



## Full wwPDB EM Validation Report ⓘ

Dec 18, 2022 – 06:19 pm GMT

PDB ID : 7ASD  
EMDB ID : EMD-11892  
Title : Structure of native royal jelly filaments  
Authors : Mattei, S.; Ban, A.; Piconi, A.; Leibundgut, M.; Glockshuber, R.;  
Boehringer, D.  
Deposited on : 2020-10-27  
Resolution : 3.50 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](mailto: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>.

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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 : 1.8.4, 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

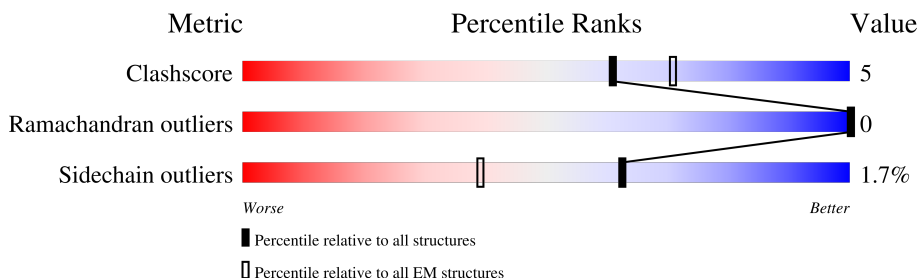
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.50 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  $\geq 3$ , 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	AA	432	
1	BA	432	
1	CA	432	
1	DA	432	
1	EA	432	
1	FA	432	
1	GA	432	
1	HA	432	

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Mol	Chain	Length	Quality of chain
2	AB	78	58% 42%
2	BB	78	56% 42%
2	CB	78	55% 42%
2	DB	78	55% 42%
2	EB	78	56% 42%
2	FB	78	58% 42%
2	GB	78	58% 42%
2	HB	78	56% 42%
3	AAc	2	50% 100%
3	BAc	2	50% 100%
3	CAc	2	50% 100%
3	DAc	2	50% 100%
3	EAc	2	50% 100%
3	FAc	2	50% 100%
3	GAc	2	50% 100%
3	HAc	2	50% 100%
4	A	3	33% 100%
4	B	3	100%
4	C	3	100%
4	D	3	33% 100%
4	E	3	100%
4	F	3	33% 100%
4	G	3	33% 100%
4	H	3	100%

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 30296 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 Major royal jelly protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	AA	413	3316	2084	562	653	17	2	0
1	BA	413	3316	2084	562	653	17	2	0
1	CA	413	3316	2084	562	653	17	2	0
1	DA	413	3316	2084	562	653	17	2	0
1	EA	413	3316	2084	562	653	17	2	0
1	FA	413	3316	2084	562	653	17	2	0
1	GA	413	3316	2084	562	653	17	2	0
1	HA	413	3316	2084	562	653	17	2	0

- Molecule 2 is a protein called Apisimin.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
2	AB	45	327	207	53	67	0	0
2	BB	45	327	207	53	67	0	0
2	CB	45	327	207	53	67	0	0
2	DB	45	327	207	53	67	0	0
2	EB	45	327	207	53	67	0	0
2	FB	45	327	207	53	67	0	0
2	GB	45	327	207	53	67	0	0

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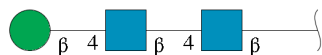
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
2	HB	45	327	207	53	67	0	0

- Molecule 3 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
			Total	C	N	O		
3	AAc	2	28	16	2	10	0	0
3	BAc	2	28	16	2	10	0	0
3	CAc	2	28	16	2	10	0	0
3	DAc	2	28	16	2	10	0	0
3	EAc	2	28	16	2	10	0	0
3	FAc	2	28	16	2	10	0	0
3	GAc	2	28	16	2	10	0	0
3	HAc	2	28	16	2	10	0	0

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



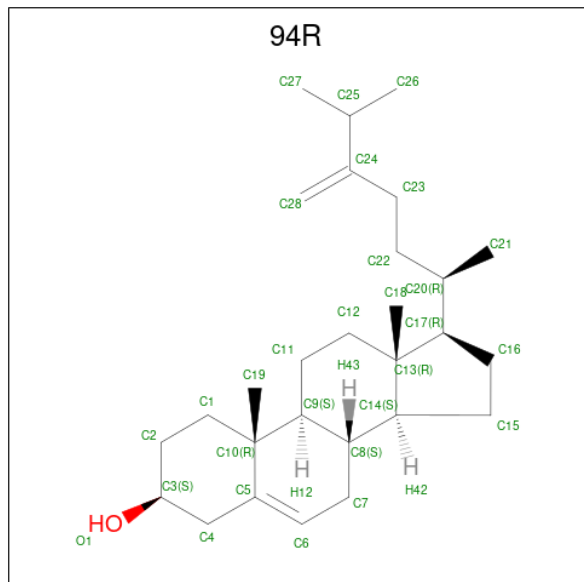
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
4	A	3	39	22	2	15	0	0
4	B	3	39	22	2	15	0	0
4	C	3	39	22	2	15	0	0

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Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
4	D	3	Total	C	N	O	0	0
			39	22	2	15		
4	E	3	Total	C	N	O	0	0
			39	22	2	15		
4	F	3	Total	C	N	O	0	0
			39	22	2	15		
4	G	3	Total	C	N	O	0	0
			39	22	2	15		
4	H	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 5 is (3beta,14beta,17alpha)-ergosta-5,24(28)-dien-3-ol (three-letter code: 94R) (formula: C<sub>28</sub>H<sub>46</sub>O).



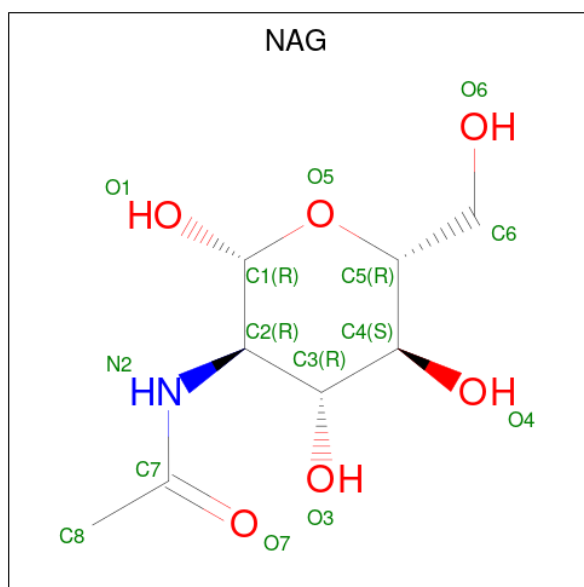
Mol	Chain	Residues	Atoms			AltConf
			Total	C	O	
5	AA	1	Total	C	O	0
			29	28	1	
5	AB	1	Total	C	O	0
			29	28	1	
5	BA	1	Total	C	O	0
			29	28	1	
5	BB	1	Total	C	O	0
			29	28	1	
5	CA	1	Total	C	O	0
			29	28	1	
5	CB	1	Total	C	O	0
			29	28	1	

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Mol	Chain	Residues	Atoms			AltConf
			Total	C	O	
5	DA	1	29	28	1	0
5	DB	1	29	28	1	0
5	EA	1	29	28	1	0
5	EB	1	29	28	1	0
5	FA	1	29	28	1	0
5	FB	1	29	28	1	0
5	GA	1	29	28	1	0
5	GB	1	29	28	1	0
5	HA	1	29	28	1	0
5	HB	1	29	28	1	0

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



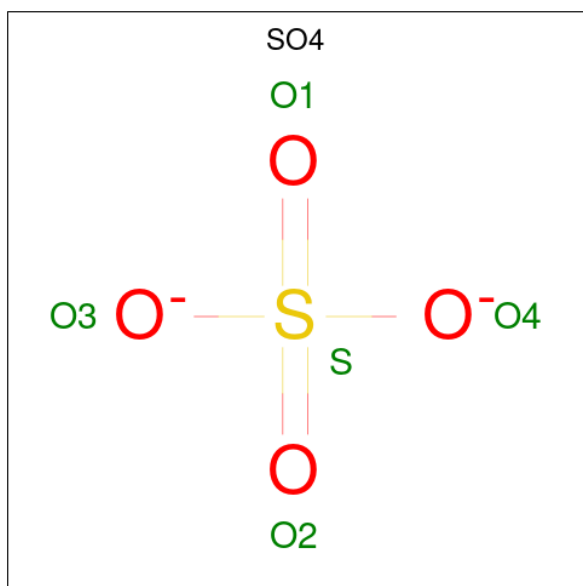
Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
6	AA	1	14	8	1	5	0

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Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
6	BA	1	Total 14	C 8	N 1	O 5	0
6	CA	1	Total 14	C 8	N 1	O 5	0
6	DA	1	Total 14	C 8	N 1	O 5	0
6	EA	1	Total 14	C 8	N 1	O 5	0
6	FA	1	Total 14	C 8	N 1	O 5	0
6	GA	1	Total 14	C 8	N 1	O 5	0
6	HA	1	Total 14	C 8	N 1	O 5	0

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			AltConf
			Total	O	S	
7	AA	1	Total 5	O 4	S 1	0
7	BA	1	Total 5	O 4	S 1	0
7	CA	1	Total 5	O 4	S 1	0
7	DA	1	Total 5	O 4	S 1	0

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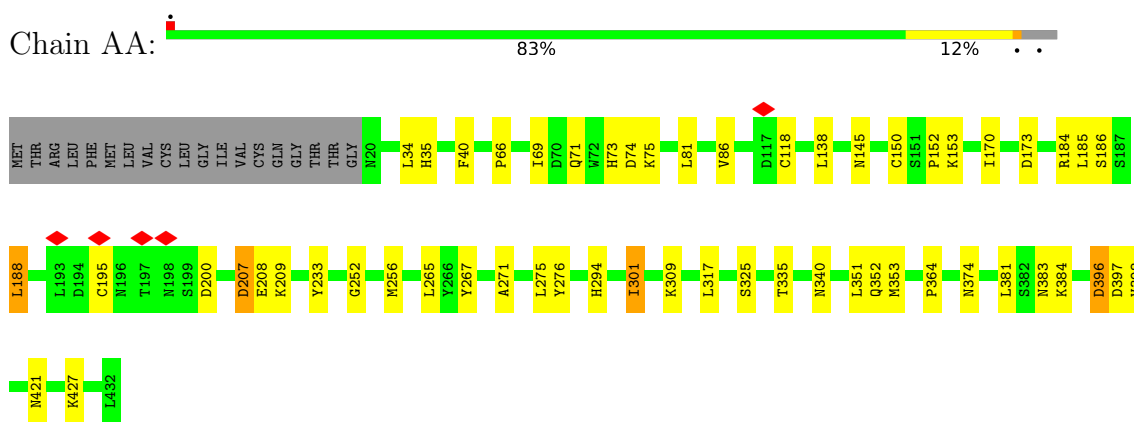
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>			<b>AltConf</b>
7	EA	1	Total 5	O 4	S 1	0
7	FA	1	Total 5	O 4	S 1	0
7	GA	1	Total 5	O 4	S 1	0
7	HA	1	Total 5	O 4	S 1	0

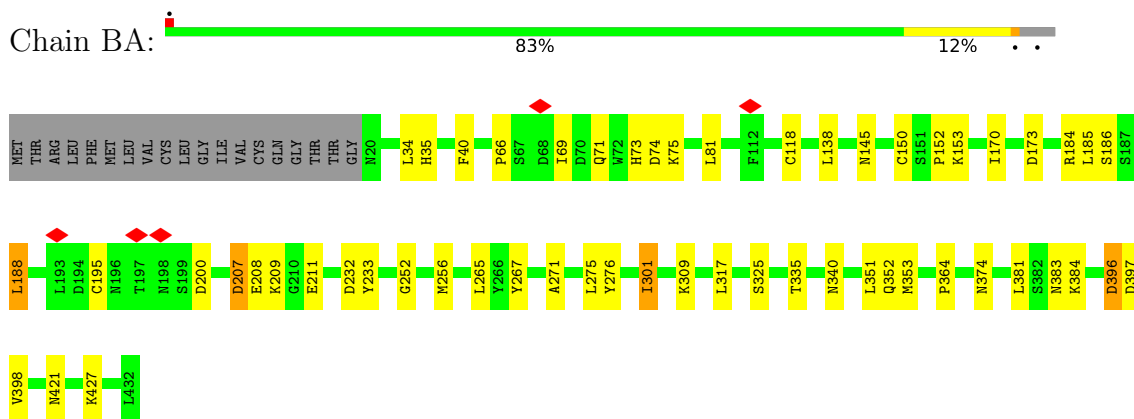
### 3 Residue-property plots [i](#)

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.

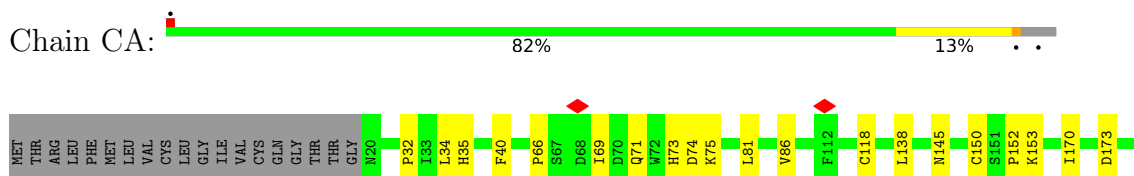
- Molecule 1: Major royal jelly protein 1

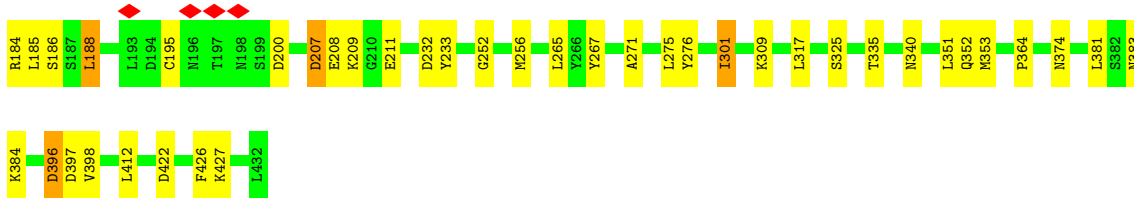


- Molecule 1: Major royal jelly protein 1

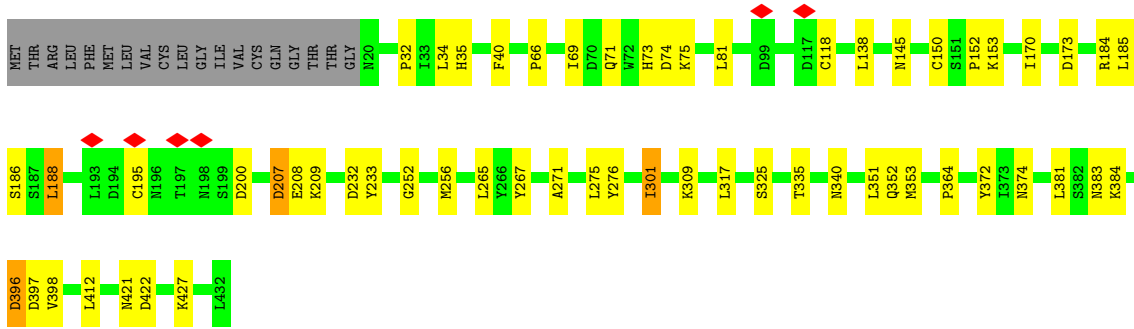
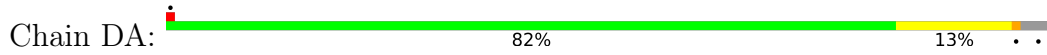


- Molecule 1: Major royal jelly protein 1

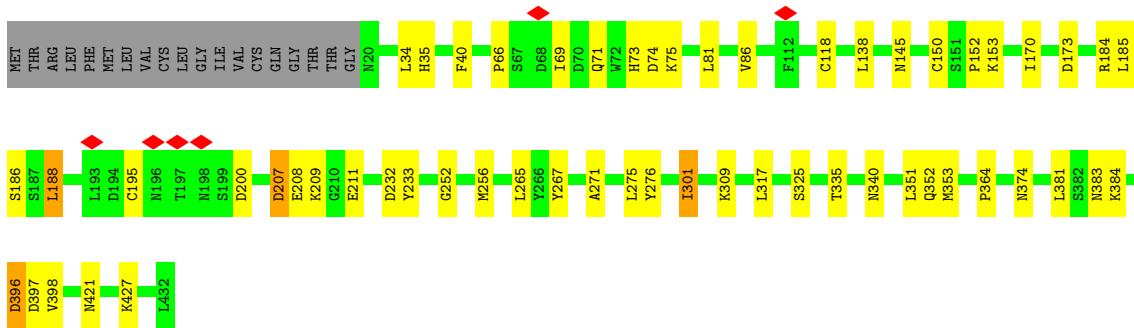
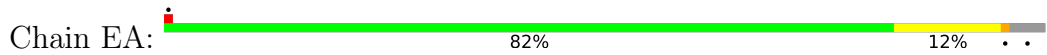




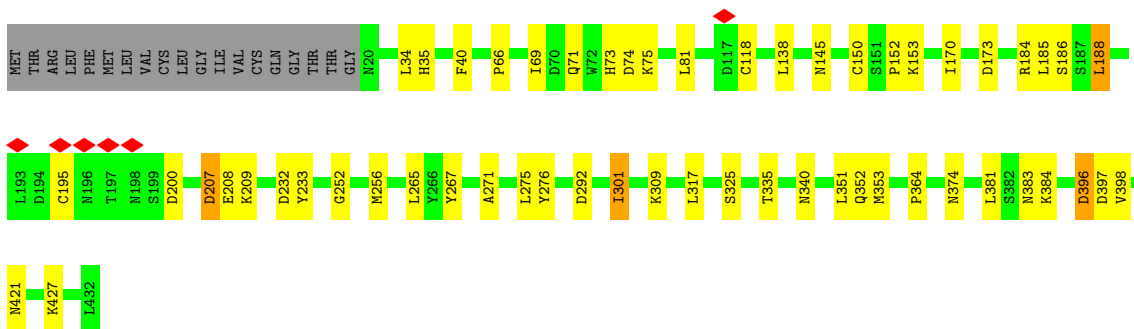
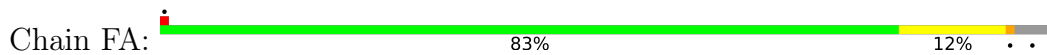
• Molecule 1: Major royal jelly protein 1



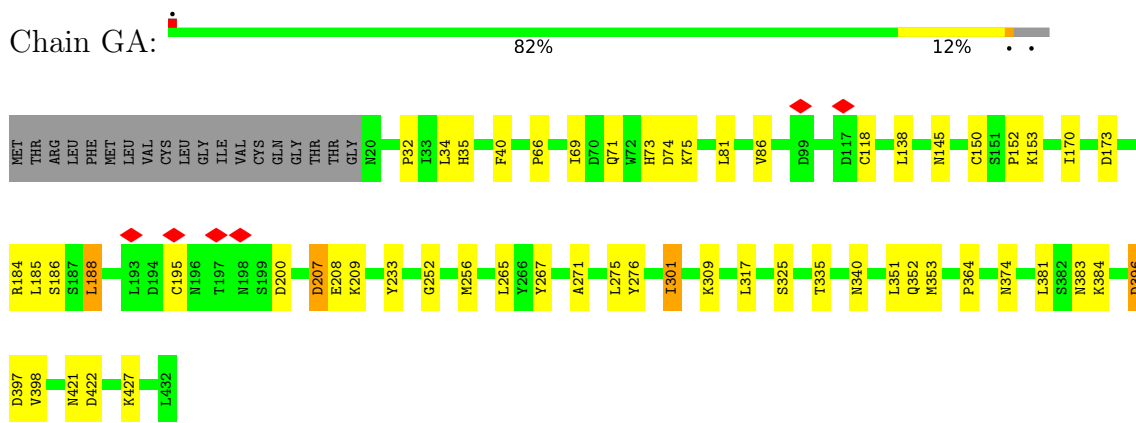
• Molecule 1: Major royal jelly protein 1



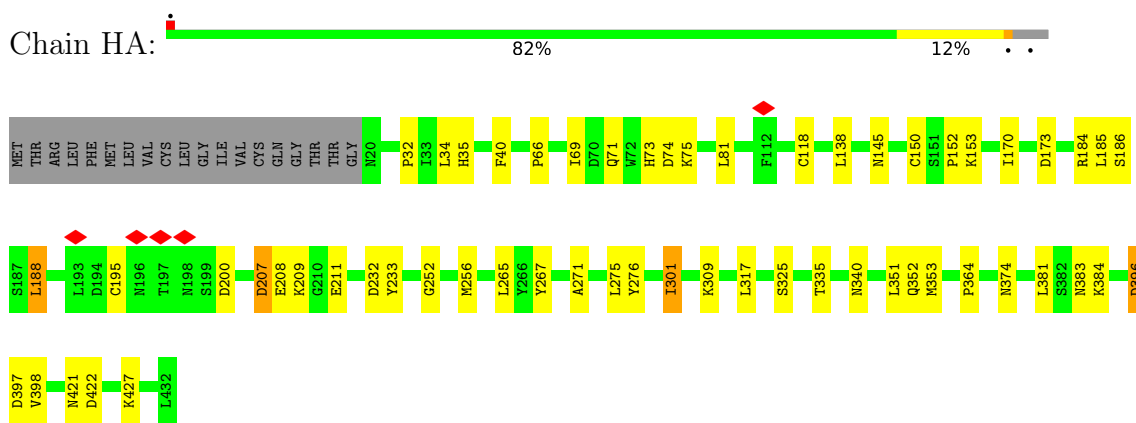
• Molecule 1: Major royal jelly protein 1



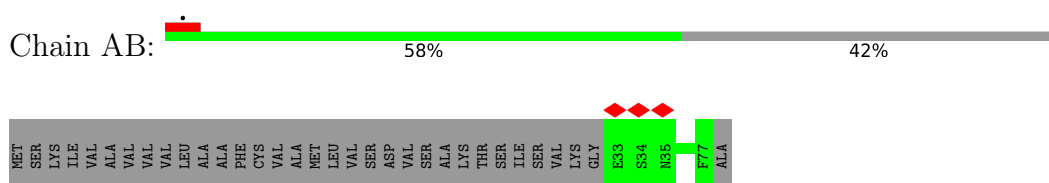
• Molecule 1: Major royal jelly protein 1



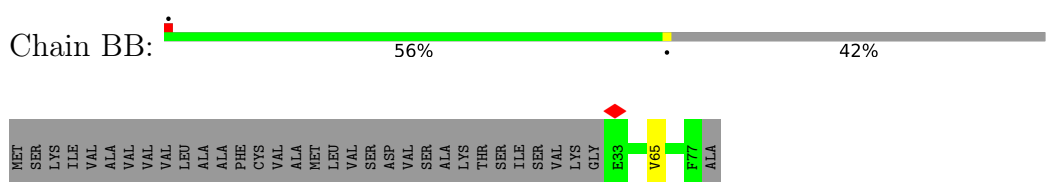
• Molecule 1: Major royal jelly protein 1



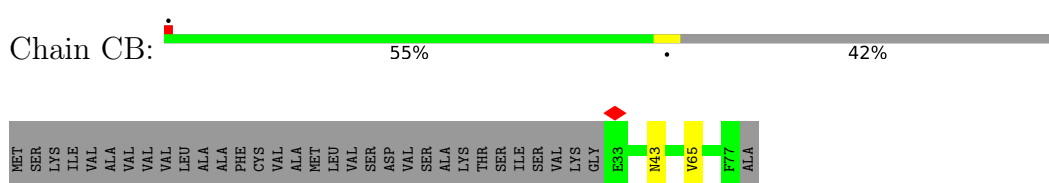
• Molecule 2: Apisimin



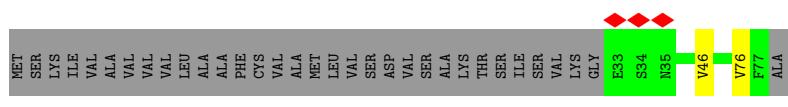
• Molecule 2: Apisimin



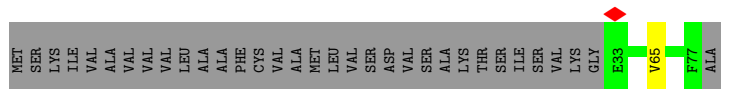
• Molecule 2: Apisimin



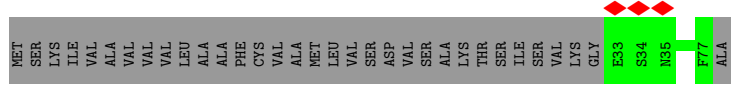
• Molecule 2: Apisimin



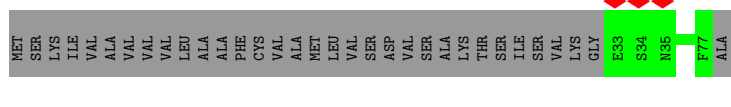
• Molecule 2: Apisimin



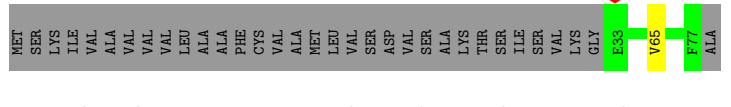
• Molecule 2: Apisimin



• Molecule 2: Apisimin



• Molecule 2: Apisimin



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain HAc:  50% 100%



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

Chain A:  33% 100%



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

Chain B:  100%



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

Chain C:  100%



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

Chain D:  33% 100%



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

Chain E:  100%



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

Chain F:  33% 100%



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

Chain G:  33% 100%



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

Chain H:  100%





## 4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=64°, rise=54 Å, axial sym=D2	Depositor
Number of segments used	240483	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{Å}^2$ )	82	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	119050	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.085	Depositor
Minimum map value	-0.035	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.015	Depositor
Map size (Å)	430.08118, 430.08118, 430.08118	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.19467, 1.19467, 1.19467	Depositor

## 5 Model quality i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, BMA, 94R, SO4

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	AA	0.24	0/3389	0.42	1/4598 (0.0%)
1	BA	0.24	0/3389	0.43	1/4598 (0.0%)
1	CA	0.24	0/3389	0.42	1/4598 (0.0%)
1	DA	0.24	0/3389	0.42	1/4598 (0.0%)
1	EA	0.24	0/3389	0.43	1/4598 (0.0%)
1	FA	0.24	0/3389	0.43	1/4598 (0.0%)
1	GA	0.24	0/3389	0.42	1/4598 (0.0%)
1	HA	0.24	0/3389	0.43	1/4598 (0.0%)
2	AB	0.23	0/327	0.33	0/448
2	BB	0.23	0/327	0.33	0/448
2	CB	0.23	0/327	0.33	0/448
2	DB	0.23	0/327	0.33	0/448
2	EB	0.23	0/327	0.33	0/448
2	FB	0.23	0/327	0.33	0/448
2	GB	0.23	0/327	0.33	0/448
2	HB	0.23	0/327	0.33	0/448
All	All	0.24	0/29728	0.42	8/40368 (0.0%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	FA	200	ASP	CB-CG-OD2	5.24	123.02	118.30
1	HA	200	ASP	CB-CG-OD2	5.23	123.01	118.30
1	CA	200	ASP	CB-CG-OD2	5.22	123.00	118.30
1	BA	200	ASP	CB-CG-OD2	5.22	122.99	118.30
1	AA	200	ASP	CB-CG-OD2	5.20	122.98	118.30
1	EA	200	ASP	CB-CG-OD2	5.17	122.96	118.30
1	DA	200	ASP	CB-CG-OD2	5.15	122.93	118.30
1	GA	200	ASP	CB-CG-OD2	5.14	122.92	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	AA	3316	0	3184	33	0
1	BA	3316	0	3184	33	0
1	CA	3316	0	3184	38	0
1	DA	3316	0	3184	36	0
1	EA	3316	0	3184	34	0
1	FA	3316	0	3184	33	0
1	GA	3316	0	3184	35	0
1	HA	3316	0	3184	35	0
2	AB	327	0	343	0	0
2	BB	327	0	343	1	0
2	CB	327	0	343	2	0
2	DB	327	0	343	2	0
2	EB	327	0	343	1	0
2	FB	327	0	343	0	0
2	GB	327	0	343	0	0
2	HB	327	0	343	1	0
3	AAc	28	0	25	0	0
3	BAc	28	0	25	0	0
3	CAc	28	0	25	0	0
3	DAc	28	0	25	0	0
3	EAc	28	0	25	0	0
3	FAc	28	0	25	0	0
3	GAc	28	0	25	0	0
3	HAc	28	0	25	0	0
4	A	39	0	34	0	0
4	B	39	0	34	0	0
4	C	39	0	34	0	0
4	D	39	0	34	0	0
4	E	39	0	34	0	0
4	F	39	0	34	0	0
4	G	39	0	34	0	0
4	H	39	0	34	0	0
5	AA	29	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	AB	29	0	0	0	0
5	BA	29	0	0	1	0
5	BB	29	0	0	0	0
5	CA	29	0	0	1	0
5	CB	29	0	0	0	0
5	DA	29	0	0	1	0
5	DB	29	0	0	0	0
5	EA	29	0	0	1	0
5	EB	29	0	0	0	0
5	FA	29	0	0	1	0
5	FB	29	0	0	0	0
5	GA	29	0	0	1	0
5	GB	29	0	0	0	0
5	HA	29	0	0	1	0
5	HB	29	0	0	0	0
6	AA	14	0	13	0	0
6	BA	14	0	13	0	0
6	CA	14	0	13	0	0
6	DA	14	0	13	0	0
6	EA	14	0	13	0	0
6	FA	14	0	13	0	0
6	GA	14	0	13	0	0
6	HA	14	0	13	0	0
7	AA	5	0	0	0	0
7	BA	5	0	0	0	0
7	CA	5	0	0	0	0
7	DA	5	0	0	0	0
7	EA	5	0	0	0	0
7	FA	5	0	0	0	0
7	GA	5	0	0	0	0
7	HA	5	0	0	0	0
All	All	30296	0	28792	273	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 (273) 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:EA:118:CYS:O	1:EA:153:LYS:NZ	2.27	0.68
1:CA:118:CYS:O	1:CA:153:LYS:NZ	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:GA:118:CYS:O	1:GA:153:LYS:NZ	2.27	0.67
1:DA:118:CYS:O	1:DA:153:LYS:NZ	2.27	0.67
1:AA:118:CYS:O	1:AA:153:LYS:NZ	2.27	0.67
1:FA:118:CYS:O	1:FA:153:LYS:NZ	2.27	0.67
1:BA:118:CYS:O	1:BA:153:LYS:NZ	2.27	0.67
1:HA:118:CYS:O	1:HA:153:LYS:NZ	2.27	0.67
1:FA:138:LEU:HB2	1:FA:188:LEU:HD13	1.78	0.65
1:BA:138:LEU:HB2	1:BA:188:LEU:HD13	1.78	0.65
1:CA:138:LEU:HB2	1:CA:188:LEU:HD13	1.78	0.65
1:DA:138:LEU:HB2	1:DA:188:LEU:HD13	1.78	0.65
1:EA:138:LEU:HB2	1:EA:188:LEU:HD13	1.78	0.65
1:HA:138:LEU:HB2	1:HA:188:LEU:HD13	1.78	0.65
1:AA:138:LEU:HB2	1:AA:188:LEU:HD13	1.78	0.65
1:GA:138:LEU:HB2	1:GA:188:LEU:HD13	1.78	0.65
1:CA:325:SER:OG	1:CA:351:LEU:O	2.15	0.64
1:EA:325:SER:OG	1:EA:351:LEU:O	2.15	0.64
1:AA:364:PRO:O	5:AA:501:94R:O1	2.17	0.62
1:GA:364:PRO:O	5:GA:501:94R:O1	2.17	0.62
1:BA:364:PRO:O	5:BA:501:94R:O1	2.17	0.62
1:HA:364:PRO:O	5:HA:501:94R:O1	2.17	0.62
1:DA:364:PRO:O	5:DA:501:94R:O1	2.17	0.62
1:DA:325:SER:OG	1:DA:351:LEU:O	2.15	0.61
1:FA:364:PRO:O	5:FA:501:94R:O1	2.17	0.61
1:EA:364:PRO:O	5:EA:501:94R:O1	2.17	0.61
1:FA:325:SER:OG	1:FA:351:LEU:O	2.15	0.61
1:BA:233:TYR:HB2	1:BA:271:ALA:HB3	1.83	0.61
1:FA:233:TYR:HB2	1:FA:271:ALA:HB3	1.83	0.61
1:CA:364:PRO:O	5:CA:501:94R:O1	2.17	0.61
1:HA:233:TYR:HB2	1:HA:271:ALA:HB3	1.83	0.61
1:DA:233:TYR:HB2	1:DA:271:ALA:HB3	1.83	0.61
1:GA:233:TYR:HB2	1:GA:271:ALA:HB3	1.83	0.61
1:AA:233:TYR:HB2	1:AA:271:ALA:HB3	1.83	0.61
1:CA:233:TYR:HB2	1:CA:271:ALA:HB3	1.83	0.60
1:HA:352:GLN:HG3	1:HA:384:LYS:HG2	1.83	0.60
1:BA:352:GLN:HG3	1:BA:384:LYS:HG2	1.83	0.60
1:EA:233:TYR:HB2	1:EA:271:ALA:HB3	1.83	0.60
1:AA:352:GLN:HG3	1:AA:384:LYS:HG2	1.83	0.60
1:GA:352:GLN:HG3	1:GA:384:LYS:HG2	1.83	0.60
1:HA:325:SER:OG	1:HA:351:LEU:O	2.15	0.60
1:BA:325:SER:OG	1:BA:351:LEU:O	2.15	0.60
1:FA:352:GLN:HG3	1:FA:384:LYS:HG2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DA:352:GLN:HG3	1:DA:384:LYS:HG2	1.83	0.59
1:AA:325:SER:OG	1:AA:351:LEU:O	2.15	0.59
1:GA:325:SER:OG	1:GA:351:LEU:O	2.15	0.59
1:CA:352:GLN:HG3	1:CA:384:LYS:HG2	1.83	0.59
1:EA:352:GLN:HG3	1:EA:384:LYS:HG2	1.83	0.59
1:BA:275:LEU:HB3	1:BA:301:ILE:HD11	1.86	0.58
1:FA:275:LEU:HB3	1:FA:301:ILE:HD11	1.86	0.58
1:GA:275:LEU:HB3	1:GA:301:ILE:HD11	1.86	0.58
1:HA:275:LEU:HB3	1:HA:301:ILE:HD11	1.86	0.58
1:AA:275:LEU:HB3	1:AA:301:ILE:HD11	1.86	0.58
2:EB:65:VAL:HG13	2:HB:65:VAL:HG13	1.86	0.58
1:CA:275:LEU:HB3	1:CA:301:ILE:HD11	1.86	0.58
1:DA:275:LEU:HB3	1:DA:301:ILE:HD11	1.86	0.58
1:EA:275:LEU:HB3	1:EA:301:ILE:HD11	1.86	0.58
1:BA:335:THR:O	1:BA:340:ASN:ND2	2.32	0.57
1:HA:335:THR:O	1:HA:340:ASN:ND2	2.33	0.57
1:GA:256:MET:HB3	1:GA:265:LEU:HD11	1.88	0.56
1:AA:256:MET:HB3	1:AA:265:LEU:HD11	1.88	0.56
1:AA:335:THR:O	1:AA:340:ASN:ND2	2.33	0.56
1:GA:335:THR:O	1:GA:340:ASN:ND2	2.33	0.56
1:EA:256:MET:HB3	1:EA:265:LEU:HD11	1.88	0.56
1:HA:256:MET:HB3	1:HA:265:LEU:HD11	1.88	0.55
1:AA:184:ARG:H	1:AA:209:LYS:HB2	1.71	0.55
1:BA:256:MET:HB3	1:BA:265:LEU:HD11	1.88	0.55
1:GA:184:ARG:H	1:GA:209:LYS:HB2	1.72	0.55
1:CA:256:MET:HB3	1:CA:265:LEU:HD11	1.88	0.55
1:GA:309:LYS:HB3	1:GA:317:LEU:HD11	1.88	0.55
1:AA:309:LYS:HB3	1:AA:317:LEU:HD11	1.88	0.55
1:DA:256:MET:HB3	1:DA:265:LEU:HD11	1.88	0.55
1:FA:256:MET:HB3	1:FA:265:LEU:HD11	1.88	0.55
1:BA:309:LYS:HB3	1:BA:317:LEU:HD11	1.88	0.54
1:HA:309:LYS:HB3	1:HA:317:LEU:HD11	1.88	0.54
1:DA:335:THR:O	1:DA:340:ASN:ND2	2.32	0.54
1:FA:335:THR:O	1:FA:340:ASN:ND2	2.32	0.54
1:DA:184:ARG:H	1:DA:209:LYS:HB2	1.71	0.54
1:FA:374:ASN:OD1	1:FA:427:LYS:NZ	2.35	0.54
1:BA:184:ARG:H	1:BA:209:LYS:HB2	1.71	0.54
1:BA:184:ARG:NH1	1:BA:208:GLU:OE1	2.41	0.54
1:FA:184:ARG:H	1:FA:209:LYS:HB2	1.72	0.54
1:HA:184:ARG:H	1:HA:209:LYS:HB2	1.71	0.54
1:AA:397:ASP:OD1	1:AA:398:VAL:N	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:309:LYS:HB3	1:CA:317:LEU:HD11	1.88	0.54
1:HA:184:ARG:NH1	1:HA:208:GLU:OE1	2.41	0.54
1:CA:184:ARG:H	1:CA:209:LYS:HB2	1.72	0.53
1:EA:309:LYS:HB3	1:EA:317:LEU:HD11	1.88	0.53
1:GA:397:ASP:OD1	1:GA:398:VAL:N	2.38	0.53
1:HA:185:LEU:HD23	1:HA:207:ASP:HB3	1.91	0.53
1:EA:184:ARG:H	1:EA:209:LYS:HB2	1.71	0.53
1:BA:185:LEU:HD23	1:BA:207:ASP:HB3	1.91	0.53
1:EA:335:THR:O	1:EA:340:ASN:ND2	2.33	0.53
1:BA:81:LEU:HD13	1:BA:145:ASN:HD21	1.73	0.53
1:DA:185:LEU:HD23	1:DA:207:ASP:HB3	1.91	0.53
1:FA:185:LEU:HD23	1:FA:207:ASP:HB3	1.91	0.53
1:DA:309:LYS:HB3	1:DA:317:LEU:HD11	1.88	0.53
1:EA:397:ASP:OD1	1:EA:398:VAL:N	2.38	0.53
1:FA:309:LYS:HB3	1:FA:317:LEU:HD11	1.88	0.53
1:GA:81:LEU:HD13	1:GA:145:ASN:HD21	1.73	0.53
1:AA:81:LEU:HD13	1:AA:145:ASN:HD21	1.73	0.53
1:CA:335:THR:O	1:CA:340:ASN:ND2	2.33	0.53
1:FA:81:LEU:HD13	1:FA:145:ASN:HD21	1.73	0.53
1:GA:184:ARG:NH1	1:GA:208:GLU:OE1	2.41	0.53
1:DA:184:ARG:NH1	1:DA:208:GLU:OE1	2.41	0.53
1:EA:185:LEU:HD23	1:EA:207:ASP:HB3	1.91	0.53
1:FA:184:ARG:NH1	1:FA:208:GLU:OE1	2.41	0.53
1:GA:185:LEU:HD23	1:GA:207:ASP:HB3	1.91	0.53
1:HA:81:LEU:HD13	1:HA:145:ASN:HD21	1.73	0.53
1:AA:184:ARG:NH1	1:AA:208:GLU:OE1	2.41	0.53
1:AA:185:LEU:HD23	1:AA:207:ASP:HB3	1.91	0.53
1:CA:397:ASP:OD1	1:CA:398:VAL:N	2.38	0.53
1:DA:81:LEU:HD13	1:DA:145:ASN:HD21	1.73	0.53
1:EA:374:ASN:OD1	1:EA:427:LYS:NZ	2.35	0.53
1:CA:185:LEU:HD23	1:CA:207:ASP:HB3	1.91	0.53
1:CA:184:ARG:NH1	1:CA:208:GLU:OE1	2.41	0.52
1:BA:397:ASP:OD1	1:BA:398:VAL:N	2.38	0.52
1:HA:397:ASP:OD1	1:HA:398:VAL:N	2.38	0.52
1:EA:184:ARG:NH1	1:EA:208:GLU:OE1	2.41	0.52
1:CA:374:ASN:OD1	1:CA:427:LYS:NZ	2.35	0.52
1:AA:40:PHE:HE2	1:AA:66:PRO:HD3	1.75	0.52
2:BB:65:VAL:HG13	2:CB:65:VAL:HG13	1.91	0.52
1:GA:40:PHE:HE2	1:GA:66:PRO:HD3	1.75	0.52
1:EA:81:LEU:HD13	1:EA:145:ASN:HD21	1.73	0.51
1:CA:81:LEU:HD13	1:CA:145:ASN:HD21	1.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DA:397:ASP:OD1	1:DA:398:VAL:N	2.38	0.51
1:CA:40:PHE:HE2	1:CA:66:PRO:HD3	1.75	0.51
1:EA:40:PHE:HE2	1:EA:66:PRO:HD3	1.75	0.51
1:GA:396:ASP:N	1:GA:396:ASP:OD1	2.43	0.51
1:CA:152:PRO:HG2	1:CA:170:ILE:HB	1.93	0.51
1:FA:397:ASP:OD1	1:FA:398:VAL:N	2.38	0.51
1:AA:396:ASP:OD1	1:AA:396:ASP:N	2.43	0.51
1:EA:152:PRO:HG2	1:EA:170:ILE:HB	1.93	0.51
1:BA:40:PHE:HE2	1:BA:66:PRO:HD3	1.75	0.51
1:HA:396:ASP:OD1	1:HA:396:ASP:N	2.43	0.51
1:BA:396:ASP:N	1:BA:396:ASP:OD1	2.43	0.51
1:HA:40:PHE:HE2	1:HA:66:PRO:HD3	1.75	0.51
1:FA:396:ASP:OD1	1:FA:396:ASP:N	2.43	0.50
1:DA:40:PHE:HE2	1:DA:66:PRO:HD3	1.75	0.50
1:DA:396:ASP:OD1	1:DA:396:ASP:N	2.43	0.50
1:FA:40:PHE:HE2	1:FA:66:PRO:HD3	1.75	0.50
1:FA:152:PRO:HG2	1:FA:170:ILE:HB	1.93	0.49
1:GA:32:PRO:HG3	1:HA:422:ASP:HB2	1.93	0.49
1:BA:152:PRO:HG2	1:BA:170:ILE:HB	1.93	0.49
1:DA:152:PRO:HG2	1:DA:170:ILE:HB	1.93	0.49
1:GA:152:PRO:HG2	1:GA:170:ILE:HB	1.93	0.49
1:HA:152:PRO:HG2	1:HA:170:ILE:HB	1.93	0.49
1:AA:152:PRO:HG2	1:AA:170:ILE:HB	1.93	0.49
1:CA:396:ASP:OD1	1:CA:396:ASP:N	2.43	0.49
1:HA:374:ASN:OD1	1:HA:427:LYS:NZ	2.35	0.49
1:EA:396:ASP:N	1:EA:396:ASP:OD1	2.43	0.49
1:BA:374:ASN:OD1	1:BA:427:LYS:NZ	2.35	0.49
1:GA:422:ASP:HB2	1:HA:32:PRO:HG3	1.94	0.48
1:HA:69:ILE:HG23	1:HA:381:LEU:HD22	1.95	0.47
1:BA:69:ILE:HG23	1:BA:381:LEU:HD22	1.95	0.47
1:FA:69:ILE:HG23	1:FA:381:LEU:HD22	1.96	0.47
1:DA:69:ILE:HG23	1:DA:381:LEU:HD22	1.95	0.47
1:AA:374:ASN:OD1	1:AA:427:LYS:NZ	2.35	0.47
1:GA:69:ILE:HG23	1:GA:381:LEU:HD22	1.95	0.46
1:AA:69:ILE:HG23	1:AA:381:LEU:HD22	1.95	0.46
1:CA:69:ILE:HG23	1:CA:381:LEU:HD22	1.95	0.46
1:CA:364:PRO:HG3	2:DB:76:VAL:HG13	1.97	0.46
1:GA:374:ASN:OD1	1:GA:427:LYS:NZ	2.35	0.46
1:CA:412:LEU:HD23	1:DA:412:LEU:HD23	1.98	0.46
1:EA:69:ILE:HG23	1:EA:381:LEU:HD22	1.95	0.46
1:EA:208:GLU:HB2	1:EA:252:GLY:H	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:208:GLU:HB2	1:CA:252:GLY:H	1.81	0.46
1:HA:208:GLU:HB2	1:HA:252:GLY:H	1.81	0.46
1:AA:208:GLU:HB2	1:AA:252:GLY:H	1.81	0.46
1:BA:208:GLU:HB2	1:BA:252:GLY:H	1.81	0.46
1:GA:208:GLU:HB2	1:GA:252:GLY:H	1.81	0.46
1:FA:208:GLU:HB2	1:FA:252:GLY:H	1.81	0.45
1:DA:208:GLU:HB2	1:DA:252:GLY:H	1.81	0.45
1:FA:173:ASP:OD1	1:FA:173:ASP:N	2.49	0.45
1:DA:173:ASP:OD1	1:DA:173:ASP:N	2.49	0.44
1:HA:173:ASP:OD1	1:HA:173:ASP:N	2.49	0.44
1:BA:173:ASP:OD1	1:BA:173:ASP:N	2.49	0.44
1:AA:173:ASP:OD1	1:AA:173:ASP:N	2.49	0.44
1:CA:184:ARG:N	1:CA:209:LYS:HB2	2.33	0.44
1:EA:184:ARG:N	1:EA:209:LYS:HB2	2.33	0.44
1:GA:173:ASP:OD1	1:GA:173:ASP:N	2.49	0.44
1:FA:186:SER:HB3	1:FA:208:GLU:HG2	2.00	0.44
1:GA:184:ARG:N	1:GA:209:LYS:HB2	2.33	0.44
1:CA:71:GLN:HA	1:CA:75:LYS:O	2.18	0.44
1:DA:186:SER:HB3	1:DA:208:GLU:HG2	2.00	0.44
1:EA:71:GLN:HA	1:EA:75:LYS:O	2.18	0.44
1:DA:71:GLN:HA	1:DA:75:LYS:O	2.18	0.43
1:FA:71:GLN:HA	1:FA:75:LYS:O	2.18	0.43
1:AA:186:SER:HB3	1:AA:208:GLU:HG2	2.00	0.43
1:BA:186:SER:HB3	1:BA:208:GLU:HG2	2.00	0.43
2:CB:43:ASN:OD1	1:DA:372:TYR:OH	2.26	0.43
1:EA:145:ASN:HB3	1:EA:184:ARG:HH21	1.83	0.43
1:GA:186:SER:HB3	1:GA:208:GLU:HG2	2.00	0.43
1:HA:186:SER:HB3	1:HA:208:GLU:HG2	2.00	0.43
1:AA:145:ASN:HB3	1:AA:184:ARG:HH21	1.84	0.43
1:BA:71:GLN:HA	1:BA:75:LYS:O	2.18	0.43
1:BA:74:ASP:N	1:BA:74:ASP:OD1	2.52	0.43
1:CA:145:ASN:HB3	1:CA:184:ARG:HH21	1.84	0.43
1:HA:71:GLN:HA	1:HA:75:LYS:O	2.18	0.43
1:HA:74:ASP:OD1	1:HA:74:ASP:N	2.52	0.43
1:GA:145:ASN:HB3	1:GA:184:ARG:HH21	1.84	0.43
1:AA:71:GLN:HA	1:AA:75:LYS:O	2.18	0.43
1:BA:353:MET:HB3	1:BA:383:ASN:O	2.19	0.43
1:CA:186:SER:HB3	1:CA:208:GLU:HG2	2.00	0.43
1:EA:186:SER:HB3	1:EA:208:GLU:HG2	2.00	0.43
1:GA:353:MET:HB3	1:GA:383:ASN:O	2.19	0.43
1:HA:353:MET:HB3	1:HA:383:ASN:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:353:MET:HB3	1:AA:383:ASN:O	2.19	0.43
1:BA:184:ARG:N	1:BA:209:LYS:HB2	2.33	0.43
1:DA:353:MET:HB3	1:DA:383:ASN:O	2.19	0.43
1:FA:353:MET:HB3	1:FA:383:ASN:O	2.19	0.43
1:HA:184:ARG:N	1:HA:209:LYS:HB2	2.33	0.43
1:DA:145:ASN:HB3	1:DA:184:ARG:HH21	1.84	0.42
1:DA:184:ARG:N	1:DA:209:LYS:HB2	2.33	0.42
1:FA:184:ARG:N	1:FA:209:LYS:HB2	2.33	0.42
1:GA:71:GLN:HA	1:GA:75:LYS:O	2.18	0.42
1:AA:294[B]:HIS:NE2	1:FA:292:ASP:OD1	2.52	0.42
1:EA:353:MET:HB3	1:EA:383:ASN:O	2.19	0.42
1:FA:145:ASN:HB3	1:FA:184:ARG:HH21	1.83	0.42
1:CA:74:ASP:N	1:CA:74:ASP:OD1	2.52	0.42
1:CA:353:MET:HB3	1:CA:383:ASN:O	2.19	0.42
1:DA:374:ASN:OD1	1:DA:427:LYS:NZ	2.35	0.42
1:DA:421:ASN:OD1	1:DA:421:ASN:N	2.53	0.42
1:FA:421:ASN:OD1	1:FA:421:ASN:N	2.53	0.42
1:DA:74:ASP:N	1:DA:74:ASP:OD1	2.52	0.42
1:EA:74:ASP:N	1:EA:74:ASP:OD1	2.52	0.42
1:BA:145:ASN:HB3	1:BA:184:ARG:HH21	1.84	0.42
1:FA:74:ASP:OD1	1:FA:74:ASP:N	2.52	0.42
1:BA:421:ASN:OD1	1:BA:421:ASN:N	2.53	0.42
1:HA:145:ASN:HB3	1:HA:184:ARG:HH21	1.84	0.42
1:EA:232:ASP:OD1	1:EA:233:TYR:N	2.53	0.42
1:HA:421:ASN:N	1:HA:421:ASN:OD1	2.53	0.42
1:AA:421:ASN:OD1	1:AA:421:ASN:N	2.53	0.42
1:CA:232:ASP:OD1	1:CA:233:TYR:N	2.53	0.42
1:DA:267:TYR:CZ	1:DA:276:TYR:HB2	2.55	0.42
1:FA:267:TYR:CZ	1:FA:276:TYR:HB2	2.55	0.42
1:EA:421:ASN:OD1	1:EA:421:ASN:N	2.53	0.42
1:GA:421:ASN:OD1	1:GA:421:ASN:N	2.53	0.42
1:FA:232:ASP:OD1	1:FA:233:TYR:N	2.53	0.41
1:GA:74:ASP:N	1:GA:74:ASP:OD1	2.52	0.41
1:AA:74:ASP:OD1	1:AA:74:ASP:N	2.52	0.41
1:CA:34:LEU:HG	1:CA:35:HIS:CD2	2.56	0.41
1:CA:267:TYR:CZ	1:CA:276:TYR:HB2	2.55	0.41
1:EA:173:ASP:OD1	1:EA:173:ASP:N	2.49	0.41
1:DA:232:ASP:OD1	1:DA:233:TYR:N	2.53	0.41
1:EA:34:LEU:HG	1:EA:35:HIS:CD2	2.56	0.41
1:EA:267:TYR:CZ	1:EA:276:TYR:HB2	2.55	0.41
1:BA:34:LEU:HG	1:BA:35:HIS:CD2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:32:PRO:HG3	1:DA:422:ASP:HB2	2.02	0.41
1:DA:34:LEU:HG	1:DA:35:HIS:CD2	2.56	0.41
1:FA:34:LEU:HG	1:FA:35:HIS:CD2	2.56	0.41
1:HA:34:LEU:HG	1:HA:35:HIS:CD2	2.56	0.41
1:HA:267:TYR:CZ	1:HA:276:TYR:HB2	2.55	0.41
1:CA:173:ASP:OD1	1:CA:173:ASP:N	2.49	0.41
1:BA:267:TYR:CZ	1:BA:276:TYR:HB2	2.55	0.41
1:AA:267:TYR:CZ	1:AA:276:TYR:HB2	2.55	0.41
1:BA:211:GLU:HG2	1:BA:232:ASP:HA	2.03	0.41
1:GA:74:ASP:OD1	1:GA:75:LYS:HG3	2.21	0.41
1:HA:211:GLU:HG2	1:HA:232:ASP:HA	2.03	0.41
1:AA:74:ASP:OD1	1:AA:75:LYS:HG3	2.21	0.41
1:AA:34:LEU:HG	1:AA:35:HIS:CD2	2.56	0.40
1:BA:232:ASP:OD1	1:BA:233:TYR:N	2.53	0.40
1:GA:267:TYR:CZ	1:GA:276:TYR:HB2	2.55	0.40
1:EA:211:GLU:HG2	1:EA:232:ASP:HA	2.03	0.40
1:HA:232:ASP:OD1	1:HA:233:TYR:N	2.53	0.40
1:AA:81:LEU:O	1:AA:86:VAL:HG21	2.22	0.40
1:CA:211:GLU:HG2	1:CA:232:ASP:HA	2.02	0.40
1:CA:422:ASP:HB2	1:DA:32:PRO:HG3	2.03	0.40
1:EA:81:LEU:O	1:EA:86:VAL:HG21	2.22	0.40
1:CA:81:LEU:O	1:CA:86:VAL:HG21	2.22	0.40
1:GA:34:LEU:HG	1:GA:35:HIS:CD2	2.56	0.40
1:GA:81:LEU:O	1:GA:86:VAL:HG21	2.22	0.40
1:CA:426:PHE:CE1	2:DB:46:VAL:HG21	2.56	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	AA	413/432 (96%)	403 (98%)	10 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	BA	413/432 (96%)	403 (98%)	10 (2%)	0	100	100
1	CA	413/432 (96%)	403 (98%)	10 (2%)	0	100	100
1	DA	413/432 (96%)	403 (98%)	10 (2%)	0	100	100
1	EA	413/432 (96%)	403 (98%)	10 (2%)	0	100	100
1	FA	413/432 (96%)	403 (98%)	10 (2%)	0	100	100
1	GA	413/432 (96%)	403 (98%)	10 (2%)	0	100	100
1	HA	413/432 (96%)	403 (98%)	10 (2%)	0	100	100
2	AB	43/78 (55%)	43 (100%)	0	0	100	100
2	BB	43/78 (55%)	43 (100%)	0	0	100	100
2	CB	43/78 (55%)	43 (100%)	0	0	100	100
2	DB	43/78 (55%)	43 (100%)	0	0	100	100
2	EB	43/78 (55%)	43 (100%)	0	0	100	100
2	FB	43/78 (55%)	43 (100%)	0	0	100	100
2	GB	43/78 (55%)	43 (100%)	0	0	100	100
2	HB	43/78 (55%)	43 (100%)	0	0	100	100
All	All	3648/4080 (89%)	3568 (98%)	80 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	AA	377/391 (96%)	370 (98%)	7 (2%)	57	80
1	BA	377/391 (96%)	370 (98%)	7 (2%)	57	80
1	CA	377/391 (96%)	370 (98%)	7 (2%)	57	80
1	DA	377/391 (96%)	370 (98%)	7 (2%)	57	80
1	EA	377/391 (96%)	370 (98%)	7 (2%)	57	80
1	FA	377/391 (96%)	370 (98%)	7 (2%)	57	80

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	GA	377/391 (96%)	370 (98%)	7 (2%)	57	80
1	HA	377/391 (96%)	370 (98%)	7 (2%)	57	80
2	AB	40/66 (61%)	40 (100%)	0	100	100
2	BB	40/66 (61%)	40 (100%)	0	100	100
2	CB	40/66 (61%)	40 (100%)	0	100	100
2	DB	40/66 (61%)	40 (100%)	0	100	100
2	EB	40/66 (61%)	40 (100%)	0	100	100
2	FB	40/66 (61%)	40 (100%)	0	100	100
2	GB	40/66 (61%)	40 (100%)	0	100	100
2	HB	40/66 (61%)	40 (100%)	0	100	100
All	All	3336/3656 (91%)	3280 (98%)	56 (2%)	62	82

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AA	73	HIS
1	AA	150	CYS
1	AA	188	LEU
1	AA	195	CYS
1	AA	207	ASP
1	AA	301	ILE
1	AA	396	ASP
1	BA	73	HIS
1	BA	150	CYS
1	BA	188	LEU
1	BA	195	CYS
1	BA	207	ASP
1	BA	301	ILE
1	BA	396	ASP
1	CA	73	HIS
1	CA	150	CYS
1	CA	188	LEU
1	CA	195	CYS
1	CA	207	ASP
1	CA	301	ILE
1	CA	396	ASP
1	DA	73	HIS
1	DA	150	CYS

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Mol	Chain	Res	Type
1	DA	188	LEU
1	DA	195	CYS
1	DA	207	ASP
1	DA	301	ILE
1	DA	396	ASP
1	EA	73	HIS
1	EA	150	CYS
1	EA	188	LEU
1	EA	195	CYS
1	EA	207	ASP
1	EA	301	ILE
1	EA	396	ASP
1	FA	73	HIS
1	FA	150	CYS
1	FA	188	LEU
1	FA	195	CYS
1	FA	207	ASP
1	FA	301	ILE
1	FA	396	ASP
1	GA	73	HIS
1	GA	150	CYS
1	GA	188	LEU
1	GA	195	CYS
1	GA	207	ASP
1	GA	301	ILE
1	GA	396	ASP
1	HA	73	HIS
1	HA	150	CYS
1	HA	188	LEU
1	HA	195	CYS
1	HA	207	ASP
1	HA	301	ILE
1	HA	396	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	AA	35	HIS
1	AA	63	ASN
1	AA	64	ASN
1	AA	145	ASN
1	AA	299	GLN

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Mol	Chain	Res	Type
1	BA	35	HIS
1	BA	63	ASN
1	BA	64	ASN
1	BA	145	ASN
1	BA	299	GLN
1	CA	35	HIS
1	CA	63	ASN
1	CA	64	ASN
1	CA	145	ASN
1	CA	299	GLN
1	DA	35	HIS
1	DA	63	ASN
1	DA	64	ASN
1	DA	299	GLN
1	EA	35	HIS
1	EA	63	ASN
1	EA	64	ASN
1	EA	145	ASN
1	EA	299	GLN
1	FA	35	HIS
1	FA	63	ASN
1	FA	64	ASN
1	FA	299	GLN
1	GA	35	HIS
1	GA	63	ASN
1	GA	64	ASN
1	GA	145	ASN
1	GA	299	GLN
1	HA	35	HIS
1	HA	63	ASN
1	HA	64	ASN
1	HA	145	ASN
1	HA	299	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

40 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	A	1	4,1	14,14,15	0.22	0	17,19,21	0.43	0
4	NAG	A	2	4	14,14,15	0.25	0	17,19,21	0.42	0
4	BMA	A	3	4	11,11,12	0.56	0	15,15,17	0.76	0
3	NAG	AAc	1	3,1	14,14,15	0.24	0	17,19,21	0.55	0
3	NAG	AAc	2	3	14,14,15	0.21	0	17,19,21	0.46	0
4	NAG	B	1	4,1	14,14,15	0.23	0	17,19,21	0.43	0
4	NAG	B	2	4	14,14,15	0.23	0	17,19,21	0.41	0
4	BMA	B	3	4	11,11,12	0.56	0	15,15,17	0.76	0
3	NAG	BAc	1	3,1	14,14,15	0.26	0	17,19,21	0.55	0
3	NAG	BAc	2	3	14,14,15	0.22	0	17,19,21	0.45	0
4	NAG	C	1	4,1	14,14,15	0.23	0	17,19,21	0.43	0
4	NAG	C	2	4	14,14,15	0.23	0	17,19,21	0.42	0
4	BMA	C	3	4	11,11,12	0.56	0	15,15,17	0.77	0
3	NAG	CAc	1	3,1	14,14,15	0.26	0	17,19,21	0.56	0
3	NAG	CAc	2	3	14,14,15	0.22	0	17,19,21	0.45	0
4	NAG	D	1	4,1	14,14,15	0.24	0	17,19,21	0.43	0
4	NAG	D	2	4	14,14,15	0.25	0	17,19,21	0.42	0
4	BMA	D	3	4	11,11,12	0.57	0	15,15,17	0.77	0
3	NAG	DAc	1	3,1	14,14,15	0.25	0	17,19,21	0.56	0
3	NAG	DAc	2	3	14,14,15	0.22	0	17,19,21	0.45	0
4	NAG	E	1	4,1	14,14,15	0.24	0	17,19,21	0.43	0
4	NAG	E	2	4	14,14,15	0.25	0	17,19,21	0.42	0
4	BMA	E	3	4	11,11,12	0.56	0	15,15,17	0.77	0
3	NAG	EAc	1	3,1	14,14,15	0.25	0	17,19,21	0.55	0
3	NAG	EAc	2	3	14,14,15	0.21	0	17,19,21	0.46	0
4	NAG	F	1	4,1	14,14,15	0.23	0	17,19,21	0.43	0
4	NAG	F	2	4	14,14,15	0.24	0	17,19,21	0.42	0
4	BMA	F	3	4	11,11,12	0.56	0	15,15,17	0.77	0
3	NAG	FAc	1	3,1	14,14,15	0.24	0	17,19,21	0.55	0
3	NAG	FAc	2	3	14,14,15	0.22	0	17,19,21	0.45	0
4	NAG	G	1	4,1	14,14,15	0.23	0	17,19,21	0.43	0
4	NAG	G	2	4	14,14,15	0.22	0	17,19,21	0.42	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	BMA	G	3	4	11,11,12	0.56	0	15,15,17	0.77	0
3	NAG	GAc	1	3,1	14,14,15	0.26	0	17,19,21	0.56	0
3	NAG	GAc	2	3	14,14,15	0.21	0	17,19,21	0.44	0
4	NAG	H	1	4,1	14,14,15	0.23	0	17,19,21	0.43	0
4	NAG	H	2	4	14,14,15	0.24	0	17,19,21	0.42	0
4	BMA	H	3	4	11,11,12	0.56	0	15,15,17	0.77	0
3	NAG	HAc	1	3,1	14,14,15	0.26	0	17,19,21	0.56	0
3	NAG	HAc	2	3	14,14,15	0.21	0	17,19,21	0.46	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	A	2	4	-	0/6/23/26	0/1/1/1
4	BMA	A	3	4	-	0/2/19/22	0/1/1/1
3	NAG	AAc	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	AAc	2	3	-	2/6/23/26	0/1/1/1
4	NAG	B	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	B	2	4	-	0/6/23/26	0/1/1/1
4	BMA	B	3	4	-	0/2/19/22	0/1/1/1
3	NAG	BAc	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	BAc	2	3	-	2/6/23/26	0/1/1/1
4	NAG	C	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	C	2	4	-	0/6/23/26	0/1/1/1
4	BMA	C	3	4	-	0/2/19/22	0/1/1/1
3	NAG	CAC	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	CAC	2	3	-	2/6/23/26	0/1/1/1
4	NAG	D	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	0/6/23/26	0/1/1/1
4	BMA	D	3	4	-	0/2/19/22	0/1/1/1
3	NAG	DAC	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	DAC	2	3	-	2/6/23/26	0/1/1/1
4	NAG	E	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	0/2/19/22	0/1/1/1
3	NAG	EAC	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	EAC	2	3	-	2/6/23/26	0/1/1/1
4	NAG	F	1	4,1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1
4	BMA	F	3	4	-	0/2/19/22	0/1/1/1
3	NAG	FAc	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	FAc	2	3	-	2/6/23/26	0/1/1/1
4	NAG	G	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
4	BMA	G	3	4	-	0/2/19/22	0/1/1/1
3	NAG	GAc	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	GAc	2	3	-	2/6/23/26	0/1/1/1
4	NAG	H	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	H	2	4	-	0/6/23/26	0/1/1/1
4	BMA	H	3	4	-	0/2/19/22	0/1/1/1
3	NAG	HAc	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	HAc	2	3	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	AAc	2	NAG	O5-C5-C6-O6
3	BAc	2	NAG	O5-C5-C6-O6
3	CAc	2	NAG	O5-C5-C6-O6
3	DAc	2	NAG	O5-C5-C6-O6
3	EAc	2	NAG	O5-C5-C6-O6
3	FAc	2	NAG	O5-C5-C6-O6
3	GAc	2	NAG	O5-C5-C6-O6
3	HAc	2	NAG	O5-C5-C6-O6
3	AAc	2	NAG	C4-C5-C6-O6
3	BAc	2	NAG	C4-C5-C6-O6
3	CAc	2	NAG	C4-C5-C6-O6
3	DAc	2	NAG	C4-C5-C6-O6
3	EAc	2	NAG	C4-C5-C6-O6
3	FAc	2	NAG	C4-C5-C6-O6
3	GAc	2	NAG	C4-C5-C6-O6
3	HAc	2	NAG	C4-C5-C6-O6
3	AAc	1	NAG	C3-C2-N2-C7
3	BAc	1	NAG	C3-C2-N2-C7
3	CAc	1	NAG	C3-C2-N2-C7

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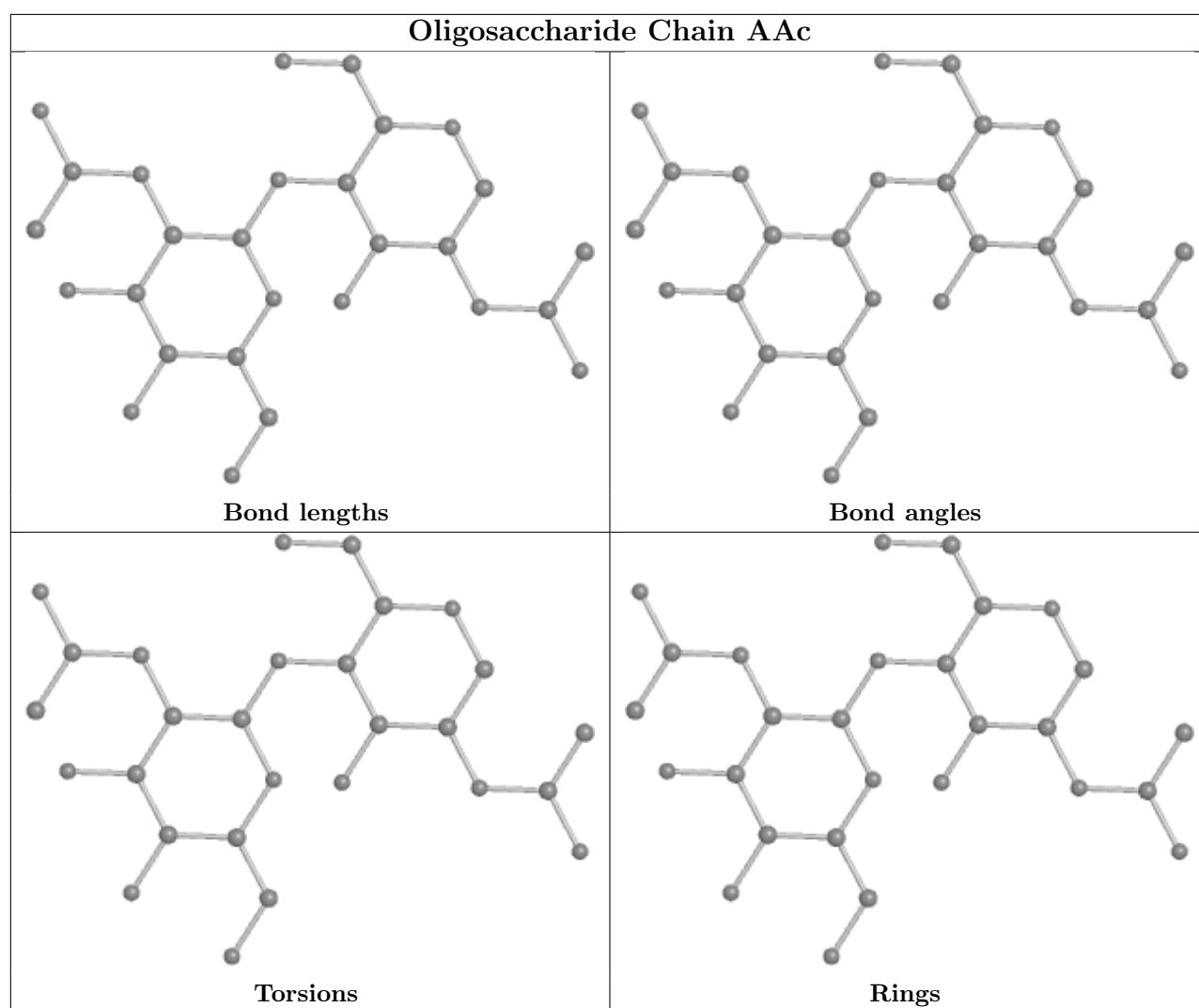
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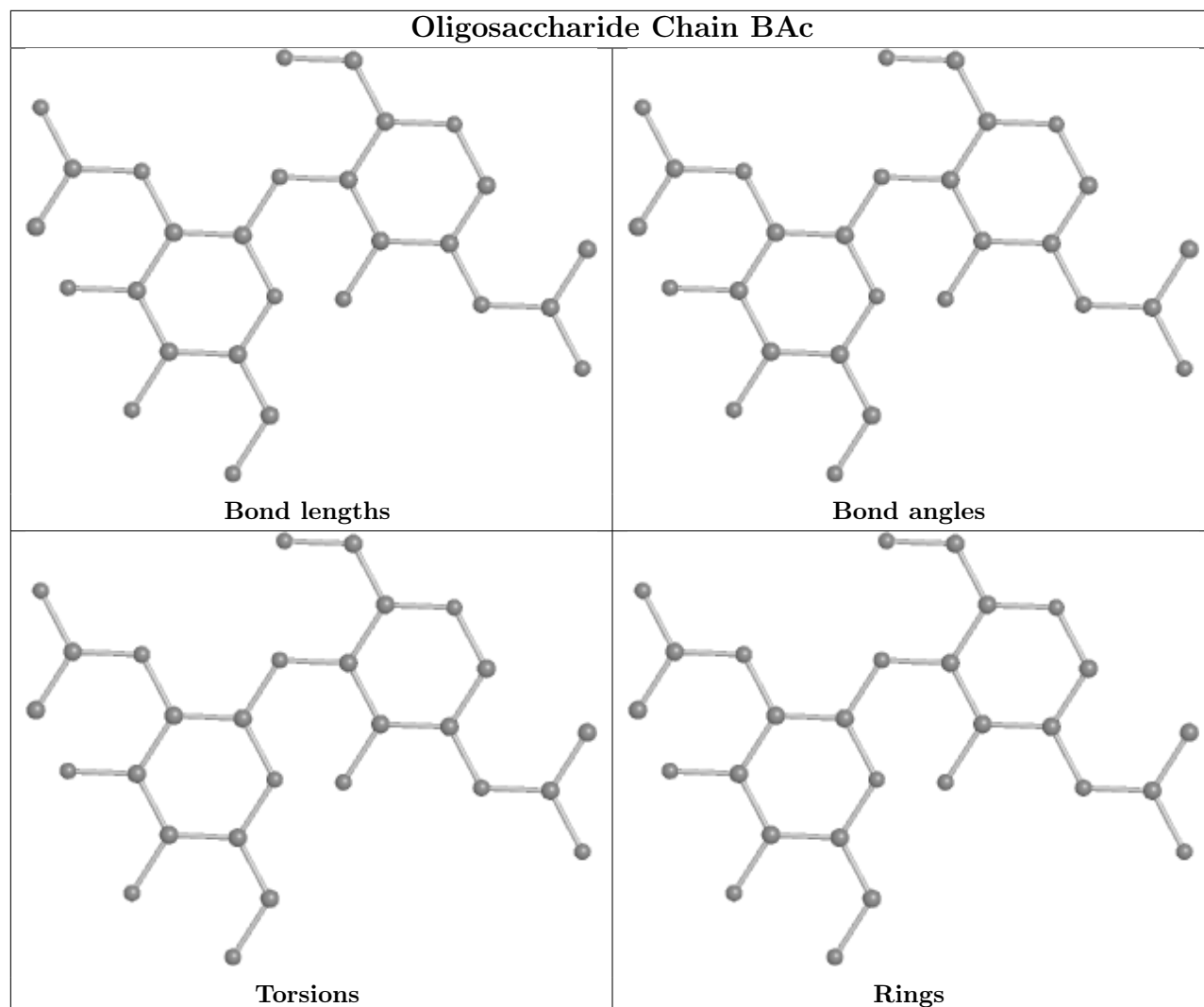
Mol	Chain	Res	Type	Atoms
3	DAc	1	NAG	C3-C2-N2-C7
3	EAc	1	NAG	C3-C2-N2-C7
3	FAc	1	NAG	C3-C2-N2-C7
3	GAc	1	NAG	C3-C2-N2-C7
3	HAc	1	NAG	C3-C2-N2-C7

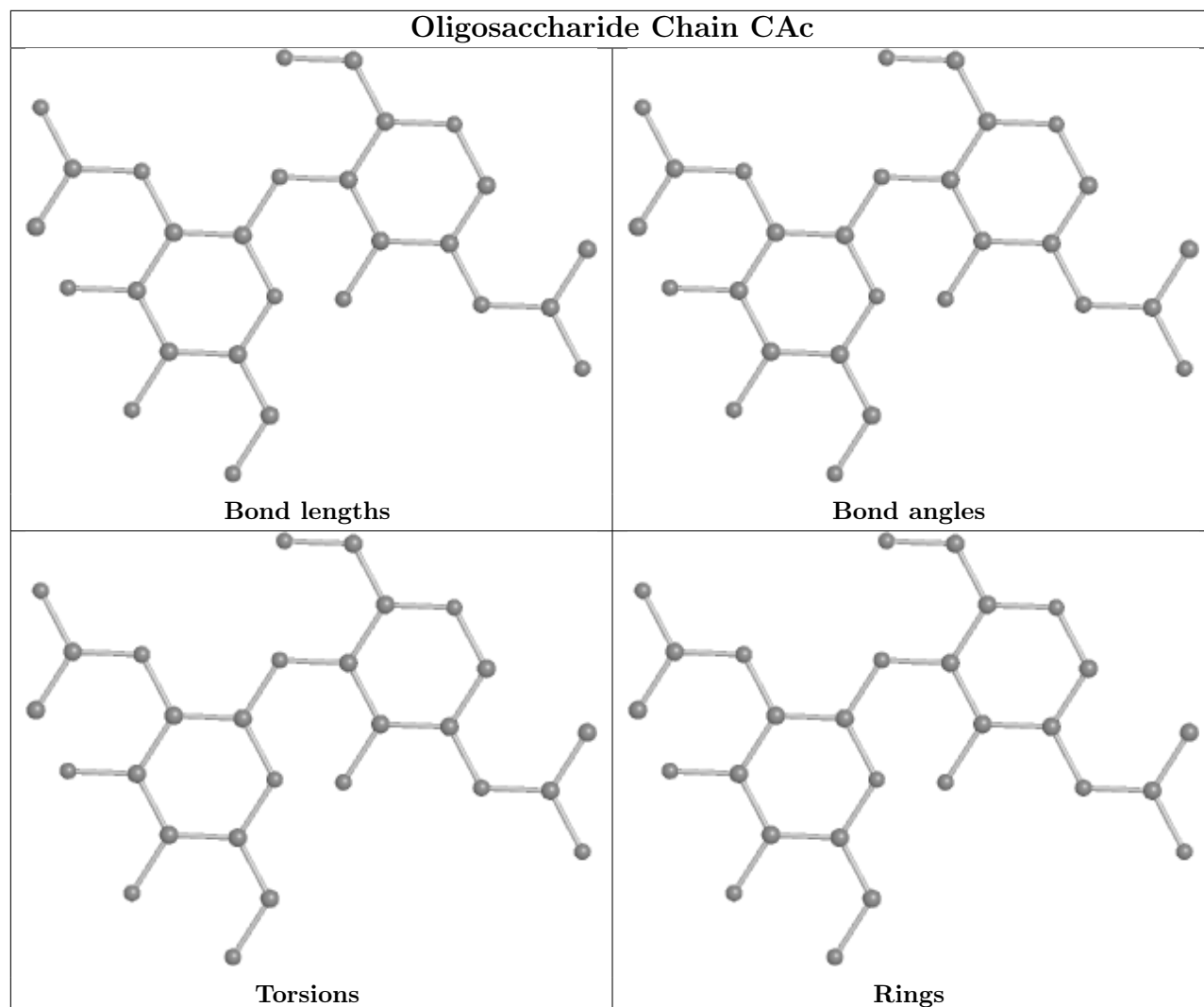
There are no ring outliers.

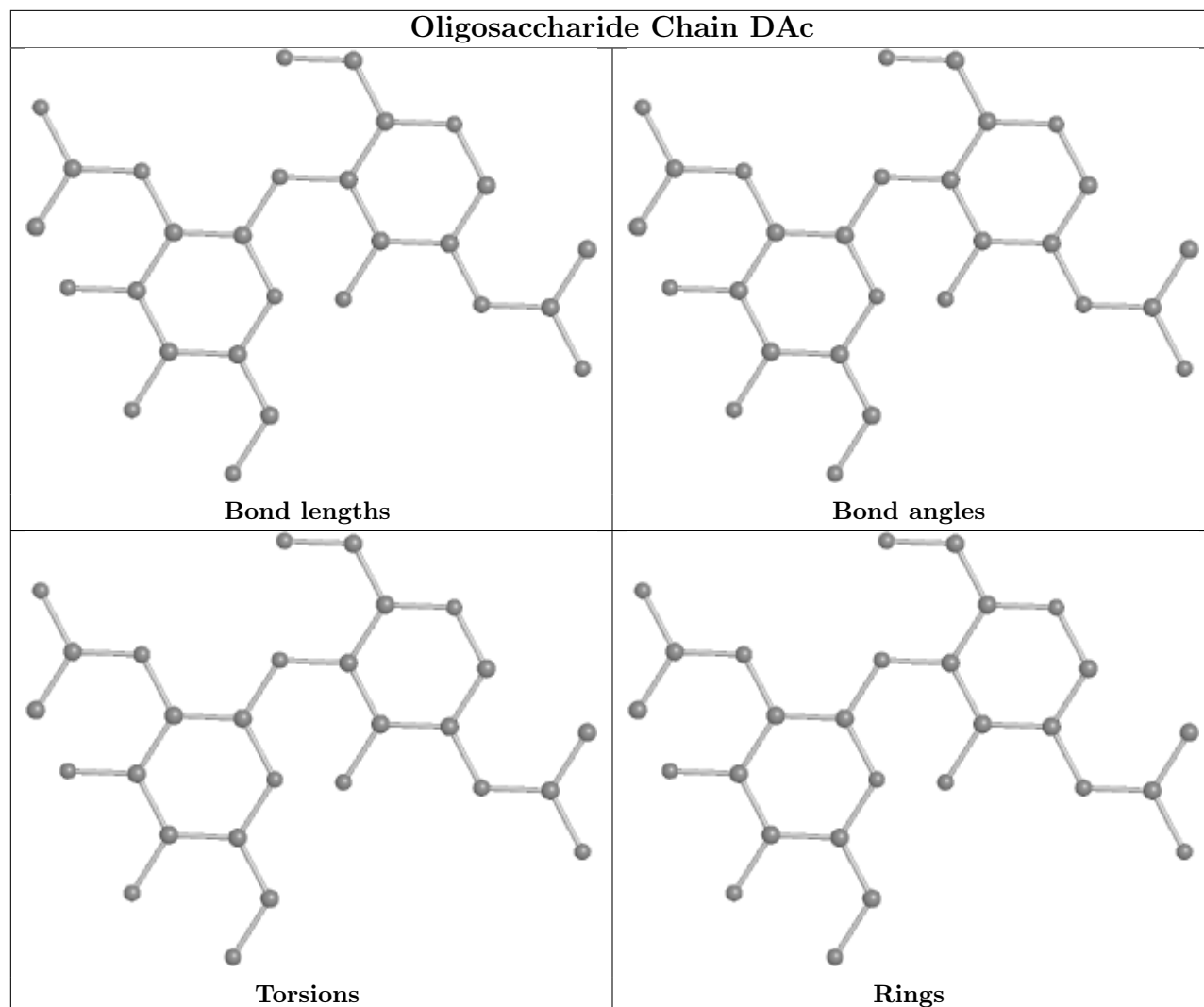
No monomer is involved in short contacts.

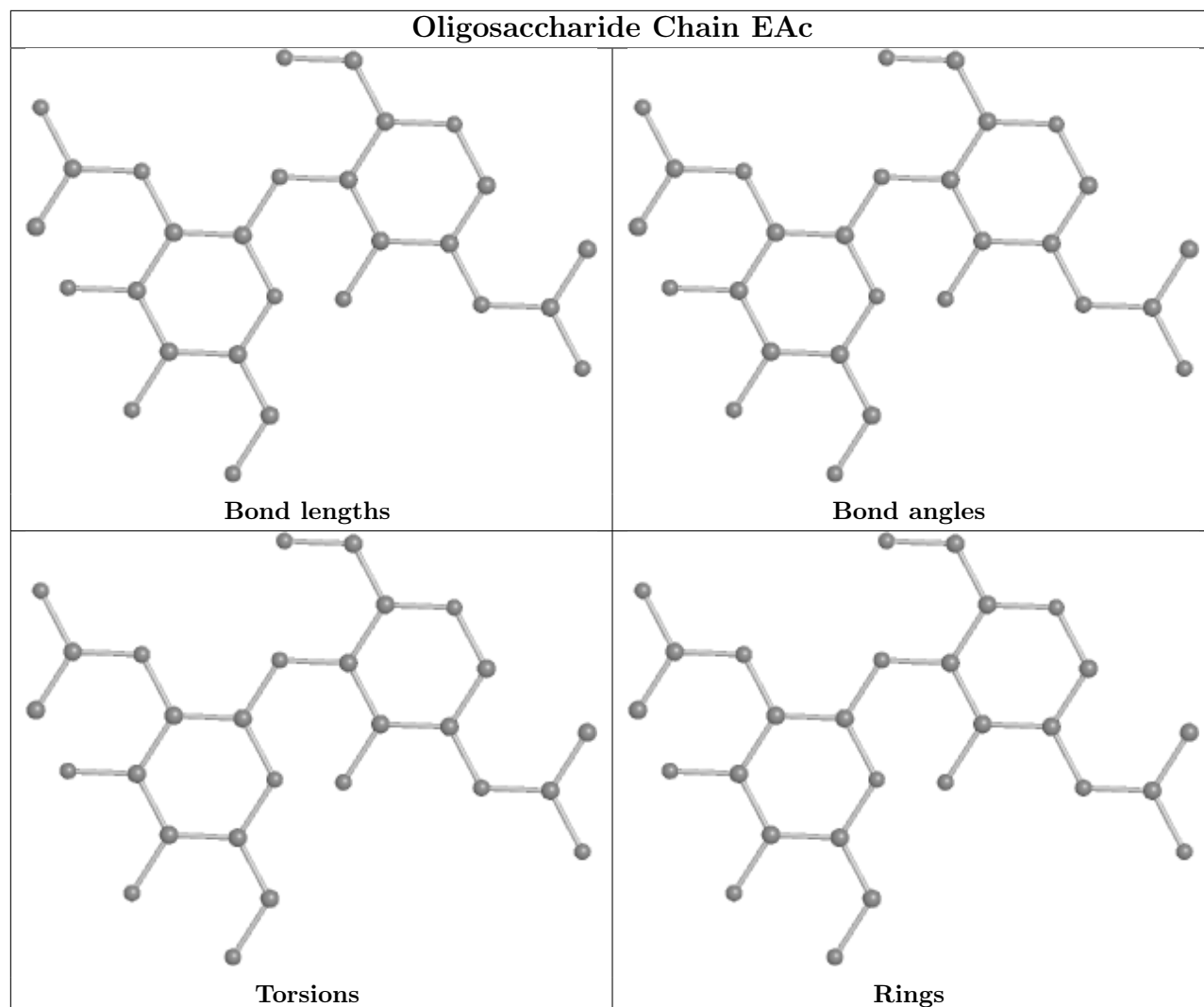
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

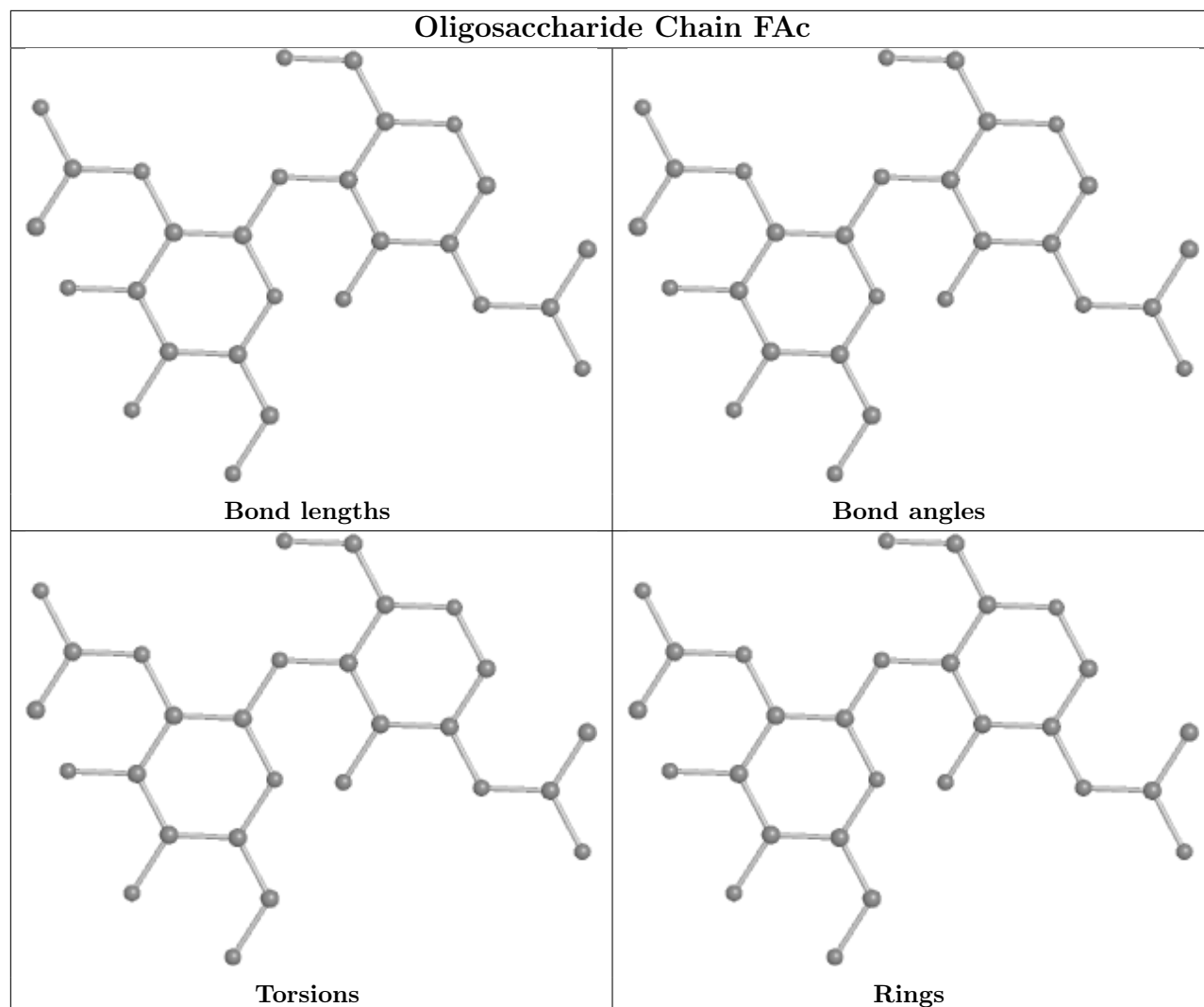




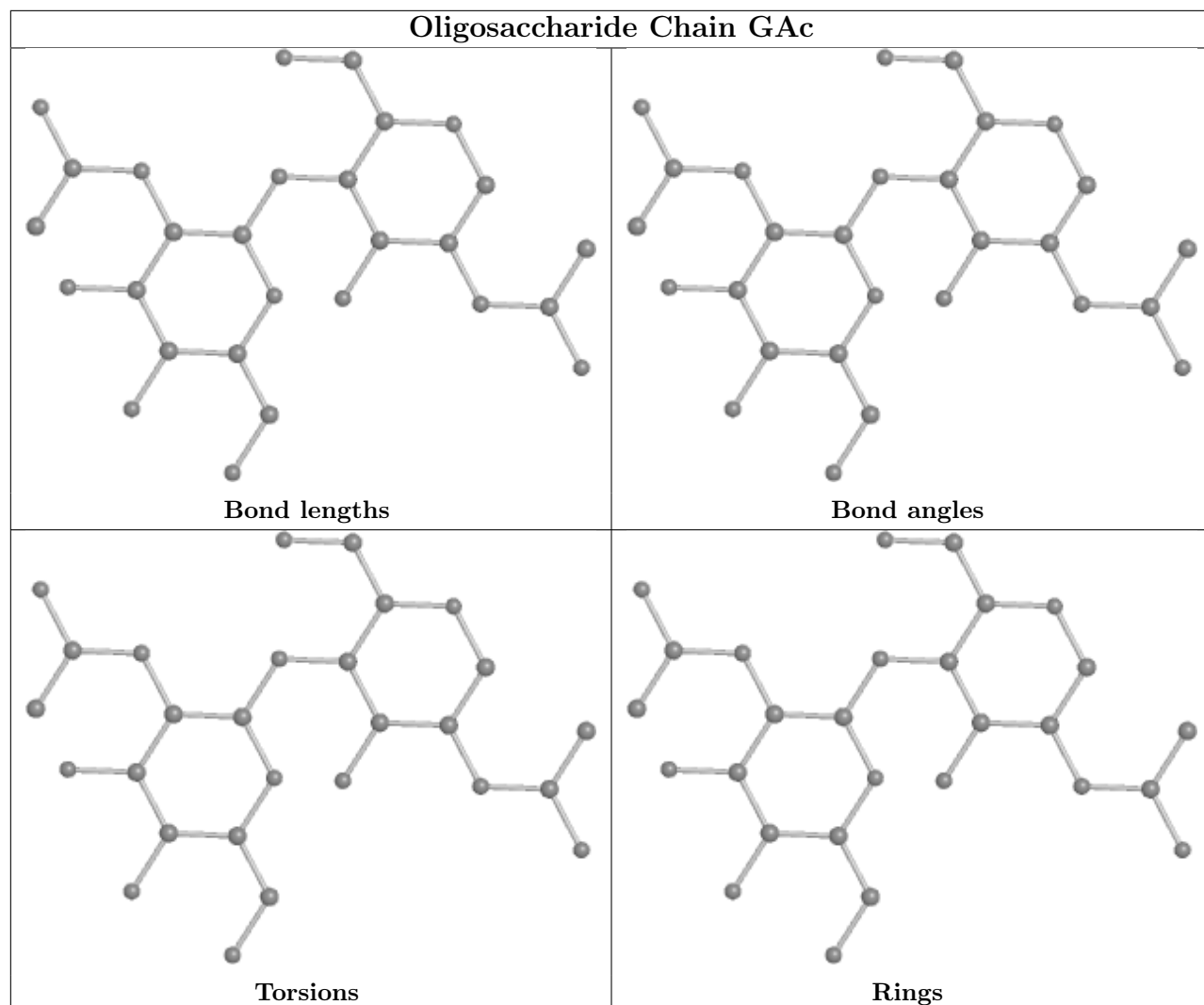


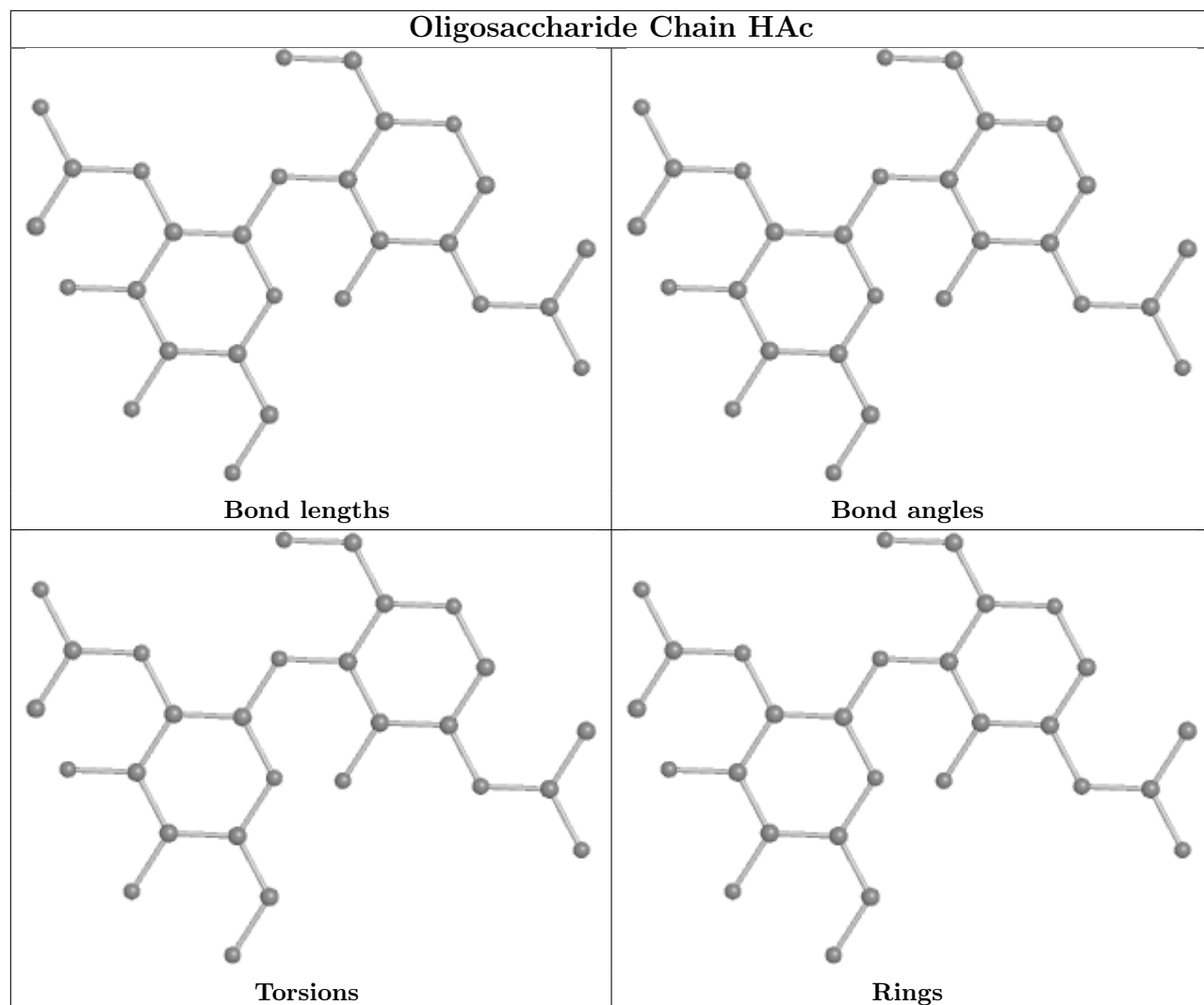


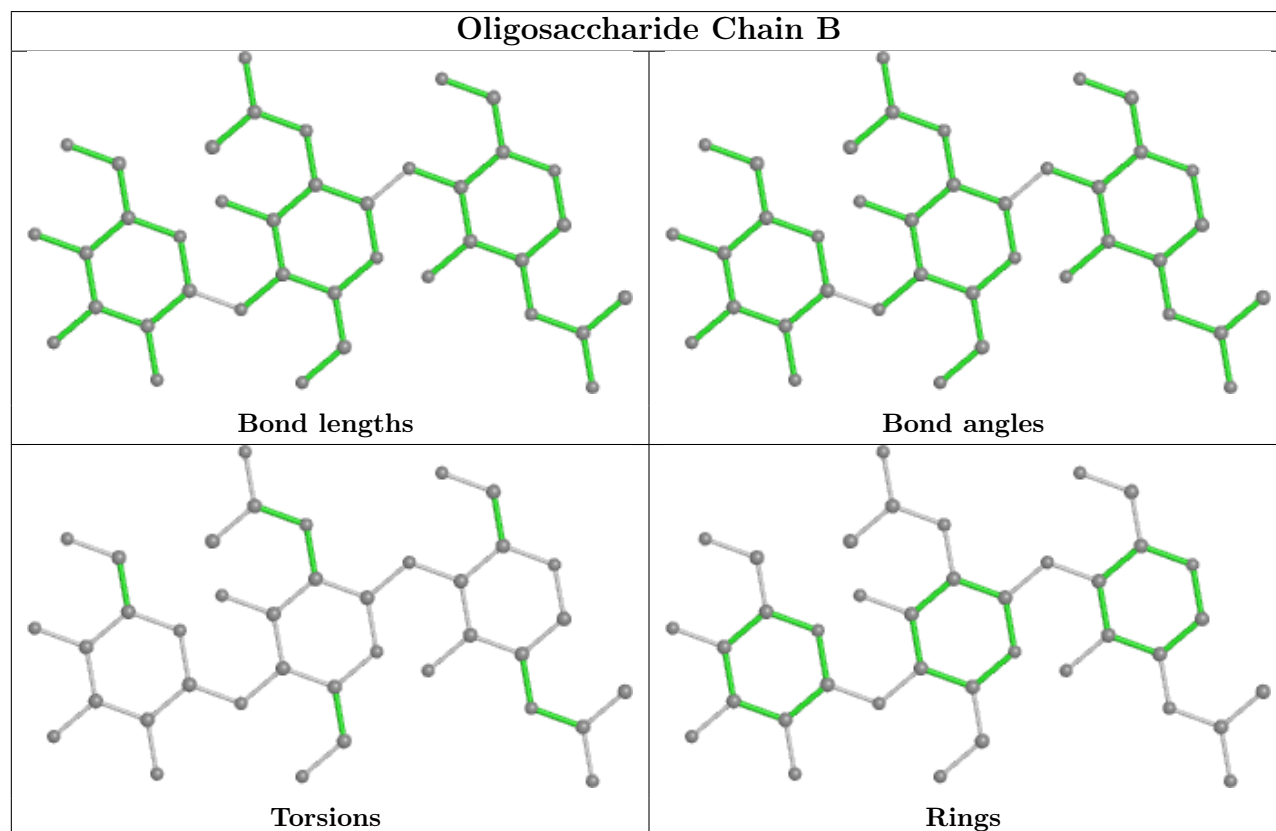
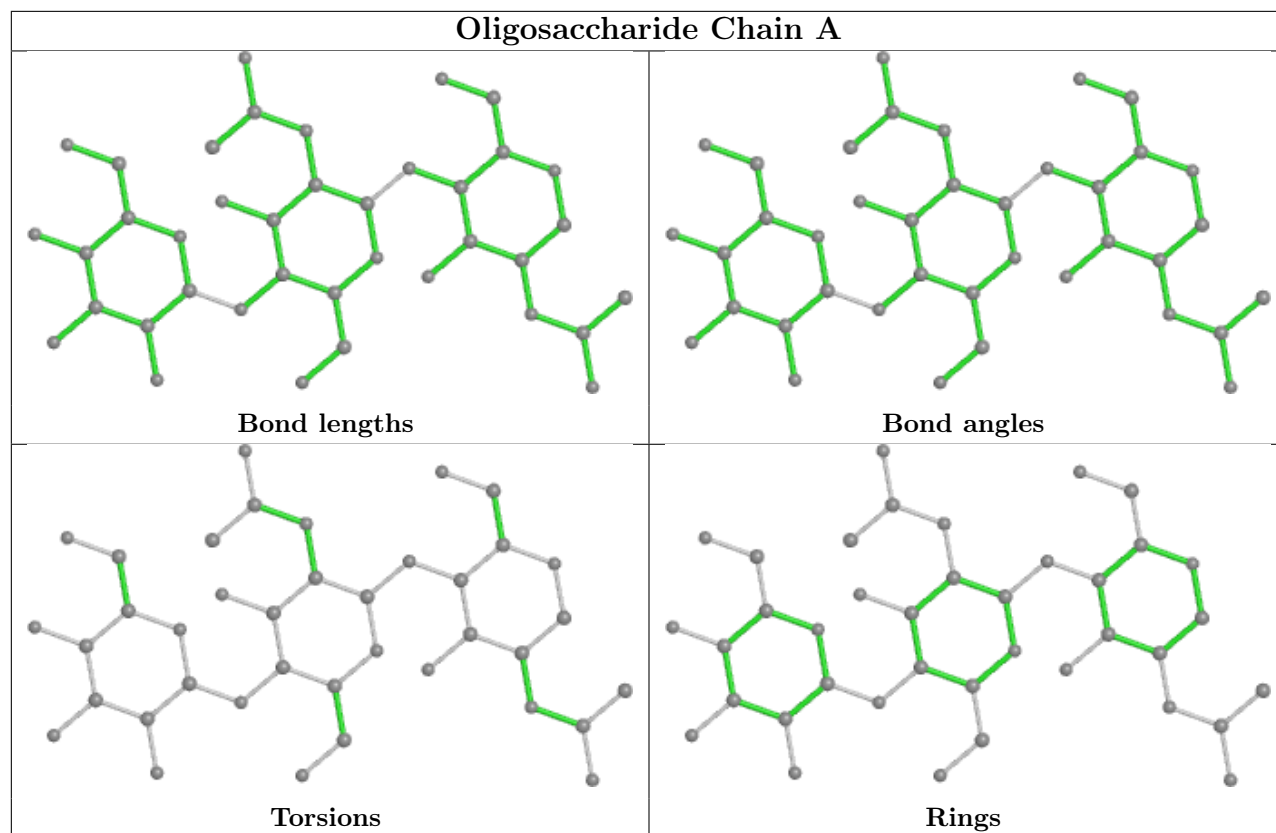


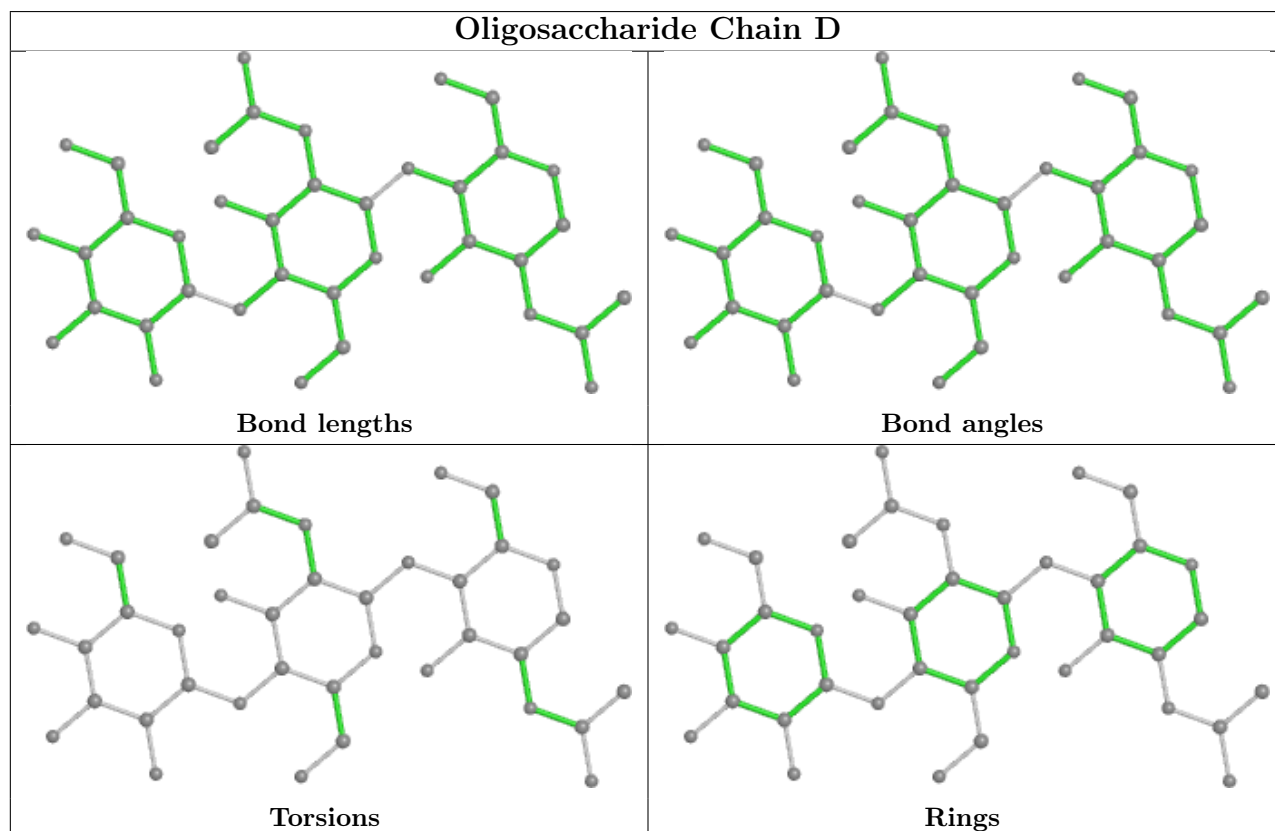
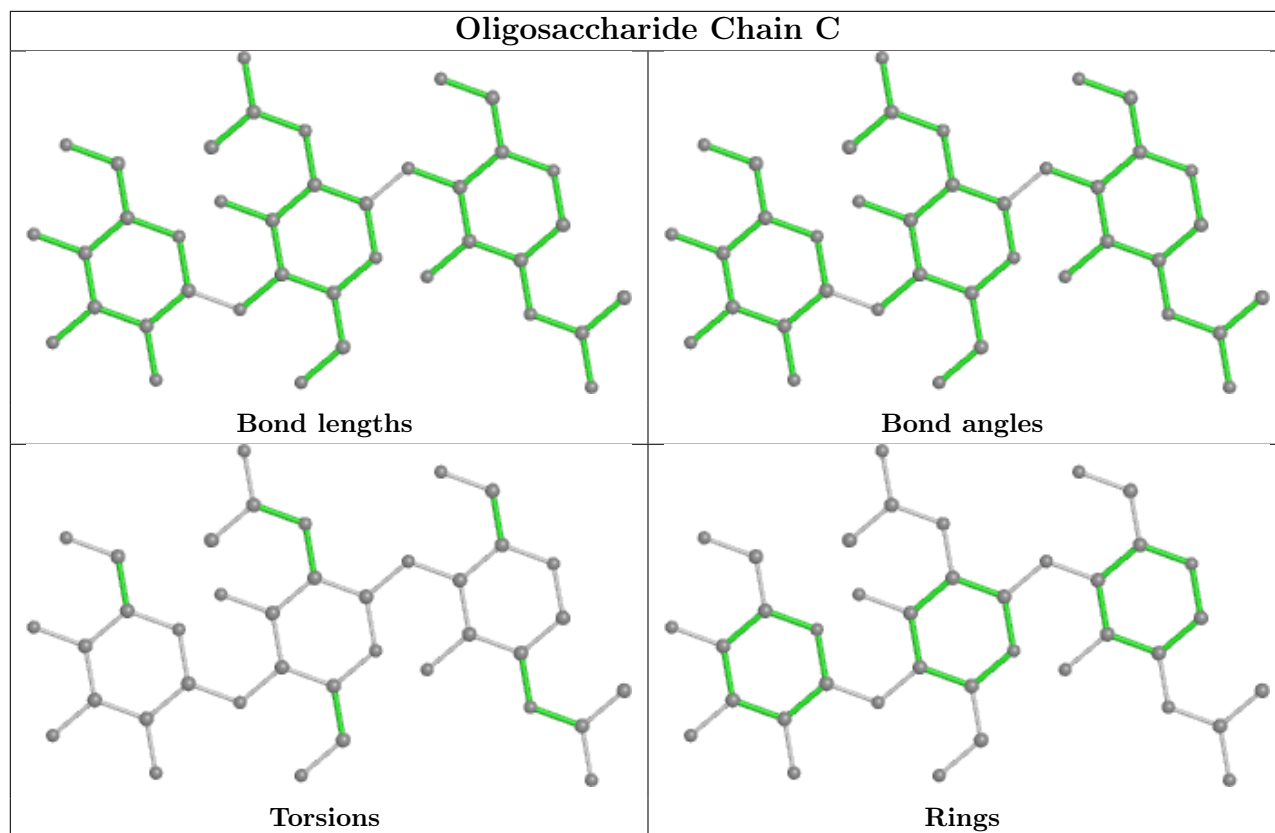


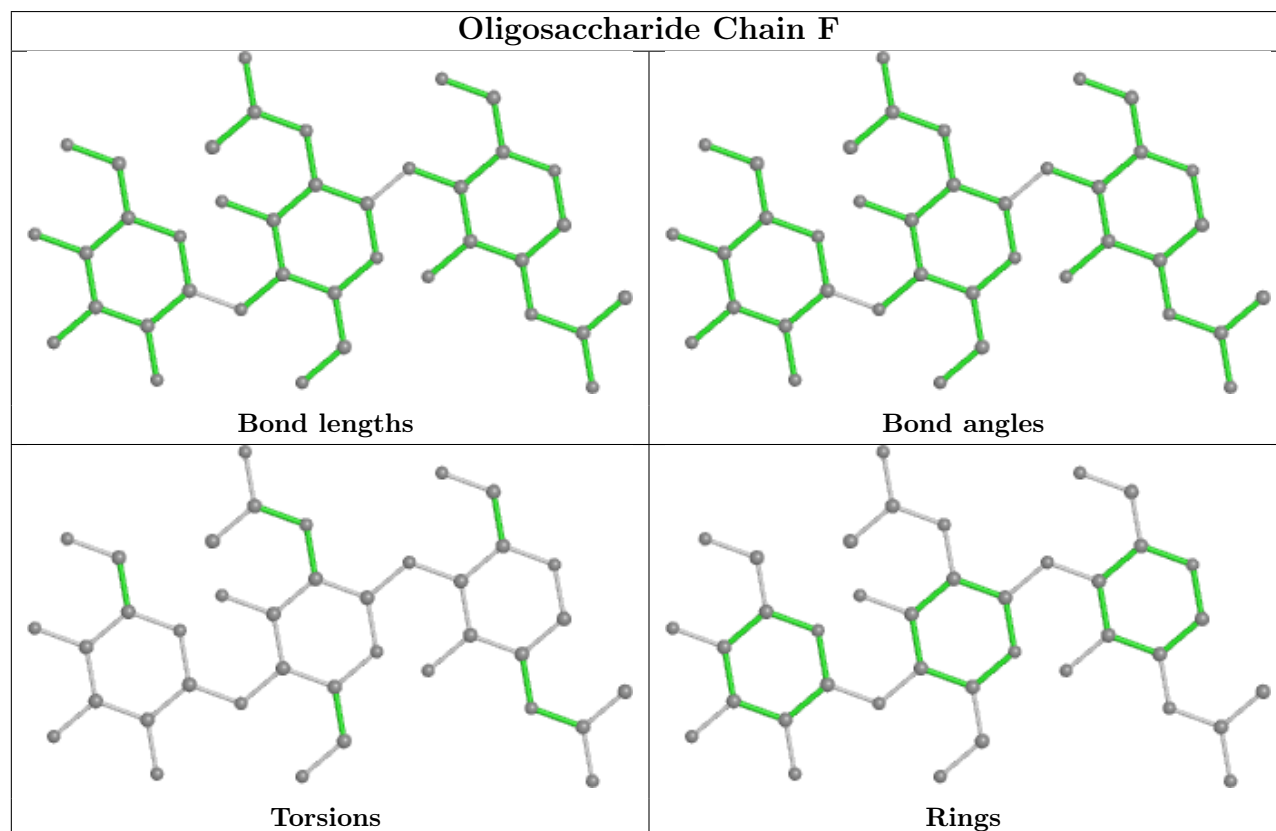
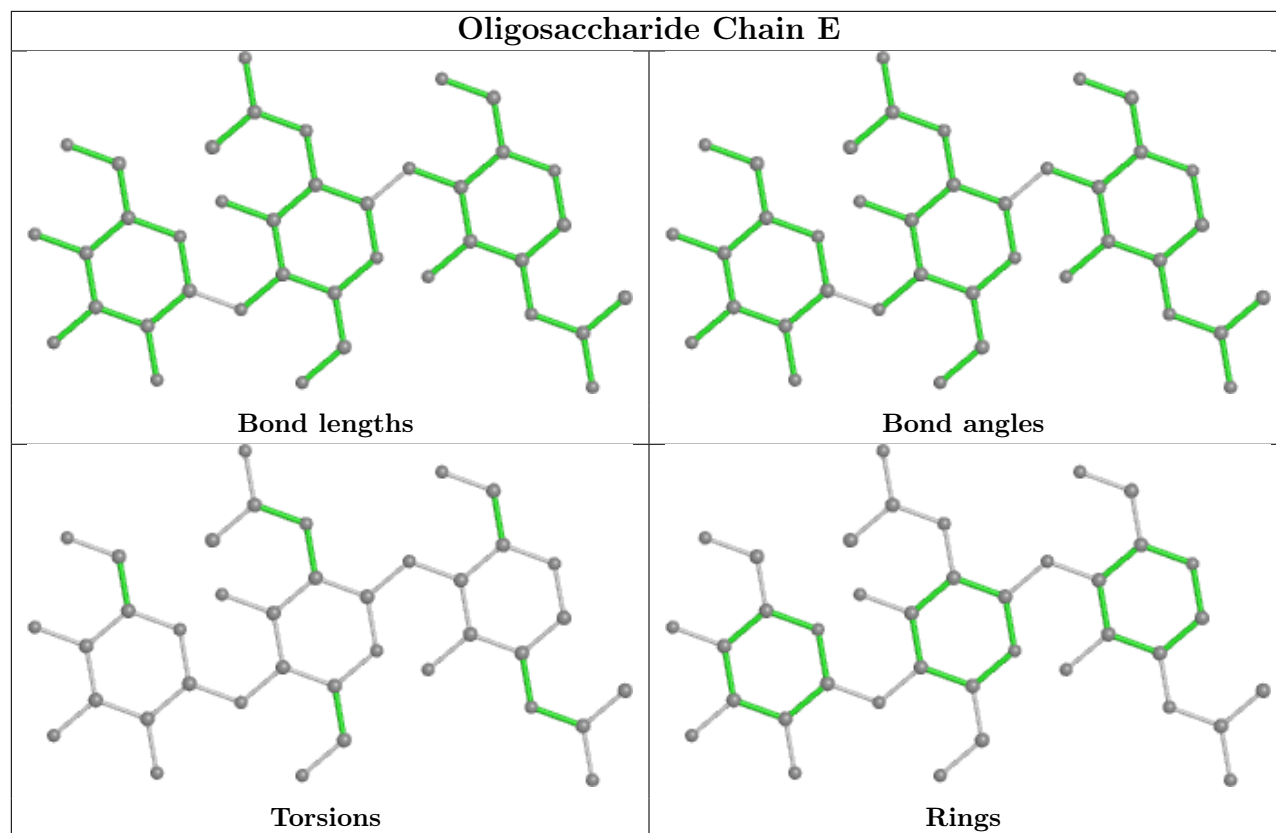


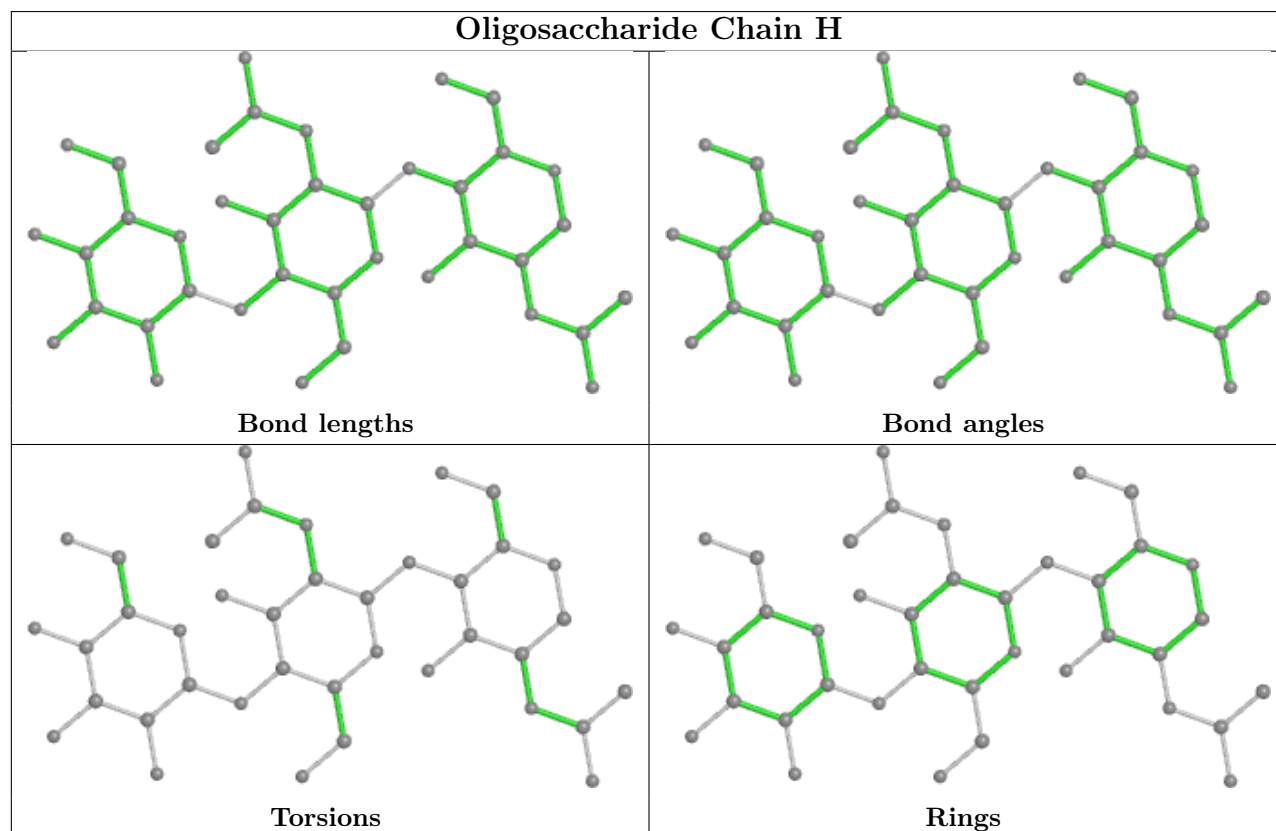
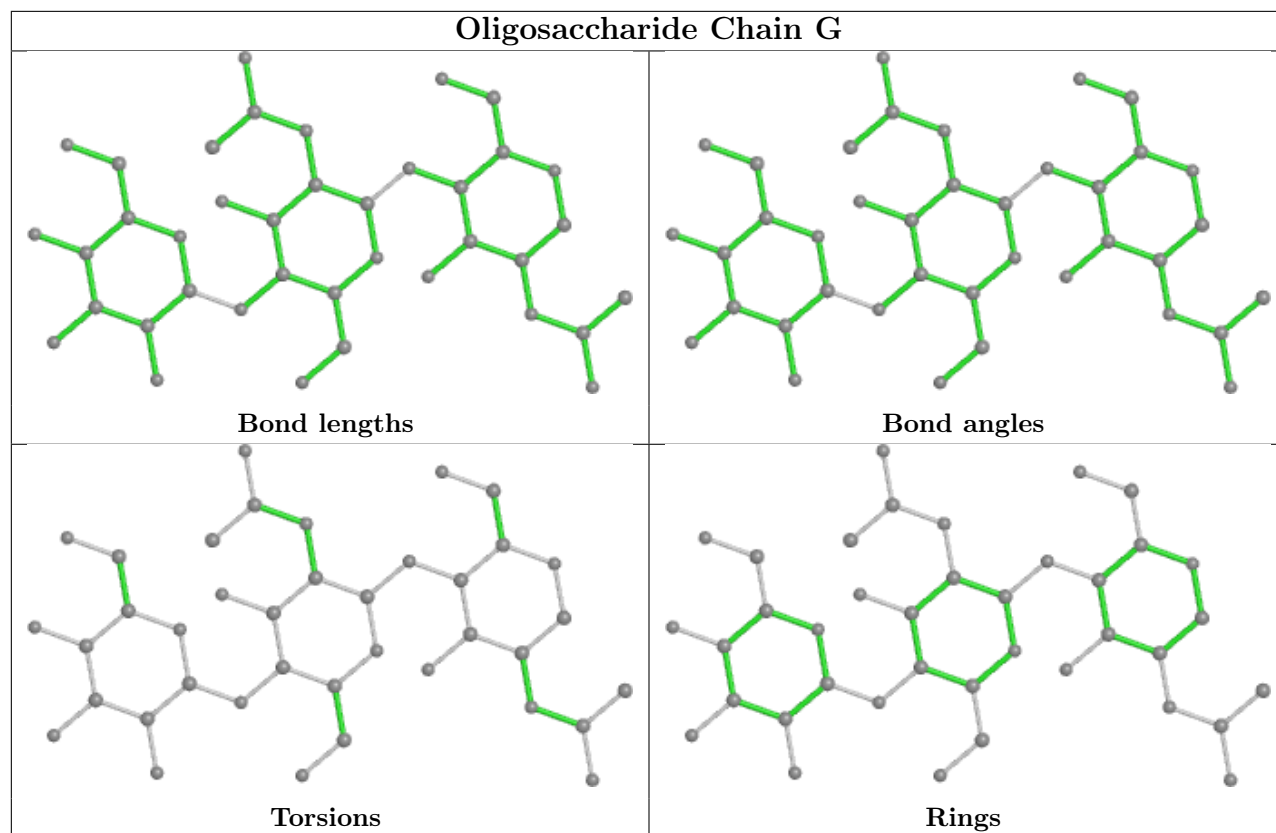












## 5.6 Ligand geometry

32 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
6	NAG	CA	502	1	14,14,15	0.21	0	17,19,21	0.39	0
7	SO4	EA	503	-	4,4,4	0.14	0	6,6,6	0.05	0
6	NAG	BA	502	1	14,14,15	0.20	0	17,19,21	0.39	0
5	94R	HA	501	-	32,32,32	0.55	0	49,50,50	1.69	11 (22%)
6	NAG	DA	502	1	14,14,15	0.21	0	17,19,21	0.39	0
6	NAG	HA	502	1	14,14,15	0.21	0	17,19,21	0.39	0
5	94R	BB	101	-	32,32,32	0.55	0	49,50,50	1.70	11 (22%)
5	94R	GB	101	-	32,32,32	0.55	0	49,50,50	1.70	11 (22%)
5	94R	EA	501	-	32,32,32	0.55	0	49,50,50	1.69	11 (22%)
7	SO4	AA	503	-	4,4,4	0.14	0	6,6,6	0.05	0
7	SO4	DA	503	-	4,4,4	0.14	0	6,6,6	0.05	0
6	NAG	FA	502	1	14,14,15	0.21	0	17,19,21	0.39	0
6	NAG	EA	502	1	14,14,15	0.21	0	17,19,21	0.38	0
7	SO4	BA	503	-	4,4,4	0.14	0	6,6,6	0.04	0
5	94R	CB	101	-	32,32,32	0.54	0	49,50,50	1.70	11 (22%)
7	SO4	FA	503	-	4,4,4	0.14	0	6,6,6	0.04	0
7	SO4	GA	503	-	4,4,4	0.14	0	6,6,6	0.04	0
5	94R	FB	101	-	32,32,32	0.55	0	49,50,50	1.69	11 (22%)
5	94R	DB	101	-	32,32,32	0.55	0	49,50,50	1.70	11 (22%)
5	94R	AA	501	-	32,32,32	0.56	0	49,50,50	1.69	12 (24%)
5	94R	EB	101	-	32,32,32	0.54	0	49,50,50	1.70	11 (22%)
6	NAG	GA	502	1	14,14,15	0.21	0	17,19,21	0.39	0
5	94R	DA	501	-	32,32,32	0.55	0	49,50,50	1.69	11 (22%)
5	94R	HB	101	-	32,32,32	0.55	0	49,50,50	1.70	11 (22%)
5	94R	GA	501	-	32,32,32	0.56	0	49,50,50	1.69	11 (22%)
6	NAG	AA	502	1	14,14,15	0.20	0	17,19,21	0.39	0
7	SO4	CA	503	-	4,4,4	0.14	0	6,6,6	0.04	0
5	94R	BA	501	-	32,32,32	0.55	0	49,50,50	1.69	11 (22%)
7	SO4	HA	503	-	4,4,4	0.14	0	6,6,6	0.05	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	94R	AB	101	-	32,32,32	0.54	0	49,50,50	1.69	11 (22%)
5	94R	FA	501	-	32,32,32	0.56	0	49,50,50	1.69	12 (24%)
5	94R	CA	501	-	32,32,32	0.55	0	49,50,50	1.69	11 (22%)

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
6	NAG	CA	502	1	-	0/6/23/26	0/1/1/1
6	NAG	BA	502	1	-	0/6/23/26	0/1/1/1
5	94R	HA	501	-	-	6/13/71/71	0/4/4/4
6	NAG	DA	502	1	-	0/6/23/26	0/1/1/1
6	NAG	HA	502	1	-	0/6/23/26	0/1/1/1
5	94R	BB	101	-	-	6/13/71/71	0/4/4/4
5	94R	GB	101	-	-	6/13/71/71	0/4/4/4
5	94R	EA	501	-	-	6/13/71/71	0/4/4/4
6	NAG	FA	502	1	-	0/6/23/26	0/1/1/1
6	NAG	EA	502	1	-	0/6/23/26	0/1/1/1
5	94R	CB	101	-	-	6/13/71/71	0/4/4/4
5	94R	FB	101	-	-	6/13/71/71	0/4/4/4
5	94R	DB	101	-	-	6/13/71/71	0/4/4/4
5	94R	AA	501	-	-	6/13/71/71	0/4/4/4
5	94R	EB	101	-	-	6/13/71/71	0/4/4/4
6	NAG	GA	502	1	-	0/6/23/26	0/1/1/1
5	94R	DA	501	-	-	6/13/71/71	0/4/4/4
5	94R	HB	101	-	-	6/13/71/71	0/4/4/4
5	94R	GA	501	-	-	6/13/71/71	0/4/4/4
6	NAG	AA	502	1	-	0/6/23/26	0/1/1/1
5	94R	BA	501	-	-	6/13/71/71	0/4/4/4
5	94R	AB	101	-	-	6/13/71/71	0/4/4/4
5	94R	FA	501	-	-	6/13/71/71	0/4/4/4
5	94R	CA	501	-	-	6/13/71/71	0/4/4/4

There are no bond length outliers.

All (178) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	HA	501	94R	C13-C17-C20	-5.51	110.86	119.49
5	GA	501	94R	C13-C17-C20	-5.51	110.86	119.49
5	BA	501	94R	C13-C17-C20	-5.51	110.86	119.49
5	AA	501	94R	C13-C17-C20	-5.50	110.87	119.49
5	EA	501	94R	C13-C17-C20	-5.50	110.87	119.49
5	FA	501	94R	C13-C17-C20	-5.49	110.88	119.49
5	DA	501	94R	C13-C17-C20	-5.49	110.89	119.49
5	CA	501	94R	C13-C17-C20	-5.48	110.90	119.49
5	AB	101	94R	C13-C17-C20	-5.48	110.91	119.49
5	HB	101	94R	C13-C17-C20	-5.46	110.93	119.49
5	EB	101	94R	C13-C17-C20	-5.46	110.94	119.49
5	GB	101	94R	C13-C17-C20	-5.46	110.94	119.49
5	CB	101	94R	C13-C17-C20	-5.45	110.94	119.49
5	DB	101	94R	C13-C17-C20	-5.45	110.96	119.49
5	BB	101	94R	C13-C17-C20	-5.44	110.96	119.49
5	FB	101	94R	C13-C17-C20	-5.43	110.98	119.49
5	AB	101	94R	C13-C14-C8	-4.37	107.91	114.38
5	BB	101	94R	C13-C14-C8	-4.36	107.93	114.38
5	EB	101	94R	C13-C14-C8	-4.36	107.93	114.38
5	CB	101	94R	C13-C14-C8	-4.35	107.94	114.38
5	HB	101	94R	C13-C14-C8	-4.35	107.94	114.38
5	FB	101	94R	C13-C14-C8	-4.34	107.95	114.38
5	DB	101	94R	C13-C14-C8	-4.33	107.96	114.38
5	GB	101	94R	C13-C14-C8	-4.33	107.97	114.38
5	GA	501	94R	C13-C14-C8	-4.31	107.99	114.38
5	AA	501	94R	C13-C14-C8	-4.30	108.02	114.38
5	CA	501	94R	C13-C14-C8	-4.29	108.03	114.38
5	BA	501	94R	C13-C14-C8	-4.29	108.03	114.38
5	DA	501	94R	C13-C14-C8	-4.29	108.03	114.38
5	EA	501	94R	C13-C14-C8	-4.28	108.04	114.38
5	FA	501	94R	C13-C14-C8	-4.28	108.04	114.38
5	HA	501	94R	C13-C14-C8	-4.26	108.07	114.38
5	FA	501	94R	C19-C10-C9	-3.30	107.75	111.68
5	DA	501	94R	C19-C10-C9	-3.29	107.76	111.68
5	GB	101	94R	C19-C10-C9	-3.28	107.77	111.68
5	AA	501	94R	C19-C10-C9	-3.28	107.77	111.68
5	FB	101	94R	C19-C10-C9	-3.27	107.78	111.68
5	EB	101	94R	C19-C10-C9	-3.27	107.79	111.68
5	HB	101	94R	C19-C10-C9	-3.27	107.79	111.68
5	DB	101	94R	C19-C10-C9	-3.26	107.79	111.68
5	HA	501	94R	C19-C10-C9	-3.26	107.79	111.68
5	AB	101	94R	C19-C10-C9	-3.26	107.79	111.68
5	BA	501	94R	C19-C10-C9	-3.26	107.80	111.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	EA	501	94R	C19-C10-C9	-3.26	107.80	111.68
5	BB	101	94R	C19-C10-C9	-3.25	107.80	111.68
5	CA	501	94R	C19-C10-C9	-3.25	107.80	111.68
5	GA	501	94R	C19-C10-C9	-3.23	107.83	111.68
5	CB	101	94R	C19-C10-C9	-3.23	107.83	111.68
5	GA	501	94R	C11-C12-C13	-2.97	107.69	112.78
5	EA	501	94R	C11-C12-C13	-2.97	107.69	112.78
5	CA	501	94R	C11-C12-C13	-2.96	107.70	112.78
5	HB	101	94R	C11-C12-C13	-2.96	107.70	112.78
5	DA	501	94R	C11-C12-C13	-2.96	107.70	112.78
5	AA	501	94R	C11-C12-C13	-2.96	107.71	112.78
5	AB	101	94R	C11-C12-C13	-2.96	107.71	112.78
5	BB	101	94R	C11-C12-C13	-2.95	107.71	112.78
5	HA	501	94R	C11-C12-C13	-2.95	107.72	112.78
5	CB	101	94R	C11-C12-C13	-2.95	107.72	112.78
5	BA	501	94R	C11-C12-C13	-2.95	107.73	112.78
5	FB	101	94R	C11-C12-C13	-2.94	107.73	112.78
5	FA	501	94R	C11-C12-C13	-2.94	107.74	112.78
5	EB	101	94R	C11-C12-C13	-2.94	107.74	112.78
5	GB	101	94R	C11-C12-C13	-2.93	107.75	112.78
5	DB	101	94R	C11-C12-C13	-2.93	107.76	112.78
5	AB	101	94R	C23-C22-C20	-2.61	109.75	114.52
5	DA	501	94R	C23-C22-C20	-2.60	109.77	114.52
5	CA	501	94R	C23-C22-C20	-2.59	109.78	114.52
5	DB	101	94R	C23-C22-C20	-2.59	109.78	114.52
5	EB	101	94R	C23-C22-C20	-2.58	109.80	114.52
5	HB	101	94R	C23-C22-C20	-2.58	109.80	114.52
5	FB	101	94R	C23-C22-C20	-2.58	109.80	114.52
5	GB	101	94R	C23-C22-C20	-2.58	109.80	114.52
5	BB	101	94R	C23-C22-C20	-2.58	109.80	114.52
5	FA	501	94R	C23-C22-C20	-2.58	109.81	114.52
5	CB	101	94R	C23-C22-C20	-2.58	109.81	114.52
5	BA	501	94R	C23-C22-C20	-2.58	109.81	114.52
5	HA	501	94R	C23-C22-C20	-2.57	109.82	114.52
5	EA	501	94R	C23-C22-C20	-2.57	109.83	114.52
5	GA	501	94R	C23-C22-C20	-2.56	109.84	114.52
5	AA	501	94R	C23-C22-C20	-2.56	109.84	114.52
5	EB	101	94R	C8-C7-C6	-2.32	109.40	112.73
5	DB	101	94R	C8-C7-C6	-2.31	109.41	112.73
5	GB	101	94R	C8-C7-C6	-2.31	109.41	112.73
5	BB	101	94R	C8-C7-C6	-2.31	109.42	112.73
5	AA	501	94R	C21-C20-C17	-2.31	109.39	112.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	DA	501	94R	C21-C20-C17	-2.30	109.40	112.92
5	CB	101	94R	C8-C7-C6	-2.30	109.43	112.73
5	FA	501	94R	C21-C20-C17	-2.29	109.41	112.92
5	FB	101	94R	C8-C7-C6	-2.29	109.44	112.73
5	BB	101	94R	C11-C9-C10	-2.29	110.06	113.08
5	HB	101	94R	C8-C7-C6	-2.29	109.44	112.73
5	FB	101	94R	C11-C9-C10	-2.29	110.07	113.08
5	CA	501	94R	C8-C7-C6	-2.29	109.45	112.73
5	EA	501	94R	C21-C20-C17	-2.29	109.42	112.92
5	GA	501	94R	C21-C20-C17	-2.28	109.42	112.92
5	HA	501	94R	C8-C7-C6	-2.28	109.45	112.73
5	EB	101	94R	C11-C9-C10	-2.28	110.08	113.08
5	HA	501	94R	C21-C20-C17	-2.27	109.44	112.92
5	EA	501	94R	C8-C7-C6	-2.27	109.47	112.73
5	CA	501	94R	C3-C4-C5	-2.27	108.17	112.03
5	BA	501	94R	C21-C20-C17	-2.27	109.45	112.92
5	DA	501	94R	C8-C7-C6	-2.27	109.48	112.73
5	CA	501	94R	C21-C20-C17	-2.27	109.45	112.92
5	DB	101	94R	C11-C9-C10	-2.26	110.10	113.08
5	GA	501	94R	C8-C7-C6	-2.26	109.48	112.73
5	EA	501	94R	C3-C4-C5	-2.26	108.19	112.03
5	DA	501	94R	C3-C4-C5	-2.26	108.19	112.03
5	BA	501	94R	C8-C7-C6	-2.26	109.48	112.73
5	AB	101	94R	C8-C7-C6	-2.26	109.49	112.73
5	FA	501	94R	C8-C7-C6	-2.26	109.49	112.73
5	AA	501	94R	C3-C4-C5	-2.26	108.20	112.03
5	AA	501	94R	C8-C7-C6	-2.26	109.49	112.73
5	GB	101	94R	C11-C9-C10	-2.26	110.11	113.08
5	CB	101	94R	C11-C9-C10	-2.25	110.11	113.08
5	GB	101	94R	C21-C20-C17	-2.25	109.47	112.92
5	HA	501	94R	C3-C4-C5	-2.25	108.21	112.03
5	AB	101	94R	C11-C9-C10	-2.25	110.12	113.08
5	GA	501	94R	C11-C9-C10	-2.25	110.12	113.08
5	DB	101	94R	C21-C20-C17	-2.24	109.49	112.92
5	GA	501	94R	C3-C4-C5	-2.24	108.22	112.03
5	HB	101	94R	C21-C20-C17	-2.24	109.49	112.92
5	BA	501	94R	C11-C9-C10	-2.24	110.13	113.08
5	FA	501	94R	C3-C4-C5	-2.24	108.23	112.03
5	DB	101	94R	C3-C4-C5	-2.23	108.23	112.03
5	BB	101	94R	C21-C20-C17	-2.23	109.50	112.92
5	HB	101	94R	C11-C9-C10	-2.23	110.14	113.08
5	BA	501	94R	C3-C4-C5	-2.23	108.24	112.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	CB	101	94R	C3-C4-C5	-2.23	108.24	112.03
5	CA	501	94R	C11-C9-C10	-2.23	110.14	113.08
5	HA	501	94R	C11-C9-C10	-2.23	110.14	113.08
5	AA	501	94R	C11-C9-C10	-2.23	110.15	113.08
5	BB	101	94R	C3-C4-C5	-2.22	108.25	112.03
5	GB	101	94R	C3-C4-C5	-2.22	108.26	112.03
5	CB	101	94R	C21-C20-C17	-2.22	109.53	112.92
5	EB	101	94R	C21-C20-C17	-2.22	109.53	112.92
5	FB	101	94R	C3-C4-C5	-2.22	108.27	112.03
5	HB	101	94R	C3-C4-C5	-2.21	108.27	112.03
5	FB	101	94R	C21-C20-C17	-2.21	109.53	112.92
5	AB	101	94R	C21-C20-C17	-2.21	109.53	112.92
5	EB	101	94R	C3-C4-C5	-2.20	108.29	112.03
5	DA	501	94R	C11-C9-C10	-2.19	110.19	113.08
5	AB	101	94R	C3-C4-C5	-2.19	108.30	112.03
5	EA	501	94R	C11-C9-C10	-2.19	110.19	113.08
5	FA	501	94R	C11-C9-C10	-2.19	110.19	113.08
5	GB	101	94R	C14-C13-C17	2.18	102.66	100.07
5	BB	101	94R	C14-C13-C17	2.17	102.65	100.07
5	DB	101	94R	C14-C13-C17	2.17	102.65	100.07
5	FB	101	94R	C14-C13-C17	2.16	102.64	100.07
5	EB	101	94R	C14-C13-C17	2.15	102.62	100.07
5	HB	101	94R	C14-C13-C17	2.13	102.60	100.07
5	CB	101	94R	C14-C13-C17	2.13	102.59	100.07
5	AB	101	94R	C14-C13-C17	2.12	102.58	100.07
5	HB	101	94R	C9-C10-C5	2.10	112.95	109.65
5	CB	101	94R	C9-C10-C5	2.10	112.94	109.65
5	GB	101	94R	C9-C10-C5	2.10	112.94	109.65
5	AB	101	94R	C9-C10-C5	2.10	112.94	109.65
5	FB	101	94R	C9-C10-C5	2.09	112.94	109.65
5	EB	101	94R	C9-C10-C5	2.08	112.92	109.65
5	DB	101	94R	C9-C10-C5	2.08	112.91	109.65
5	FA	501	94R	C14-C13-C17	2.07	102.53	100.07
5	BB	101	94R	C9-C10-C5	2.06	112.88	109.65
5	DA	501	94R	C9-C10-C5	2.06	112.88	109.65
5	GA	501	94R	C9-C10-C5	2.05	112.87	109.65
5	EA	501	94R	C9-C10-C5	2.05	112.86	109.65
5	AA	501	94R	C14-C13-C17	2.05	102.50	100.07
5	EA	501	94R	C14-C13-C17	2.05	102.50	100.07
5	FA	501	94R	C9-C10-C5	2.04	112.86	109.65
5	DA	501	94R	C14-C13-C17	2.04	102.49	100.07
5	CA	501	94R	C14-C13-C17	2.04	102.49	100.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	HA	501	94R	C9-C10-C5	2.04	112.85	109.65
5	HA	501	94R	C14-C13-C17	2.04	102.49	100.07
5	BA	501	94R	C14-C13-C17	2.03	102.48	100.07
5	GA	501	94R	C14-C13-C17	2.02	102.47	100.07
5	AA	501	94R	C9-C10-C5	2.02	112.81	109.65
5	CA	501	94R	C9-C10-C5	2.01	112.81	109.65
5	BA	501	94R	C9-C10-C5	2.01	112.81	109.65
5	AA	501	94R	C15-C14-C13	-2.01	101.42	103.84
5	FA	501	94R	C15-C14-C13	-2.00	101.43	103.84

There are no chirality outliers.

All (96) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	AA	501	94R	C20-C22-C23-C24
5	BA	501	94R	C20-C22-C23-C24
5	CA	501	94R	C20-C22-C23-C24
5	DA	501	94R	C20-C22-C23-C24
5	EA	501	94R	C20-C22-C23-C24
5	FA	501	94R	C20-C22-C23-C24
5	GA	501	94R	C20-C22-C23-C24
5	HA	501	94R	C20-C22-C23-C24
5	AA	501	94R	C13-C17-C20-C21
5	BA	501	94R	C13-C17-C20-C21
5	CA	501	94R	C13-C17-C20-C21
5	DA	501	94R	C13-C17-C20-C21
5	EA	501	94R	C13-C17-C20-C21
5	FA	501	94R	C13-C17-C20-C21
5	GA	501	94R	C13-C17-C20-C21
5	HA	501	94R	C13-C17-C20-C21
5	AA	501	94R	C21-C20-C22-C23
5	AB	101	94R	C21-C20-C22-C23
5	BA	501	94R	C21-C20-C22-C23
5	CA	501	94R	C21-C20-C22-C23
5	DA	501	94R	C21-C20-C22-C23
5	DB	101	94R	C21-C20-C22-C23
5	EA	501	94R	C21-C20-C22-C23
5	FA	501	94R	C21-C20-C22-C23
5	GA	501	94R	C21-C20-C22-C23
5	HA	501	94R	C21-C20-C22-C23
5	BB	101	94R	C21-C20-C22-C23
5	CB	101	94R	C21-C20-C22-C23

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Mol	Chain	Res	Type	Atoms
5	EB	101	94R	C21-C20-C22-C23
5	FB	101	94R	C21-C20-C22-C23
5	GB	101	94R	C21-C20-C22-C23
5	HB	101	94R	C21-C20-C22-C23
5	AA	501	94R	C13-C17-C20-C22
5	BA	501	94R	C13-C17-C20-C22
5	CA	501	94R	C13-C17-C20-C22
5	DA	501	94R	C13-C17-C20-C22
5	EA	501	94R	C13-C17-C20-C22
5	FA	501	94R	C13-C17-C20-C22
5	GA	501	94R	C13-C17-C20-C22
5	HA	501	94R	C13-C17-C20-C22
5	AB	101	94R	C13-C17-C20-C21
5	BB	101	94R	C13-C17-C20-C21
5	CB	101	94R	C13-C17-C20-C21
5	DB	101	94R	C13-C17-C20-C21
5	EB	101	94R	C13-C17-C20-C21
5	FB	101	94R	C13-C17-C20-C21
5	GB	101	94R	C13-C17-C20-C21
5	HB	101	94R	C13-C17-C20-C21
5	AB	101	94R	C13-C17-C20-C22
5	BB	101	94R	C13-C17-C20-C22
5	CB	101	94R	C13-C17-C20-C22
5	DB	101	94R	C13-C17-C20-C22
5	EB	101	94R	C13-C17-C20-C22
5	FB	101	94R	C13-C17-C20-C22
5	GB	101	94R	C13-C17-C20-C22
5	HB	101	94R	C13-C17-C20-C22
5	AB	101	94R	C20-C22-C23-C24
5	BB	101	94R	C20-C22-C23-C24
5	CB	101	94R	C20-C22-C23-C24
5	DB	101	94R	C20-C22-C23-C24
5	EB	101	94R	C20-C22-C23-C24
5	FB	101	94R	C20-C22-C23-C24
5	GB	101	94R	C20-C22-C23-C24
5	HB	101	94R	C20-C22-C23-C24
5	AA	501	94R	C16-C17-C20-C21
5	BA	501	94R	C16-C17-C20-C21
5	CA	501	94R	C16-C17-C20-C21
5	DA	501	94R	C16-C17-C20-C21
5	EA	501	94R	C16-C17-C20-C21
5	FA	501	94R	C16-C17-C20-C21

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Mol	Chain	Res	Type	Atoms
5	GA	501	94R	C16-C17-C20-C21
5	HA	501	94R	C16-C17-C20-C21
5	AB	101	94R	C16-C17-C20-C21
5	BB	101	94R	C16-C17-C20-C21
5	CB	101	94R	C16-C17-C20-C21
5	DB	101	94R	C16-C17-C20-C21
5	EB	101	94R	C16-C17-C20-C21
5	FB	101	94R	C16-C17-C20-C21
5	GB	101	94R	C16-C17-C20-C21
5	HB	101	94R	C16-C17-C20-C21
5	AA	501	94R	C16-C17-C20-C22
5	BA	501	94R	C16-C17-C20-C22
5	CA	501	94R	C16-C17-C20-C22
5	DA	501	94R	C16-C17-C20-C22
5	FA	501	94R	C16-C17-C20-C22
5	HA	501	94R	C16-C17-C20-C22
5	EA	501	94R	C16-C17-C20-C22
5	GA	501	94R	C16-C17-C20-C22
5	AB	101	94R	C16-C17-C20-C22
5	BB	101	94R	C16-C17-C20-C22
5	CB	101	94R	C16-C17-C20-C22
5	DB	101	94R	C16-C17-C20-C22
5	EB	101	94R	C16-C17-C20-C22
5	FB	101	94R	C16-C17-C20-C22
5	GB	101	94R	C16-C17-C20-C22
5	HB	101	94R	C16-C17-C20-C22

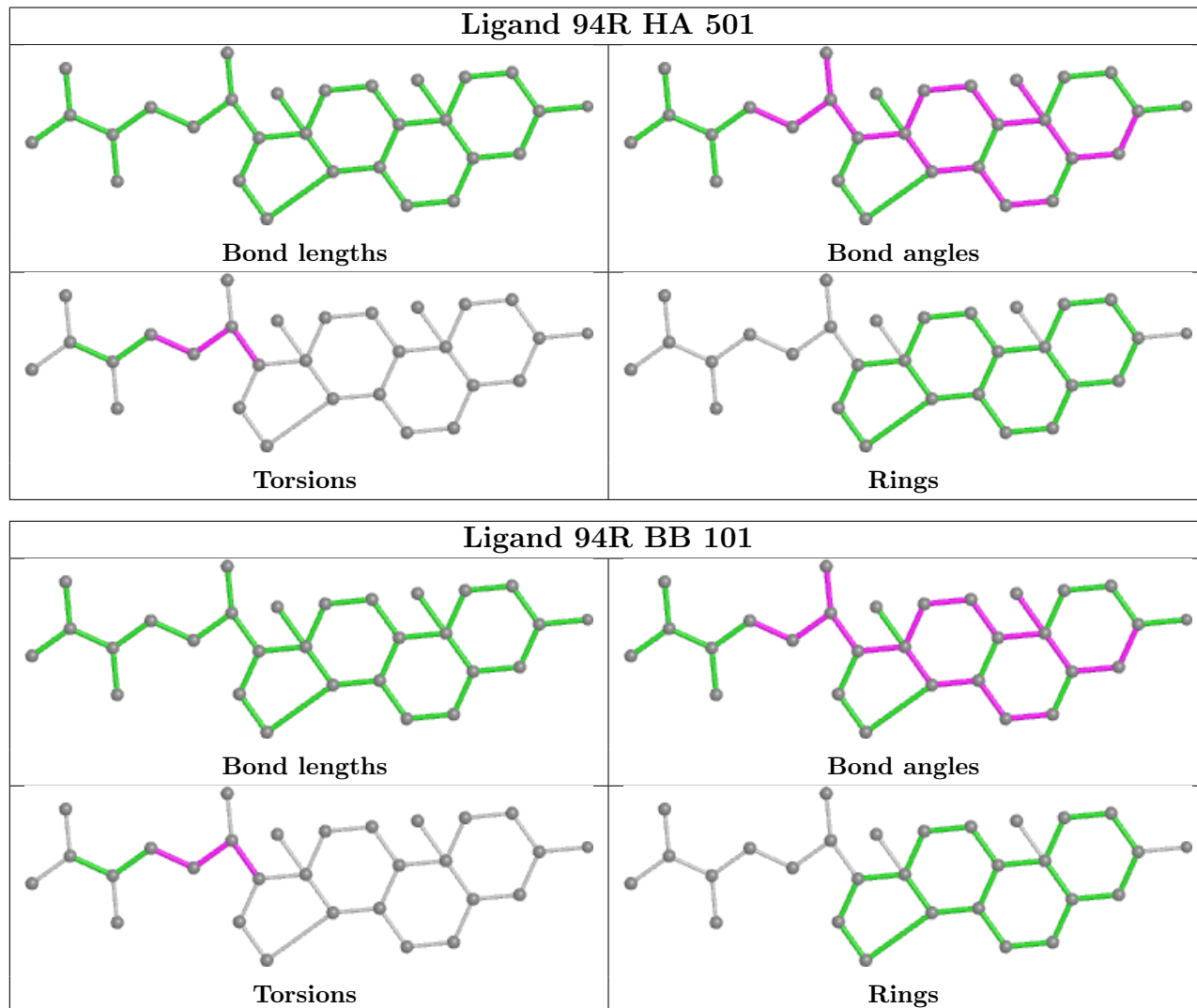
There are no ring outliers.

8 monomers are involved in 8 short contacts:

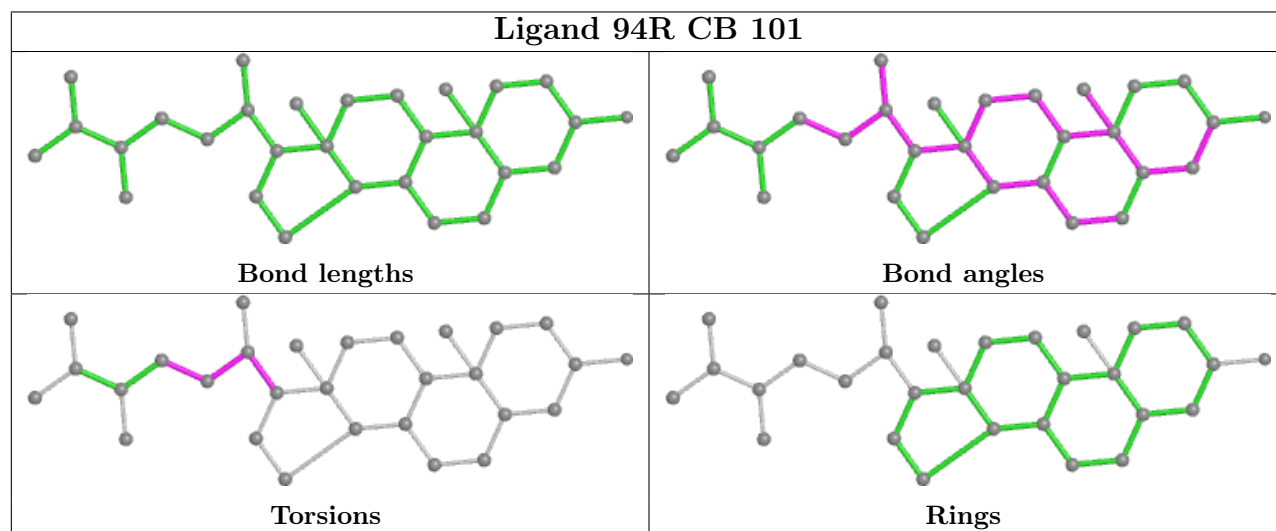
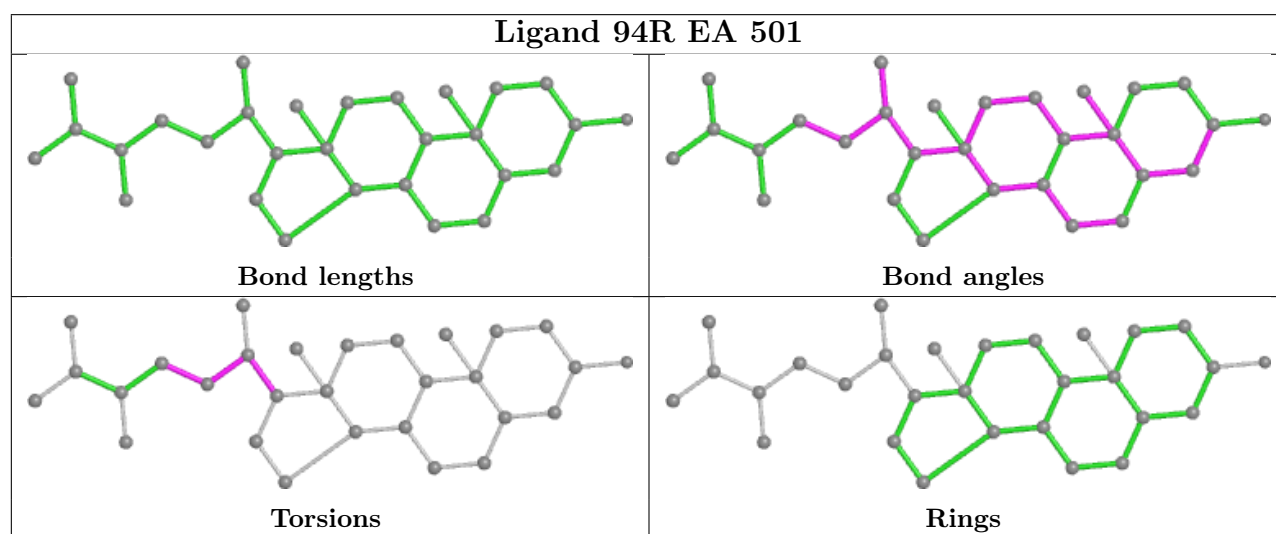
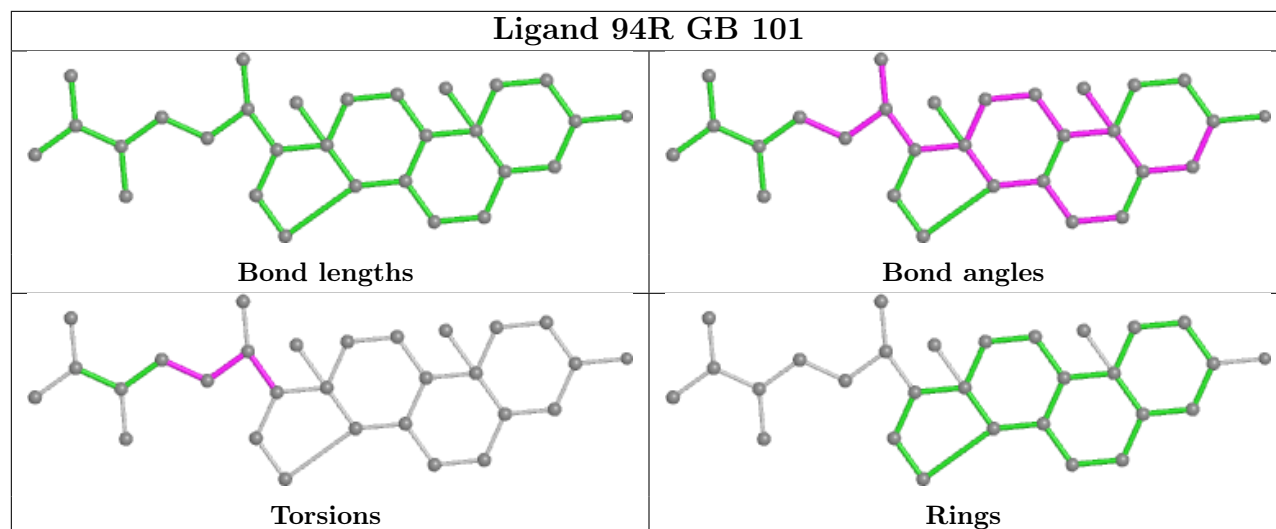
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	HA	501	94R	1	0
5	EA	501	94R	1	0
5	AA	501	94R	1	0
5	DA	501	94R	1	0
5	GA	501	94R	1	0
5	BA	501	94R	1	0
5	FA	501	94R	1	0
5	CA	501	94R	1	0

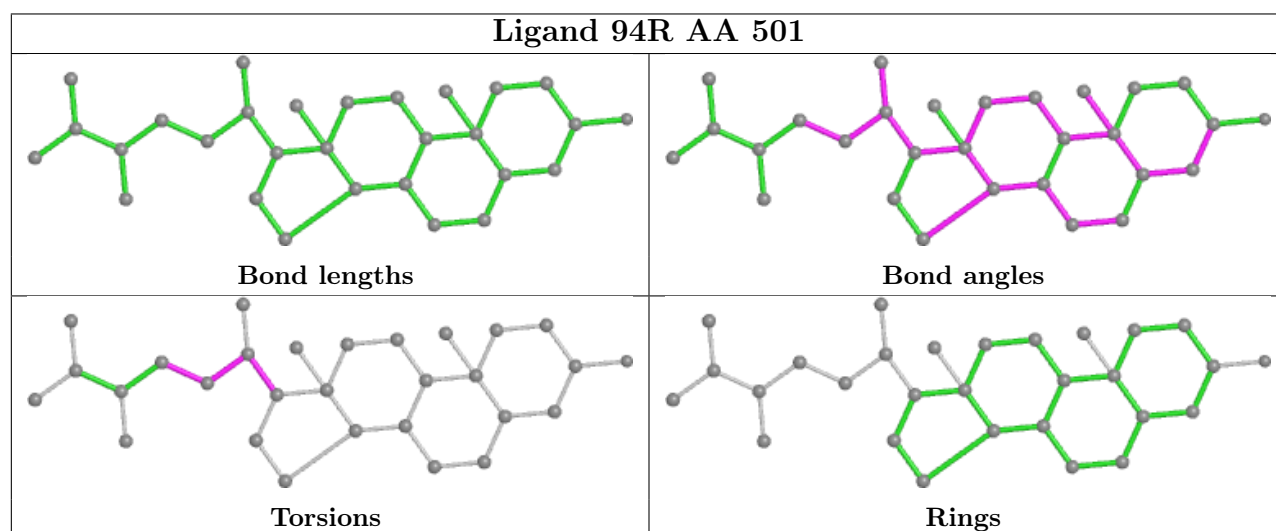
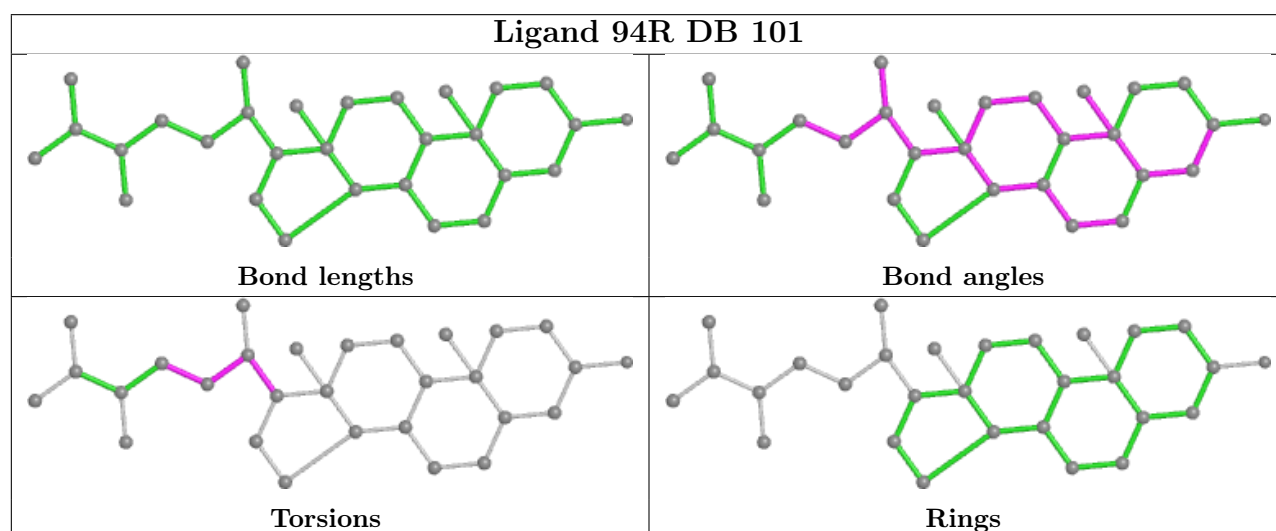
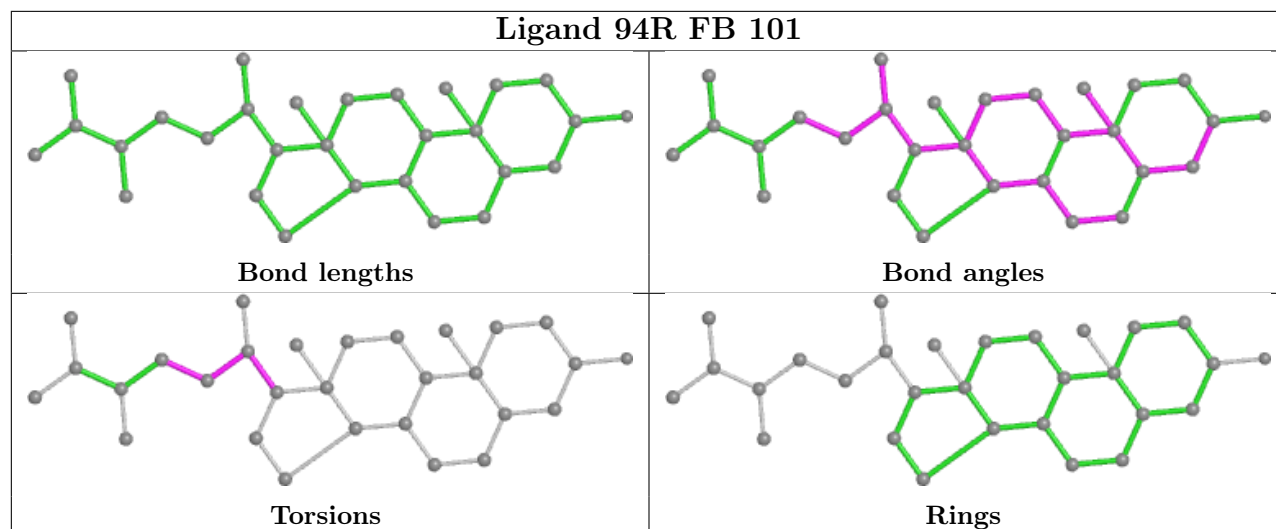
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

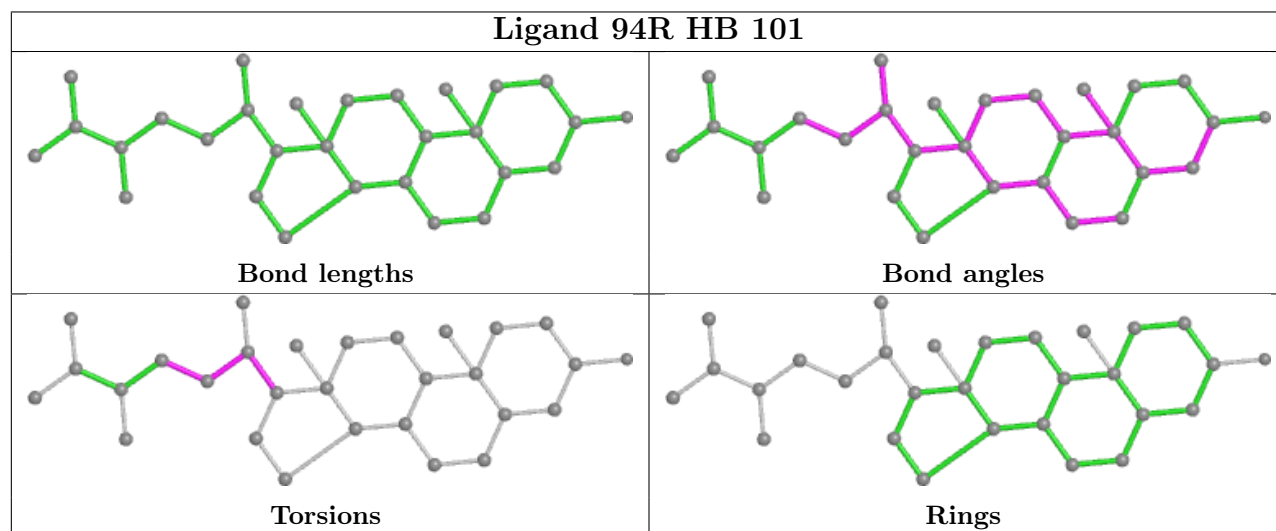
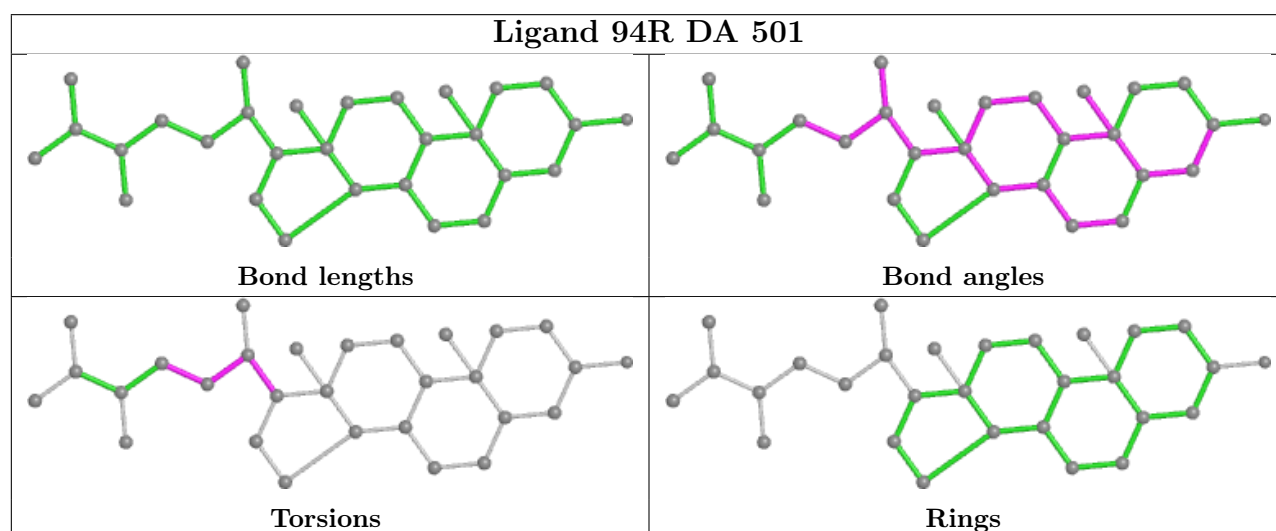
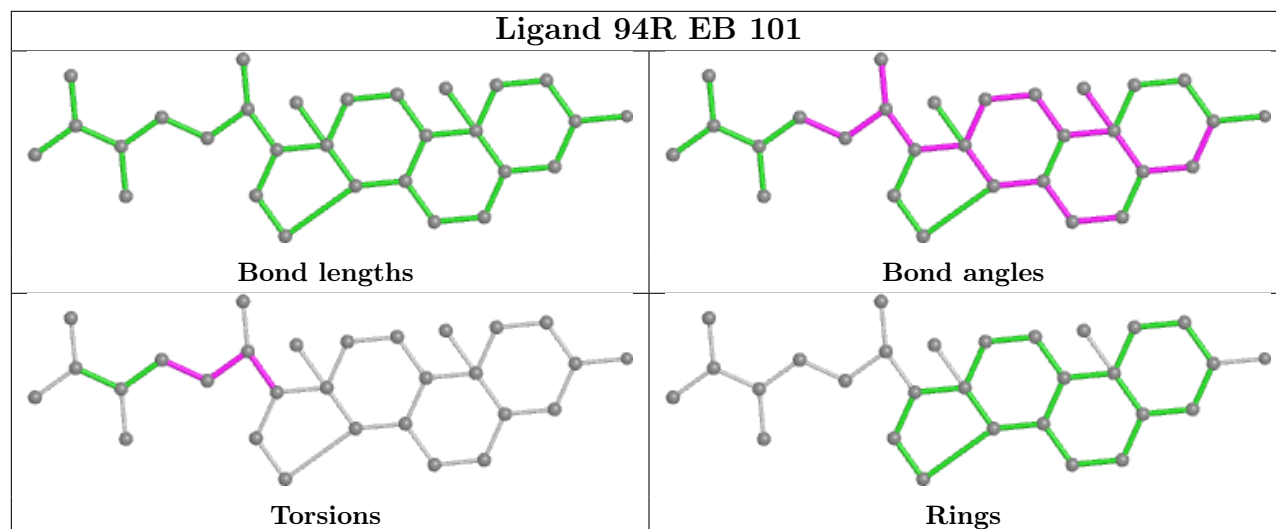
addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.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

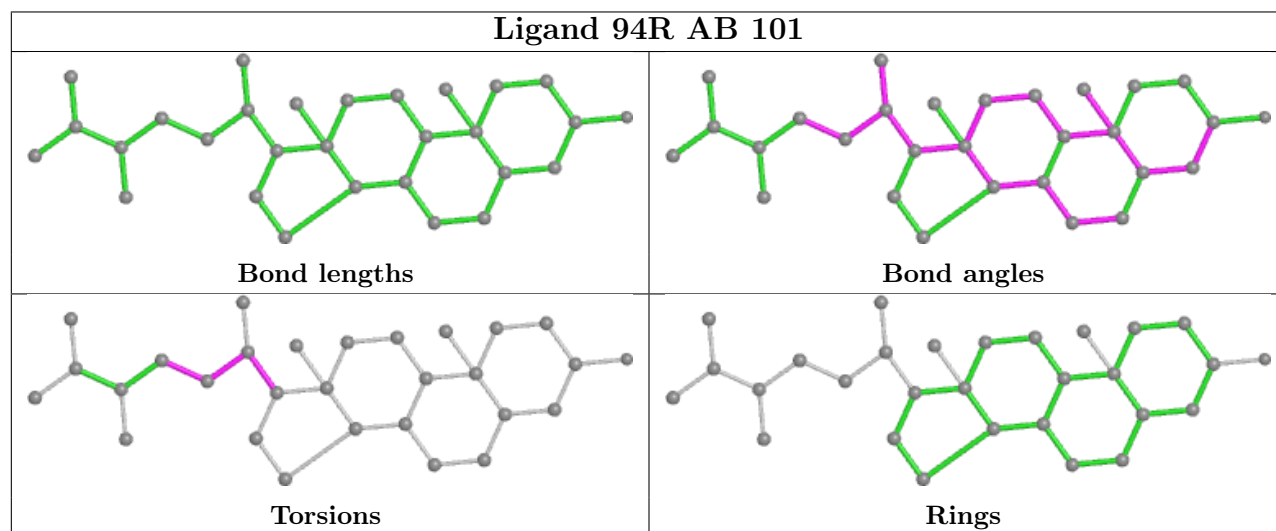
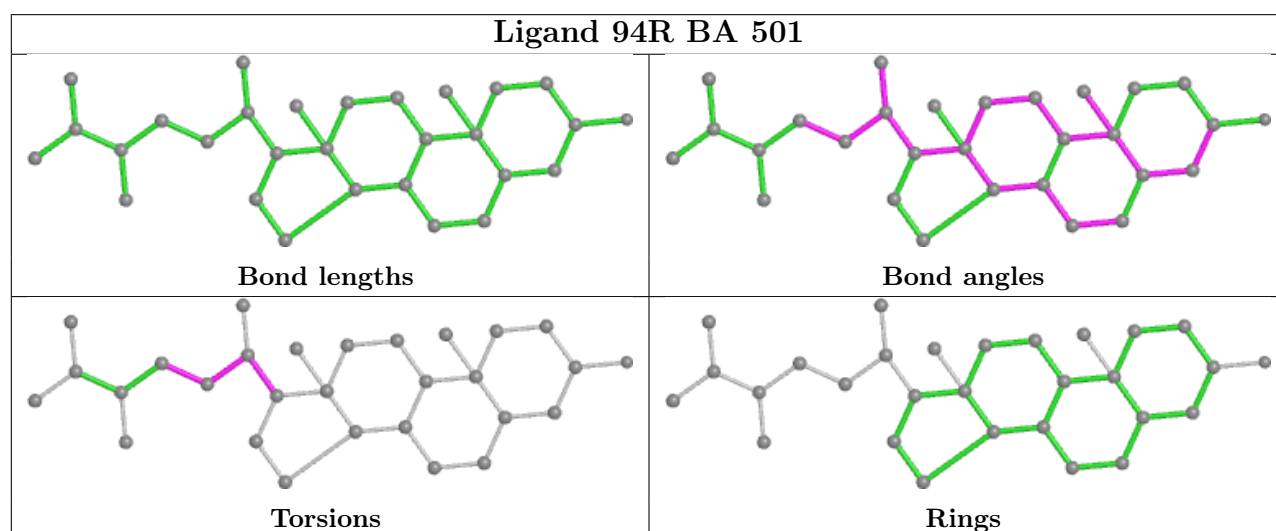
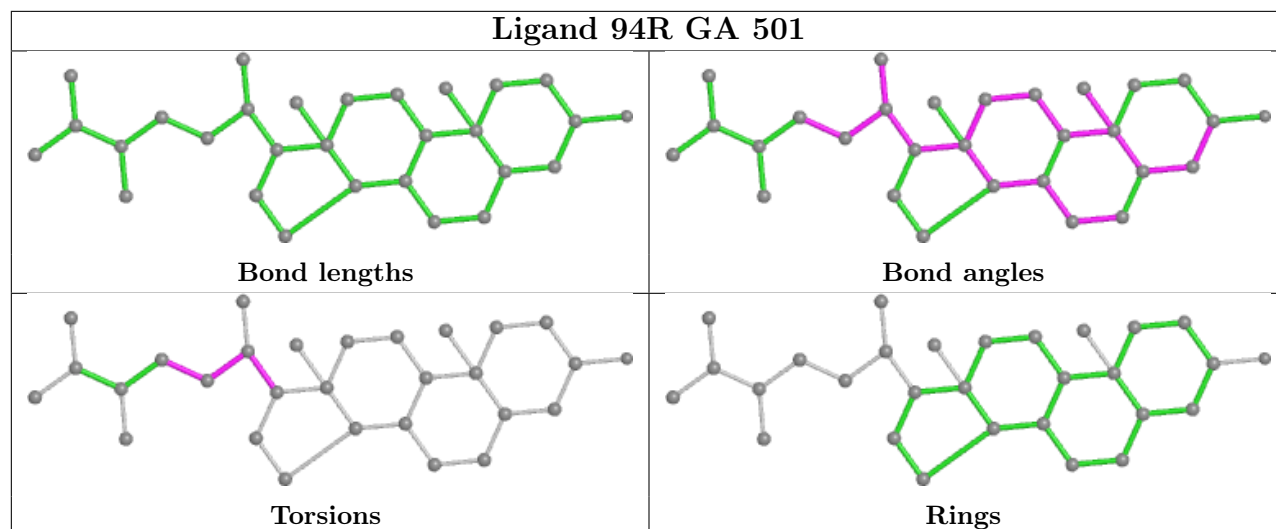


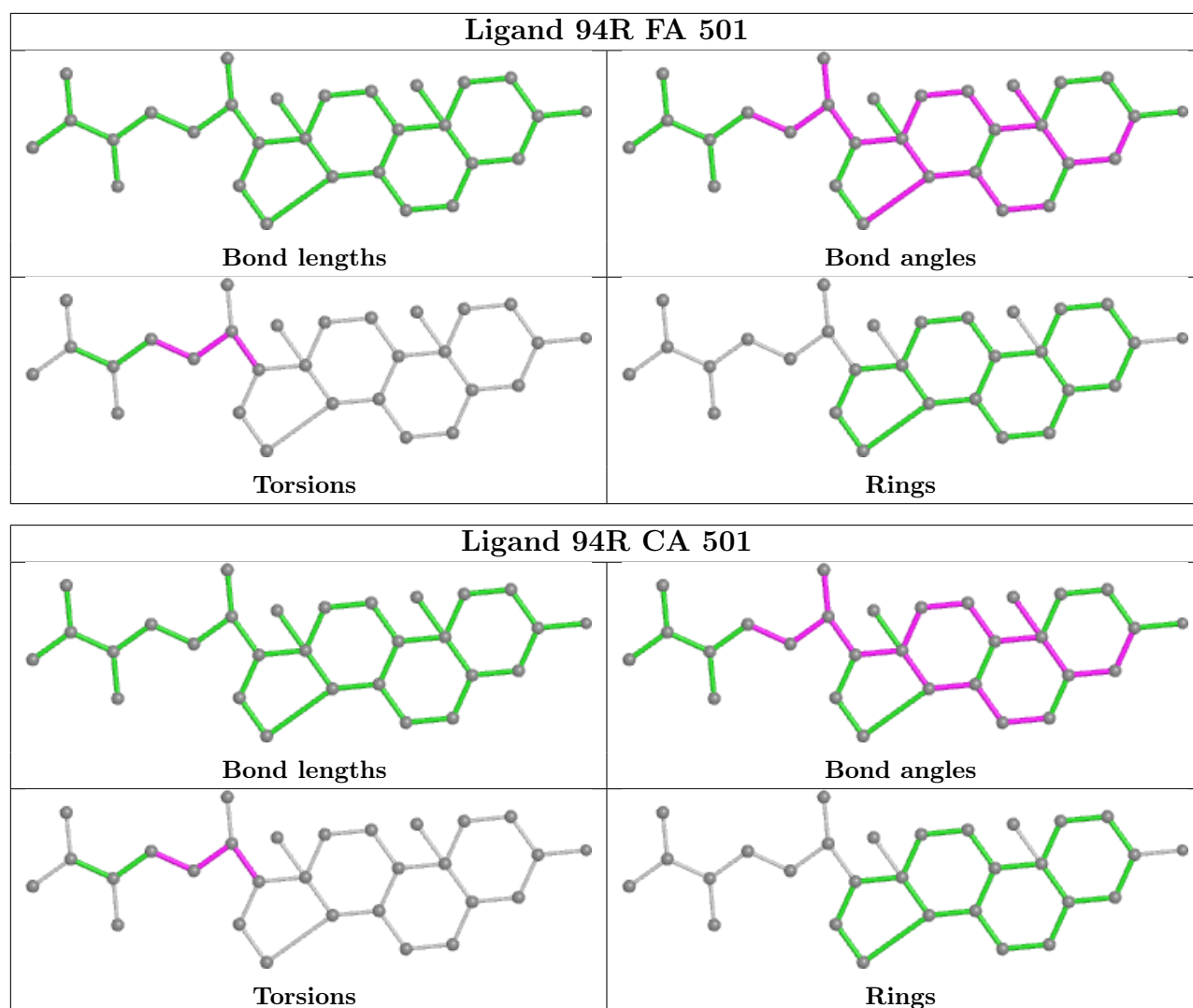












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

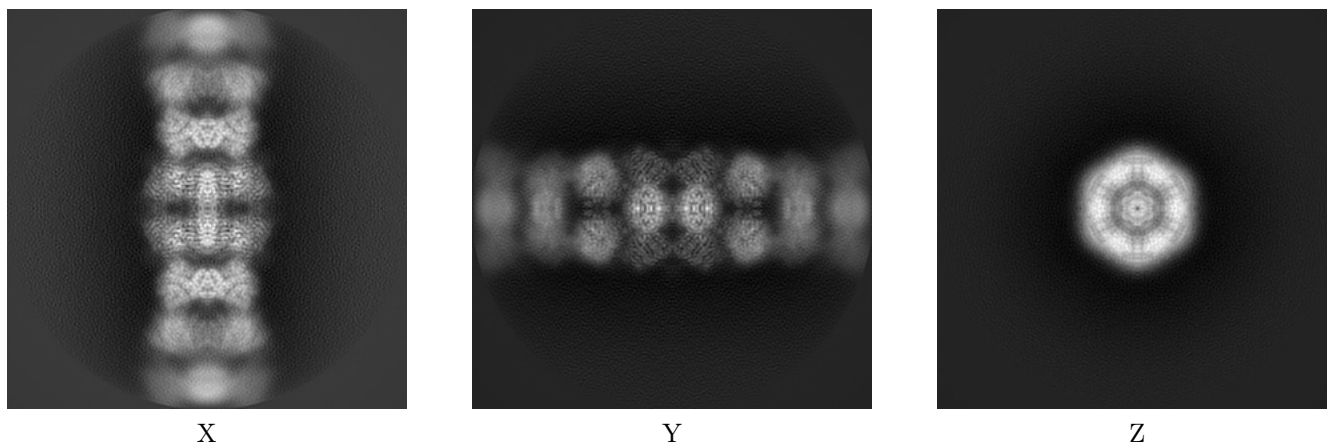
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11892. These allow visual inspection of the internal detail of the map and identification of artifacts.

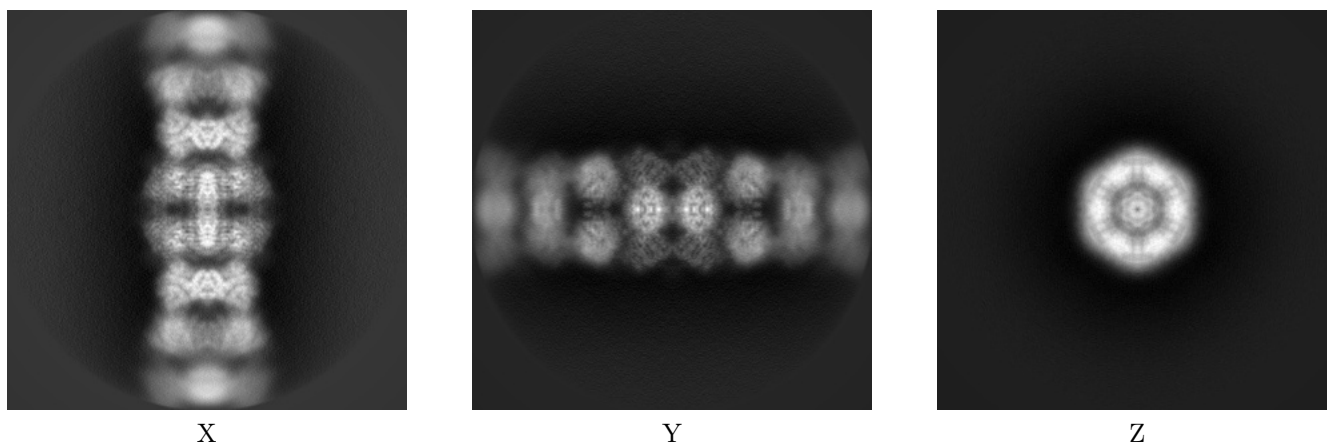
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

#### 6.1.1 Primary map



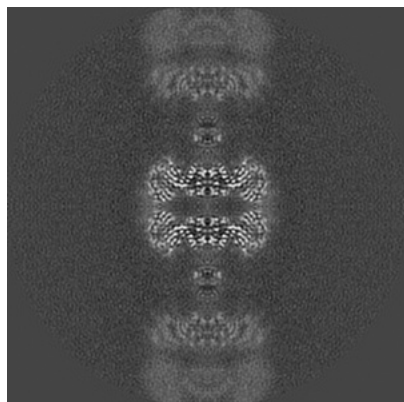
#### 6.1.2 Raw map



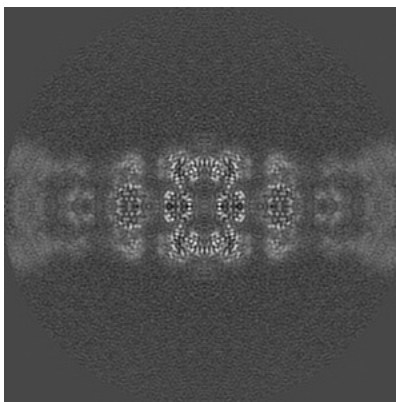
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

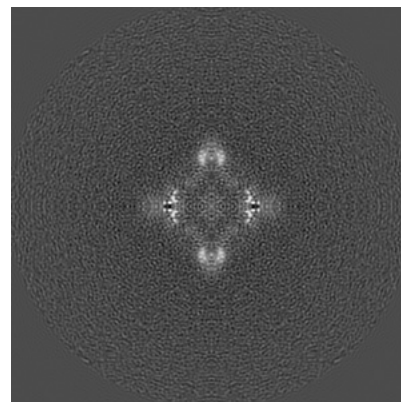
### 6.2.1 Primary map



X Index: 180

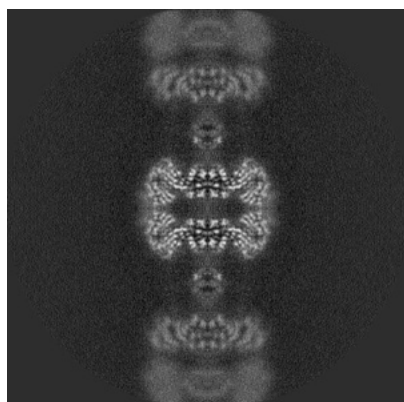


Y Index: 180

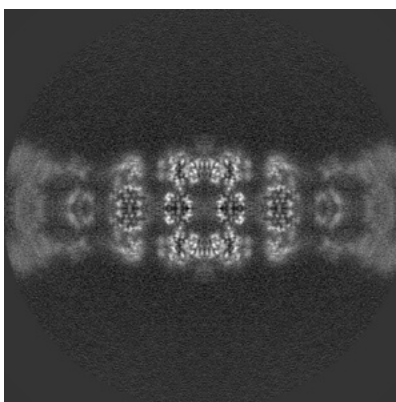


Z Index: 180

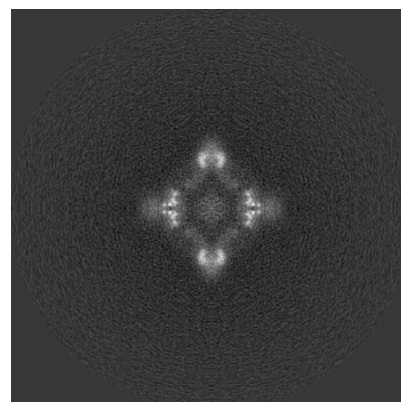
### 6.2.2 Raw map



X Index: 180



Y Index: 180

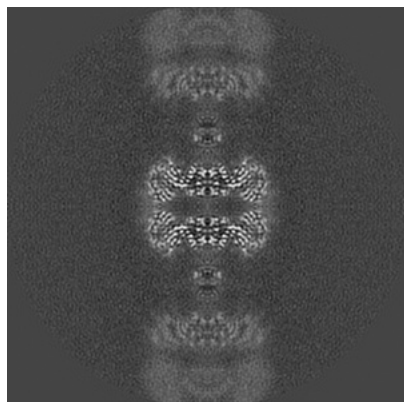


Z Index: 180

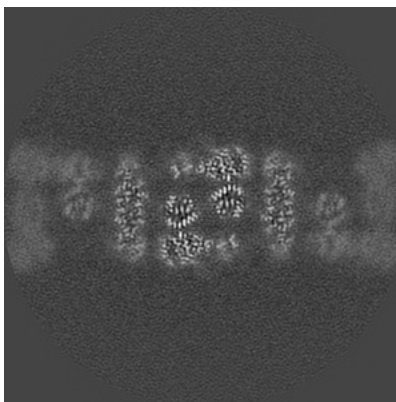
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

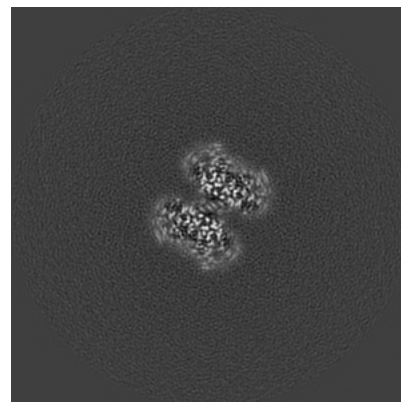
### 6.3.1 Primary map



X Index: 180

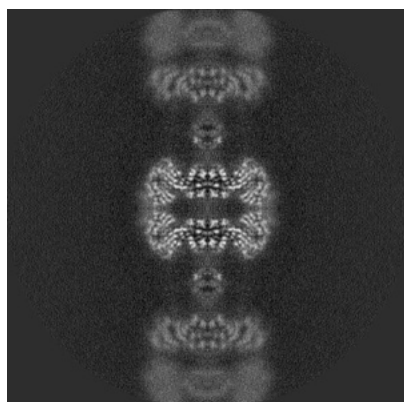


Y Index: 174

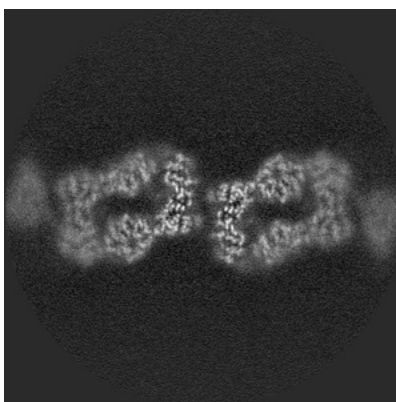


Z Index: 157

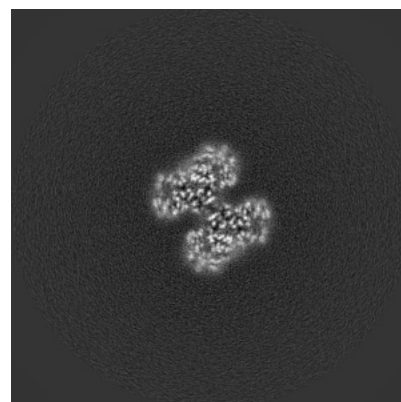
### 6.3.2 Raw map



X Index: 180



Y Index: 208



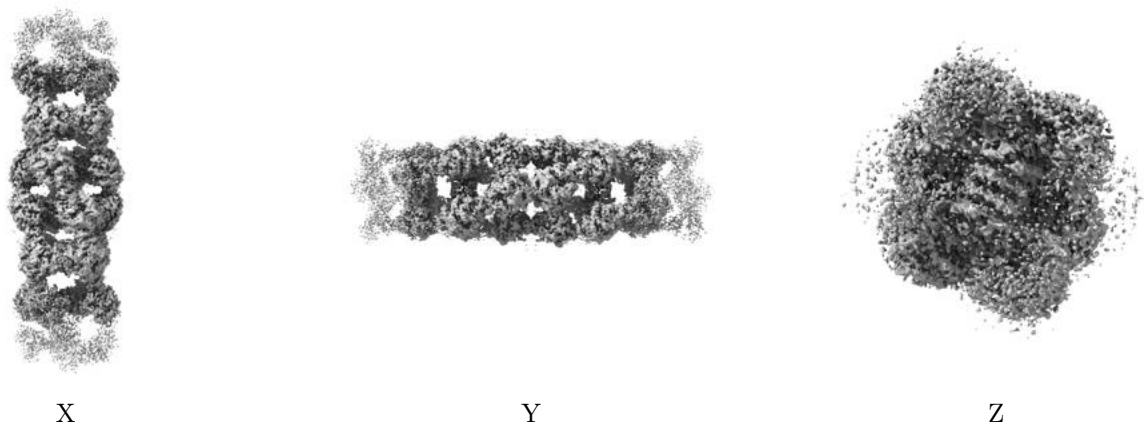
Z Index: 200

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



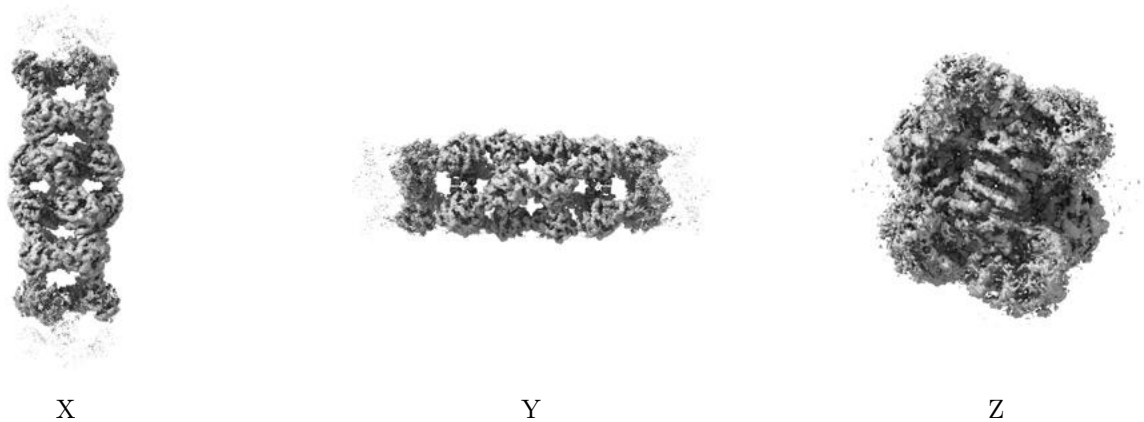
## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.015. 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.

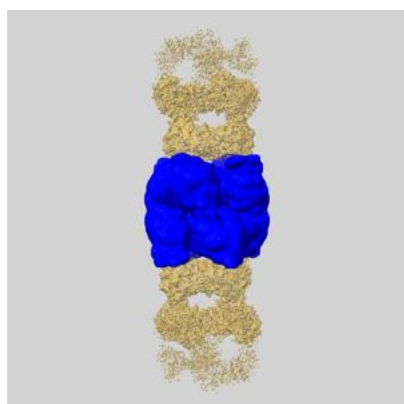
## 6.5 Mask visualisation [i](#)

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

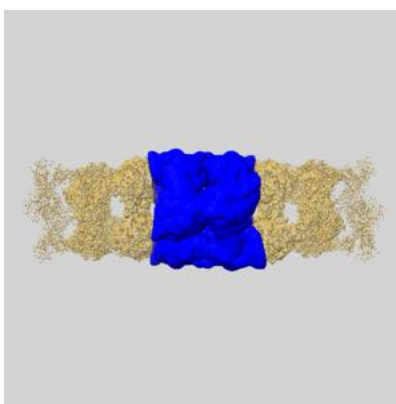
A mask typically either:

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

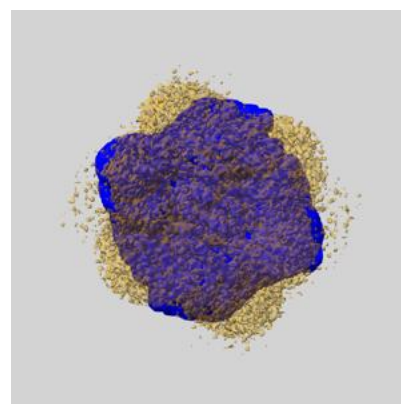
### 6.5.1 emd\_11892\_msk\_1.map [i](#)



X



Y

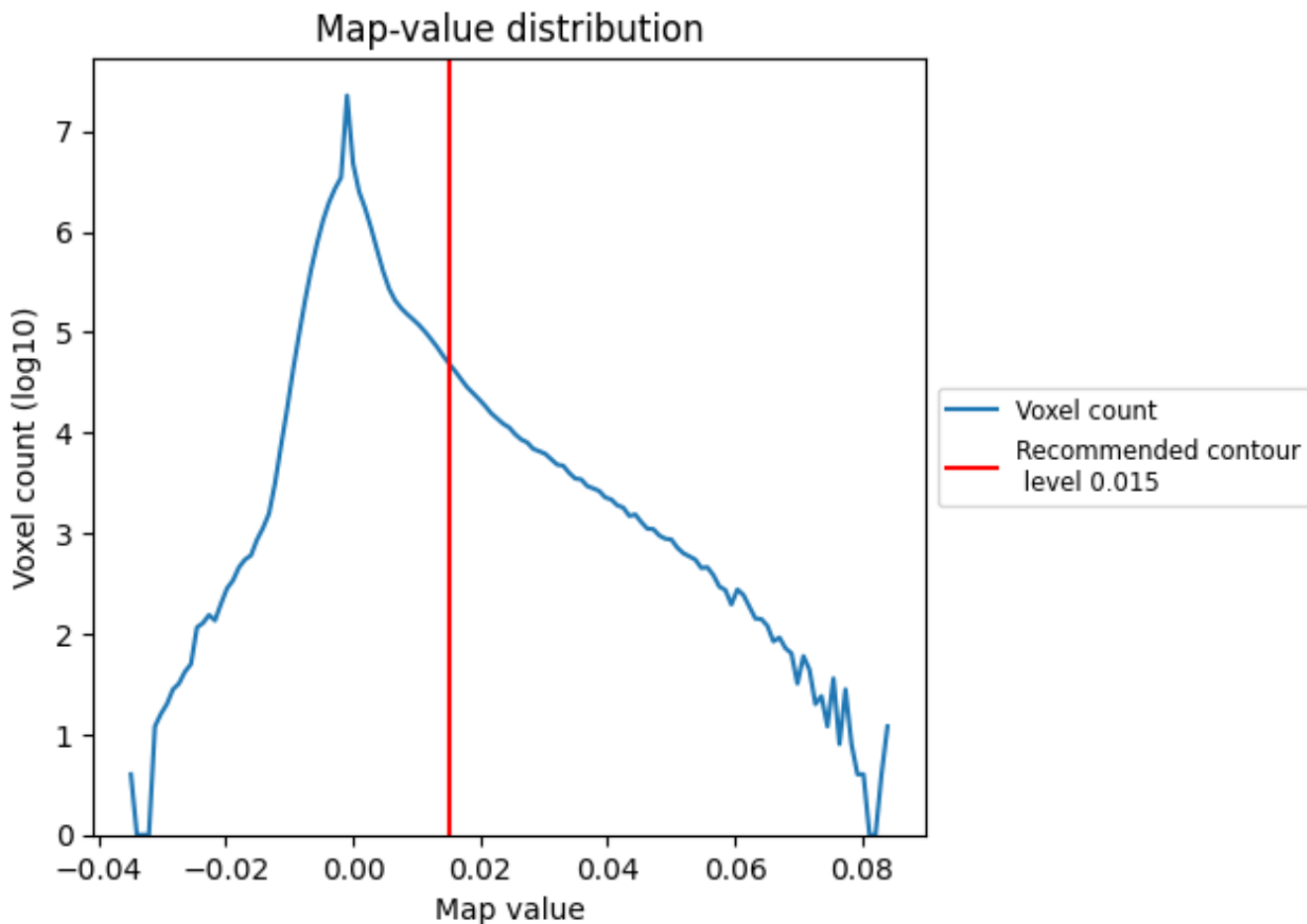


Z

## 7 Map analysis [i](#)

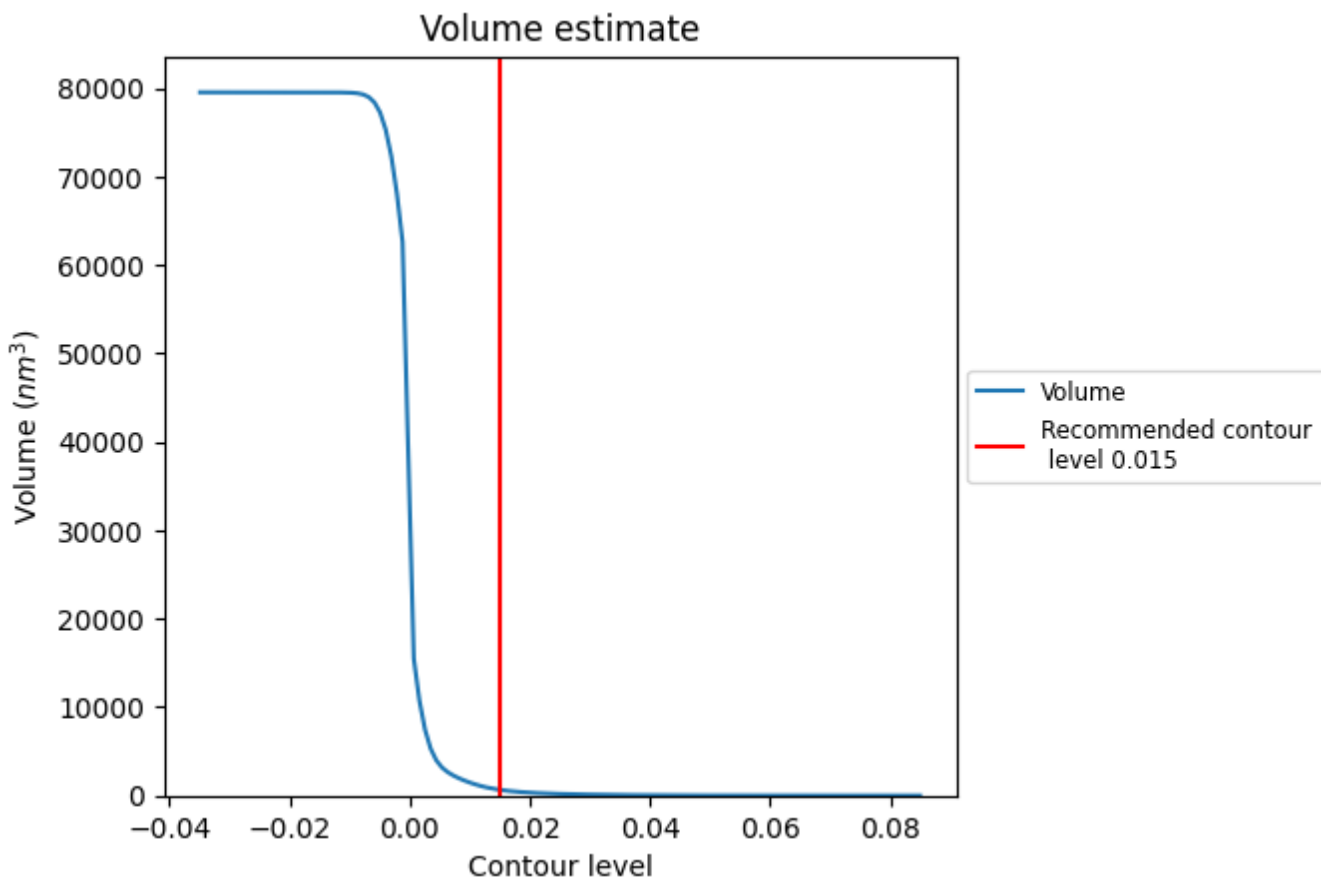
This section contains the results of statistical analysis of the map.

### 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.

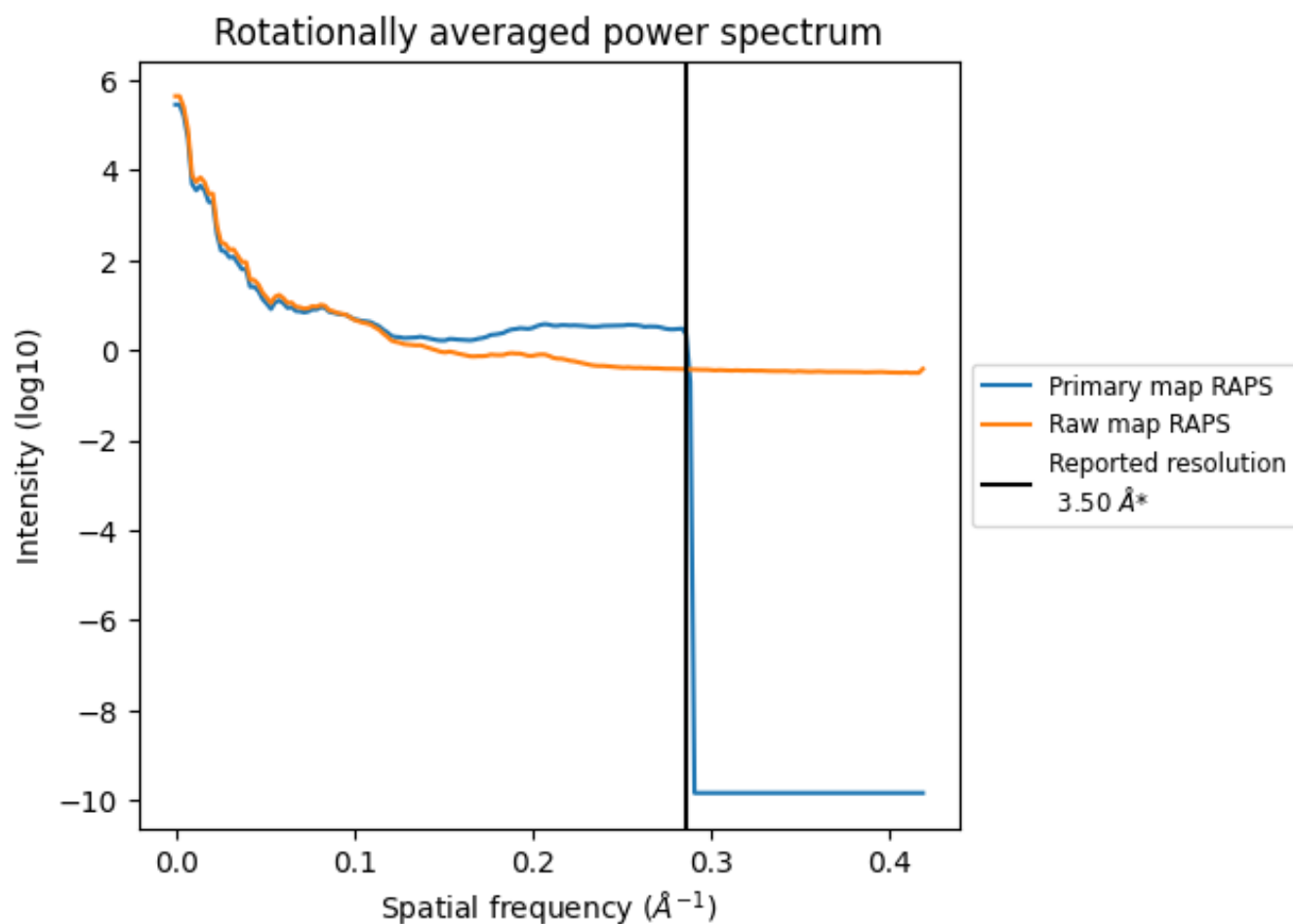
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 649 nm<sup>3</sup>; this corresponds to an approximate mass of 587 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 i

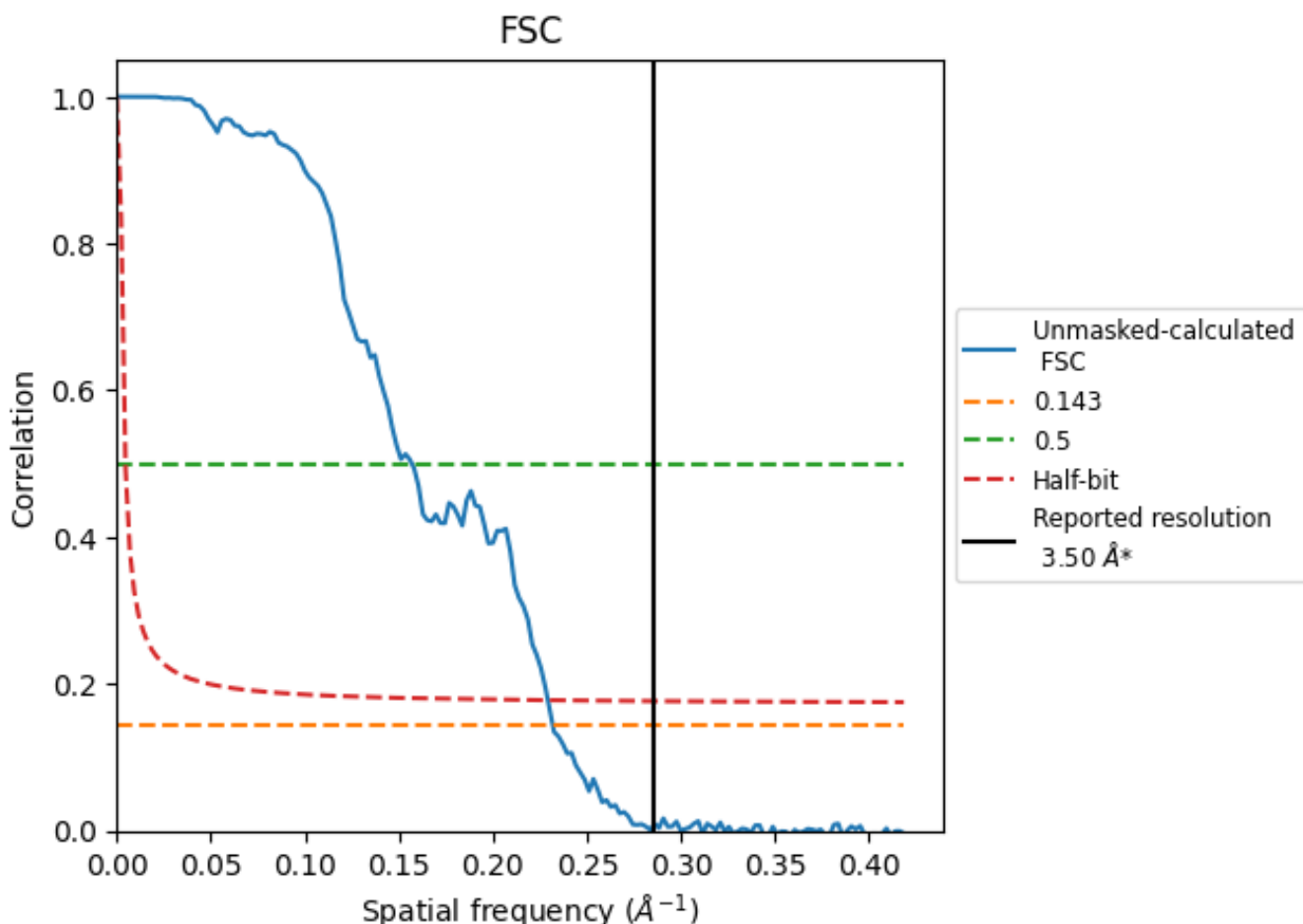


\*Reported resolution corresponds to spatial frequency of 0.286 Å<sup>-1</sup>

## 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 0.286 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

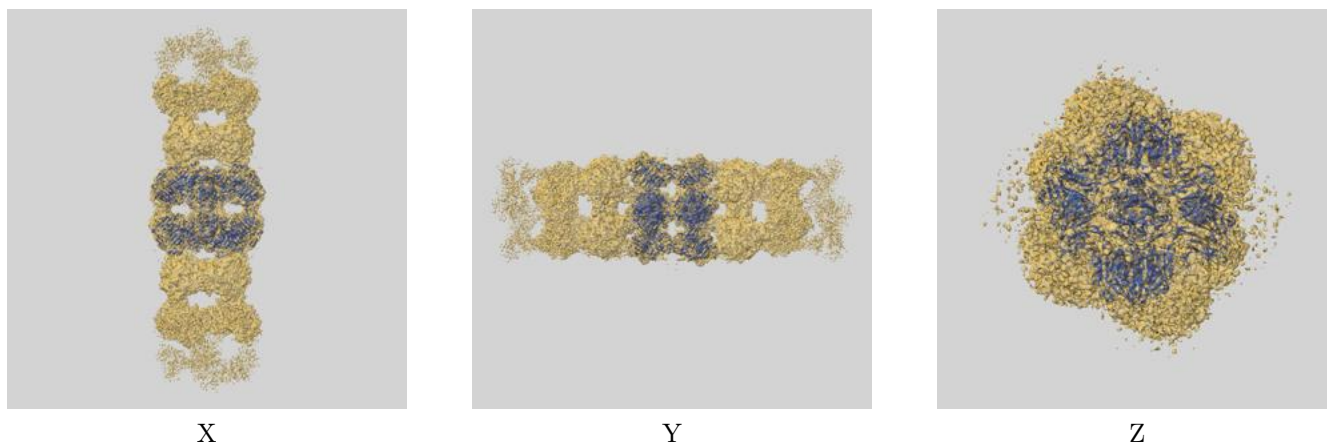
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.31	6.38	4.36

\*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 4.31 differs from the reported value 3.5 by more than 10 %

## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-11892 and PDB model 7ASD. Per-residue inclusion information can be found in section 3 on page 10.

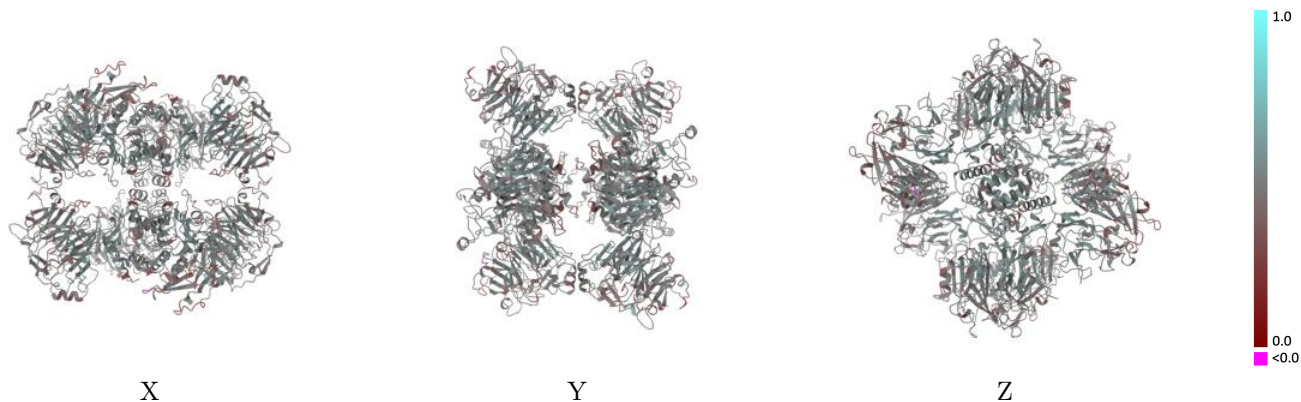
### 9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.015 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.

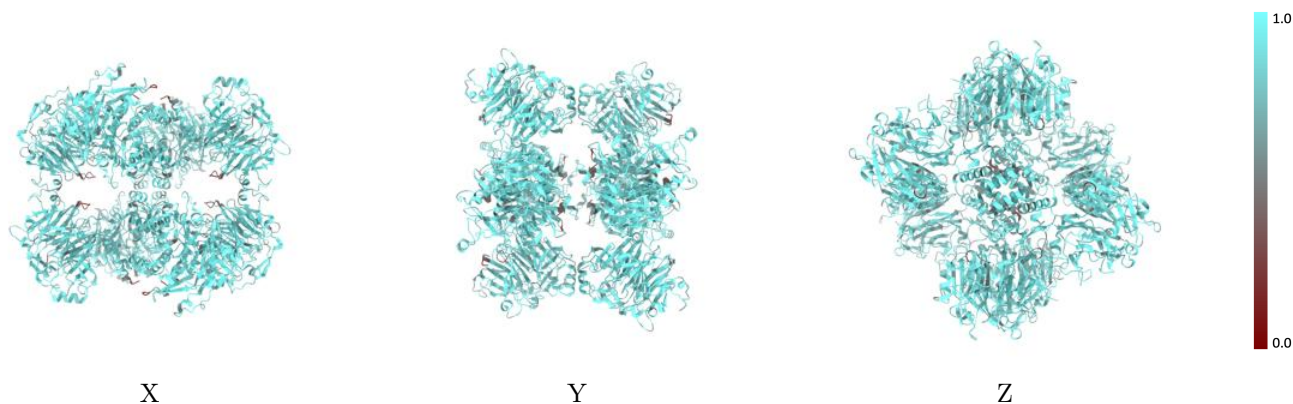


## 9.2 Q-score mapped to coordinate model [\(i\)](#)



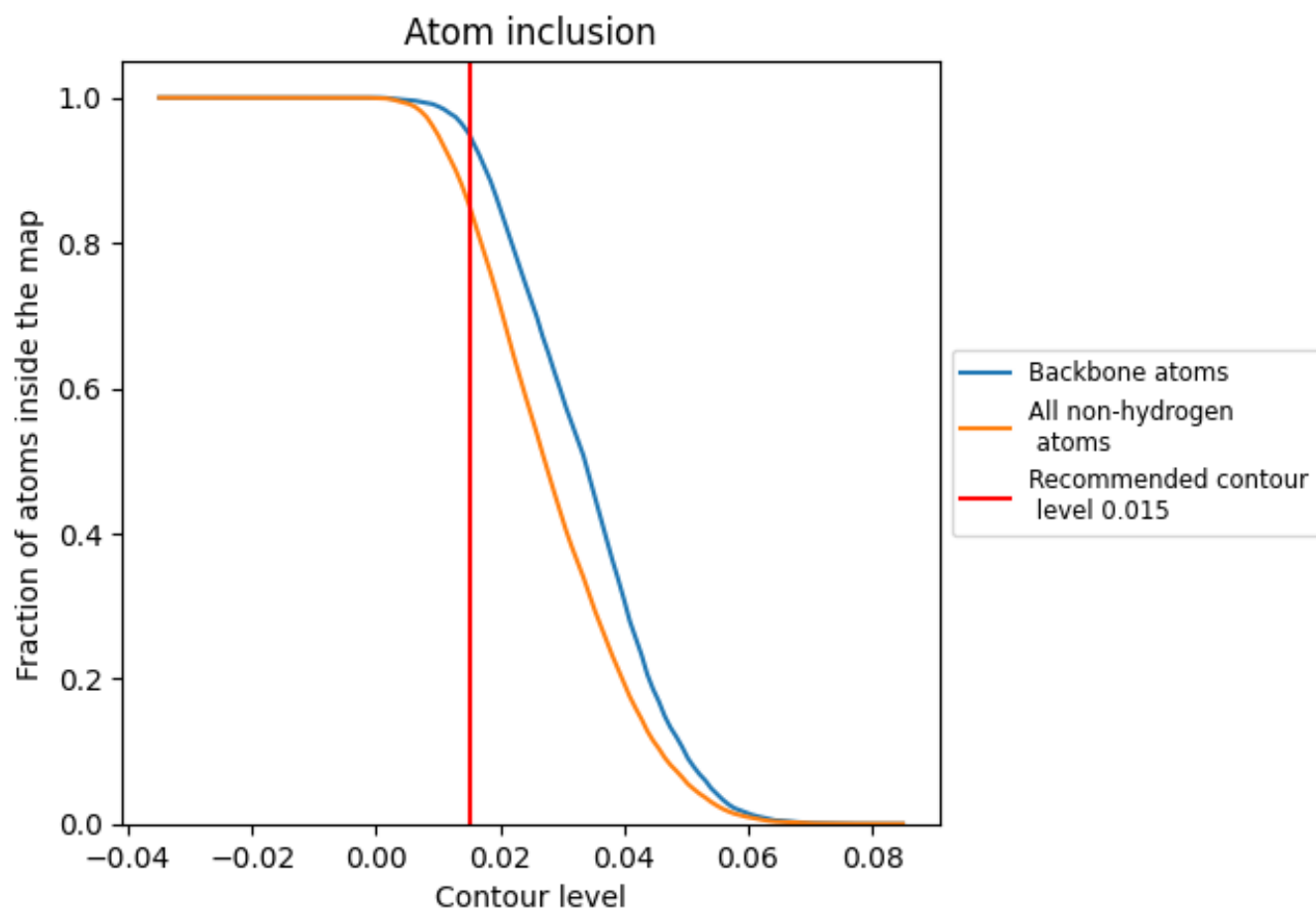
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.015).



































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 85% 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 (0.015) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8526	 0.4640
A	 0.6667	 0.3400
AA	 0.8589	 0.4700
AAc	 0.3571	 0.3430
AB	 0.8315	 0.4890
B	 0.6923	 0.3360
BA	 0.8622	 0.4570
BAc	 0.3929	 0.3780
BB	 0.8567	 0.4820
C	 0.6923	 0.3370
CA	 0.8622	 0.4590
CAC	 0.4286	 0.3590
CB	 0.8567	 0.4860
D	 0.6667	 0.3470
DA	 0.8577	 0.4690
DAC	 0.2857	 0.3260
DB	 0.8343	 0.4870
E	 0.7179	 0.3340
EA	 0.8595	 0.4570
EAC	 0.4286	 0.3720
EB	 0.8567	 0.4820
F	 0.6667	 0.3560
FA	 0.8610	 0.4710
FAC	 0.3571	 0.3440
FB	 0.8315	 0.4870
G	 0.6667	 0.3380
GA	 0.8540	 0.4700
GAC	 0.3214	 0.3380
GB	 0.8371	 0.4920
H	 0.6667	 0.3570
HA	 0.8616	 0.4610
HAC	 0.4286	 0.3730
HB	 0.8511	 0.4890

