



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 02:42 AM UTC

PDB ID : 9LOY / pdb_00009loy
EMDB ID : EMD-63263
Title : Cryo-EM structure of SARS-CoV-2 JN.1 spike glycoprotein in complex with F61R2-780 Fab
Authors : Wang, X.; Li, X.
Deposited on : 2025-01-23
Resolution : 2.66 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

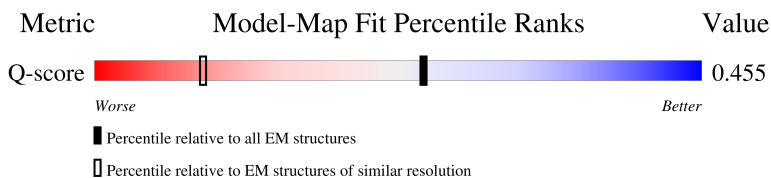
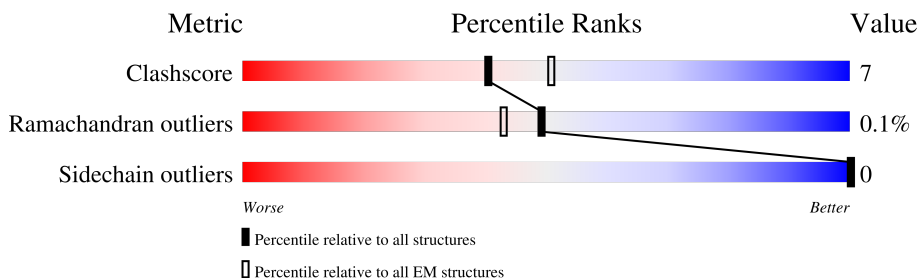
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data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

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.66 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	9119 (2.16 - 3.16)

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	1200	15% 70% 14% 16%
1	B	1200	55% 13% 32%
1	C	1200	17% 71% 13% 15%
2	E	117	100% 79% 21%

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Mol	Chain	Length	Quality of chain
2	H	117	99%
			76%24%
3	F	106	94%
			78%16%6%
3	I	106	94%
			78%16%6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 25971 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	1002	Total	C	N	O	S	0	0
			7884	5071	1300	1477	36		
1	B	818	Total	C	N	O	S	0	0
			6375	4082	1049	1217	27		
1	C	1021	Total	C	N	O	S	0	0
			8006	5136	1326	1508	36		

There are 207 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ILE	THR	variant	UNP P0DTC2
A	24	THR	ARG	conflict	UNP P0DTC2
A	?	-	LEU	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	27	SER	ALA	conflict	UNP P0DTC2
A	50	LEU	SER	conflict	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	127	PHE	VAL	conflict	UNP P0DTC2
A	142	ASP	GLY	variant	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	157	SER	PHE	conflict	UNP P0DTC2
A	158	GLY	ARG	conflict	UNP P0DTC2
A	?	-	ASN	deletion	UNP P0DTC2
A	212	ILE	LEU	conflict	UNP P0DTC2
A	213	GLY	VAL	variant	UNP P0DTC2
A	216	PHE	LEU	conflict	UNP P0DTC2
A	245	ASN	HIS	conflict	UNP P0DTC2
A	264	ASP	ALA	conflict	UNP P0DTC2
A	332	VAL	ILE	conflict	UNP P0DTC2
A	339	HIS	GLY	variant	UNP P0DTC2
A	356	THR	LYS	conflict	UNP P0DTC2
A	371	PHE	SER	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	376	ALA	THR	variant	UNP P0DTC2
A	403	LYS	ARG	conflict	UNP P0DTC2
A	405	ASN	ASP	variant	UNP P0DTC2
A	408	SER	ARG	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	445	HIS	VAL	conflict	UNP P0DTC2
A	446	SER	GLY	variant	UNP P0DTC2
A	450	ASP	ASN	conflict	UNP P0DTC2
A	452	TRP	LEU	conflict	UNP P0DTC2
A	455	SER	LEU	conflict	UNP P0DTC2
A	460	LYS	ASN	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	481	LYS	ASN	conflict	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	484	LYS	GLU	variant	UNP P0DTC2
A	486	PRO	PHE	variant	UNP P0DTC2
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2
A	554	LYS	GLU	conflict	UNP P0DTC2
A	570	VAL	ALA	conflict	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	621	SER	PRO	conflict	UNP P0DTC2
A	655	TYR	HIS	variant	UNP P0DTC2
A	679	LYS	ASN	variant	UNP P0DTC2
A	681	ARG	PRO	variant	UNP P0DTC2
A	682	GLY	ARG	conflict	UNP P0DTC2
A	683	SER	ARG	conflict	UNP P0DTC2
A	685	SER	ARG	conflict	UNP P0DTC2
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	817	PRO	PHE	engineered mutation	UNP P0DTC2
A	892	PRO	ALA	engineered mutation	UNP P0DTC2
A	899	PRO	ALA	engineered mutation	UNP P0DTC2
A	939	PHE	SER	conflict	UNP P0DTC2
A	942	PRO	ALA	engineered mutation	UNP P0DTC2
A	954	HIS	GLN	variant	UNP P0DTC2
A	969	LYS	ASN	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	986	PRO	LYS	engineered mutation	UNP P0DTC2
A	987	PRO	VAL	engineered mutation	UNP P0DTC2
A	1143	LEU	PRO	conflict	UNP P0DTC2
B	22	ILE	THR	variant	UNP P0DTC2
B	24	THR	ARG	conflict	UNP P0DTC2
B	?	-	LEU	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	27	SER	ALA	conflict	UNP P0DTC2
B	50	LEU	SER	conflict	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	127	PHE	VAL	conflict	UNP P0DTC2
B	143	ASP	GLY	variant	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	157	SER	PHE	conflict	UNP P0DTC2
B	158	GLY	ARG	conflict	UNP P0DTC2
B	?	-	ASN	deletion	UNP P0DTC2
B	211	ILE	LEU	conflict	UNP P0DTC2
B	213	GLY	VAL	variant	UNP P0DTC2
B	216	PHE	LEU	conflict	UNP P0DTC2
B	245	ASN	HIS	conflict	UNP P0DTC2
B	264	ASP	ALA	conflict	UNP P0DTC2
B	333	VAL	ILE	conflict	UNP P0DTC2
B	340	HIS	GLY	variant	UNP P0DTC2
B	357	THR	LYS	conflict	UNP P0DTC2
B	372	PHE	SER	variant	UNP P0DTC2
B	374	PRO	SER	variant	UNP P0DTC2
B	376	PHE	SER	variant	UNP P0DTC2
B	377	ALA	THR	variant	UNP P0DTC2
B	404	LYS	ARG	conflict	UNP P0DTC2
B	406	ASN	ASP	variant	UNP P0DTC2
B	409	SER	ARG	variant	UNP P0DTC2
B	418	ASN	LYS	variant	UNP P0DTC2
B	441	LYS	ASN	variant	UNP P0DTC2
B	446	HIS	VAL	conflict	UNP P0DTC2
B	447	SER	GLY	variant	UNP P0DTC2
B	451	ASP	ASN	conflict	UNP P0DTC2
B	453	TRP	LEU	conflict	UNP P0DTC2
B	456	SER	LEU	conflict	UNP P0DTC2
B	461	LYS	ASN	variant	UNP P0DTC2
B	478	ASN	SER	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	479	LYS	THR	variant	UNP P0DTC2
B	482	LYS	ASN	conflict	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	484	LYS	GLU	variant	UNP P0DTC2
B	486	PRO	PHE	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	505	HIS	TYR	variant	UNP P0DTC2
B	554	LYS	GLU	conflict	UNP P0DTC2
B	570	VAL	ALA	conflict	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	621	SER	PRO	conflict	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	679	LYS	ASN	variant	UNP P0DTC2
B	681	ARG	PRO	variant	UNP P0DTC2
B	682	GLY	ARG	conflict	UNP P0DTC2
B	683	SER	ARG	conflict	UNP P0DTC2
B	685	SER	ARG	conflict	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	817	PRO	PHE	engineered mutation	UNP P0DTC2
B	892	PRO	ALA	engineered mutation	UNP P0DTC2
B	899	PRO	ALA	engineered mutation	UNP P0DTC2
B	939	PHE	SER	conflict	UNP P0DTC2
B	942	PRO	ALA	engineered mutation	UNP P0DTC2
B	954	HIS	GLN	variant	UNP P0DTC2
B	969	LYS	ASN	variant	UNP P0DTC2
B	986	PRO	LYS	engineered mutation	UNP P0DTC2
B	987	PRO	VAL	engineered mutation	UNP P0DTC2
B	1143	LEU	PRO	conflict	UNP P0DTC2
C	22	ILE	THR	variant	UNP P0DTC2
C	24	THR	ARG	conflict	UNP P0DTC2
C	?	-	LEU	deletion	UNP P0DTC2
C	?	-	PRO	deletion	UNP P0DTC2
C	?	-	PRO	deletion	UNP P0DTC2
C	27	SER	ALA	conflict	UNP P0DTC2
C	50	LEU	SER	conflict	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	127	PHE	VAL	conflict	UNP P0DTC2
C	142	ASP	GLY	variant	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	157	SER	PHE	conflict	UNP P0DTC2
C	158	GLY	ARG	conflict	UNP P0DTC2
C	?	-	ASN	deletion	UNP P0DTC2
C	211	ILE	LEU	conflict	UNP P0DTC2
C	213	GLY	VAL	variant	UNP P0DTC2
C	216	PHE	LEU	conflict	UNP P0DTC2
C	245	ASN	HIS	conflict	UNP P0DTC2
C	264	ASP	ALA	conflict	UNP P0DTC2
C	332	VAL	ILE	conflict	UNP P0DTC2
C	339	HIS	GLY	variant	UNP P0DTC2
C	356	THR	LYS	conflict	UNP P0DTC2
C	371	PHE	SER	variant	UNP P0DTC2
C	373	PRO	SER	variant	UNP P0DTC2
C	375	PHE	SER	variant	UNP P0DTC2
C	376	ALA	THR	variant	UNP P0DTC2
C	403	LYS	ARG	conflict	UNP P0DTC2
C	405	ASN	ASP	variant	UNP P0DTC2
C	408	SER	ARG	variant	UNP P0DTC2
C	417	ASN	LYS	variant	UNP P0DTC2
C	440	LYS	ASN	variant	UNP P0DTC2
C	445	HIS	VAL	conflict	UNP P0DTC2
C	446	SER	GLY	variant	UNP P0DTC2
C	450	ASP	ASN	conflict	UNP P0DTC2
C	452	TRP	LEU	conflict	UNP P0DTC2
C	455	SER	LEU	conflict	UNP P0DTC2
C	460	LYS	ASN	variant	UNP P0DTC2
C	477	ASN	SER	variant	UNP P0DTC2
C	478	LYS	THR	variant	UNP P0DTC2
C	481	LYS	ASN	conflict	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	484	LYS	GLU	variant	UNP P0DTC2
C	486	PRO	PHE	variant	UNP P0DTC2
C	498	ARG	GLN	variant	UNP P0DTC2
C	501	TYR	ASN	variant	UNP P0DTC2
C	505	HIS	TYR	variant	UNP P0DTC2
C	554	LYS	GLU	conflict	UNP P0DTC2
C	570	VAL	ALA	conflict	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	621	SER	PRO	conflict	UNP P0DTC2
C	655	TYR	HIS	variant	UNP P0DTC2
C	679	LYS	ASN	variant	UNP P0DTC2
C	681	ARG	PRO	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	682	GLY	ARG	conflict	UNP P0DTC2
C	683	SER	ARG	conflict	UNP P0DTC2
C	685	SER	ARG	conflict	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2
C	817	PRO	PHE	engineered mutation	UNP P0DTC2
C	892	PRO	ALA	engineered mutation	UNP P0DTC2
C	899	PRO	ALA	engineered mutation	UNP P0DTC2
C	939	PHE	SER	conflict	UNP P0DTC2
C	942	PRO	ALA	engineered mutation	UNP P0DTC2
C	954	HIS	GLN	variant	UNP P0DTC2
C	969	LYS	ASN	variant	UNP P0DTC2
C	986	PRO	LYS	engineered mutation	UNP P0DTC2
C	987	PRO	VAL	engineered mutation	UNP P0DTC2
C	1143	LEU	PRO	conflict	UNP P0DTC2

- Molecule 2 is a protein called F61R2-780 heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	117	Total	C	N	O	S	0	0
			914	576	155	178	5		
2	H	117	Total	C	N	O	S	0	0
			914	576	155	178	5		

- Molecule 3 is a protein called F61R2-780 light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	F	100	Total	C	N	O	S	0	0
			764	483	130	149	2		
3	I	100	Total	C	N	O	S	0	0
			764	483	130	149	2		

- Molecule 4 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
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	

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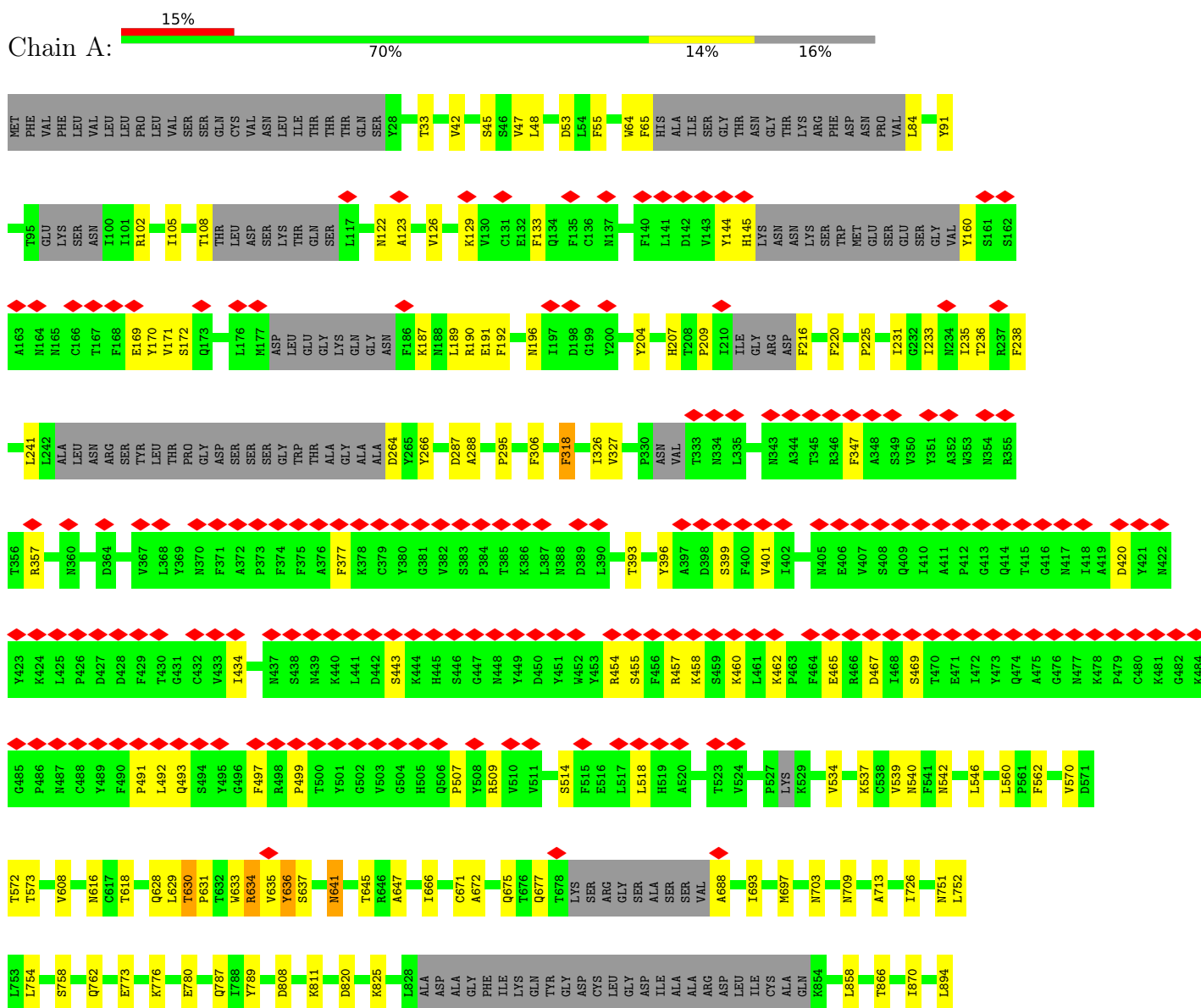
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Mol	Chain	Residues	Atoms				AltConf
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C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

3 Residue-property plots

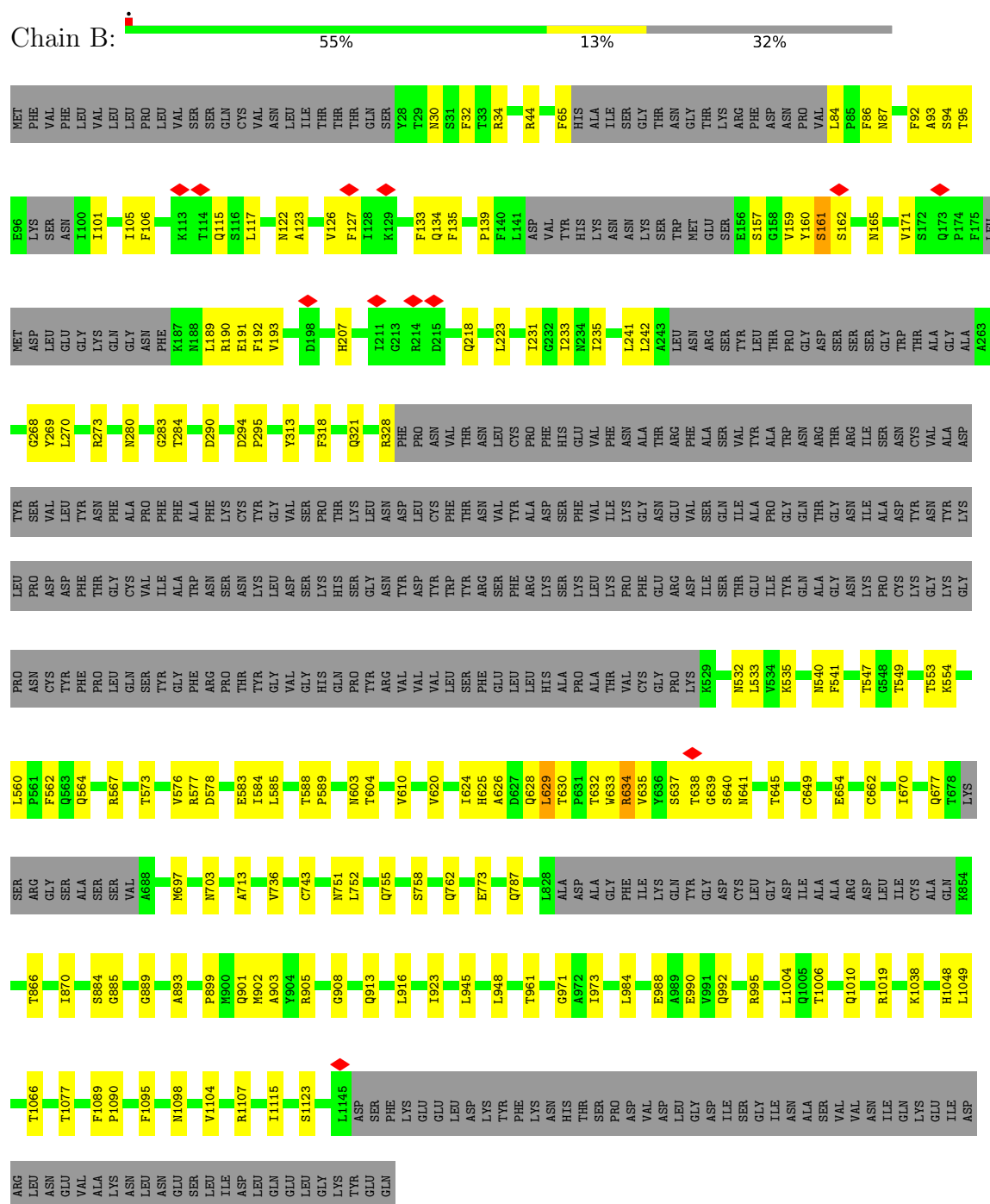
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Spike glycoprotein



TYR
GLU
GLN

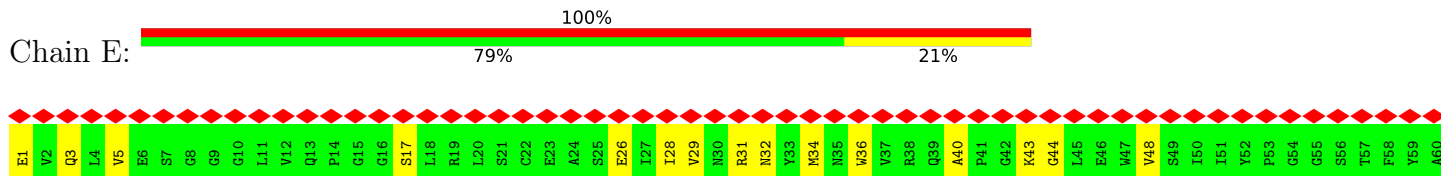
- Molecule 1: Spike glycoprotein

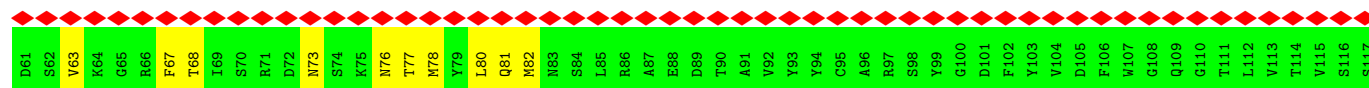


Chain C:

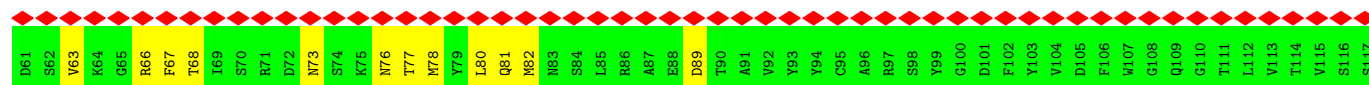
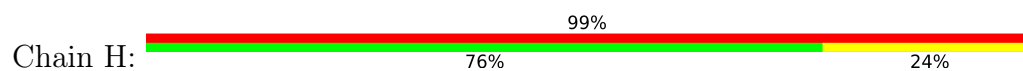


Chain E:

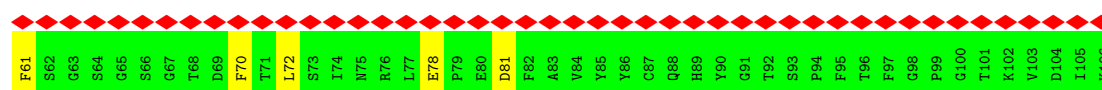
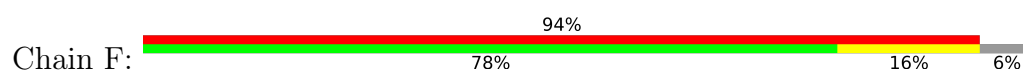




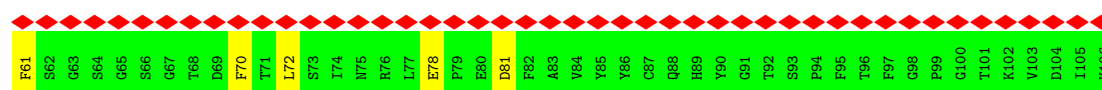
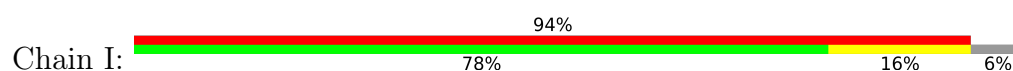
• Molecule 2: F61R2-780 heavy chain



• Molecule 3: F61R2-780 light chain



• Molecule 3: F61R2-780 light chain



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	499045	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	4.509	Depositor
Minimum map value	-2.598	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.068	Depositor
Recommended contour level	0.286	Depositor
Map size ($\text{\AA}$)	388.0, 388.0, 388.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.97, 0.97, 0.97	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.54	0/8077	0.66	12/10988 (0.1%)
1	B	0.63	3/6513 (0.0%)	0.76	16/8861 (0.2%)
1	C	0.61	0/8199	0.75	14/11157 (0.1%)
2	E	0.41	0/934	0.67	1/1264 (0.1%)
2	H	0.41	0/934	0.67	1/1264 (0.1%)
3	F	0.29	0/785	0.53	2/1066 (0.2%)
3	I	0.29	0/785	0.53	2/1066 (0.2%)
All	All	0.57	3/26227 (0.0%)	0.71	48/35666 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	899	PRO	C-O	-8.42	1.13	1.24
1	B	295	PRO	C-O	-6.28	1.16	1.24
1	B	313	TYR	C-O	-5.36	1.17	1.24

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	631	PRO	N-CA-C	12.63	131.53	113.47
1	C	631	PRO	N-CA-C	-8.61	101.56	113.53
1	C	83	VAL	CA-C-N	8.14	133.06	120.68
1	C	83	VAL	C-N-CA	8.14	133.06	120.68
1	A	641	ASN	CB-CA-C	7.42	123.46	110.85
1	A	635	VAL	CA-C-N	-7.35	112.76	123.05
1	A	635	VAL	C-N-CA	-7.35	112.76	123.05
1	B	743	CYS	CB-CA-C	-7.10	99.59	111.02
1	B	630	THR	CB-CA-C	6.87	123.71	110.17
1	C	631	PRO	CB-CA-C	6.83	123.90	112.62
1	C	318	PHE	CA-CB-CG	6.69	120.49	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	634	ARG	N-CA-C	-6.52	100.52	110.30
1	B	649	CYS	CB-CA-C	-6.36	99.25	109.75
1	A	637	SER	N-CA-C	-6.21	97.58	110.80
1	C	887	THR	CB-CA-C	6.18	121.36	110.85
1	B	751	ASN	CB-CA-C	6.13	121.28	110.85
1	C	1098	ASN	CA-CB-CG	6.12	118.72	112.60
1	B	318	PHE	CA-CB-CG	6.03	119.83	113.80
1	B	161	SER	N-CA-C	-5.91	106.23	113.50
2	H	73	ASN	CB-CA-C	-5.64	101.26	110.85
1	B	638	THR	N-CA-C	-5.64	105.22	111.36
2	E	73	ASN	CB-CA-C	-5.62	101.29	110.85
1	C	914	ASN	CA-CB-CG	5.60	118.20	112.60
1	A	1038	LYS	N-CA-C	-5.52	106.54	113.72
1	C	748	GLU	N-CA-C	-5.51	105.36	111.36
1	C	219	GLY	CA-C-O	-5.50	118.49	122.45
1	C	801	ASN	CA-CB-CG	5.48	118.08	112.60
1	B	157	SER	N-CA-C	-5.45	106.00	113.30
1	B	294	ASP	CA-CB-CG	5.43	118.03	112.60
1	B	1098	ASN	CA-CB-CG	5.41	118.01	112.60
1	B	899	PRO	N-CA-CB	-5.38	97.41	103.33
1	A	709	ASN	CA-CB-CG	5.37	117.97	112.60
1	B	634	ARG	CA-C-N	-5.35	112.95	120.77
1	B	634	ARG	C-N-CA	-5.35	112.95	120.77
1	A	318	PHE	CA-CB-CG	5.26	119.06	113.80
1	C	294	ASP	CA-CB-CG	5.26	117.86	112.60
1	A	636	TYR	CB-CA-C	5.16	118.65	110.19
1	C	294	ASP	CB-CA-C	5.09	120.19	110.17
1	B	629	LEU	CA-C-O	-5.08	116.14	121.88
1	B	637	SER	CA-C-N	-5.07	113.09	120.29
1	B	637	SER	C-N-CA	-5.07	113.09	120.29
1	C	904	TYR	CA-CB-CG	5.07	123.02	113.90
1	A	630	THR	CA-C-N	5.03	124.20	118.97
1	A	630	THR	C-N-CA	5.03	124.20	118.97
3	I	37	GLN	CA-C-N	-5.02	113.36	120.49
3	I	37	GLN	C-N-CA	-5.02	113.36	120.49
3	F	37	GLN	CA-C-N	-5.01	113.38	120.49
3	F	37	GLN	C-N-CA	-5.01	113.38	120.49

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	7884	0	7694	100	0
1	B	6375	0	6269	103	0
1	C	8006	0	7824	112	0
2	E	914	0	876	20	0
2	H	914	0	876	23	0
3	F	764	0	735	10	0
3	I	764	0	735	9	0
4	A	126	0	117	0	0
4	B	98	0	91	0	0
4	C	126	0	117	3	0
All	All	25971	0	25334	359	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 7.

All (359) 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:B:93:ALA:HB1	1:B:189:LEU:HD11	1.62	0.80
1:A:645:THR:HG22	1:A:647:ALA:H	1.48	0.78
1:B:901:GLN:HE21	1:B:905:ARG:HH21	1.32	0.75
1:C:128:ILE:HD13	1:C:170:TYR:HD2	1.51	0.75
1:C:1101:HIS:CD2	4:C:1306:NAG:H5	2.23	0.74
1:B:578:ASP:HB3	1:B:583:GLU:H	1.51	0.73
1:B:577:ARG:HE	1:B:584:ILE:HD11	1.52	0.73
1:B:984:LEU:HD21	1:B:988:GLU:HG3	1.71	0.72
1:B:626:ALA:HA	1:B:633:TRP:CH2	2.23	0.72
1:B:639:GLY:C	1:B:641:ASN:H	1.97	0.72
1:A:1011:GLN:OE1	1:A:1014:ARG:NH2	2.23	0.72
1:B:1104:VAL:HG13	1:B:1115:ILE:HG12	1.72	0.71
3:F:53:ARG:NH2	3:F:61:PHE:O	2.24	0.71
3:I:53:ARG:NH2	3:I:61:PHE:O	2.24	0.70
1:B:626:ALA:HA	1:B:633:TRP:HH2	1.57	0.69
1:C:454:ARG:NH2	1:C:469:SER:O	2.25	0.69
1:A:454:ARG:NH2	1:A:469:SER:O	2.25	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:TYR:OH	1:A:191:GLU:OE2	2.06	0.67
1:A:677:GLN:O	1:A:688:ALA:N	2.26	0.67
1:A:971:GLY:HA3	1:A:995:ARG:HH21	1.58	0.67
1:C:64:TRP:HE1	1:C:264:ASP:HB2	1.60	0.67
1:C:1091:ARG:NH1	1:C:1118:ASP:O	2.28	0.66
1:C:1101:HIS:HD2	4:C:1306:NAG:H5	1.59	0.66
1:A:752:LEU:HD11	1:A:990:GLU:HG2	1.76	0.66
1:B:328:ARG:HD3	1:B:578:ASP:OD1	1.97	0.65
1:B:65:PHE:CE2	1:B:84:LEU:HD11	2.32	0.65
1:A:787:GLN:OE1	1:B:703:ASN:ND2	2.30	0.64
1:C:544:ASN:HD21	1:C:579:PRO:HB3	1.60	0.64
1:B:885:GLY:HA2	1:B:901:GLN:NE2	2.11	0.64
1:C:176:LEU:O	1:C:207:HIS:NE2	2.31	0.64
1:A:811:LYS:NZ	1:A:820:ASP:OD2	2.24	0.64
1:B:736:VAL:HG11	1:B:1004:LEU:HD11	1.80	0.64
1:C:105:ILE:HD11	1:C:241:LEU:HD21	1.78	0.64
2:H:68:THR:HB	2:H:81:GLN:HB3	1.79	0.64
1:A:546:LEU:HD21	1:A:573:THR:HG21	1.79	0.63
1:B:135:PHE:CE1	1:B:159:VAL:HG13	2.34	0.62
1:A:108:THR:O	1:A:236:THR:OG1	2.17	0.62
2:E:68:THR:HB	2:E:81:GLN:HB3	1.79	0.62
1:A:773:GLU:OE2	1:A:1019:ARG:NH1	2.30	0.62
1:C:626:ALA:CB	1:C:628:GLN:HE21	2.11	0.62
1:C:973:ILE:HD12	1:C:983:ARG:HD2	1.82	0.62
1:B:32:PHE:CD2	1:B:218:GLN:HG2	2.35	0.61
1:B:787:GLN:OE1	1:C:703:ASN:ND2	2.32	0.61
1:C:57:PRO:HB3	1:C:273:ARG:HH12	1.66	0.61
1:A:900:MET:HE1	1:B:1077:THR:HG23	1.83	0.61
1:C:457:ARG:NH1	1:C:467:ASP:OD2	2.34	0.61
1:C:57:PRO:HB3	1:C:273:ARG:NH1	2.16	0.60
1:C:175:PHE:O	1:C:177:MET:HG2	2.01	0.60
1:C:90:VAL:HG21	1:C:238:PHE:CE1	2.37	0.60
1:A:457:ARG:NH1	1:A:467:ASP:OD2	2.34	0.60
1:B:135:PHE:HE1	1:B:159:VAL:HG13	1.64	0.60
1:C:1028:LYS:NZ	1:C:1042:PHE:O	2.34	0.59
1:A:1141:LEU:O	1:A:1144:GLU:HG2	2.03	0.59
1:B:117:LEU:HD12	1:B:231:ILE:HG21	1.84	0.59
1:B:122:ASN:HD22	1:B:127:PHE:HE2	1.50	0.59
1:C:32:PHE:CD2	1:C:218:GLN:HG2	2.37	0.59
1:C:916:LEU:C	1:C:916:LEU:HD23	2.27	0.59
2:H:28:ILE:HD13	2:H:31:ARG:HH12	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:VAL:HG23	1:C:174:PRO:HA	1.82	0.59
1:C:91:TYR:OH	1:C:191:GLU:OE1	2.18	0.58
1:B:773:GLU:OE2	1:B:1019:ARG:NH1	2.34	0.58
2:H:17:SER:OG	2:H:82:MET:O	2.18	0.58
1:A:983:ARG:O	1:A:984:LEU:HD23	2.05	0.57
3:I:19:LEU:HD13	3:I:72:LEU:HD23	1.87	0.57
2:E:28:ILE:HD13	2:E:31:ARG:HH12	1.68	0.57
1:A:231:ILE:HG22	1:A:233:ILE:HG23	1.85	0.57
1:B:567:ARG:HG2	1:B:573:THR:HA	1.87	0.57
1:A:825:LYS:NZ	1:A:941:THR:O	2.38	0.57
1:C:735:SER:HB3	1:C:859:THR:OG1	2.04	0.56
1:C:65:PHE:CZ	1:C:84:LEU:HD21	2.40	0.56
1:B:532:ASN:OD1	1:B:533:LEU:N	2.39	0.56
1:B:1006:THR:O	1:B:1010:GLN:HG2	2.05	0.56
1:A:1028:LYS:NZ	1:A:1042:PHE:O	2.39	0.56
1:A:570:VAL:HG23	1:A:572:THR:HG23	1.88	0.56
1:B:115:GLN:HE22	1:B:165:ASN:HD22	1.53	0.55
3:F:19:LEU:HD13	3:F:72:LEU:HD23	1.87	0.55
1:B:973:ILE:HG23	1:B:992:GLN:OE1	2.07	0.55
1:C:196:ASN:HD21	1:C:235:ILE:HD12	1.71	0.55
1:C:971:GLY:C	1:C:995:ARG:HH21	2.15	0.55
1:C:321:GLN:HE21	1:C:630:THR:HB	1.71	0.55
1:A:190:ARG:HB3	1:A:192:PHE:HE1	1.71	0.55
1:B:639:GLY:C	1:B:641:ASN:N	2.64	0.55
1:B:190:ARG:HG2	1:B:207:HIS:ND1	2.22	0.55
1:A:189:LEU:O	1:A:207:HIS:HA	2.07	0.55
1:A:64:TRP:HE1	1:A:264:ASP:CG	2.16	0.54
1:C:347:PHE:CE2	1:C:399:SER:HB2	2.42	0.54
1:A:347:PHE:CE2	1:A:399:SER:HB2	2.42	0.54
1:C:231:ILE:HG22	1:C:233:ILE:HG23	1.88	0.54
1:C:626:ALA:O	1:C:627:ASP:OD1	2.26	0.54
1:A:671:CYS:SG	1:A:697:MET:HE2	2.48	0.54
1:B:133:PHE:HA	1:B:162:SER:O	2.08	0.54
3:F:32:LEU:HD22	3:F:70:PHE:CG	2.43	0.54
1:B:95:THR:HG23	1:B:189:LEU:HD13	1.90	0.54
1:C:327:VAL:HG12	1:C:327:VAL:O	2.07	0.54
3:I:32:LEU:HD22	3:I:70:PHE:CG	2.43	0.54
1:C:984:LEU:HB3	1:C:988:GLU:HG3	1.91	0.53
1:B:971:GLY:C	1:B:995:ARG:HE	2.17	0.53
1:A:144:TYR:HD2	1:A:145:HIS:ND1	2.07	0.53
1:A:462:LYS:N	1:A:465:GLU:OE2	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:PHE:HA	1:A:630:THR:OG1	2.09	0.52
1:A:443:SER:HB3	1:A:499:PRO:HD3	1.91	0.52
1:A:751:ASN:HA	1:A:754:LEU:HD12	1.90	0.52
1:B:92:PHE:HB3	1:B:192:PHE:HB2	1.91	0.52
1:C:1104:VAL:HG13	1:C:1115:ILE:HG12	1.91	0.52
1:A:455:SER:HG	1:A:493:GLN:CD	2.17	0.52
1:C:462:LYS:N	1:C:465:GLU:OE2	2.42	0.52
2:E:63:VAL:HB	2:E:67:PHE:CD1	2.44	0.52
1:A:789:TYR:HA	1:B:703:ASN:O	2.09	0.52
1:B:752:LEU:HD11	1:B:990:GLU:HG2	1.91	0.52
1:C:715:PRO:HA	1:C:1072:GLU:HA	1.92	0.52
2:H:63:VAL:HB	2:H:67:PHE:CD1	2.44	0.52
1:C:30:ASN:HB3	1:C:32:PHE:CE1	2.44	0.52
1:C:44:ARG:O	1:C:283:GLY:HA2	2.10	0.52
1:B:134:GLN:HB3	1:B:161:SER:HB2	1.92	0.52
3:F:60:ARG:HH21	3:F:81:ASP:CG	2.18	0.52
1:B:106:PHE:HB3	1:B:235:ILE:HD13	1.91	0.52
1:B:662:CYS:HB2	1:B:697:MET:HE3	1.91	0.52
1:B:588:THR:CG2	1:B:589:PRO:HD2	2.40	0.51
1:B:752:LEU:O	1:B:755:GLN:HG2	2.11	0.51
1:A:675:GLN:HG3	1:A:693:ILE:HD11	1.91	0.51
1:B:65:PHE:CZ	1:B:84:LEU:HD11	2.45	0.51
1:C:443:SER:HB3	1:C:499:PRO:HD3	1.91	0.51
2:E:80:LEU:HD23	2:E:82:MET:HE2	1.93	0.51
2:H:80:LEU:HD23	2:H:82:MET:HE2	1.93	0.51
3:I:60:ARG:HH21	3:I:81:ASP:CG	2.18	0.51
2:E:40:ALA:HB3	2:E:43:LYS:HD2	1.93	0.51
2:H:40:ALA:HB3	2:H:43:LYS:HD2	1.93	0.51
1:B:945:LEU:HD12	1:B:948:LEU:HD12	1.93	0.51
1:A:42:VAL:HG11	1:B:567:ARG:HE	1.76	0.51
1:C:95:THR:HG22	1:C:189:LEU:HD13	1.92	0.51
1:C:396:TYR:HB2	1:C:514:SER:HB3	1.92	0.51
1:A:973:ILE:HD12	1:A:983:ARG:NH2	2.26	0.50
1:B:193:VAL:HG13	1:B:270:LEU:HD11	1.92	0.50
1:B:588:THR:HG23	1:B:589:PRO:HD2	1.92	0.50
1:A:396:TYR:HB2	1:A:514:SER:HB3	1.92	0.50
1:C:130:VAL:HG21	1:C:231:ILE:HD12	1.94	0.50
1:B:603:ASN:OD1	1:B:604:THR:N	2.45	0.50
1:C:87:ASN:ND2	1:C:269:TYR:CD2	2.78	0.50
1:A:1090:PRO:HD3	1:A:1095:PHE:CE2	2.47	0.50
1:C:455:SER:HG	1:C:493:GLN:CD	2.17	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:632:THR:C	1:B:634:ARG:N	2.66	0.50
1:A:808:ASP:HB3	1:A:811:LYS:HE2	1.93	0.49
1:B:560:LEU:O	1:B:562:PHE:N	2.45	0.49
1:B:884:SER:OG	1:B:893:ALA:HB1	2.11	0.49
1:C:645:THR:HG23	1:C:670:ILE:HG13	1.94	0.49
1:A:773:GLU:OE2	1:A:1019:ARG:HD3	2.12	0.49
1:C:128:ILE:HD13	1:C:170:TYR:CD2	2.40	0.49
1:C:196:ASN:ND2	1:C:235:ILE:HB	2.27	0.49
1:B:971:GLY:HA3	1:B:995:ARG:HH21	1.78	0.49
1:A:187:LYS:O	1:A:209:PRO:HA	2.13	0.49
1:C:87:ASN:ND2	1:C:269:TYR:CE2	2.81	0.49
1:B:273:ARG:NH1	1:B:290:ASP:OD1	2.46	0.49
1:B:533:LEU:HD21	1:B:585:LEU:HD11	1.95	0.49
1:C:130:VAL:HG21	1:C:231:ILE:HG23	1.95	0.49
1:B:654:GLU:OE2	1:B:677:GLN:HG3	2.13	0.49
1:C:748:GLU:O	1:C:752:LEU:HD12	2.12	0.48
2:E:17:SER:OG	2:E:82:MET:O	2.18	0.48
1:B:758:SER:O	1:B:762:GLN:HG3	2.12	0.48
1:B:916:LEU:HD12	1:B:923:ILE:HD13	1.96	0.48
1:C:393:THR:HG21	1:C:518:LEU:HB2	1.95	0.48
1:A:633:TRP:O	1:A:634:ARG:C	2.54	0.48
1:B:628:GLN:HB3	1:B:633:TRP:HE1	1.78	0.48
1:B:268:GLY:C	1:B:269:TYR:HD1	2.22	0.48
3:I:11:LEU:HD12	3:I:15:GLU:HG3	1.96	0.48
1:C:190:ARG:HG3	1:C:207:HIS:NE2	2.29	0.48
1:C:328:ARG:NH2	1:C:578:ASP:OD2	2.47	0.48
1:A:393:THR:HG21	1:A:518:LEU:HB2	1.95	0.48
1:A:204:TYR:CE2	1:A:225:PRO:HG3	2.49	0.47
1:A:636:TYR:C	1:A:641:ASN:HD21	2.22	0.47
1:A:1121:PHE:HE2	1:C:914:ASN:HD22	1.61	0.47
1:B:135:PHE:HE1	1:B:159:VAL:CG1	2.27	0.47
1:C:532:ASN:OD1	1:C:533:LEU:N	2.45	0.47
1:A:42:VAL:HG21	1:B:567:ARG:HH21	1.79	0.47
1:B:105:ILE:HD11	1:B:241:LEU:HD21	1.96	0.47
2:E:29:VAL:HG22	2:E:34:MET:HE2	1.96	0.47
1:B:135:PHE:HB3	1:B:139:PRO:HG3	1.95	0.47
3:F:78:GLU:HB2	3:F:81:ASP:CG	2.39	0.47
1:C:1090:PRO:HD3	1:C:1095:PHE:CE2	2.50	0.47
3:F:11:LEU:HD12	3:F:15:GLU:HG3	1.96	0.47
2:H:29:VAL:HG22	2:H:34:MET:HE2	1.95	0.47
3:I:78:GLU:HB2	3:I:81:ASP:CG	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:458:LYS:HD3	2:H:31:ARG:HD3	1.96	0.47
1:A:295:PRO:HB2	1:A:608:VAL:HG21	1.96	0.47
1:A:894:LEU:HB3	1:B:713:ALA:HB3	1.97	0.47
1:C:196:ASN:OD1	1:C:201:PHE:HD1	1.98	0.47
1:C:605:SER:OG	1:C:607:GLN:HG3	2.15	0.47
1:A:170:TYR:CE2	1:A:172:SER:HB2	2.49	0.47
1:A:628:GLN:O	1:A:629:LEU:HD23	2.15	0.47
1:C:328:ARG:NH1	1:C:533:LEU:HB2	2.30	0.47
1:A:703:ASN:ND2	1:C:787:GLN:OE1	2.46	0.47
1:C:458:LYS:HD3	2:E:31:ARG:HD3	1.96	0.47
1:C:786:LYS:HG3	1:C:787:GLN:HG3	1.96	0.47
1:A:401:VAL:HG22	1:A:509:ARG:HG2	1.97	0.46
1:A:1006:THR:O	1:A:1010:GLN:HG2	2.15	0.46
1:B:533:LEU:HD23	1:B:535:LYS:NZ	2.29	0.46
1:C:295:PRO:HG3	1:C:633:TRP:CE3	2.49	0.46
1:A:560:LEU:O	1:A:562:PHE:N	2.48	0.46
1:A:985:ASP:HB2	1:A:988:GLU:OE1	2.14	0.46
1:B:30:ASN:HB3	1:B:32:PHE:CE1	2.51	0.46
1:B:44:ARG:O	1:B:283:GLY:HA2	2.15	0.46
1:B:93:ALA:C	1:B:94:SER:HG	2.24	0.46
1:C:196:ASN:HD21	1:C:235:ILE:HB	1.81	0.46
1:C:401:VAL:HG22	1:C:509:ARG:HG2	1.97	0.46
1:A:65:PHE:HE2	1:A:84:LEU:HD11	1.79	0.46
1:A:327:VAL:HG23	1:A:542:ASN:HB3	1.98	0.46
1:C:662:CYS:HB2	1:C:697:MET:HE3	1.96	0.46
1:C:1006:THR:O	1:C:1010:GLN:OE1	2.34	0.46
1:A:126:VAL:O	1:A:171:VAL:HA	2.15	0.46
1:A:216:PHE:HB2	1:A:266:TYR:CZ	2.51	0.46
1:C:294:ASP:OD1	1:C:297:SER:OG	2.30	0.46
1:A:53:ASP:HB3	1:A:55:PHE:CE2	2.50	0.46
1:B:280:ASN:OD1	1:B:284:THR:OG1	2.24	0.46
1:B:866:THR:O	1:B:870:ILE:HG12	2.16	0.46
3:F:12:SER:O	3:F:15:GLU:HG2	2.16	0.46
3:I:12:SER:O	3:I:15:GLU:HG2	2.16	0.46
1:A:497:PHE:CE2	1:A:507:PRO:HB3	2.51	0.46
1:B:634:ARG:O	1:B:635:VAL:C	2.56	0.46
1:C:36:VAL:HG21	1:C:220:PHE:HE1	1.81	0.46
1:C:458:LYS:CD	2:E:31:ARG:HD3	2.46	0.46
1:C:326:ILE:HD11	1:C:534:VAL:HB	1.98	0.46
1:C:497:PHE:CE2	1:C:507:PRO:HB3	2.51	0.46
1:C:914:ASN:OD1	1:C:915:VAL:HG23	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1101:HIS:HD2	4:C:1306:NAG:C5	2.28	0.46
1:A:420:ASP:OD1	1:A:460:LYS:HE3	2.17	0.45
1:B:101:ILE:HG12	1:B:242:LEU:HD12	1.97	0.45
1:B:903:ALA:HB1	1:B:913:GLN:HB2	1.97	0.45
1:C:193:VAL:HG13	1:C:270:LEU:HD11	1.98	0.45
1:A:916:LEU:HD12	1:A:923:ILE:HD13	1.97	0.45
1:C:187:LYS:O	1:C:210:ILE:N	2.32	0.45
1:A:458:LYS:CD	2:H:31:ARG:HD3	2.46	0.45
1:B:629:LEU:HD12	1:B:629:LEU:HA	1.75	0.45
1:C:165:ASN:O	1:C:167:THR:N	2.50	0.45
1:C:420:ASP:OD1	1:C:460:LYS:HE3	2.17	0.45
1:B:620:VAL:O	1:B:624:ILE:HG12	2.16	0.45
1:B:639:GLY:O	1:B:641:ASN:N	2.49	0.45
1:C:165:ASN:O	1:C:167:THR:HG23	2.17	0.45
1:B:541:PHE:O	1:B:547:THR:HA	2.17	0.44
1:B:902:MET:HE1	1:B:1049:LEU:HD13	2.00	0.44
1:A:102:ARG:HB3	1:A:241:LEU:HB2	1.99	0.44
1:A:628:GLN:O	1:A:628:GLN:HG3	2.18	0.44
1:C:377:PHE:CD1	1:C:434:ILE:HG12	2.53	0.44
1:C:645:THR:HG22	1:C:647:ALA:H	1.82	0.44
1:A:64:TRP:C	1:A:64:TRP:CD1	2.95	0.44
1:B:328:ARG:CZ	1:B:533:LEU:HB2	2.48	0.44
1:B:1019:ARG:NH2	1:C:1017:GLU:HG2	2.33	0.44
1:A:33:THR:HG22	1:A:220:PHE:HD1	1.82	0.44
1:A:858:LEU:HD13	1:A:959:LEU:HD22	1.98	0.44
1:C:188:ASN:OD1	1:C:190:ARG:NH1	2.51	0.44
1:C:551:VAL:HB	1:C:588:THR:CG2	2.47	0.44
1:A:491:PRO:O	1:A:492:LEU:HD23	2.17	0.44
1:B:328:ARG:NH1	1:B:533:LEU:HB2	2.33	0.44
1:C:491:PRO:O	1:C:492:LEU:HD23	2.17	0.44
3:I:36:GLN:HB2	3:I:46:LEU:HD11	2.00	0.44
1:C:916:LEU:HD23	1:C:916:LEU:O	2.18	0.44
1:B:86:PHE:O	1:B:87:ASN:C	2.61	0.44
1:A:105:ILE:O	1:A:238:PHE:HA	2.18	0.44
1:A:713:ALA:HB3	1:C:894:LEU:HB3	2.00	0.44
1:A:122:ASN:OD1	1:A:123:ALA:N	2.47	0.43
1:C:328:ARG:HH12	1:C:533:LEU:HB2	1.83	0.43
3:F:36:GLN:HB2	3:F:46:LEU:HD11	2.00	0.43
1:B:126:VAL:O	1:B:171:VAL:HA	2.17	0.43
2:H:78:MET:HE3	2:H:78:MET:HB2	1.89	0.43
1:B:122:ASN:OD1	1:B:123:ALA:N	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1090:PRO:HD3	1:B:1095:PHE:CE2	2.53	0.43
1:C:638:THR:O	1:C:639:GLY:C	2.60	0.43
2:E:28:ILE:O	2:E:32:ASN:ND2	2.51	0.43
1:A:377:PHE:CD1	1:A:434:ILE:HG12	2.53	0.43
1:A:904:TYR:CZ	1:B:1107:ARG:HD3	2.53	0.43
1:B:231:ILE:HG22	1:B:233:ILE:HG23	1.99	0.43
2:H:28:ILE:O	2:H:32:ASN:ND2	2.51	0.43
1:B:645:THR:HG23	1:B:670:ILE:HG13	2.00	0.43
1:A:48:LEU:HD22	1:A:306:PHE:CE1	2.54	0.43
1:A:825:LYS:HE3	1:A:945:LEU:HD23	2.01	0.43
1:C:115:GLN:HG2	1:C:132:GLU:HG3	2.00	0.43
1:B:901:GLN:NE2	1:B:905:ARG:HH21	2.06	0.43
1:A:196:ASN:OD1	1:A:235:ILE:HD12	2.19	0.42
1:C:733:LYS:HG2	1:C:771:ALA:HA	2.01	0.42
1:C:908:GLY:O	1:C:1038:LYS:HE3	2.18	0.42
3:F:12:SER:N	3:F:15:GLU:OE2	2.52	0.42
2:H:28:ILE:HA	2:H:76:ASN:HD21	1.84	0.42
3:F:28:PRO:HB2	3:F:30:THR:HG22	2.01	0.42
2:H:36:TRP:O	2:H:48:VAL:HB	2.19	0.42
2:H:43:LYS:HG2	2:H:44:GLY:N	2.34	0.42
1:A:866:THR:O	1:A:870:ILE:HG12	2.19	0.42
1:A:190:ARG:HB3	1:A:192:PHE:CE1	2.52	0.42
1:A:326:ILE:N	1:A:540:ASN:O	2.43	0.42
1:A:758:SER:O	1:A:762:GLN:HG3	2.19	0.42
1:C:630:THR:O	1:C:631:PRO:C	2.63	0.42
2:E:36:TRP:O	2:E:48:VAL:HB	2.19	0.42
2:H:63:VAL:HB	2:H:67:PHE:CE1	2.55	0.42
3:I:28:PRO:HB2	3:I:30:THR:HG22	2.01	0.42
1:C:31:SER:OG	1:C:60:SER:OG	2.34	0.42
1:C:902:MET:HB3	1:C:916:LEU:CD1	2.50	0.42
1:C:64:TRP:NE1	1:C:264:ASP:HB2	2.32	0.42
1:C:726:ILE:HG12	1:C:1061:VAL:HG22	2.01	0.42
1:C:854:LYS:HD3	1:C:854:LYS:HA	1.91	0.42
2:E:36:TRP:HE1	2:E:78:MET:CG	2.33	0.42
2:E:43:LYS:HG2	2:E:44:GLY:N	2.34	0.42
1:B:32:PHE:HD2	1:B:218:GLN:HG2	1.83	0.42
1:B:540:ASN:OD1	1:B:549:THR:HG22	2.20	0.42
1:B:553:THR:HG22	1:B:554:LYS:N	2.34	0.42
1:B:889:GLY:O	1:C:1045:LYS:HE3	2.19	0.42
1:A:287:ASP:OD1	1:A:288:ALA:N	2.46	0.42
1:A:984:LEU:HB3	1:A:988:GLU:HG2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:66:ARG:NH1	2:H:89:ASP:OD2	2.40	0.42
1:A:945:LEU:HD12	1:A:948:LEU:HD12	2.02	0.41
1:C:328:ARG:HD2	1:C:328:ARG:HA	1.87	0.41
1:C:357:ARG:HG3	1:C:396:TYR:CE1	2.55	0.41
2:E:28:ILE:HA	2:E:76:ASN:HD21	1.84	0.41
2:E:63:VAL:HB	2:E:67:PHE:CE1	2.55	0.41
1:A:84:LEU:HB2	1:A:238:PHE:O	2.21	0.41
1:B:610:VAL:CG2	1:B:635:VAL:HG11	2.50	0.41
1:C:294:ASP:OD1	1:C:297:SER:CB	2.68	0.41
1:C:626:ALA:HB1	1:C:628:GLN:HE21	1.84	0.41
1:C:629:LEU:HD12	1:C:629:LEU:HA	1.88	0.41
1:A:45:SER:O	1:A:47:VAL:HG13	2.19	0.41
1:B:191:GLU:OE1	1:B:223:LEU:HD11	2.21	0.41
1:C:569:ILE:HG13	1:C:570:VAL:HG13	2.02	0.41
2:H:77:THR:HG22	2:H:78:MET:H	1.85	0.41
1:B:610:VAL:HG23	1:B:635:VAL:HG11	2.01	0.41
1:C:36:VAL:HG21	1:C:220:PHE:CE1	2.55	0.41
1:C:971:GLY:HA3	1:C:995:ARG:NH2	2.34	0.41
1:A:616:ASN:HB3	1:A:618:THR:HG22	2.02	0.41
2:E:77:THR:HG22	2:E:78:MET:H	1.85	0.41
1:A:534:VAL:HG11	1:A:539:VAL:HG11	2.02	0.41
1:B:908:GLY:O	1:B:1038:LYS:HE3	2.21	0.41
1:B:1089:PHE:CE1	1:B:1123:SER:HB3	2.56	0.41
1:C:462:LYS:HB2	1:C:462:LYS:HE2	1.73	0.41
1:A:762:GLN:NE2	1:B:961:THR:HG21	2.36	0.41
1:B:160:TYR:O	1:B:160:TYR:CD1	2.72	0.41
1:B:321:GLN:HE21	1:B:629:LEU:HB2	1.84	0.41
1:C:106:PHE:HB3	1:C:235:ILE:HD13	2.02	0.41
2:H:26:GLU:HG3	2:H:26:GLU:O	2.21	0.41
1:A:133:PHE:HD1	1:A:160:TYR:HA	1.86	0.41
1:A:1092:GLU:O	1:A:1092:GLU:HG2	2.21	0.41
1:B:564:GLN:HE22	1:B:577:ARG:HD2	1.86	0.41
1:C:551:VAL:HB	1:C:588:THR:HG22	2.03	0.41
1:C:1139:ASP:HB3	1:C:1142:GLN:HE21	1.86	0.41
1:A:357:ARG:HG3	1:A:396:TYR:CE1	2.55	0.40
1:A:537:LYS:HE2	1:A:537:LYS:HB2	1.90	0.40
1:A:666:ILE:HD11	1:A:672:ALA:HB2	2.03	0.40
1:B:34:ARG:NH2	1:B:191:GLU:OE2	2.51	0.40
1:B:576:VAL:O	1:B:584:ILE:HA	2.21	0.40
2:E:26:GLU:O	2:E:26:GLU:HG3	2.21	0.40
2:H:36:TRP:HE1	2:H:78:MET:CG	2.33	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:LEU:HD23	1:A:84:LEU:HA	1.87	0.40
1:A:726:ILE:HG12	1:A:1061:VAL:HG22	2.03	0.40
1:C:379:CYS:HA	1:C:432:CYS:HA	2.03	0.40
2:E:5:VAL:O	2:E:5:VAL:HG23	2.22	0.40
2:H:36:TRP:HE1	2:H:78:MET:HG2	1.86	0.40
1:B:1048:HIS:HA	1:B:1066:THR:HG22	2.03	0.40
1:C:200:TYR:HB3	1:C:228:ASP:OD1	2.22	0.40
2:H:1:GLU:HG2	2:H:2:VAL:HG23	2.04	0.40
1:A:129:LYS:HG2	1:A:169:GLU:OE1	2.21	0.40
1:A:776:LYS:HE2	1:A:780:GLU:OE1	2.20	0.40
2:E:1:GLU:N	2:E:3:GLN:HE21	2.20	0.40
2:E:36:TRP:HE1	2:E:78:MET:HG2	1.86	0.40
2:H:5:VAL:O	2:H:5:VAL:HG23	2.22	0.40
2:H:1:GLU:N	2:H:3:GLN:HE21	2.20	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	978/1200 (82%)	941 (96%)	37 (4%)	0	100	100
1	B	800/1200 (67%)	767 (96%)	31 (4%)	2 (0%)	36	52
1	C	1003/1200 (84%)	962 (96%)	40 (4%)	1 (0%)	48	66
2	E	115/117 (98%)	108 (94%)	7 (6%)	0	100	100
2	H	115/117 (98%)	108 (94%)	7 (6%)	0	100	100
3	F	98/106 (92%)	92 (94%)	6 (6%)	0	100	100
3	I	98/106 (92%)	92 (94%)	6 (6%)	0	100	100
All	All	3207/4046 (79%)	3070 (96%)	134 (4%)	3 (0%)	49	66

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	640	SER
1	B	625	HIS
1	C	632	THR

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	198/1050 (19%)	198 (100%)	0	100	100
1	B	615/1050 (59%)	615 (100%)	0	100	100
1	C	894/1050 (85%)	894 (100%)	0	100	100
2	E	98/98 (100%)	98 (100%)	0	100	100
2	H	98/98 (100%)	98 (100%)	0	100	100
3	F	82/88 (93%)	82 (100%)	0	100	100
3	I	82/88 (93%)	82 (100%)	0	100	100
All	All	2067/3522 (59%)	2067 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	960	ASN
1	A	1074	ASN
1	A	1106	GLN
1	B	49	HIS
1	B	87	ASN
1	B	115	GLN
1	B	607	GLN
1	B	625	HIS
1	B	677	GLN
1	B	751	ASN

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Mol	Chain	Res	Type
1	B	901	GLN
1	B	907	ASN
1	B	1010	GLN
1	C	49	HIS
1	C	61	ASN
1	C	87	ASN
1	C	196	ASN
1	C	321	GLN
1	C	542	ASN
1	C	603	ASN
1	C	607	GLN
1	C	625	HIS
1	C	628	GLN
1	C	690	GLN
1	C	751	ASN
1	C	901	GLN
1	C	992	GLN
1	C	1101	HIS
1	C	1106	GLN
2	E	3	GLN
2	E	39	GLN
3	F	37	GLN
3	F	75	ASN
2	H	3	GLN
2	H	39	GLN
3	I	37	GLN
3	I	75	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

25 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
4	NAG	A	1302	1	14,14,15	0.40	0	17,19,21	1.66	1 (5%)
4	NAG	A	1308	1	14,14,15	0.23	0	17,19,21	0.88	1 (5%)
4	NAG	C	1303	1	14,14,15	0.55	0	17,19,21	1.59	4 (23%)
4	NAG	C	1305	1	14,14,15	0.58	0	17,19,21	1.69	4 (23%)
4	NAG	A	1304	1	14,14,15	0.67	0	17,19,21	1.53	3 (17%)
4	NAG	B	1303	1	14,14,15	0.44	0	17,19,21	1.24	1 (5%)
4	NAG	B	1302	1	14,14,15	0.43	0	17,19,21	0.63	0
4	NAG	C	1307	1	14,14,15	0.43	0	17,19,21	1.05	1 (5%)
4	NAG	C	1301	1	14,14,15	0.41	0	17,19,21	0.70	0
4	NAG	A	1303	1	14,14,15	0.51	0	17,19,21	1.22	1 (5%)
4	NAG	A	1305	1	14,14,15	0.38	0	17,19,21	0.85	0
4	NAG	B	1305	1	14,14,15	0.67	0	17,19,21	1.10	1 (5%)
4	NAG	A	1309	1	14,14,15	0.40	0	17,19,21	1.10	2 (11%)
4	NAG	B	1307	1	14,14,15	0.52	0	17,19,21	1.00	1 (5%)
4	NAG	C	1306	1	14,14,15	0.40	0	17,19,21	1.33	4 (23%)
4	NAG	A	1301	1	14,14,15	0.40	0	17,19,21	1.27	3 (17%)
4	NAG	B	1301	1	14,14,15	0.25	0	17,19,21	1.31	2 (11%)
4	NAG	A	1306	1	14,14,15	0.32	0	17,19,21	1.03	2 (11%)
4	NAG	A	1307	1	14,14,15	0.34	0	17,19,21	1.49	2 (11%)
4	NAG	B	1304	1	14,14,15	0.71	0	17,19,21	0.94	0
4	NAG	C	1308	1	14,14,15	0.33	0	17,19,21	1.48	2 (11%)
4	NAG	C	1309	1	14,14,15	0.22	0	17,19,21	0.88	1 (5%)
4	NAG	B	1306	1	14,14,15	0.37	0	17,19,21	1.70	5 (29%)
4	NAG	C	1302	1	14,14,15	0.39	0	17,19,21	0.77	1 (5%)
4	NAG	C	1304	1	14,14,15	0.64	0	17,19,21	1.51	2 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1302	1	-	3/6/23/26	0/1/1/1
4	NAG	A	1308	1	-	4/6/23/26	0/1/1/1
4	NAG	C	1303	1	-	4/6/23/26	0/1/1/1
4	NAG	C	1305	1	-	4/6/23/26	0/1/1/1
4	NAG	A	1304	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1303	1	-	5/6/23/26	0/1/1/1
4	NAG	B	1302	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1307	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1301	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1303	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1305	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1305	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1309	1	-	3/6/23/26	0/1/1/1
4	NAG	B	1307	1	-	4/6/23/26	0/1/1/1
4	NAG	C	1306	1	-	3/6/23/26	0/1/1/1
4	NAG	A	1301	1	-	3/6/23/26	0/1/1/1
4	NAG	B	1301	1	-	3/6/23/26	0/1/1/1
4	NAG	A	1306	1	-	5/6/23/26	0/1/1/1
4	NAG	A	1307	1	-	3/6/23/26	0/1/1/1
4	NAG	B	1304	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1308	1	-	3/6/23/26	0/1/1/1
4	NAG	C	1309	1	-	4/6/23/26	0/1/1/1
4	NAG	B	1306	1	-	4/6/23/26	0/1/1/1
4	NAG	C	1302	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1304	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1302	NAG	C2-N2-C7	5.42	130.16	122.90
4	C	1303	NAG	C1-O5-C5	4.46	118.17	112.19
4	B	1303	NAG	C1-O5-C5	4.04	117.59	112.19
4	C	1304	NAG	O5-C1-C2	-4.03	105.05	111.29
4	A	1307	NAG	C2-N2-C7	3.89	128.11	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1308	NAG	C2-N2-C7	3.87	128.09	122.90
4	C	1308	NAG	C1-C2-N2	3.56	116.05	110.43
4	B	1306	NAG	O5-C1-C2	-3.56	105.79	111.29
4	A	1307	NAG	C1-C2-N2	3.55	116.03	110.43
4	B	1307	NAG	C1-O5-C5	3.46	116.83	112.19
4	C	1305	NAG	C2-N2-C7	3.39	127.44	122.90
4	B	1306	NAG	C2-N2-C7	3.33	127.36	122.90
4	A	1304	NAG	O5-C1-C2	-3.25	106.26	111.29
4	C	1305	NAG	O5-C1-C2	-3.21	106.33	111.29
4	A	1304	NAG	C1-O5-C5	3.05	116.28	112.19
4	A	1304	NAG	C2-N2-C7	-3.02	118.85	122.90
4	B	1301	NAG	C1-O5-C5	2.98	116.18	112.19
4	C	1305	NAG	C1-O5-C5	2.87	116.04	112.19
4	A	1301	NAG	C4-C3-C2	-2.87	106.81	111.02
4	C	1306	NAG	O5-C1-C2	-2.79	106.98	111.29
4	A	1301	NAG	C2-N2-C7	-2.74	119.23	122.90
4	B	1301	NAG	C2-N2-C7	2.68	126.50	122.90
4	C	1306	NAG	C1-C2-N2	2.67	114.64	110.43
4	A	1309	NAG	C2-N2-C7	-2.61	119.40	122.90
4	B	1306	NAG	C1-C2-N2	2.58	114.50	110.43
4	A	1303	NAG	O5-C1-C2	-2.53	107.38	111.29
4	A	1306	NAG	C1-O5-C5	2.45	115.47	112.19
4	B	1305	NAG	O5-C1-C2	-2.41	107.56	111.29
4	C	1302	NAG	C4-C3-C2	-2.34	107.58	111.02
4	A	1306	NAG	C2-N2-C7	2.34	126.03	122.90
4	C	1303	NAG	C2-N2-C7	2.32	126.01	122.90
4	C	1306	NAG	C4-C3-C2	-2.31	107.63	111.02
4	B	1306	NAG	O5-C5-C6	2.30	112.13	107.66
4	C	1307	NAG	O5-C1-C2	-2.26	107.79	111.29
4	A	1309	NAG	C4-C3-C2	-2.20	107.79	111.02
4	A	1301	NAG	C1-O5-C5	-2.12	109.34	112.19
4	C	1305	NAG	C1-C2-N2	2.11	113.76	110.43
4	B	1306	NAG	C3-C4-C5	-2.08	106.46	110.23
4	C	1304	NAG	C1-O5-C5	2.08	114.97	112.19
4	C	1306	NAG	C2-N2-C7	2.06	125.66	122.90
4	C	1303	NAG	C1-C2-N2	-2.04	107.22	110.43
4	C	1303	NAG	O5-C5-C6	-2.04	103.70	107.66
4	C	1309	NAG	C2-N2-C7	2.03	125.62	122.90
4	A	1308	NAG	C2-N2-C7	2.02	125.60	122.90

There are no chirality outliers.

All (65) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1301	NAG	C8-C7-N2-C2
4	A	1301	NAG	O7-C7-N2-C2
4	A	1302	NAG	C3-C2-N2-C7
4	A	1306	NAG	C1-C2-N2-C7
4	A	1307	NAG	C1-C2-N2-C7
4	A	1307	NAG	C8-C7-N2-C2
4	A	1307	NAG	O7-C7-N2-C2
4	A	1308	NAG	C1-C2-N2-C7
4	B	1303	NAG	C1-C2-N2-C7
4	B	1303	NAG	C8-C7-N2-C2
4	B	1303	NAG	O7-C7-N2-C2
4	B	1306	NAG	C1-C2-N2-C7
4	B	1307	NAG	C8-C7-N2-C2
4	B	1307	NAG	O7-C7-N2-C2
4	C	1302	NAG	C8-C7-N2-C2
4	C	1302	NAG	O7-C7-N2-C2
4	C	1306	NAG	C1-C2-N2-C7
4	C	1308	NAG	C1-C2-N2-C7
4	C	1308	NAG	C8-C7-N2-C2
4	C	1308	NAG	O7-C7-N2-C2
4	C	1309	NAG	C1-C2-N2-C7
4	A	1302	NAG	C8-C7-N2-C2
4	A	1302	NAG	O7-C7-N2-C2
4	B	1302	NAG	C8-C7-N2-C2
4	B	1302	NAG	O7-C7-N2-C2
4	B	1304	NAG	C8-C7-N2-C2
4	B	1304	NAG	O7-C7-N2-C2
4	A	1306	NAG	O7-C7-N2-C2
4	A	1308	NAG	O7-C7-N2-C2
4	C	1309	NAG	O7-C7-N2-C2
4	A	1306	NAG	O5-C5-C6-O6
4	A	1306	NAG	C8-C7-N2-C2
4	A	1308	NAG	C8-C7-N2-C2
4	C	1309	NAG	C8-C7-N2-C2
4	C	1303	NAG	O5-C5-C6-O6
4	A	1306	NAG	C4-C5-C6-O6
4	B	1306	NAG	C8-C7-N2-C2
4	C	1303	NAG	C4-C5-C6-O6
4	B	1306	NAG	O7-C7-N2-C2
4	B	1303	NAG	O5-C5-C6-O6
4	B	1303	NAG	C4-C5-C6-O6
4	C	1305	NAG	O5-C5-C6-O6
4	C	1305	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	A	1303	NAG	C8-C7-N2-C2
4	B	1305	NAG	O5-C5-C6-O6
4	B	1307	NAG	O5-C5-C6-O6
4	A	1308	NAG	O5-C5-C6-O6
4	C	1309	NAG	O5-C5-C6-O6
4	B	1307	NAG	C4-C5-C6-O6
4	A	1309	NAG	C8-C7-N2-C2
4	B	1301	NAG	O5-C5-C6-O6
4	A	1301	NAG	O5-C5-C6-O6
4	A	1303	NAG	O7-C7-N2-C2
4	B	1305	NAG	C4-C5-C6-O6
4	A	1309	NAG	O7-C7-N2-C2
4	B	1301	NAG	C1-C2-N2-C7
4	C	1303	NAG	C1-C2-N2-C7
4	C	1305	NAG	C1-C2-N2-C7
4	C	1306	NAG	C8-C7-N2-C2
4	B	1301	NAG	C3-C2-N2-C7
4	C	1303	NAG	C3-C2-N2-C7
4	C	1306	NAG	O7-C7-N2-C2
4	A	1309	NAG	C4-C5-C6-O6
4	B	1306	NAG	C4-C5-C6-O6
4	C	1305	NAG	C3-C2-N2-C7

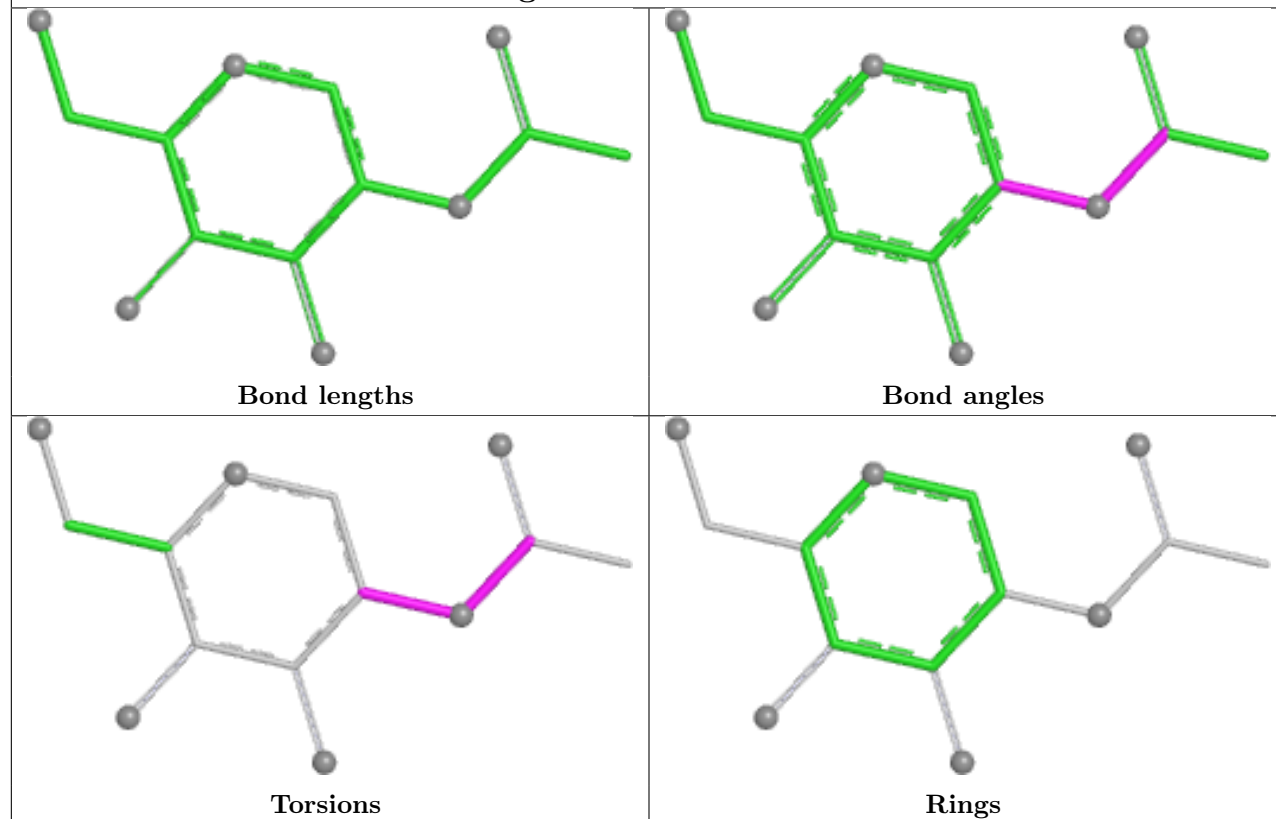
There are no ring outliers.

1 monomer is involved in 3 short contacts:

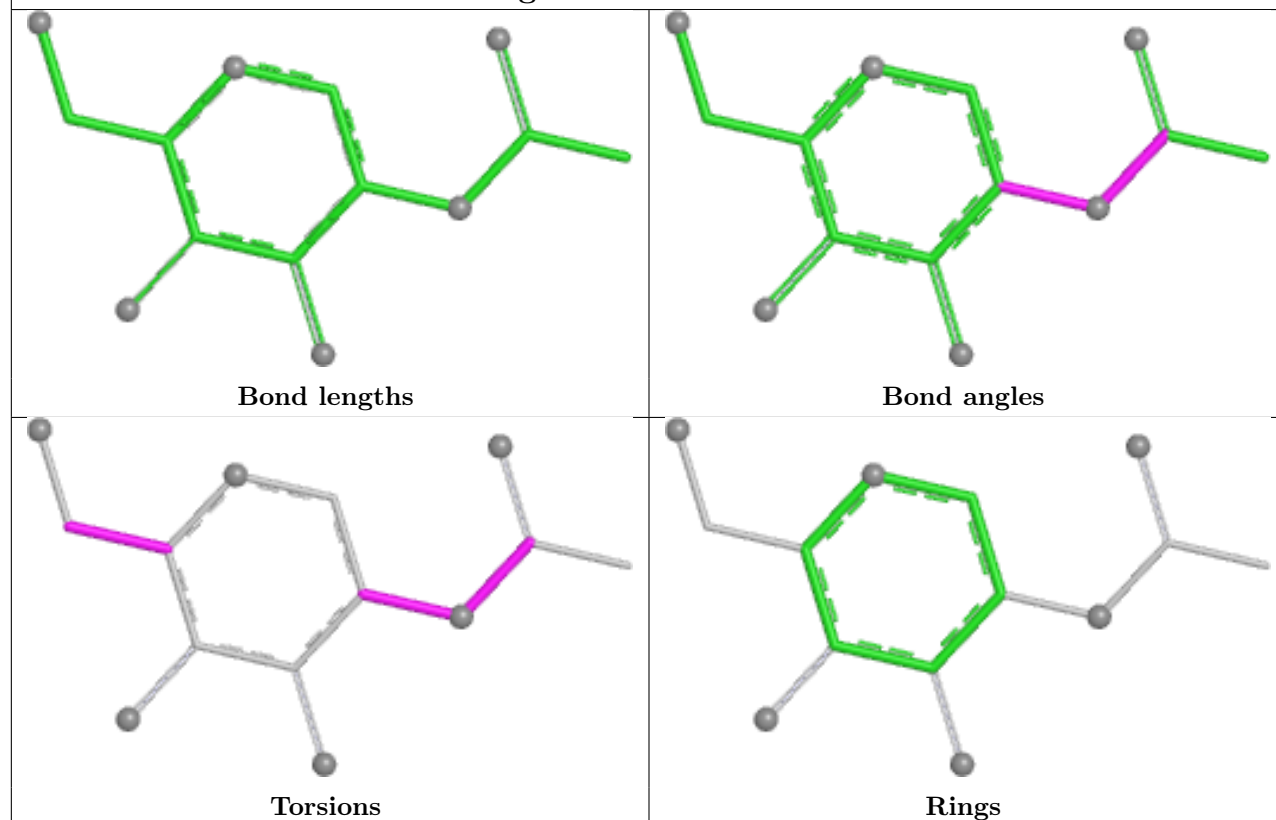
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1306	NAG	3	0

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.

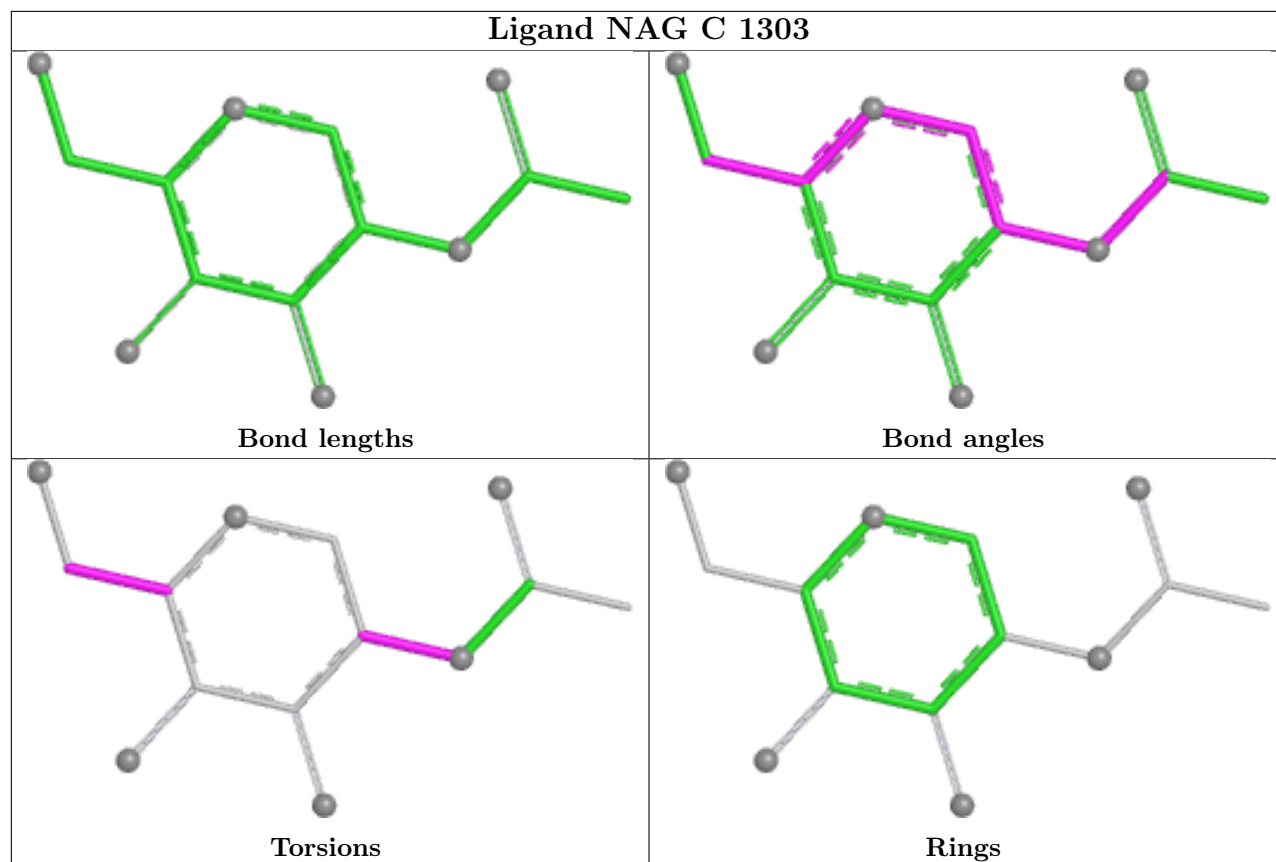
Ligand NAG A 1302



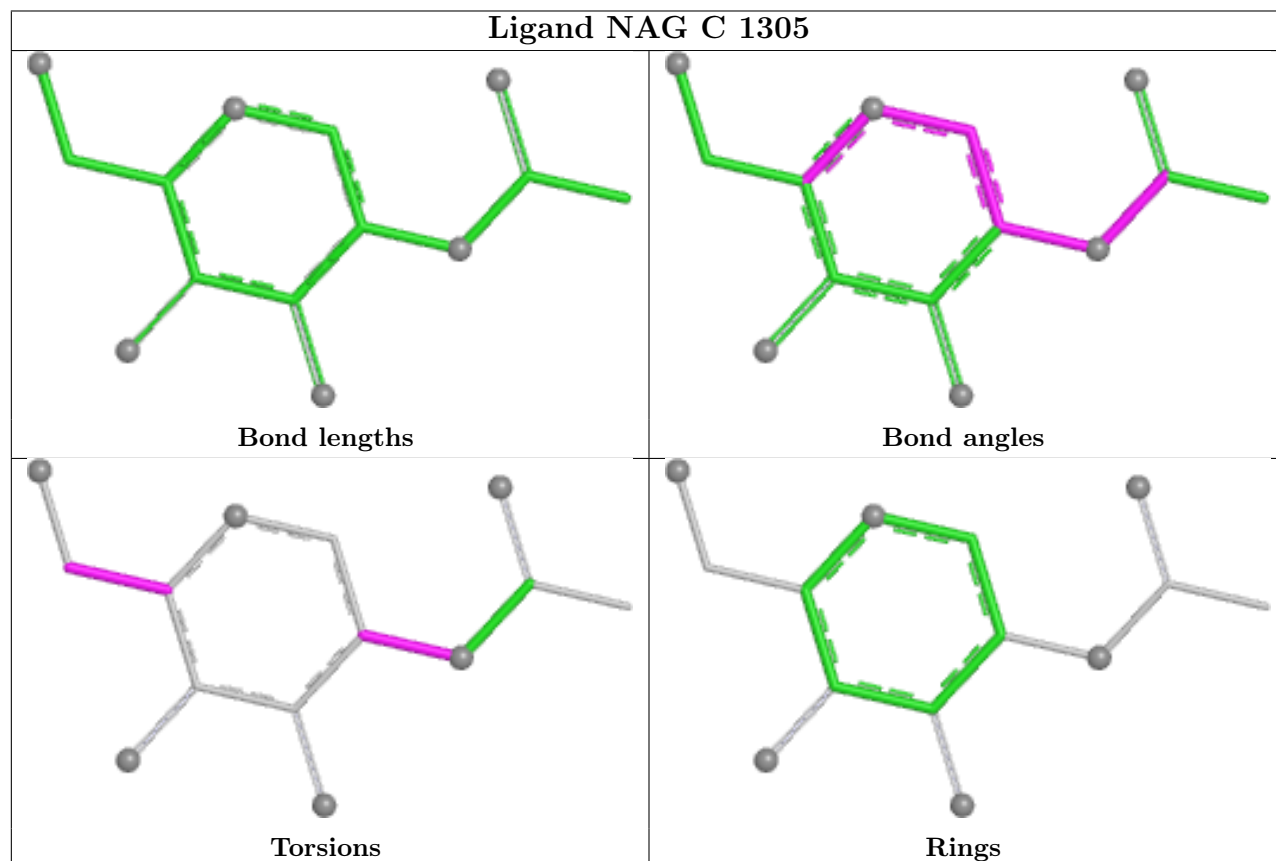
Ligand NAG A 1308

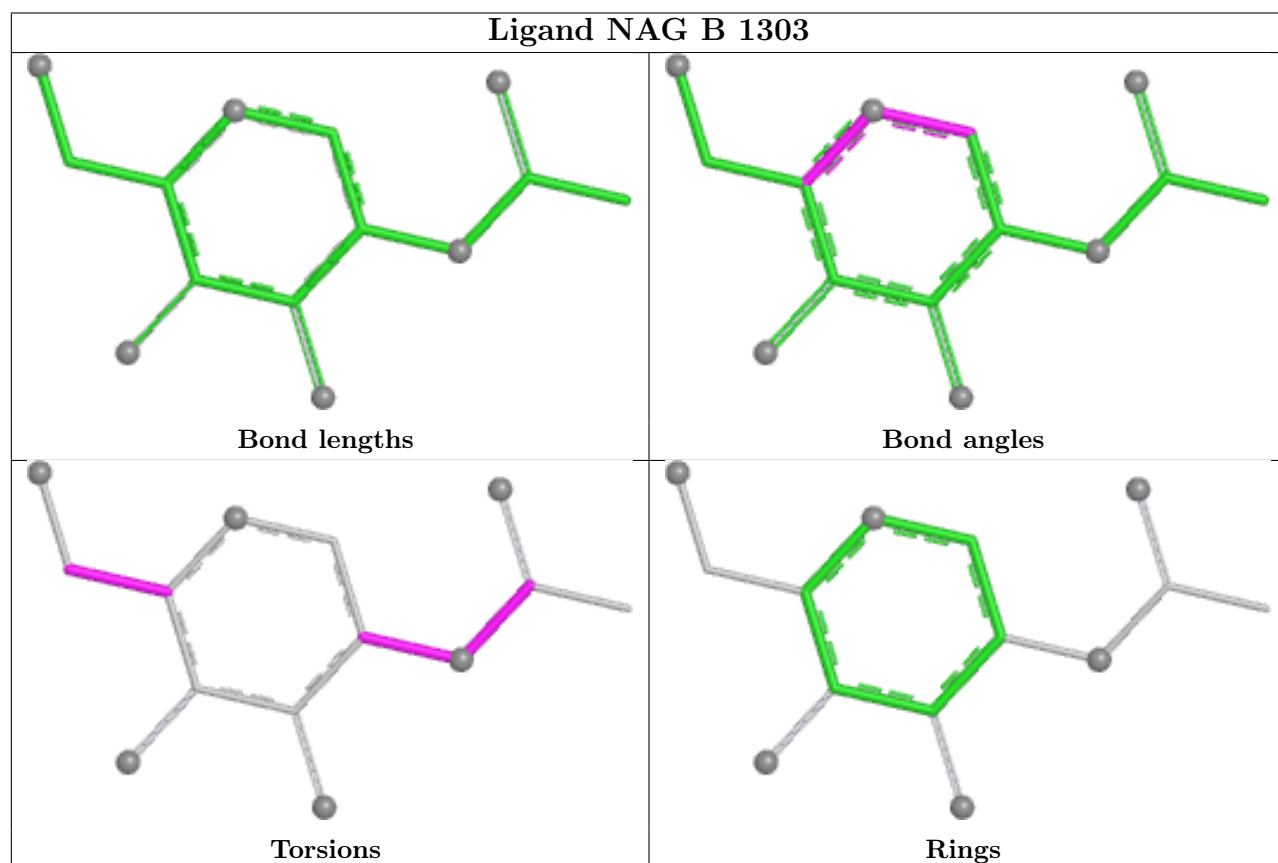
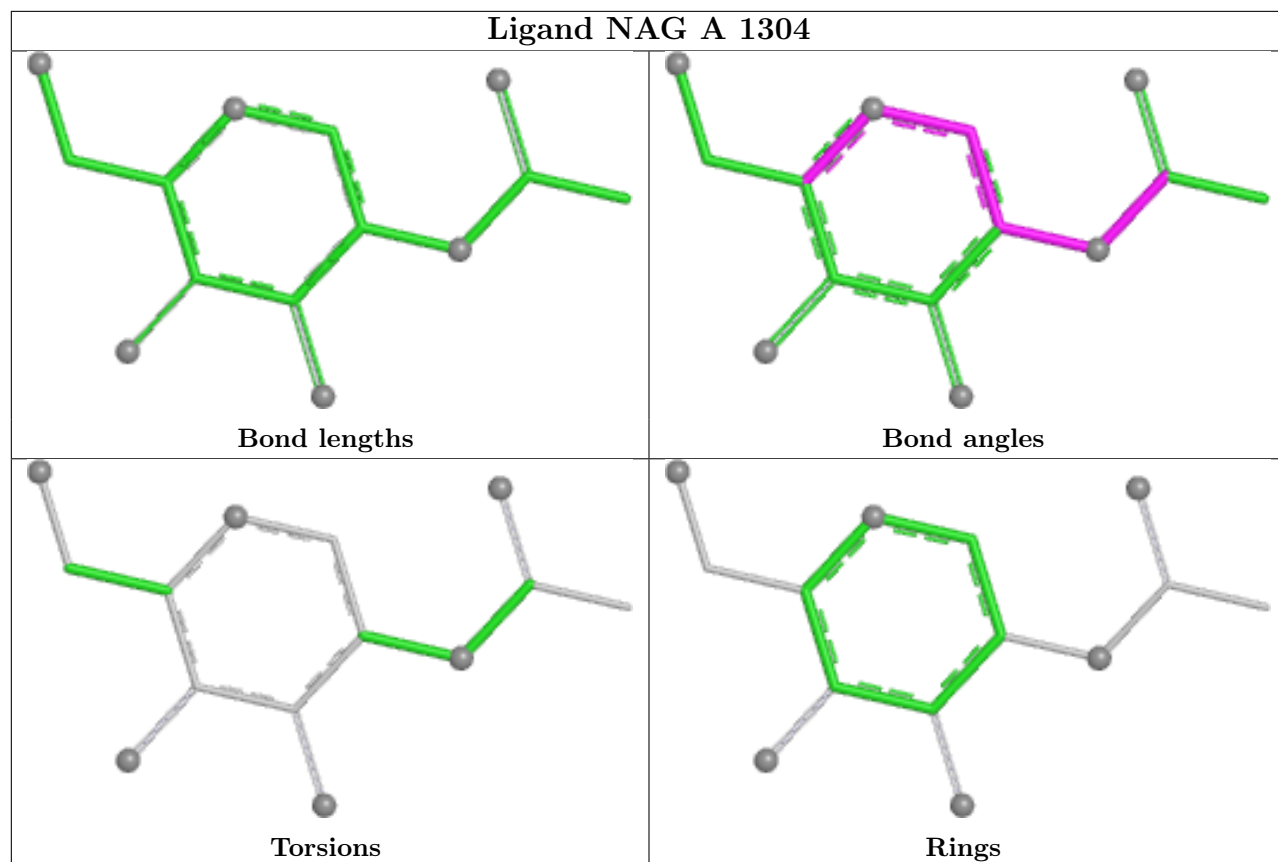


Ligand NAG C 1303

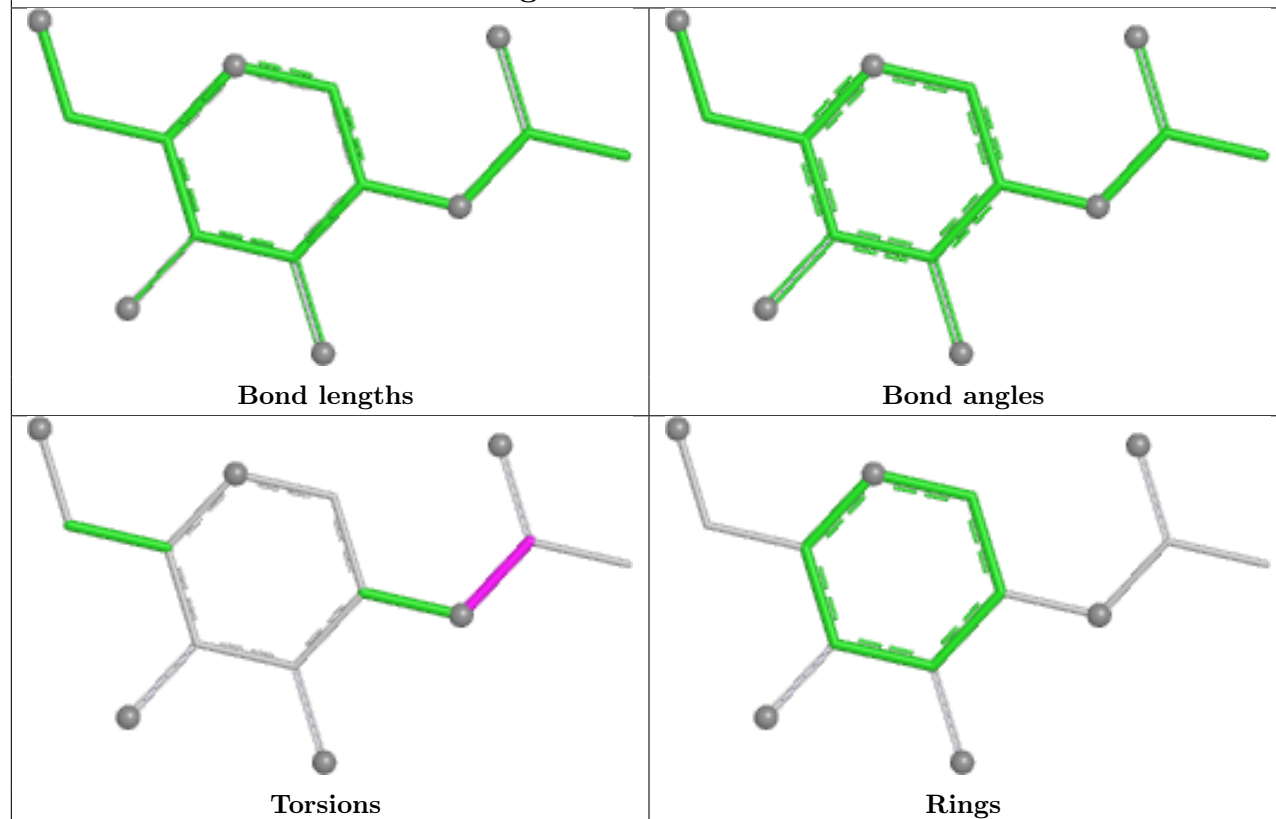


Ligand NAG C 1305

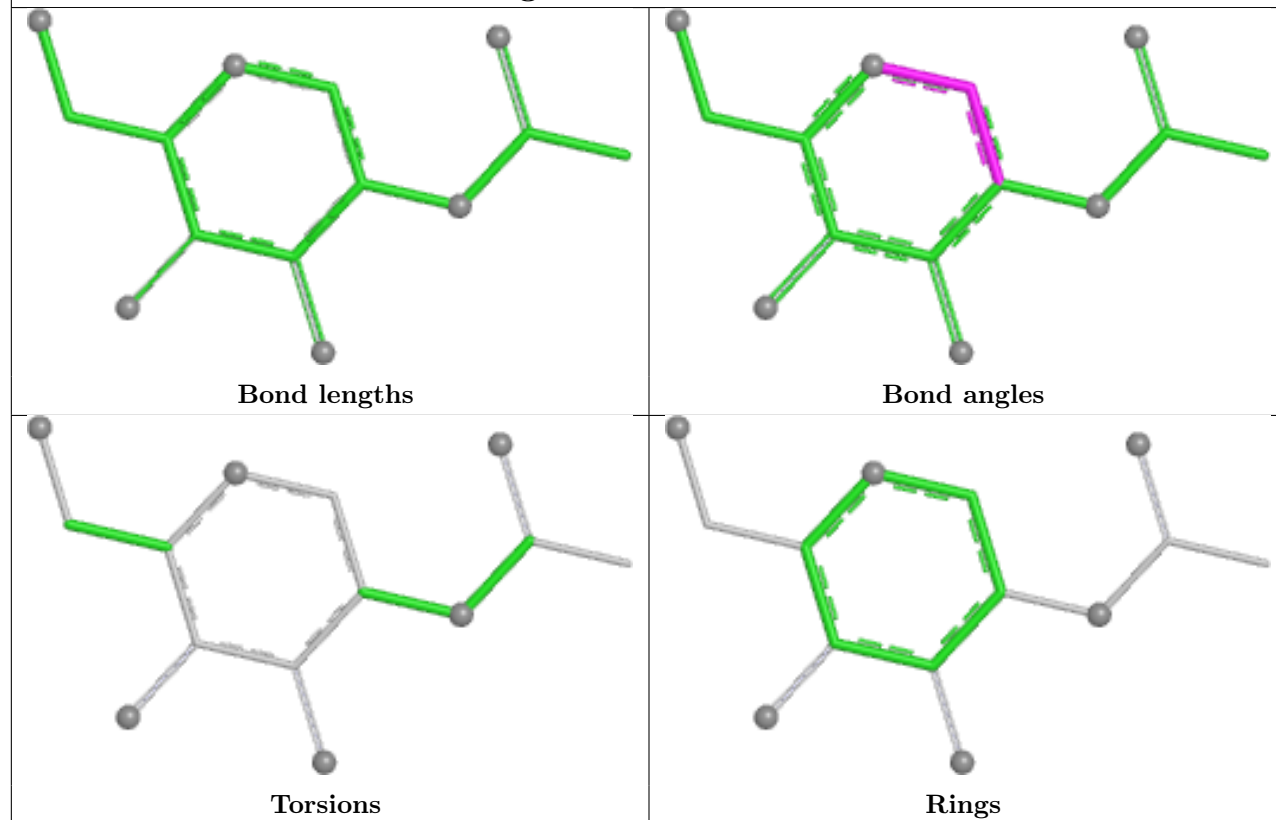


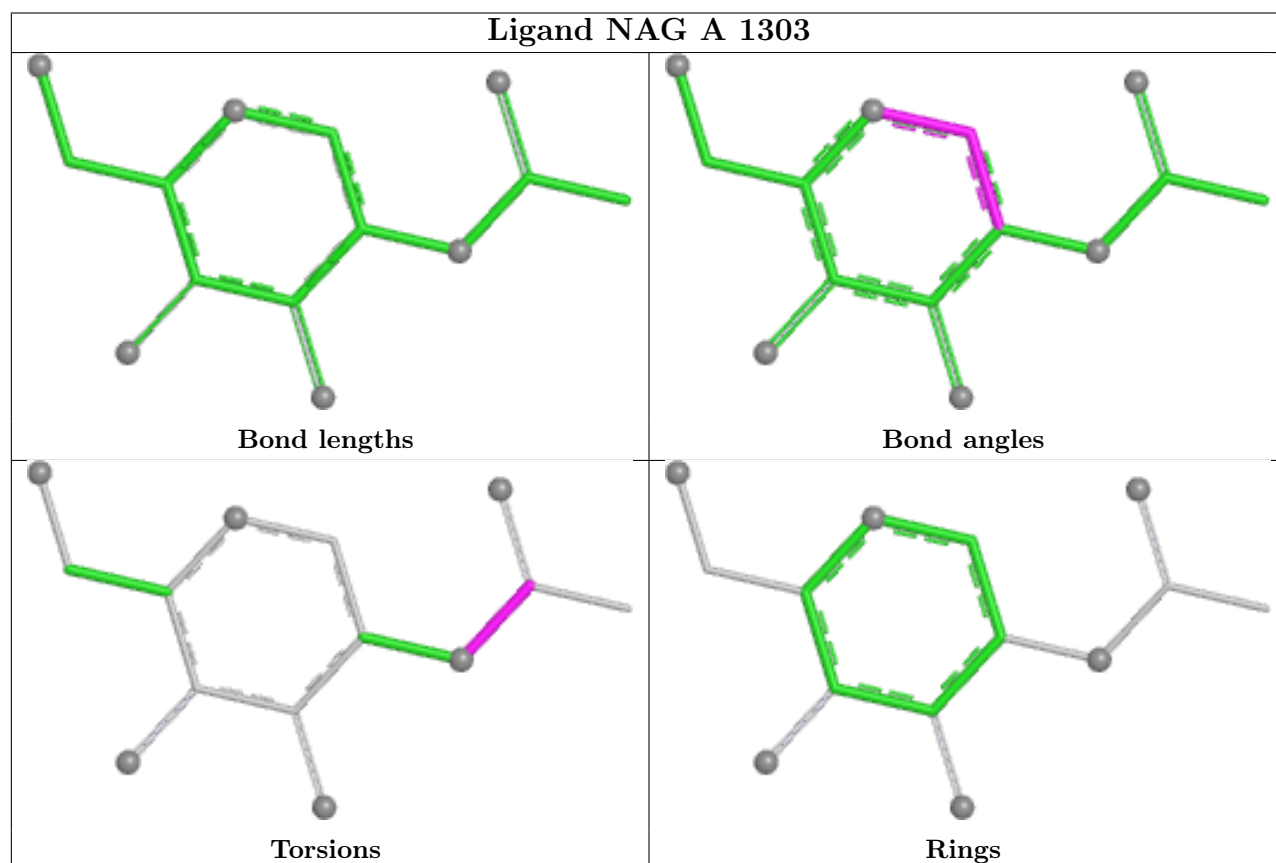
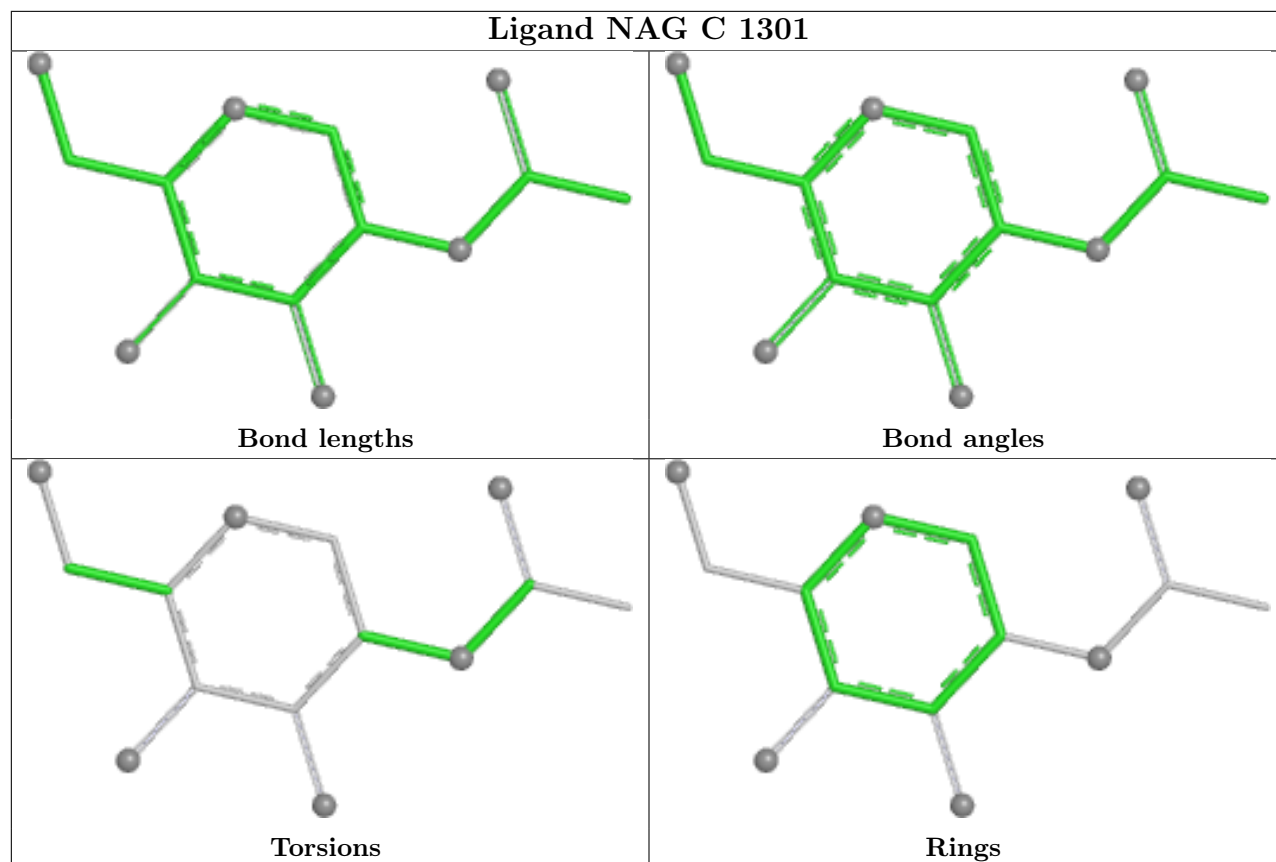


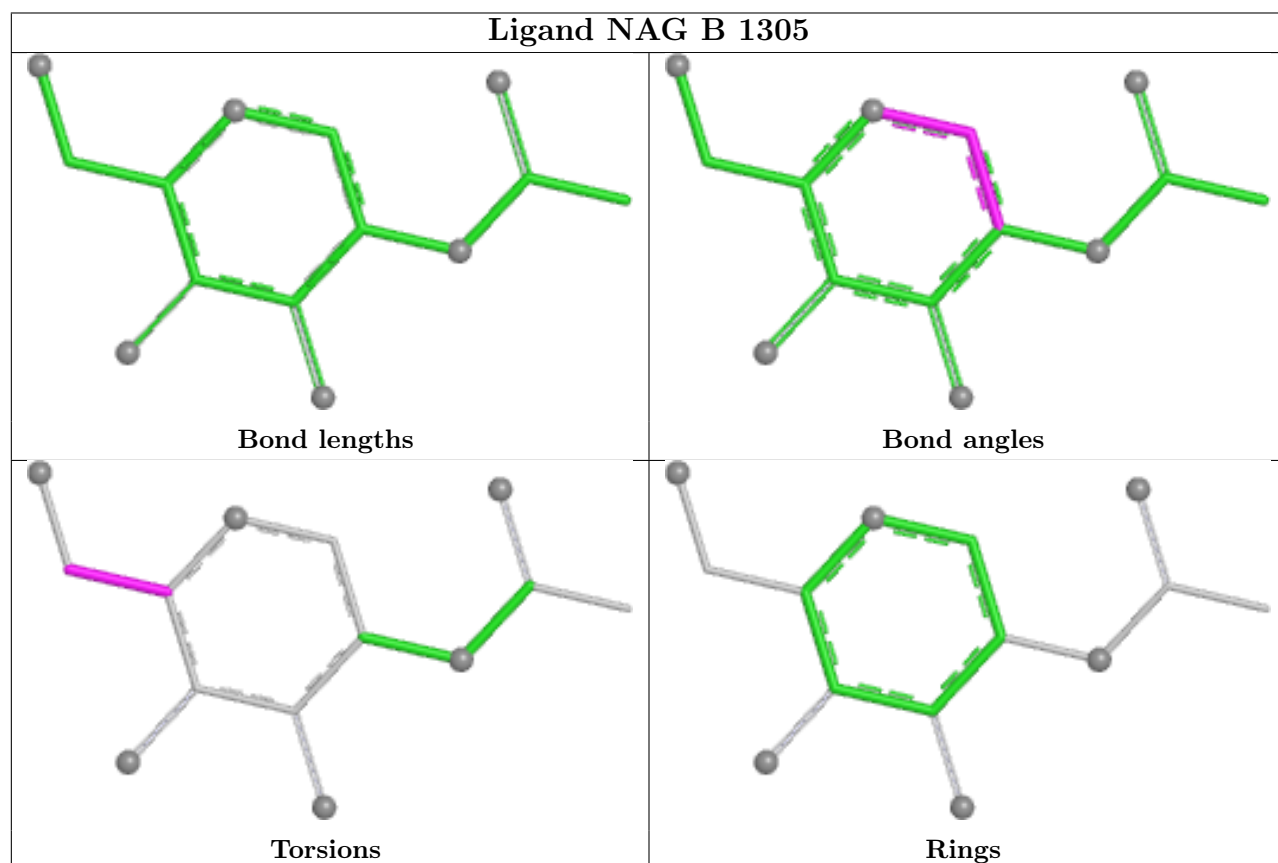
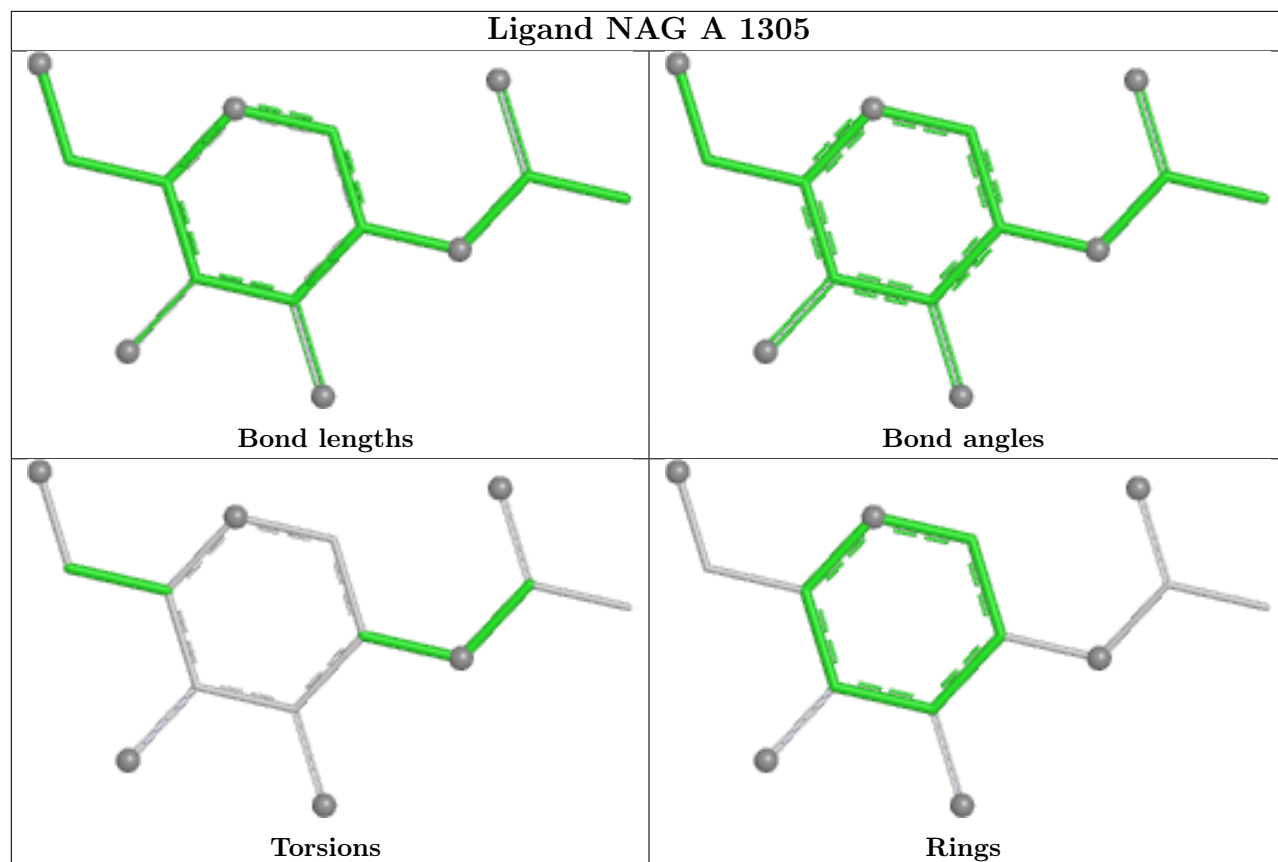
Ligand NAG B 1302

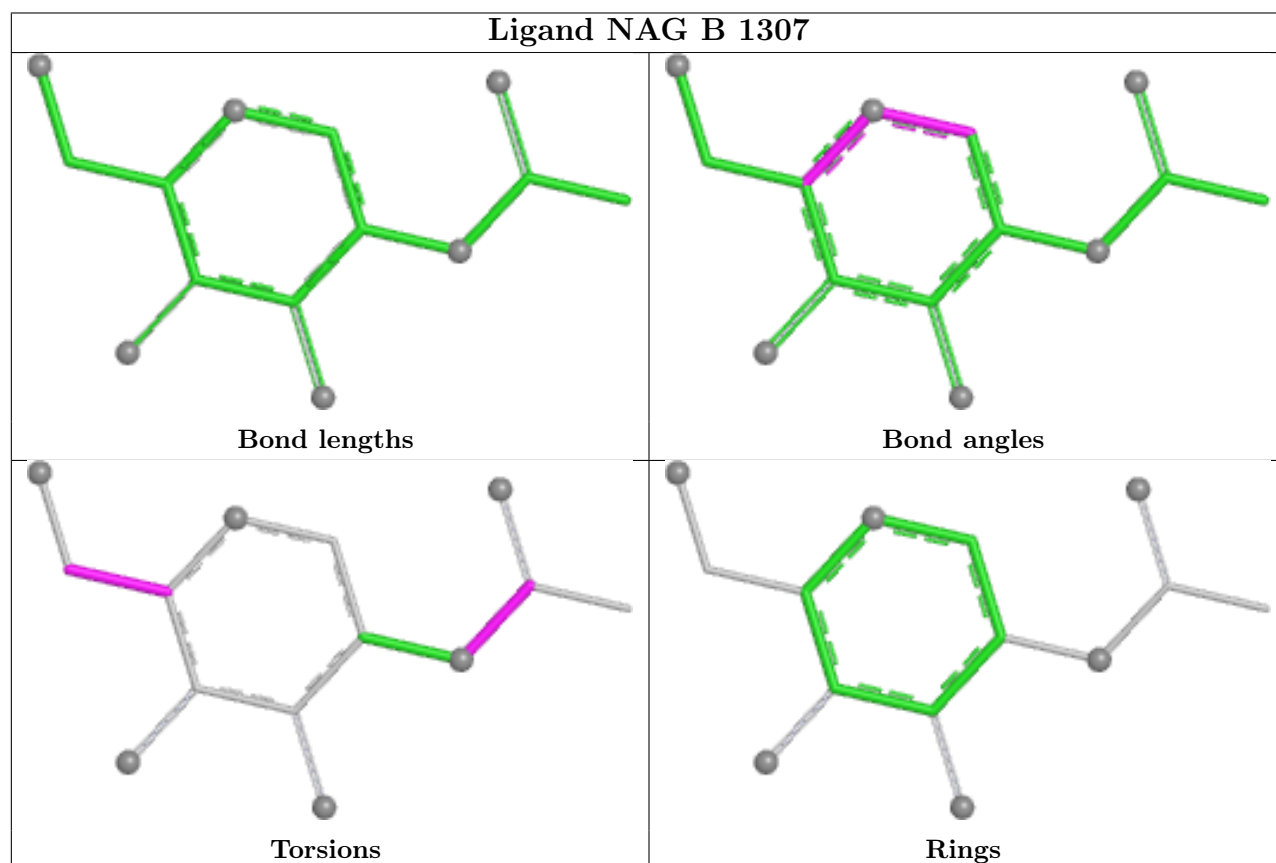
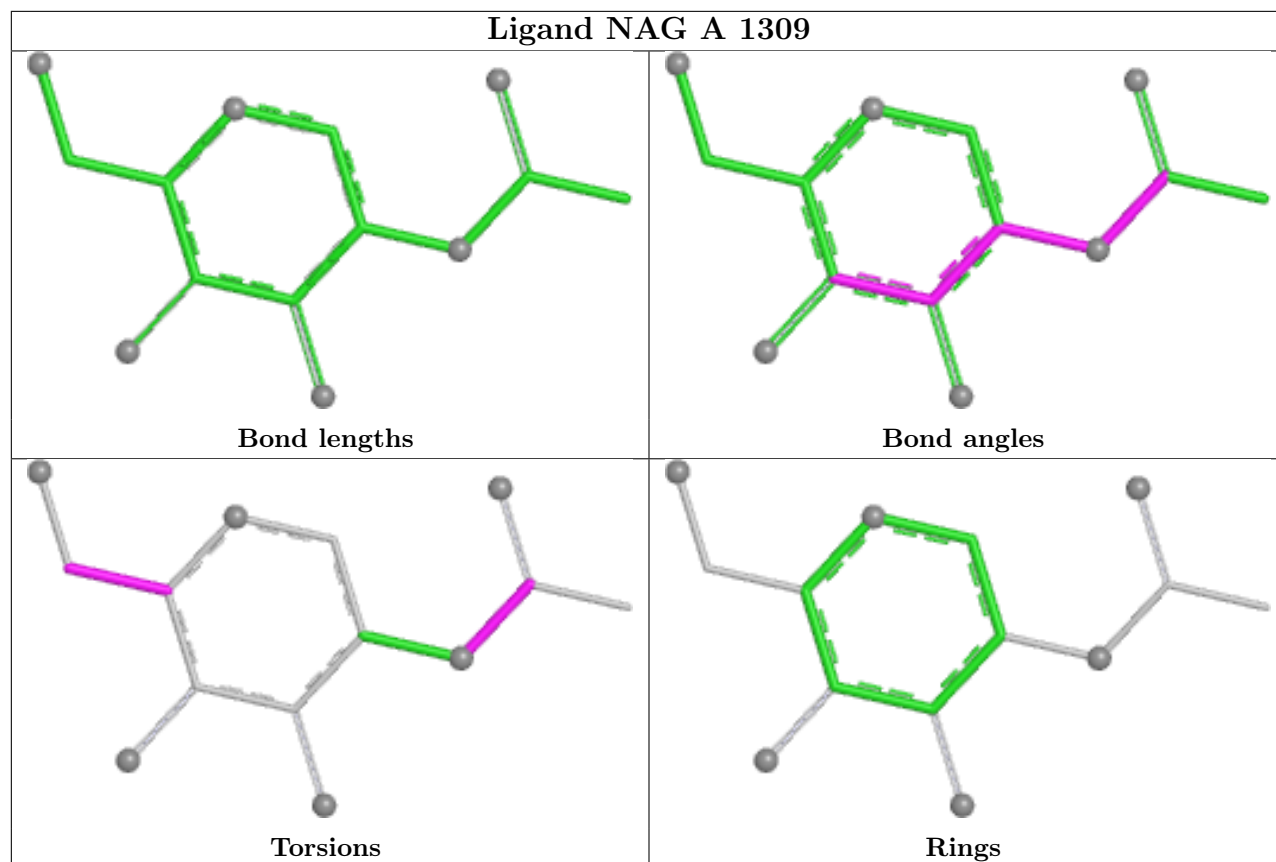


Ligand NAG C 1307

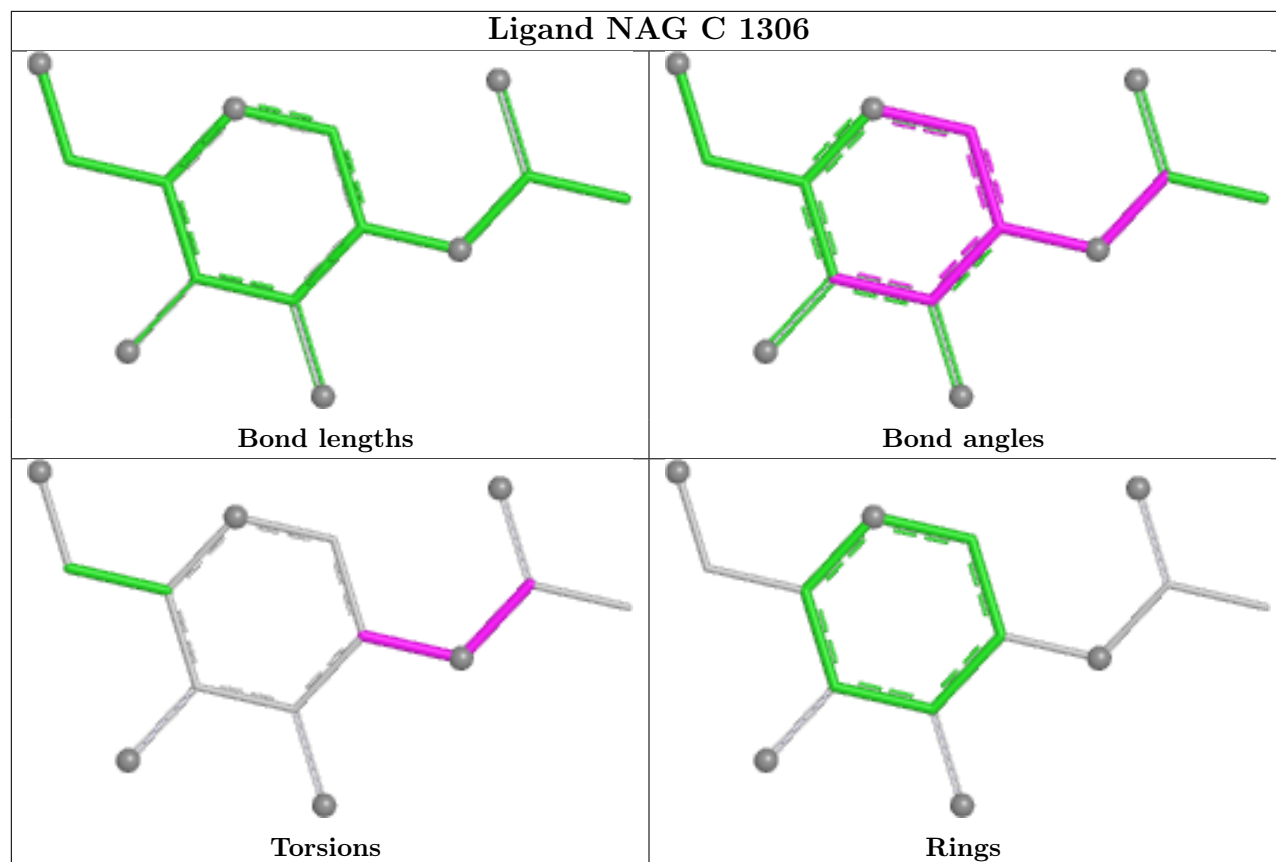




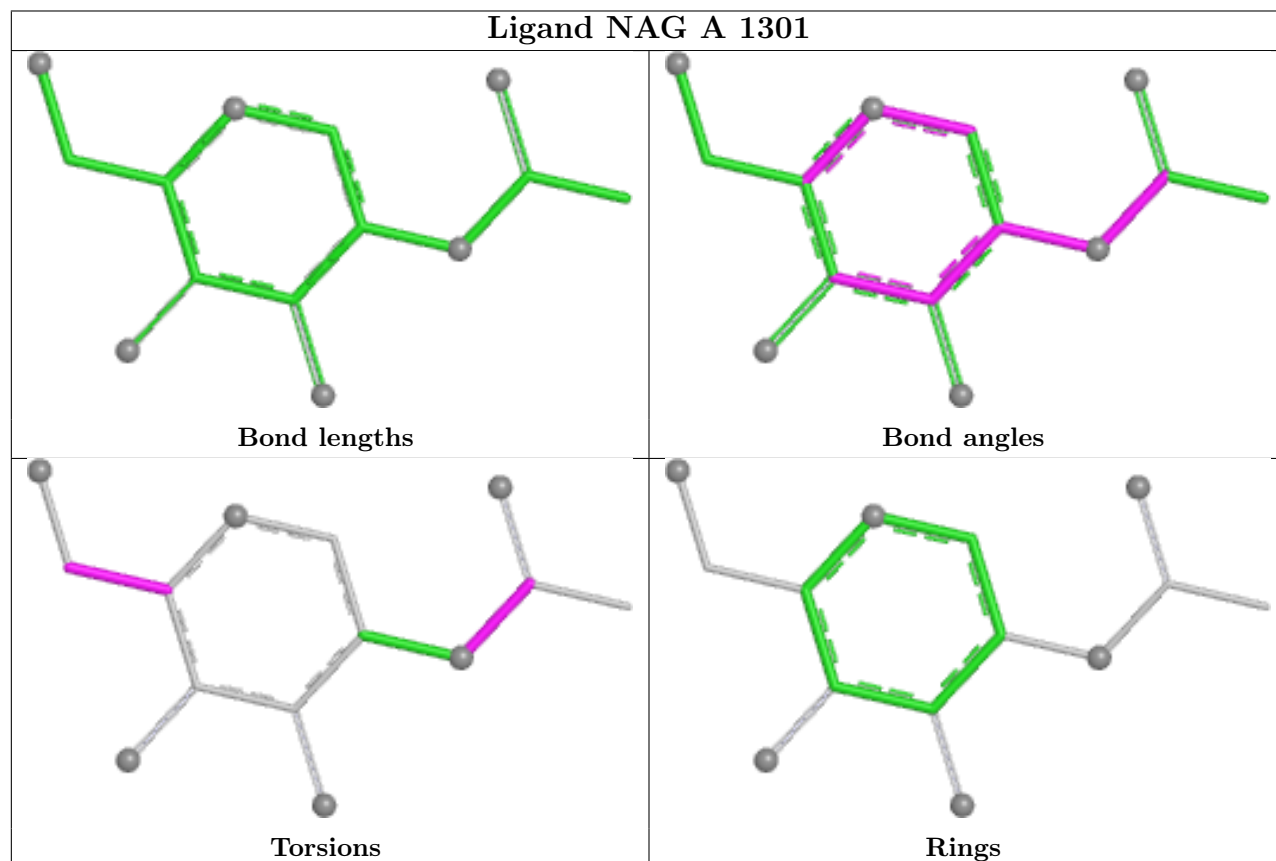




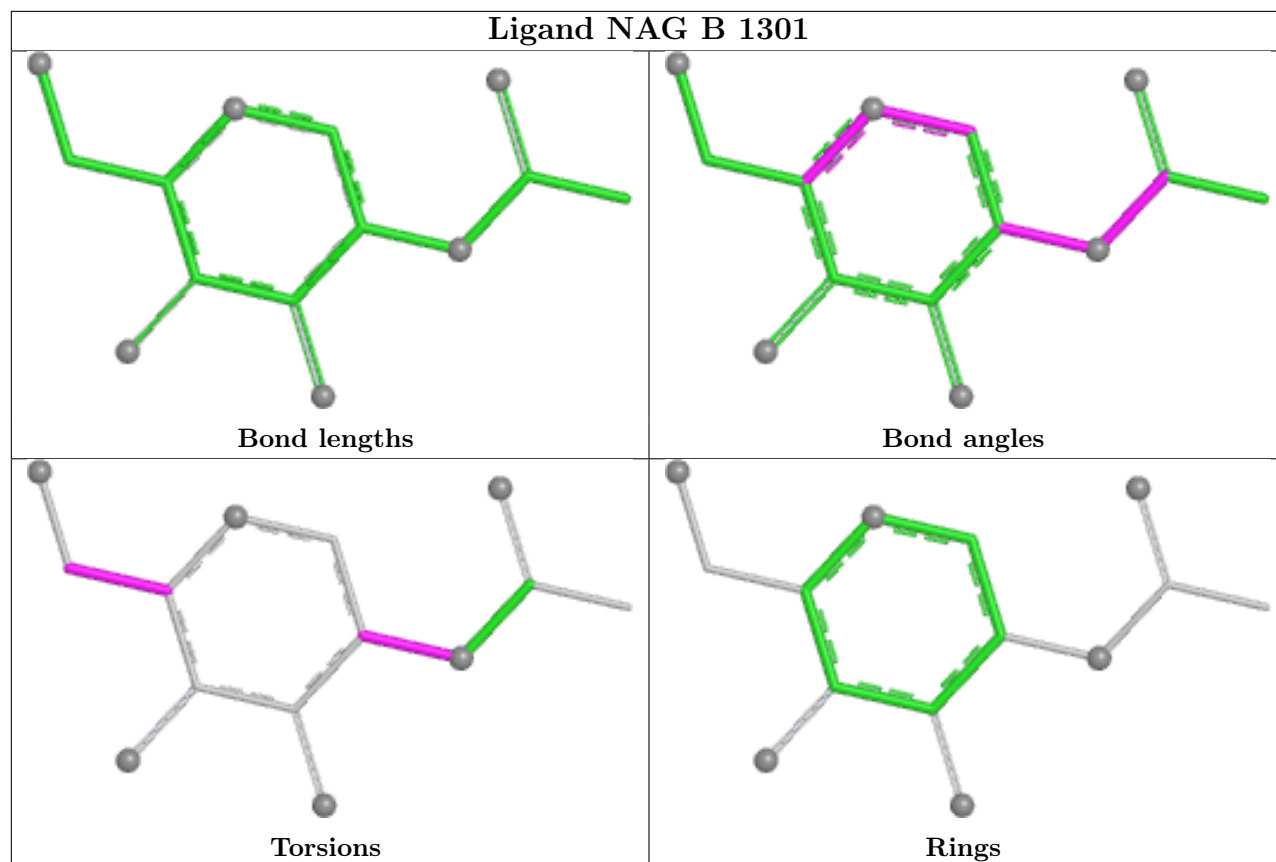
Ligand NAG C 1306



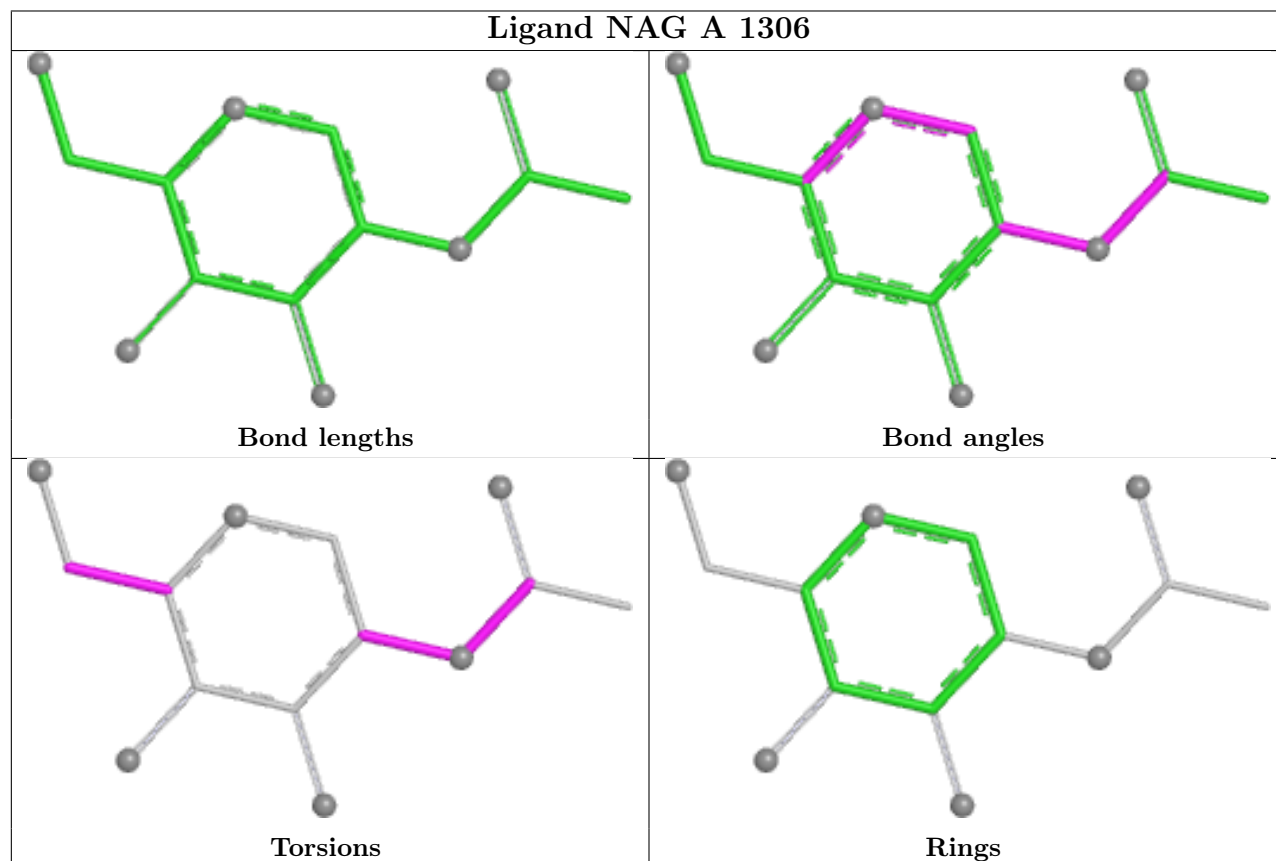
Ligand NAG A 1301



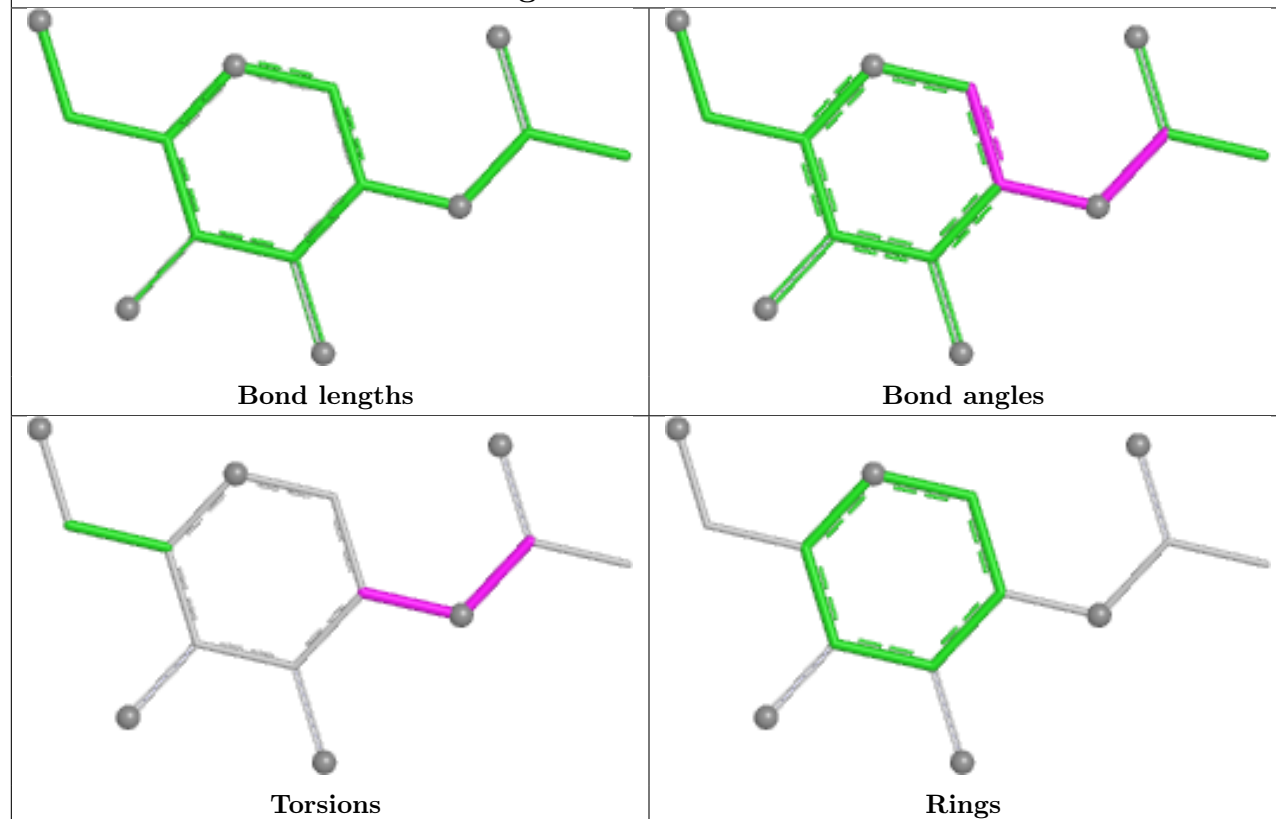
Ligand NAG B 1301



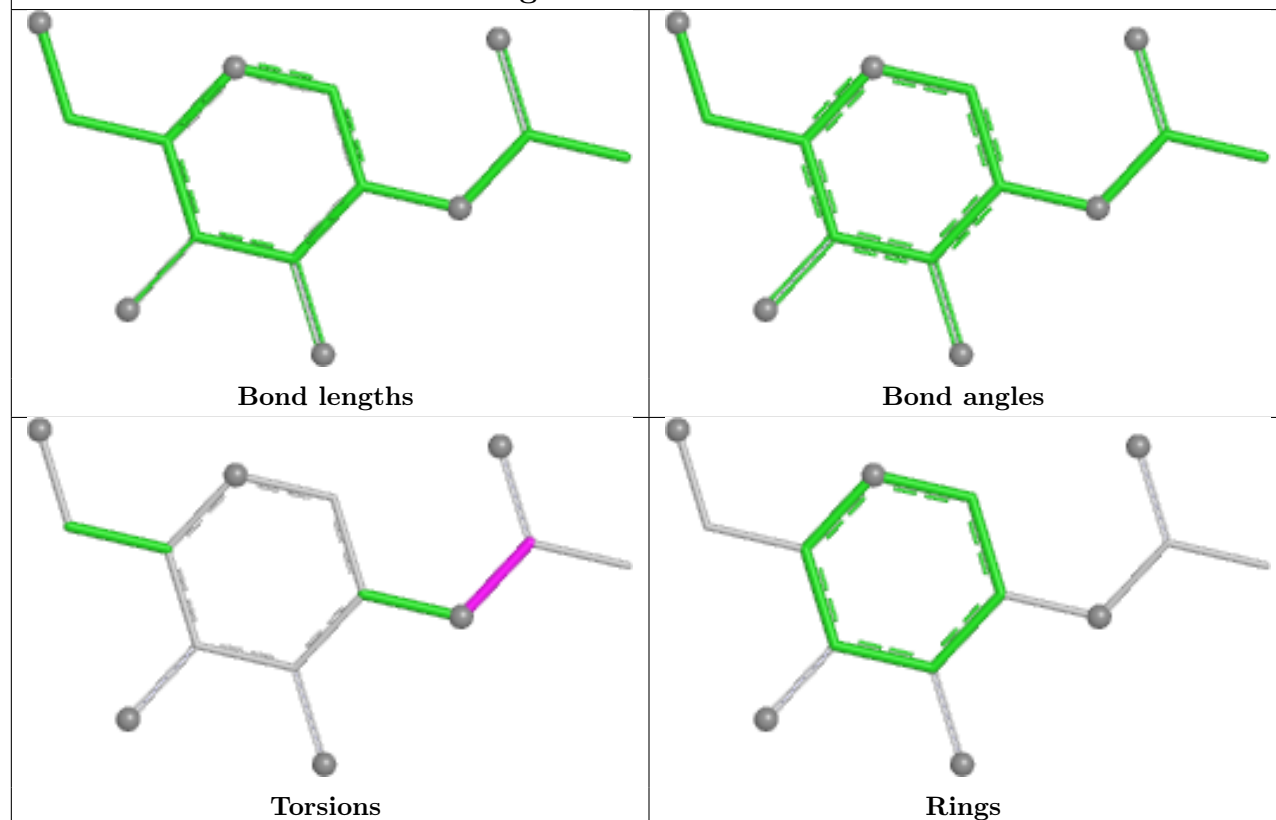
Ligand NAG A 1306

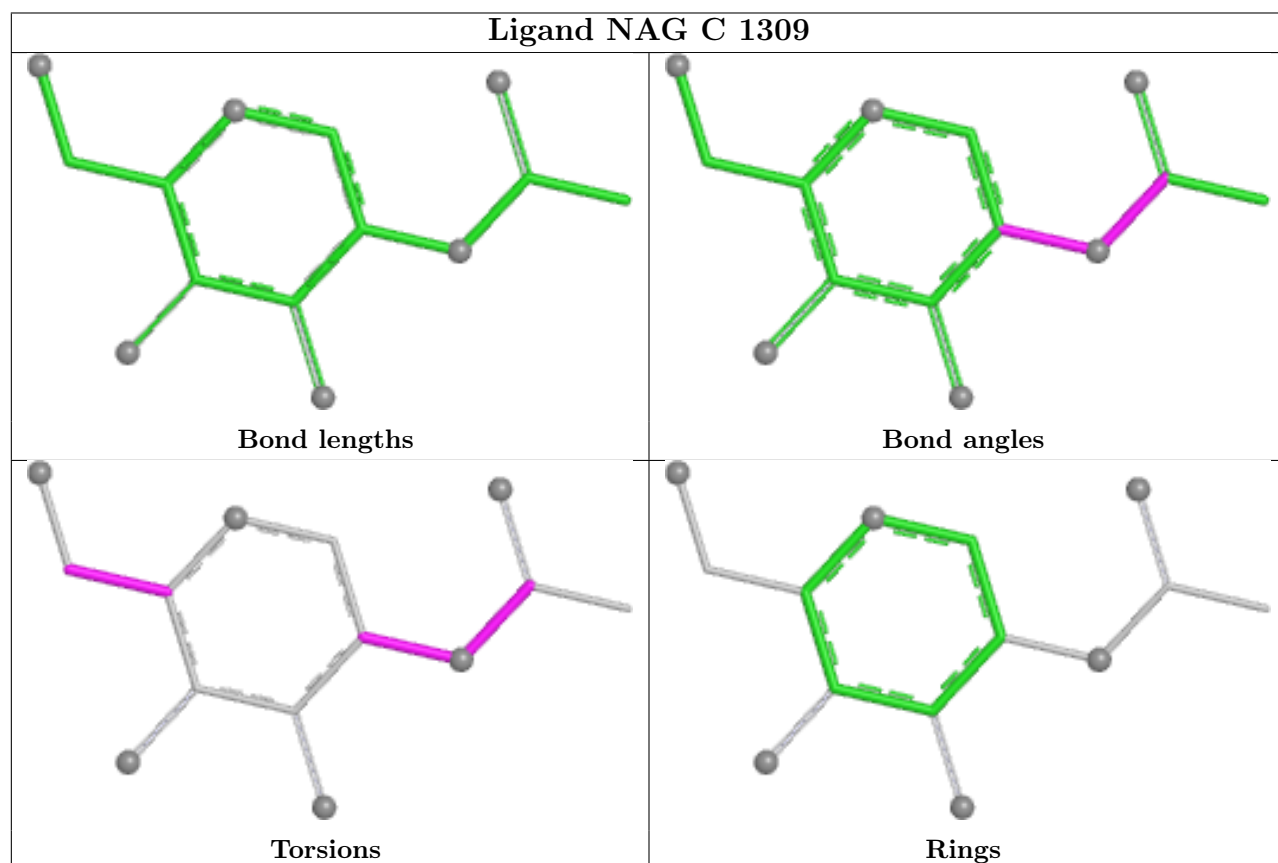
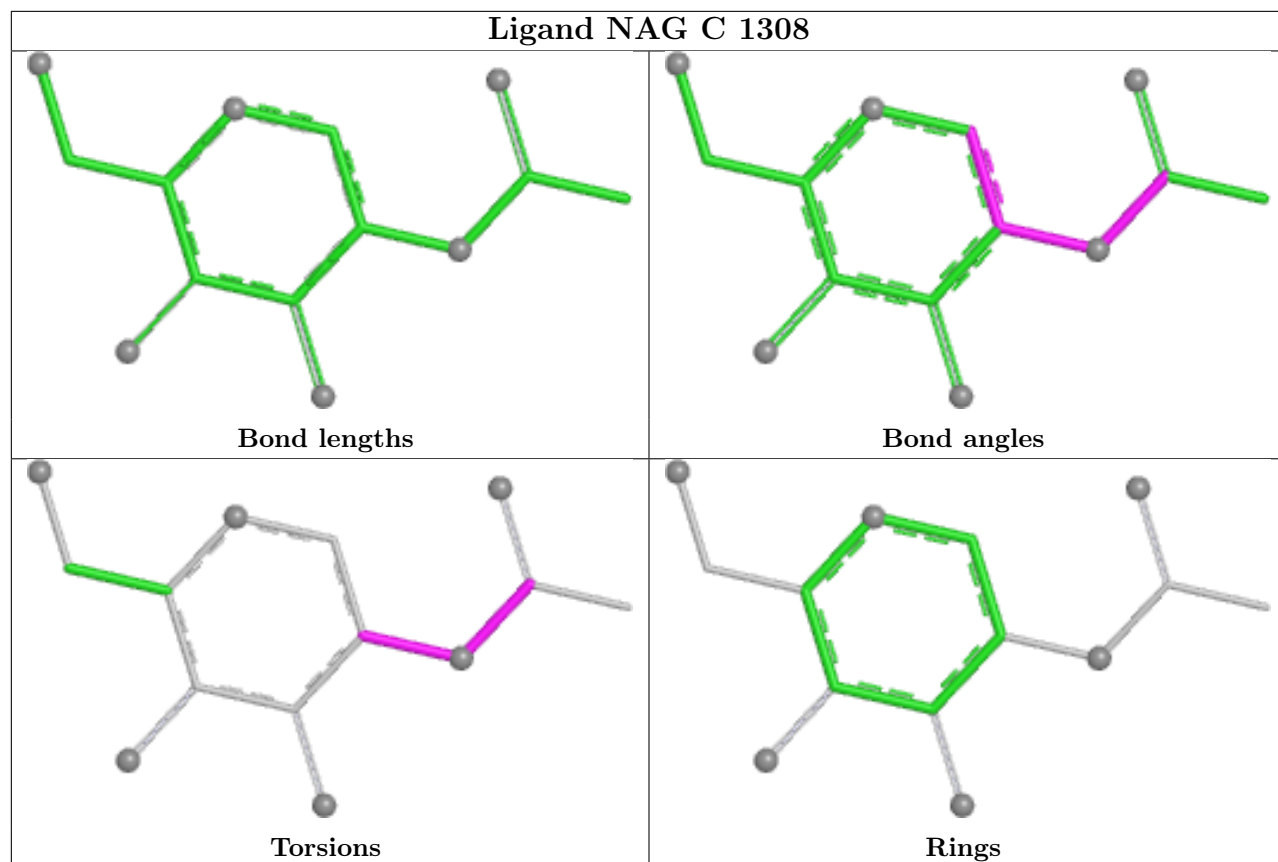


Ligand NAG A 1307

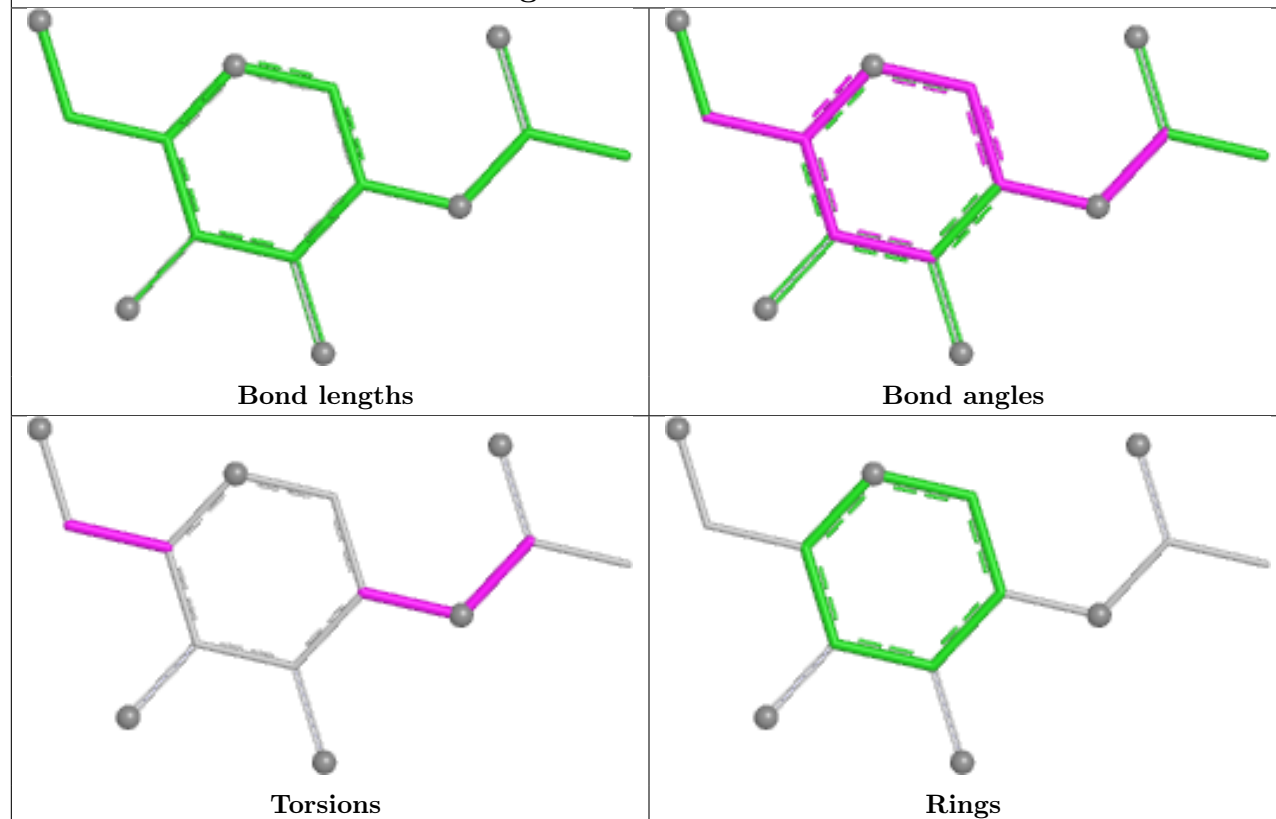


Ligand NAG B 1304

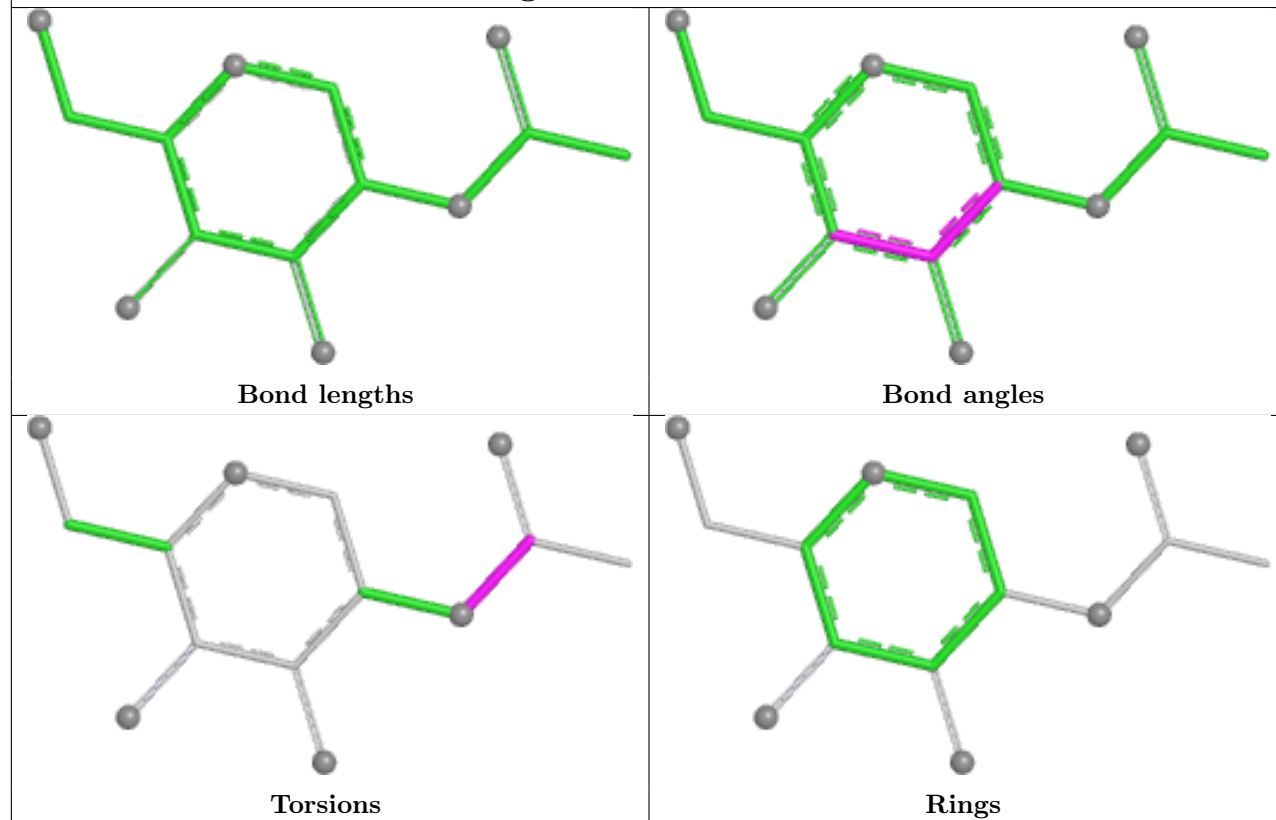


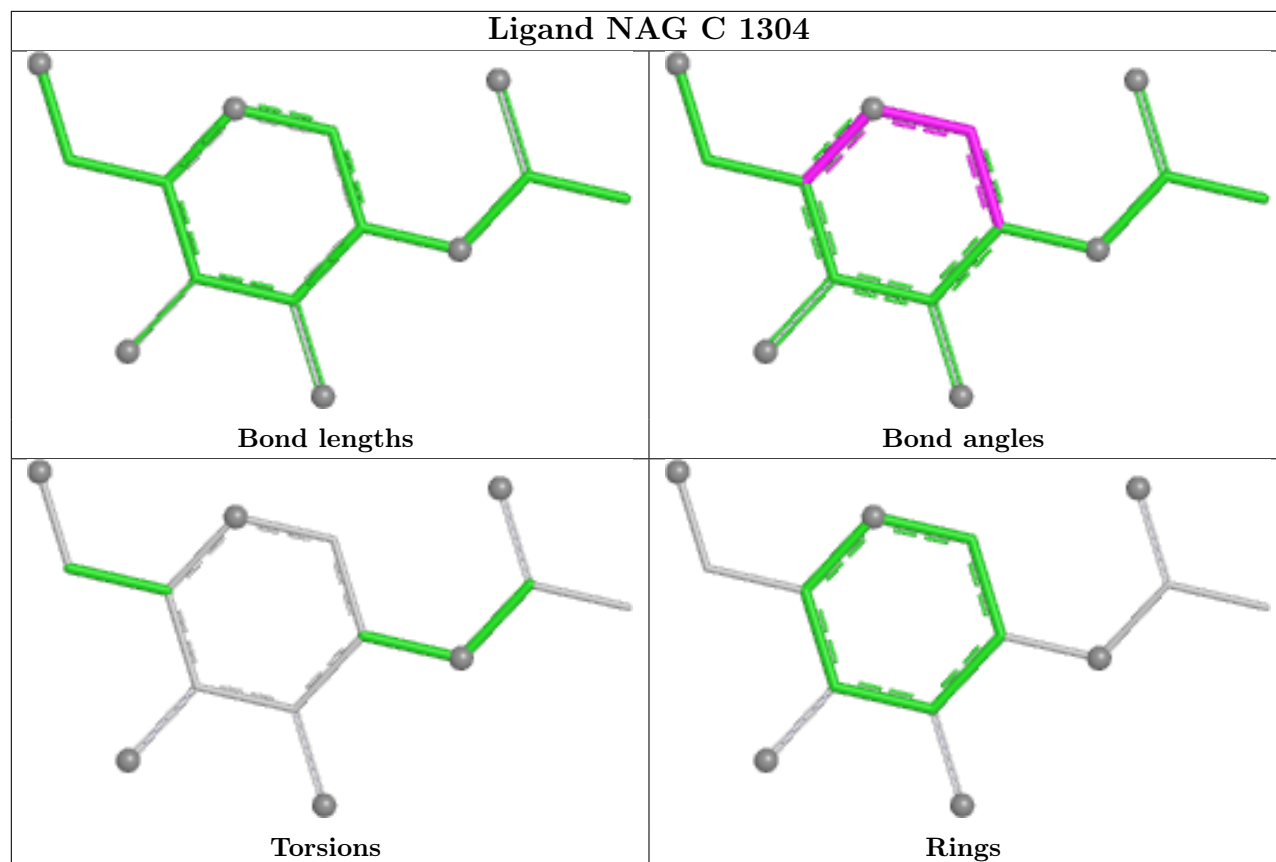


Ligand NAG B 1306



Ligand NAG C 1302





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.

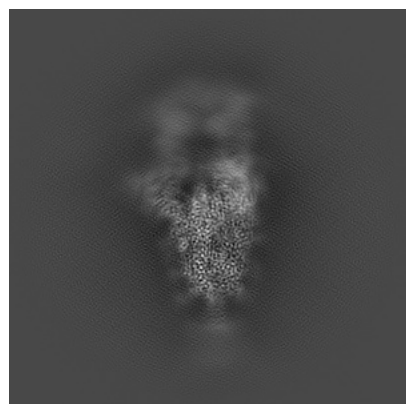
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-63263. These allow visual inspection of the internal detail of the map and identification of artifacts.

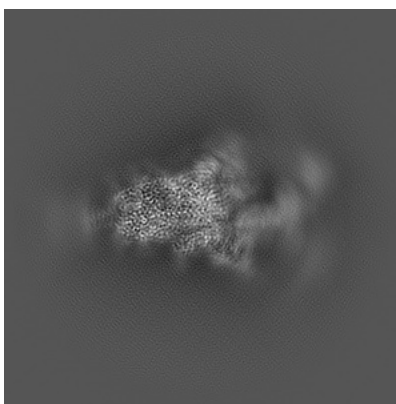
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

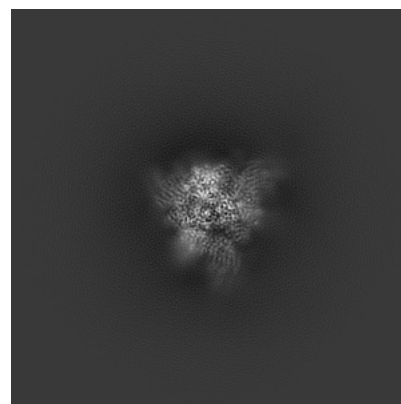
6.1.1 Primary map



X

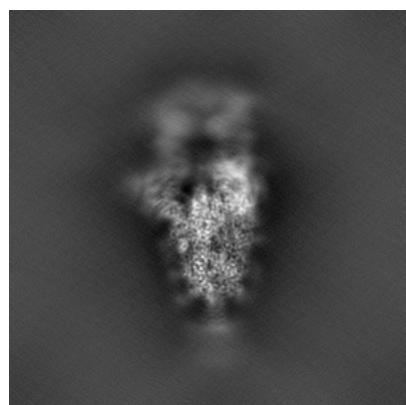


Y

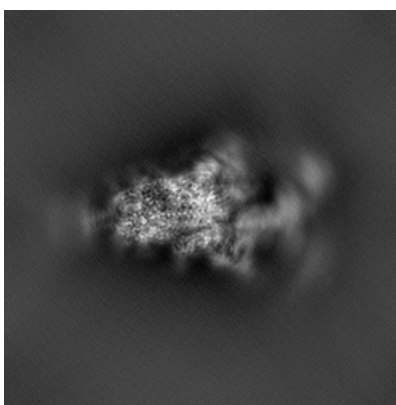


Z

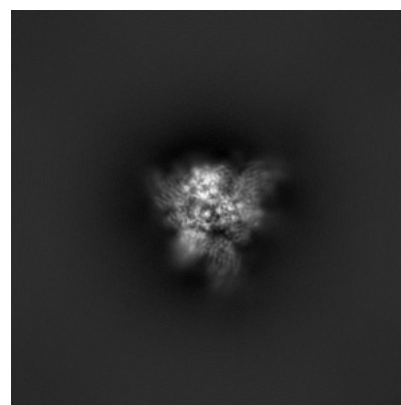
6.1.2 Raw map



X



Y

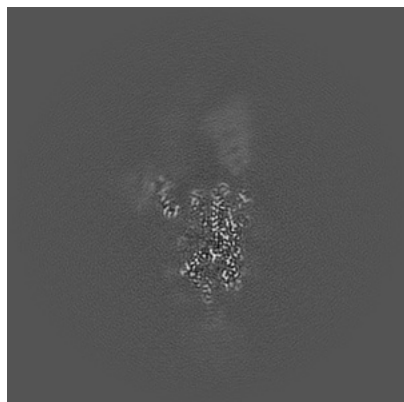


Z

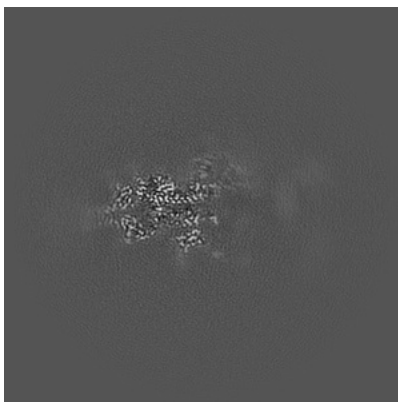
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

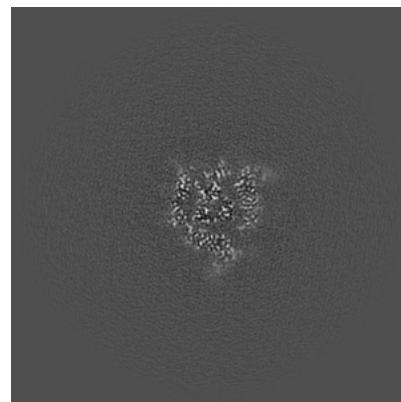
6.2.1 Primary map



X Index: 200

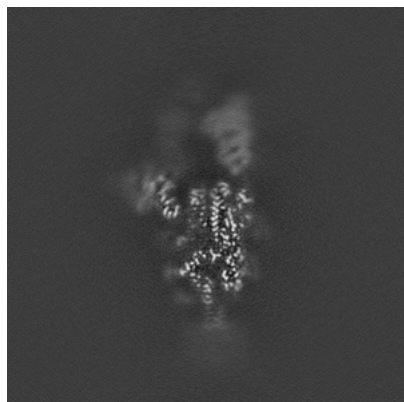


Y Index: 200

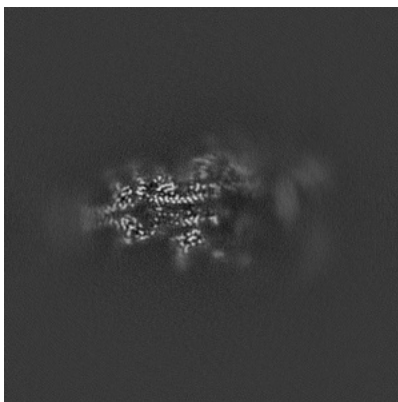


Z Index: 200

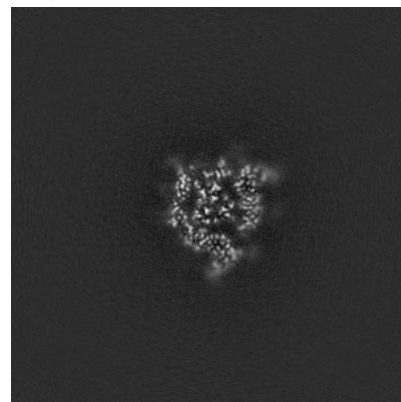
6.2.2 Raw map



X Index: 200



Y Index: 200

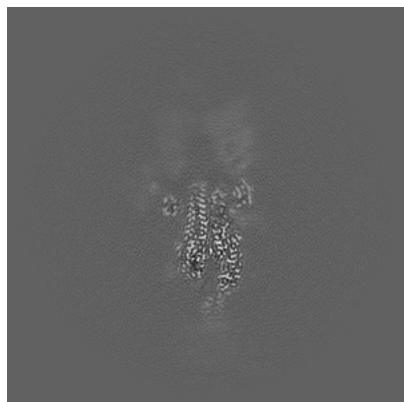


Z Index: 200

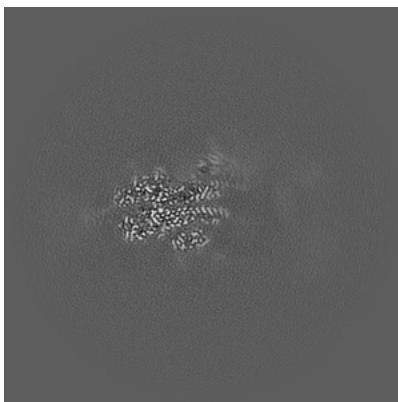
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

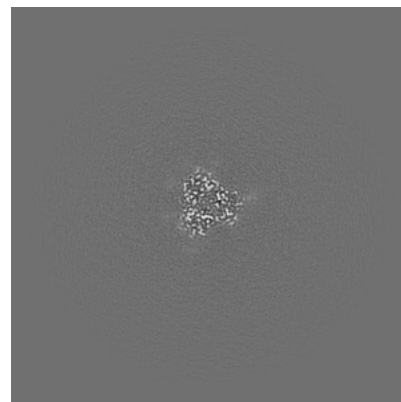
6.3.1 Primary map



X Index: 195

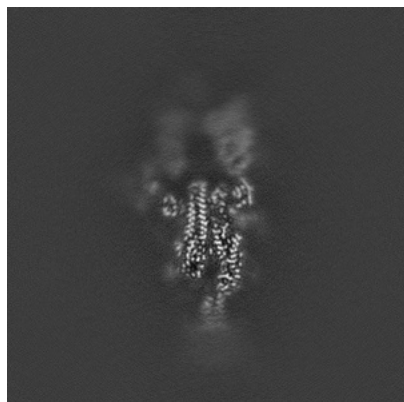


Y Index: 196

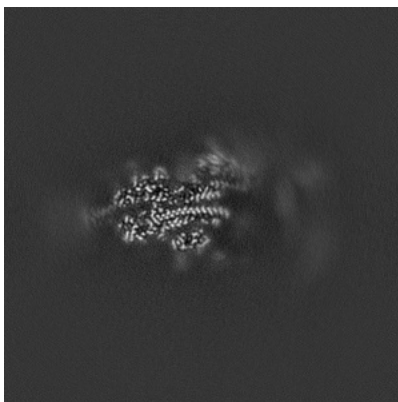


Z Index: 146

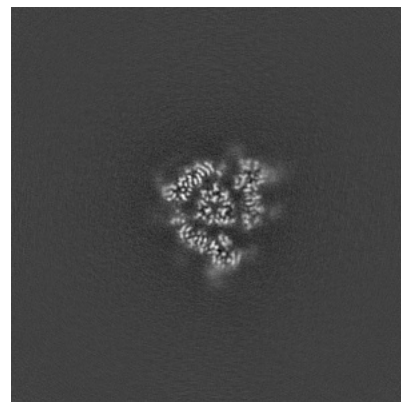
6.3.2 Raw map



X Index: 195



Y Index: 196

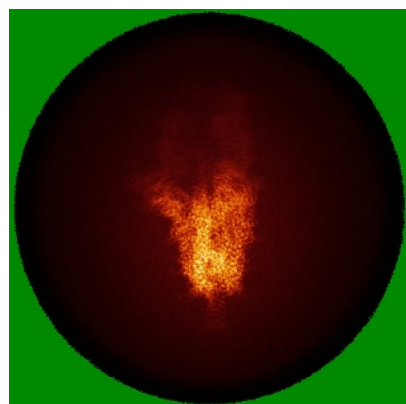


Z Index: 205

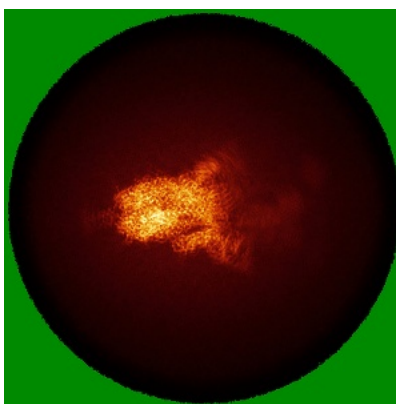
The images above show the largest variance slices of the map in three orthogonal directions.

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

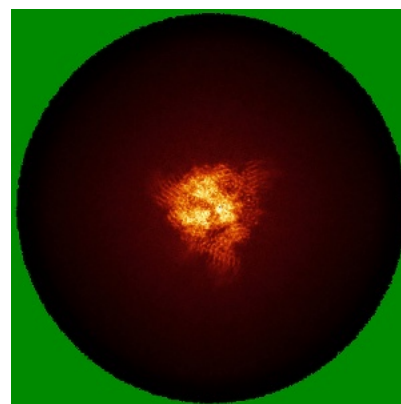
6.4.1 Primary map



X

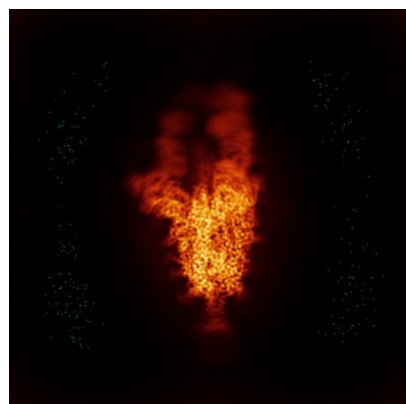


Y

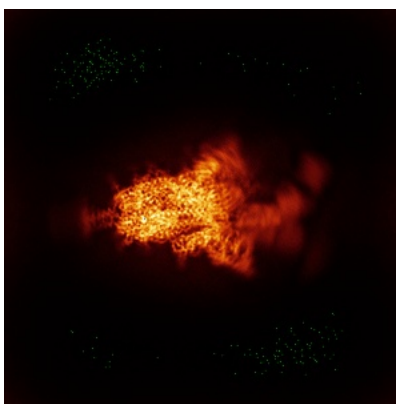


Z

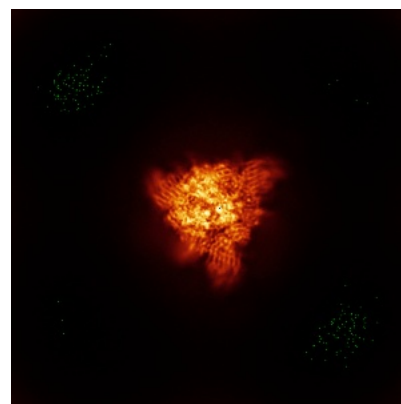
6.4.2 Raw map



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



X



Y



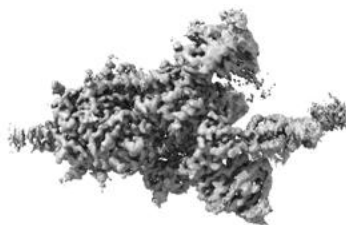
Z

The images above show the 3D surface view of the map at the recommended contour level 0.286. 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



X



Y



Z

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.

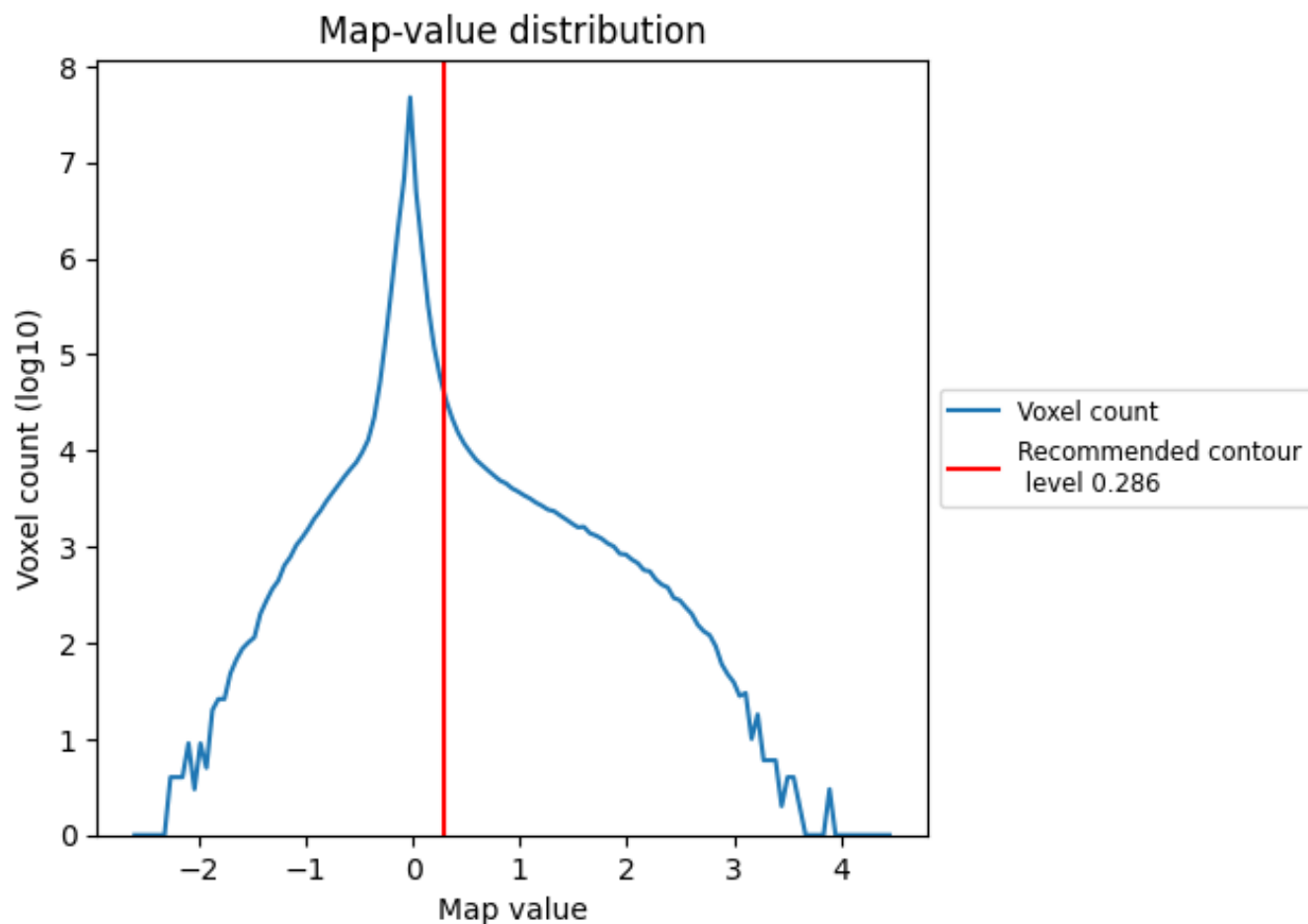
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

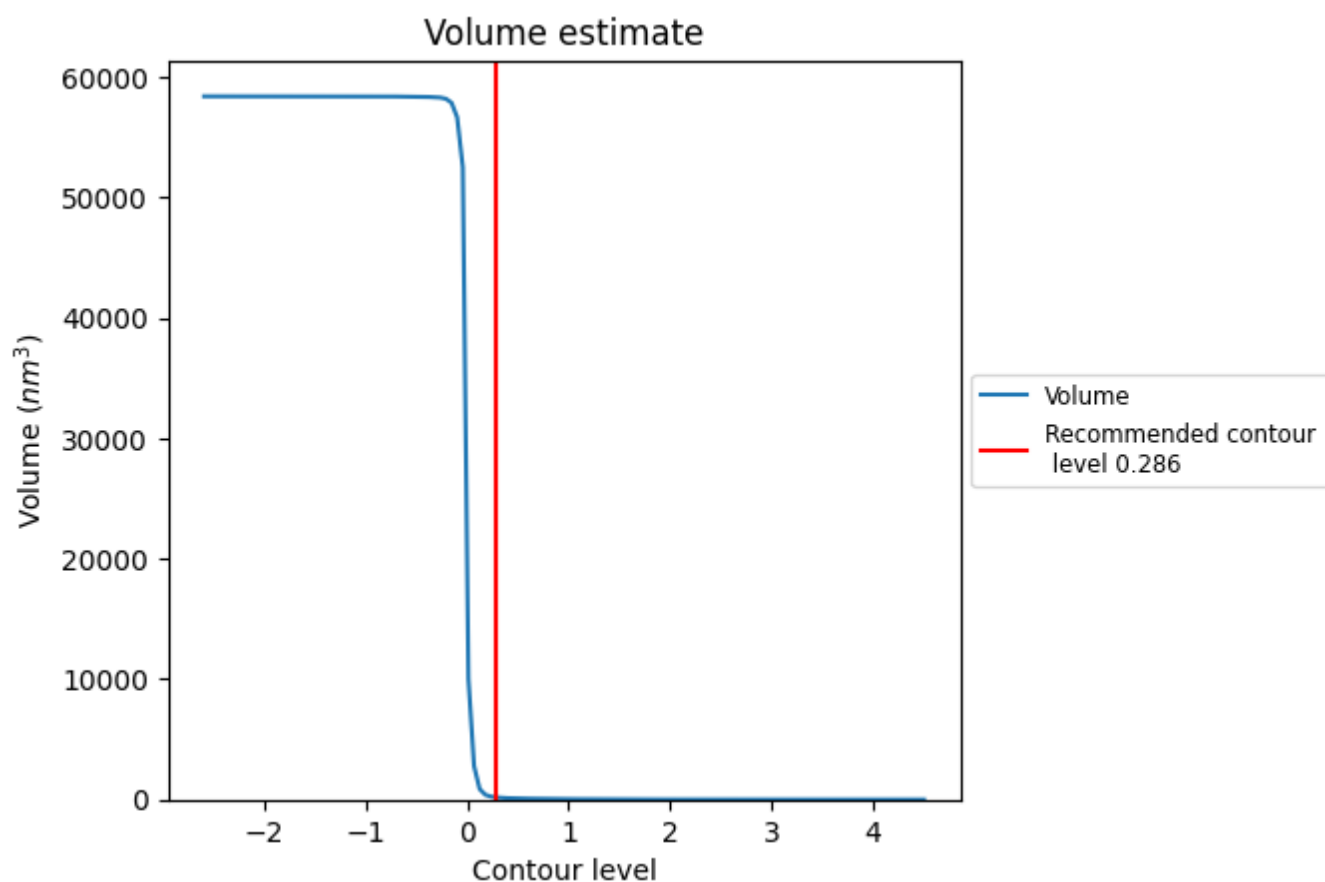
This section contains the results of statistical analysis of the map.

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.

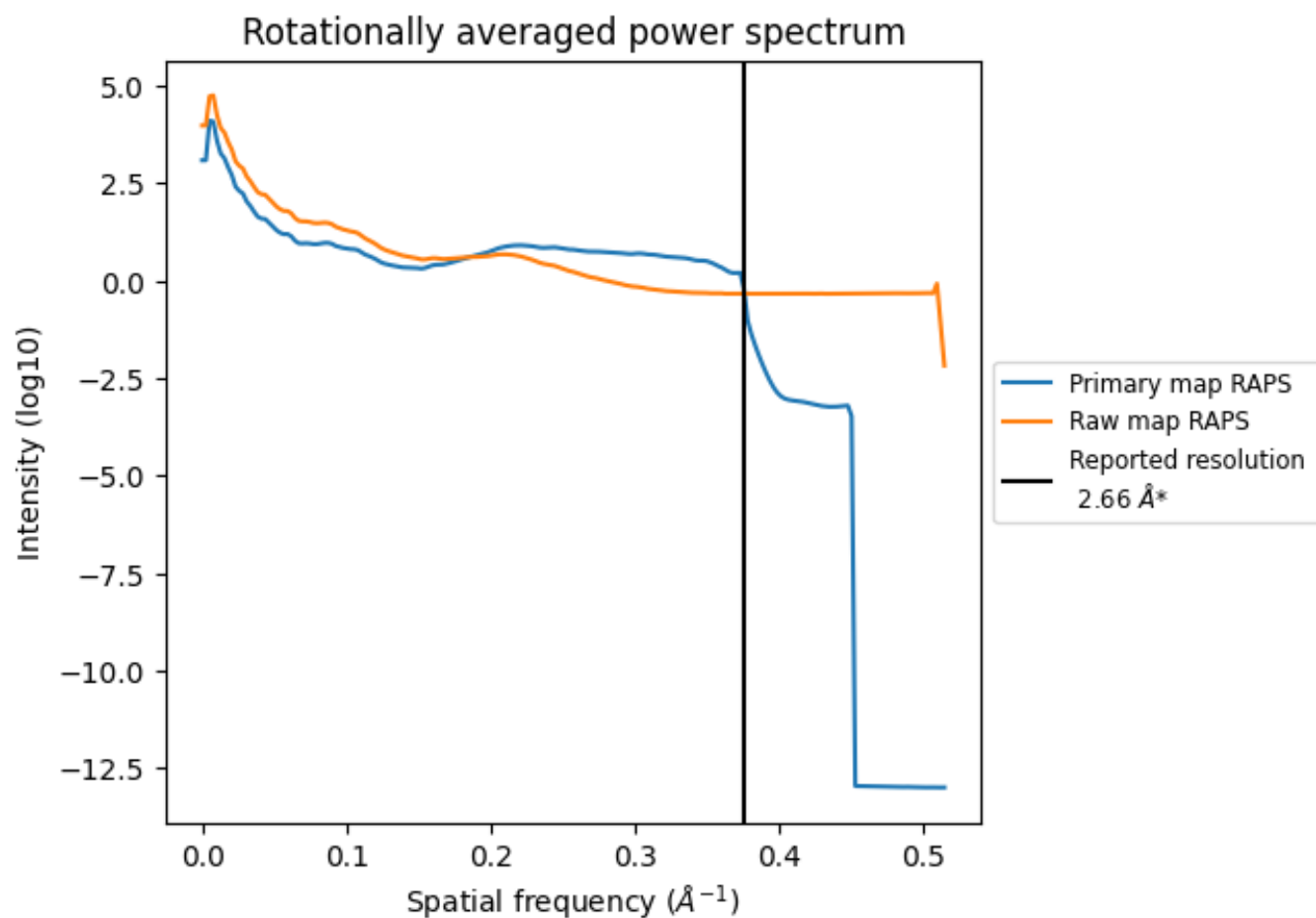
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 182 nm³; this corresponds to an approximate mass of 164 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 ⓘ

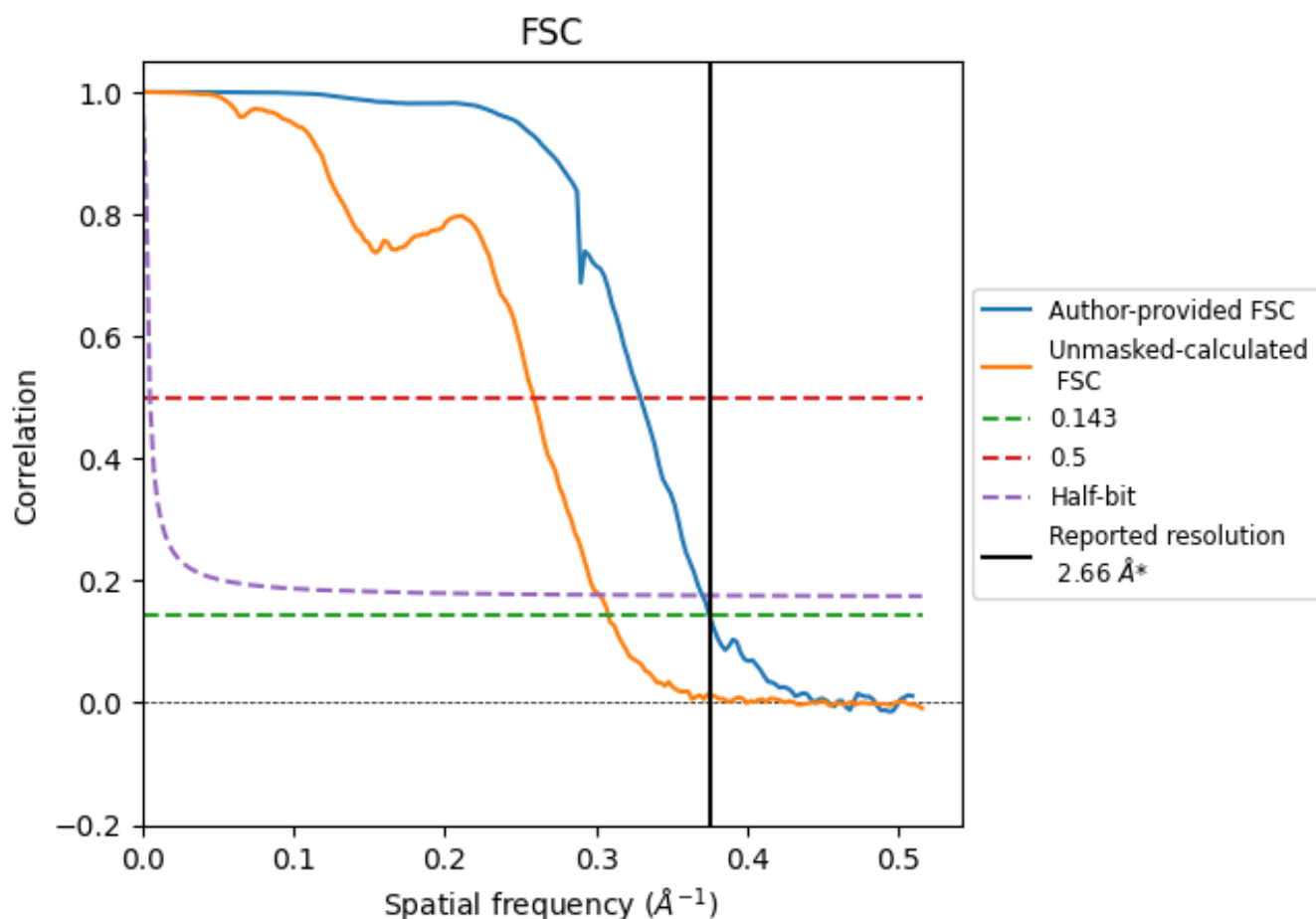


*Reported resolution corresponds to spatial frequency of 0.376 Å⁻¹

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 0.376 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.66	-	-
Author-provided FSC curve	2.67	3.04	2.70
Unmasked-calculated*	3.25	3.86	3.31

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.25 differs from the reported value 2.66 by more than 10 %

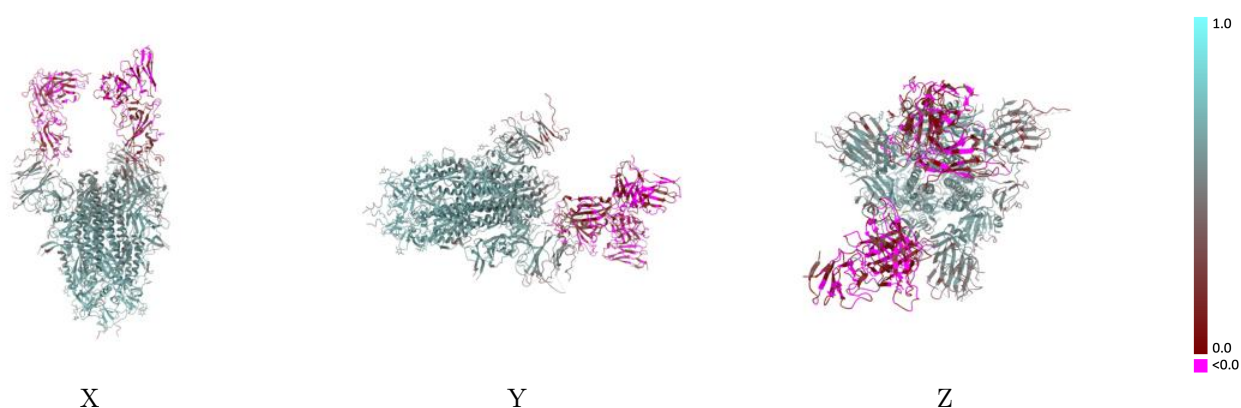
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-63263 and PDB model 9LOY. Per-residue inclusion information can be found in section 3 on page 12.

9.1 Map-model overlay [i](#)

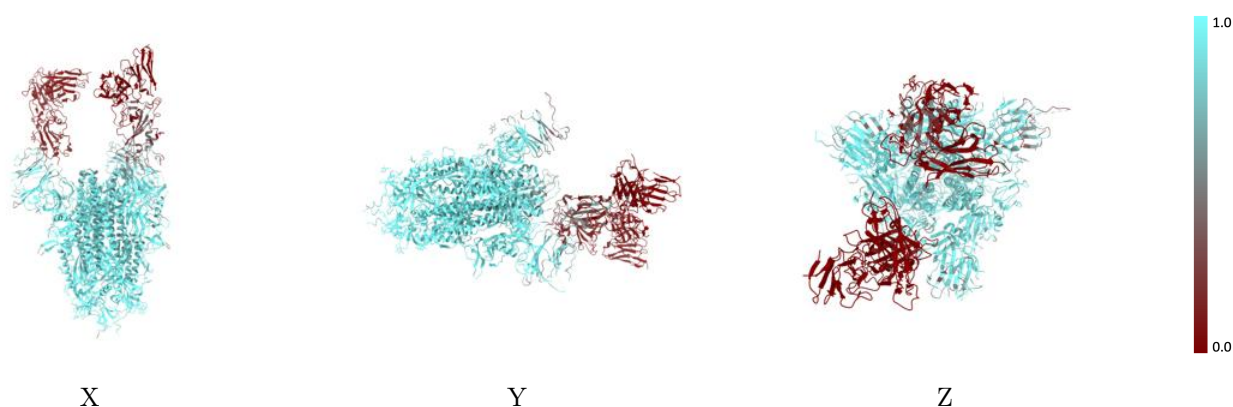
This section was not generated.

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



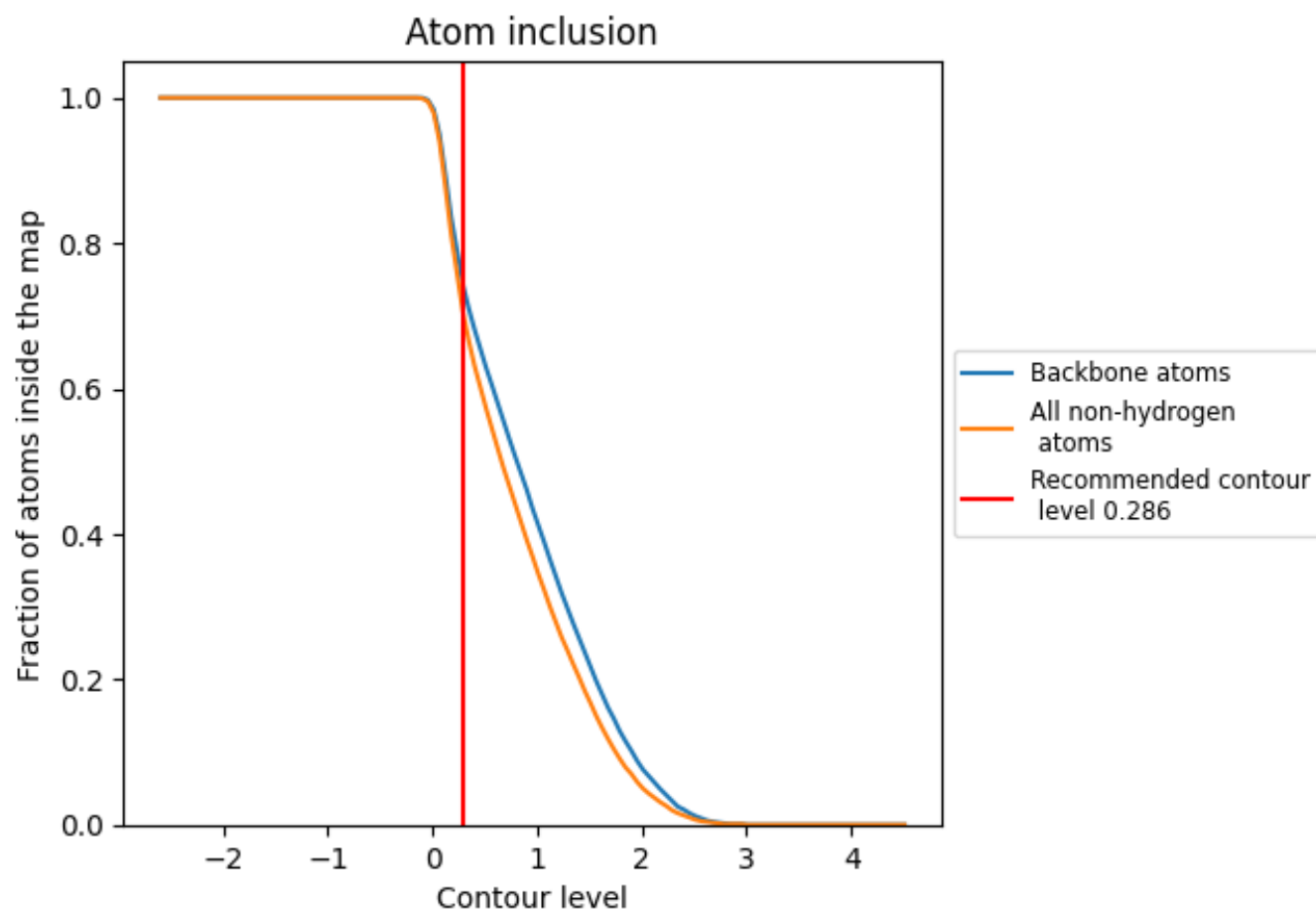
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.286).

9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 71% 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 (0.286) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7080	0.4550
A	0.7740	0.4940
B	0.9120	0.5820
C	0.7610	0.4890
E	0.0030	0.0370
F	0.0030	0.0300
H	0.0290	0.0360
I	0.0430	0.0290

1.0

0.0

<0.0