



Full wwPDB EM Validation Report ⓘ

Nov 21, 2022 – 05:26 AM EST

PDB ID : 7U0P
EMDB ID : EMD-26262
Title : SARS-Cov2 S protein structure in complex with neutralizing monoclonal antibody 002-S21F2
Authors : Patel, A.; Ortlund, E.
Deposited on : 2022-02-18
Resolution : 3.76 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.dev43
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

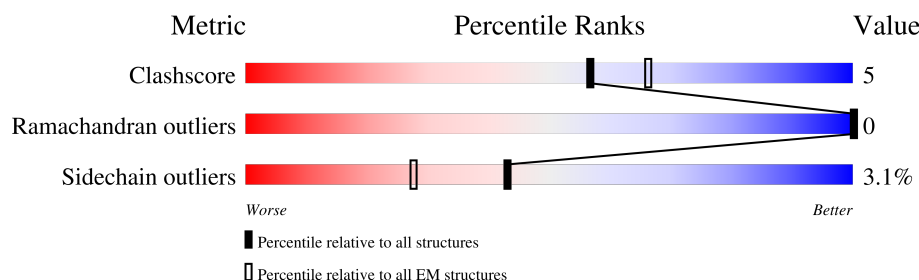
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.76 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	1208	<div> <div>22%</div> <div>75%</div> <div>12%</div> <div>12%</div> </div>
1	B	1208	<div> <div>13%</div> <div>77%</div> <div>12%</div> <div>10%</div> </div>
1	C	1208	<div> <div>10%</div> <div>74%</div> <div>13%</div> <div>12%</div> </div>
2	D	449	<div> <div>27%</div> <div>41%</div> <div>8%</div> <div>50%</div> </div>
2	E	449	<div> <div>30%</div> <div>43%</div> <div>7%</div> <div>50%</div> </div>
3	F	215	<div> <div>63%</div> <div>83%</div> <div>16%</div> <div>•</div> </div>
3	I	215	<div> <div>54%</div> <div>80%</div> <div>20%</div> </div>
4	B1	2	<div> <div>50%</div> <div>100%</div> </div>

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Mol	Chain	Length	Quality of chain
4	G	2	<div><div>50%</div><div><div></div><div></div><div></div></div><div>50%</div><div>50%</div></div>
4	H	2	<div><div>100%</div><div><div></div></div></div>
4	J	2	<div><div>100%</div><div><div></div></div></div>
4	K	2	<div><div>50%</div><div><div></div><div></div><div></div></div><div>100%</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 32371 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	1069	Total	C	N	O	S	0	0
			8366	5339	1397	1591	39		
1	B	1088	Total	C	N	O	S	0	0
			8515	5436	1421	1620	38		
1	C	1067	Total	C	N	O	S	0	0
			8350	5327	1395	1589	39		

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	682	GLY	ARG	conflict	UNP P0DTC2
A	683	SER	ARG	conflict	UNP P0DTC2
A	685	SER	ARG	conflict	UNP P0DTC2
A	817	PRO	PHE	conflict	UNP P0DTC2
A	892	PRO	ALA	conflict	UNP P0DTC2
A	899	PRO	ALA	conflict	UNP P0DTC2
A	942	PRO	ALA	conflict	UNP P0DTC2
A	986	PRO	LYS	conflict	UNP P0DTC2
A	987	PRO	VAL	conflict	UNP P0DTC2
B	682	GLY	ARG	conflict	UNP P0DTC2
B	683	SER	ARG	conflict	UNP P0DTC2
B	685	SER	ARG	conflict	UNP P0DTC2
B	817	PRO	PHE	conflict	UNP P0DTC2
B	892	PRO	ALA	conflict	UNP P0DTC2
B	899	PRO	ALA	conflict	UNP P0DTC2
B	942	PRO	ALA	conflict	UNP P0DTC2
B	986	PRO	LYS	conflict	UNP P0DTC2
B	987	PRO	VAL	conflict	UNP P0DTC2
C	682	GLY	ARG	conflict	UNP P0DTC2
C	683	SER	ARG	conflict	UNP P0DTC2
C	685	SER	ARG	conflict	UNP P0DTC2
C	817	PRO	PHE	conflict	UNP P0DTC2
C	892	PRO	ALA	conflict	UNP P0DTC2
C	899	PRO	ALA	conflict	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	942	PRO	ALA	conflict	UNP P0DTC2
C	986	PRO	LYS	conflict	UNP P0DTC2
C	987	PRO	VAL	conflict	UNP P0DTC2

- Molecule 2 is a protein called mAb 002-S21F2 Heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	224	Total	C	N	O	S	0	0
			1684	1069	275	331	9		
2	D	224	Total	C	N	O	S	0	0
			1684	1069	275	331	9		

- Molecule 3 is a protein called mAb 002_S21F2 light chain.

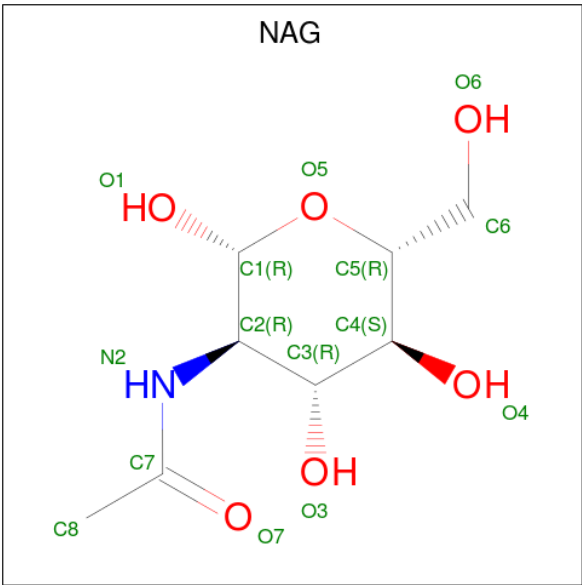
Mol	Chain	Residues	Atoms					AltConf	Trace
3	I	215	Total	C	N	O	S	0	0
			1662	1040	275	341	6		
3	F	215	Total	C	N	O	S	0	0
			1662	1040	275	341	6		

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



Mol	Chain	Residues	Atoms				AltConf	Trace
4	B1	2	Total	C	N	O	0	0
			28	16	2	10		
4	G	2	Total	C	N	O	0	0
			28	16	2	10		
4	H	2	Total	C	N	O	0	0
			28	16	2	10		
4	J	2	Total	C	N	O	0	0
			28	16	2	10		
4	K	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			112	64	8	40	
5	A	1	Total	C	N	O	0
			112	64	8	40	
5	A	1	Total	C	N	O	0
			112	64	8	40	
5	A	1	Total	C	N	O	0
			112	64	8	40	
5	A	1	Total	C	N	O	0
			112	64	8	40	
5	A	1	Total	C	N	O	0
			112	64	8	40	
5	A	1	Total	C	N	O	0
			112	64	8	40	
5	B	1	Total	C	N	O	0
			84	48	6	30	
5	B	1	Total	C	N	O	0
			84	48	6	30	
5	B	1	Total	C	N	O	0
			84	48	6	30	
5	B	1	Total	C	N	O	0
			84	48	6	30	
5	B	1	Total	C	N	O	0
			84	48	6	30	
5	B	1	Total	C	N	O	0
			84	48	6	30	

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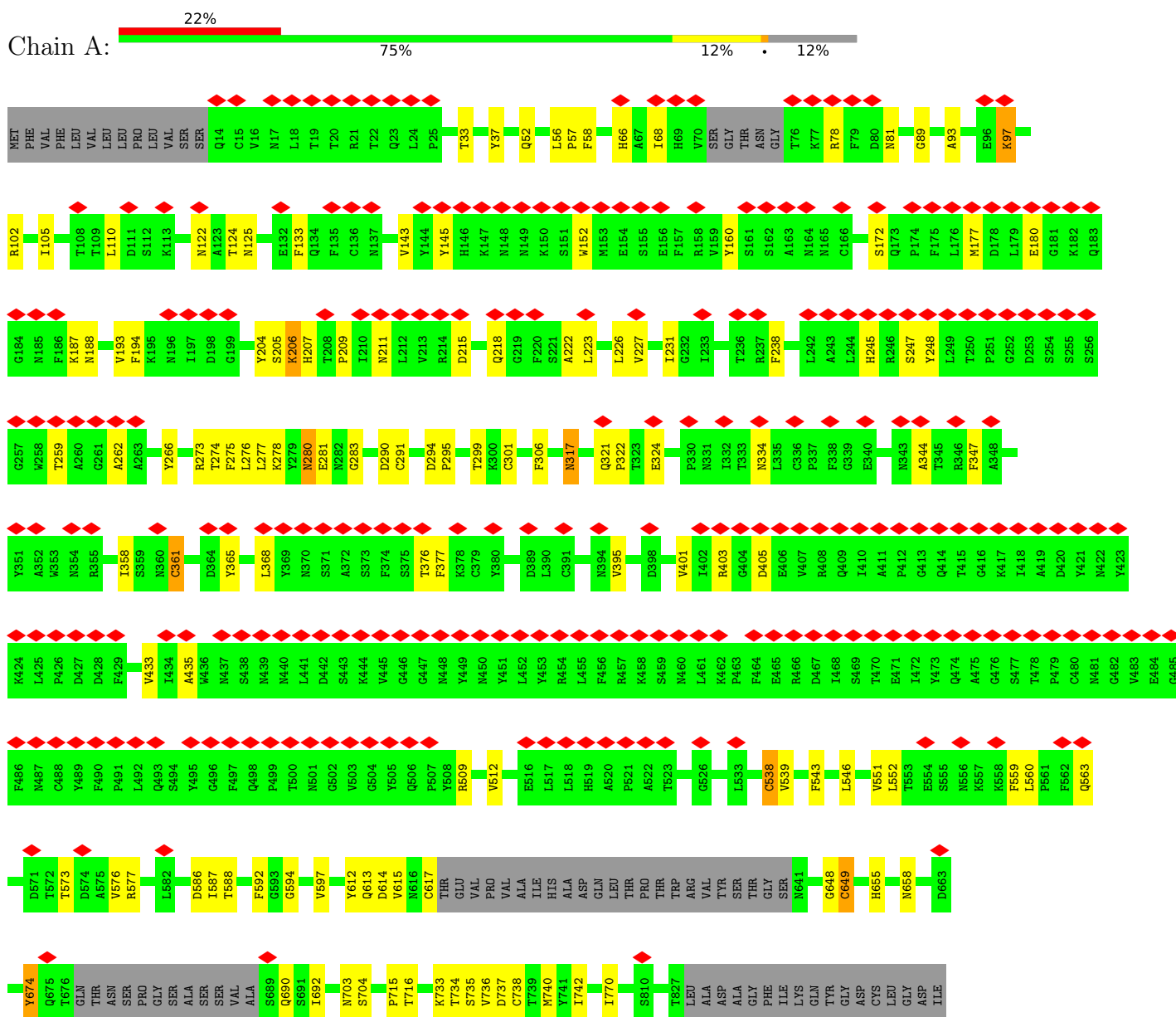
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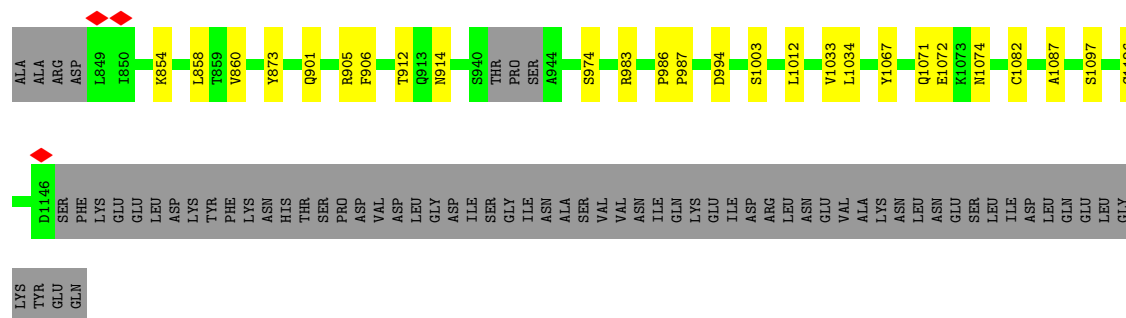
Mol	Chain	Residues	Atoms				AltConf
5	C	1	Total	C	N	O	0
			112	64	8	40	
5	C	1	Total	C	N	O	0
			112	64	8	40	
5	C	1	Total	C	N	O	0
			112	64	8	40	
5	C	1	Total	C	N	O	0
			112	64	8	40	
5	C	1	Total	C	N	O	0
			112	64	8	40	
5	C	1	Total	C	N	O	0
			112	64	8	40	
5	C	1	Total	C	N	O	0
			112	64	8	40	

3 Residue-property plots

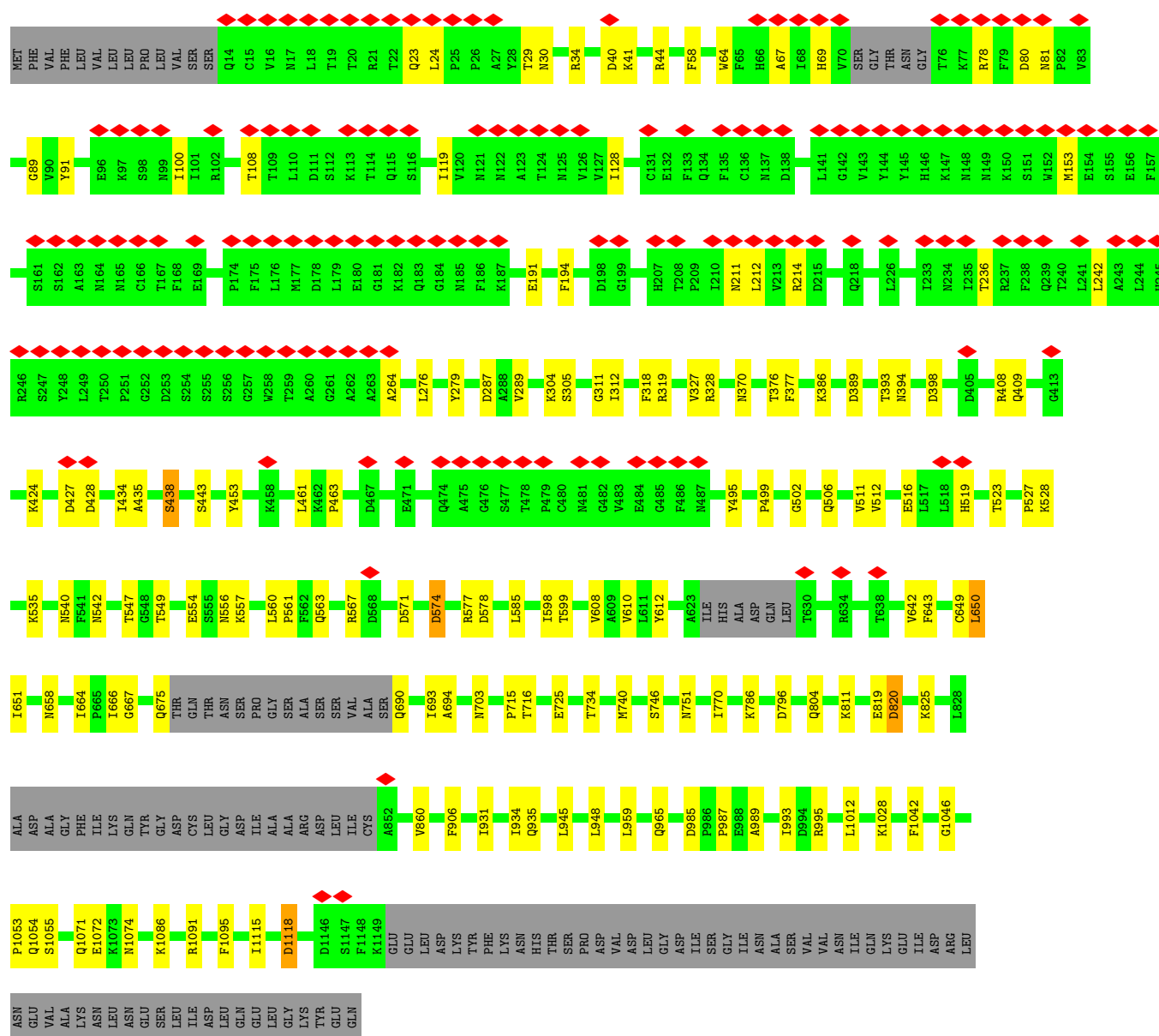
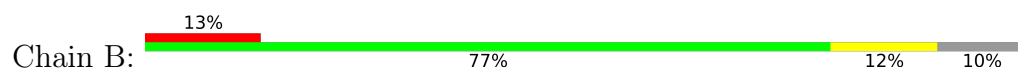
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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

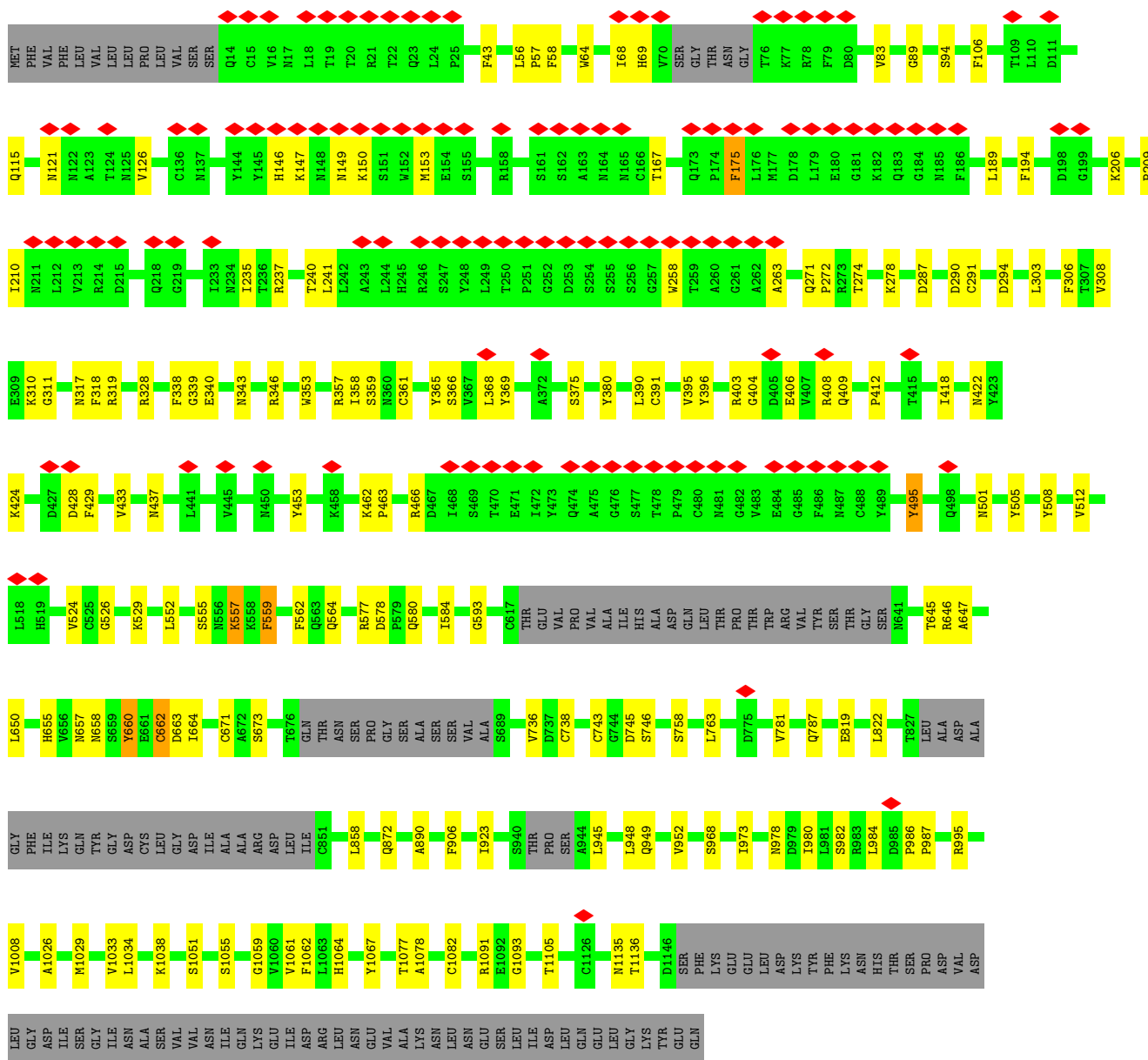
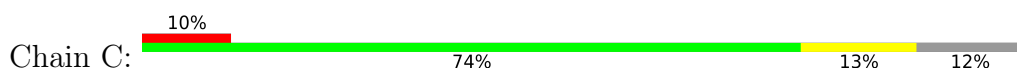




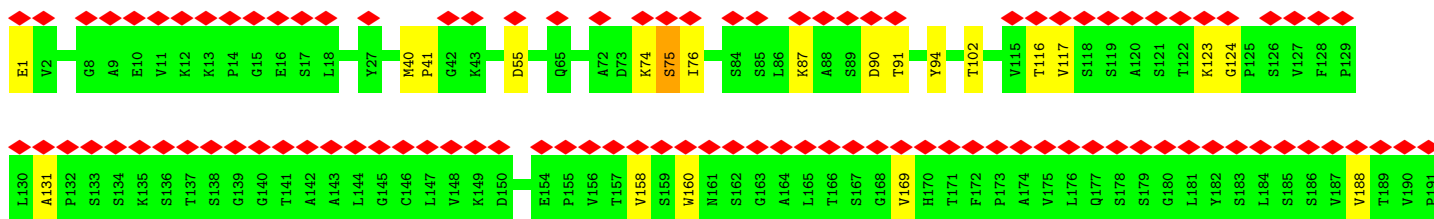
• Molecule 1: Spike glycoprotein

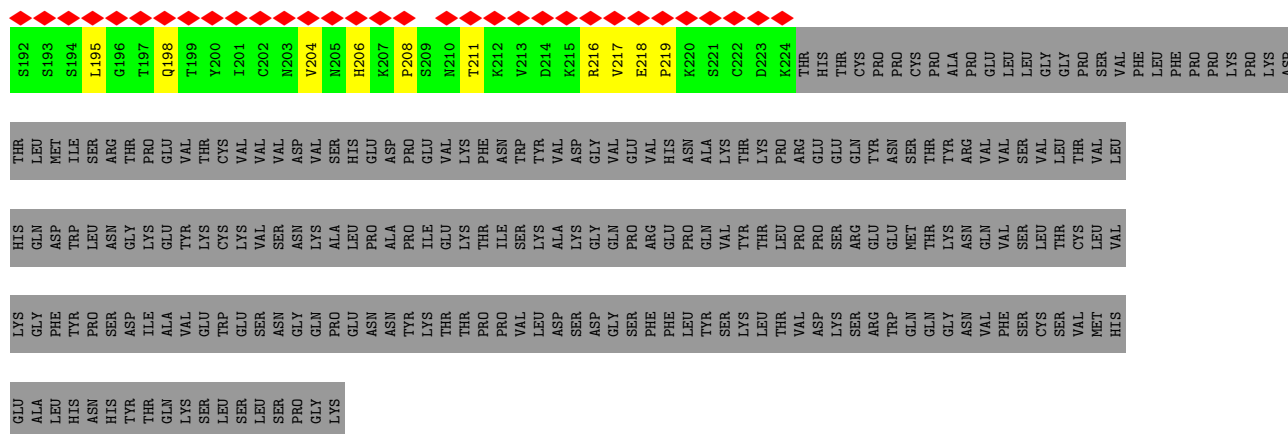


• Molecule 1: Spike glycoprotein

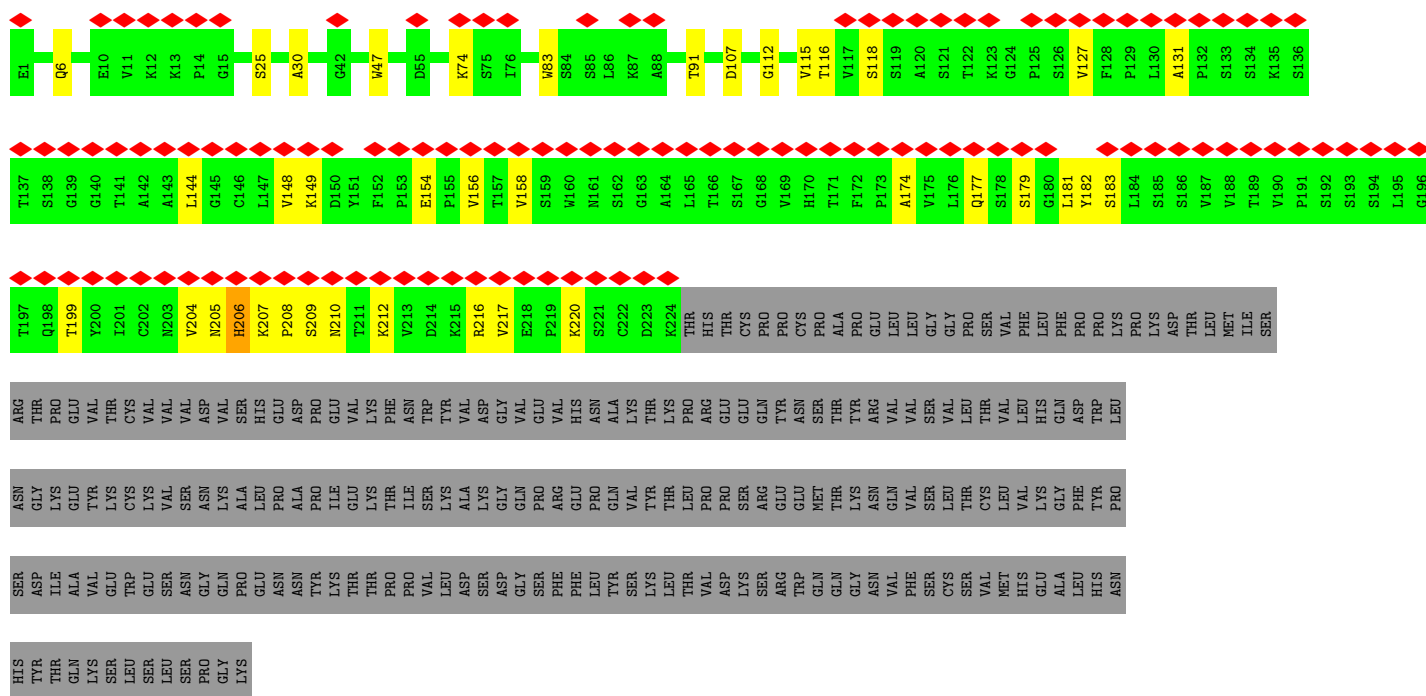
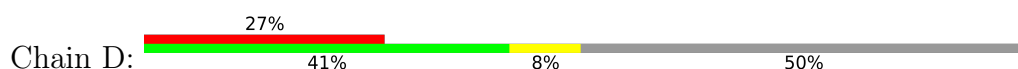


• Molecule 2: mAb 002-S21F2 Heavy chain

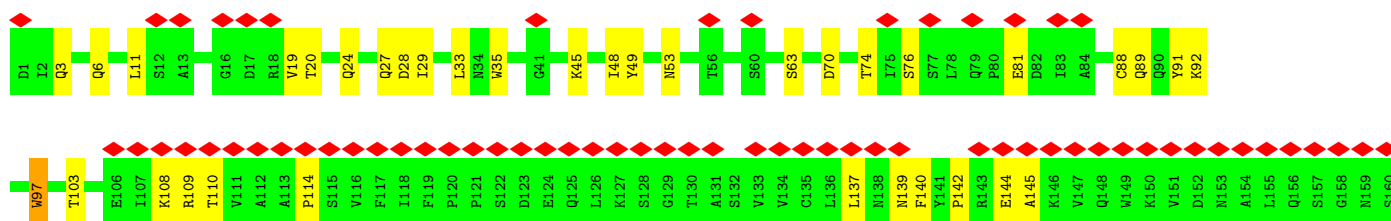
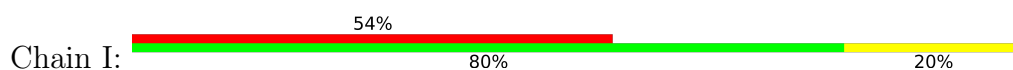


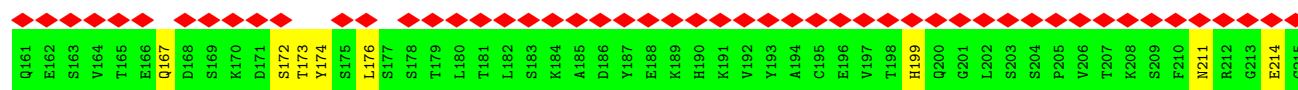


• Molecule 2: mAb 002-S21F2 Heavy chain

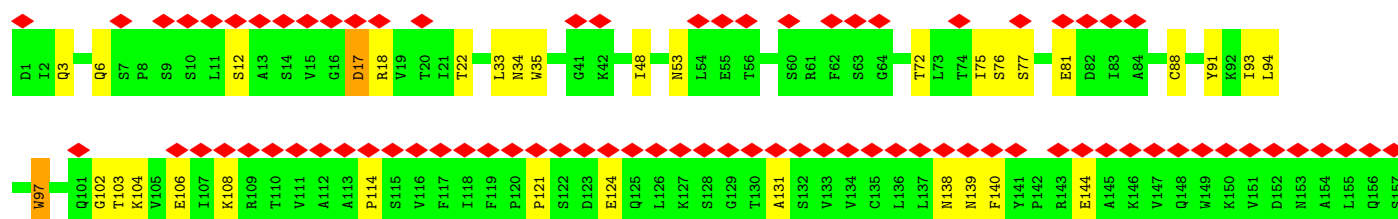
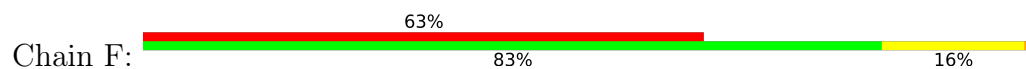


• Molecule 3: mAb 002_S21F2 light chain





- Molecule 3: mAb 002_S21F2 light chain



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



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



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



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



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



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	129931	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	51	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	79000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	4.865	Depositor
Minimum map value	-2.416	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.089	Depositor
Recommended contour level	0.66	Depositor
Map size (Å)	440.0, 440.0, 440.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	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	0.25	0/8562	0.47	0/11654
1	B	0.25	0/8719	0.48	1/11874 (0.0%)
1	C	0.30	2/8546 (0.0%)	0.54	4/11632 (0.0%)
2	D	0.27	0/1728	0.52	0/2352
2	E	0.26	0/1728	0.49	0/2352
3	F	0.27	0/1696	0.49	0/2303
3	I	0.26	0/1696	0.49	0/2303
All	All	0.27	2/32675 (0.0%)	0.50	5/44470 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	209	PRO	CB-CG	-11.22	0.93	1.50
1	C	209	PRO	CG-CD	-9.44	1.19	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	209	PRO	N-CD-CG	-17.52	76.92	103.20
1	C	209	PRO	CA-CB-CG	-15.57	74.42	104.00
1	C	209	PRO	CB-CG-CD	12.01	153.36	106.50
1	C	209	PRO	CA-N-CD	-6.88	101.86	111.50
1	B	1118	ASP	CB-CG-OD1	5.44	123.20	118.30

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	8366	0	8171	81	0
1	B	8515	0	8324	79	0
1	C	8350	0	8145	93	0
2	D	1684	0	1649	24	0
2	E	1684	0	1649	19	0
3	F	1662	0	1617	21	0
3	I	1662	0	1617	23	0
4	B1	28	0	25	0	0
4	G	28	0	25	0	0
4	H	28	0	25	0	0
4	J	28	0	25	0	0
4	K	28	0	25	0	0
5	A	112	0	104	0	0
5	B	84	0	78	0	0
5	C	112	0	104	0	0
All	All	32371	0	31583	326	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 5.

All (326) 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:560:LEU:HD12	1:B:561:PRO:HD2	1.60	0.83
1:A:674:TYR:HB2	1:A:692:ILE:HD13	1.67	0.76
2:D:131:ALA:H	2:D:220:LYS:HE3	1.53	0.74
1:A:89:GLY:HA2	1:A:194:PHE:O	1.90	0.71
1:A:93:ALA:HB3	1:A:266:TYR:HB2	1.73	0.69
1:B:212:LEU:HD23	1:B:214:ARG:H	1.56	0.69
2:D:177:GLN:HE21	2:D:181:LEU:HB2	1.56	0.69
1:C:564:GLN:HE22	1:C:577:ARG:HH21	1.42	0.68
3:F:3:GLN:OE1	3:F:3:GLN:N	2.26	0.68
1:A:740:MET:SD	1:C:319:ARG:NH2	2.67	0.68
3:F:144:GLU:N	3:F:144:GLU:OE2	2.27	0.67
1:A:983:ARG:HG2	1:C:390:LEU:HD11	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:LYS:HA	1:A:211:ASN:HB3	1.77	0.66
1:C:271:GLN:OE1	1:C:272:PRO:HD2	1.94	0.66
1:C:353:TRP:O	1:C:466:ARG:NH1	2.29	0.66
2:E:158:VAL:HG12	2:E:204:VAL:HG22	1.79	0.65
3:F:22:THR:HG22	3:F:72:THR:HG22	1.77	0.65
1:B:547:THR:OG1	1:C:978:ASN:ND2	2.30	0.65
2:D:154:GLU:N	2:D:154:GLU:OE1	2.30	0.65
1:B:394:ASN:ND2	1:B:516:GLU:OE1	2.30	0.65
1:B:675:GLN:O	1:B:690:GLN:N	2.30	0.64
3:I:63:SER:OG	3:I:74:THR:OG1	2.13	0.64
1:C:1093:GLY:HA3	1:C:1105:THR:O	1.97	0.64
2:D:149:LYS:HB3	2:D:183:SER:HA	1.78	0.64
1:B:386:LYS:NZ	1:C:982:SER:O	2.31	0.64
1:C:973:ILE:HB	1:C:980:ILE:HD11	1.81	0.63
1:B:563:GLN:O	1:B:577:ARG:NH2	2.32	0.63
1:C:1135:ASN:OD1	1:C:1136:THR:N	2.32	0.62
1:B:703:ASN:ND2	1:C:787:GLN:OE1	2.33	0.62
3:F:6:GLN:OE1	3:F:102:GLY:N	2.27	0.61
3:I:139:ASN:HD22	3:I:173:THR:HG21	1.65	0.61
2:D:177:GLN:NE2	2:D:179:SER:OG	2.34	0.61
1:A:546:LEU:HD21	1:A:573:THR:HG21	1.81	0.60
1:C:328:ARG:NH2	1:C:580:GLN:OE1	2.34	0.60
3:F:124:GLU:OE1	3:F:124:GLU:N	2.24	0.60
1:A:66:HIS:O	1:A:78:ARG:NH1	2.35	0.60
1:A:612:TYR:O	1:A:648:GLY:HA3	2.02	0.59
1:C:290:ASP:OD1	1:C:291:CYS:N	2.35	0.59
1:A:401:VAL:HG22	1:A:509:ARG:HG2	1.84	0.59
1:A:737:ASP:OD2	1:C:317:ASN:ND2	2.35	0.59
1:B:770:ILE:HD11	1:B:1012:LEU:HA	1.84	0.59
1:B:69:HIS:HB3	1:B:78:ARG:HB2	1.83	0.59
1:B:89:GLY:HA2	1:B:194:PHE:O	2.02	0.59
1:B:574:ASP:OD2	1:B:574:ASP:N	2.35	0.59
1:A:324:GLU:H	1:A:539:VAL:HG12	1.68	0.59
1:A:152:TRP:NE1	1:A:180:GLU:OE1	2.36	0.58
1:B:535:LYS:NZ	1:B:554:GLU:OE2	2.35	0.58
1:A:358:ILE:HB	1:A:395:VAL:HB	1.86	0.58
2:D:206:HIS:CE1	2:D:209:SER:HB2	2.38	0.58
1:B:67:ALA:HA	1:B:80:ASP:HB3	1.85	0.58
1:B:666:ILE:HG22	1:B:667:GLY:H	1.69	0.58
1:A:281:GLU:OE1	1:A:281:GLU:N	2.27	0.58
1:A:290:ASP:OD1	1:A:291:CYS:N	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:391:CYS:SG	1:C:526:GLY:N	2.77	0.57
1:A:716:THR:OG1	1:A:1071:GLN:O	2.23	0.57
3:I:144:GLU:OE1	3:I:144:GLU:N	2.28	0.57
1:A:122:ASN:ND2	1:A:124:THR:OG1	2.37	0.57
1:A:736:VAL:HG22	1:A:858:LEU:HD22	1.86	0.57
2:E:123:LYS:NZ	2:E:124:GLY:O	2.36	0.57
1:B:811:LYS:NZ	1:B:820:ASP:OD2	2.36	0.57
1:B:610:VAL:HG12	1:B:651:ILE:HB	1.87	0.57
3:F:91:TYR:HB2	3:F:97:TRP:CE3	2.40	0.57
1:C:406:GLU:OE2	1:C:409:GLN:HG3	2.04	0.56
1:C:645:THR:HG22	1:C:647:ALA:H	1.70	0.56
1:A:715:PRO:HA	1:A:1072:GLU:HA	1.87	0.56
3:F:12:SER:HB2	3:F:106:GLU:HB2	1.86	0.56
1:C:424:LYS:HB3	1:C:463:PRO:HA	1.87	0.56
1:A:33:THR:OG1	1:A:218:GLN:O	2.22	0.56
1:B:23:GLN:NE2	1:B:80:ASP:OD1	2.39	0.55
1:B:393:THR:O	1:B:523:THR:OG1	2.21	0.55
3:F:114:PRO:HB3	3:F:140:PHE:HB3	1.87	0.55
1:B:643:PHE:HB3	1:B:650:LEU:HG	1.88	0.55
1:B:398:ASP:HB2	1:B:512:VAL:HG22	1.87	0.55
1:B:725:GLU:OE1	1:B:1028:LYS:NZ	2.40	0.55
2:E:91:THR:HG22	2:E:116:THR:HA	1.89	0.55
3:I:3:GLN:N	3:I:3:GLN:OE1	2.40	0.55
1:A:52:GLN:HB2	1:A:274:THR:HG22	1.87	0.55
1:B:804:GLN:NE2	1:B:935:GLN:OE1	2.37	0.55
2:E:195:LEU:HD13	2:E:219:PRO:HG3	1.89	0.55
1:A:901:GLN:HE21	1:A:905:ARG:HE	1.55	0.55
1:C:357:ARG:HG2	1:C:396:TYR:HE1	1.71	0.54
1:A:273:ARG:HB2	1:A:275:PHE:HE2	1.73	0.54
1:C:189:LEU:HB2	1:C:210:ILE:HD11	1.88	0.54
1:A:538:CYS:HB2	1:A:551:VAL:HG12	1.89	0.54
1:A:125:ASN:ND2	1:A:172:SER:O	2.38	0.54
1:A:334:ASN:HB2	1:A:361:CYS:HA	1.89	0.54
1:C:346:ARG:NH1	2:E:102:THR:O	2.41	0.54
1:C:763:LEU:HD22	1:C:1008:VAL:HG21	1.89	0.54
1:A:299:THR:HG21	1:A:597:VAL:HG21	1.90	0.54
1:C:358:ILE:HB	1:C:395:VAL:HB	1.89	0.54
1:A:280:ASN:OD1	1:A:283:GLY:N	2.41	0.53
1:C:555:SER:OG	1:C:557:LYS:NZ	2.40	0.53
1:C:340:GLU:N	1:C:340:GLU:OE1	2.41	0.53
1:C:428:ASP:OD1	1:C:428:ASP:N	2.40	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:114:PRO:HB3	3:I:140:PHE:HB3	1.91	0.53
1:C:404:GLY:HA2	1:C:508:TYR:CD2	2.43	0.52
2:D:30:ALA:HB2	2:D:74:LYS:HD2	1.92	0.52
2:D:156:VAL:HG23	2:D:206:HIS:HB3	1.91	0.52
3:I:137:LEU:HB2	3:I:176:LEU:HB3	1.91	0.52
1:B:434:ILE:HB	1:B:511:VAL:HG12	1.90	0.52
2:E:206:HIS:HB3	2:E:211:THR:HG22	1.90	0.52
2:D:199:THR:OG1	2:D:216:ARG:NE	2.41	0.52
1:B:24:LEU:HD11	1:B:78:ARG:HD3	1.91	0.52
1:A:317:ASN:HA	1:A:594:GLY:HA2	1.90	0.52
1:B:438:SER:O	1:B:438:SER:OG	2.28	0.51
1:C:359:SER:HA	1:C:524:VAL:HG23	1.93	0.51
1:A:551:VAL:HG22	1:A:588:THR:HB	1.93	0.51
2:D:174:ALA:HB1	2:D:182:TYR:HB3	1.92	0.51
1:C:819:GLU:OE2	1:C:1055:SER:OG	2.27	0.51
1:A:295:PRO:O	1:A:299:THR:HG23	2.11	0.51
1:C:89:GLY:HA2	1:C:194:PHE:O	2.11	0.50
1:C:106:PHE:HD2	1:C:235:ILE:HG21	1.77	0.50
1:C:736:VAL:HG22	1:C:858:LEU:HD23	1.94	0.50
2:D:47:TRP:HE1	3:I:89:GLN:NE2	2.08	0.50
1:B:312:ILE:HG22	1:B:598:ILE:HG12	1.94	0.50
1:B:327:VAL:HG12	1:B:542:ASN:HB3	1.93	0.50
3:I:81:GLU:OE1	3:I:81:GLU:N	2.37	0.50
2:D:6:GLN:OE1	2:D:112:GLY:N	2.35	0.50
1:C:83:VAL:HG21	1:C:237:ARG:HH21	1.76	0.50
1:C:68:ILE:HG12	1:C:263:ALA:H	1.76	0.50
1:B:108:THR:HA	1:B:236:THR:HG22	1.93	0.50
1:B:931:ILE:O	1:B:934:ILE:HG22	2.12	0.50
1:C:115:GLN:NE2	1:C:167:THR:OG1	2.42	0.50
2:E:91:THR:HG22	2:E:117:VAL:H	1.76	0.50
1:A:365:TYR:O	1:A:368:LEU:HB2	2.12	0.49
1:B:91:TYR:OH	1:B:191:GLU:OE2	2.30	0.49
2:D:144:LEU:HD21	2:D:217:VAL:HG11	1.94	0.49
3:I:20:THR:HG23	3:I:74:THR:HG22	1.93	0.49
1:B:100:ILE:HG22	1:B:242:LEU:HD12	1.94	0.49
1:C:294:ASP:OD1	1:C:294:ASP:N	2.45	0.49
1:C:949:GLN:O	1:C:952:VAL:HG22	2.12	0.49
1:A:703:ASN:OD1	1:A:704:SER:N	2.46	0.49
1:A:563:GLN:O	1:A:577:ARG:NE	2.38	0.49
1:B:64:TRP:HE1	1:B:264:ALA:HB1	1.77	0.49
1:B:1091:ARG:NH1	1:B:1118:ASP:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1053:PRO:O	1:B:1054:GLN:NE2	2.42	0.49
1:A:247:SER:HB2	1:A:259:THR:HG23	1.94	0.49
1:B:734:THR:HG22	1:B:860:VAL:HG12	1.95	0.49
3:I:167:GLN:OE1	3:I:174:TYR:OH	2.21	0.49
2:D:207:LYS:N	2:D:208:PRO:HD2	2.27	0.48
3:I:145:ALA:HB2	3:I:199:HIS:HD2	1.78	0.48
1:A:278:LYS:HD2	1:A:306:PHE:HE2	1.79	0.48
1:B:540:ASN:HB3	1:B:549:THR:HG22	1.95	0.48
1:B:945:LEU:HD12	1:B:948:LEU:HD12	1.95	0.48
3:F:35:TRP:CZ3	3:F:88:CYS:HB3	2.49	0.48
1:A:205:SER:HB3	1:A:226:LEU:HD22	1.94	0.48
1:B:599:THR:HB	1:B:608:VAL:HG22	1.94	0.48
1:C:126:VAL:HG11	1:C:175:PHE:HE2	1.77	0.48
1:C:303:LEU:HD12	1:C:308:VAL:HG12	1.95	0.48
1:C:403:ARG:HB3	1:C:495:TYR:HE1	1.79	0.48
3:F:33:LEU:O	3:F:34:ASN:ND2	2.44	0.48
1:C:1077:THR:OG1	1:C:1078:ALA:N	2.46	0.48
1:C:658:ASN:ND2	1:C:660:TYR:OH	2.47	0.48
1:B:693:ILE:HD12	1:B:694:ALA:H	1.79	0.48
3:F:138:ASN:OD1	3:F:139:ASN:HB2	2.14	0.48
1:C:559:PHE:HB2	1:C:584:ILE:HG13	1.95	0.48
1:A:226:LEU:HG	1:A:227:VAL:HG23	1.95	0.47
2:D:127:VAL:HA	2:D:148:VAL:HG22	1.96	0.47
1:A:344:ALA:HB3	1:A:347:PHE:HE1	1.79	0.47
1:B:642:VAL:HG23	1:B:651:ILE:HD11	1.95	0.47
1:B:825:LYS:HD2	1:B:825:LYS:HA	1.58	0.47
1:C:346:ARG:NH2	3:F:91:TYR:OH	2.47	0.47
1:C:1033:VAL:HG12	1:C:1034:LEU:HD23	1.95	0.47
3:I:142:PRO:HG2	3:I:144:GLU:OE2	2.14	0.47
2:E:91:THR:HG23	2:E:117:VAL:HG12	1.97	0.47
3:I:211:ASN:HB3	3:I:214:GLU:HB2	1.96	0.47
1:A:206:LYS:NZ	1:A:222:ALA:O	2.41	0.47
1:B:304:LYS:HE2	1:B:304:LYS:HB3	1.69	0.47
1:A:145:TYR:O	1:A:248:TYR:OH	2.27	0.47
3:I:24:GLN:NE2	3:I:70:ASP:OD1	2.47	0.47
2:D:158:VAL:HG22	2:D:204:VAL:HG22	1.96	0.47
1:A:89:GLY:CA	1:A:194:PHE:O	2.62	0.46
1:A:614:ASP:OD1	1:A:614:ASP:N	2.44	0.46
3:I:109:ARG:HE	3:I:172:SER:HB2	1.80	0.46
1:C:1051:SER:OG	1:C:1064:HIS:ND1	2.26	0.46
1:A:994:ASP:OD2	1:C:995:ARG:NH2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:547:THR:O	1:C:978:ASN:ND2	2.48	0.46
1:C:146:HIS:O	1:C:150:LYS:N	2.48	0.46
2:E:206:HIS:CD2	2:E:208:PRO:HD2	2.50	0.46
1:C:278:LYS:NZ	1:C:287:ASP:OD2	2.37	0.46
1:C:146:HIS:HB3	1:C:149:ASN:HB2	1.97	0.46
2:E:75:SER:OG	2:E:76:ILE:HD12	2.16	0.46
1:C:662:CYS:HB2	1:C:671:CYS:HB2	1.47	0.46
3:I:91:TYR:HB2	3:I:97:TRP:CE3	2.50	0.46
1:A:321:GLN:HG2	1:A:322:PRO:HD2	1.98	0.46
1:C:339:GLY:O	1:C:343:ASN:HB2	2.16	0.46
2:E:131:ALA:HB2	2:E:217:VAL:HG12	1.98	0.45
1:B:376:THR:HB	1:B:435:ALA:HB3	1.99	0.45
1:C:56:LEU:HD12	1:C:57:PRO:HD2	1.97	0.45
1:C:462:LYS:HE2	1:C:462:LYS:HA	1.98	0.45
1:B:985:ASP:HB2	1:B:987:PRO:HD2	1.98	0.45
1:A:206:LYS:HB2	1:A:223:LEU:HG	1.98	0.45
1:A:294:ASP:OD1	1:A:294:ASP:N	2.49	0.45
1:B:443:SER:HB3	1:B:499:PRO:HG3	1.99	0.45
1:C:406:GLU:OE1	1:C:418:ILE:HG12	2.17	0.45
1:A:552:LEU:HD13	1:A:587:ILE:HG12	1.98	0.45
1:A:1074:ASN:OD1	1:A:1074:ASN:N	2.44	0.45
1:B:40:ASP:OD1	1:B:41:LYS:N	2.46	0.45
1:A:81:ASN:OD1	1:A:81:ASN:N	2.49	0.44
1:B:612:TYR:HB2	1:B:649:CYS:SG	2.57	0.44
1:C:562:PHE:O	1:C:564:GLN:NE2	2.48	0.44
1:A:543:PHE:CD2	1:A:576:VAL:HG11	2.51	0.44
1:B:81:ASN:N	1:B:81:ASN:OD1	2.49	0.44
1:B:819:GLU:OE2	1:B:1055:SER:OG	2.30	0.44
1:A:433:VAL:HG22	1:A:512:VAL:HG22	1.99	0.44
1:B:989:ALA:O	1:B:993:ILE:HG12	2.18	0.44
1:A:560:LEU:HB2	1:A:563:GLN:HB2	1.99	0.44
3:I:6:GLN:NE2	3:I:103:THR:OG1	2.50	0.44
1:B:319:ARG:NH1	1:C:745:ASP:OD1	2.49	0.44
1:A:617:CYS:HB2	1:A:649:CYS:HB2	1.43	0.44
1:A:854:LYS:HA	1:A:854:LYS:HD3	1.67	0.44
1:B:995:ARG:H	1:B:995:ARG:HG2	1.63	0.44
1:C:318:PHE:HB3	1:C:593:GLY:HA3	2.00	0.44
3:I:35:TRP:CZ3	3:I:88:CYS:HB3	2.53	0.44
1:B:428:ASP:OD1	1:B:428:ASP:N	2.51	0.44
1:C:365:TYR:HD1	1:C:368:LEU:HD12	1.82	0.44
1:A:613:GLN:O	1:A:615:VAL:N	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:240:THR:OG1	1:C:241:LEU:N	2.51	0.43
2:D:206:HIS:O	2:D:206:HIS:ND1	2.49	0.43
1:A:1087:ALA:HB2	1:A:1126:CYS:HB3	2.00	0.43
1:B:516:GLU:OE2	1:B:519:HIS:ND1	2.51	0.43
2:E:169:VAL:HG22	2:E:188:VAL:HG12	2.01	0.43
1:A:193:VAL:HB	1:A:204:TYR:HB2	2.00	0.43
1:A:674:TYR:O	1:A:674:TYR:CG	2.71	0.43
1:B:716:THR:HG22	1:B:1071:GLN:O	2.18	0.43
2:E:74:LYS:HD3	2:E:74:LYS:C	2.39	0.43
2:D:205:ASN:HB2	2:D:212:LYS:HE2	1.99	0.43
3:I:28:ASP:OD2	3:I:29:ILE:N	2.51	0.43
1:C:380:TYR:HE2	1:C:412:PRO:HD2	1.83	0.43
1:C:948:LEU:HD21	1:C:1059:GLY:HA3	2.01	0.43
1:C:986:PRO:N	1:C:987:PRO:HD2	2.33	0.43
3:F:48:ILE:HG23	3:F:53:ASN:O	2.18	0.43
1:A:68:ILE:HD13	1:A:262:ALA:HB1	2.00	0.43
1:B:119:ILE:HG12	1:B:128:ILE:HG12	2.00	0.43
1:B:311:GLY:HA2	1:B:664:ILE:HD12	2.00	0.43
1:C:437:ASN:HA	1:C:508:TYR:HD1	1.84	0.43
1:A:658:ASN:N	1:A:658:ASN:OD1	2.52	0.43
1:B:29:THR:OG1	1:B:30:ASN:N	2.51	0.43
1:B:1046:GLY:HA2	1:C:890:ALA:HB1	2.01	0.43
3:I:109:ARG:HG2	3:I:110:THR:H	1.82	0.43
3:F:6:GLN:NE2	3:F:103:THR:OG1	2.52	0.43
1:B:1095:PHE:HE1	1:B:1115:ILE:HD12	1.85	0.42
1:A:105:ILE:HD11	1:A:110:LEU:HD13	2.01	0.42
1:C:317:ASN:OD1	1:C:317:ASN:N	2.53	0.42
1:B:276:LEU:HB3	1:B:289:VAL:HG13	2.01	0.42
1:C:274:THR:HG23	1:C:291:CYS:HB2	2.00	0.42
1:C:357:ARG:HG2	1:C:396:TYR:CE1	2.53	0.42
1:A:143:VAL:HG22	1:A:245:HIS:HA	2.01	0.42
1:A:738:CYS:O	1:A:742:ILE:HG12	2.20	0.42
1:A:912:THR:OG1	1:A:914:ASN:OD1	2.28	0.42
1:C:310:LYS:NZ	1:C:663:ASP:OD2	2.37	0.42
1:C:645:THR:HG22	1:C:647:ALA:N	2.33	0.42
1:A:1033:VAL:HG12	1:A:1034:LEU:HD23	2.01	0.42
1:C:328:ARG:HH21	1:C:580:GLN:HB2	1.83	0.42
2:D:206:HIS:HE1	2:D:209:SER:HB2	1.82	0.42
3:I:33:LEU:HD23	3:I:33:LEU:HA	1.86	0.42
1:C:206:LYS:HD2	1:C:206:LYS:HA	1.82	0.42
2:E:40:MET:HG3	2:E:41:PRO:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:VAL:HG21	1:C:237:ARG:NH2	2.35	0.42
1:C:433:VAL:HG23	1:C:512:VAL:HG22	2.01	0.42
1:C:984:LEU:HD23	1:C:984:LEU:HA	1.93	0.42
1:A:188:ASN:HA	1:A:209:PRO:HA	2.02	0.42
1:C:311:GLY:HA2	1:C:664:ILE:HD12	2.01	0.42
1:B:527:PRO:O	1:B:528:LYS:HG2	2.20	0.41
1:C:418:ILE:HG22	1:C:422:ASN:HD21	1.84	0.41
1:C:501:ASN:HB3	1:C:505:TYR:HB2	2.02	0.41
1:C:822:LEU:HD11	1:C:945:LEU:HD21	2.02	0.41
2:E:76:ILE:HD12	2:E:76:ILE:N	2.36	0.41
1:B:377:PHE:HD1	1:B:434:ILE:HG12	1.85	0.41
2:E:87:LYS:HA	2:E:87:LYS:HD3	1.80	0.41
2:D:91:THR:OG1	2:D:116:THR:HA	2.20	0.41
3:F:17:ASP:N	3:F:17:ASP:OD1	2.53	0.41
3:F:18:ARG:HA	3:F:75:ILE:O	2.20	0.41
2:E:216:ARG:HH12	2:E:218:GLU:HB3	1.85	0.41
1:A:37:TYR:HB3	1:A:223:LEU:HB2	2.02	0.41
1:C:763:LEU:HD23	1:C:763:LEU:HA	1.89	0.41
1:A:56:LEU:HD12	1:A:57:PRO:HD2	2.02	0.41
1:A:102:ARG:HD3	1:A:102:ARG:HA	1.78	0.41
1:A:276:LEU:C	1:A:277:LEU:HD23	2.40	0.41
1:A:403:ARG:HE	1:A:405:ASP:HB2	1.85	0.41
1:C:1038:LYS:HE2	1:C:1038:LYS:HB3	1.88	0.41
2:D:91:THR:HA	2:D:115:VAL:O	2.21	0.41
1:A:215:ASP:OD2	1:A:266:TYR:OH	2.39	0.41
1:C:69:HIS:CD2	1:C:258:TRP:HB2	2.56	0.41
1:C:646:ARG:H	1:C:646:ARG:HG3	1.69	0.41
3:F:93:ILE:HG22	3:F:94:LEU:H	1.86	0.41
1:C:147:LYS:H	1:C:147:LYS:HG2	1.64	0.41
2:D:207:LYS:O	2:D:210:ASN:ND2	2.54	0.41
3:I:11:LEU:HD11	3:I:19:VAL:HG13	2.03	0.41
1:A:734:THR:HG22	1:A:860:VAL:HG12	2.03	0.41
1:B:44:ARG:HB2	1:B:279:TYR:CD2	2.56	0.41
1:B:424:LYS:HB3	1:B:463:PRO:HA	2.03	0.41
1:B:424:LYS:HA	1:B:424:LYS:HD2	1.80	0.41
1:B:557:LYS:HZ1	1:C:43:PHE:HE2	1.69	0.41
1:C:781:VAL:HG22	1:C:1026:ALA:HB2	2.02	0.41
1:C:923:ILE:H	1:C:923:ILE:HG12	1.67	0.41
2:D:206:HIS:HD1	2:D:206:HIS:C	2.24	0.41
3:I:48:ILE:HG23	3:I:53:ASN:O	2.21	0.41
3:F:81:GLU:N	3:F:81:GLU:OE1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:ARG:H	1:B:34:ARG:HG2	1.59	0.41
1:B:502:GLY:O	1:B:506:GLN:HG3	2.22	0.41
1:B:959:LEU:HD23	1:B:959:LEU:HA	1.89	0.41
2:E:91:THR:CG2	2:E:117:VAL:H	2.34	0.41
3:F:193:TYR:HB2	3:F:210:PHE:CE1	2.56	0.41
1:A:133:PHE:HB3	1:A:160:TYR:HB3	2.02	0.40
1:B:1074:ASN:OD1	1:B:1074:ASN:N	2.54	0.40
1:B:1086:LYS:H	1:B:1086:LYS:HG3	1.71	0.40
1:A:376:THR:HB	1:A:435:ALA:HB3	2.02	0.40
1:B:715:PRO:HA	1:B:1072:GLU:HA	2.02	0.40
1:B:786:LYS:HA	1:B:786:LYS:HD3	1.79	0.40
1:C:366:SER:HA	1:C:369:TYR:CZ	2.56	0.40
1:A:231:ILE:HD13	1:A:231:ILE:HA	1.88	0.40
1:A:770:ILE:HD11	1:A:1012:LEU:HA	2.03	0.40
1:B:211:ASN:OD1	1:B:211:ASN:N	2.54	0.40
1:B:556:ASN:OD1	1:B:556:ASN:N	2.52	0.40
1:C:328:ARG:NE	1:C:578:ASP:OD2	2.55	0.40
1:C:1029:MET:HB2	1:C:1062:PHE:HZ	1.86	0.40
3:F:76:SER:OG	3:F:77:SER:N	2.53	0.40
3:F:121:PRO:HG2	3:F:131:ALA:HB1	2.02	0.40
1:A:97:LYS:HB2	1:A:177:MET:HE1	2.03	0.40
1:A:986:PRO:HB2	1:A:987:PRO:HD3	2.03	0.40
1:B:328:ARG:NH2	1:B:578:ASP:OD2	2.52	0.40
1:C:403:ARG:HB3	1:C:495:TYR:CE1	2.56	0.40
1:C:822:LEU:HD21	1:C:1061:VAL:HG21	2.03	0.40
2:E:90:ASP:OD1	2:E:94:TYR:OH	2.40	0.40
2:D:149:LYS:HD3	2:D:183:SER:HB2	2.03	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	1057/1208 (88%)	1019 (96%)	38 (4%)	0	100	100
1	B	1078/1208 (89%)	1046 (97%)	32 (3%)	0	100	100
1	C	1055/1208 (87%)	1019 (97%)	36 (3%)	0	100	100
2	D	222/449 (49%)	211 (95%)	11 (5%)	0	100	100
2	E	222/449 (49%)	216 (97%)	6 (3%)	0	100	100
3	F	213/215 (99%)	201 (94%)	12 (6%)	0	100	100
3	I	213/215 (99%)	205 (96%)	8 (4%)	0	100	100
All	All	4060/4952 (82%)	3917 (96%)	143 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	937/1056 (89%)	910 (97%)	27 (3%)	42	67
1	B	954/1056 (90%)	926 (97%)	28 (3%)	42	67
1	C	935/1056 (88%)	901 (96%)	34 (4%)	35	63
2	D	188/396 (48%)	183 (97%)	5 (3%)	44	69
2	E	188/396 (48%)	183 (97%)	5 (3%)	44	69
3	F	191/191 (100%)	187 (98%)	4 (2%)	53	74
3	I	191/191 (100%)	184 (96%)	7 (4%)	34	62
All	All	3584/4342 (82%)	3474 (97%)	110 (3%)	43	65

All (110) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	PHE
1	A	97	LYS
1	A	206	LYS
1	A	207	HIS

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Mol	Chain	Res	Type
1	A	238	PHE
1	A	280	ASN
1	A	301	CYS
1	A	317	ASN
1	A	361	CYS
1	A	377	PHE
1	A	538	CYS
1	A	559	PHE
1	A	586	ASP
1	A	592	PHE
1	A	649	CYS
1	A	655	HIS
1	A	674	TYR
1	A	690	GLN
1	A	733	LYS
1	A	735	SER
1	A	873	TYR
1	A	906	PHE
1	A	974	SER
1	A	1003	SER
1	A	1067	TYR
1	A	1082	CYS
1	A	1097	SER
1	B	58	PHE
1	B	153	MET
1	B	287	ASP
1	B	305	SER
1	B	318	PHE
1	B	370	ASN
1	B	389	ASP
1	B	408	ARG
1	B	409	GLN
1	B	427	ASP
1	B	438	SER
1	B	453	TYR
1	B	461	LEU
1	B	495	TYR
1	B	567	ARG
1	B	571	ASP
1	B	574	ASP
1	B	585	LEU
1	B	650	LEU

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Mol	Chain	Res	Type
1	B	658	ASN
1	B	740	MET
1	B	746	SER
1	B	751	ASN
1	B	796	ASP
1	B	820	ASP
1	B	906	PHE
1	B	965	GLN
1	B	1042	PHE
1	C	58	PHE
1	C	64	TRP
1	C	94	SER
1	C	121	ASN
1	C	153	MET
1	C	175	PHE
1	C	306	PHE
1	C	338	PHE
1	C	361	CYS
1	C	375	SER
1	C	408	ARG
1	C	429	PHE
1	C	453	TYR
1	C	495	TYR
1	C	529	LYS
1	C	552	LEU
1	C	557	LYS
1	C	559	PHE
1	C	650	LEU
1	C	655	HIS
1	C	657	ASN
1	C	660	TYR
1	C	662	CYS
1	C	673	SER
1	C	738	CYS
1	C	743	CYS
1	C	746	SER
1	C	758	SER
1	C	872	GLN
1	C	906	PHE
1	C	968	SER
1	C	1067	TYR
1	C	1082	CYS

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Mol	Chain	Res	Type
1	C	1091	ARG
2	E	1	GLU
2	E	55	ASP
2	E	75	SER
2	E	160	TRP
2	E	198	GLN
2	D	25	SER
2	D	83	TRP
2	D	107	ASP
2	D	118	SER
2	D	206	HIS
3	I	27	GLN
3	I	45	LYS
3	I	49	TYR
3	I	76	SER
3	I	92	LYS
3	I	97	TRP
3	I	108	LYS
3	F	17	ASP
3	F	97	TRP
3	F	104	LYS
3	F	108	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	ASN
1	A	115	GLN
1	A	121	ASN
1	A	122	ASN
1	A	207	HIS
1	A	239	GLN
1	A	901	GLN
1	A	907	ASN
1	A	935	GLN
1	A	1135	ASN
1	B	30	ASN
1	B	99	ASN
1	B	115	GLN
1	B	239	GLN
1	B	360	ASN
1	B	422	ASN

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Mol	Chain	Res	Type
1	B	563	GLN
1	B	658	ASN
1	B	954	GLN
1	C	30	ASN
1	C	115	GLN
1	C	207	HIS
1	C	239	GLN
1	C	439	ASN
1	C	498	GLN
1	C	564	GLN
1	C	658	ASN
1	C	913	GLN
1	C	920	GLN
1	C	978	ASN
2	E	111	GLN
2	D	177	GLN
3	I	89	GLN
3	I	139	ASN
3	F	31	ASN
3	F	34	ASN
3	F	38	GLN
3	F	53	ASN
3	F	79	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 ⓘ

10 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	B1	1	1,4	14,14,15	0.26	0	17,19,21	0.53	0
4	NAG	B1	2	4	14,14,15	0.27	0	17,19,21	0.50	0
4	NAG	G	1	1,4	14,14,15	0.40	0	17,19,21	0.72	1 (5%)
4	NAG	G	2	4	14,14,15	0.20	0	17,19,21	0.47	0
4	NAG	H	1	1,4	14,14,15	0.22	0	17,19,21	0.41	0
4	NAG	H	2	4	14,14,15	0.20	0	17,19,21	0.43	0
4	NAG	J	1	1,4	14,14,15	0.28	0	17,19,21	0.54	0
4	NAG	J	2	4	14,14,15	0.32	0	17,19,21	0.57	0
4	NAG	K	1	1,4	14,14,15	0.38	0	17,19,21	1.78	2 (11%)
4	NAG	K	2	4	14,14,15	0.42	0	17,19,21	1.04	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	B1	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	B1	2	4	-	3/6/23/26	0/1/1/1
4	NAG	G	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
4	NAG	H	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1
4	NAG	J	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	J	2	4	-	3/6/23/26	0/1/1/1
4	NAG	K	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	K	2	4	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	K	1	NAG	C1-O5-C5	6.36	120.80	112.19
4	K	2	NAG	C1-O5-C5	3.20	116.53	112.19
4	K	1	NAG	C3-C4-C5	2.53	114.76	110.24
4	K	2	NAG	C3-C4-C5	2.11	114.00	110.24
4	G	1	NAG	C1-O5-C5	2.00	114.90	112.19

There are no chirality outliers.

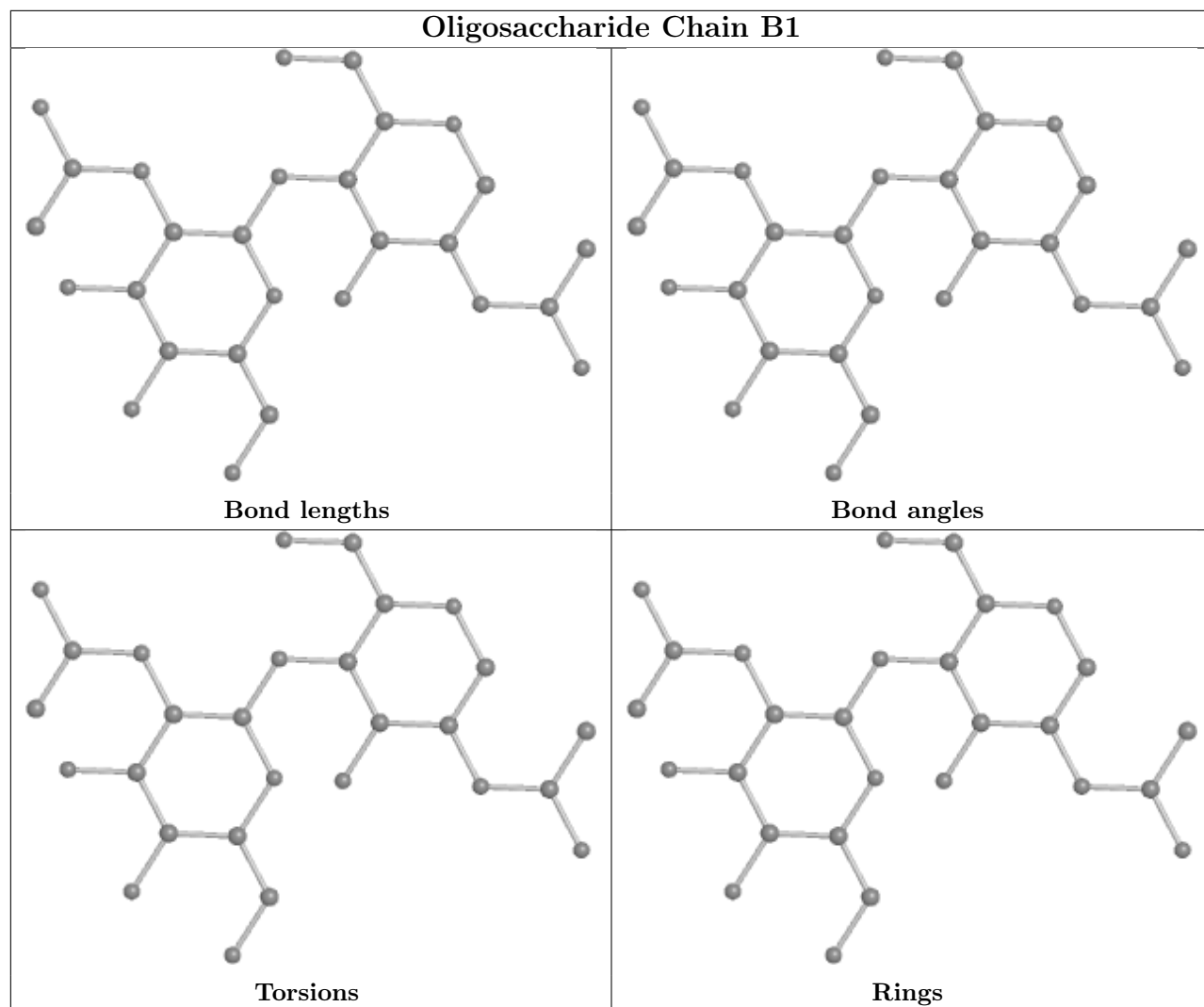
All (22) torsion outliers are listed below:

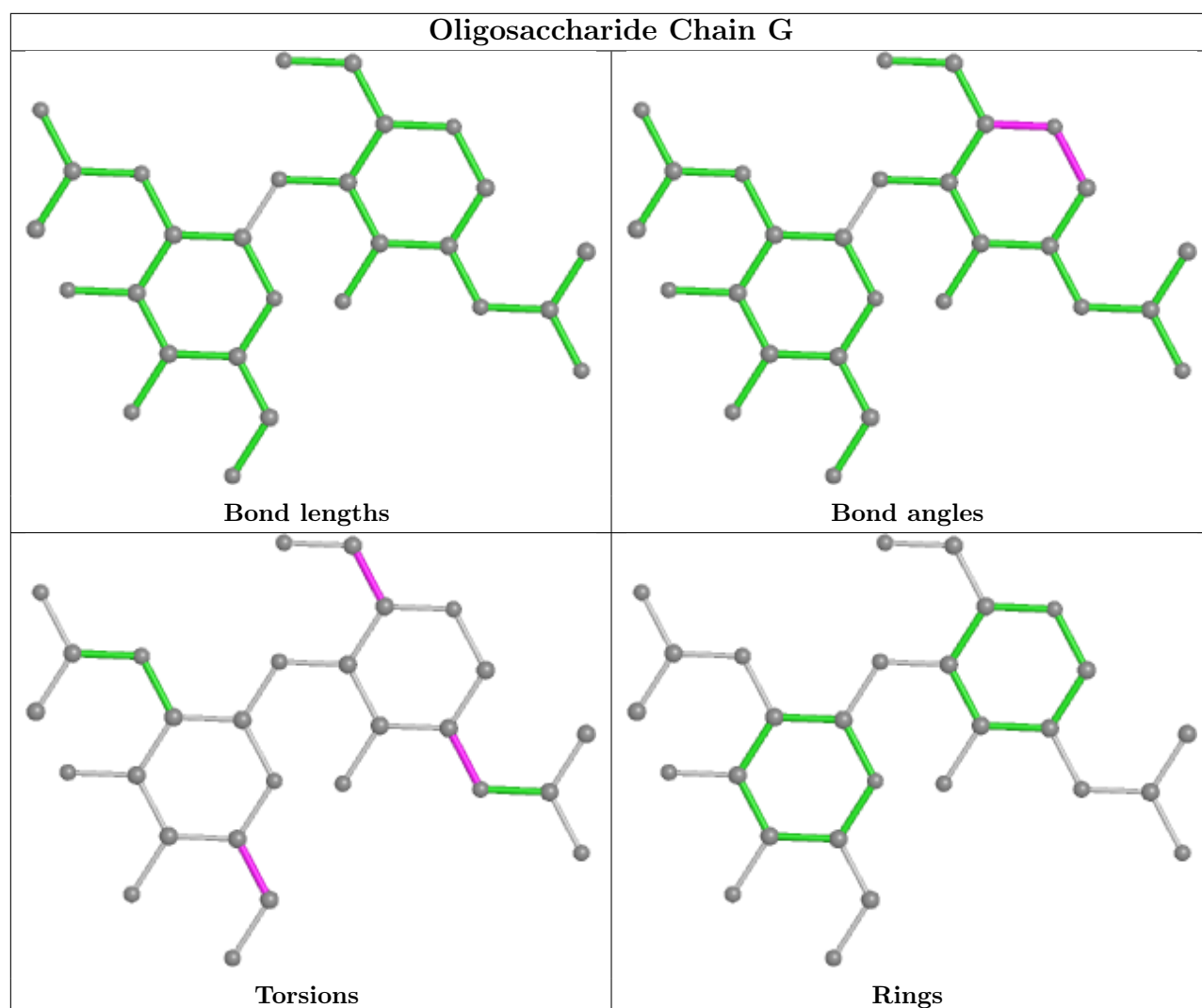
Mol	Chain	Res	Type	Atoms
4	B1	2	NAG	O5-C5-C6-O6
4	K	1	NAG	O5-C5-C6-O6
4	K	1	NAG	C4-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	B1	2	NAG	C4-C5-C6-O6
4	H	2	NAG	C4-C5-C6-O6
4	H	2	NAG	O5-C5-C6-O6
4	K	2	NAG	C8-C7-N2-C2
4	K	2	NAG	O7-C7-N2-C2
4	G	2	NAG	C4-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
4	J	1	NAG	C4-C5-C6-O6
4	J	1	NAG	O5-C5-C6-O6
4	B1	1	NAG	C3-C2-N2-C7
4	B1	2	NAG	C3-C2-N2-C7
4	G	1	NAG	C3-C2-N2-C7
4	J	1	NAG	C3-C2-N2-C7
4	J	2	NAG	C3-C2-N2-C7
4	K	1	NAG	C3-C2-N2-C7
4	J	2	NAG	C4-C5-C6-O6
4	J	2	NAG	O5-C5-C6-O6

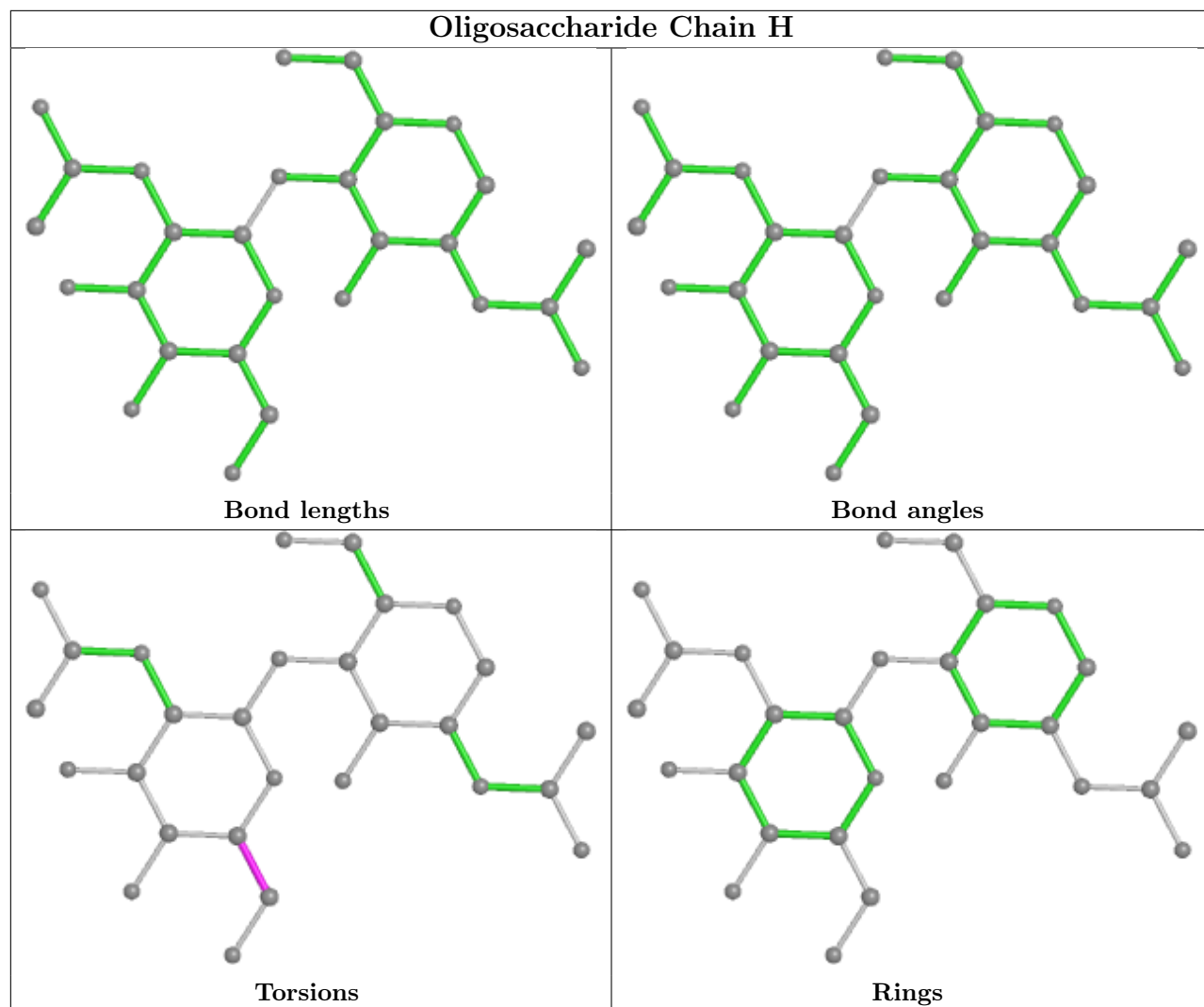
There are no ring outliers.

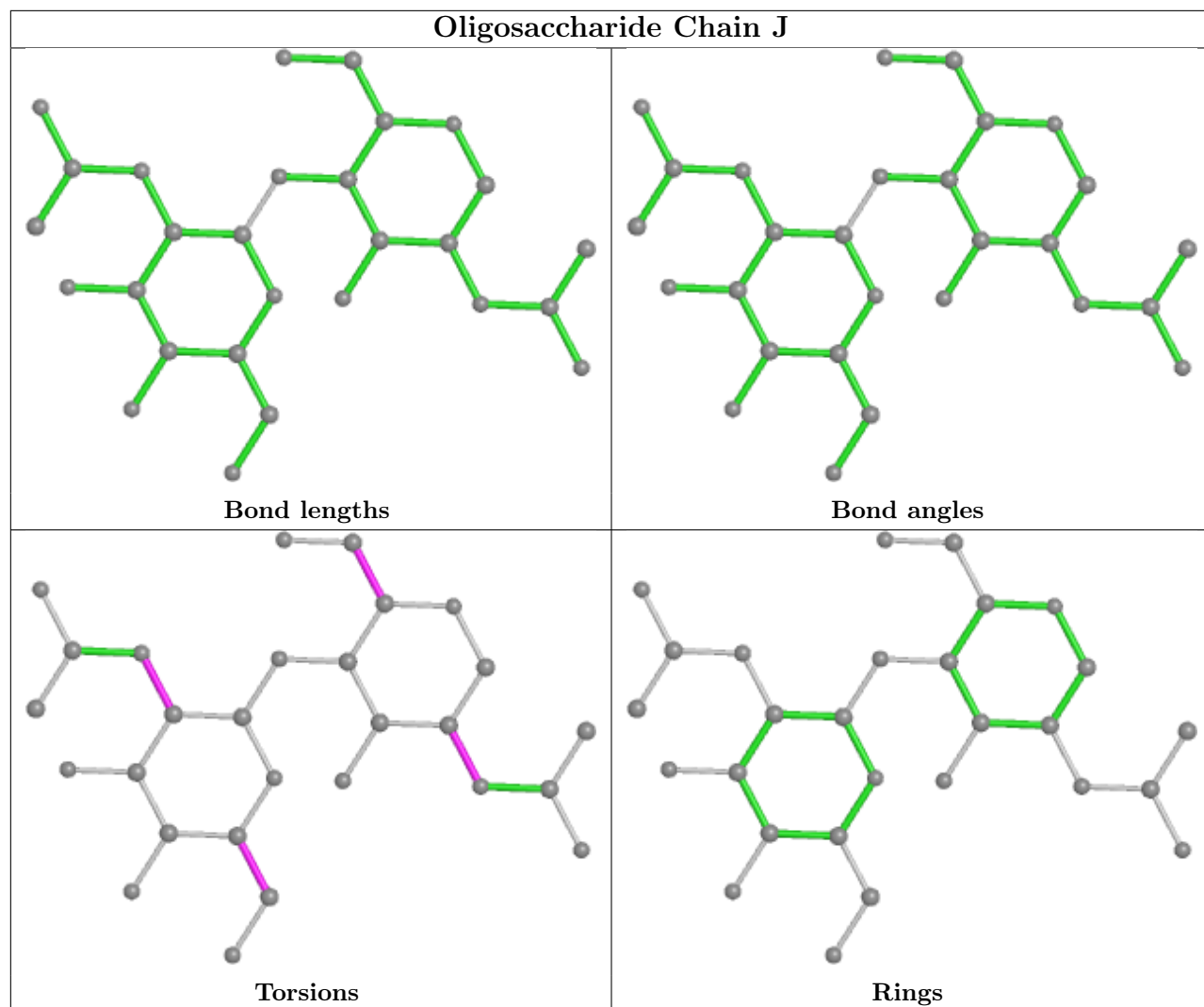
No monomer is involved in short contacts.

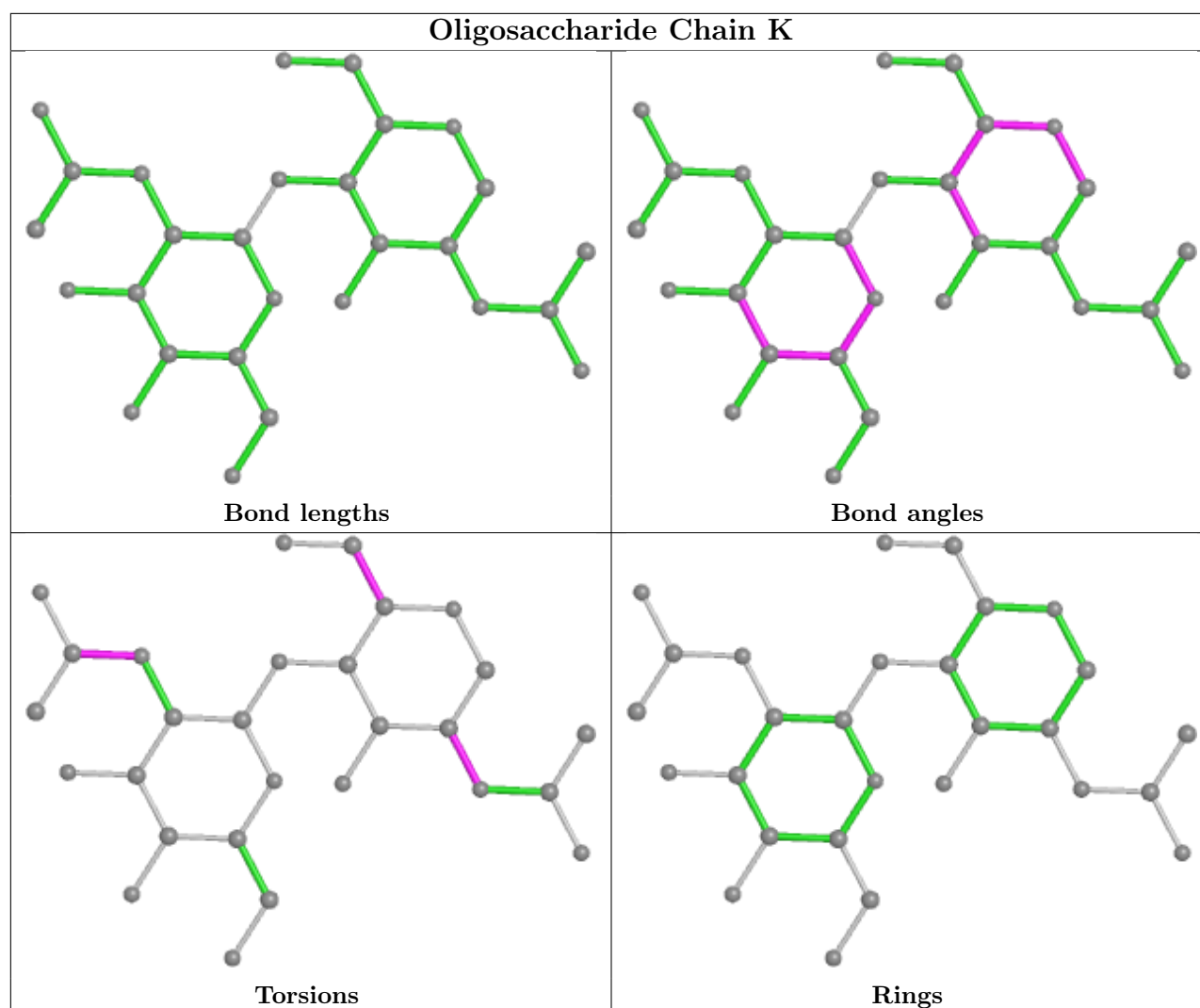
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

22 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	$\# Z > 2$
5	NAG	A	1304	1	14,14,15	0.29	0	17,19,21	0.50	0
5	NAG	C	1306	1	14,14,15	0.24	0	17,19,21	0.46	0
5	NAG	B	1302	1	14,14,15	0.25	0	17,19,21	0.51	0
5	NAG	B	1304	1	14,14,15	0.30	0	17,19,21	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	A	1308	1	14,14,15	0.21	0	17,19,21	0.43	0
5	NAG	B	1305	1	14,14,15	0.27	0	17,19,21	0.45	0
5	NAG	B	1303	1	14,14,15	0.24	0	17,19,21	0.47	0
5	NAG	C	1307	1	14,14,15	0.19	0	17,19,21	0.43	0
5	NAG	C	1305	1	14,14,15	0.21	0	17,19,21	0.42	0
5	NAG	B	1301	1	14,14,15	0.85	1 (7%)	17,19,21	0.97	1 (5%)
5	NAG	C	1302	1	14,14,15	0.27	0	17,19,21	0.55	0
5	NAG	A	1305	1	14,14,15	0.21	0	17,19,21	0.43	0
5	NAG	C	1303	1	14,14,15	0.23	0	17,19,21	0.45	0
5	NAG	C	1301	1	14,14,15	0.65	1 (7%)	17,19,21	0.75	0
5	NAG	A	1306	1	14,14,15	0.27	0	17,19,21	0.44	0
5	NAG	B	1306	1	14,14,15	0.20	0	17,19,21	0.41	0
5	NAG	C	1308	1	14,14,15	0.22	0	17,19,21	0.43	0
5	NAG	A	1303	1	14,14,15	0.25	0	17,19,21	0.52	0
5	NAG	C	1304	1	14,14,15	0.30	0	17,19,21	0.51	0
5	NAG	A	1301	1	14,14,15	0.83	1 (7%)	17,19,21	1.02	1 (5%)
5	NAG	A	1307	1	14,14,15	0.32	0	17,19,21	0.42	0
5	NAG	A	1302	1	14,14,15	0.22	0	17,19,21	0.43	0

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

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

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1301	NAG	O5-C1	2.75	1.48	1.43
5	A	1301	NAG	O5-C1	2.57	1.47	1.43
5	C	1301	NAG	O5-C1	-2.23	1.40	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1301	NAG	C1-O5-C5	3.80	117.34	112.19
5	B	1301	NAG	C1-O5-C5	3.76	117.29	112.19

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	1308	NAG	O5-C5-C6-O6
5	A	1301	NAG	O5-C5-C6-O6
5	B	1301	NAG	O5-C5-C6-O6
5	C	1301	NAG	O5-C5-C6-O6
5	C	1308	NAG	C4-C5-C6-O6
5	B	1301	NAG	C4-C5-C6-O6
5	C	1301	NAG	C4-C5-C6-O6
5	A	1301	NAG	C4-C5-C6-O6
5	A	1306	NAG	O5-C5-C6-O6
5	B	1306	NAG	C4-C5-C6-O6
5	C	1308	NAG	C8-C7-N2-C2
5	C	1308	NAG	O7-C7-N2-C2
5	A	1306	NAG	C4-C5-C6-O6
5	A	1304	NAG	O5-C5-C6-O6
5	B	1305	NAG	O5-C5-C6-O6

Continued on next page...

Continued from previous page...

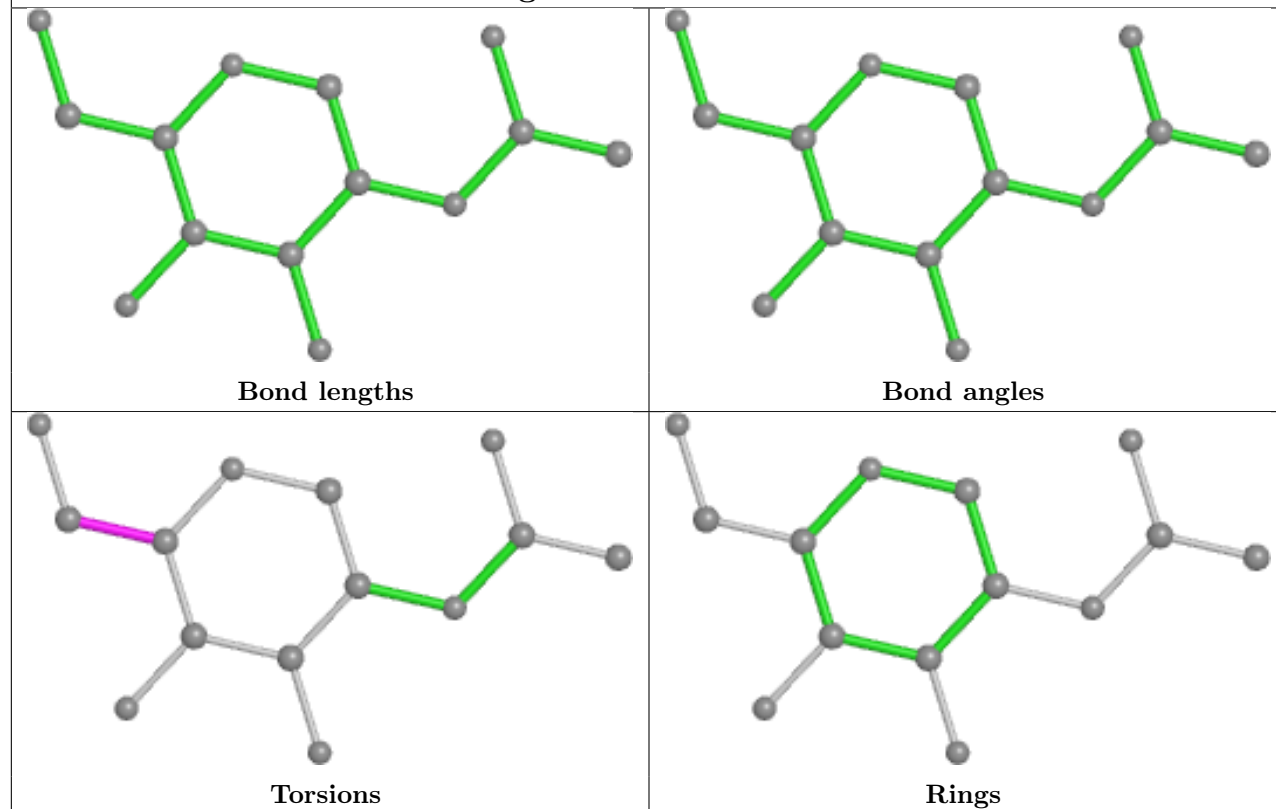
Mol	Chain	Res	Type	Atoms
5	C	1305	NAG	O5-C5-C6-O6
5	A	1304	NAG	C4-C5-C6-O6
5	B	1306	NAG	O5-C5-C6-O6
5	B	1305	NAG	C4-C5-C6-O6
5	B	1303	NAG	C4-C5-C6-O6
5	A	1307	NAG	O5-C5-C6-O6
5	B	1302	NAG	O5-C5-C6-O6
5	C	1302	NAG	C4-C5-C6-O6
5	C	1307	NAG	C4-C5-C6-O6
5	C	1307	NAG	O5-C5-C6-O6
5	B	1303	NAG	O5-C5-C6-O6
5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

There are no ring outliers.

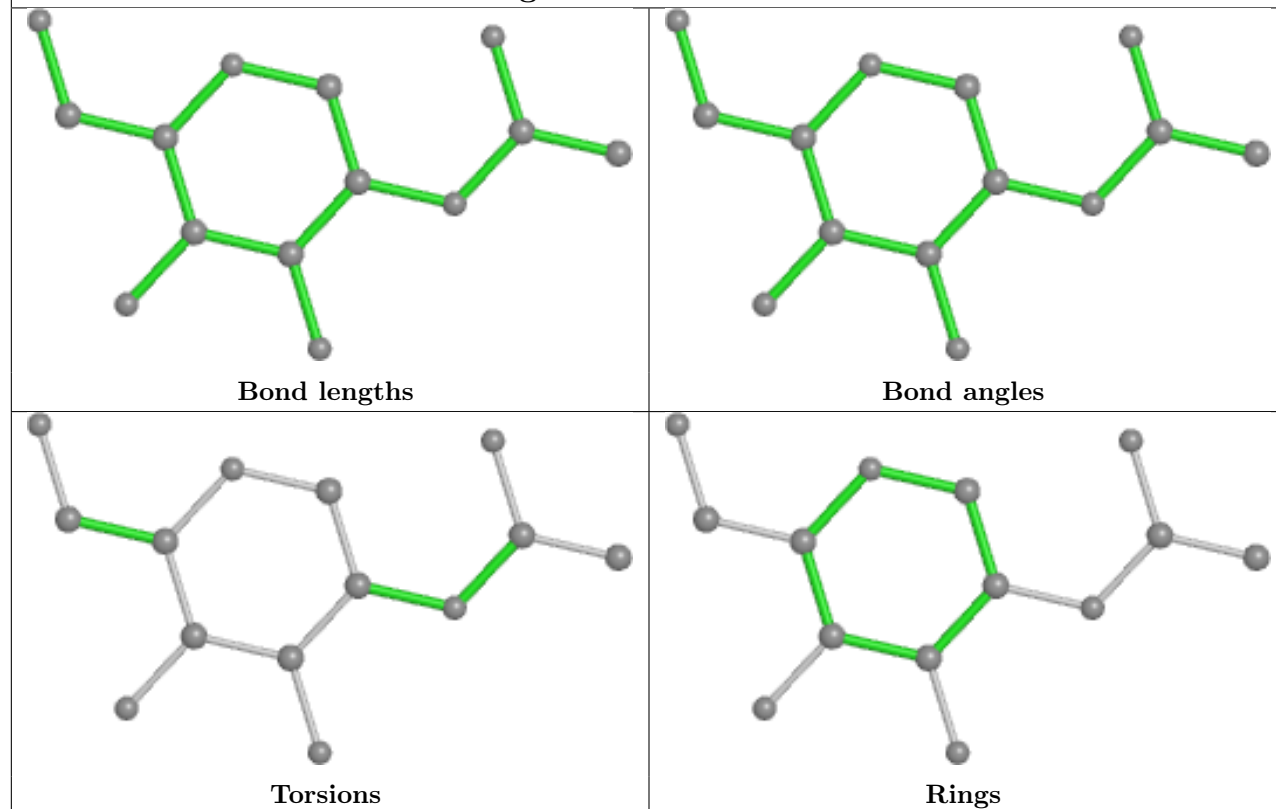
No monomer is involved in short contacts.

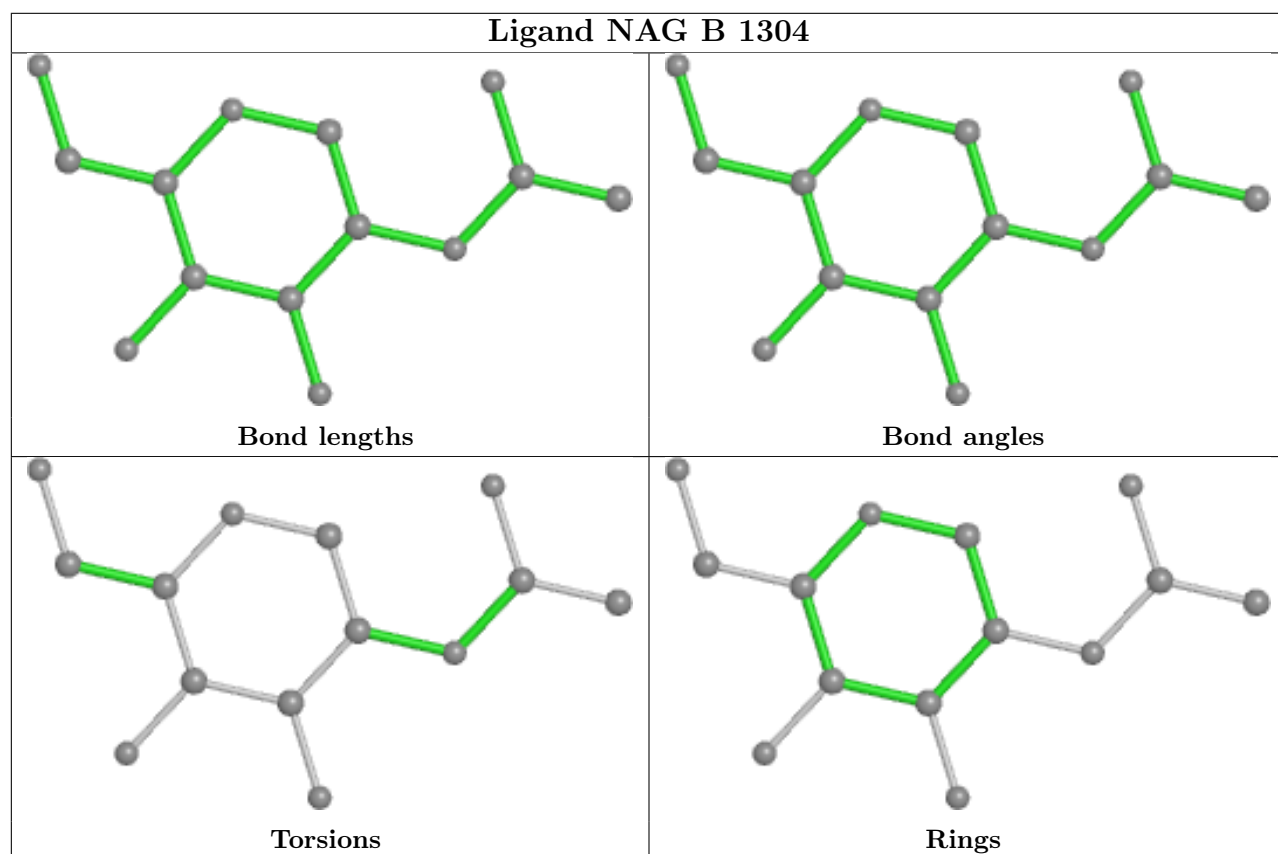
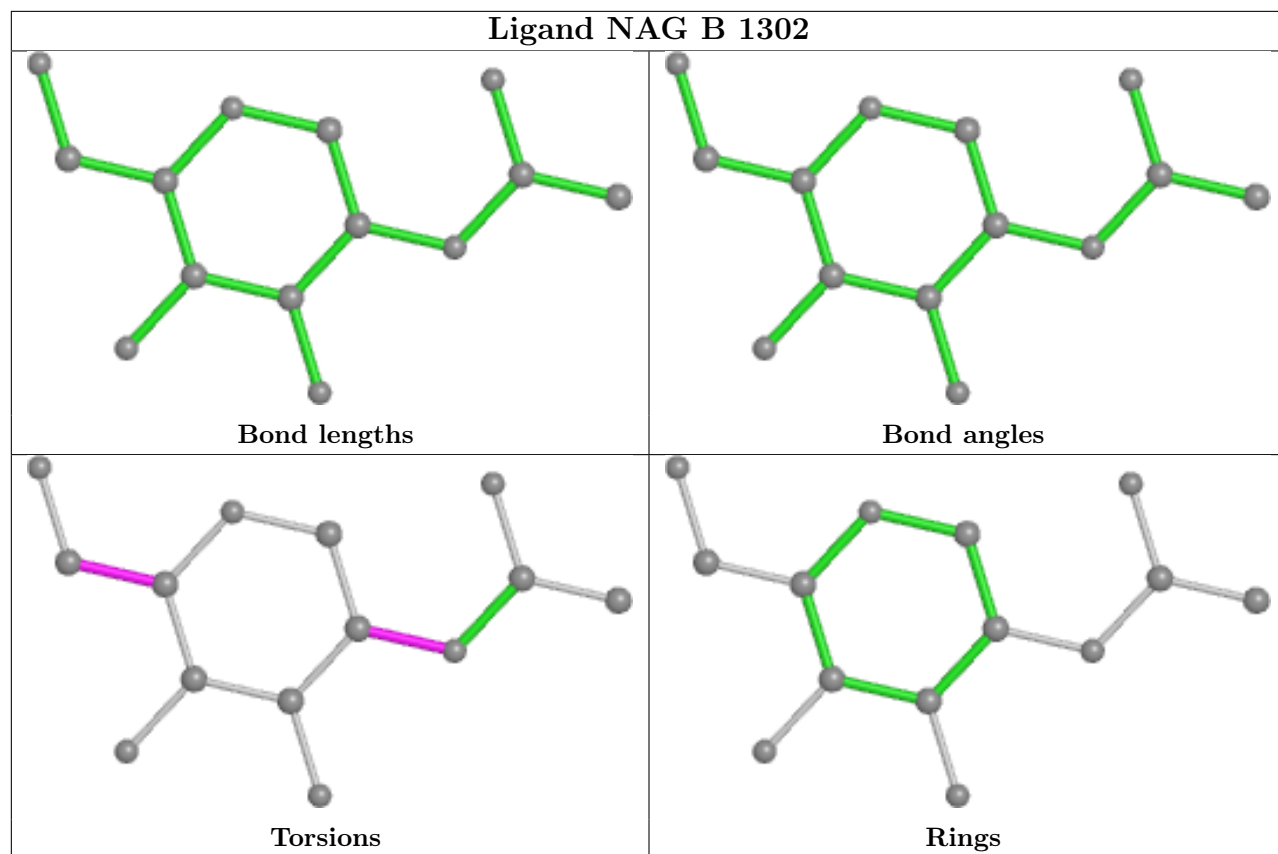
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

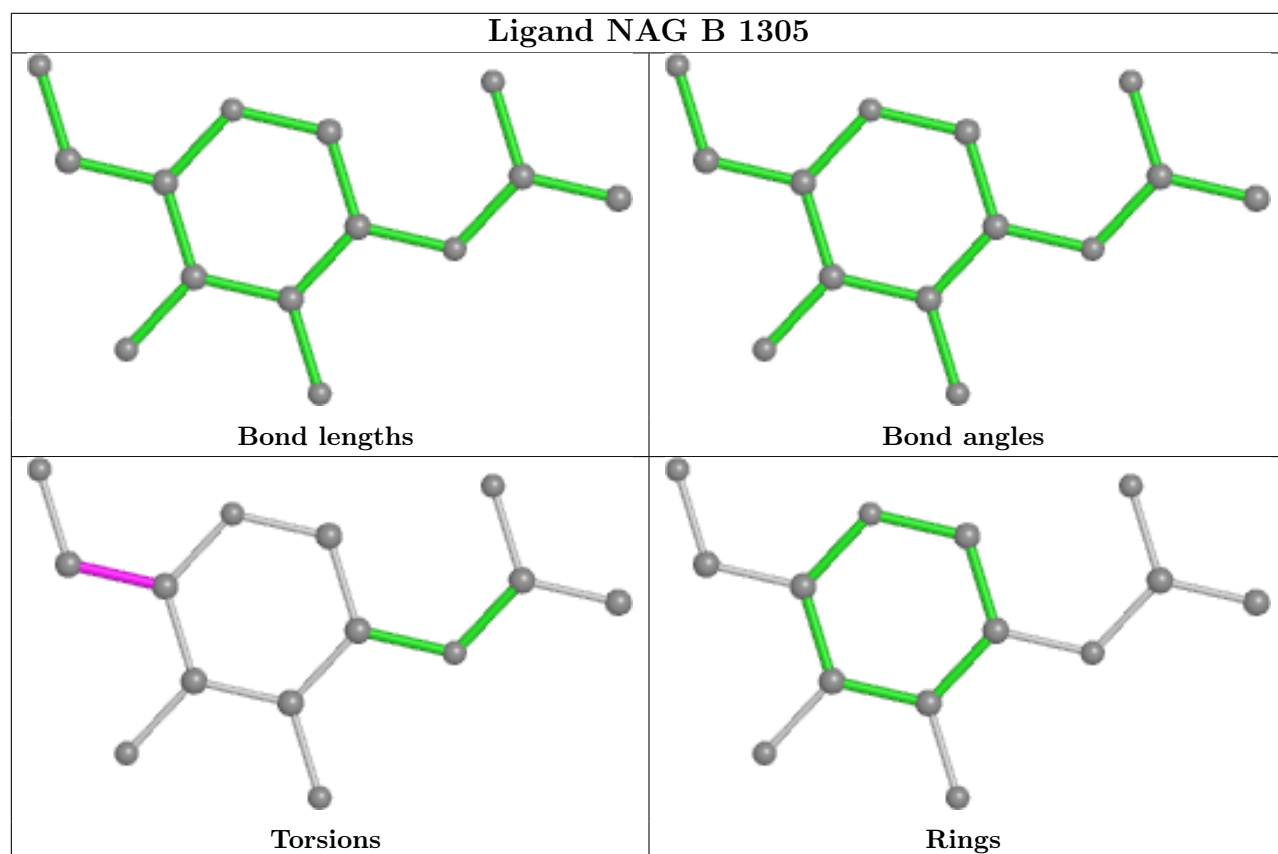
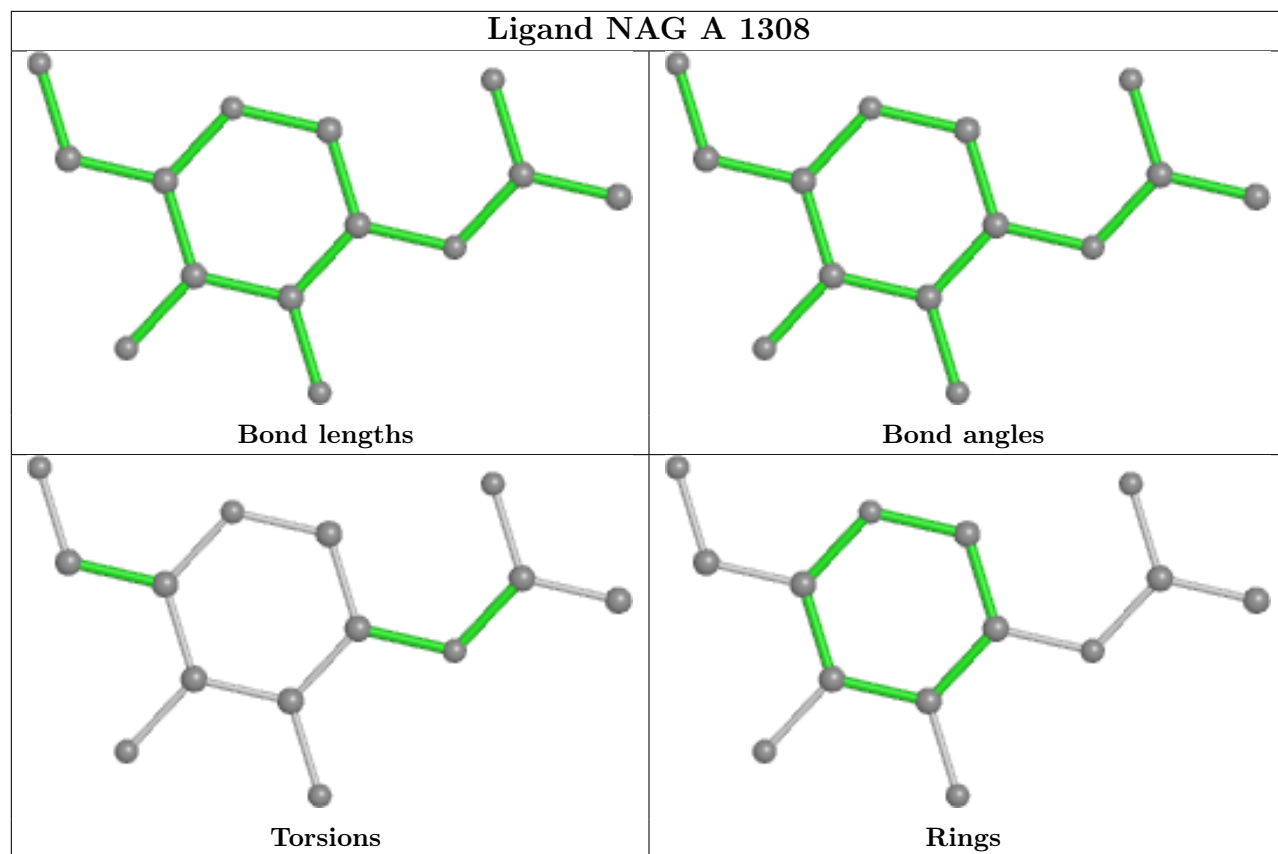
Ligand NAG A 1304

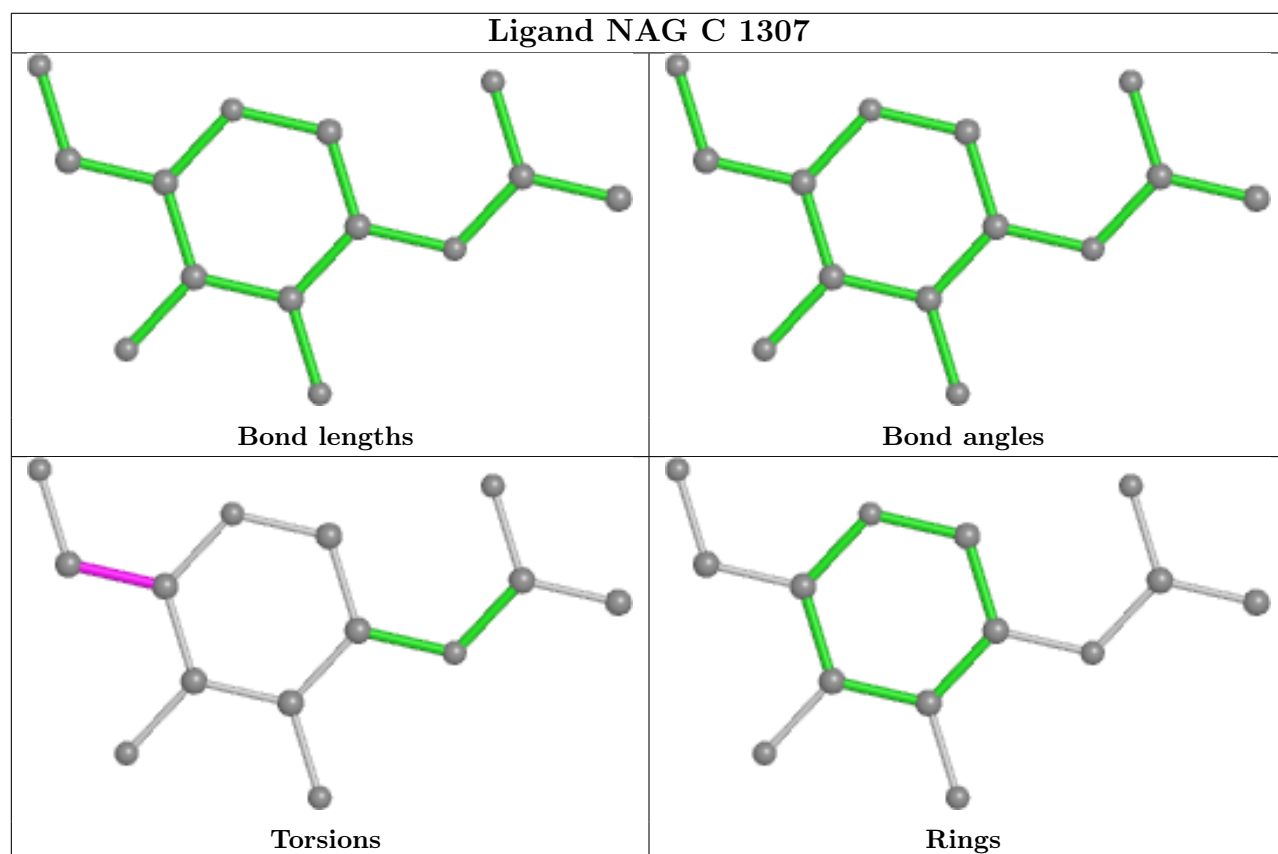
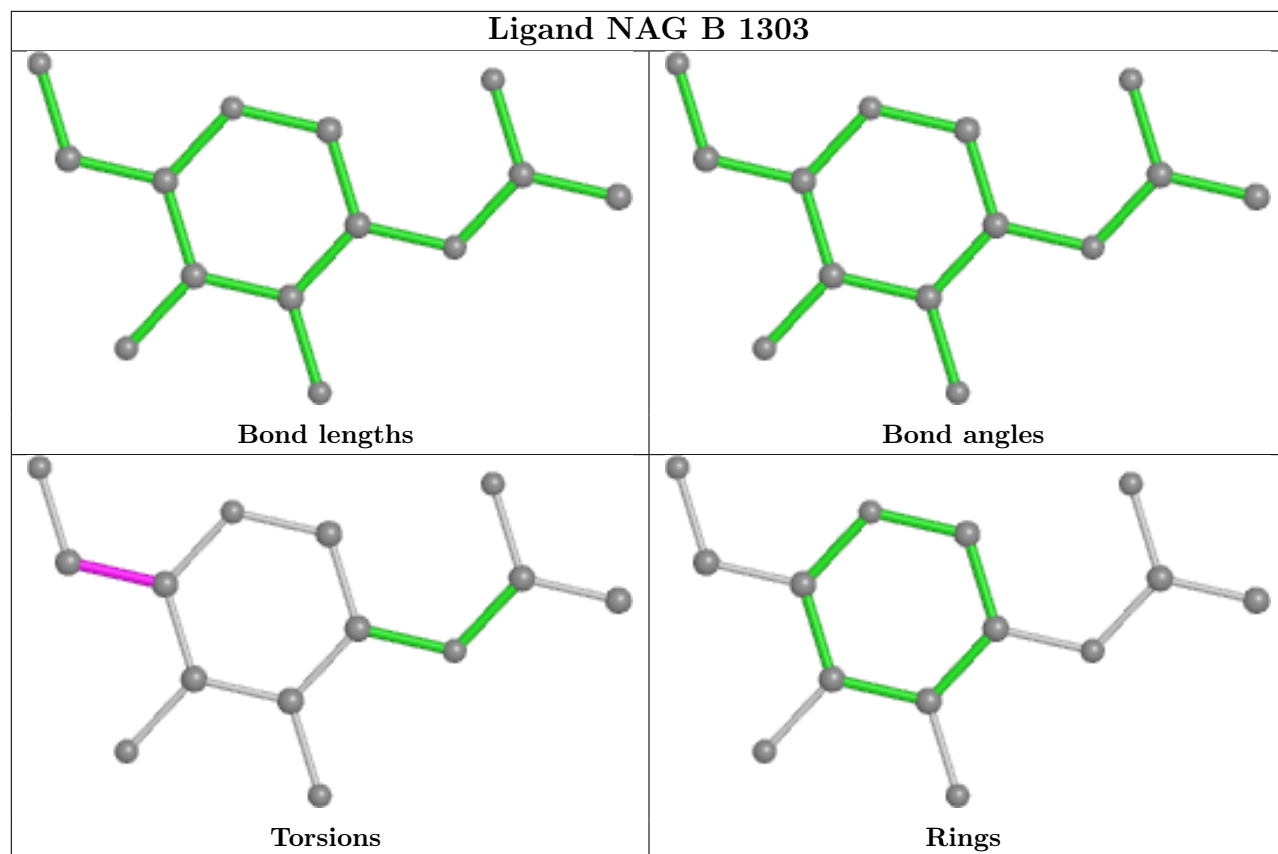


Ligand NAG C 1306

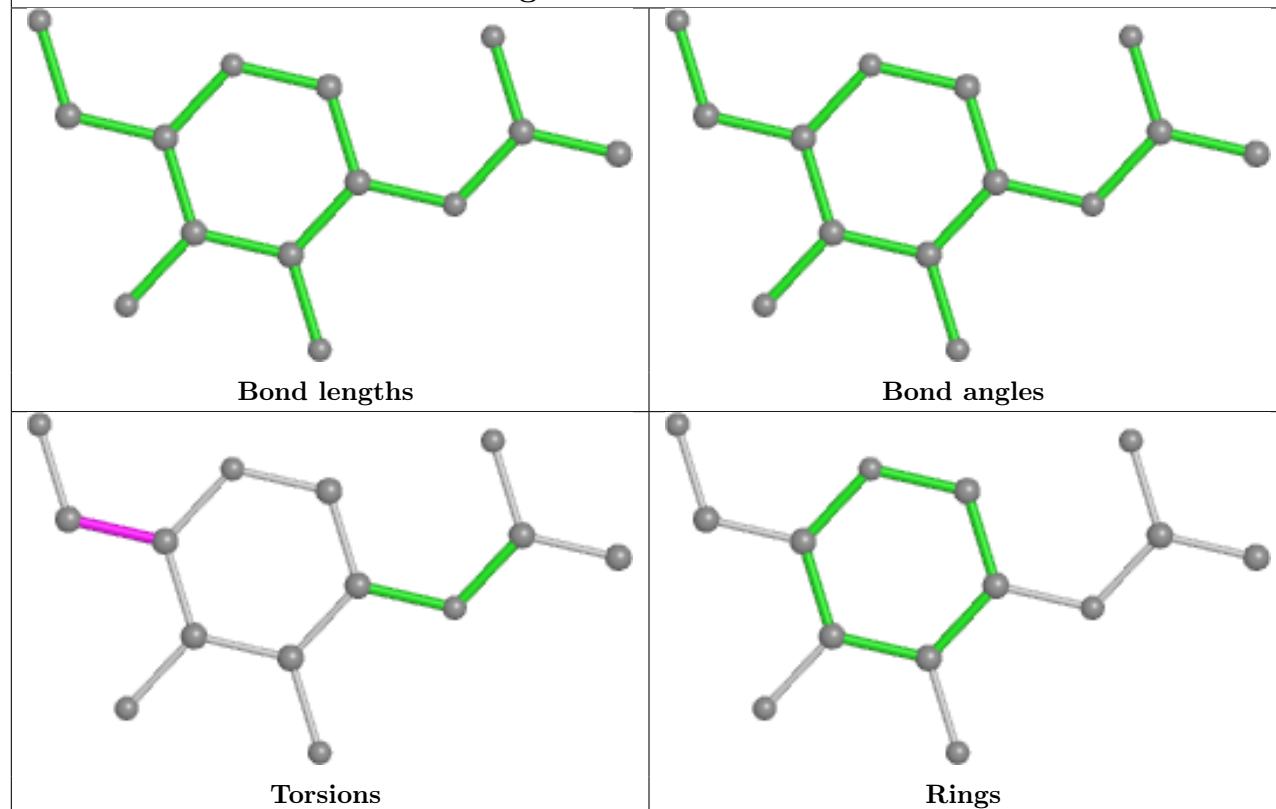




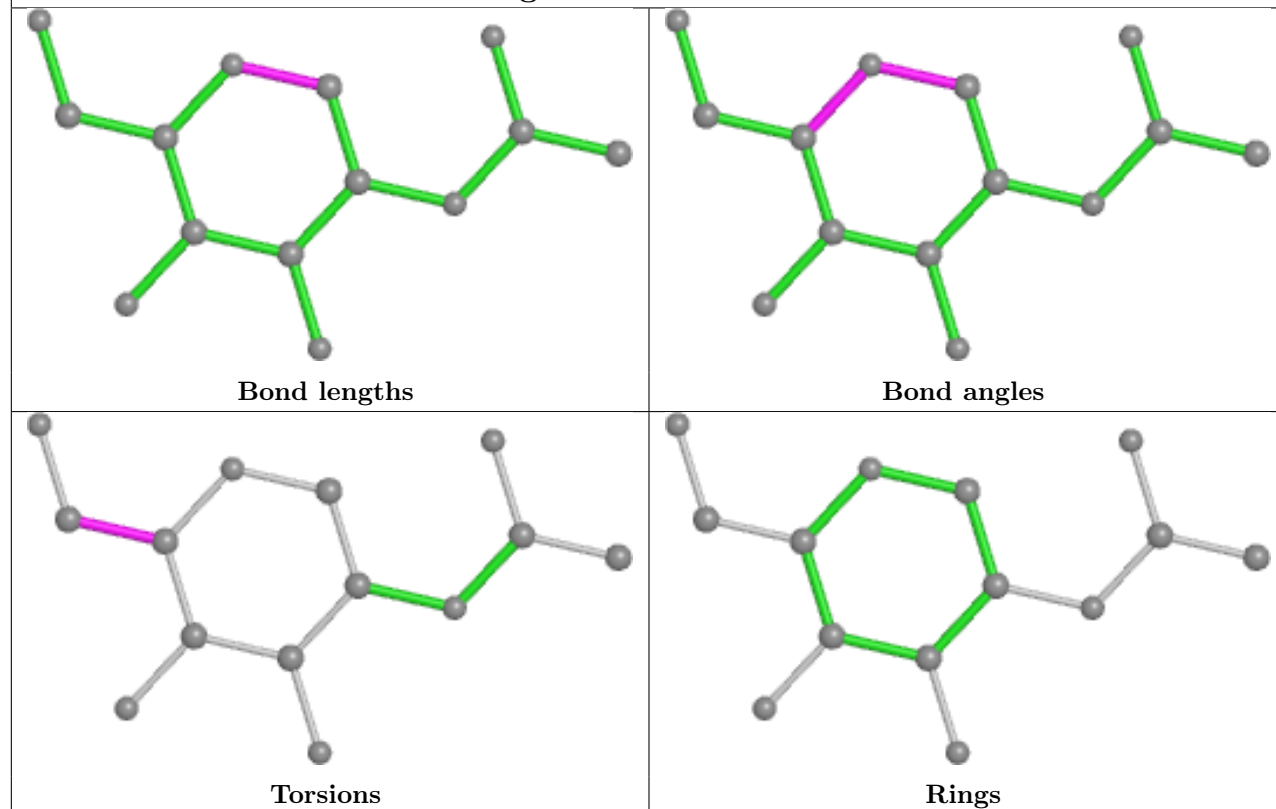




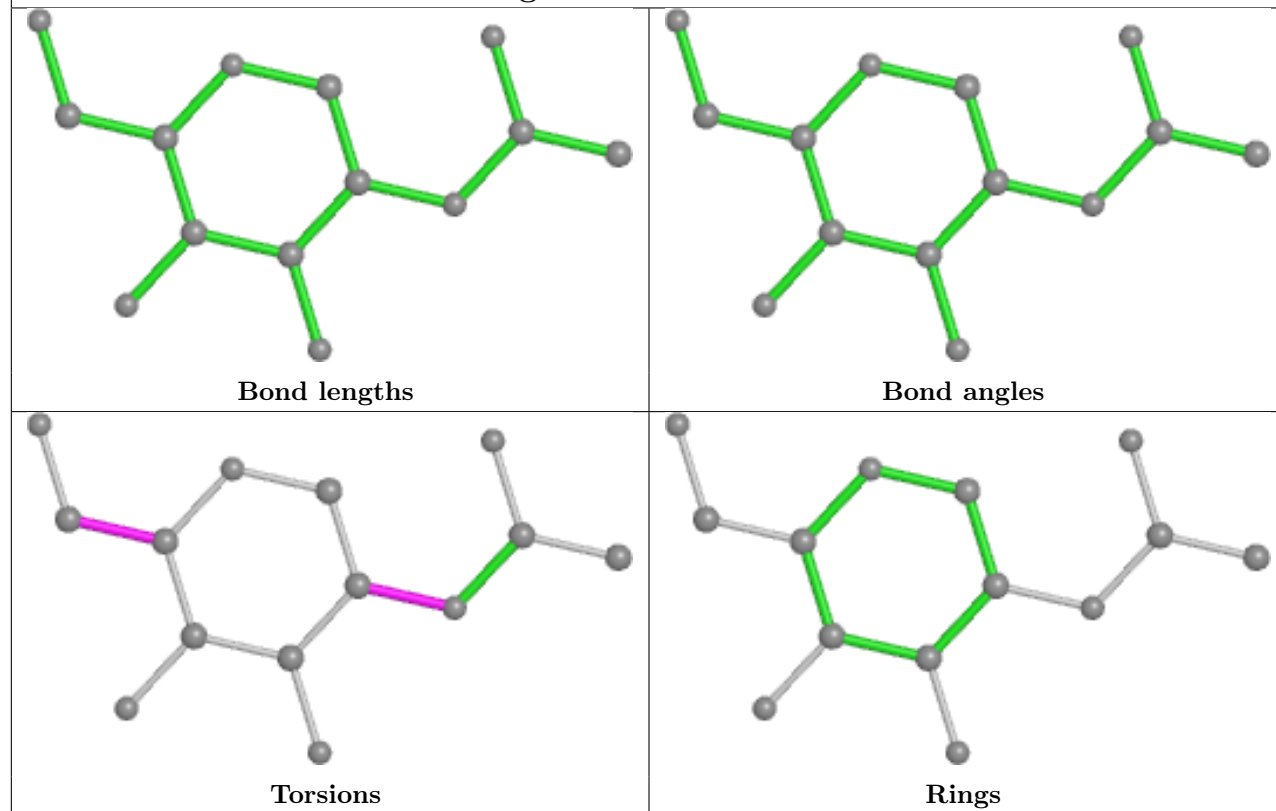
Ligand NAG C 1305



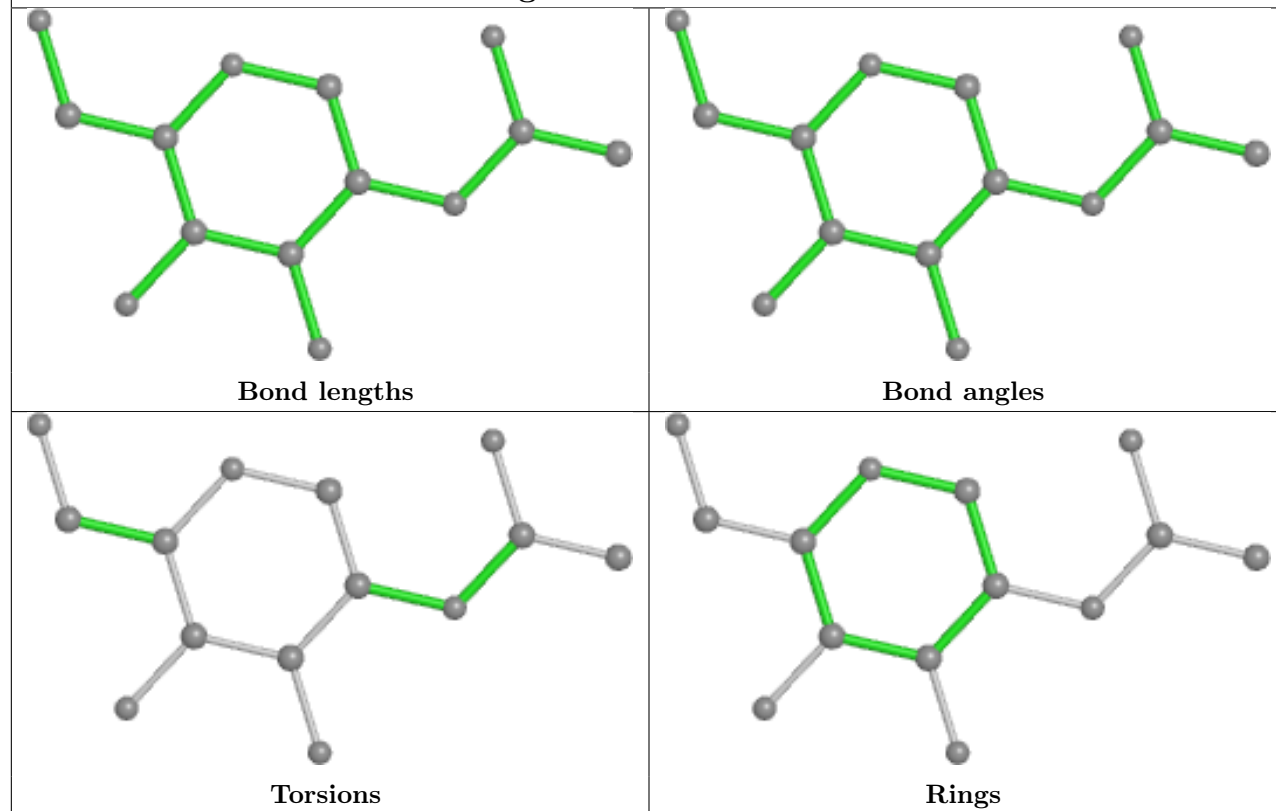
Ligand NAG B 1301

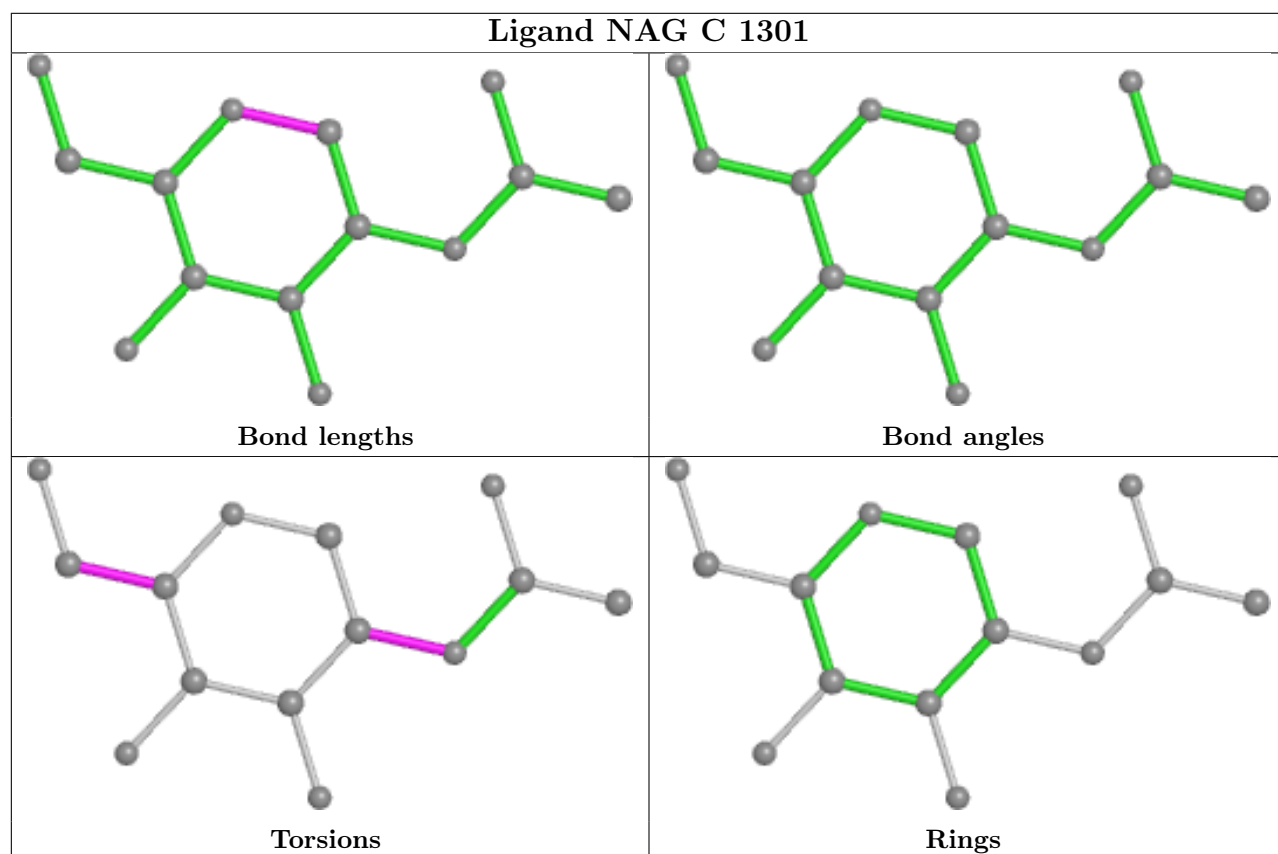
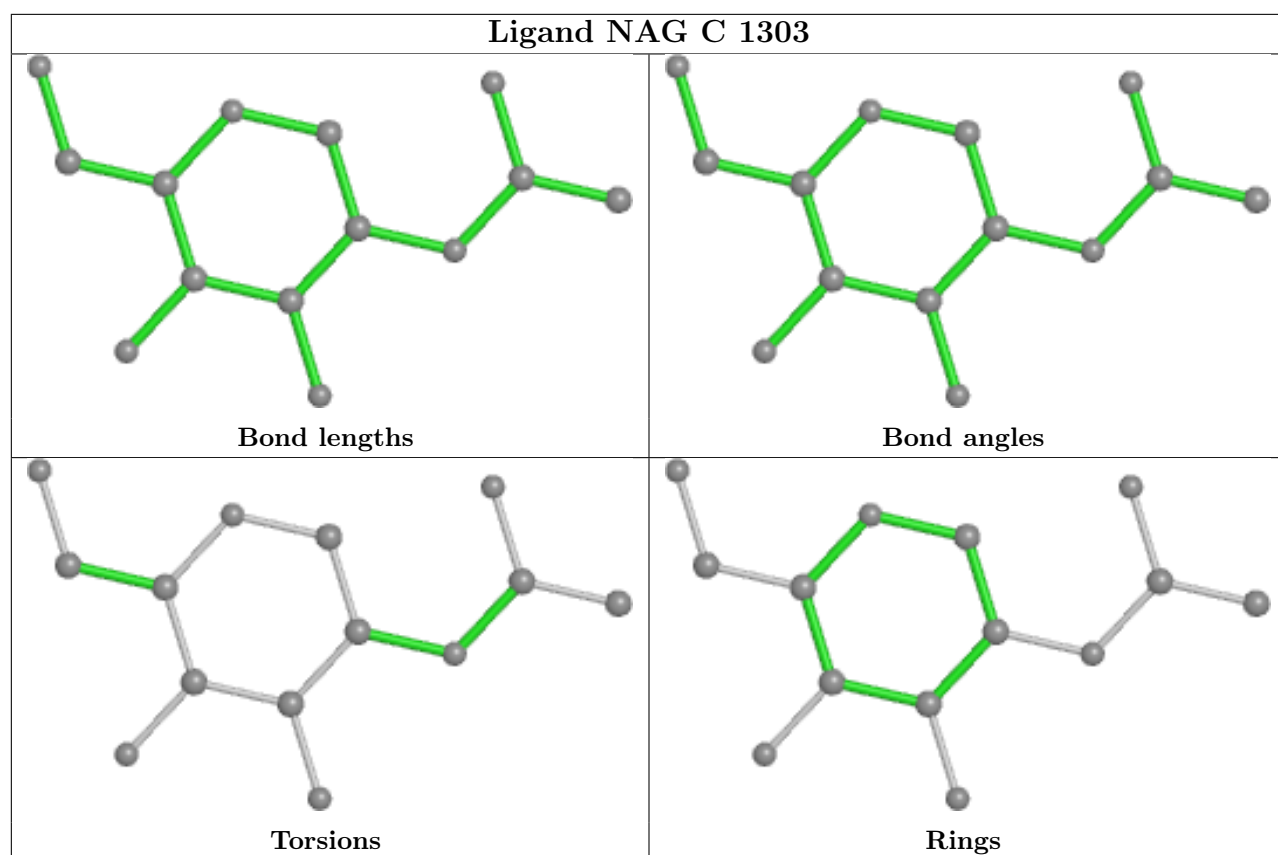


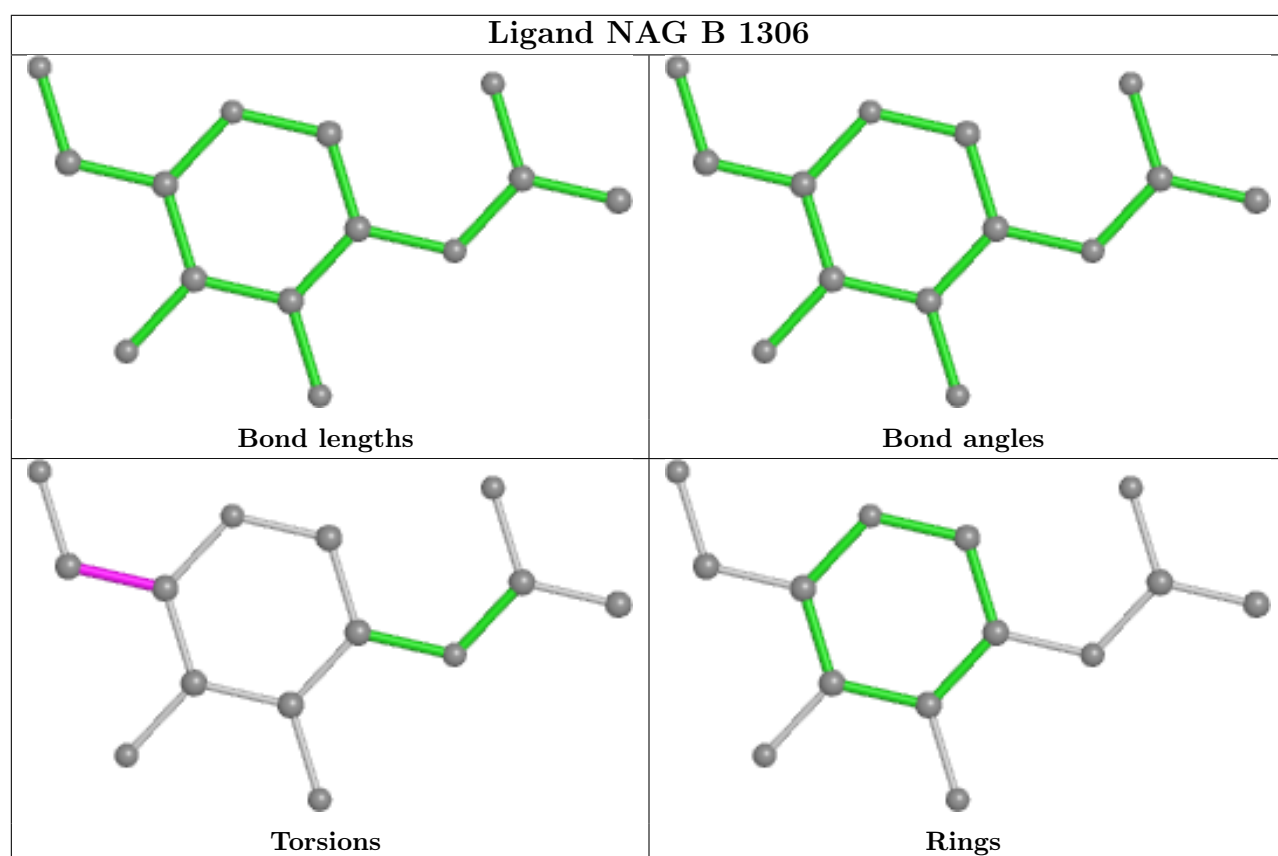
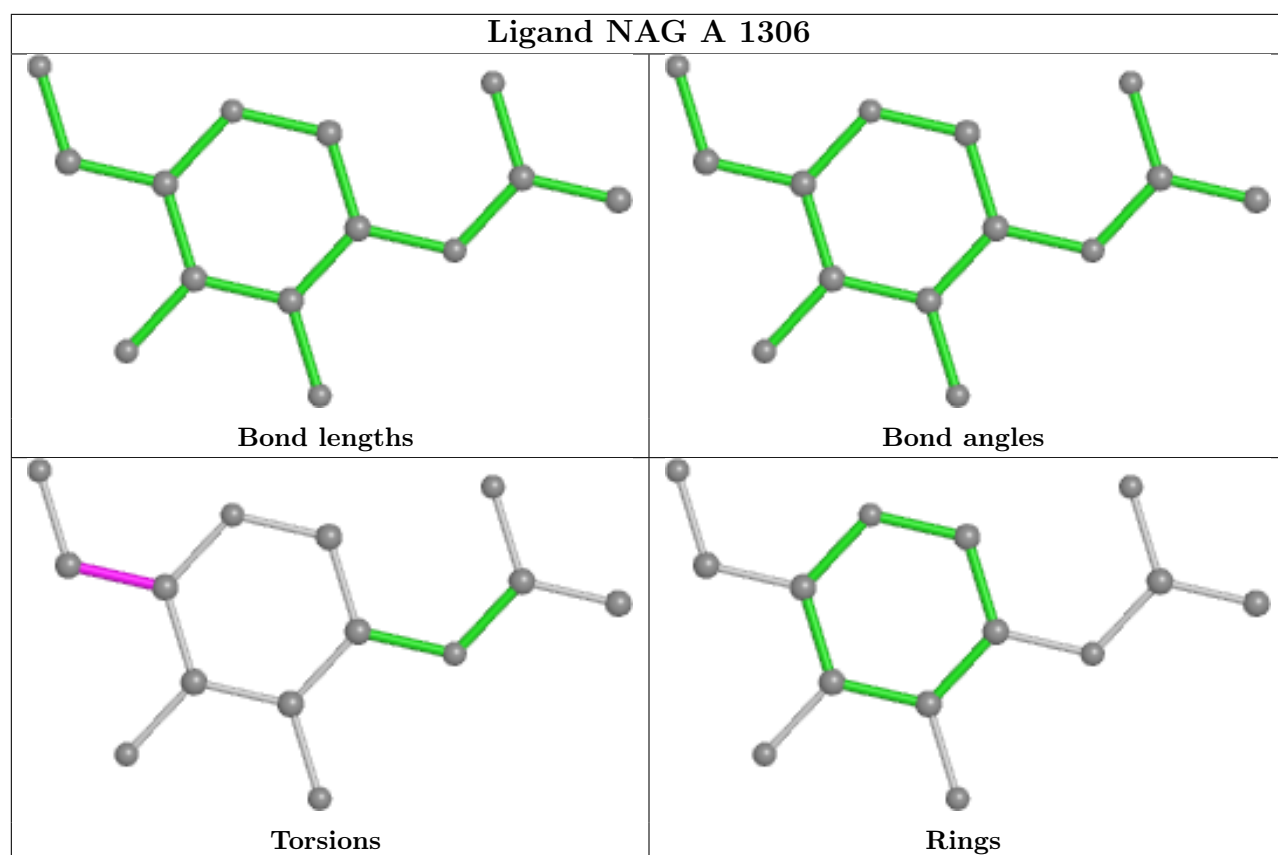
Ligand NAG C 1302



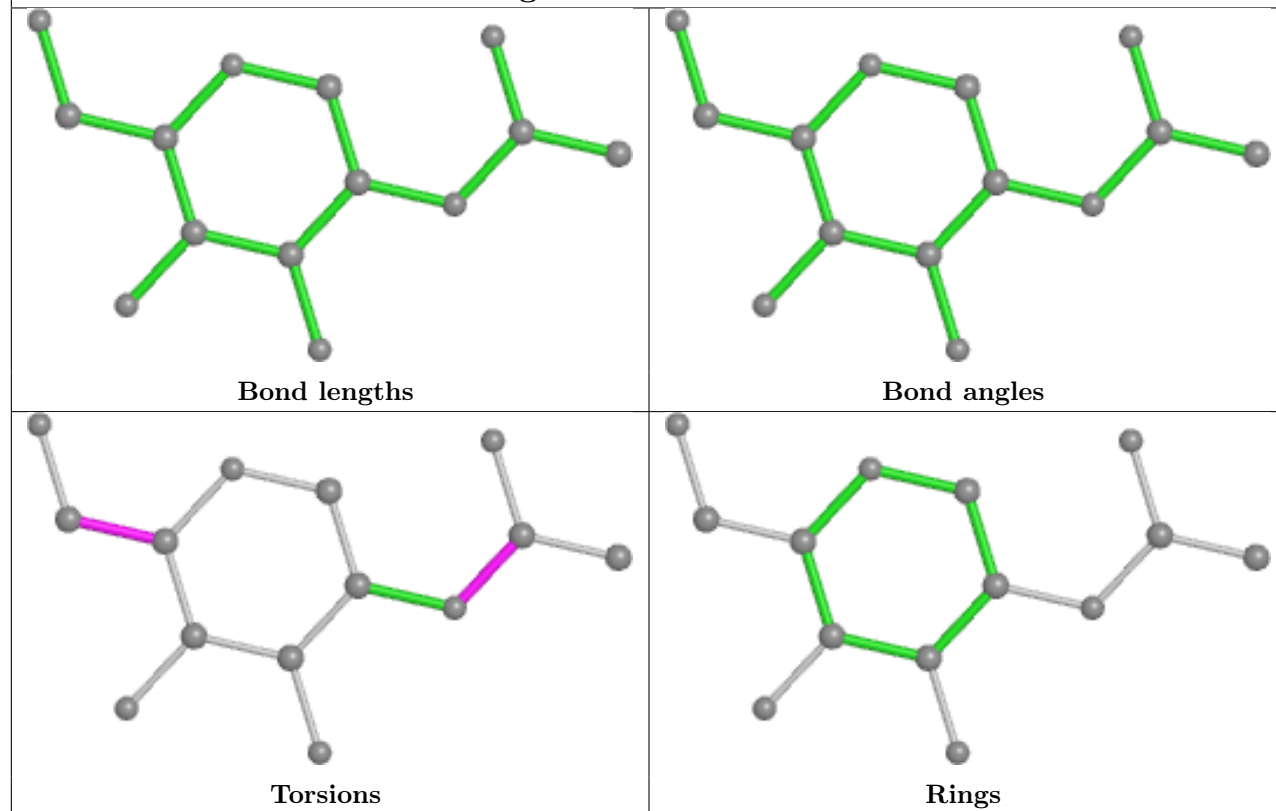
Ligand NAG A 1305



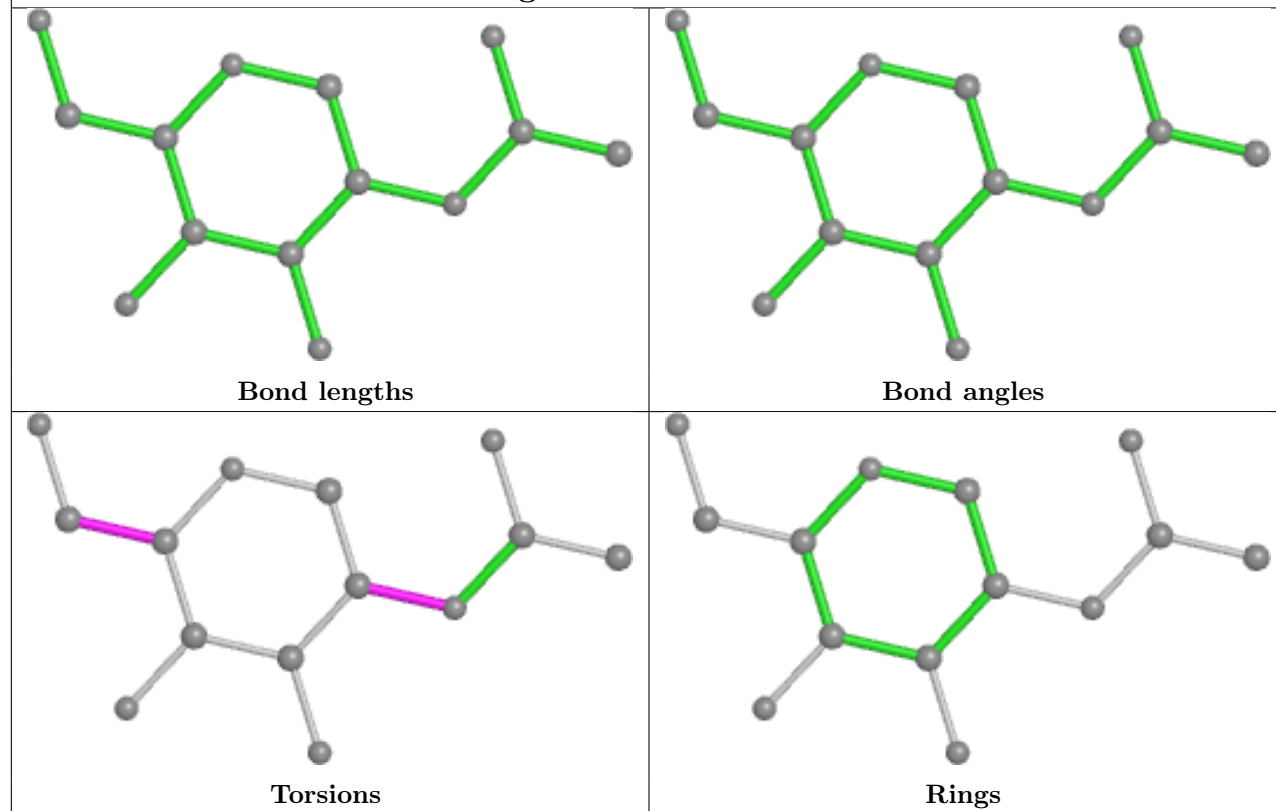




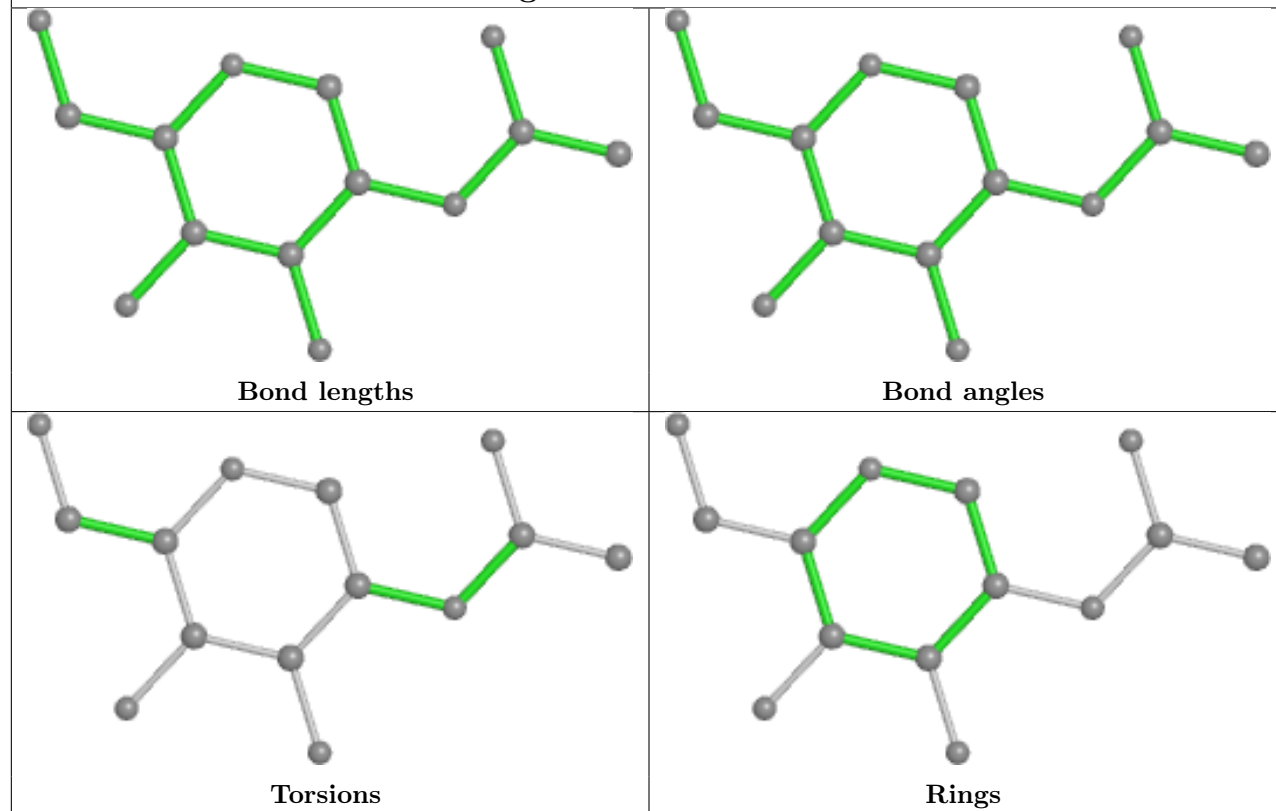
Ligand NAG C 1308



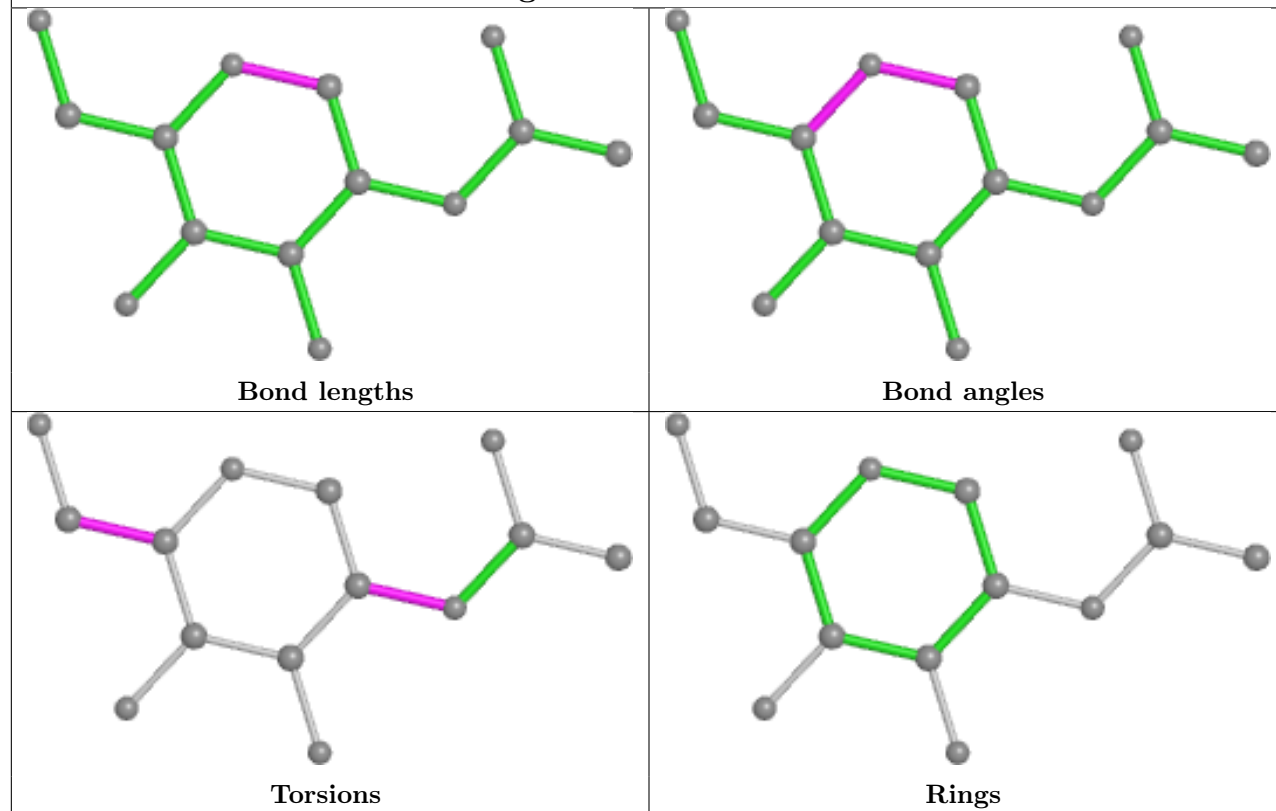
Ligand NAG A 1303

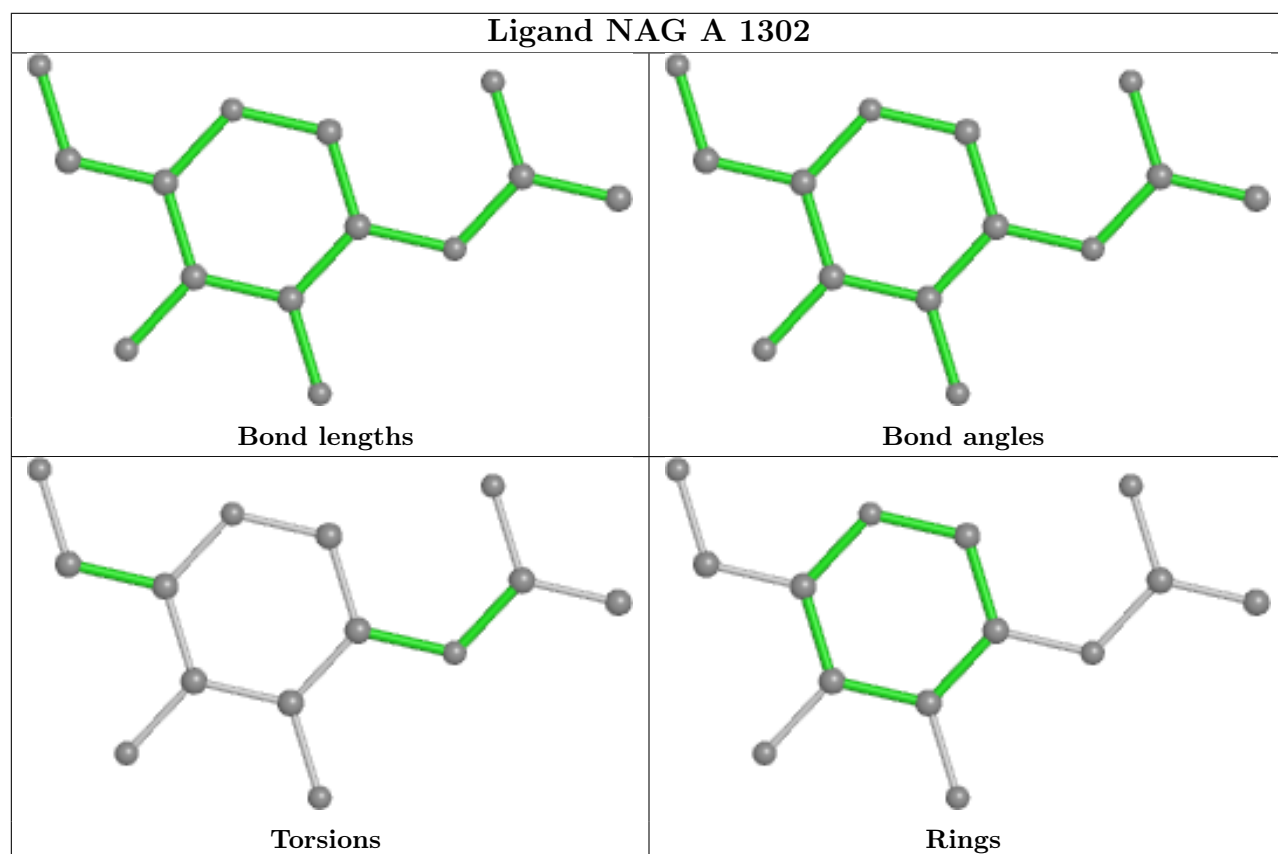
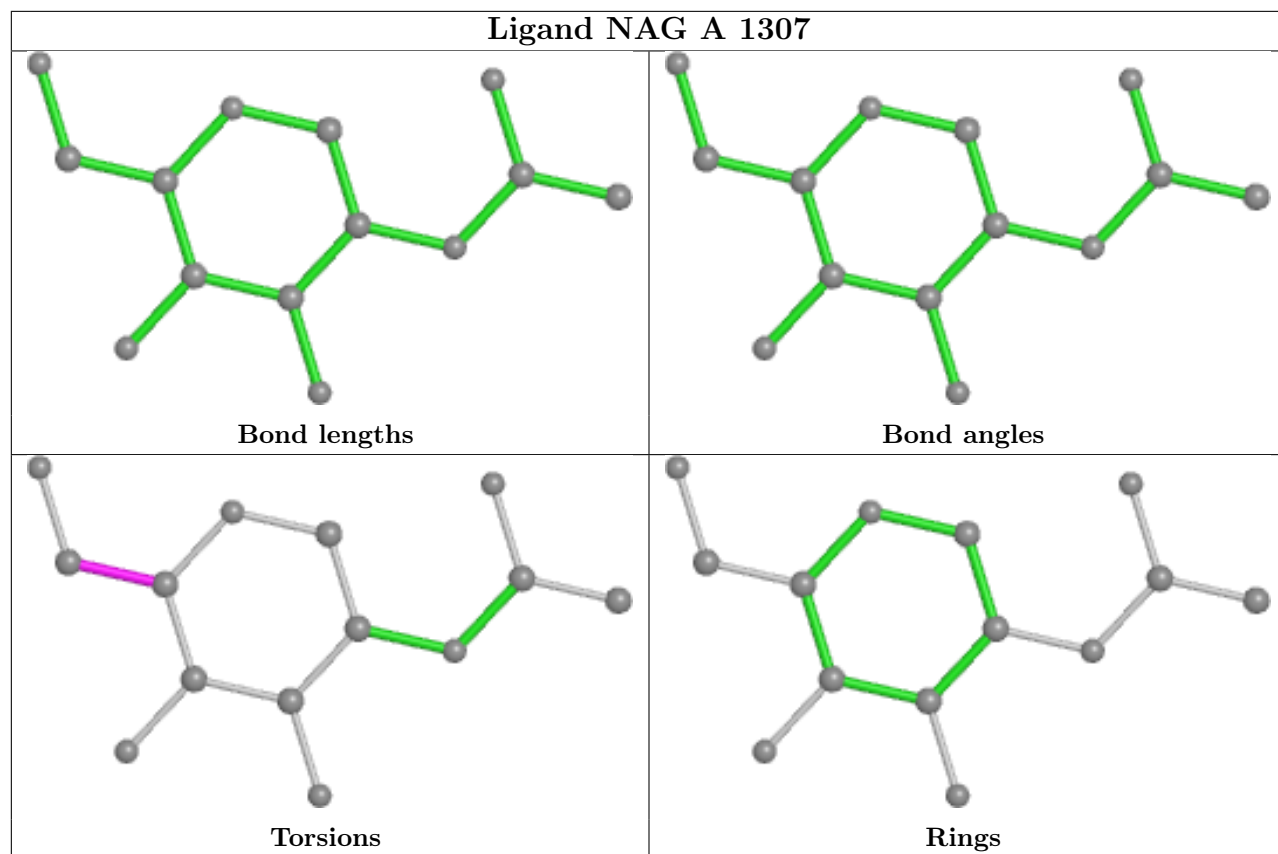


Ligand NAG C 1304



Ligand NAG A 1301





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.

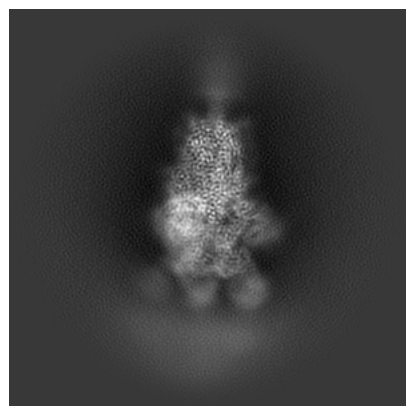
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26262. These allow visual inspection of the internal detail of the map and identification of artifacts.

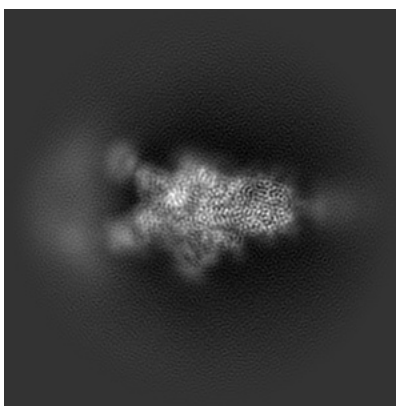
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

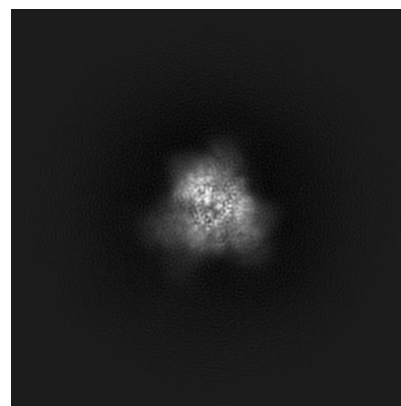
6.1.1 Primary map



X

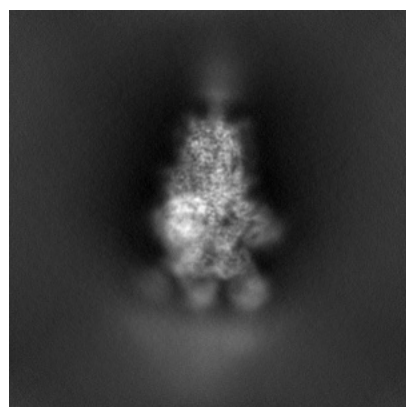


Y

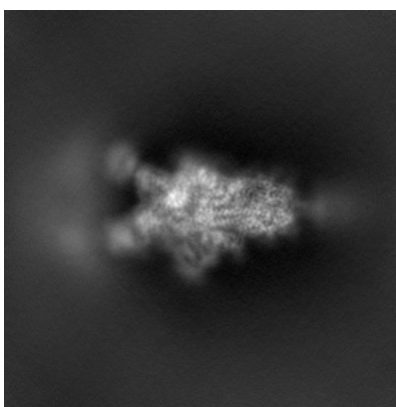


Z

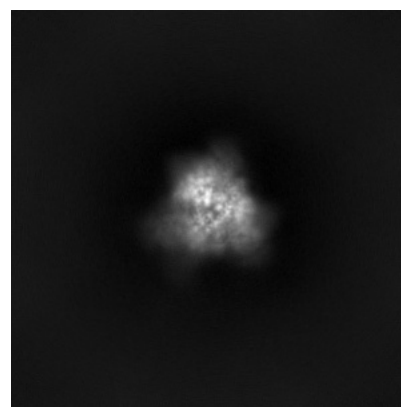
6.1.2 Raw map



X



Y

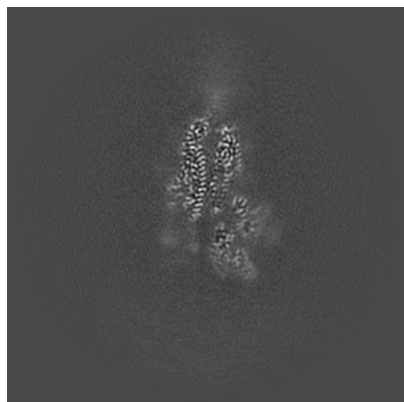


Z

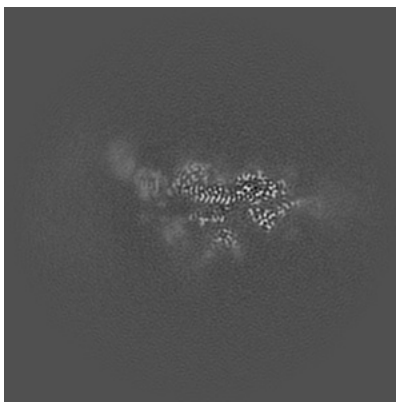
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

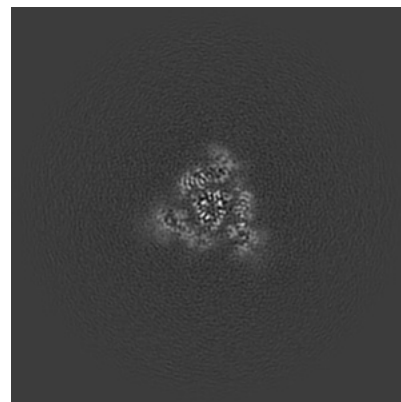
6.2.1 Primary map



X Index: 200

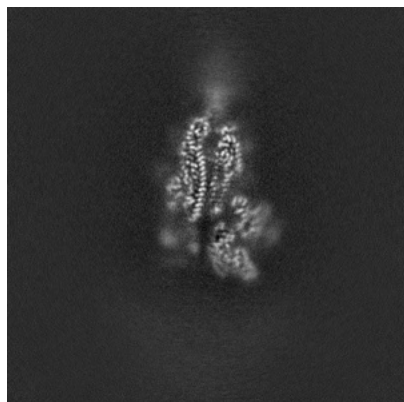


Y Index: 200

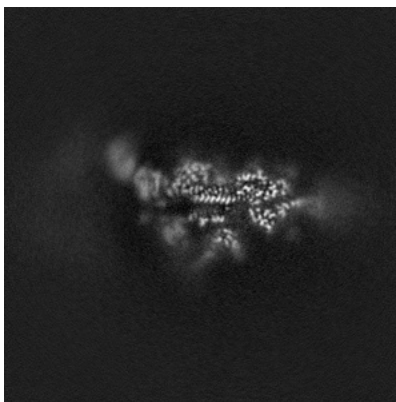


Z Index: 200

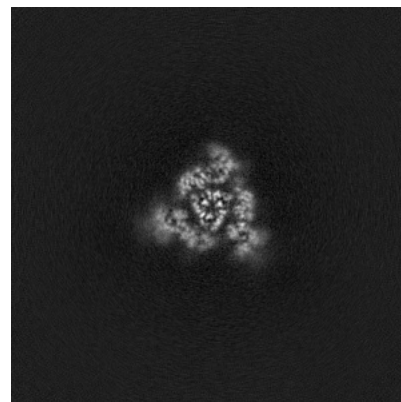
6.2.2 Raw map



X Index: 200



Y Index: 200

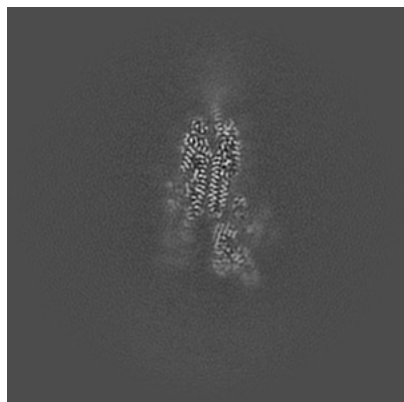


Z Index: 200

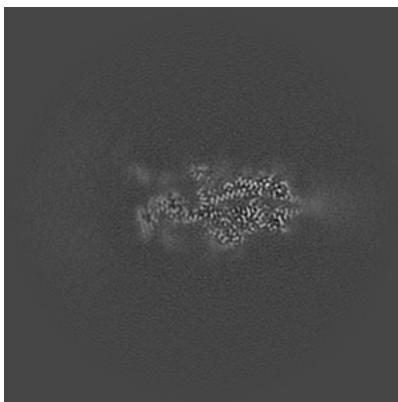
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

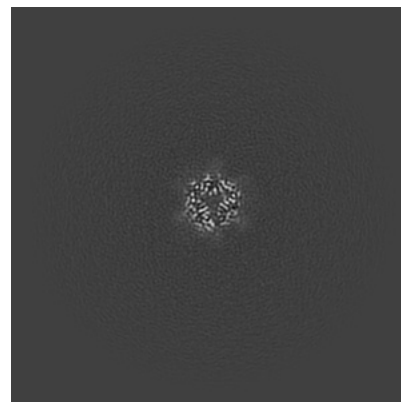
6.3.1 Primary map



X Index: 196

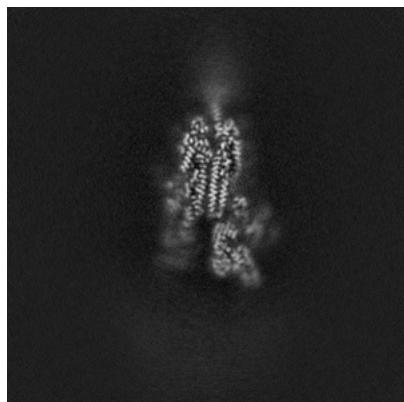


Y Index: 211

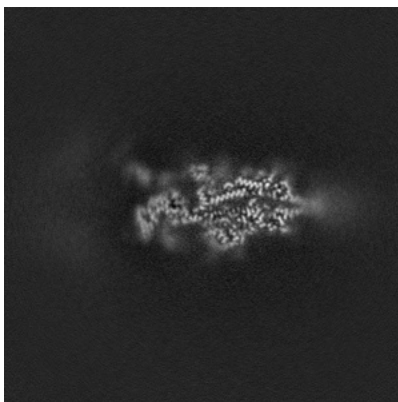


Z Index: 264

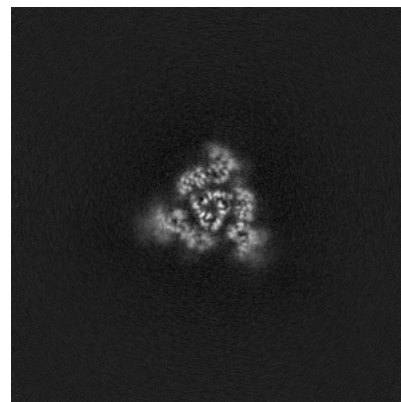
6.3.2 Raw map



X Index: 196



Y Index: 211

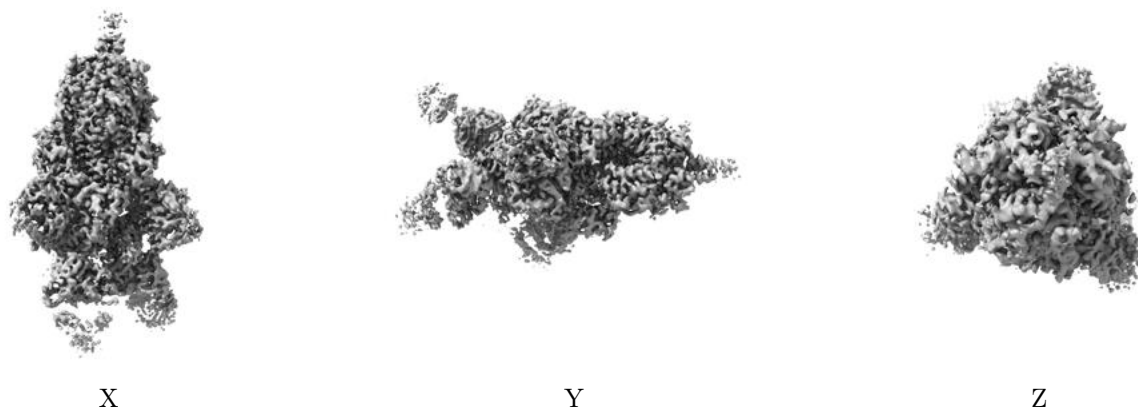


Z Index: 199

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

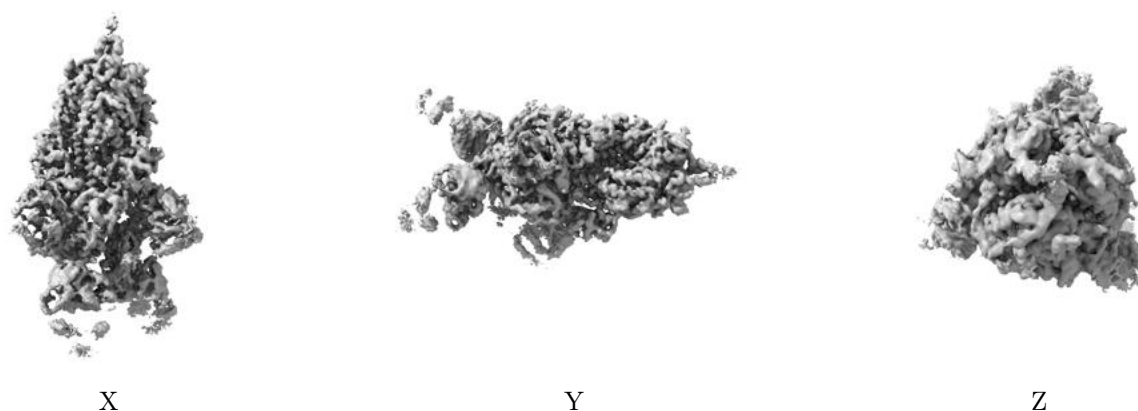
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.66. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 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.

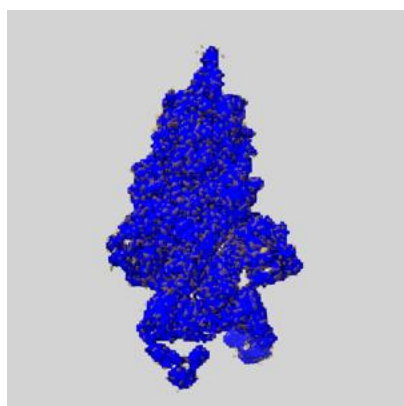
6.5 Mask visualisation [i](#)

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

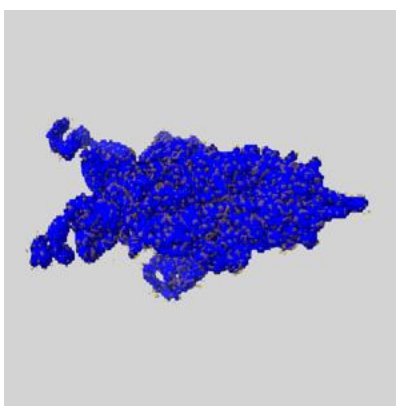
A mask typically either:

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

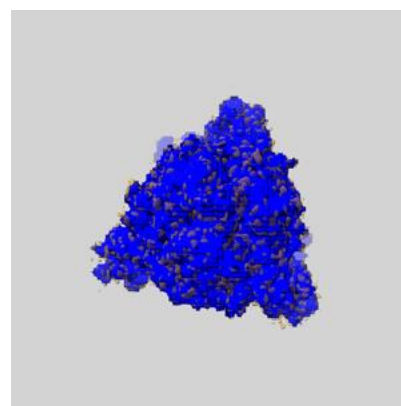
6.5.1 emd_26262_msk_1.map [i](#)



X



Y

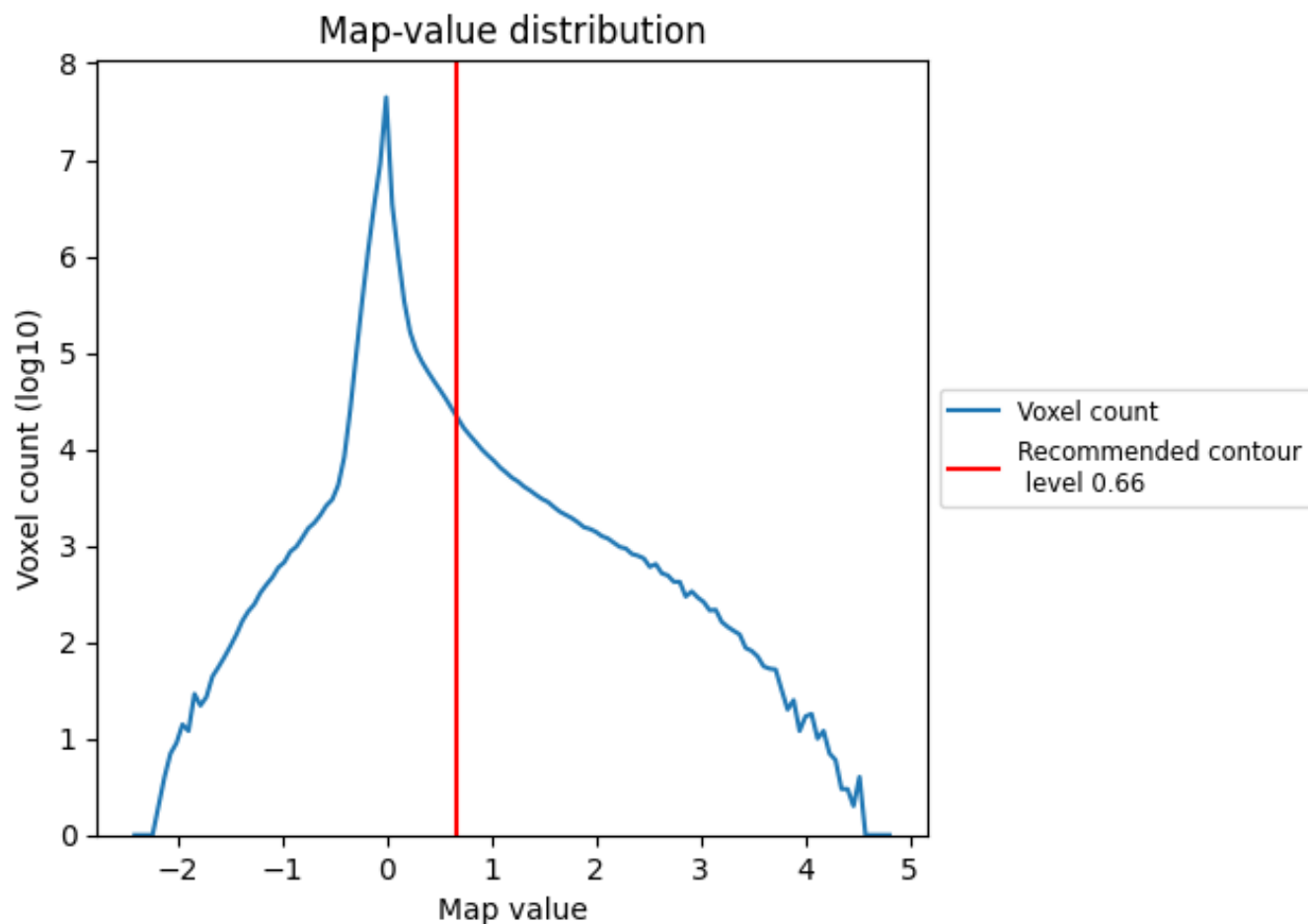


Z

7 Map analysis [i](#)

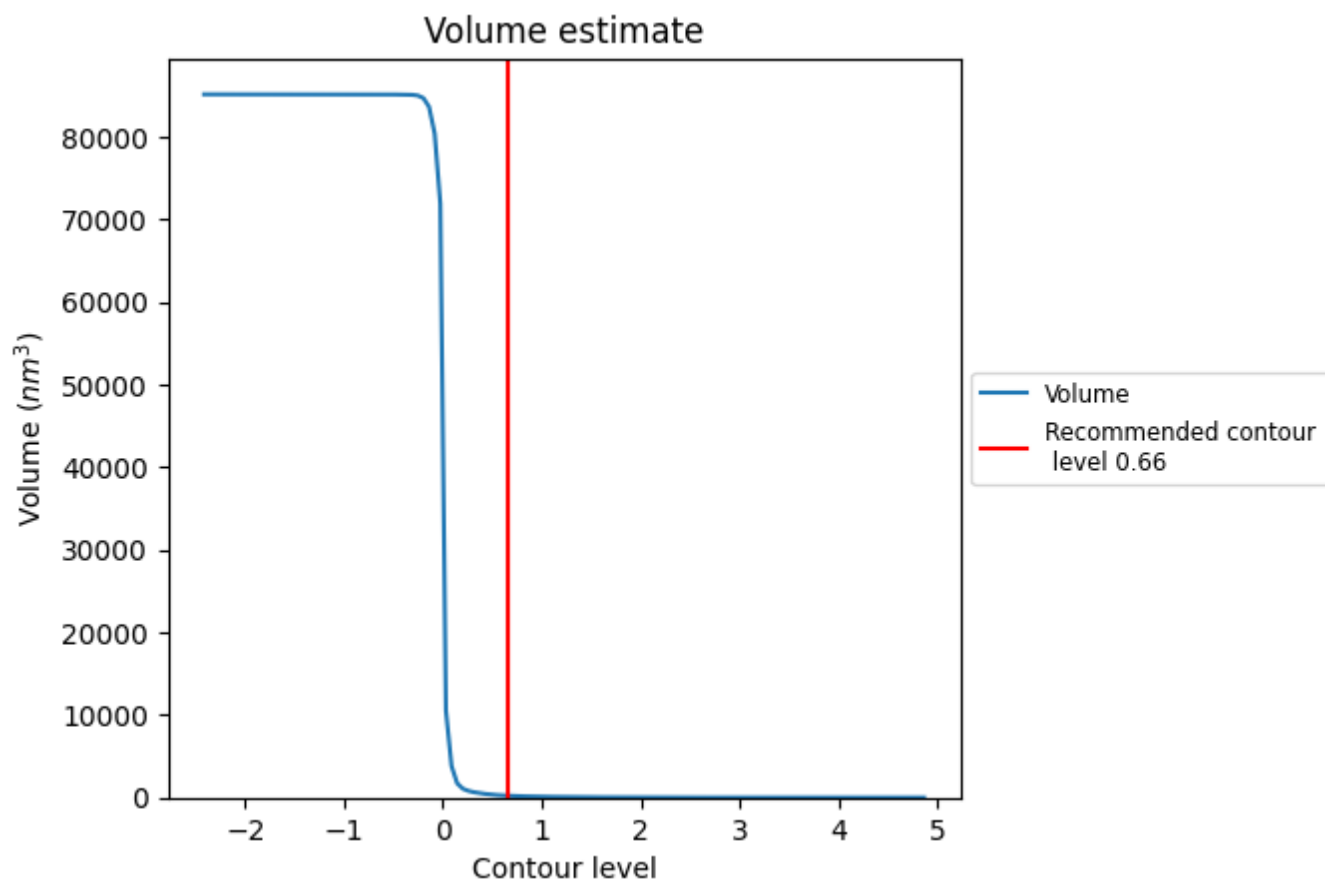
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

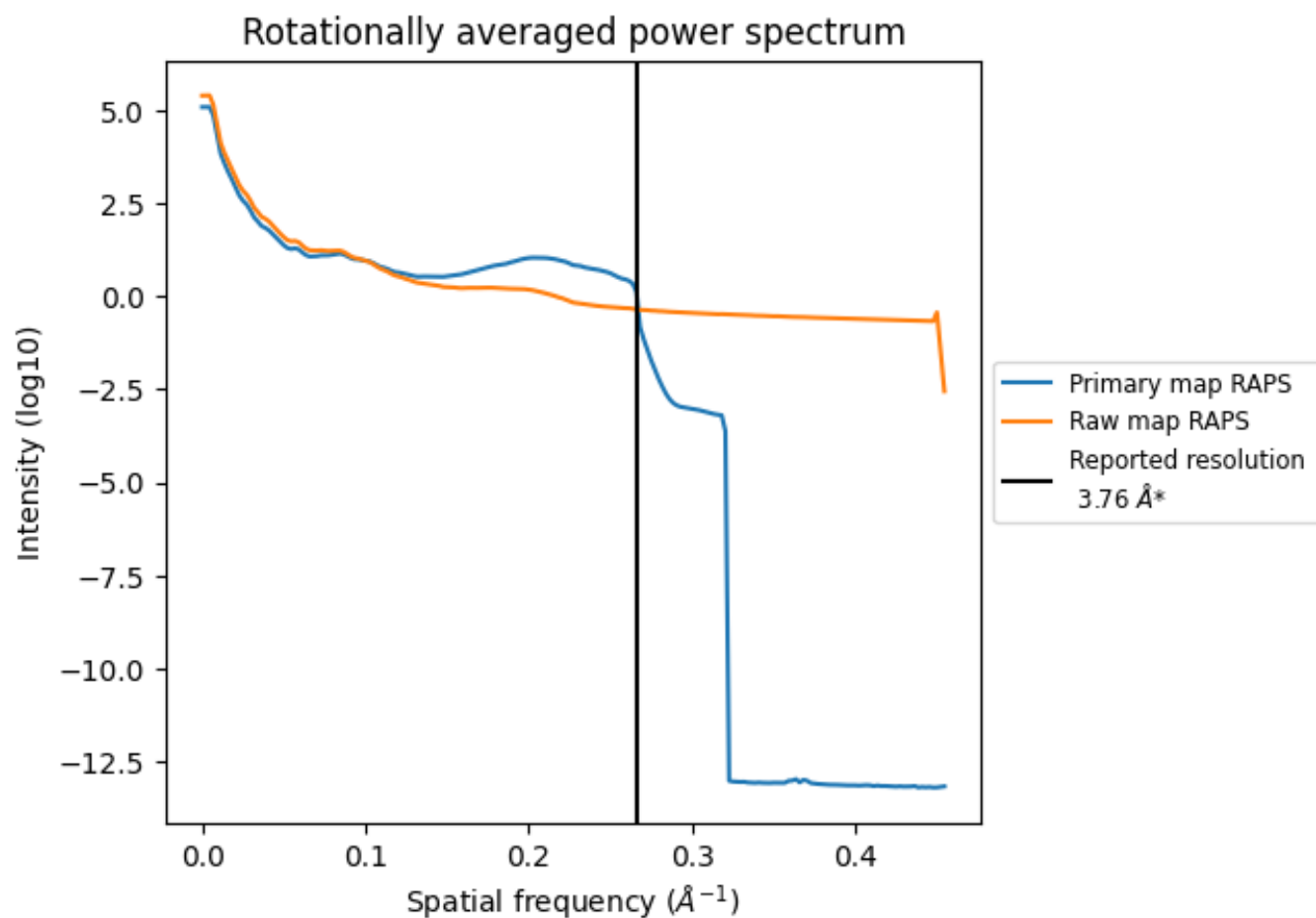
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 219 nm³; this corresponds to an approximate mass of 198 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 ⓘ

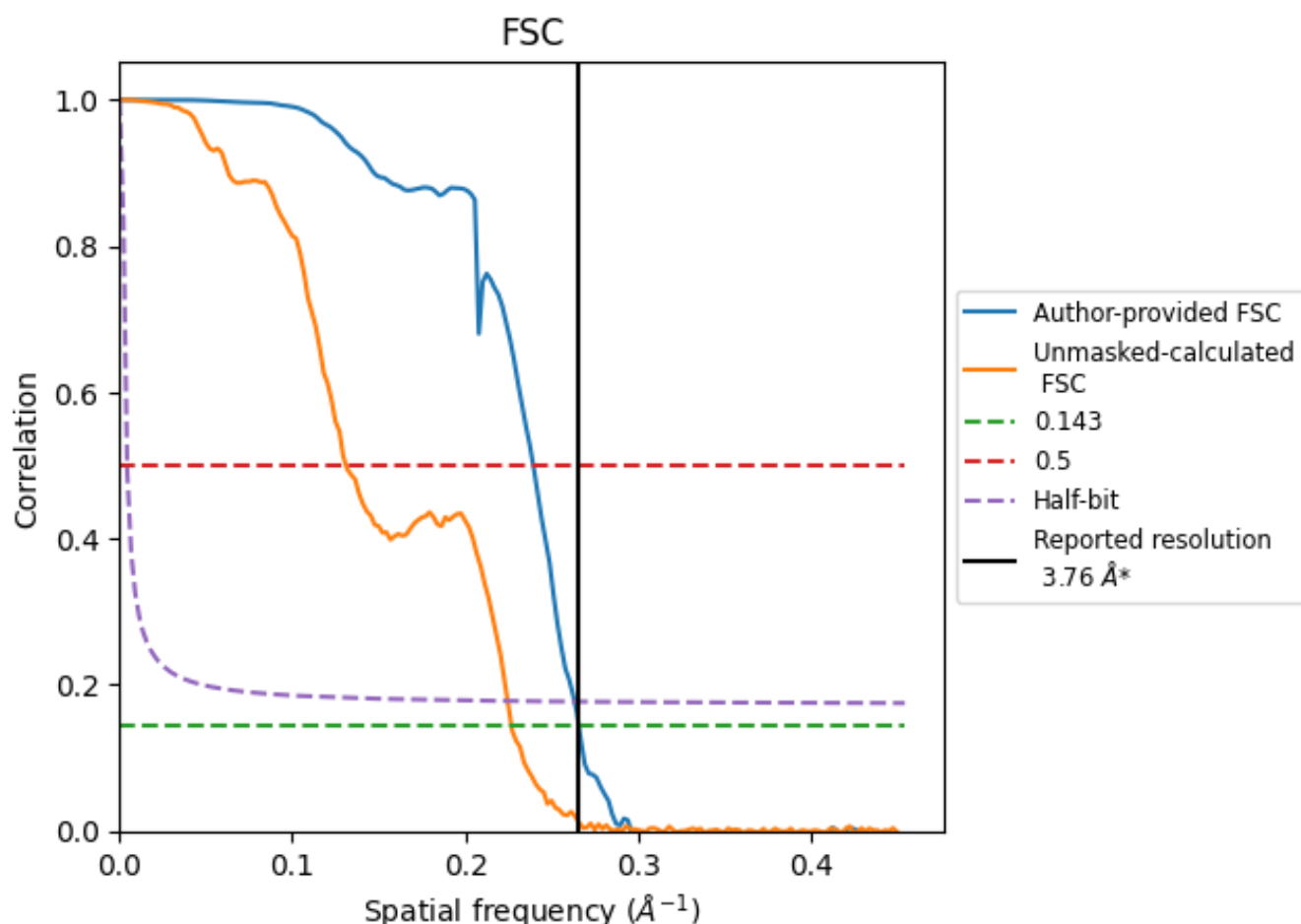


*Reported resolution corresponds to spatial frequency of 0.266 Å⁻¹

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 0.266 Å⁻¹

8.2 Resolution estimates [i](#)

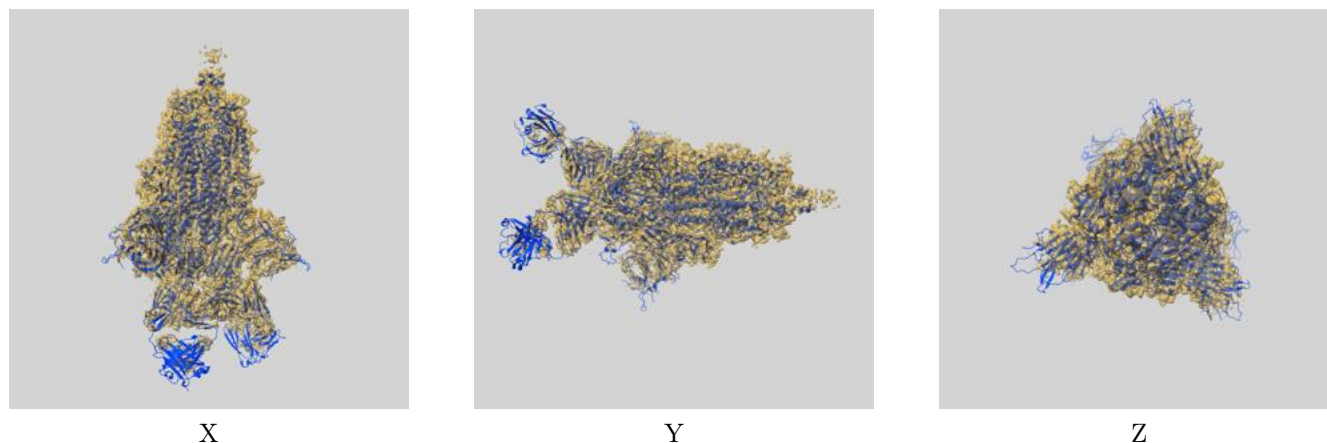
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.76	-	-
Author-provided FSC curve	3.76	4.18	3.80
Unmasked-calculated*	4.41	7.63	4.45

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.41 differs from the reported value 3.76 by more than 10 %

9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-26262 and PDB model 7U0P. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



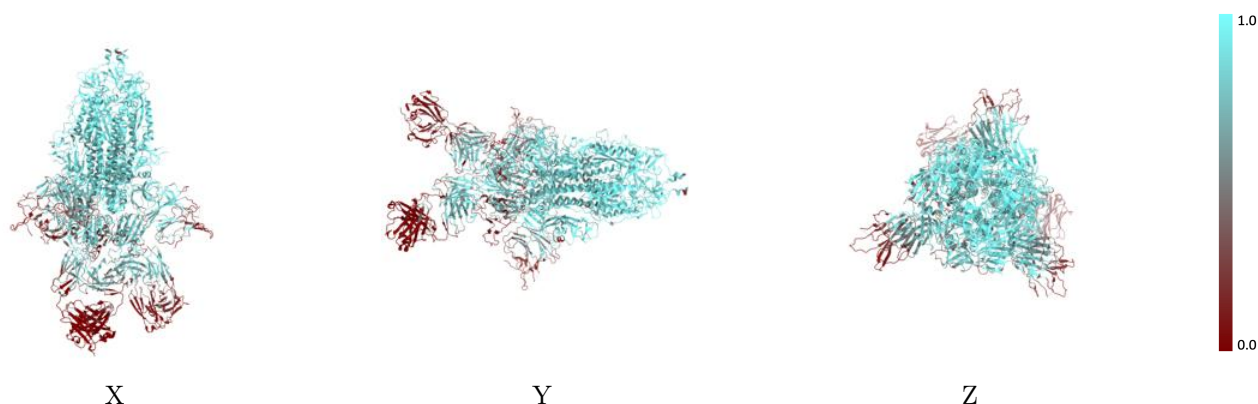
The images above show the 3D surface view of the map at the recommended contour level 0.66 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



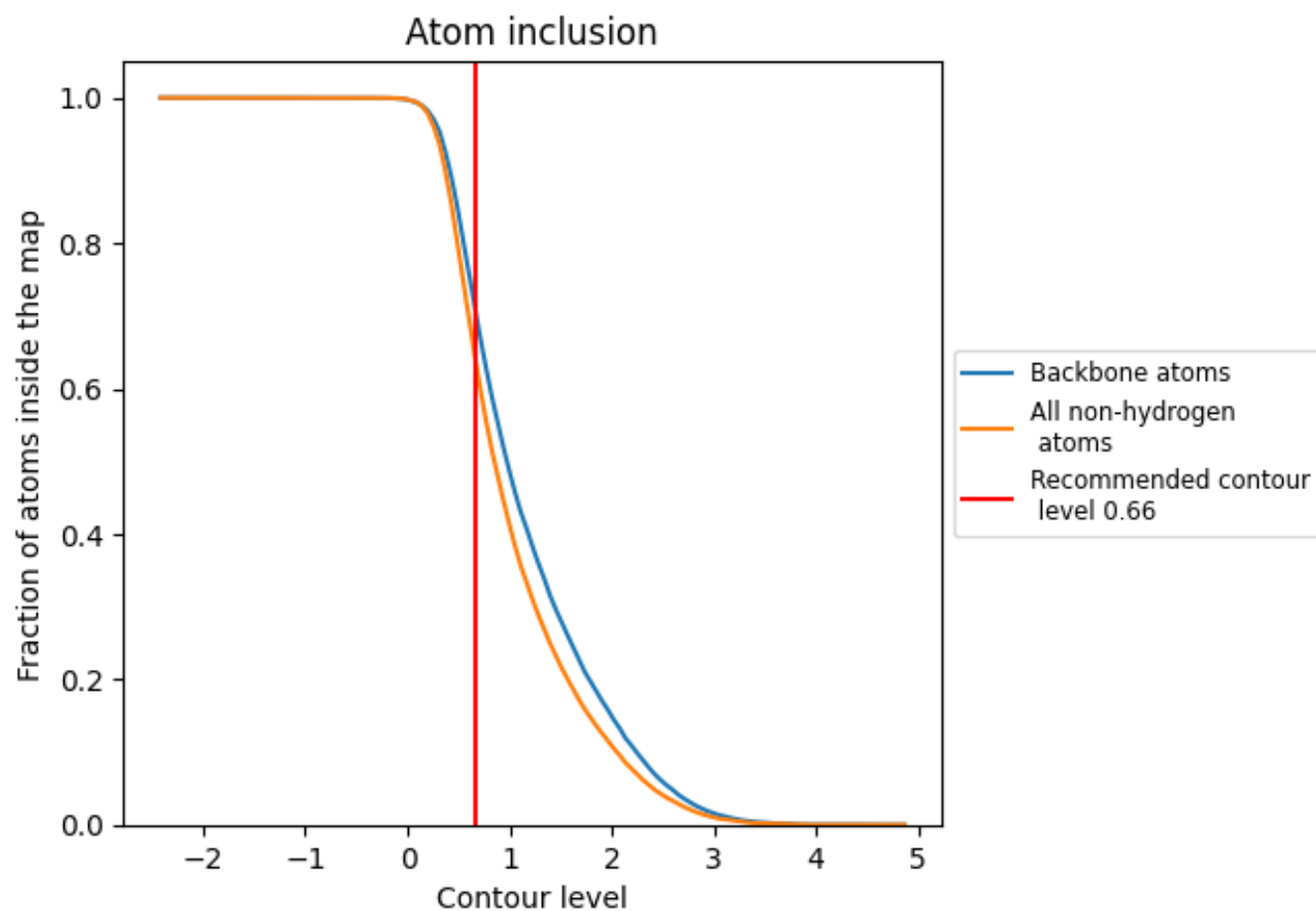
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.66).

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 64% 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 (0.66) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.6441	<div></div> 0.3660
A	<div></div> 0.6635	<div></div> 0.3970
B	<div></div> 0.7313	<div></div> 0.4210
B1	<div></div> 0.5357	<div></div> 0.3620
C	<div></div> 0.7708	<div></div> 0.4240
D	<div></div> 0.3926	<div></div> 0.1990
E	<div></div> 0.3281	<div></div> 0.1760
F	<div></div> 0.2977	<div></div> 0.1670
G	<div></div> 0.5357	<div></div> 0.2610
H	<div></div> 0.8214	<div></div> 0.4630
I	<div></div> 0.3728	<div></div> 0.1970
J	<div></div> 0.7143	<div></div> 0.4310
K	<div></div> 0.4643	<div></div> 0.2210

1.0

0.0

<0.0