.26	119.76
1	A	203	LEU	C-N-CA	6.50	126.26	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1058	ALA	CA-C-N	6.48	126.11	119.56
1	B	1058	ALA	C-N-CA	6.48	126.11	119.56
1	C	405	LEU	CA-C-N	6.45	126.41	120.03
1	C	405	LEU	C-N-CA	6.45	126.41	120.03
1	C	106	ILE	N-CA-C	6.45	117.14	108.11
1	B	363	SER	CA-C-N	6.43	126.56	119.87
1	B	363	SER	C-N-CA	6.43	126.56	119.87
1	A	301	GLN	CA-C-N	6.42	126.33	119.85
1	A	301	GLN	C-N-CA	6.42	126.33	119.85
1	B	707	LEU	CA-C-N	6.41	126.36	119.76
1	B	707	LEU	C-N-CA	6.41	126.36	119.76
1	A	405	LEU	CA-C-N	6.40	126.77	119.92
1	A	405	LEU	C-N-CA	6.40	126.77	119.92
1	A	788	ASP	CA-C-N	6.40	126.52	119.87
1	A	788	ASP	C-N-CA	6.40	126.52	119.87
1	B	71	LEU	CA-C-N	6.37	126.08	119.64
1	B	71	LEU	C-N-CA	6.37	126.08	119.64
1	C	309	PHE	CA-C-N	6.36	126.37	119.89
1	C	309	PHE	C-N-CA	6.36	126.37	119.89
1	C	1048	VAL	CA-C-N	6.36	126.31	119.76
1	C	1048	VAL	C-N-CA	6.36	126.31	119.76
1	C	68	ASN	CA-C-N	6.32	126.08	119.76
1	C	68	ASN	C-N-CA	6.32	126.08	119.76
1	B	1091	GLU	CA-C-N	6.29	126.80	120.14
1	B	1091	GLU	C-N-CA	6.29	126.80	120.14
1	C	56	LEU	CA-C-N	6.25	126.28	119.90
1	C	56	LEU	C-N-CA	6.25	126.28	119.90
1	B	203	LEU	CA-C-N	6.24	126.21	119.78
1	B	203	LEU	C-N-CA	6.24	126.21	119.78
1	B	442	LYS	CA-C-N	6.20	126.17	119.78
1	B	442	LYS	C-N-CA	6.20	126.17	119.78
1	B	644	ILE	CA-C-N	6.18	125.94	119.76
1	B	644	ILE	C-N-CA	6.18	125.94	119.76
1	C	363	SER	CA-C-N	6.17	125.79	119.56
1	C	363	SER	C-N-CA	6.17	125.79	119.56
1	C	707	LEU	CA-C-N	6.14	126.08	119.76
1	C	707	LEU	C-N-CA	6.14	126.08	119.76
1	B	486	GLN	CA-C-N	6.13	126.05	119.85
1	B	486	GLN	C-N-CA	6.13	126.05	119.85
1	B	558	ASP	CA-C-N	6.13	125.81	119.56
1	B	558	ASP	C-N-CA	6.13	125.81	119.56
1	C	1091	GLU	CA-C-N	6.12	126.32	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1091	GLU	C-N-CA	6.12	126.32	119.90
1	A	707	LEU	CA-C-N	6.11	126.06	119.76
1	A	707	LEU	C-N-CA	6.11	126.06	119.76
1	B	391	ALA	CA-C-N	6.11	126.29	119.32
1	B	391	ALA	C-N-CA	6.11	126.29	119.32
1	A	1091	GLU	CA-C-N	6.08	127.44	119.84
1	A	1091	GLU	C-N-CA	6.08	127.44	119.84
1	C	644	ILE	CA-C-N	6.07	126.39	119.83
1	C	644	ILE	C-N-CA	6.07	126.39	119.83
1	C	694	ILE	CA-C-N	6.07	126.03	119.78
1	C	694	ILE	C-N-CA	6.07	126.03	119.78
1	A	786	LEU	CA-C-N	6.04	125.98	119.76
1	A	786	LEU	C-N-CA	6.04	125.98	119.76
1	B	871	GLY	CA-C-N	6.02	126.52	120.14
1	B	871	GLY	C-N-CA	6.02	126.52	120.14
1	B	579	THR	CA-C-N	6.02	126.48	120.52
1	B	579	THR	C-N-CA	6.02	126.48	120.52
1	A	500	ALA	CA-C-N	5.99	126.01	119.90
1	A	500	ALA	C-N-CA	5.99	126.01	119.90
1	C	274	ASP	CA-C-N	5.98	126.41	119.47
1	C	274	ASP	C-N-CA	5.98	126.41	119.47
1	C	921	THR	CA-C-N	5.96	126.38	119.47
1	C	921	THR	C-N-CA	5.96	126.38	119.47
1	A	274	ASP	CA-C-N	5.92	127.25	119.84
1	A	274	ASP	C-N-CA	5.92	127.25	119.84
1	A	871	GLY	CA-C-N	5.92	125.83	119.85
1	A	871	GLY	C-N-CA	5.92	125.83	119.85
1	C	391	ALA	CA-C-N	5.92	125.83	119.85
1	C	391	ALA	C-N-CA	5.92	125.83	119.85
1	B	316	CYS	CA-C-N	5.91	125.60	119.64
1	B	316	CYS	C-N-CA	5.91	125.60	119.64
1	B	1048	VAL	CA-C-N	5.85	125.76	119.85
1	B	1048	VAL	C-N-CA	5.85	125.76	119.85
1	A	644	ILE	CA-C-N	5.83	126.02	119.90
1	A	644	ILE	C-N-CA	5.83	126.02	119.90
1	B	211	GLU	CA-C-N	5.82	125.58	119.76
1	B	211	GLU	C-N-CA	5.82	125.58	119.76
1	C	1054	ASN	CA-CB-CG	5.82	118.42	112.60
1	A	1048	VAL	CA-C-N	5.80	125.71	119.85
1	A	1048	VAL	C-N-CA	5.80	125.71	119.85
1	C	871	GLY	CA-C-N	5.80	125.70	119.85
1	C	871	GLY	C-N-CA	5.80	125.70	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	251	GLN	CA-C-N	5.79	127.08	119.84
1	B	251	GLN	C-N-CA	5.79	127.08	119.84
1	B	301	GLN	CA-C-N	5.78	126.12	119.93
1	B	301	GLN	C-N-CA	5.78	126.12	119.93
1	A	694	ILE	CA-C-N	5.76	125.70	119.76
1	A	694	ILE	C-N-CA	5.76	125.70	119.76
1	B	500	ALA	CA-C-N	5.74	125.69	119.78
1	B	500	ALA	C-N-CA	5.74	125.69	119.78
1	C	24	LEU	CA-C-N	5.73	126.28	120.38
1	C	24	LEU	C-N-CA	5.73	126.28	120.38
1	A	771	THR	CA-C-N	5.71	126.26	120.38
1	A	771	THR	C-N-CA	5.71	126.26	120.38
1	A	540	LEU	CA-C-N	5.67	126.93	119.84
1	A	540	LEU	C-N-CA	5.67	126.93	119.84
1	B	694	ILE	CA-C-N	5.66	125.42	119.76
1	B	694	ILE	C-N-CA	5.66	125.42	119.76
1	B	786	LEU	CA-C-N	5.66	125.59	119.76
1	B	786	LEU	C-N-CA	5.66	125.59	119.76
1	C	1069	PHE	CA-C-N	5.65	125.60	119.78
1	C	1069	PHE	C-N-CA	5.65	125.60	119.78
1	C	216	LEU	CA-C-N	5.64	126.89	119.84
1	C	216	LEU	C-N-CA	5.64	126.89	119.84
1	B	540	LEU	CA-C-N	5.59	126.82	119.84
1	B	540	LEU	C-N-CA	5.59	126.82	119.84
1	A	579	THR	CA-C-N	5.58	126.95	120.98
1	A	579	THR	C-N-CA	5.58	126.95	120.98
1	C	579	THR	CA-C-N	5.56	126.93	120.98
1	C	579	THR	C-N-CA	5.56	126.93	120.98
1	C	442	LYS	CA-C-N	5.56	125.88	119.93
1	C	442	LYS	C-N-CA	5.56	125.88	119.93
1	C	771	THR	CA-C-N	5.56	126.11	120.38
1	C	771	THR	C-N-CA	5.56	126.11	120.38
1	C	1122	GLN	CA-C-N	5.53	125.20	119.56
1	C	1122	GLN	C-N-CA	5.53	125.20	119.56
1	B	771	THR	CA-C-N	5.52	126.06	120.38
1	B	771	THR	C-N-CA	5.52	126.06	120.38
1	C	786	LEU	CA-C-N	5.49	125.50	119.90
1	C	786	LEU	C-N-CA	5.49	125.50	119.90
1	A	125	ASP	CA-C-N	5.49	125.39	119.85
1	A	125	ASP	C-N-CA	5.49	125.39	119.85
1	A	1036	ALA	CA-C-N	5.47	126.68	119.84
1	A	1036	ALA	C-N-CA	5.47	126.68	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1122	GLN	CA-C-N	5.46	125.13	119.56
1	B	1122	GLN	C-N-CA	5.46	125.13	119.56
1	B	458	THR	CA-C-N	5.45	125.77	119.93
1	B	458	THR	C-N-CA	5.45	125.77	119.93
1	A	841	LEU	CA-C-N	5.44	125.98	120.38
1	A	841	LEU	C-N-CA	5.44	125.98	120.38
1	C	568	THR	CA-C-N	5.43	125.33	119.85
1	C	568	THR	C-N-CA	5.43	125.33	119.85
1	A	309	PHE	CA-C-N	5.41	126.60	119.84
1	A	309	PHE	C-N-CA	5.41	126.60	119.84
1	C	251	GLN	CA-C-N	5.41	125.72	119.93
1	C	251	GLN	C-N-CA	5.41	125.72	119.93
1	A	1122	GLN	CA-C-N	5.40	125.02	119.56
1	A	1122	GLN	C-N-CA	5.40	125.02	119.56
1	C	841	LEU	CA-C-N	5.39	125.94	120.38
1	C	841	LEU	C-N-CA	5.39	125.94	120.38
1	B	1036	ALA	CA-C-N	5.39	126.58	119.84
1	B	1036	ALA	C-N-CA	5.39	126.58	119.84
1	B	876	ILE	CA-C-N	5.35	125.36	119.90
1	B	876	ILE	C-N-CA	5.35	125.36	119.90
1	B	405	LEU	CA-C-N	5.34	126.52	119.84
1	B	405	LEU	C-N-CA	5.34	126.52	119.84
1	A	478	GLN	CA-C-N	5.34	125.65	119.93
1	A	478	GLN	C-N-CA	5.34	125.65	119.93
1	A	1069	PHE	CA-C-N	5.34	125.10	119.76
1	A	1069	PHE	C-N-CA	5.34	125.10	119.76
1	A	160	GLN	CA-C-N	5.33	126.50	119.84
1	A	160	GLN	C-N-CA	5.33	126.50	119.84
1	A	966	LYS	CA-C-N	5.28	124.95	119.56
1	A	966	LYS	C-N-CA	5.28	124.95	119.56
1	A	32	PHE	CA-C-N	5.25	131.16	121.70
1	A	32	PHE	C-N-CA	5.25	131.16	121.70
1	C	788	ASP	CA-C-N	5.23	126.38	119.84
1	C	788	ASP	C-N-CA	5.23	126.38	119.84
1	B	106	ILE	N-CA-C	5.22	115.99	108.48
1	C	966	LYS	CA-C-N	5.21	124.83	119.56
1	C	966	LYS	C-N-CA	5.21	124.83	119.56
1	B	1032	PHE	CA-C-N	5.21	125.66	120.14
1	B	1032	PHE	C-N-CA	5.21	125.66	120.14
1	A	44	ARG	N-CA-C	5.21	117.36	111.11
1	C	32	PHE	CA-C-N	5.21	131.07	121.70
1	C	32	PHE	C-N-CA	5.21	131.07	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	540	LEU	CA-C-N	5.20	126.34	119.84
1	C	540	LEU	C-N-CA	5.20	126.34	119.84
1	C	1032	PHE	CA-C-N	5.17	125.62	120.14
1	C	1032	PHE	C-N-CA	5.17	125.62	120.14
1	A	442	LYS	CA-C-N	5.17	125.41	119.83
1	A	442	LYS	C-N-CA	5.17	125.41	119.83
1	A	106	ILE	N-CA-C	5.12	116.28	109.37
1	C	470	SER	CA-C-N	5.10	125.10	120.21
1	C	470	SER	C-N-CA	5.10	125.10	120.21
1	B	1069	PHE	CA-C-N	5.08	124.98	119.85
1	B	1069	PHE	C-N-CA	5.08	124.98	119.85

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	689	ASN	Sidechain
1	B	921	THR	Peptide

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	8265	0	8084	31	0
1	B	8262	0	8090	48	0
1	C	8254	0	8074	40	0
2	A	182	0	169	2	0
2	B	168	0	156	0	0
2	C	182	0	169	2	0
All	All	25313	0	24742	117	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 2.

All (117) 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:560:GLN:HB3	2:A:1213:NAG:H83	1.72	0.72
1:B:125:ASP:N	1:B:126:PRO:CD	2.65	0.60
1:B:138:SER:HG	1:B:139:TRP:CD1	2.20	0.59
1:B:337:ARG:HH12	1:C:155:PHE:H	1.50	0.59
1:A:466:PHE:CG	1:A:467:ASN:N	2.72	0.58
1:B:98:ASP:OD2	1:B:100:LYS:NZ	2.37	0.56
1:A:71:LEU:N	1:A:72:PRO:CD	2.68	0.56
1:C:71:LEU:N	1:C:72:PRO:HD3	2.20	0.56
1:C:33:THR:OG1	1:C:204:PRO:O	2.24	0.56
1:A:486:GLN:N	1:A:487:PRO:CD	2.69	0.56
1:B:920:SER:O	1:B:921:THR:C	2.49	0.55
1:B:40:ASP:N	1:B:40:ASP:OD1	2.37	0.55
1:C:173:PHE:H	1:C:174:LYS:HA	1.72	0.55
1:A:1103:SER:OG	1:A:1104:GLY:N	2.35	0.55
1:B:1090:TYR:O	1:B:1091:GLU:C	2.50	0.55
1:A:160:GLN:HB2	1:A:161:PRO:CD	2.35	0.54
1:A:183:ASN:HB3	1:A:184:ILE:HA	1.88	0.54
1:A:71:LEU:N	1:A:72:PRO:HD3	2.23	0.54
1:A:125:ASP:N	1:A:126:PRO:CD	2.70	0.54
1:B:71:LEU:N	1:B:72:PRO:CD	2.71	0.54
1:B:183:ASN:HB3	1:B:184:ILE:HA	1.90	0.54
1:C:71:LEU:N	1:C:72:PRO:CD	2.71	0.54
1:C:1122:GLN:N	1:C:1123:PRO:CD	2.71	0.54
1:C:173:PHE:N	1:C:174:LYS:HA	2.24	0.53
1:C:485:TYR:O	1:C:486:GLN:C	2.51	0.52
1:B:133:HIS:O	1:B:136:ASN:ND2	2.43	0.52
1:C:160:GLN:N	1:C:161:PRO:CD	2.73	0.52
1:B:52:GLN:O	1:B:53:ASP:CB	2.58	0.51
1:A:486:GLN:N	1:A:487:PRO:HD3	2.26	0.51
1:B:1063:HIS:CG	1:B:1064:ASP:H	2.29	0.50
1:B:160:GLN:N	1:B:161:PRO:CD	2.74	0.50
1:B:1063:HIS:CG	1:B:1064:ASP:N	2.80	0.50
1:C:270:ASP:OD1	1:C:271:CYS:N	2.44	0.50
1:C:183:ASN:CB	1:C:184:ILE:HA	2.42	0.48
1:B:71:LEU:N	1:B:72:PRO:HD3	2.29	0.48
1:B:1122:GLN:N	1:B:1123:PRO:CD	2.76	0.48
1:B:337:ARG:NH2	1:B:376:TYR:OH	2.46	0.48
1:B:125:ASP:N	1:B:126:PRO:HD2	2.28	0.48
1:C:470:SER:HB3	1:C:471:PRO:HD3	1.95	0.47
1:B:160:GLN:N	1:B:160:GLN:OE1	2.48	0.47
1:B:1008:LYS:NZ	1:B:1022:PHE:O	2.47	0.47
1:C:1103:SER:OG	1:C:1104:GLY:N	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:PHE:N	1:A:174:LYS:HA	2.29	0.47
1:B:139:TRP:O	1:B:140:MET:HB2	2.15	0.47
1:B:520:ASN:C	1:B:520:ASN:OD1	2.56	0.47
1:B:470:SER:N	1:B:471:PRO:HD2	2.28	0.47
1:C:183:ASN:HB2	1:C:184:ILE:HA	1.95	0.47
1:B:215:ASP:OD1	1:B:215:ASP:N	2.48	0.47
1:B:333:TRP:CD1	1:B:333:TRP:H	2.33	0.47
1:B:1057:THR:OG1	1:B:1058:ALA:N	2.48	0.47
1:C:125:ASP:N	1:C:126:PRO:CD	2.79	0.46
1:B:173:PHE:N	1:B:174:LYS:HA	2.31	0.46
1:B:236:SER:OG	1:B:237:GLY:N	2.50	0.45
1:B:432:GLN:O	1:B:433:TYR:C	2.59	0.45
1:C:920:SER:O	1:C:921:THR:C	2.58	0.45
1:A:33:THR:H	1:A:206:GLY:H	1.63	0.45
1:C:552:THR:O	1:C:553:THR:C	2.58	0.45
1:A:1122:GLN:N	1:A:1123:PRO:HD2	2.31	0.45
1:A:32:PHE:HA	1:A:33:THR:HB	1.99	0.45
1:C:1008:LYS:NZ	1:C:1022:PHE:O	2.49	0.45
1:A:449:SER:OG	1:A:450:THR:N	2.49	0.45
1:B:183:ASN:CB	1:B:184:ILE:HA	2.47	0.45
1:A:864:SER:O	1:A:881:GLN:NE2	2.49	0.45
1:C:847:ASP:O	1:C:848:GLU:HB2	2.17	0.45
1:C:84:LYS:NZ	1:C:168:GLY:O	2.47	0.45
1:C:647:GLY:O	1:C:649:GLY:N	2.50	0.45
1:C:864:SER:OG	1:C:867:THR:OG1	2.35	0.45
1:A:560:GLN:CB	2:A:1213:NAG:H83	2.43	0.44
1:C:949:ASN:OD1	1:C:955:SER:N	2.50	0.44
1:A:183:ASN:CB	1:A:184:ILE:HA	2.48	0.44
1:B:548:ASP:OD1	1:B:548:ASP:C	2.61	0.44
1:B:920:SER:OG	1:B:921:THR:N	2.51	0.43
1:A:966:LYS:N	1:A:967:PRO:HD2	2.34	0.43
1:A:888:GLY:O	1:A:1018:LYS:NZ	2.51	0.43
1:B:173:PHE:H	1:B:174:LYS:HA	1.83	0.43
1:B:966:LYS:HB3	1:B:967:PRO:HD3	2.00	0.43
1:C:846:THR:H	1:C:849:MET:HE3	1.84	0.43
1:A:553:THR:C	1:A:555:ALA:H	2.27	0.43
1:C:486:GLN:N	1:C:487:PRO:HD3	2.33	0.43
1:B:70:VAL:HA	1:B:71:LEU:HB2	2.00	0.43
1:C:133:HIS:ND1	1:C:134:LYS:N	2.63	0.43
1:B:842:PRO:HA	1:B:843:PRO:HD3	1.89	0.42
1:C:61:ASN:OD1	2:C:1201:NAG:N2	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:LEU:HD23	1:B:54:LEU:H	1.84	0.42
1:C:470:SER:HB3	1:C:471:PRO:CD	2.50	0.42
1:A:37:TYR:CD1	1:A:37:TYR:C	2.97	0.42
1:A:470:SER:N	1:A:471:PRO:HD2	2.33	0.42
1:B:791:LYS:NZ	1:B:800:ASP:OD2	2.40	0.42
1:A:139:TRP:O	1:A:140:MET:HB3	2.19	0.42
1:A:441:LEU:O	1:A:442:LYS:HB3	2.20	0.42
1:A:457:SER:OG	1:A:458:THR:N	2.52	0.42
1:B:966:LYS:HB3	1:B:967:PRO:CD	2.50	0.42
1:C:966:LYS:N	1:C:967:PRO:HD2	2.34	0.42
1:A:273:LEU:O	1:A:274:ASP:C	2.63	0.42
1:B:378:ASP:OD1	1:B:378:ASP:C	2.61	0.42
1:C:89:ARG:NH2	1:C:130:VAL:O	2.53	0.42
1:C:125:ASP:N	1:C:126:PRO:HD2	2.35	0.42
1:A:173:PHE:H	1:A:174:LYS:HA	1.85	0.42
1:B:470:SER:N	1:B:471:PRO:CD	2.82	0.42
1:C:307:VAL:H	1:C:511:THR:HG22	1.84	0.42
1:B:337:ARG:NH1	1:C:155:PHE:H	2.17	0.41
1:C:120:PHE:CZ	1:C:122:PHE:HB2	2.55	0.41
1:C:72:PRO:HG2	1:C:74:ASN:H	1.85	0.41
1:A:442:LYS:HA	1:A:443:PRO:HD3	1.93	0.41
1:A:474:SER:OG	1:A:475:TYR:N	2.53	0.41
1:B:55:PHE:HB2	1:B:255:PHE:CE1	2.55	0.41
1:A:1057:THR:OG1	1:A:1058:ALA:N	2.54	0.40
1:B:1056:THR:O	1:B:1077:SER:N	2.52	0.40
1:C:153:CYS:SG	1:C:154:THR:N	2.94	0.40
1:B:771:THR:HA	1:B:772:PRO:HD3	1.96	0.40
1:B:449:SER:OG	1:B:450:THR:N	2.54	0.40
1:C:269:VAL:HG23	1:C:286:PHE:CE2	2.57	0.40
1:B:32:PHE:CD2	1:B:62:VAL:CG2	3.04	0.40
1:B:160:GLN:N	1:B:161:PRO:HD2	2.36	0.40
1:C:61:ASN:OD1	2:C:1201:NAG:C2	2.69	0.40
1:C:351:SER:OG	1:C:352:ALA:N	2.54	0.40
1:C:441:LEU:O	1:C:442:LYS:C	2.65	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	1050/1188 (88%)	979 (93%)	68 (6%)	3 (0%)	36	66
1	B	1049/1188 (88%)	980 (93%)	63 (6%)	6 (1%)	21	55
1	C	1048/1188 (88%)	970 (93%)	72 (7%)	6 (1%)	21	55
All	All	3147/3564 (88%)	2929 (93%)	203 (6%)	15 (0%)	26	58

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	44	ARG
1	C	648	ALA
1	B	53	ASP
1	C	442	LYS
1	C	542	PHE
1	C	44	ARG
1	C	553	THR
1	A	138	SER
1	A	888	GLY
1	B	274	ASP
1	B	406	PRO
1	B	457	SER
1	C	552	THR
1	B	140	MET
1	A	442	LYS

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	927/1039 (89%)	926 (100%)	1 (0%)	88	87
1	B	926/1039 (89%)	924 (100%)	2 (0%)	87	85
1	C	926/1039 (89%)	925 (100%)	1 (0%)	88	87
All	All	2779/3117 (89%)	2775 (100%)	4 (0%)	87	87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1054	ASN
1	B	596	ASN
1	B	1078	ASN
1	C	1054	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	49	HIS
1	A	52	GLN
1	A	108	ASN
1	A	172	ASN
1	A	198	ASN
1	A	990	GLN
1	A	1038	HIS
1	B	49	HIS
1	B	86	ASN
1	B	121	GLN
1	B	175	ASN
1	B	226	GLN
1	B	368	ASN
1	B	586	ASN
1	B	893	GLN
1	B	991	GLN
1	C	102	GLN
1	C	124	ASN
1	C	560	GLN
1	C	586	ASN
1	C	670	GLN
1	C	1038	HIS
1	C	1088	ASN
1	C	1122	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

38 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	B	1208	1	14,14,15	1.33	1 (7%)	17,19,21	1.01	1 (5%)
2	NAG	C	1213	1	14,14,15	1.39	1 (7%)	17,19,21	1.07	1 (5%)
2	NAG	B	1211	1	14,14,15	1.35	2 (14%)	17,19,21	1.03	1 (5%)
2	NAG	C	1201	1	14,14,15	1.53	1 (7%)	17,19,21	1.16	1 (5%)
2	NAG	A	1213	1	14,14,15	1.53	2 (14%)	17,19,21	1.20	2 (11%)
2	NAG	A	1201	1	14,14,15	1.39	2 (14%)	17,19,21	1.05	1 (5%)
2	NAG	A	1205	1	14,14,15	1.31	2 (14%)	17,19,21	1.03	1 (5%)
2	NAG	B	1201	1	14,14,15	1.39	1 (7%)	17,19,21	0.84	0
2	NAG	C	1208	1	14,14,15	1.36	2 (14%)	17,19,21	1.01	1 (5%)
2	NAG	C	1209	1	14,14,15	1.35	2 (14%)	17,19,21	1.01	1 (5%)
2	NAG	C	1212	1	14,14,15	1.36	1 (7%)	17,19,21	1.06	1 (5%)
2	NAG	A	1208	1	14,14,15	1.37	2 (14%)	17,19,21	1.01	1 (5%)
2	NAG	A	1212	1	14,14,15	1.40	2 (14%)	17,19,21	0.98	1 (5%)
2	NAG	B	1210	1	14,14,15	1.29	2 (14%)	17,19,21	0.99	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	1207	1	14,14,15	1.47	2 (14%)	17,19,21	0.88	1 (5%)
2	NAG	B	1209	1	14,14,15	1.36	2 (14%)	17,19,21	1.04	1 (5%)
2	NAG	C	1206	1	14,14,15	1.34	1 (7%)	17,19,21	1.09	1 (5%)
2	NAG	A	1203	1	14,14,15	1.36	1 (7%)	17,19,21	1.08	1 (5%)
2	NAG	B	1202	1	14,14,15	1.36	2 (14%)	17,19,21	1.10	1 (5%)
2	NAG	C	1203	1	14,14,15	1.40	2 (14%)	17,19,21	1.07	1 (5%)
2	NAG	B	1204	1	14,14,15	1.38	1 (7%)	17,19,21	1.03	1 (5%)
2	NAG	A	1206	1	14,14,15	1.36	1 (7%)	17,19,21	1.07	1 (5%)
2	NAG	B	1206	1	14,14,15	1.25	2 (14%)	17,19,21	1.02	1 (5%)
2	NAG	A	1209	1	14,14,15	1.33	2 (14%)	17,19,21	1.02	1 (5%)
2	NAG	B	1205	1	14,14,15	1.39	2 (14%)	17,19,21	0.97	1 (5%)
2	NAG	A	1202	1	14,14,15	1.37	2 (14%)	17,19,21	1.03	1 (5%)
2	NAG	A	1211	1	14,14,15	1.48	2 (14%)	17,19,21	0.93	1 (5%)
2	NAG	B	1207	1	14,14,15	1.36	1 (7%)	17,19,21	1.14	1 (5%)
2	NAG	C	1211	1	14,14,15	1.33	2 (14%)	17,19,21	1.07	1 (5%)
2	NAG	A	1210	1	14,14,15	1.40	1 (7%)	17,19,21	1.00	1 (5%)
2	NAG	C	1202	1	14,14,15	1.41	2 (14%)	17,19,21	1.06	2 (11%)
2	NAG	B	1203	1	14,14,15	1.49	1 (7%)	17,19,21	1.04	1 (5%)
2	NAG	C	1204	1	14,14,15	1.40	2 (14%)	17,19,21	1.07	1 (5%)
2	NAG	C	1205	1	14,14,15	1.28	1 (7%)	17,19,21	1.67	2 (11%)
2	NAG	A	1204	1	14,14,15	1.27	2 (14%)	17,19,21	1.08	1 (5%)
2	NAG	B	1212	1	14,14,15	1.38	1 (7%)	17,19,21	1.04	1 (5%)
2	NAG	A	1207	1	14,14,15	1.39	2 (14%)	17,19,21	0.95	1 (5%)
2	NAG	C	1210	1	14,14,15	1.34	2 (14%)	17,19,21	1.00	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
2	NAG	B	1208	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1213	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1211	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1201	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1213	1	-	2/6/23/26	0/1/1/1
2	NAG	A	1201	1	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	1205	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1201	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1208	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1209	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1212	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1208	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1212	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1210	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1207	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1209	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1206	1	-	2/6/23/26	0/1/1/1
2	NAG	A	1203	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1202	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1203	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1204	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1206	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1206	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1209	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1205	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1202	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1211	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1207	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1211	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1210	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1202	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1203	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1204	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1205	1	-	3/6/23/26	0/1/1/1
2	NAG	A	1204	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1212	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1207	1	-	2/6/23/26	0/1/1/1
2	NAG	C	1210	1	-	1/6/23/26	0/1/1/1

All (62) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1201	NAG	C1-C2	4.76	1.58	1.52
2	A	1211	NAG	C1-C2	4.60	1.58	1.52
2	B	1203	NAG	C1-C2	4.58	1.58	1.52
2	A	1213	NAG	C1-C2	4.53	1.58	1.52
2	C	1207	NAG	C1-C2	4.47	1.58	1.52
2	A	1210	NAG	C1-C2	4.30	1.58	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1201	NAG	C1-C2	4.23	1.58	1.52
2	A	1212	NAG	C1-C2	4.21	1.58	1.52
2	C	1213	NAG	C1-C2	4.21	1.58	1.52
2	B	1205	NAG	C1-C2	4.20	1.58	1.52
2	A	1203	NAG	C1-C2	4.19	1.58	1.52
2	C	1204	NAG	C1-C2	4.19	1.58	1.52
2	C	1203	NAG	C1-C2	4.17	1.58	1.52
2	B	1212	NAG	C1-C2	4.16	1.58	1.52
2	C	1202	NAG	C1-C2	4.16	1.58	1.52
2	B	1204	NAG	C1-C2	4.15	1.58	1.52
2	B	1207	NAG	C1-C2	4.08	1.57	1.52
2	A	1201	NAG	C1-C2	4.08	1.57	1.52
2	C	1212	NAG	C1-C2	4.07	1.57	1.52
2	A	1202	NAG	C1-C2	4.06	1.57	1.52
2	C	1208	NAG	C1-C2	4.04	1.57	1.52
2	A	1206	NAG	C1-C2	4.04	1.57	1.52
2	A	1208	NAG	C1-C2	4.02	1.57	1.52
2	C	1206	NAG	C1-C2	4.01	1.57	1.52
2	A	1207	NAG	C1-C2	3.99	1.57	1.52
2	B	1208	NAG	C1-C2	3.96	1.57	1.52
2	B	1202	NAG	C1-C2	3.95	1.57	1.52
2	B	1209	NAG	C1-C2	3.95	1.57	1.52
2	B	1211	NAG	C1-C2	3.92	1.57	1.52
2	C	1211	NAG	C1-C2	3.91	1.57	1.52
2	C	1209	NAG	C1-C2	3.88	1.57	1.52
2	C	1210	NAG	C1-C2	3.88	1.57	1.52
2	A	1209	NAG	C1-C2	3.85	1.57	1.52
2	B	1210	NAG	C1-C2	3.82	1.57	1.52
2	C	1205	NAG	C1-C2	3.78	1.57	1.52
2	A	1205	NAG	C1-C2	3.77	1.57	1.52
2	A	1204	NAG	C1-C2	3.70	1.57	1.52
2	B	1206	NAG	C1-C2	3.65	1.57	1.52
2	A	1207	NAG	O5-C5	2.26	1.47	1.43
2	C	1207	NAG	O5-C5	2.25	1.47	1.43
2	B	1209	NAG	O5-C5	2.21	1.47	1.43
2	A	1209	NAG	O5-C5	2.19	1.47	1.43
2	C	1211	NAG	O5-C5	2.18	1.47	1.43
2	C	1209	NAG	O5-C5	2.16	1.47	1.43
2	A	1211	NAG	O5-C5	2.14	1.47	1.43
2	A	1205	NAG	O5-C5	2.12	1.47	1.43
2	A	1212	NAG	O5-C5	2.12	1.47	1.43
2	C	1210	NAG	O5-C5	2.12	1.47	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1202	NAG	O5-C5	2.12	1.47	1.43
2	A	1201	NAG	O5-C5	2.11	1.47	1.43
2	B	1211	NAG	O5-C5	2.11	1.47	1.43
2	A	1208	NAG	O5-C5	2.11	1.47	1.43
2	A	1204	NAG	O5-C5	2.10	1.47	1.43
2	C	1204	NAG	O5-C5	2.10	1.47	1.43
2	C	1203	NAG	O5-C5	2.10	1.47	1.43
2	A	1202	NAG	O5-C5	2.07	1.47	1.43
2	C	1208	NAG	O5-C5	2.06	1.47	1.43
2	A	1213	NAG	O5-C5	2.06	1.47	1.43
2	B	1205	NAG	O5-C5	2.06	1.47	1.43
2	B	1202	NAG	O5-C5	2.05	1.47	1.43
2	B	1206	NAG	O5-C5	2.03	1.47	1.43
2	B	1210	NAG	O5-C5	2.02	1.47	1.43

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1205	NAG	C8-C7-N2	5.30	124.91	116.12
2	C	1205	NAG	O7-C7-N2	-3.27	116.21	121.98
2	A	1213	NAG	C8-C7-N2	2.95	121.01	116.12
2	B	1207	NAG	C8-C7-N2	2.70	120.59	116.12
2	C	1213	NAG	C8-C7-N2	2.68	120.57	116.12
2	C	1203	NAG	C8-C7-N2	2.61	120.44	116.12
2	C	1204	NAG	C8-C7-N2	2.59	120.41	116.12
2	A	1204	NAG	C8-C7-N2	2.58	120.39	116.12
2	A	1201	NAG	C8-C7-N2	2.57	120.38	116.12
2	C	1201	NAG	C8-C7-N2	2.57	120.37	116.12
2	A	1206	NAG	C8-C7-N2	2.55	120.34	116.12
2	B	1202	NAG	C8-C7-N2	2.53	120.32	116.12
2	C	1211	NAG	C8-C7-N2	2.48	120.23	116.12
2	C	1206	NAG	C8-C7-N2	2.47	120.22	116.12
2	B	1204	NAG	C8-C7-N2	2.46	120.19	116.12
2	A	1203	NAG	C8-C7-N2	2.43	120.15	116.12
2	B	1209	NAG	C8-C7-N2	2.43	120.15	116.12
2	A	1205	NAG	C8-C7-N2	2.41	120.12	116.12
2	C	1212	NAG	C8-C7-N2	2.41	120.12	116.12
2	A	1208	NAG	C8-C7-N2	2.41	120.11	116.12
2	C	1209	NAG	C8-C7-N2	2.41	120.11	116.12
2	B	1211	NAG	C8-C7-N2	2.40	120.11	116.12
2	C	1208	NAG	C8-C7-N2	2.39	120.09	116.12
2	C	1202	NAG	C8-C7-N2	2.39	120.08	116.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1208	NAG	C8-C7-N2	2.38	120.07	116.12
2	A	1202	NAG	C8-C7-N2	2.38	120.06	116.12
2	B	1212	NAG	C8-C7-N2	2.36	120.03	116.12
2	A	1207	NAG	C8-C7-N2	2.36	120.03	116.12
2	A	1209	NAG	C8-C7-N2	2.35	120.02	116.12
2	C	1210	NAG	C8-C7-N2	2.35	120.02	116.12
2	B	1205	NAG	C8-C7-N2	2.34	120.00	116.12
2	B	1203	NAG	C8-C7-N2	2.34	120.00	116.12
2	A	1210	NAG	C8-C7-N2	2.33	119.98	116.12
2	B	1206	NAG	C8-C7-N2	2.31	119.95	116.12
2	B	1210	NAG	C8-C7-N2	2.27	119.89	116.12
2	C	1202	NAG	C2-N2-C7	-2.21	119.93	122.90
2	A	1212	NAG	C8-C7-N2	2.20	119.77	116.12
2	C	1207	NAG	C8-C7-N2	2.19	119.75	116.12
2	A	1213	NAG	O7-C7-C8	-2.19	118.16	122.05
2	A	1211	NAG	C8-C7-N2	2.11	119.62	116.12

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1205	NAG	C8-C7-N2-C2
2	C	1205	NAG	O7-C7-N2-C2
2	C	1206	NAG	O5-C5-C6-O6
2	B	1212	NAG	O5-C5-C6-O6
2	C	1209	NAG	O5-C5-C6-O6
2	A	1205	NAG	O5-C5-C6-O6
2	B	1203	NAG	O5-C5-C6-O6
2	B	1204	NAG	O5-C5-C6-O6
2	C	1213	NAG	O5-C5-C6-O6
2	C	1212	NAG	O5-C5-C6-O6
2	C	1203	NAG	O5-C5-C6-O6
2	C	1211	NAG	O5-C5-C6-O6
2	B	1205	NAG	O5-C5-C6-O6
2	B	1209	NAG	O5-C5-C6-O6
2	B	1208	NAG	O5-C5-C6-O6
2	C	1208	NAG	O5-C5-C6-O6
2	A	1208	NAG	O5-C5-C6-O6
2	A	1209	NAG	O5-C5-C6-O6
2	A	1211	NAG	O5-C5-C6-O6
2	B	1211	NAG	O5-C5-C6-O6
2	A	1210	NAG	O5-C5-C6-O6

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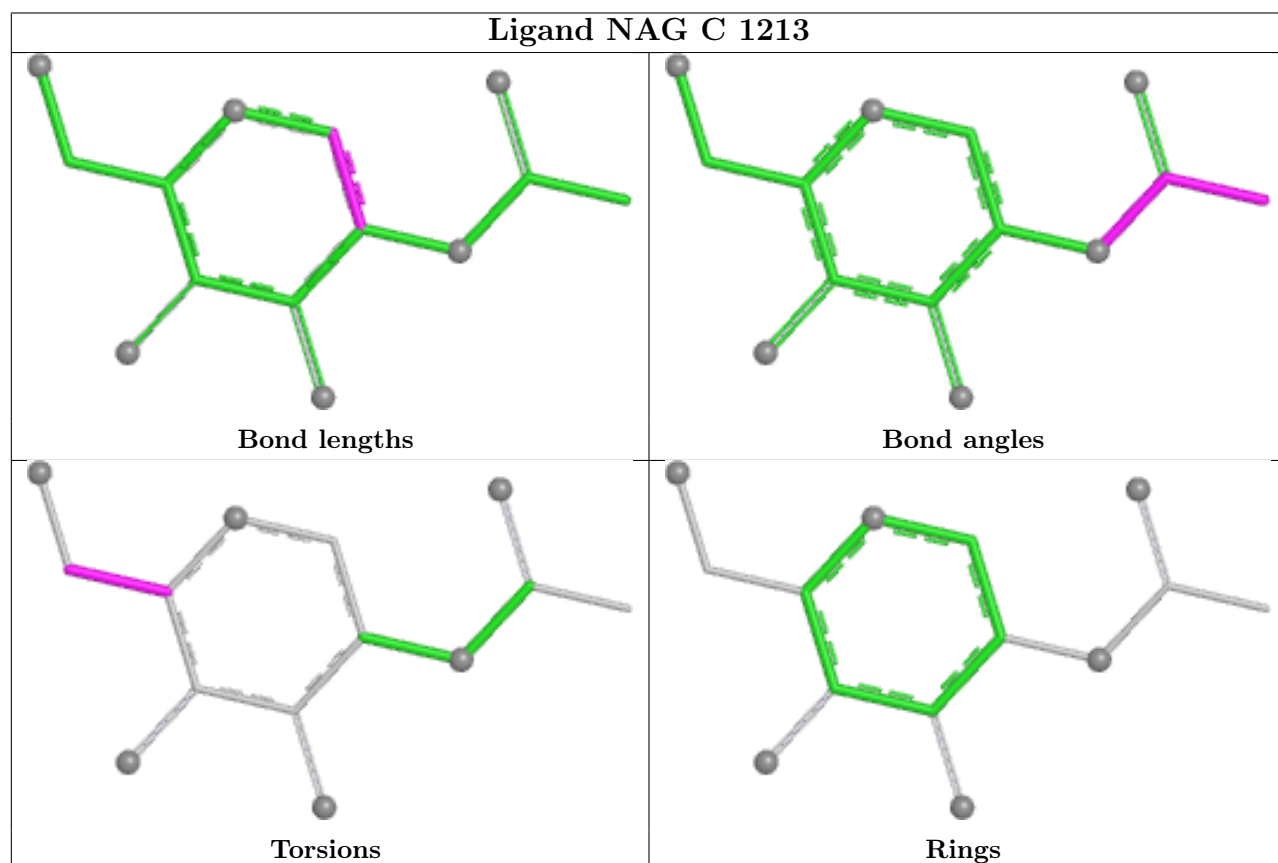
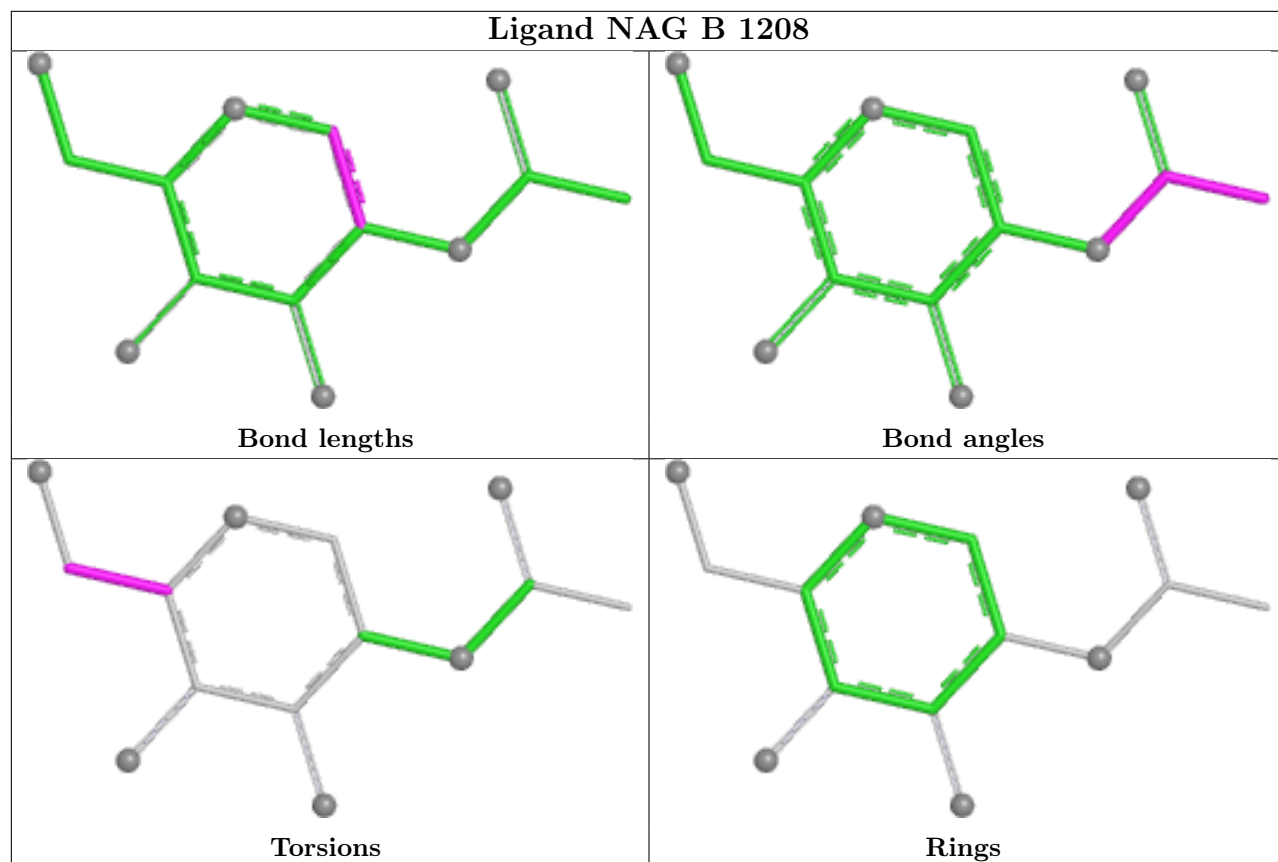
Mol	Chain	Res	Type	Atoms
2	C	1205	NAG	O5-C5-C6-O6
2	C	1210	NAG	O5-C5-C6-O6
2	B	1210	NAG	O5-C5-C6-O6
2	C	1202	NAG	O5-C5-C6-O6
2	A	1212	NAG	O5-C5-C6-O6
2	B	1201	NAG	O5-C5-C6-O6
2	A	1201	NAG	O5-C5-C6-O6
2	A	1207	NAG	C3-C2-N2-C7
2	A	1213	NAG	C3-C2-N2-C7
2	C	1206	NAG	C4-C5-C6-O6
2	A	1207	NAG	C1-C2-N2-C7
2	A	1213	NAG	C1-C2-N2-C7

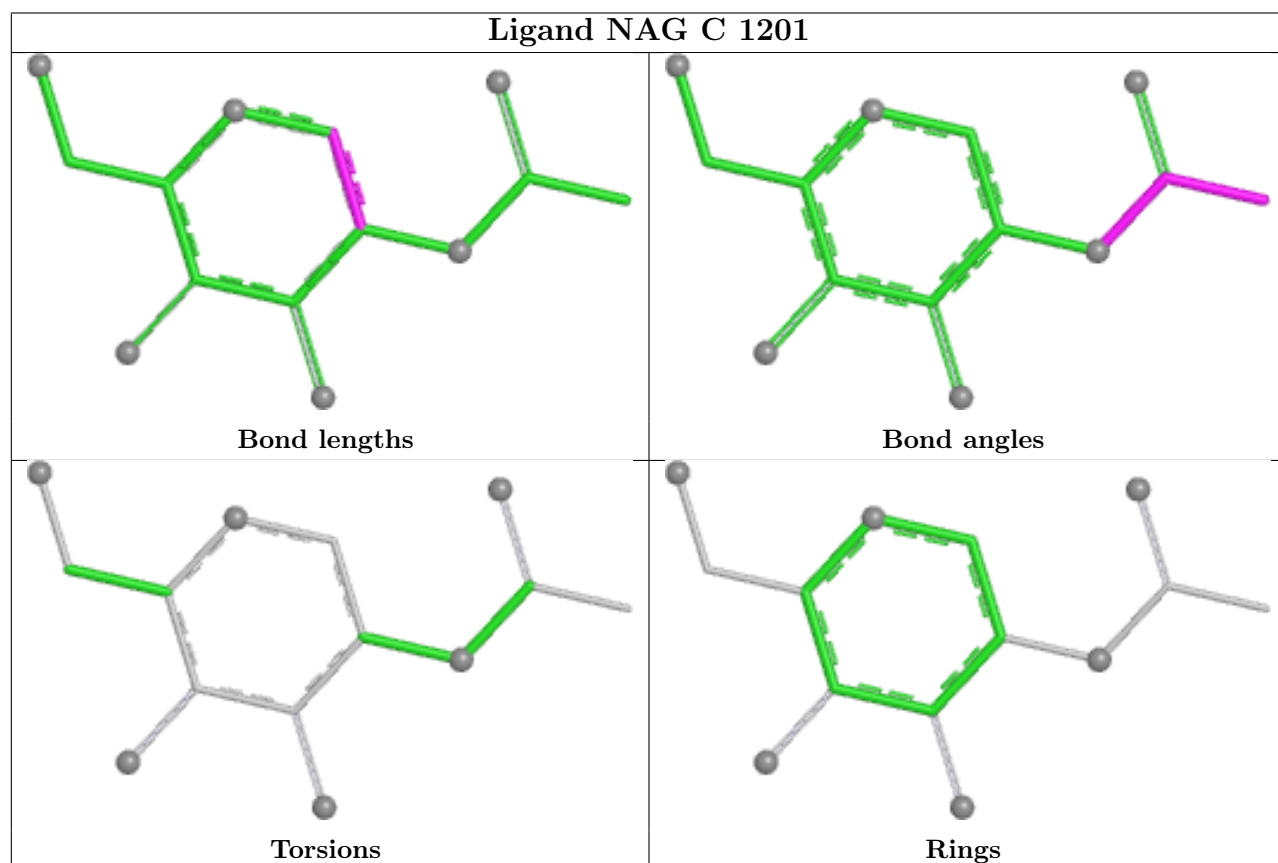
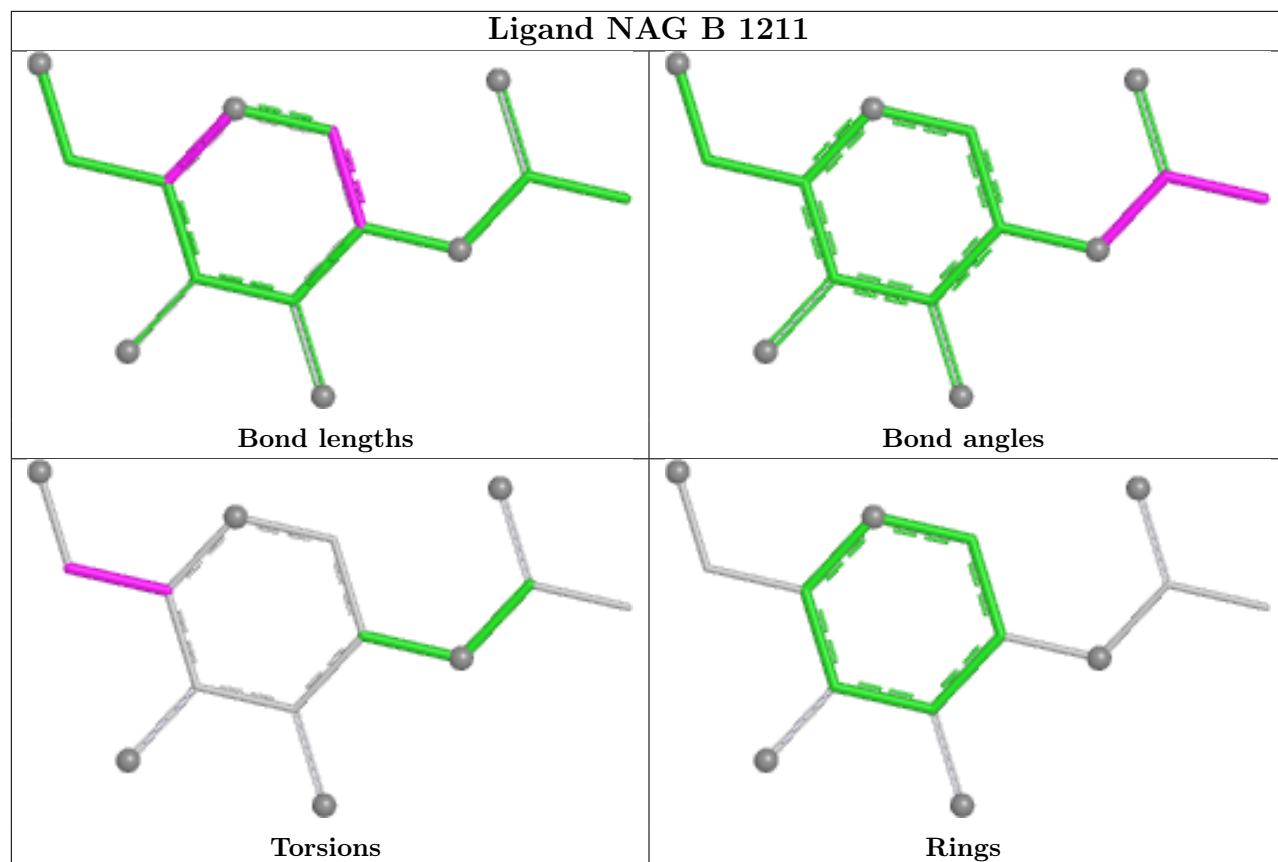
There are no ring outliers.

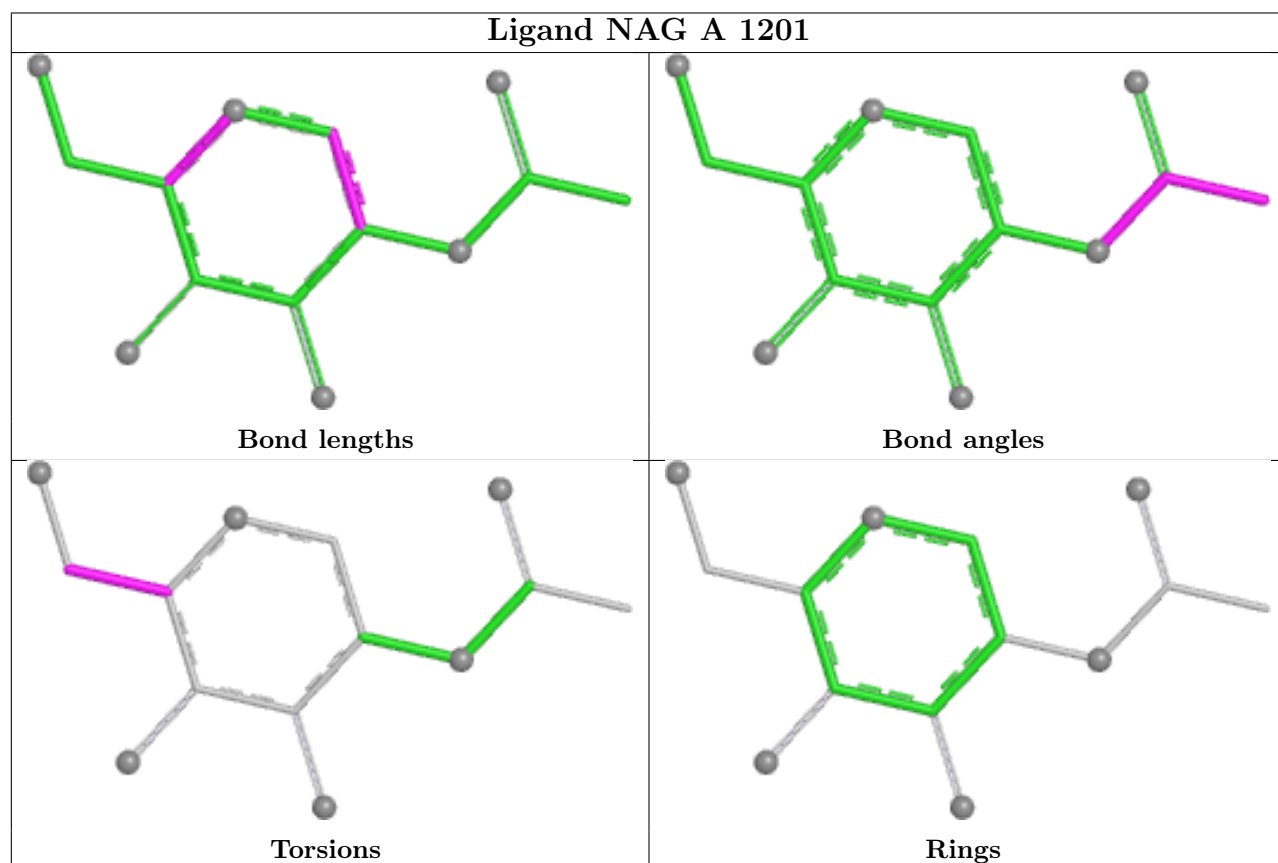
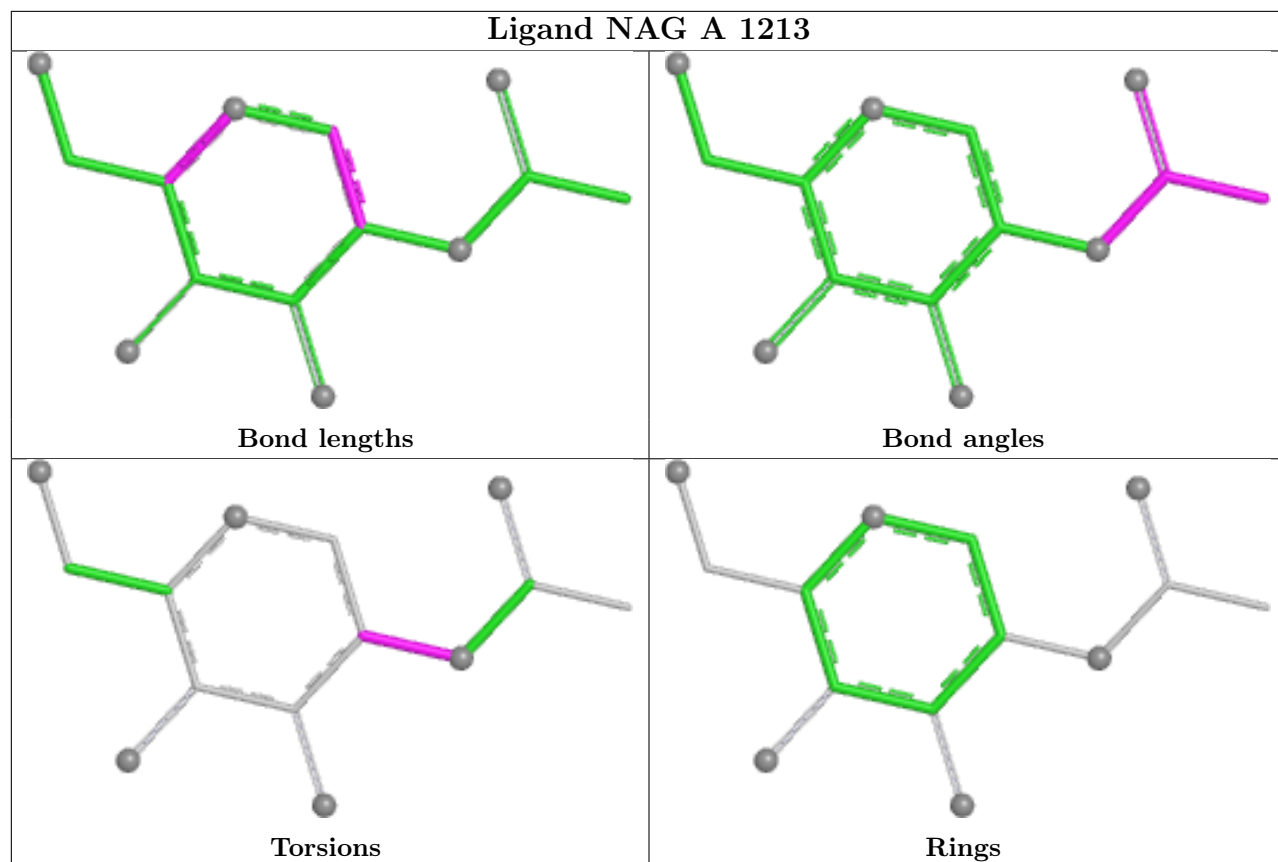
2 monomers are involved in 4 short contacts:

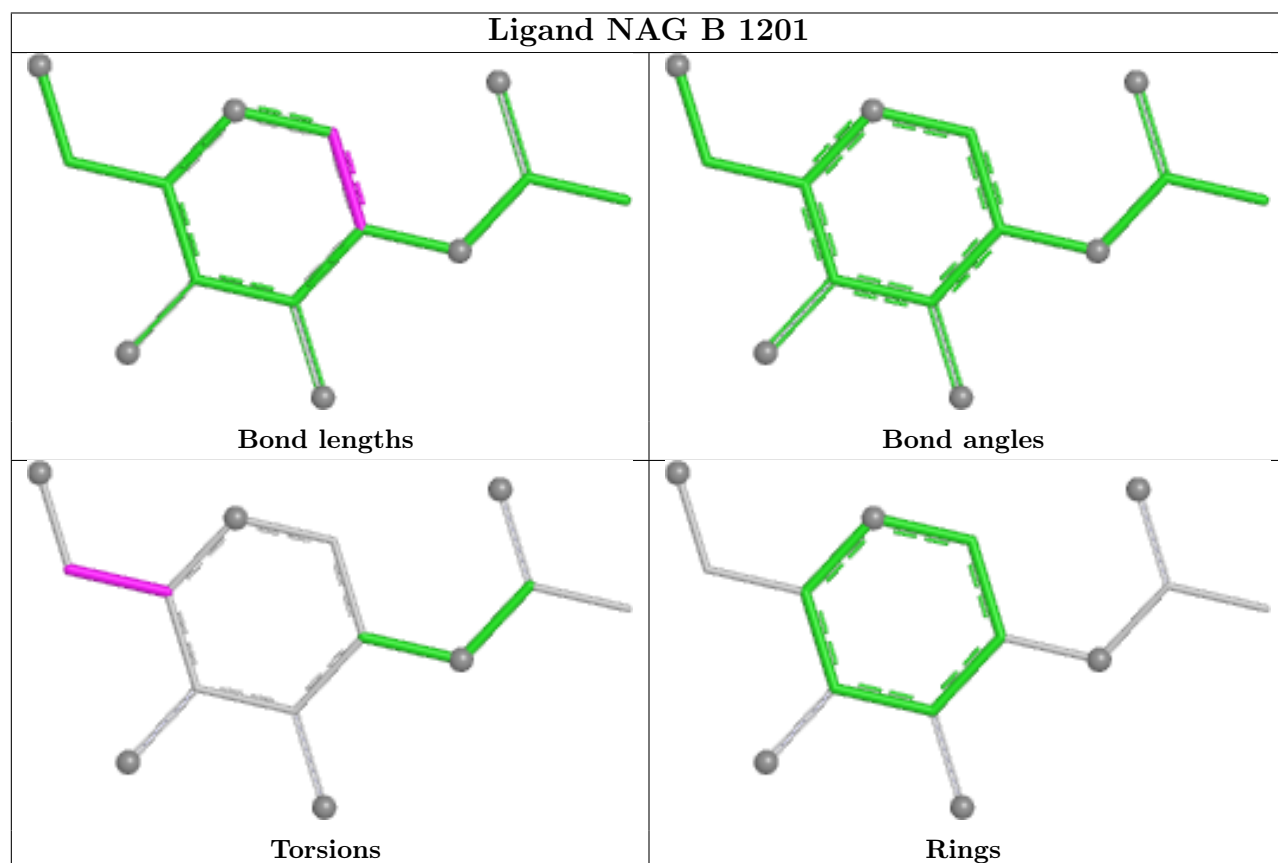
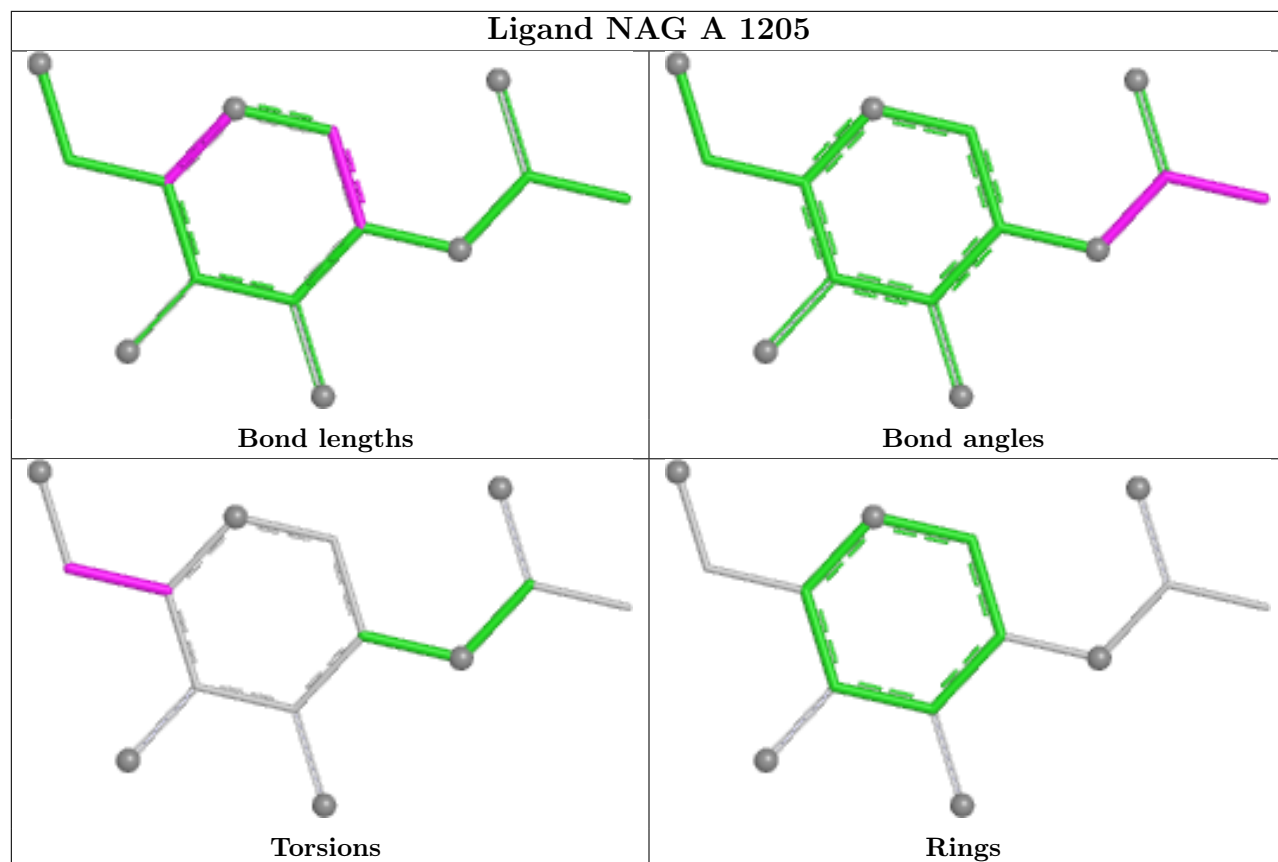
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1201	NAG	2	0
2	A	1213	NAG	2	0

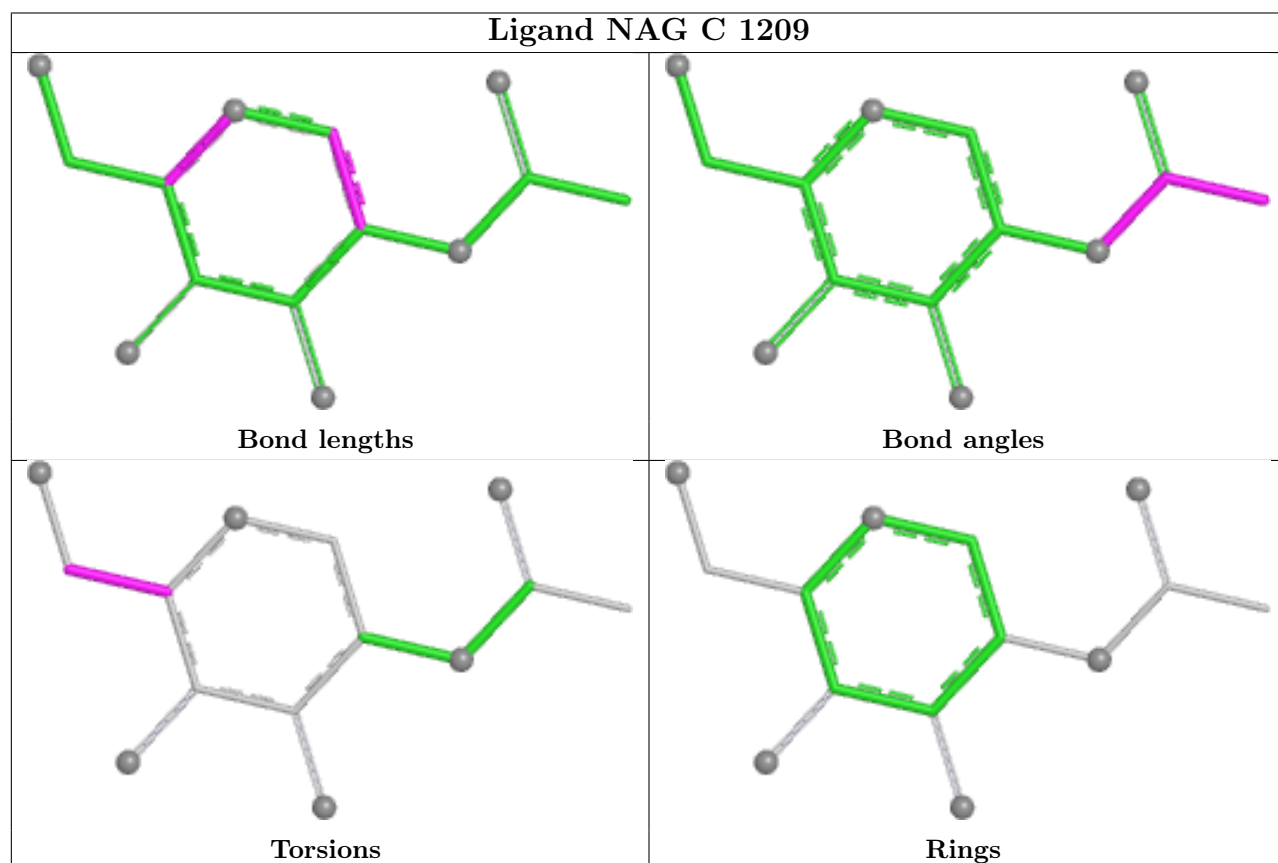
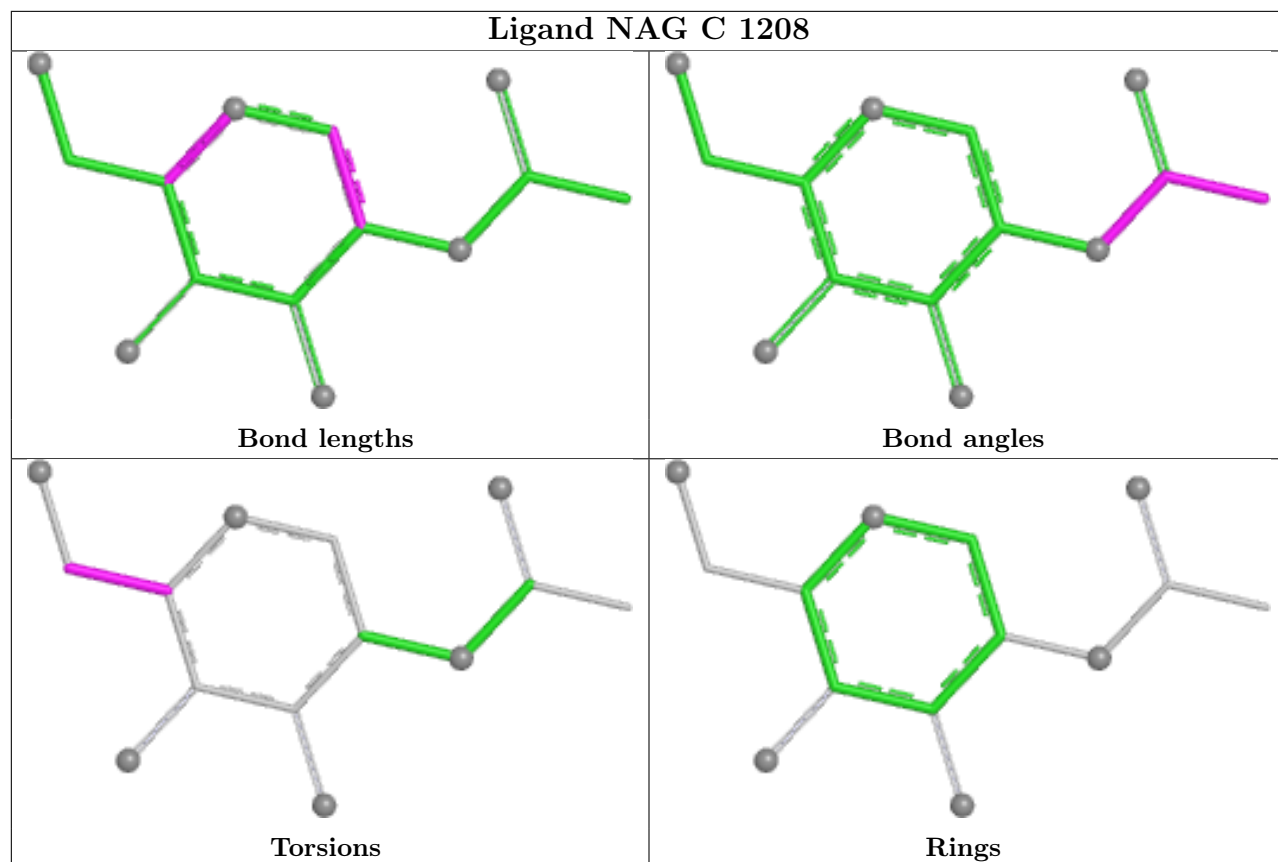
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



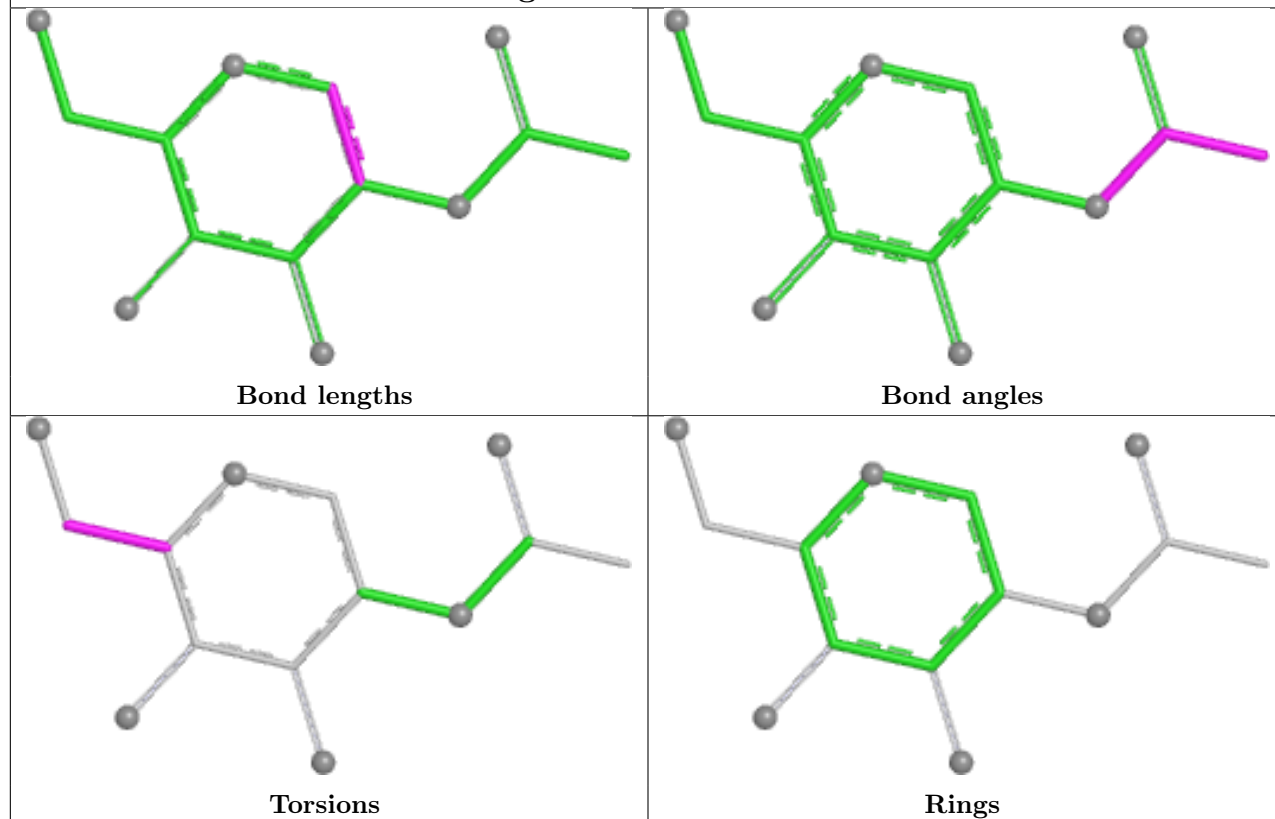




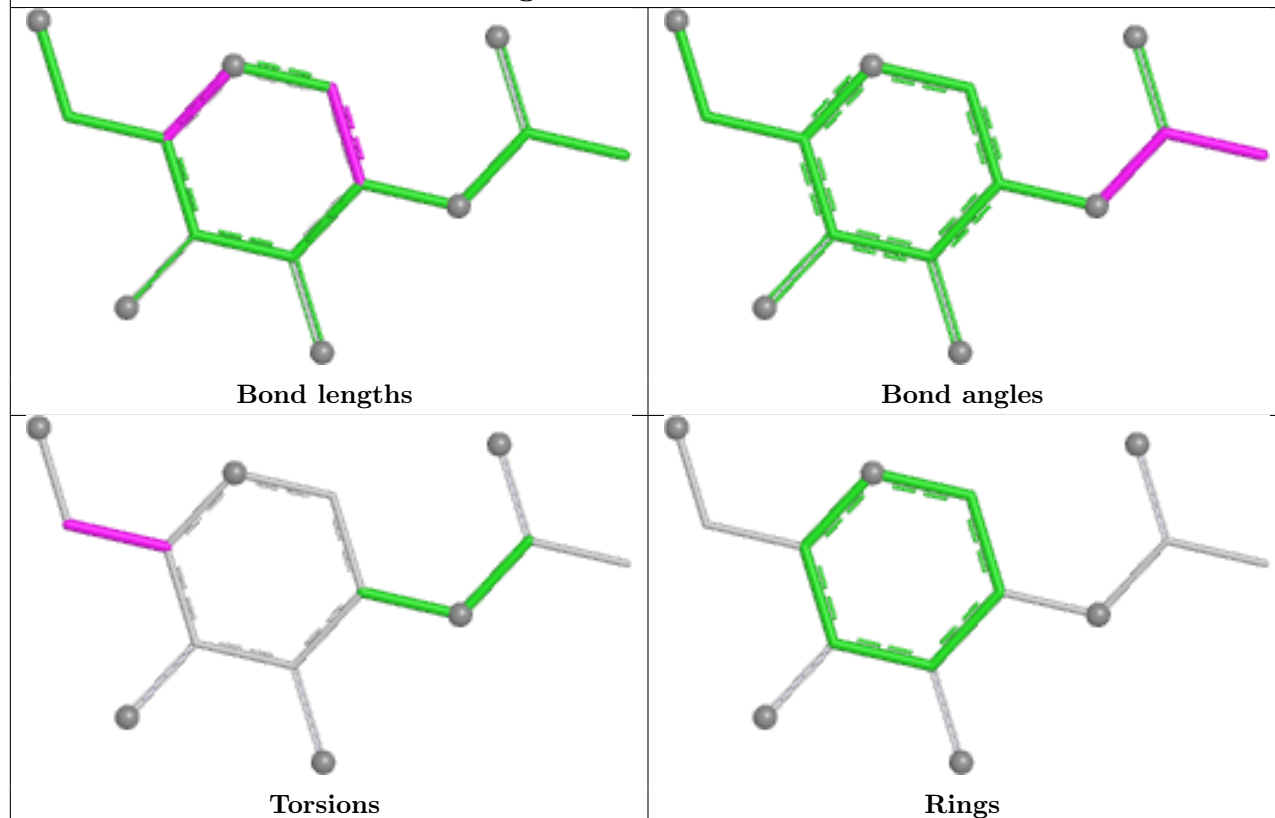




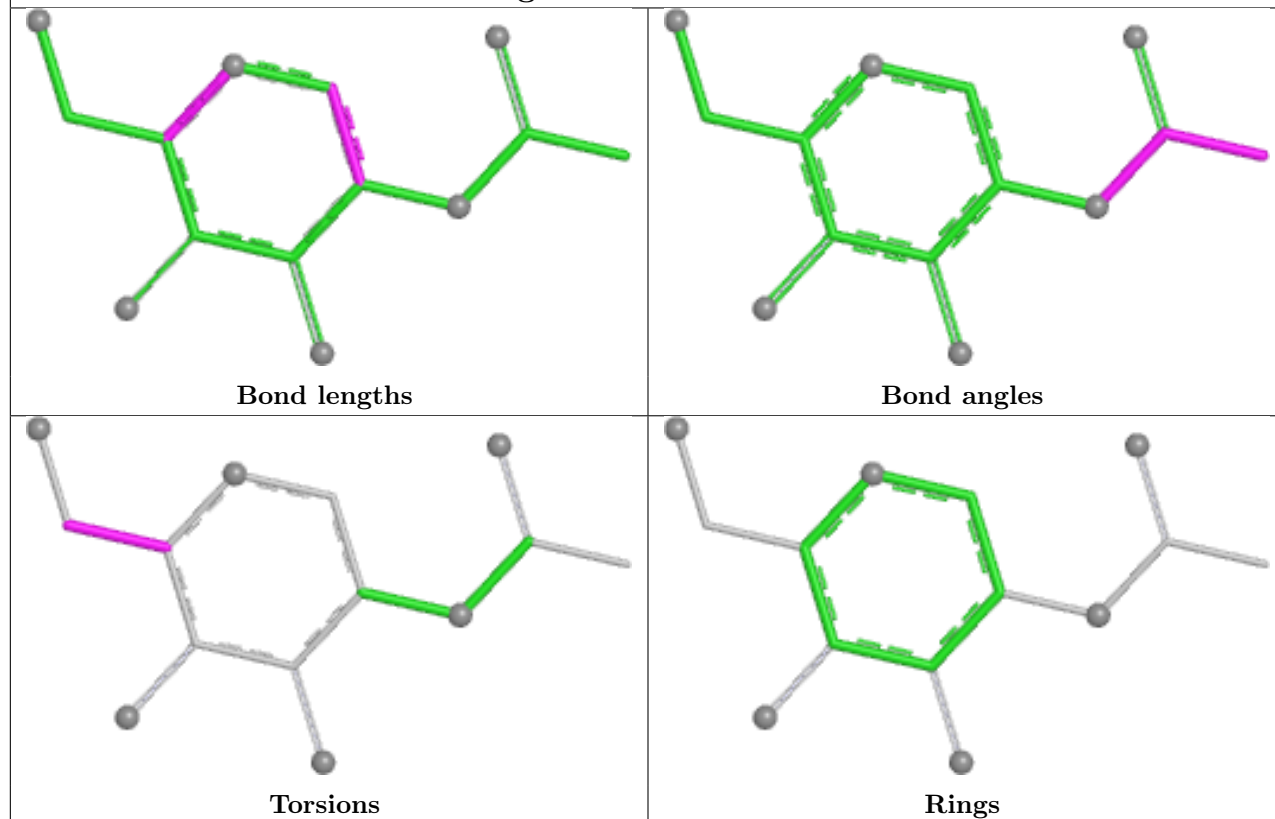
Ligand NAG C 1212



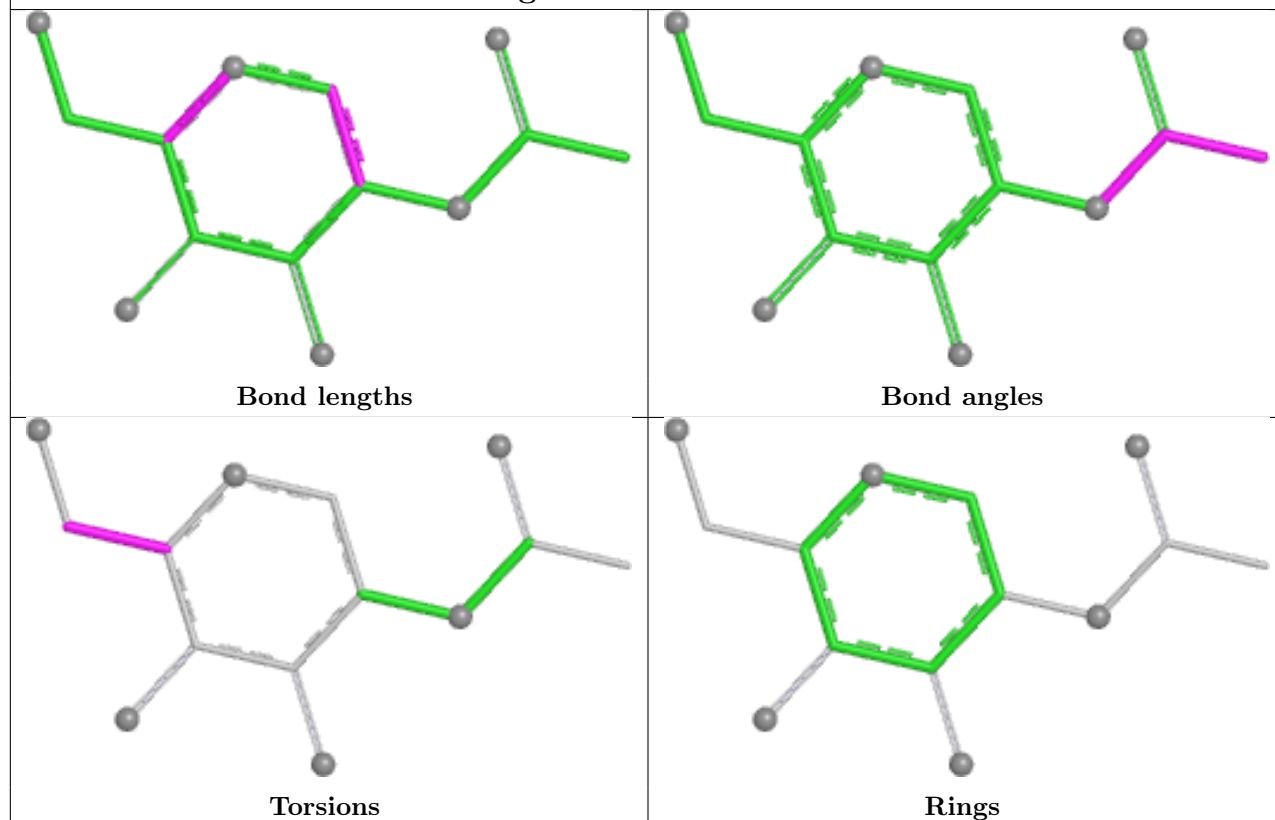
Ligand NAG A 1208

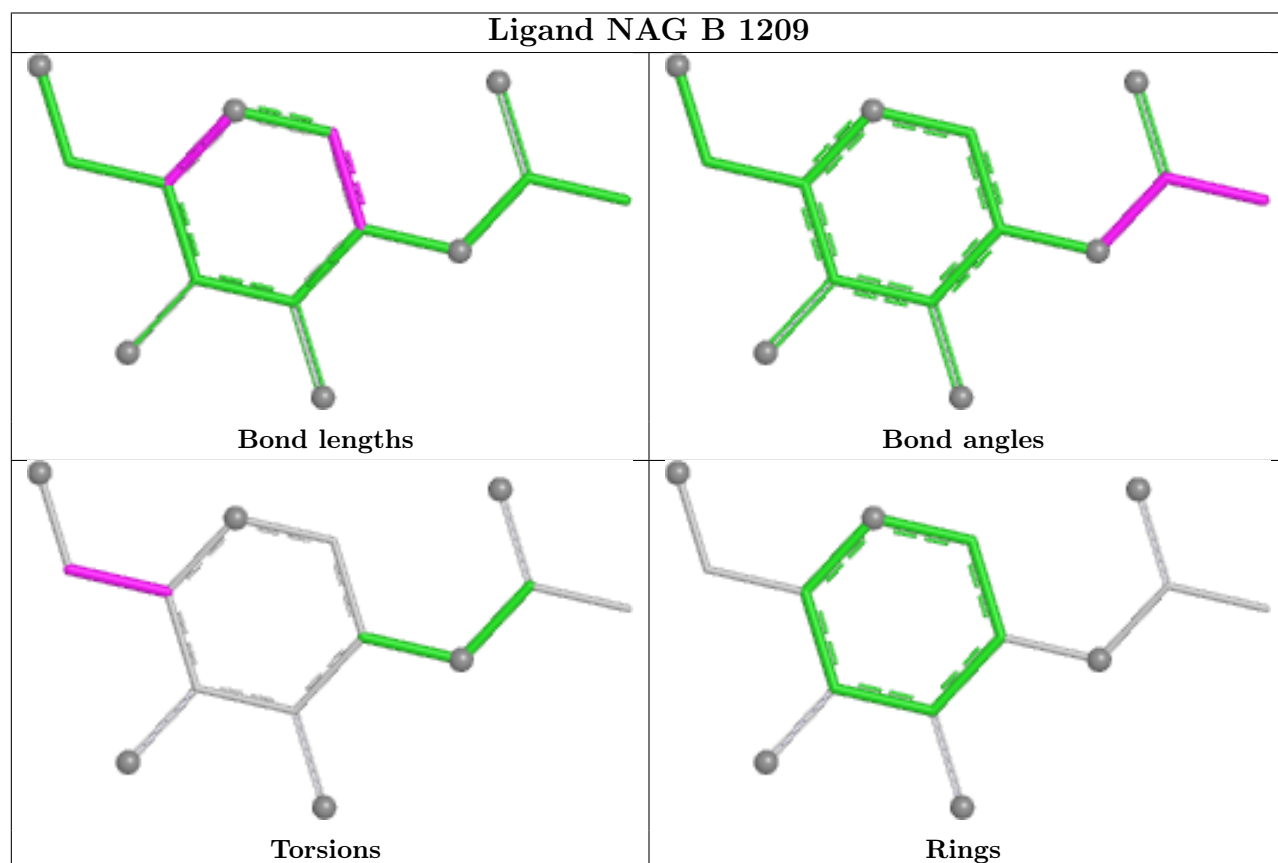
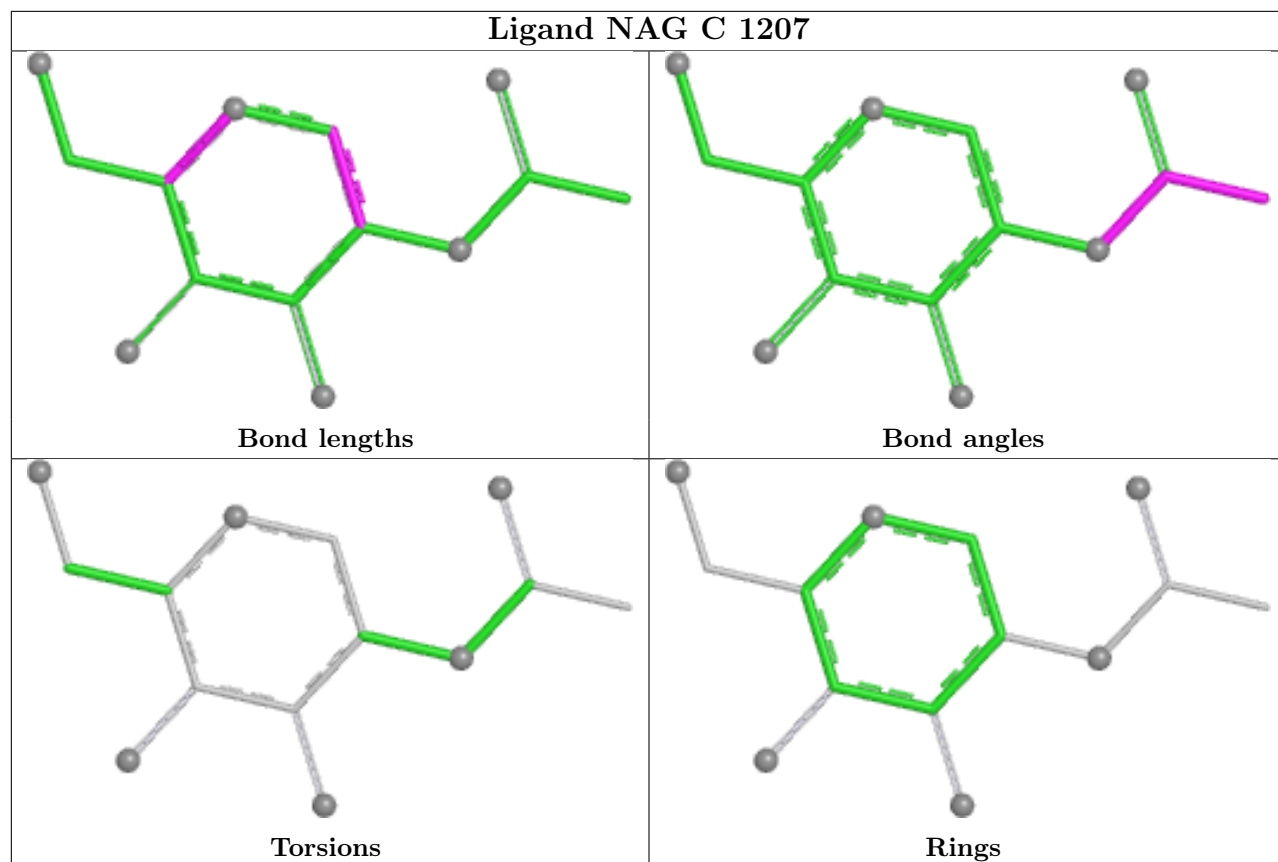


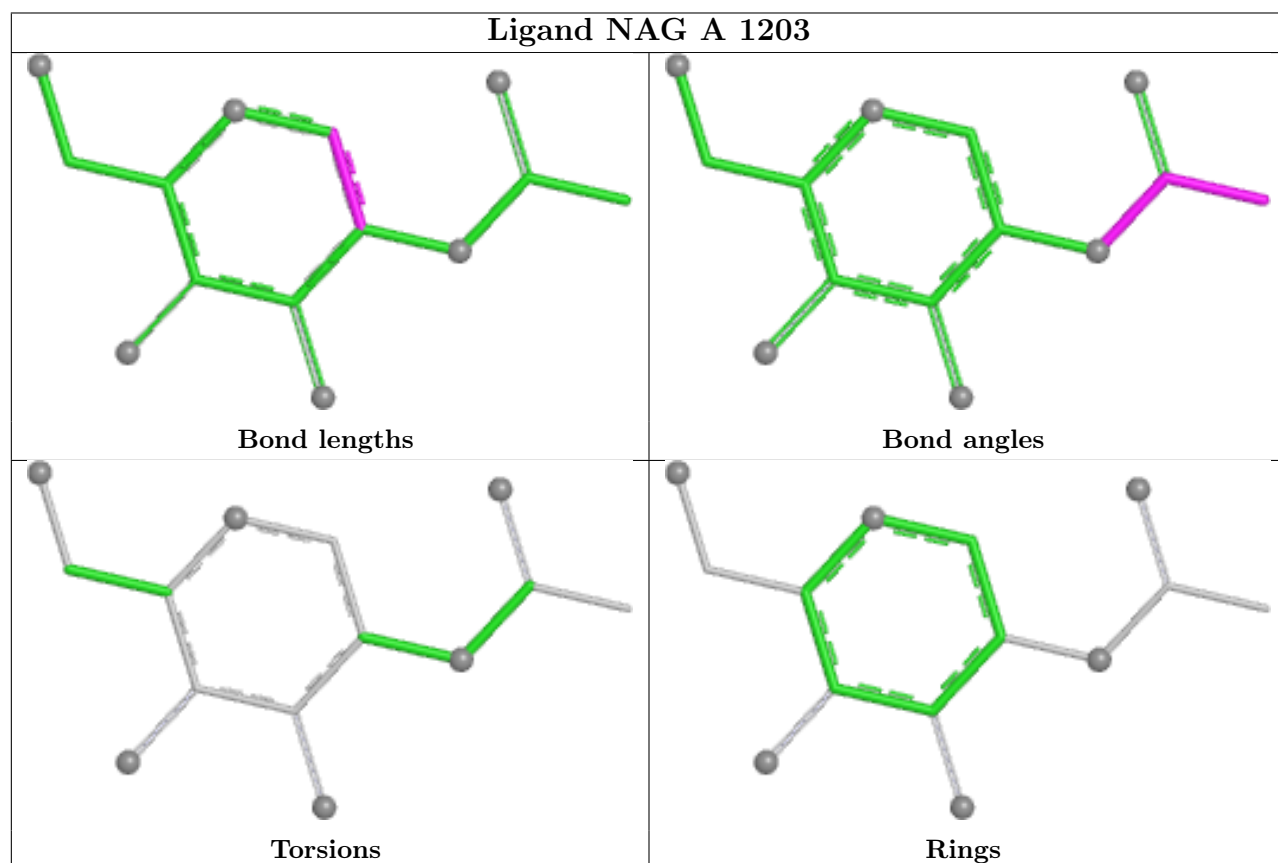
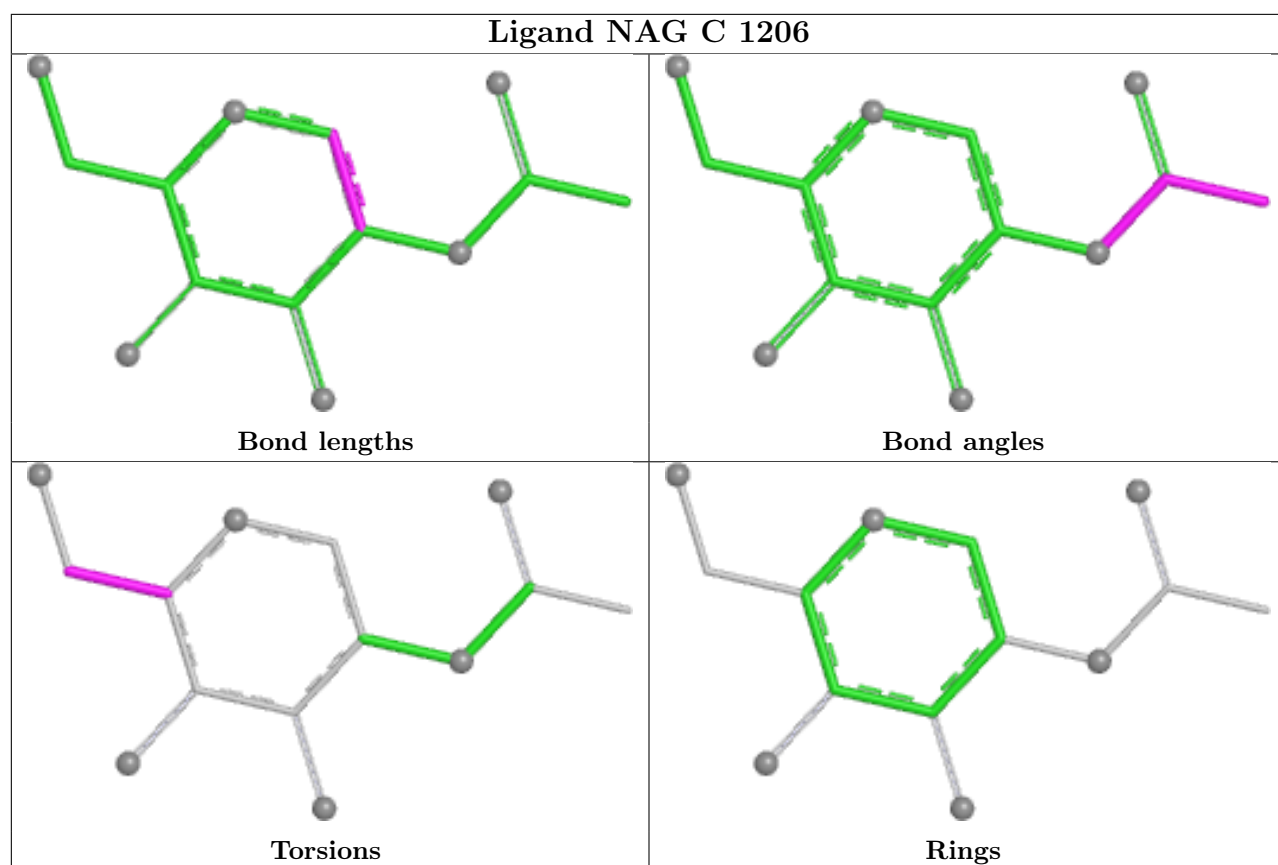
Ligand NAG A 1212

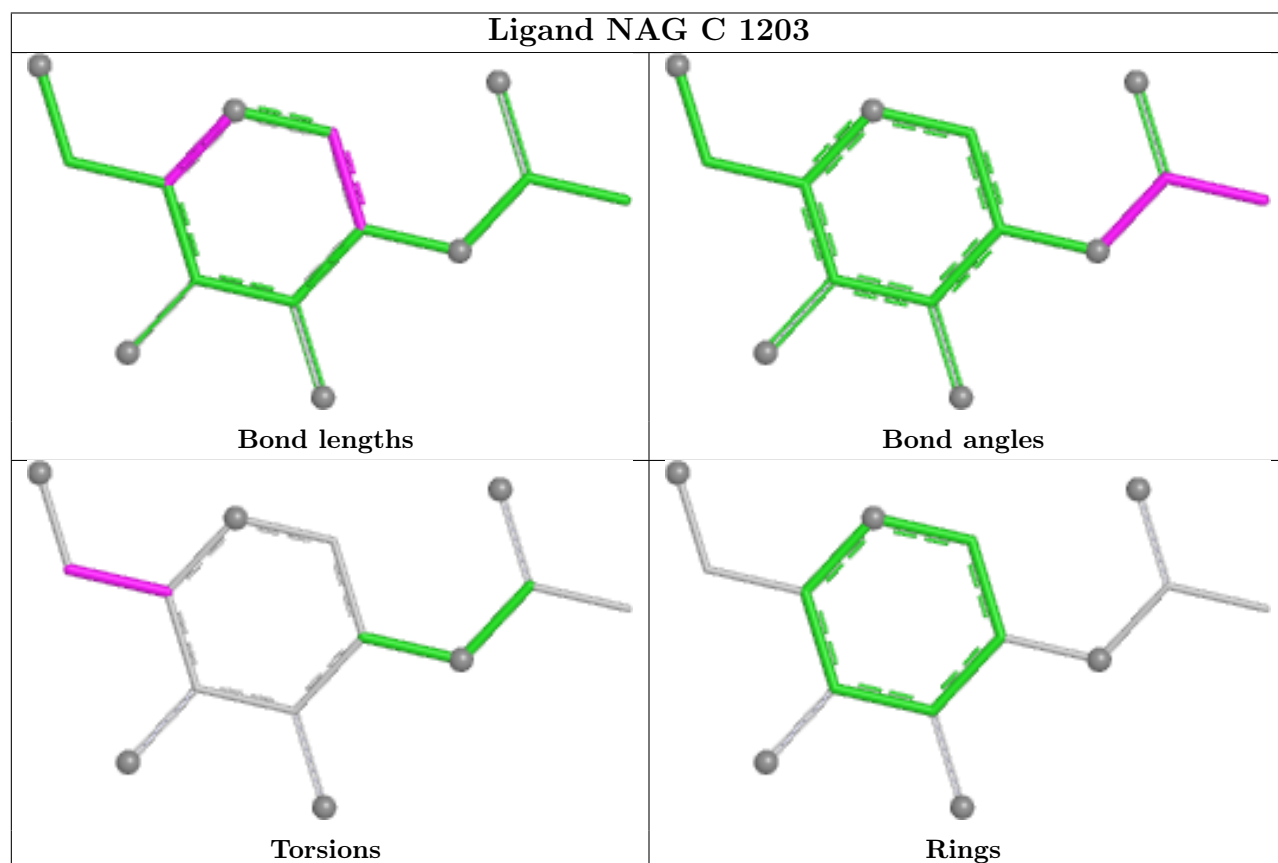
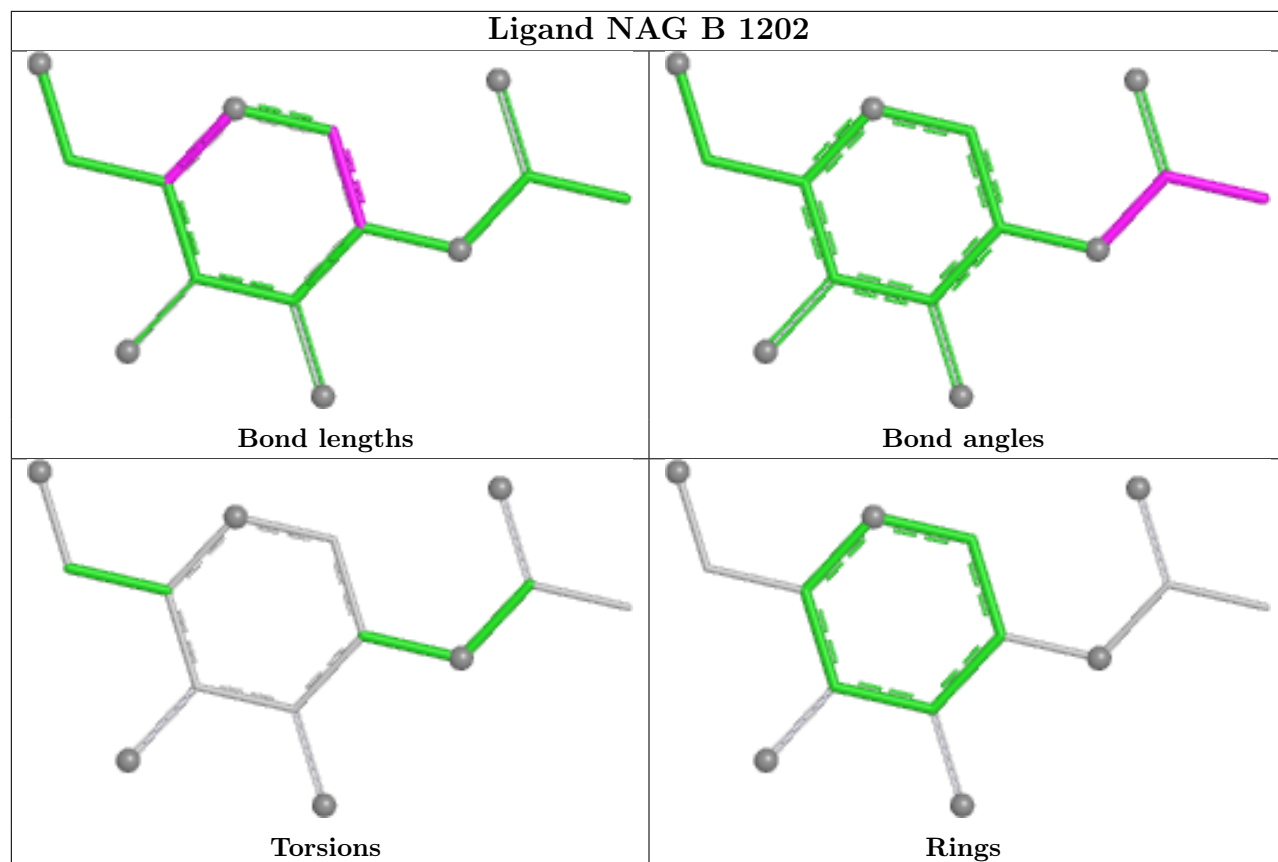


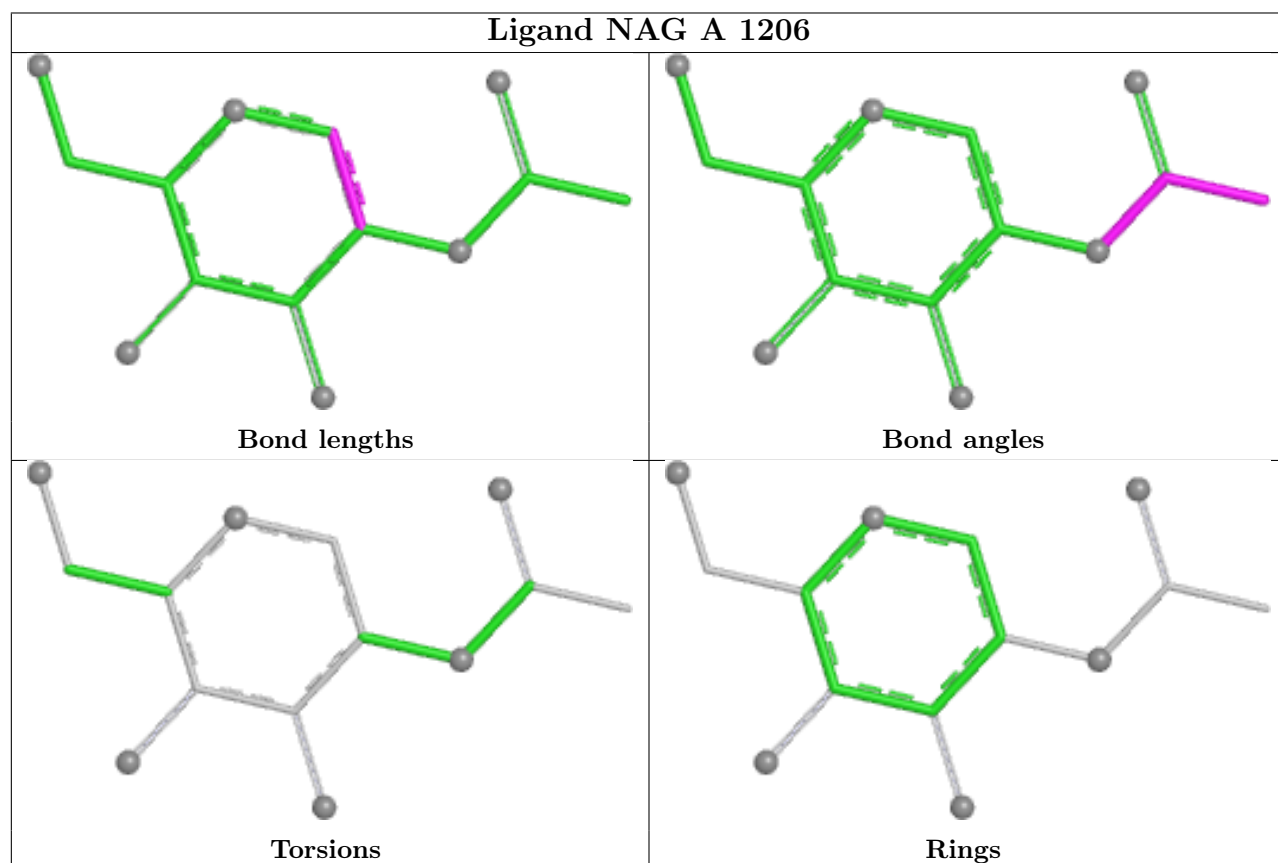
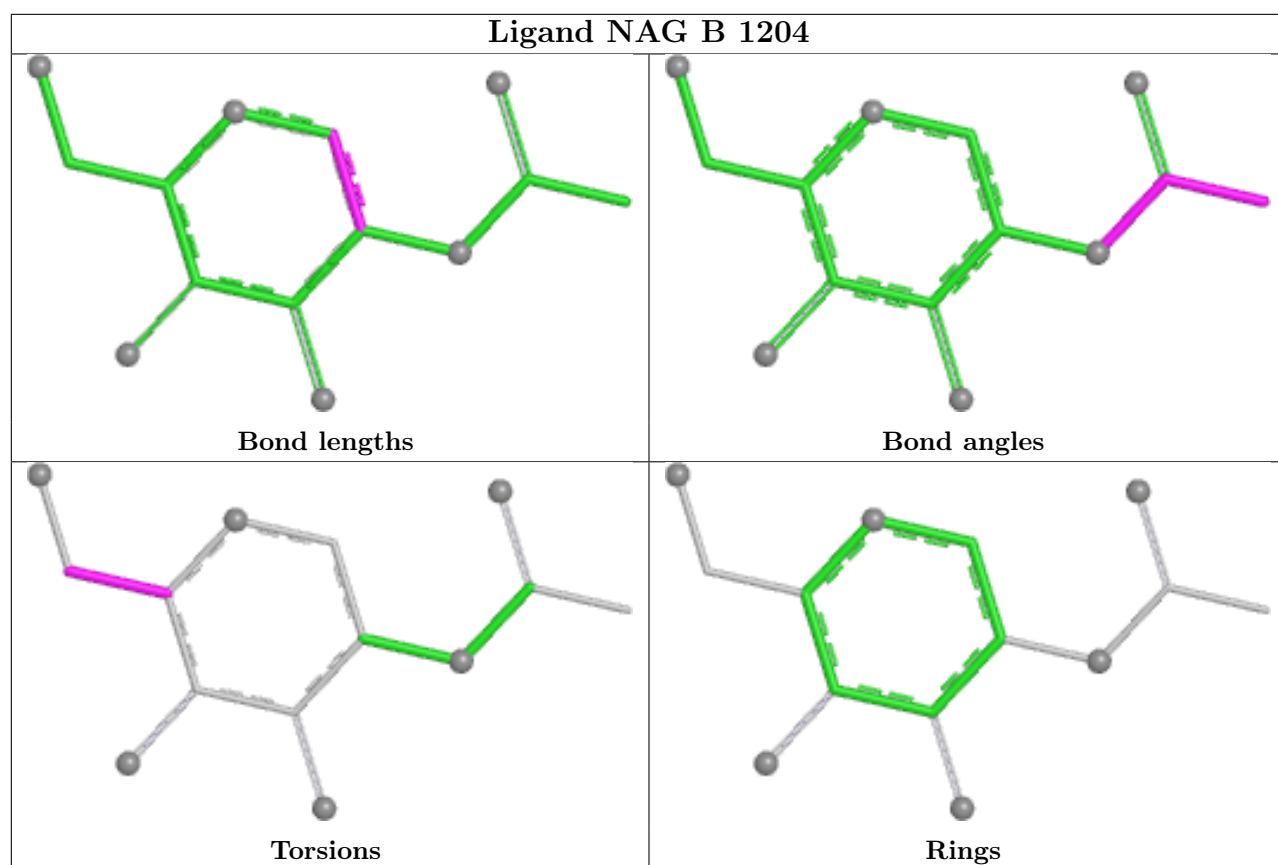
Ligand NAG B 1210

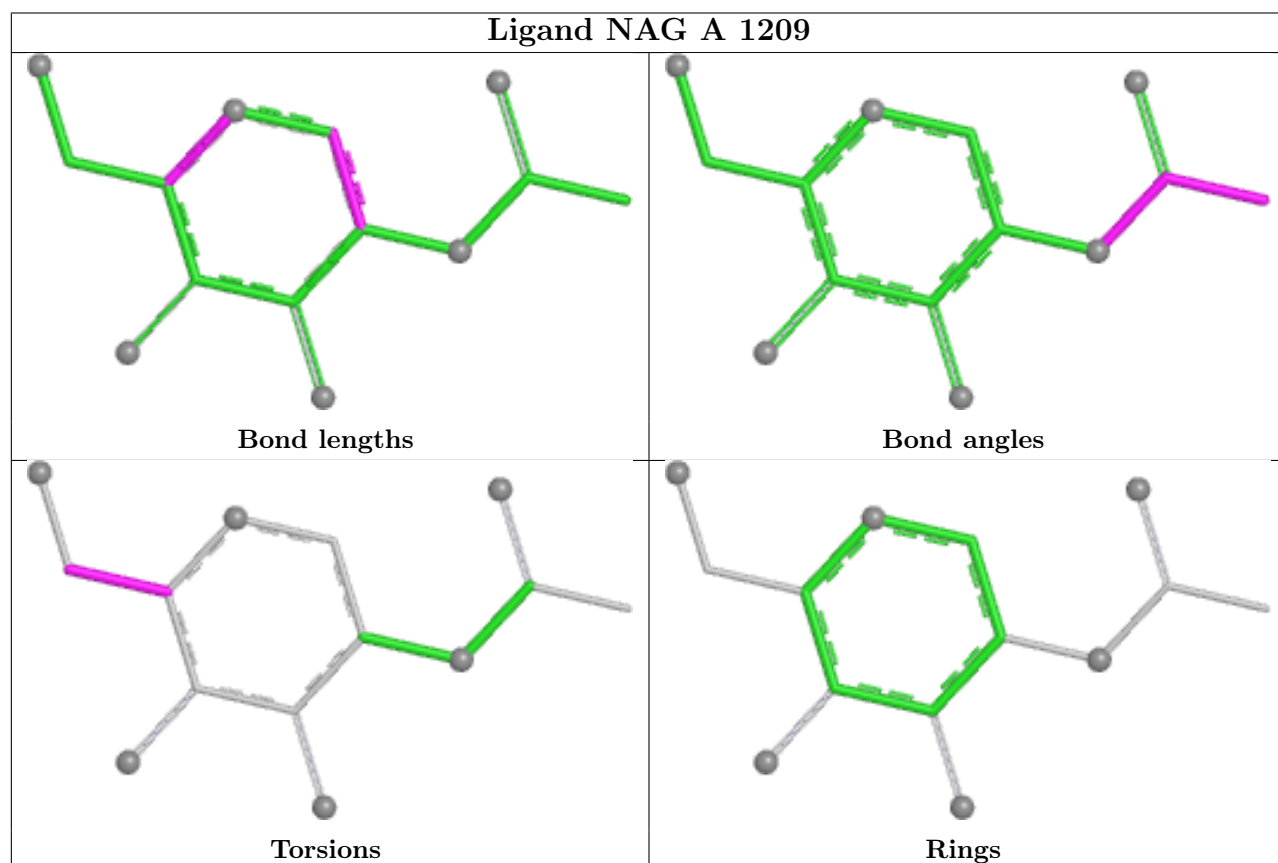
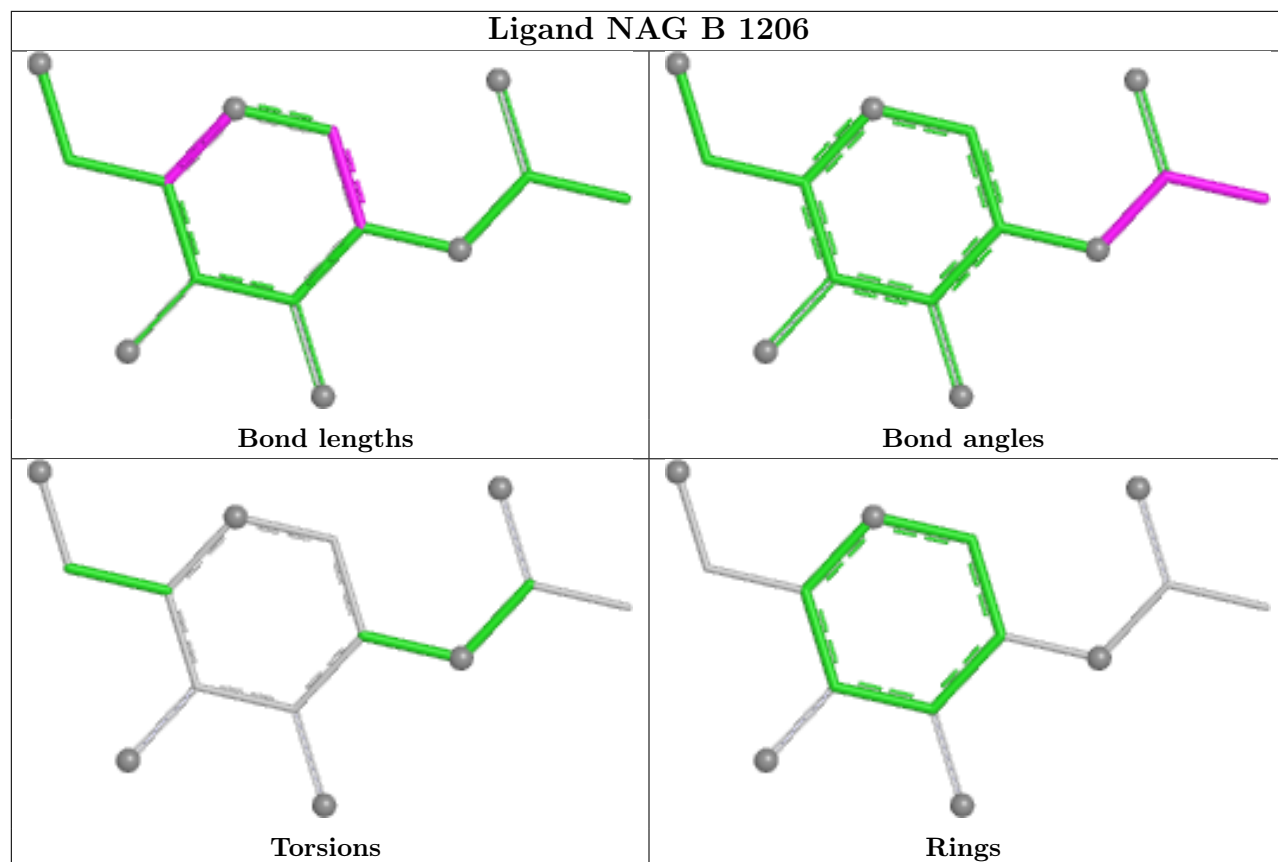




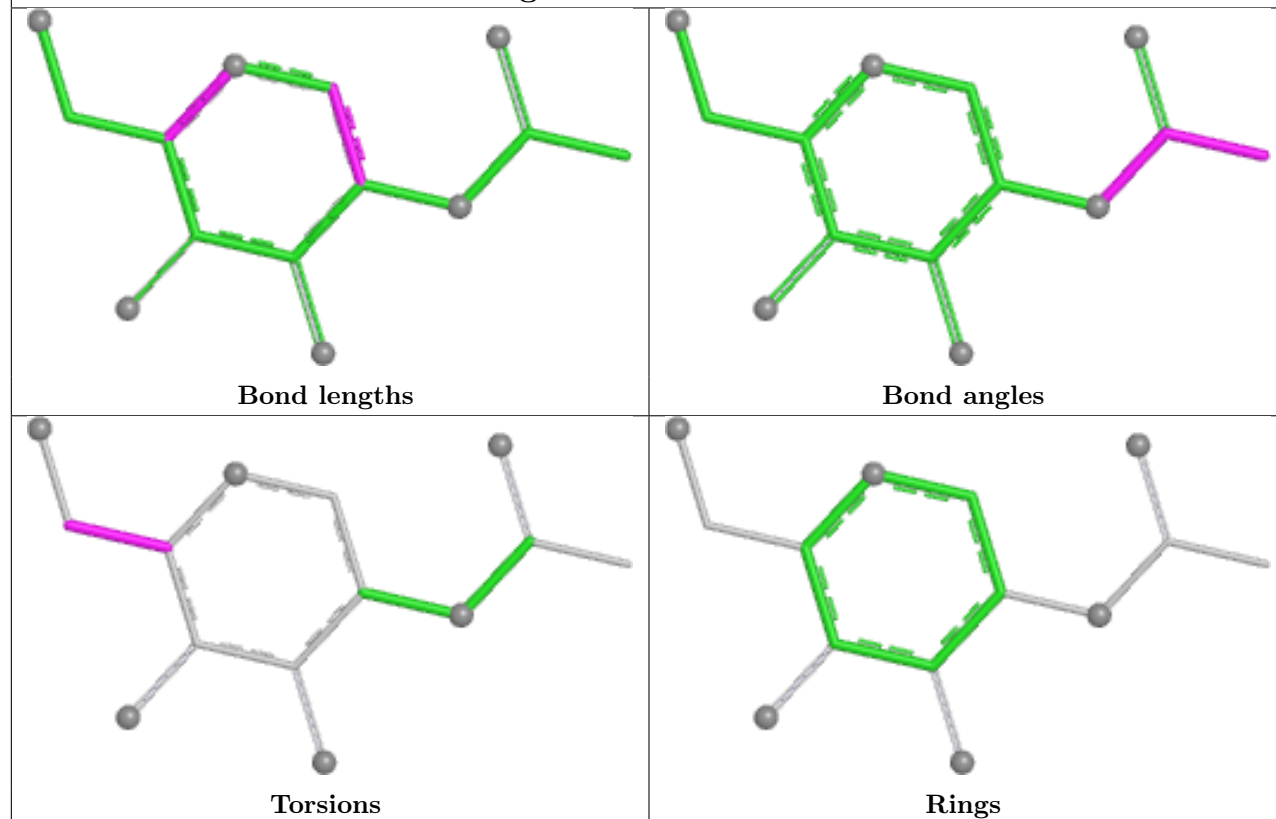




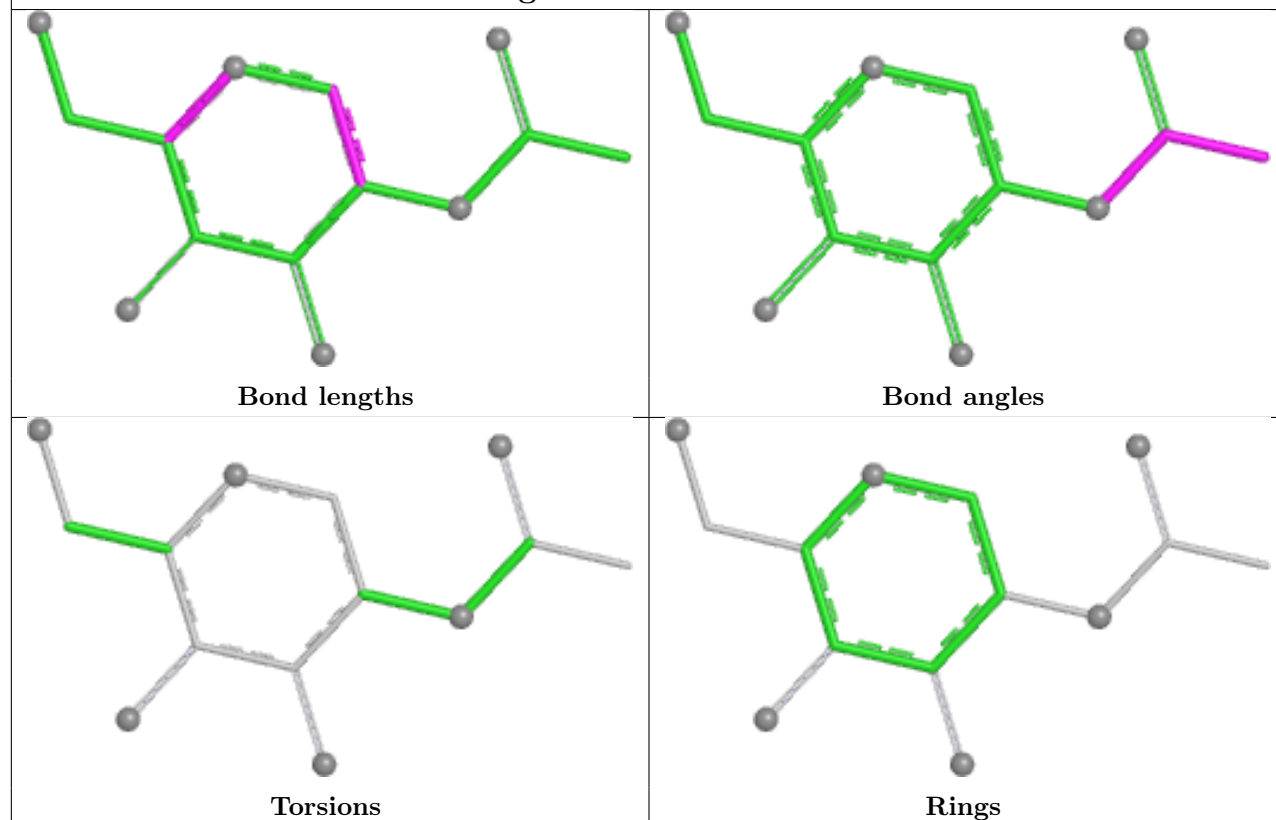


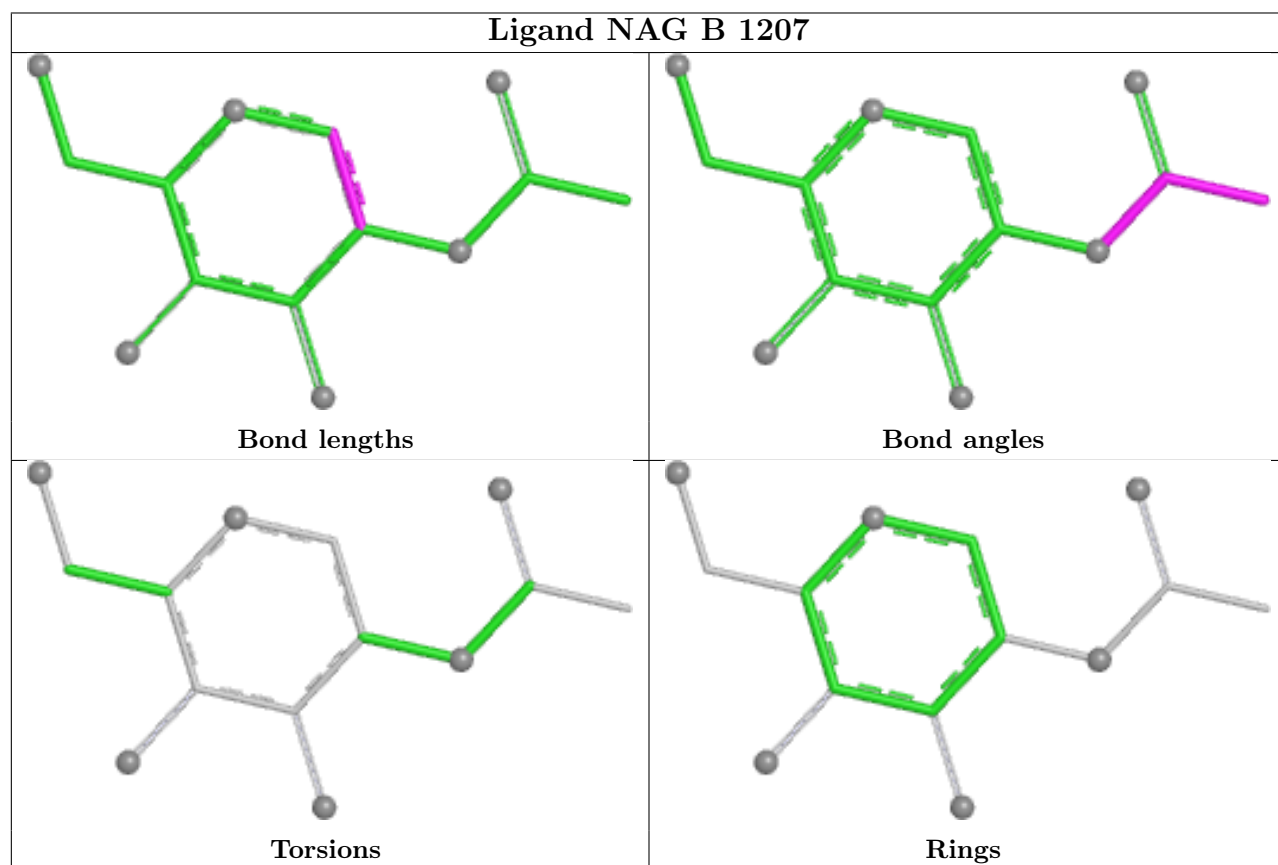
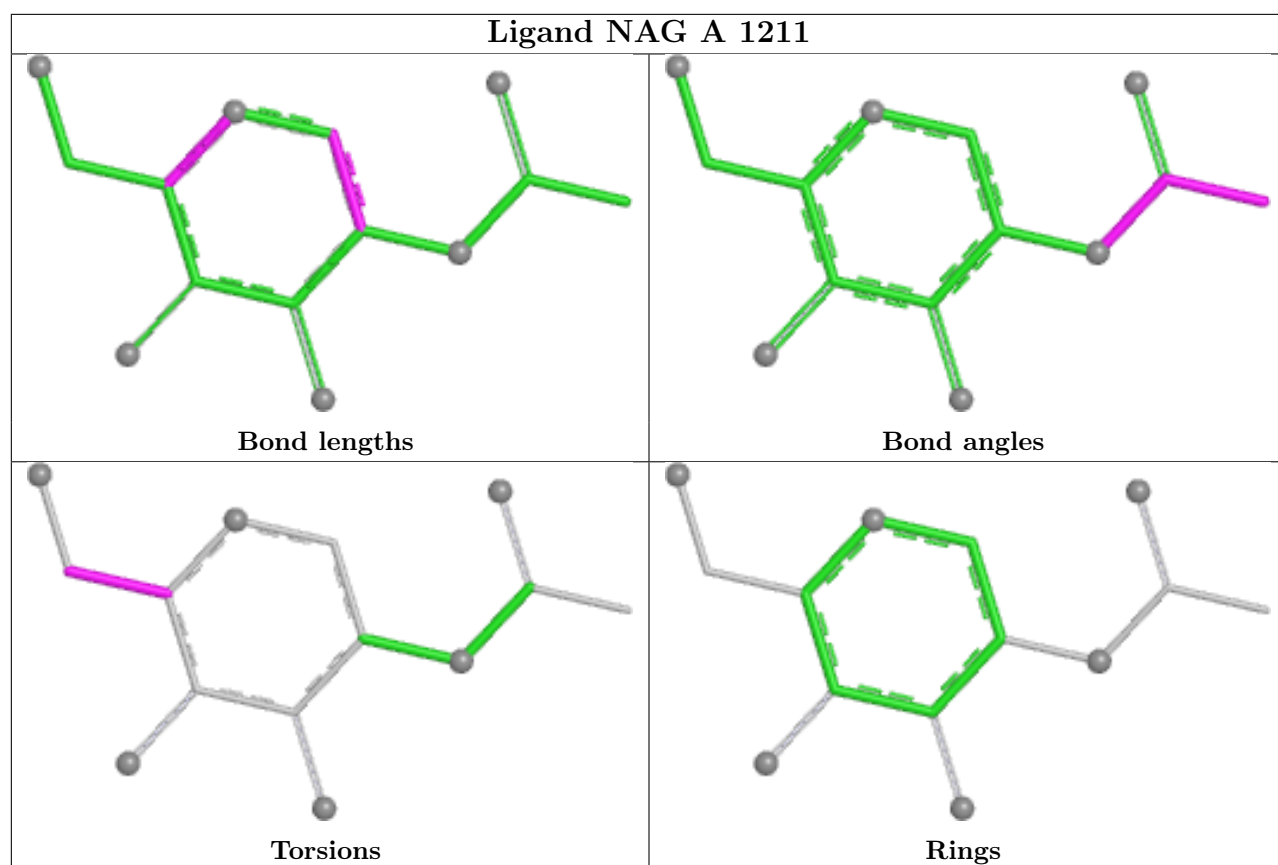


Ligand NAG B 1205

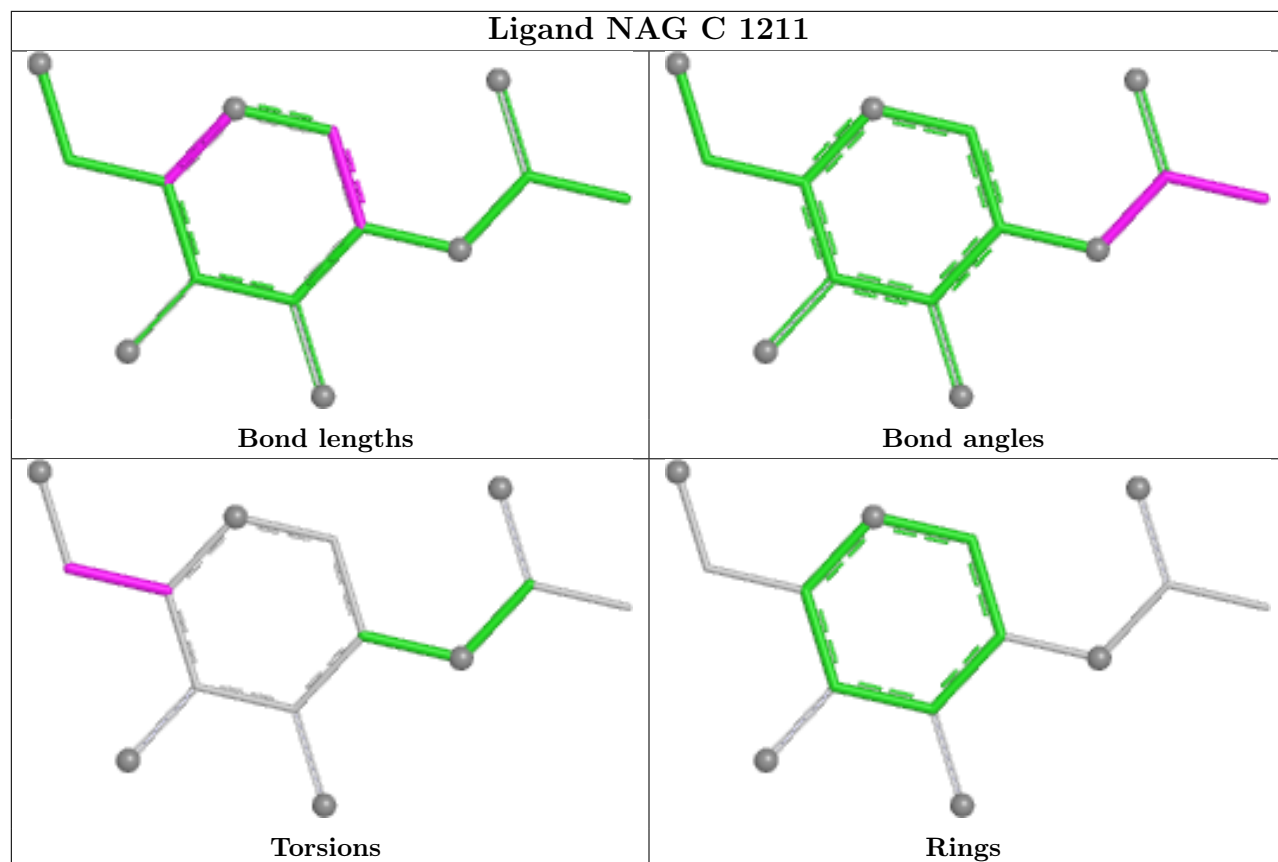


Ligand NAG A 1202

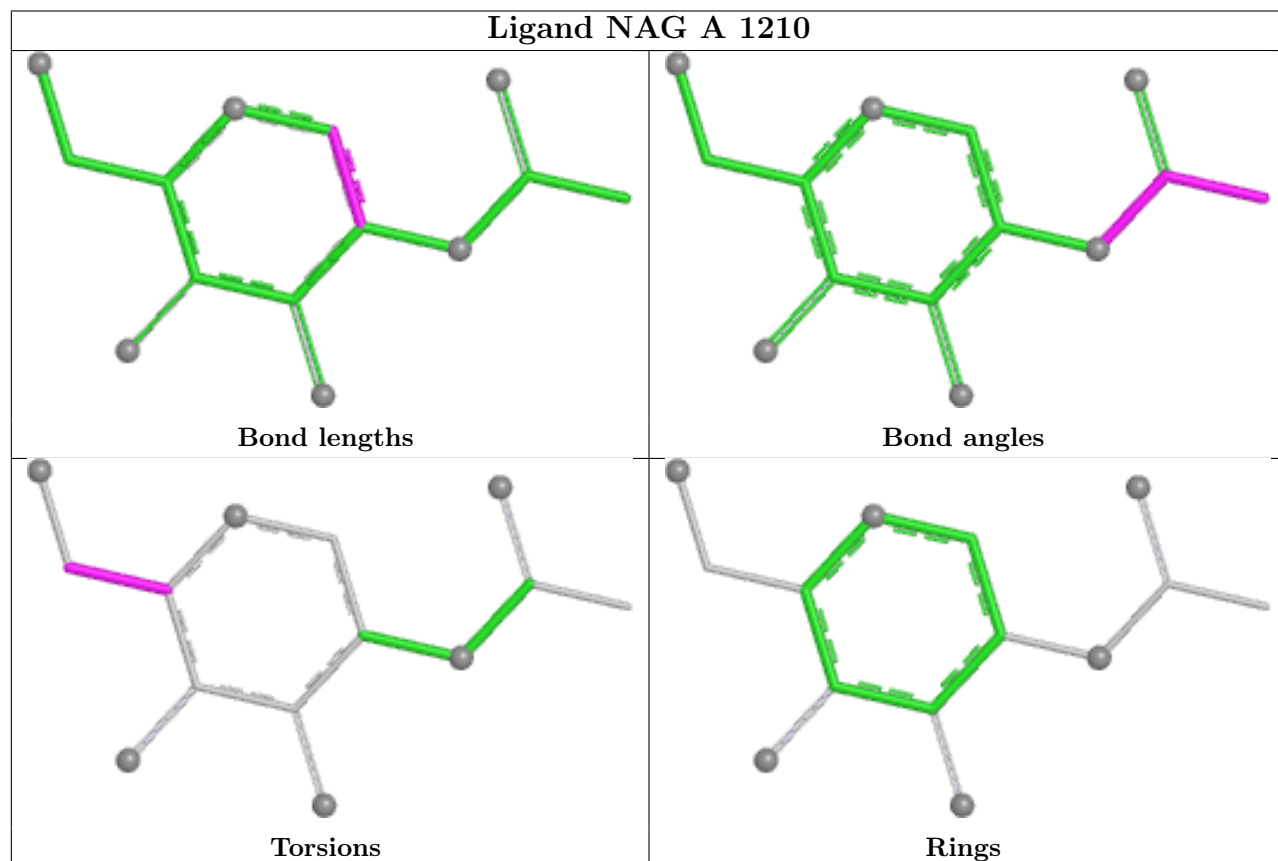




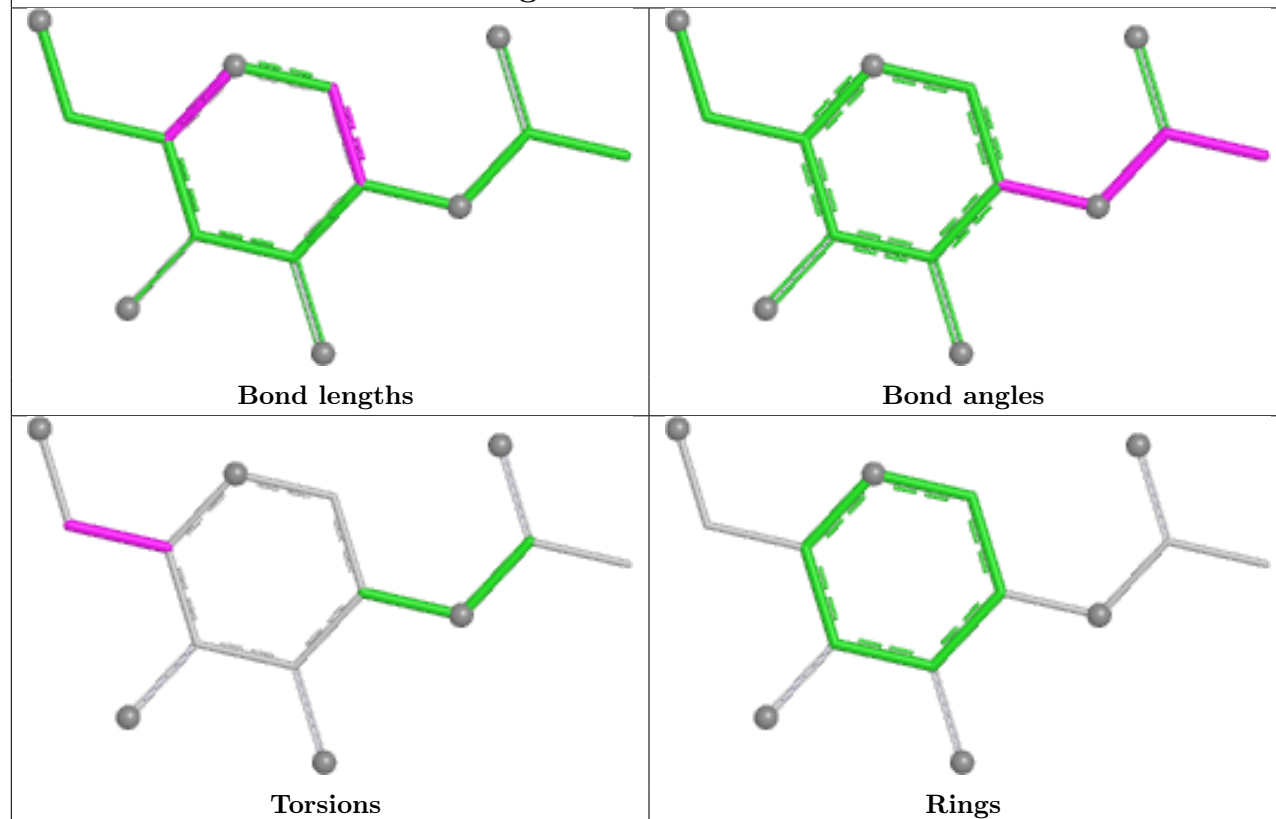
Ligand NAG C 1211



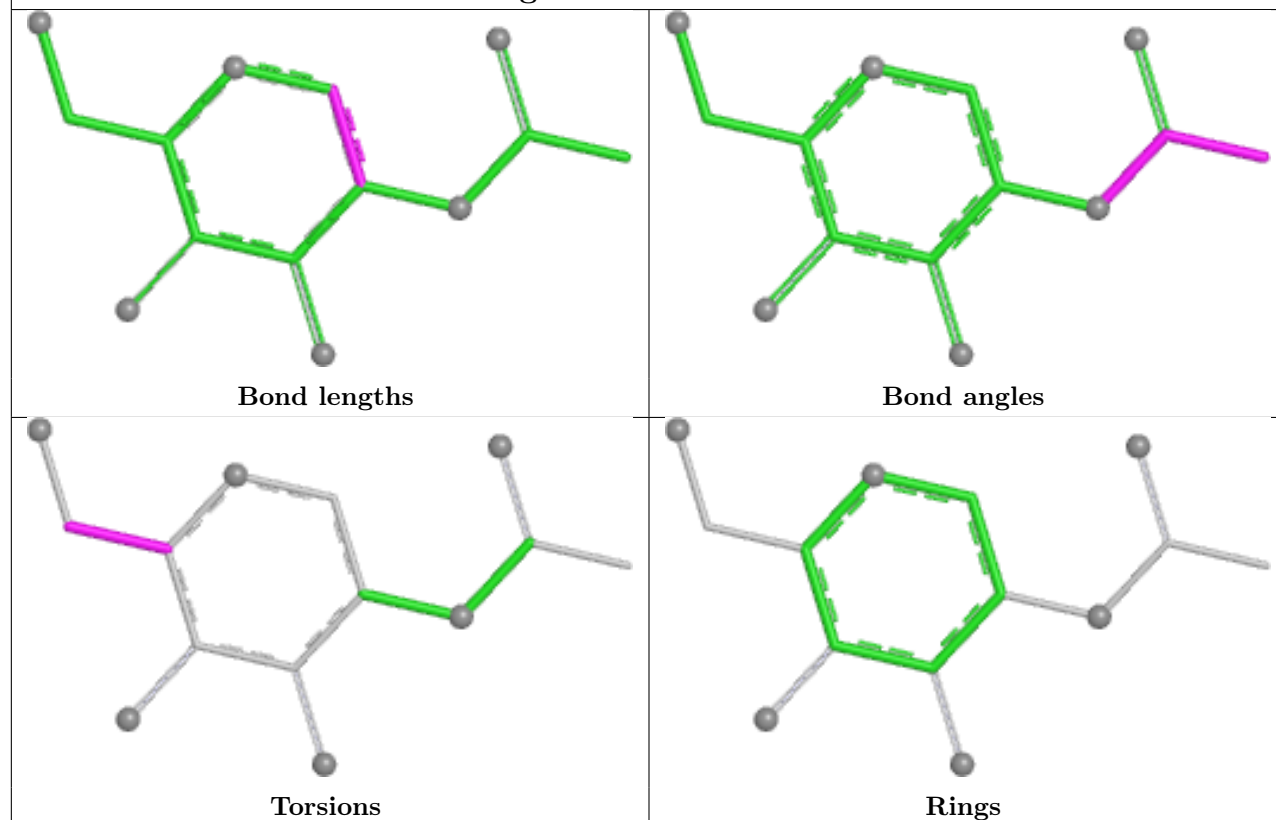
Ligand NAG A 1210



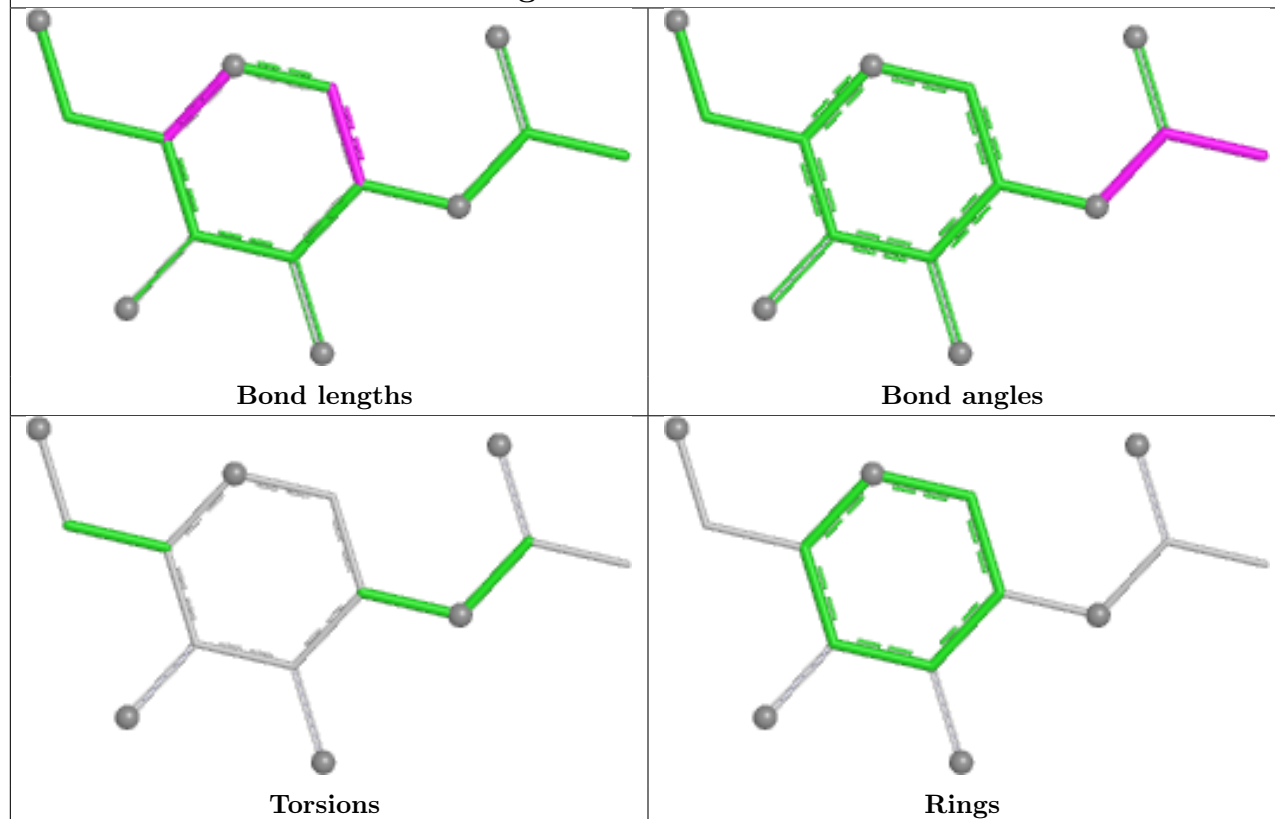
Ligand NAG C 1202



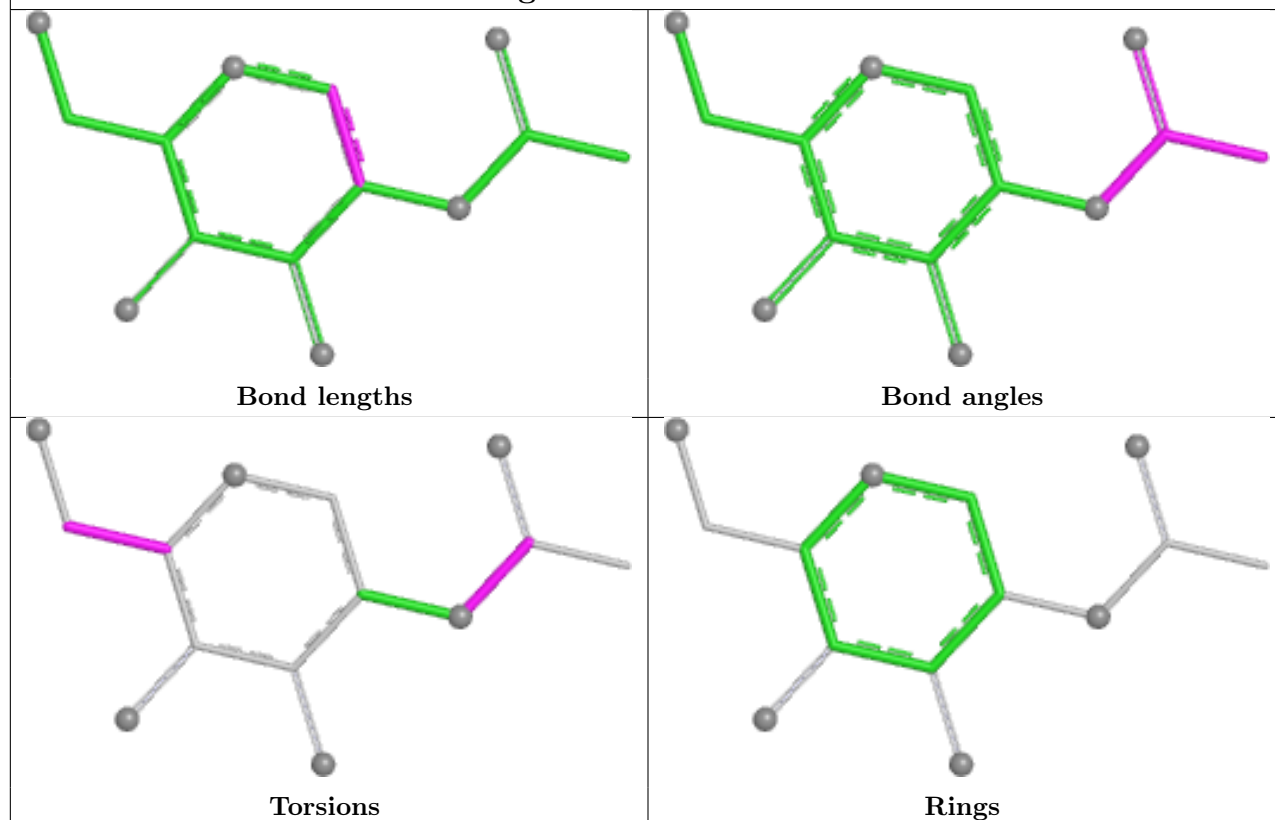
Ligand NAG B 1203

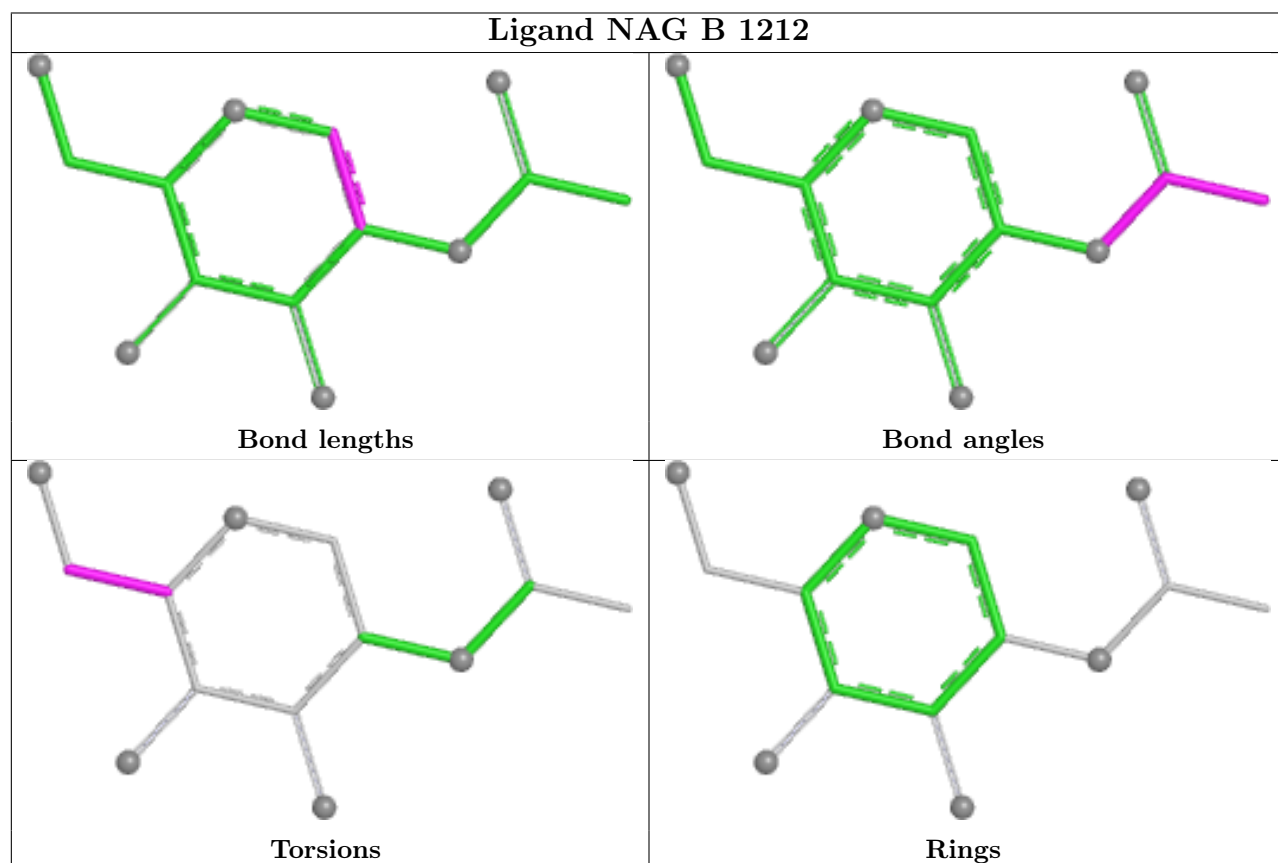
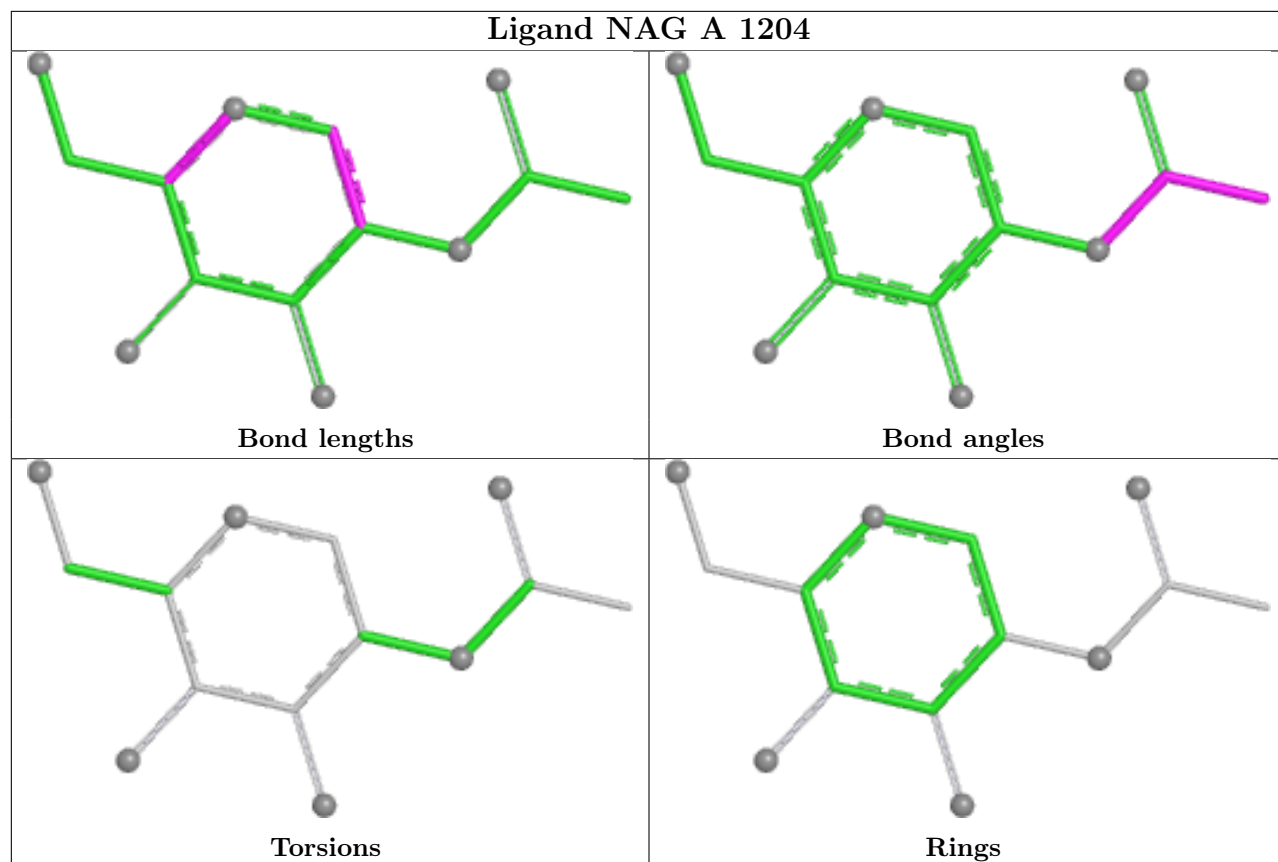


Ligand NAG C 1204

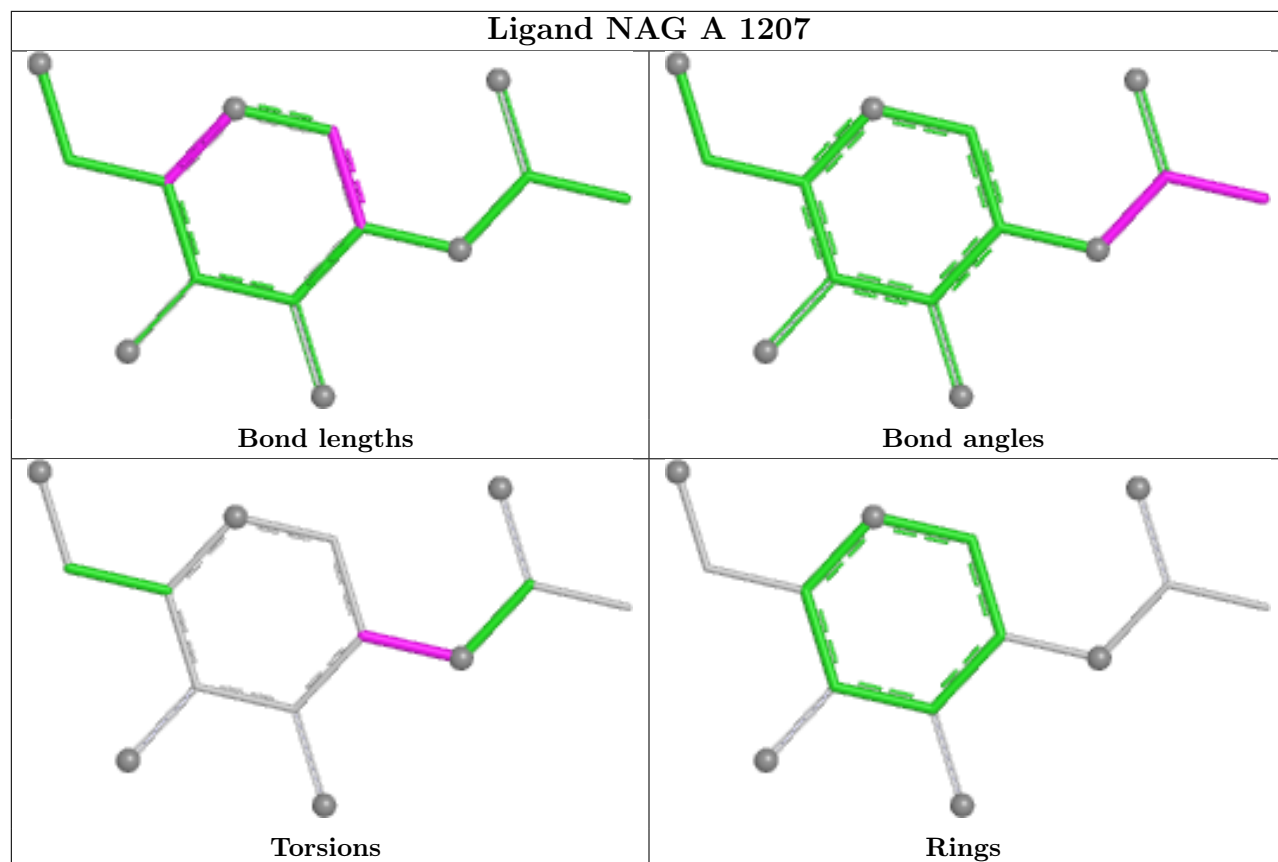


Ligand NAG C 1205

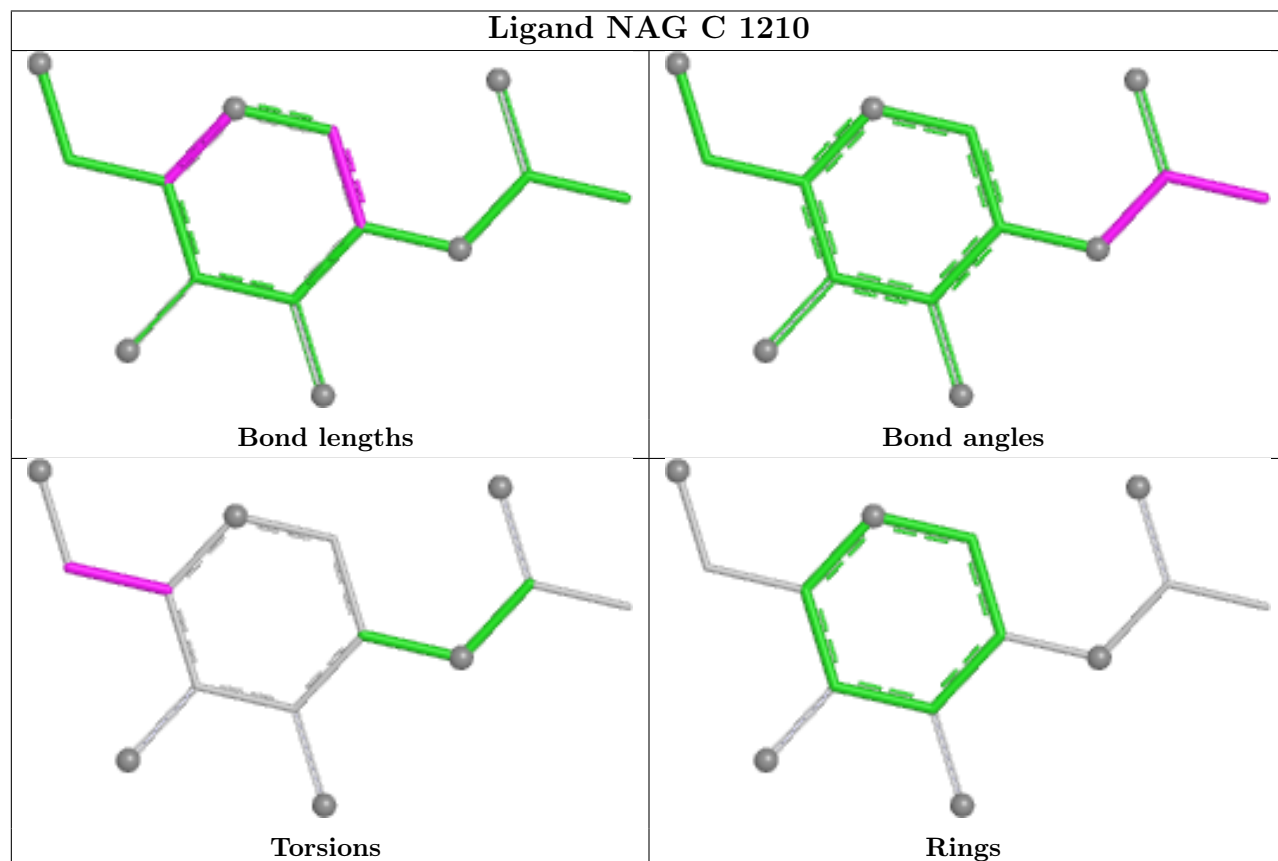




Ligand NAG A 1207



Ligand NAG C 1210



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

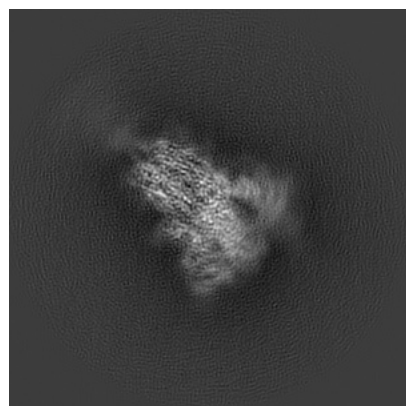
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29456. These allow visual inspection of the internal detail of the map and identification of artifacts.

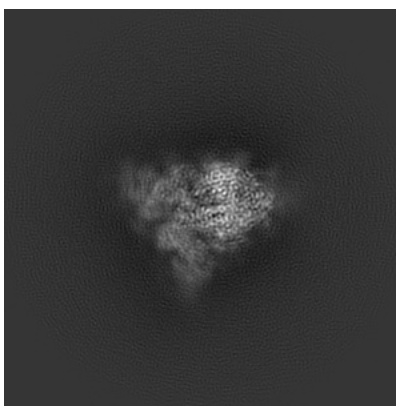
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

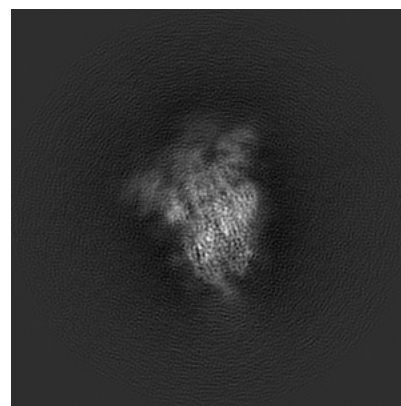
6.1.1 Primary map



X

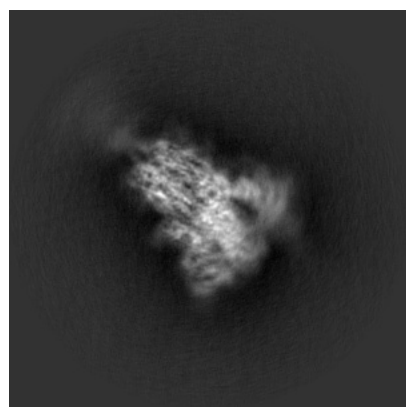


Y

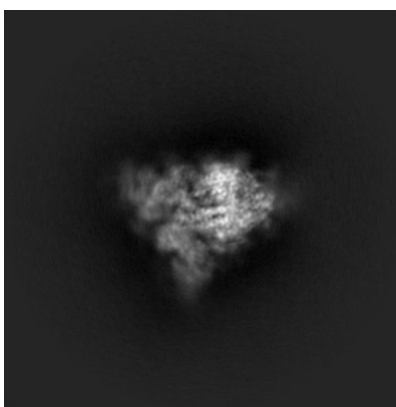


Z

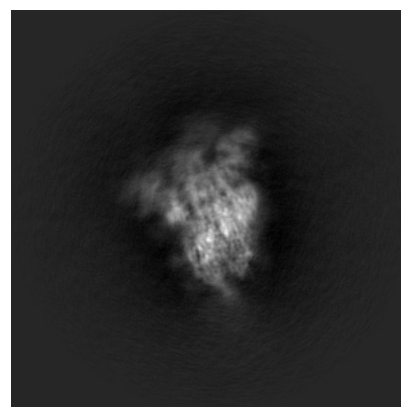
6.1.2 Raw map



X



Y

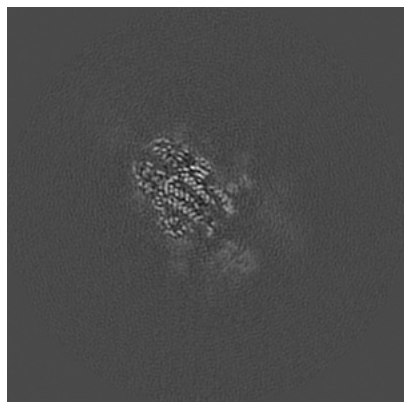


Z

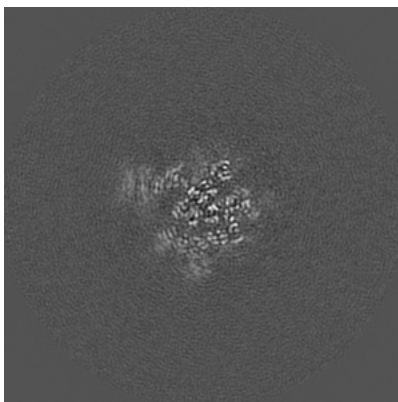
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

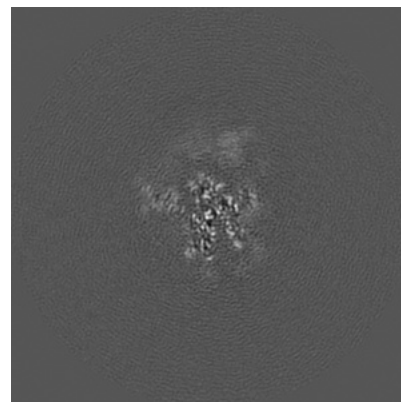
6.2.1 Primary map



X Index: 220

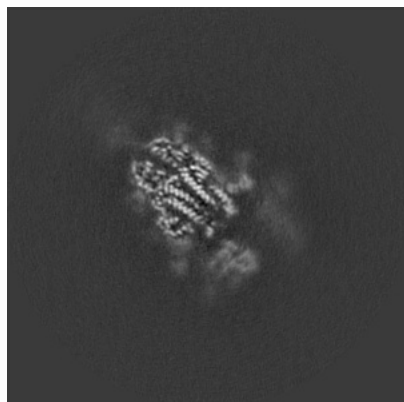


Y Index: 220

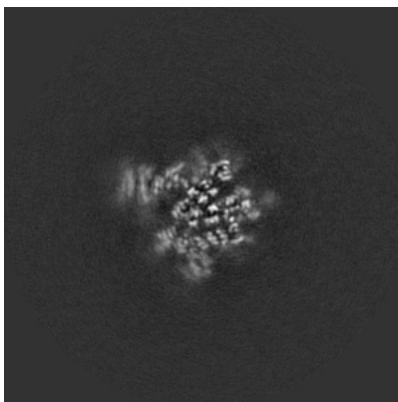


Z Index: 220

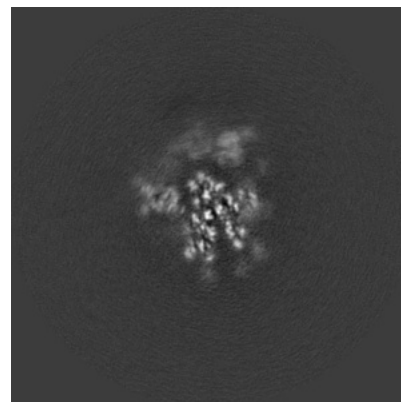
6.2.2 Raw map



X Index: 220



Y Index: 220

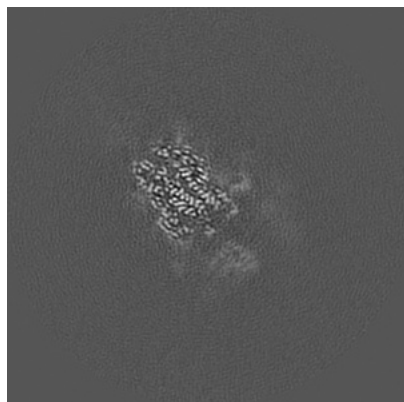


Z Index: 220

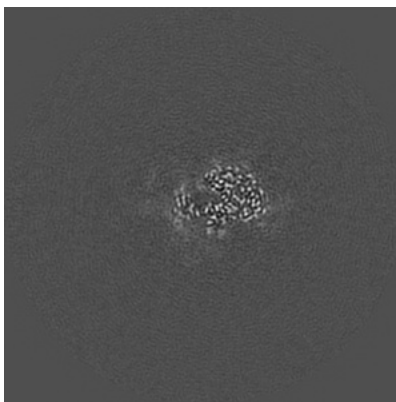
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

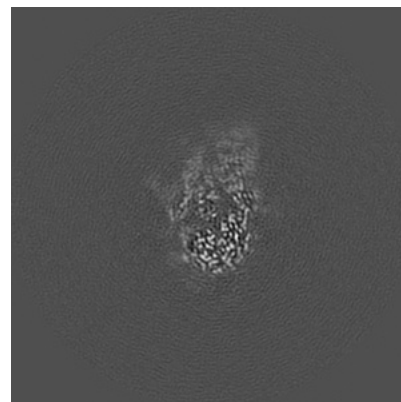
6.3.1 Primary map



X Index: 218

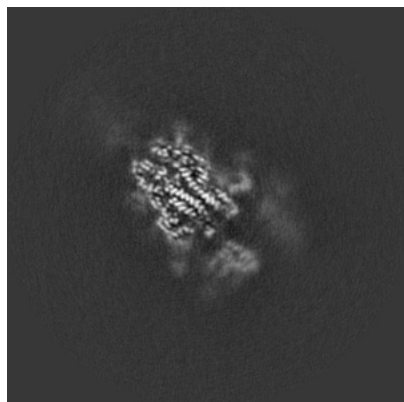


Y Index: 187

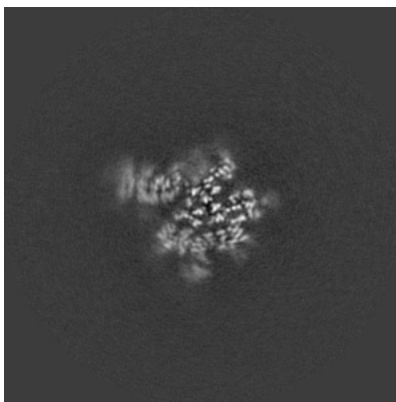


Z Index: 238

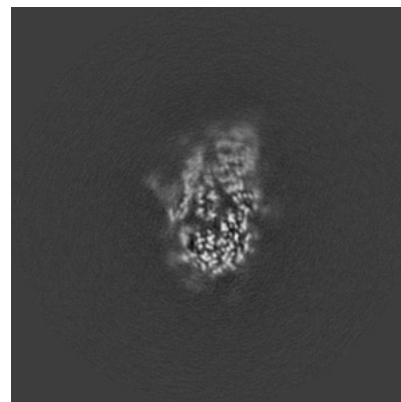
6.3.2 Raw map



X Index: 218



Y Index: 216

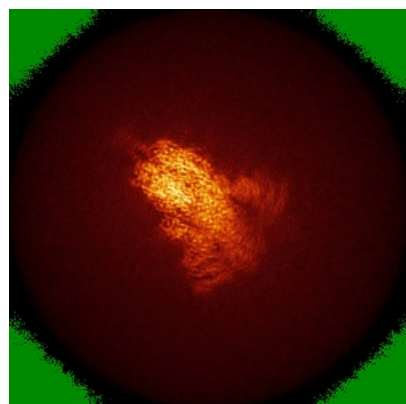


Z Index: 238

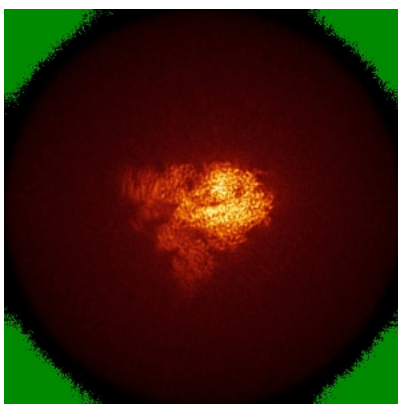
The images above show the largest variance slices of the map in three orthogonal directions.

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

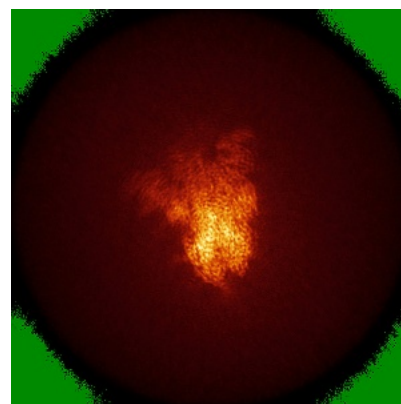
6.4.1 Primary map



X

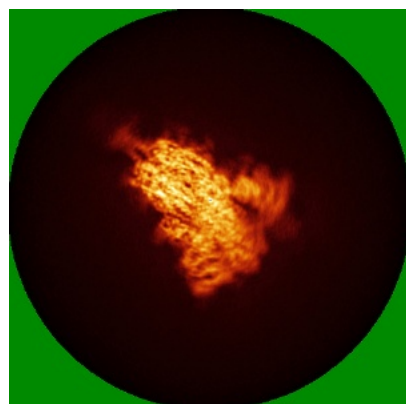


Y

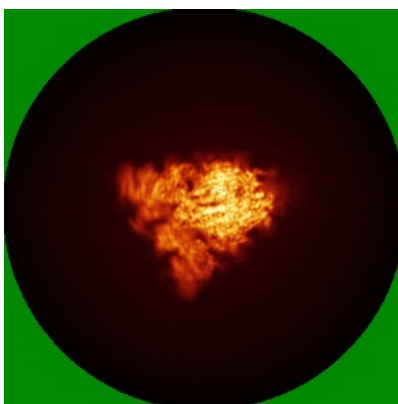


Z

6.4.2 Raw map



X



Y

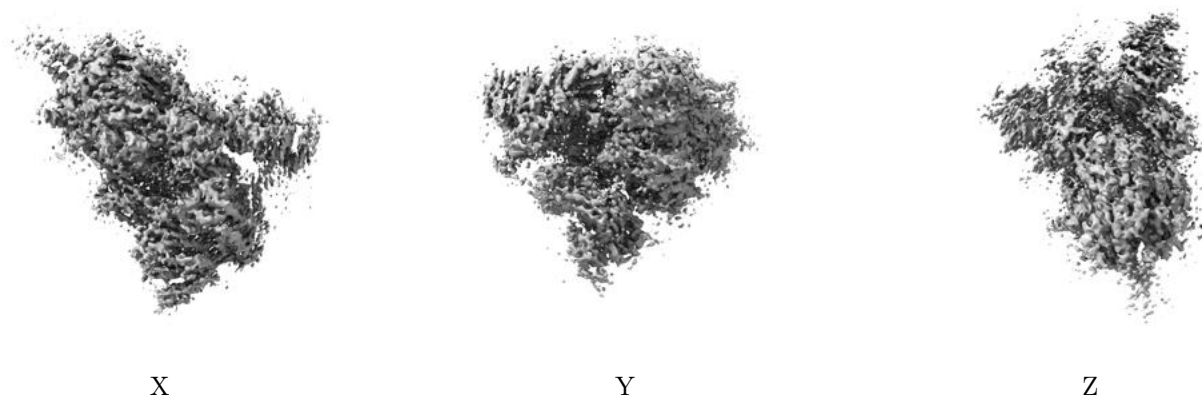


Z

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 0.007. 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.

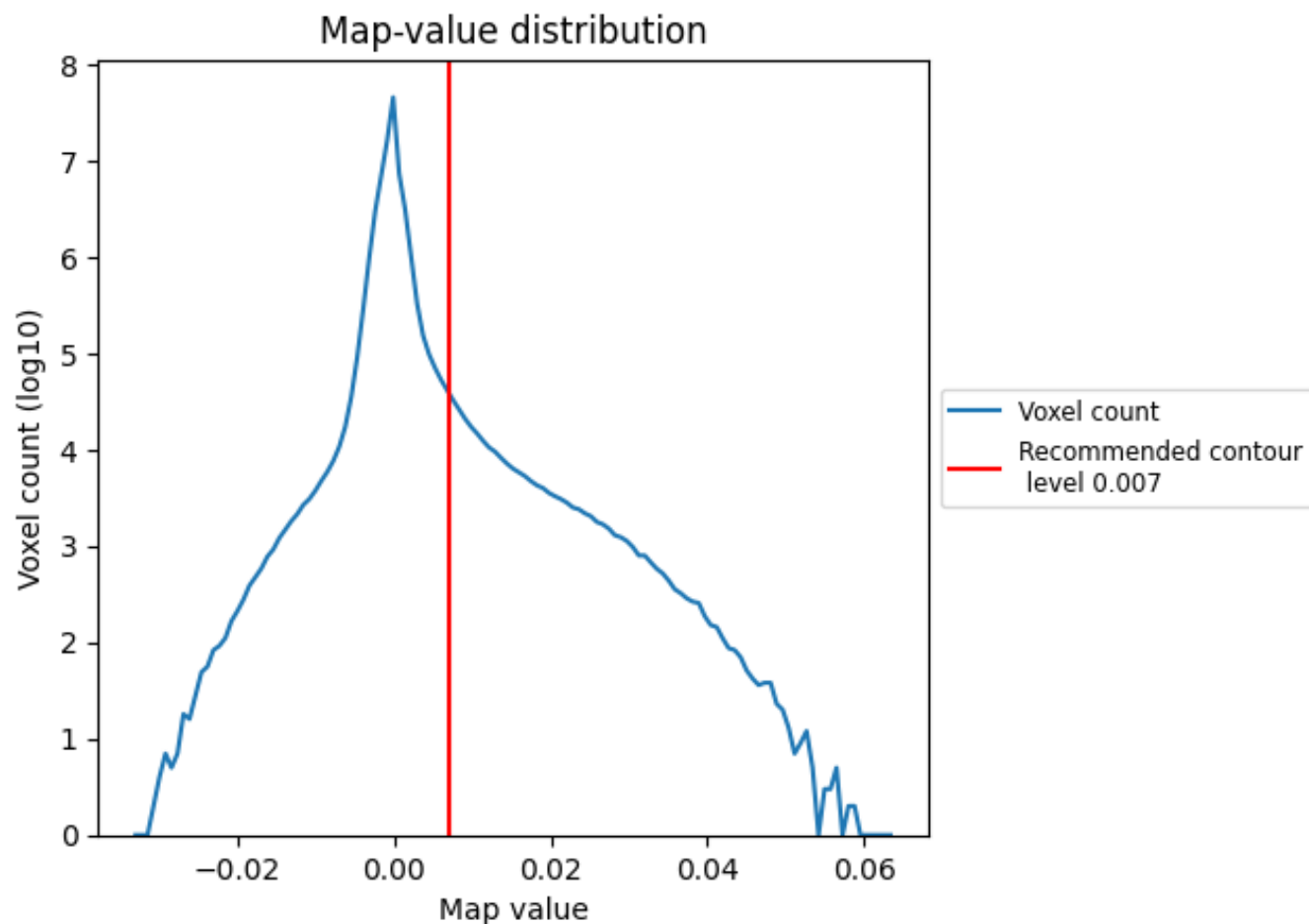
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

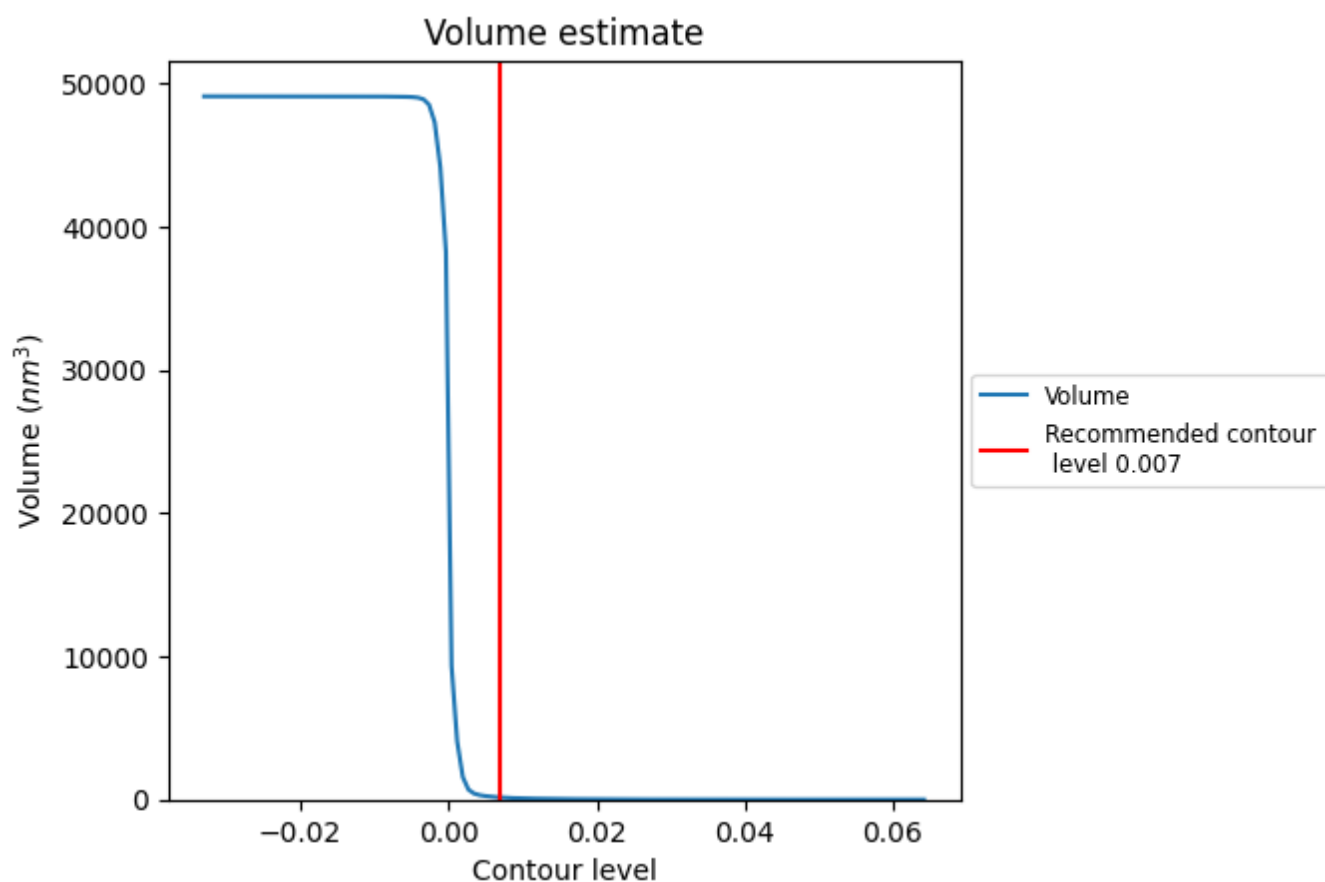
This section contains the results of statistical analysis of the map.

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.

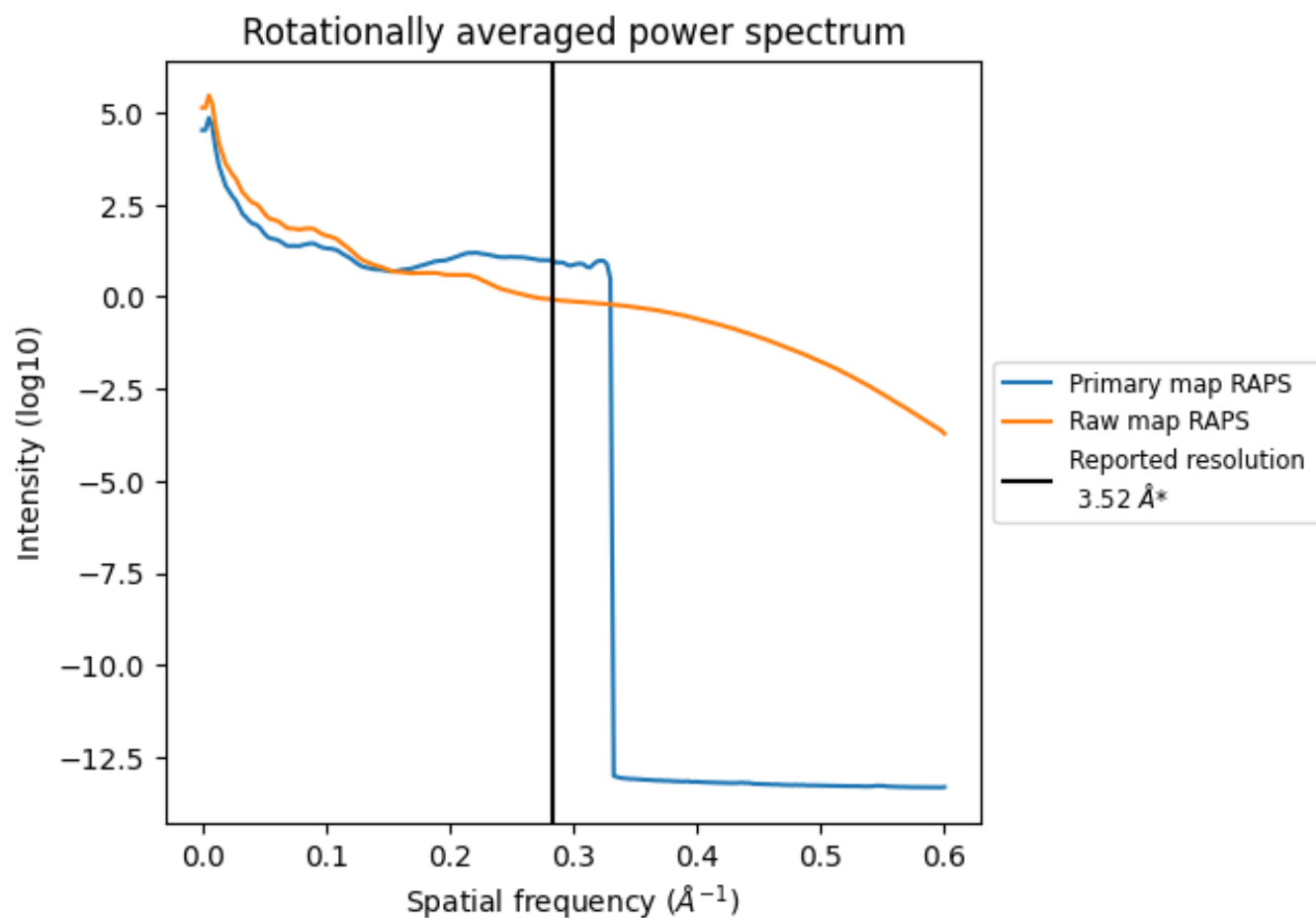
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 148 nm³; this corresponds to an approximate mass of 134 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 ⓘ

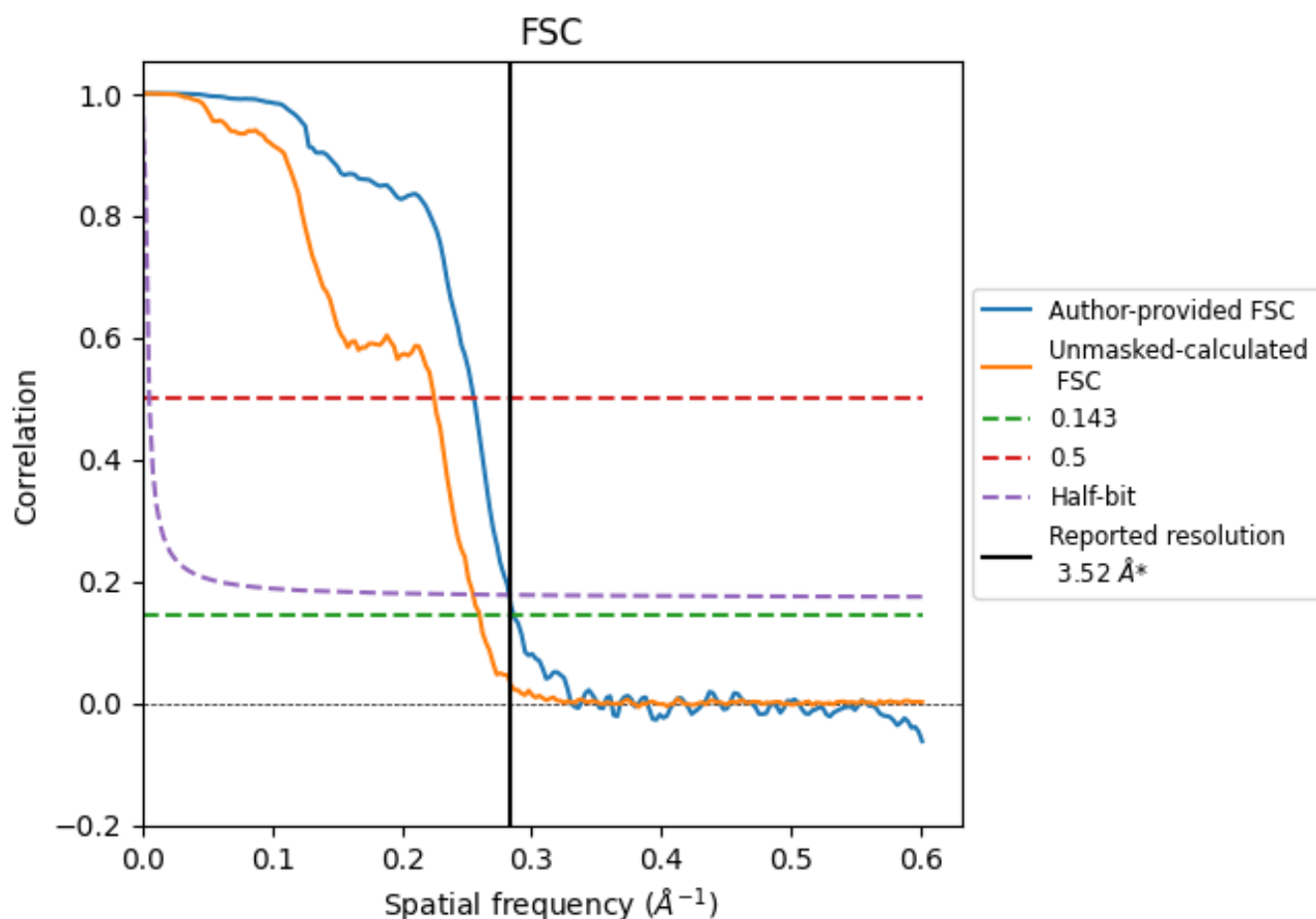


*Reported resolution corresponds to spatial frequency of 0.284 Å⁻¹

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 0.284 Å⁻¹

8.2 Resolution estimates [i](#)

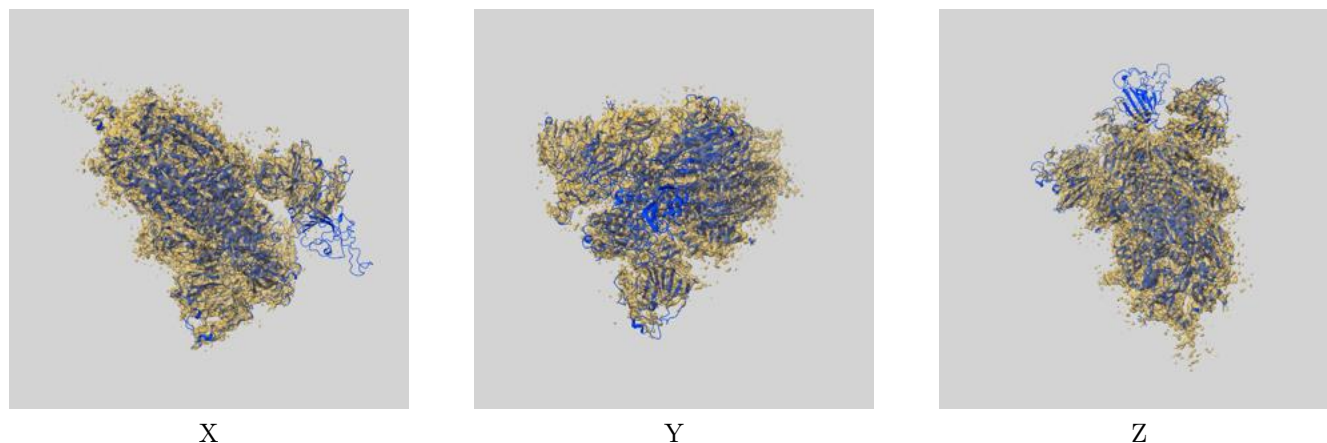
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.52	-	-
Author-provided FSC curve	3.49	3.91	3.54
Unmasked-calculated*	3.85	4.45	3.92

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

9 Map-model fit [i](#)

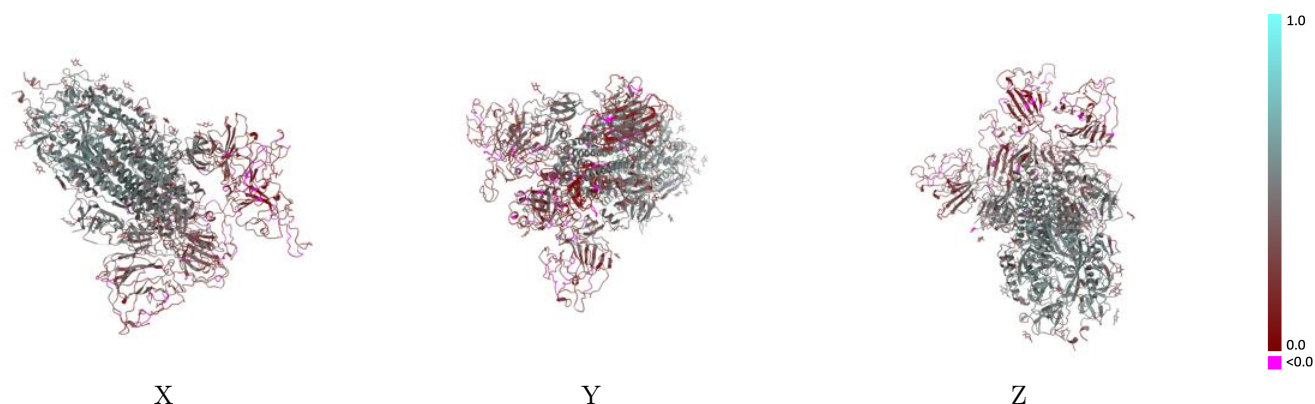
This section contains information regarding the fit between EMDB map EMD-29456 and PDB model 8FU9. Per-residue inclusion information can be found in section [3](#) on page [9](#).

9.1 Map-model overlay [i](#)



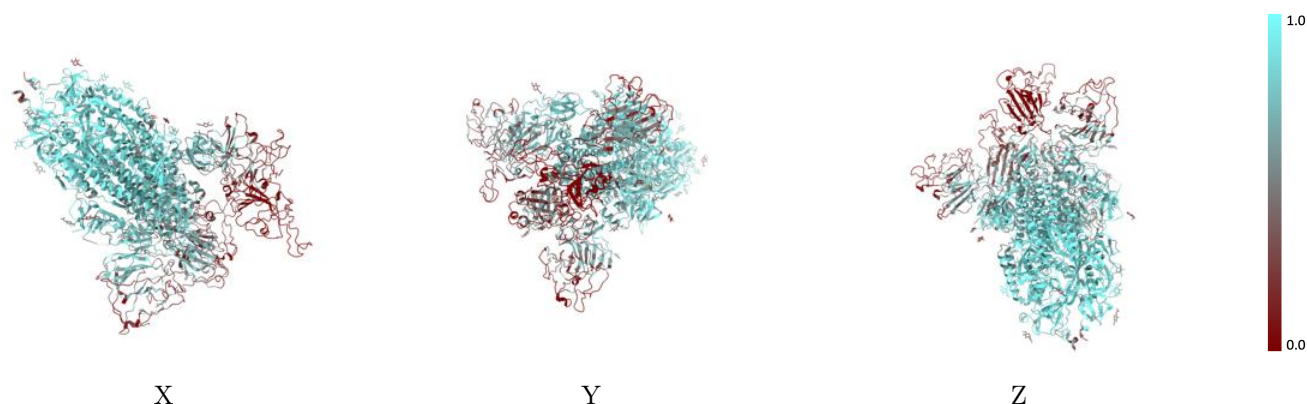
The images above show the 3D surface view of the map at the recommended contour level 0.007 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



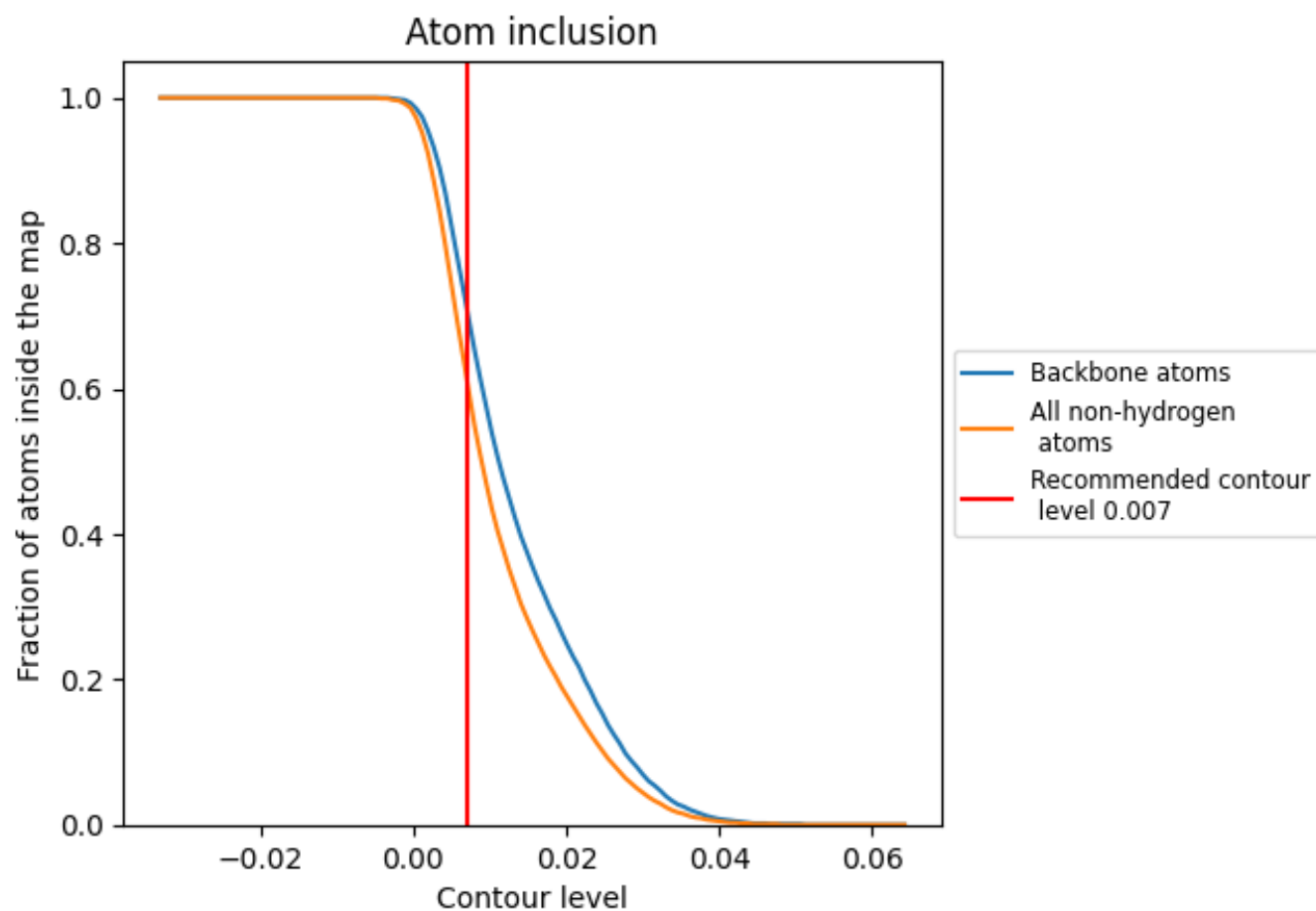
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.007).

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 61% 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 (0.007) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6110	0.3650
A	0.6370	0.3730
B	0.5570	0.3490
C	0.6400	0.3750

