



## Full wwPDB EM Validation Report ⓘ

Jun 3, 2024 – 01:48 pm BST

PDB ID : 8P4T  
EMDB ID : EMD-17428  
Title : The spike complex of the Lujo Virus  
Authors : Eilon-Ashkenazy, M.; Diskin, R.  
Deposited on : 2023-05-23  
Resolution : 2.96 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.dev92  
Mogul : 1.8.4, CSD as541be (2020)  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36.2

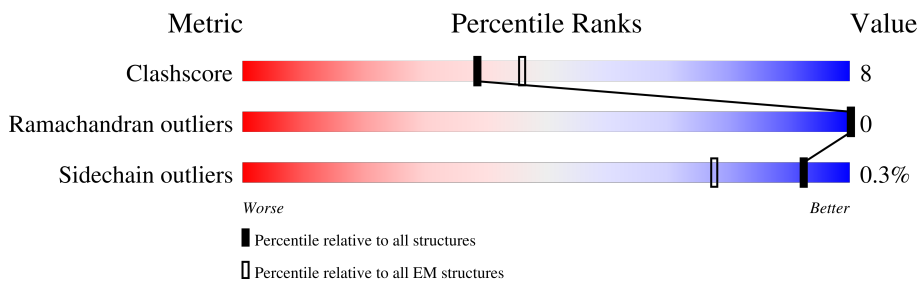
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 2.96 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	A	163	79% (green), 21% (yellow), 5% (red)
1	B	163	77% (green), 22% (yellow), 5% (red)
1	C	163	77% (green), 22% (yellow), 5% (red)
2	a	247	72% (green), 28% (grey), 5% (red)
2	b	247	72% (green), 28% (grey), 5% (red)
2	c	247	72% (green), 28% (grey), 5% (red)
3	SA	58	55% (green), 41% (grey), 38% (red), 5% (yellow)
3	SB	58	55% (green), 41% (grey), 43% (red), 5% (yellow)

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Mol	Chain	Length	Quality of chain
3	SC	58	
4	D	5	
4	G	5	
4	I	5	
5	E	7	
5	H	7	
5	J	7	
6	F	2	
6	K	2	
6	L	2	

## 2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 9976 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 Glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	162	1330	838	235	250	7	0	0
1	B	162	1330	838	235	250	7	0	0
1	C	162	1330	838	235	250	7	0	0

- Molecule 2 is a protein called Glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	a	178	1461	939	244	264	14	0	0
2	b	178	1461	939	244	264	14	0	0
2	c	178	1461	939	244	264	14	0	0

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
a	455	GLY	-	expression tag	UNP C5ILC1
a	456	GLY	-	expression tag	UNP C5ILC1
a	457	ASP	-	expression tag	UNP C5ILC1
a	458	TYR	-	expression tag	UNP C5ILC1
a	459	LYS	-	expression tag	UNP C5ILC1
a	460	ASP	-	expression tag	UNP C5ILC1
a	461	ASP	-	expression tag	UNP C5ILC1
a	462	ASP	-	expression tag	UNP C5ILC1
a	463	ASP	-	expression tag	UNP C5ILC1
a	464	LYS	-	expression tag	UNP C5ILC1
a	465	GLY	-	expression tag	UNP C5ILC1
a	466	SER	-	expression tag	UNP C5ILC1
a	467	GLY	-	expression tag	UNP C5ILC1

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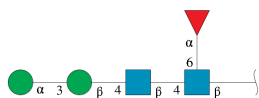
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Chain	Residue	Modelled	Actual	Comment	Reference
a	468	THR	-	expression tag	UNP C5ILC1
b	455	GLY	-	expression tag	UNP C5ILC1
b	456	GLY	-	expression tag	UNP C5ILC1
b	457	ASP	-	expression tag	UNP C5ILC1
b	458	TYR	-	expression tag	UNP C5ILC1
b	459	LYS	-	expression tag	UNP C5ILC1
b	460	ASP	-	expression tag	UNP C5ILC1
b	461	ASP	-	expression tag	UNP C5ILC1
b	462	ASP	-	expression tag	UNP C5ILC1
b	463	ASP	-	expression tag	UNP C5ILC1
b	464	LYS	-	expression tag	UNP C5ILC1
b	465	GLY	-	expression tag	UNP C5ILC1
b	466	SER	-	expression tag	UNP C5ILC1
b	467	GLY	-	expression tag	UNP C5ILC1
b	468	THR	-	expression tag	UNP C5ILC1
c	455	GLY	-	expression tag	UNP C5ILC1
c	456	GLY	-	expression tag	UNP C5ILC1
c	457	ASP	-	expression tag	UNP C5ILC1
c	458	TYR	-	expression tag	UNP C5ILC1
c	459	LYS	-	expression tag	UNP C5ILC1
c	460	ASP	-	expression tag	UNP C5ILC1
c	461	ASP	-	expression tag	UNP C5ILC1
c	462	ASP	-	expression tag	UNP C5ILC1
c	463	ASP	-	expression tag	UNP C5ILC1
c	464	LYS	-	expression tag	UNP C5ILC1
c	465	GLY	-	expression tag	UNP C5ILC1
c	466	SER	-	expression tag	UNP C5ILC1
c	467	GLY	-	expression tag	UNP C5ILC1
c	468	THR	-	expression tag	UNP C5ILC1

- Molecule 3 is a protein called Glycoprotein.

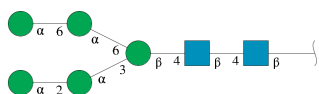
Mol	Chain	Residues	Atoms					AltConf	Trace
3	SA	34	Total	C	N	O	S	0	0
			265	182	39	43	1		
3	SB	34	Total	C	N	O	S	0	0
			265	182	39	43	1		
3	SC	34	Total	C	N	O	S	0	0
			265	182	39	43	1		

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-beta-D-mannopyranos e-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acet amido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
4	D	5	60	34	2	24	0	0
4	G	5	60	34	2	24	0	0
4	I	5	60	34	2	24	0	0

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



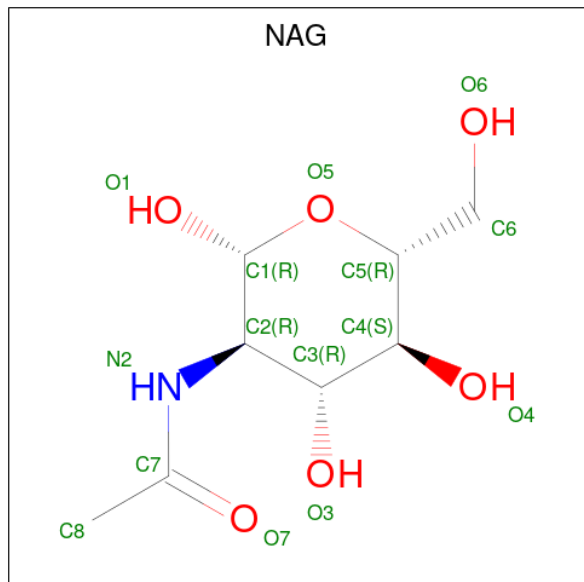
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
5	E	7	83	46	2	35	0	0
5	H	7	83	46	2	35	0	0
5	J	7	83	46	2	35	0	0

- Molecule 6 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		
6	F	2	28	16	2	10	0	0
6	K	2	28	16	2	10	0	0
6	L	2	28	16	2	10	0	0

- Molecule 7 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	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	a	1	Total	C	N	O	0
			14	8	1	5	
7	a	1	Total	C	N	O	0
			14	8	1	5	
7	a	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	C	1	Total	C	N	O	0
			14	8	1	5	

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Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
7	C	1	Total 14	C 8	N 1	O 5	0
7	C	1	Total 14	C 8	N 1	O 5	0
7	C	1	Total 14	C 8	N 1	O 5	0
7	b	1	Total 14	C 8	N 1	O 5	0
7	b	1	Total 14	C 8	N 1	O 5	0
7	b	1	Total 14	C 8	N 1	O 5	0
7	c	1	Total 14	C 8	N 1	O 5	0
7	c	1	Total 14	C 8	N 1	O 5	0
7	c	1	Total 14	C 8	N 1	O 5	0

- Molecule 8 is SODIUM ION (three-letter code: NA) (formula: Na).


Mol	Chain	Residues	Atoms		AltConf
8	A	1	Total 1	Na 1	0

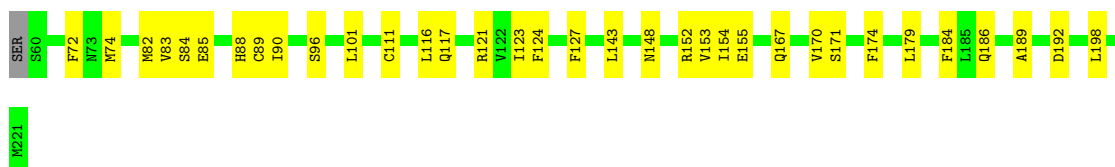


### 3 Residue-property plots


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.

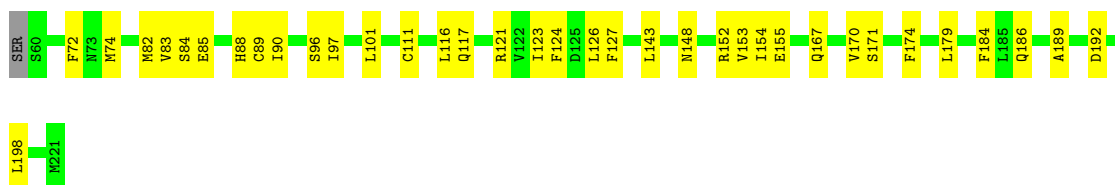
- Molecule 1: Glycoprotein

Chain A:  79% 21%




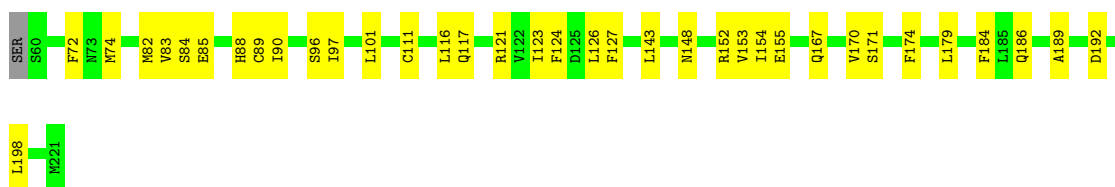
- Molecule 1: Glycoprotein

Chain B:  77% 22%



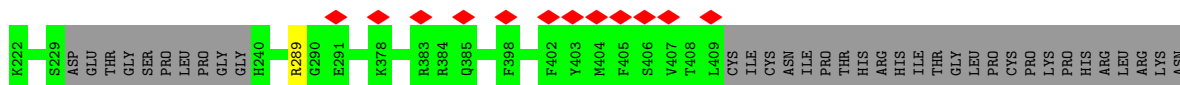
- Molecule 1: Glycoprotein

Chain C:  77% 22%



- Molecule 2: Glycoprotein

Chain a:  5% 72% 28%



GLY THR CYS ALA CYS CYS PHE PHE LYS SER ILE ASN ARG SER THR GLY TRP ALA LYS HIS GLY ASP TYR LYS ASP ASP ASP ASP LYS LYS GLY SER GLY THR

• Molecule 2: Glycoprotein



R222 S229 ASP GLU THR GLY SER PRO LYS PRO ILE ASN ARG SER GLY H240 R289 K378 R383 R384 Q385 F398 F402 Y403 M404 F405 S406 V407 T408 L409 CYS ILE ASN CYS THR PRO HIS ARG HIS ILE THR GLY LEU PRO PRO LYS HIS ARG LEU LYS ASN GLY THR

CYS ALA GLY PHE LYS SER ILE ARG ASN PRO ILE THR GLY TRP ALA LYS HIS GLY ASP TYR LYS ASP ASP ASP LYS GLY SER GLY THR

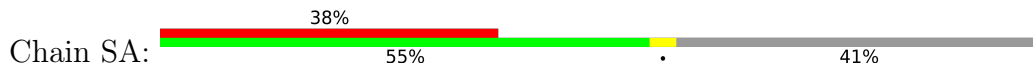
• Molecule 2: Glycoprotein



R222 S229 ASP GLU THR GLY SER PRO LYS PRO ILE ASN ARG SER H240 R289 K378 R383 R384 Q385 L392 F398 F402 Y403 M404 F405 S406 V407 T408 L409 CYS ILE CYS ASN ILE THR THR HIS ARG HIS ILE THR GLY LEU PRO PRO CYS PRO HIS ARG LEU ARG LYS ASN

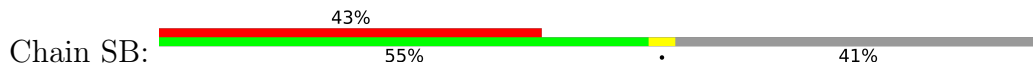
GLY THR CYS ALA CYS CYS PHE PHE LYS SER ILE ARG ASN SER THR TRP ALA LYS HIS GLY ASP TYR LYS ASP ASP ASP LYS GLY SER GLY THR

• Molecule 3: Glycoprotein



MET V2 A3 V4 F5 Q6 A7 I8 P9 E10 I11 I12 N13 E14 A15 I18 I23 I24 M25 F26 T27 L28 I29 K30 G31 V32 F33 N34 L35 TYR LYS SER GLY PHE GLN LEU VAL ILE PHE LEU LEU LEU CYS GLY ARG CYS ASP

• Molecule 3: Glycoprotein



MET V2 A3 V4 F5 Q6 A7 I8 P9 E10 I11 I12 N13 E14 A15 I18 I21 V22 I23 I24 M25 F26 T27 L28 I29 K30 G31 V32 F33 N34 L35 TYR LYS SER GLY PHE GLN LEU VAL ILE PHE LEU LEU LEU CYS GLY ARG CYS ASP

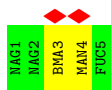
• Molecule 3: Glycoprotein



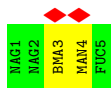
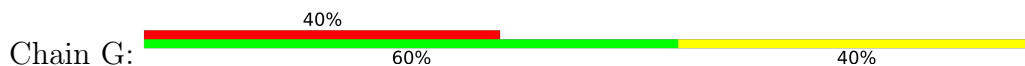
MET V2 A3 V4 F5 Q6 A7 I8 P9 E10 I11 I12 N13 E14 A15 I23 F26 T27 L28 I29 K30 G31 V32 F33 N34 L35 TYR LYS SER GLY PHE GLN LEU VAL ILE PHE LEU LEU LEU CYS GLY ARG CYS ASP

• Molecule 4: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

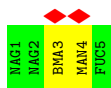
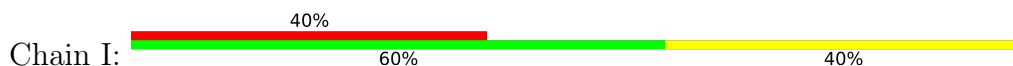




- Molecule 4: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



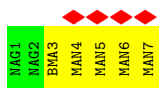
- Molecule 4: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

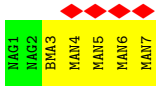


- Molecule 5: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%



- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  100%



- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  100%



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	87868	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	41	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.474	Depositor
Minimum map value	-0.250	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.017	Depositor
Recommended contour level	0.08	Depositor
Map size ( $\text{\AA}$ )	210.944, 210.944, 210.944	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.824, 0.824, 0.824	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, FUC, MAN, NA, BMA

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	0/1359	0.50	0/1835
1	B	0.25	0/1359	0.50	0/1835
1	C	0.25	0/1359	0.50	0/1835
2	a	0.24	0/1492	0.46	0/2006
2	b	0.24	0/1492	0.46	0/2006
2	c	0.24	0/1492	0.46	0/2006
3	SA	0.24	0/268	0.33	0/365
3	SB	0.24	0/268	0.33	0/365
3	SC	0.24	0/268	0.33	0/365
All	All	0.24	0/9357	0.47	0/12618

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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	1330	0	1295	22	0
1	B	1330	0	1295	23	0
1	C	1330	0	1295	25	0
2	a	1461	0	1454	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	b	1461	0	1454	0	0
2	c	1461	0	1454	0	0
3	SA	265	0	295	1	0
3	SB	265	0	295	1	0
3	SC	265	0	295	1	0
4	D	60	0	52	0	0
4	G	60	0	52	0	0
4	I	60	0	52	0	0
5	E	83	0	70	0	0
5	H	83	0	70	0	0
5	J	83	0	70	0	0
6	F	28	0	25	0	0
6	K	28	0	25	0	0
6	L	28	0	25	0	0
7	A	56	0	52	0	0
7	B	56	0	52	0	0
7	C	56	0	52	0	0
7	a	42	0	39	0	0
7	b	42	0	39	0	0
7	c	42	0	39	0	0
8	A	1	0	0	0	0
All	All	9976	0	9846	73	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 8.

All (73) 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:124:PHE:HA	1:B:127:PHE:CE1	2.35	0.62
1:C:124:PHE:HA	1:C:127:PHE:CE1	2.35	0.62
1:A:124:PHE:HA	1:A:127:PHE:CE1	2.34	0.62
1:A:155:GLU:HG3	1:A:184:PHE:HB3	1.87	0.57
1:C:155:GLU:HG3	1:C:184:PHE:HB3	1.87	0.56
1:B:155:GLU:HG3	1:B:184:PHE:HB3	1.87	0.55
1:A:72:PHE:HB2	1:A:74:MET:HE2	1.90	0.53
1:B:153:VAL:HG13	1:B:186:GLN:HG3	1.91	0.52
1:B:170:VAL:O	1:B:174:PHE:N	2.36	0.52
1:A:153:VAL:HG13	1:A:186:GLN:HG3	1.91	0.51
1:C:153:VAL:HG13	1:C:186:GLN:HG3	1.91	0.51
1:A:101:LEU:HD13	1:A:179:LEU:HD22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:72:PHE:HB2	1:C:74:MET:HE2	1.93	0.51
1:C:170:VAL:O	1:C:174:PHE:N	2.36	0.51
1:A:170:VAL:O	1:A:174:PHE:N	2.36	0.50
1:A:88:HIS:HA	1:A:174:PHE:CE1	2.47	0.50
1:B:101:LEU:HD13	1:B:179:LEU:HD22	1.93	0.50
1:B:88:HIS:HA	1:B:174:PHE:CE1	2.47	0.50
1:C:88:HIS:HA	1:C:174:PHE:CE1	2.47	0.49
1:C:111:CYS:SG	1:C:143:LEU:HG	2.53	0.49
1:A:84:SER:OG	1:A:89:CYS:HB3	2.13	0.49
1:A:111:CYS:SG	1:A:143:LEU:HG	2.53	0.49
1:B:84:SER:OG	1:B:89:CYS:HB3	2.13	0.49
1:C:101:LEU:HD13	1:C:179:LEU:HD22	1.93	0.49
1:B:111:CYS:SG	1:B:143:LEU:HG	2.52	0.49
1:A:89:CYS:SG	1:A:96:SER:HB2	2.53	0.48
1:C:84:SER:OG	1:C:89:CYS:HB3	2.13	0.48
1:B:116:LEU:HD21	1:B:154:ILE:HD11	1.96	0.48
1:C:89:CYS:SG	1:C:96:SER:HB2	2.53	0.48
1:B:72:PHE:HB2	1:B:74:MET:HE2	1.95	0.48
1:A:167:GLN:NE2	1:A:171:SER:OG	2.47	0.47
1:C:116:LEU:HD21	1:C:154:ILE:HD11	1.96	0.47
1:C:167:GLN:NE2	1:C:171:SER:OG	2.47	0.47
1:B:89:CYS:SG	1:B:96:SER:HB2	2.53	0.47
1:A:116:LEU:HD21	1:A:154:ILE:HD11	1.96	0.47
1:B:167:GLN:NE2	1:B:171:SER:OG	2.47	0.46
1:B:167:GLN:HA	1:B:170:VAL:HG22	1.98	0.46
1:C:152:ARG:HB2	1:C:189:ALA:HB3	1.98	0.46
1:B:85:GLU:HA	1:B:174:PHE:HD1	1.82	0.45
1:A:167:GLN:HA	1:A:170:VAL:HG22	1.98	0.45
1:A:85:GLU:HA	1:A:174:PHE:HD1	1.82	0.45
1:B:152:ARG:HB2	1:B:189:ALA:HB3	1.98	0.45
1:B:123:ILE:O	1:B:127:PHE:CD1	2.70	0.44
1:C:167:GLN:HA	1:C:170:VAL:HG22	1.98	0.44
1:C:85:GLU:HA	1:C:174:PHE:HD1	1.82	0.44
1:A:152:ARG:HB2	1:A:189:ALA:HB3	1.98	0.44
1:A:123:ILE:O	1:A:127:PHE:CD1	2.70	0.44
3:SB:6:GLN:O	3:SB:9:PRO:HD2	2.18	0.43
3:SA:6:GLN:O	3:SA:9:PRO:HD2	2.18	0.43
1:C:123:ILE:O	1:C:127:PHE:CD1	2.70	0.43
3:SC:6:GLN:O	3:SC:9:PRO:HD2	2.18	0.43
1:A:148:ASN:O	1:A:152:ARG:HG2	2.19	0.43
1:C:117:GLN:HG2	1:C:121:ARG:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:VAL:HG22	1:C:90:ILE:HG12	2.01	0.42
1:B:82:MET:CB	1:B:198:LEU:HB3	2.50	0.42
1:B:148:ASN:O	1:B:152:ARG:HG2	2.19	0.42
1:C:148:ASN:O	1:C:152:ARG:HG2	2.19	0.42
1:B:83:VAL:HG22	1:B:90:ILE:HG12	2.01	0.42
1:A:82:MET:CB	1:A:198:LEU:HB3	2.50	0.42
1:A:117:GLN:HG2	1:A:121:ARG:HD3	2.01	0.42
1:A:83:VAL:HG22	1:A:90:ILE:HG12	2.01	0.41
1:B:192:ASP:N	1:B:192:ASP:OD1	2.53	0.41
1:C:192:ASP:N	1:C:192:ASP:OD1	2.53	0.41
1:A:192:ASP:OD1	1:A:192:ASP:N	2.53	0.41
1:B:170:VAL:HG21	1:B:179:LEU:HG	2.02	0.41
1:C:82:MET:CB	1:C:198:LEU:HB3	2.50	0.41
1:B:117:GLN:HG2	1:B:121:ARG:HD3	2.01	0.41
1:C:97:ILE:HD13	1:C:126:LEU:HB3	2.03	0.41
1:B:97:ILE:HD13	1:B:126:LEU:HB3	2.02	0.41
1:A:170:VAL:HG21	1:A:179:LEU:HG	2.02	0.40
1:C:170:VAL:HG21	1:C:179:LEU:HG	2.02	0.40
1:C:85:GLU:HA	1:C:174:PHE:CD1	2.57	0.40
1:C:123:ILE:HD12	1:C:189:ALA:HB2	2.04	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	160/163 (98%)	156 (98%)	4 (2%)	0	100	100
1	B	160/163 (98%)	156 (98%)	4 (2%)	0	100	100
1	C	160/163 (98%)	156 (98%)	4 (2%)	0	100	100
2	a	174/247 (70%)	170 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	b	174/247 (70%)	170 (98%)	4 (2%)	0	100	100
2	c	174/247 (70%)	170 (98%)	4 (2%)	0	100	100
3	SA	32/58 (55%)	31 (97%)	1 (3%)	0	100	100
3	SB	32/58 (55%)	31 (97%)	1 (3%)	0	100	100
3	SC	32/58 (55%)	31 (97%)	1 (3%)	0	100	100
All	All	1098/1404 (78%)	1071 (98%)	27 (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	A	156/157 (99%)	156 (100%)	0	100	100
1	B	156/157 (99%)	156 (100%)	0	100	100
1	C	156/157 (99%)	156 (100%)	0	100	100
2	a	164/220 (74%)	163 (99%)	1 (1%)	86	94
2	b	164/220 (74%)	163 (99%)	1 (1%)	86	94
2	c	164/220 (74%)	163 (99%)	1 (1%)	86	94
3	SA	30/51 (59%)	30 (100%)	0	100	100
3	SB	30/51 (59%)	30 (100%)	0	100	100
3	SC	30/51 (59%)	30 (100%)	0	100	100
All	All	1050/1284 (82%)	1047 (100%)	3 (0%)	92	97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	a	289	ARG
2	b	289	ARG
2	c	289	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	a	385	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](#)

42 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	D	1	1,4	14,14,15	0.77	0	17,19,21	0.90	0
4	NAG	D	2	4	14,14,15	0.72	0	17,19,21	0.80	0
4	BMA	D	3	4	11,11,12	0.84	0	15,15,17	2.13	4 (26%)
4	MAN	D	4	4	11,11,12	0.69	0	15,15,17	1.25	1 (6%)
4	FUC	D	5	4	10,10,11	0.81	0	14,14,16	0.94	0
5	NAG	E	1	5,1	14,14,15	0.75	0	17,19,21	0.85	0
5	NAG	E	2	5	14,14,15	0.75	0	17,19,21	0.80	0
5	BMA	E	3	5	11,11,12	0.82	0	15,15,17	2.34	5 (33%)
5	MAN	E	4	5	11,11,12	0.70	0	15,15,17	1.22	1 (6%)
5	MAN	E	5	5	11,11,12	0.74	0	15,15,17	1.07	1 (6%)
5	MAN	E	6	5	11,11,12	0.78	0	15,15,17	1.19	1 (6%)
5	MAN	E	7	5	11,11,12	0.69	0	15,15,17	1.29	1 (6%)
6	NAG	F	1	6,2	14,14,15	0.71	0	17,19,21	0.97	0
6	NAG	F	2	6	14,14,15	0.71	0	17,19,21	0.83	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NAG	G	1	1,4	14,14,15	0.77	0	17,19,21	0.89	0
4	NAG	G	2	4	14,14,15	0.72	0	17,19,21	0.80	0
4	BMA	G	3	4	11,11,12	0.84	0	15,15,17	2.13	4 (26%)
4	MAN	G	4	4	11,11,12	0.69	0	15,15,17	1.25	1 (6%)
4	FUC	G	5	4	10,10,11	0.81	0	14,14,16	0.94	0
5	NAG	H	1	5,1	14,14,15	0.75	0	17,19,21	0.85	0
5	NAG	H	2	5	14,14,15	0.75	0	17,19,21	0.80	0
5	BMA	H	3	5	11,11,12	0.82	0	15,15,17	2.34	5 (33%)
5	MAN	H	4	5	11,11,12	0.70	0	15,15,17	1.22	1 (6%)
5	MAN	H	5	5	11,11,12	0.74	0	15,15,17	1.07	1 (6%)
5	MAN	H	6	5	11,11,12	0.78	0	15,15,17	1.19	1 (6%)
5	MAN	H	7	5	11,11,12	0.69	0	15,15,17	1.29	1 (6%)
4	NAG	I	1	1,4	14,14,15	0.77	0	17,19,21	0.89	0
4	NAG	I	2	4	14,14,15	0.72	0	17,19,21	0.80	0
4	BMA	I	3	4	11,11,12	0.84	0	15,15,17	2.13	4 (26%)
4	MAN	I	4	4	11,11,12	0.69	0	15,15,17	1.25	1 (6%)
4	FUC	I	5	4	10,10,11	0.81	0	14,14,16	0.94	0
5	NAG	J	1	5,1	14,14,15	0.75	0	17,19,21	0.85	0
5	NAG	J	2	5	14,14,15	0.75	0	17,19,21	0.80	0
5	BMA	J	3	5	11,11,12	0.82	0	15,15,17	2.34	5 (33%)
5	MAN	J	4	5	11,11,12	0.70	0	15,15,17	1.22	1 (6%)
5	MAN	J	5	5	11,11,12	0.74	0	15,15,17	1.07	1 (6%)
5	MAN	J	6	5	11,11,12	0.78	0	15,15,17	1.19	1 (6%)
5	MAN	J	7	5	11,11,12	0.69	0	15,15,17	1.29	1 (6%)
6	NAG	K	1	6,2	14,14,15	0.71	0	17,19,21	0.97	0
6	NAG	K	2	6	14,14,15	0.71	0	17,19,21	0.83	0
6	NAG	L	1	6,2	14,14,15	0.71	0	17,19,21	0.97	0
6	NAG	L	2	6	14,14,15	0.71	0	17,19,21	0.83	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
4	NAG	D	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	D	2	4	-	0/6/23/26	0/1/1/1
4	BMA	D	3	4	-	1/2/19/22	0/1/1/1
4	MAN	D	4	4	-	1/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FUC	D	5	4	-	-	0/1/1/1
5	NAG	E	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	E	2	5	-	0/6/23/26	0/1/1/1
5	BMA	E	3	5	-	1/2/19/22	0/1/1/1
5	MAN	E	4	5	-	2/2/19/22	0/1/1/1
5	MAN	E	5	5	-	1/2/19/22	0/1/1/1
5	MAN	E	6	5	-	0/2/19/22	0/1/1/1
5	MAN	E	7	5	-	0/2/19/22	0/1/1/1
6	NAG	F	1	6,2	-	0/6/23/26	0/1/1/1
6	NAG	F	2	6	-	0/6/23/26	0/1/1/1
4	NAG	G	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
4	BMA	G	3	4	-	1/2/19/22	0/1/1/1
4	MAN	G	4	4	-	1/2/19/22	0/1/1/1
4	FUC	G	5	4	-	-	0/1/1/1
5	NAG	H	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	H	2	5	-	0/6/23/26	0/1/1/1
5	BMA	H	3	5	-	1/2/19/22	0/1/1/1
5	MAN	H	4	5	-	2/2/19/22	0/1/1/1
5	MAN	H	5	5	-	1/2/19/22	0/1/1/1
5	MAN	H	6	5	-	0/2/19/22	0/1/1/1
5	MAN	H	7	5	-	0/2/19/22	0/1/1/1
4	NAG	I	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	I	2	4	-	0/6/23/26	0/1/1/1
4	BMA	I	3	4	-	1/2/19/22	0/1/1/1
4	MAN	I	4	4	-	1/2/19/22	0/1/1/1
4	FUC	I	5	4	-	-	0/1/1/1
5	NAG	J	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	J	2	5	-	0/6/23/26	0/1/1/1
5	BMA	J	3	5	-	1/2/19/22	0/1/1/1
5	MAN	J	4	5	-	2/2/19/22	0/1/1/1
5	MAN	J	5	5	-	1/2/19/22	0/1/1/1
5	MAN	J	6	5	-	0/2/19/22	0/1/1/1
5	MAN	J	7	5	-	0/2/19/22	0/1/1/1
6	NAG	K	1	6,2	-	0/6/23/26	0/1/1/1
6	NAG	K	2	6	-	0/6/23/26	0/1/1/1
6	NAG	L	1	6,2	-	0/6/23/26	0/1/1/1
6	NAG	L	2	6	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	3	BMA	C1-O5-C5	6.79	121.39	112.19
5	J	3	BMA	C1-O5-C5	6.79	121.39	112.19
5	H	3	BMA	C1-O5-C5	6.79	121.39	112.19
4	D	3	BMA	C1-O5-C5	6.28	120.70	112.19
4	I	3	BMA	C1-O5-C5	6.28	120.69	112.19
4	G	3	BMA	C1-O5-C5	6.27	120.69	112.19
5	H	7	MAN	C1-O5-C5	4.07	117.70	112.19
5	J	7	MAN	C1-O5-C5	4.06	117.70	112.19
5	E	7	MAN	C1-O5-C5	4.06	117.69	112.19
4	G	4	MAN	C1-O5-C5	3.88	117.45	112.19
4	I	4	MAN	C1-O5-C5	3.88	117.45	112.19
4	D	4	MAN	C1-O5-C5	3.88	117.45	112.19
5	H	4	MAN	C1-O5-C5	3.37	116.75	112.19
5	J	4	MAN	C1-O5-C5	3.37	116.75	112.19
5	E	4	MAN	C1-O5-C5	3.37	116.75	112.19
5	E	3	BMA	C3-C4-C5	3.10	115.76	110.24
5	H	3	BMA	C3-C4-C5	3.10	115.76	110.24
5	J	3	BMA	C3-C4-C5	3.09	115.76	110.24
5	J	5	MAN	C1-O5-C5	2.82	116.02	112.19
5	E	5	MAN	C1-O5-C5	2.82	116.02	112.19
5	H	5	MAN	C1-O5-C5	2.82	116.01	112.19
5	E	6	MAN	C1-O5-C5	2.66	115.79	112.19
5	J	6	MAN	C1-O5-C5	2.65	115.79	112.19
5	H	6	MAN	C1-O5-C5	2.65	115.79	112.19
5	H	3	BMA	O4-C4-C3	-2.34	104.94	110.35
5	E	3	BMA	O4-C4-C3	-2.34	104.94	110.35
5	J	3	BMA	O4-C4-C3	-2.34	104.94	110.35
4	I	3	BMA	C3-C4-C5	2.30	114.35	110.24
4	D	3	BMA	C3-C4-C5	2.30	114.34	110.24
4	G	3	BMA	C3-C4-C5	2.30	114.34	110.24
4	G	3	BMA	C2-C3-C4	2.25	114.80	110.89
4	I	3	BMA	O3-C3-C2	-2.25	105.68	109.99
4	D	3	BMA	C2-C3-C4	2.25	114.80	110.89
4	D	3	BMA	O3-C3-C2	-2.25	105.68	109.99
4	G	3	BMA	O3-C3-C2	-2.25	105.68	109.99
4	I	3	BMA	C2-C3-C4	2.25	114.79	110.89
5	E	3	BMA	O3-C3-C4	2.08	115.16	110.35
5	J	3	BMA	C2-C3-C4	2.08	114.49	110.89
5	J	3	BMA	O3-C3-C4	2.08	115.15	110.35
5	H	3	BMA	O3-C3-C4	2.08	115.15	110.35
5	H	3	BMA	C2-C3-C4	2.08	114.49	110.89
5	E	3	BMA	C2-C3-C4	2.08	114.49	110.89

There are no chirality outliers.

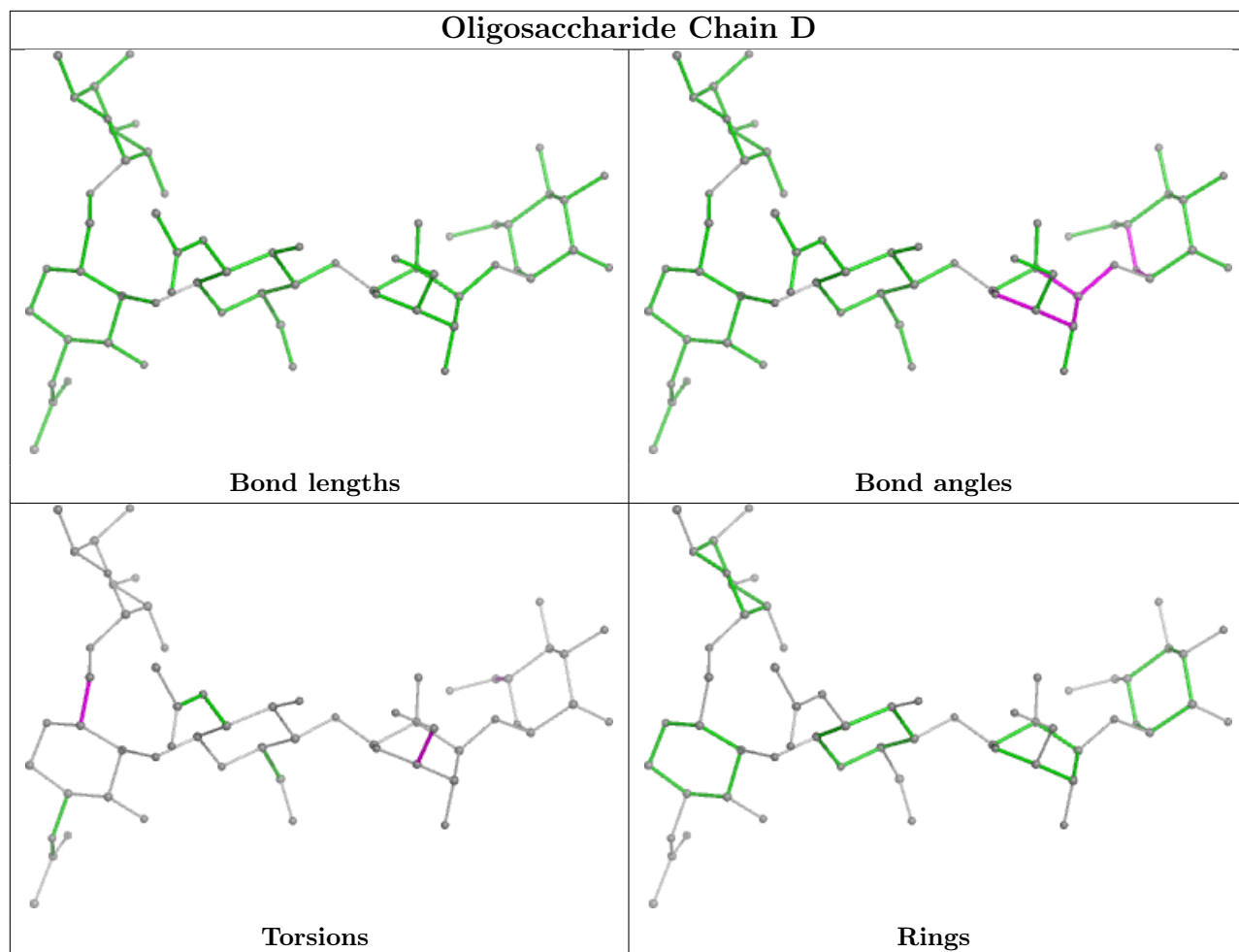
All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	4	MAN	O5-C5-C6-O6
5	H	4	MAN	O5-C5-C6-O6
5	J	4	MAN	O5-C5-C6-O6
4	D	1	NAG	C4-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
4	I	1	NAG	C4-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6
4	D	1	NAG	O5-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
5	E	3	BMA	O5-C5-C6-O6
5	H	3	BMA	O5-C5-C6-O6
5	J	3	BMA	O5-C5-C6-O6
4	D	4	MAN	C4-C5-C6-O6
4	G	4	MAN	C4-C5-C6-O6
4	I	4	MAN	C4-C5-C6-O6
5	E	4	MAN	C4-C5-C6-O6
5	H	4	MAN	C4-C5-C6-O6
5	J	4	MAN	C4-C5-C6-O6
5	E	5	MAN	C4-C5-C6-O6
5	H	5	MAN	C4-C5-C6-O6
5	J	5	MAN	C4-C5-C6-O6
4	D	3	BMA	O5-C5-C6-O6
4	G	3	BMA	O5-C5-C6-O6
4	I	3	BMA	O5-C5-C6-O6

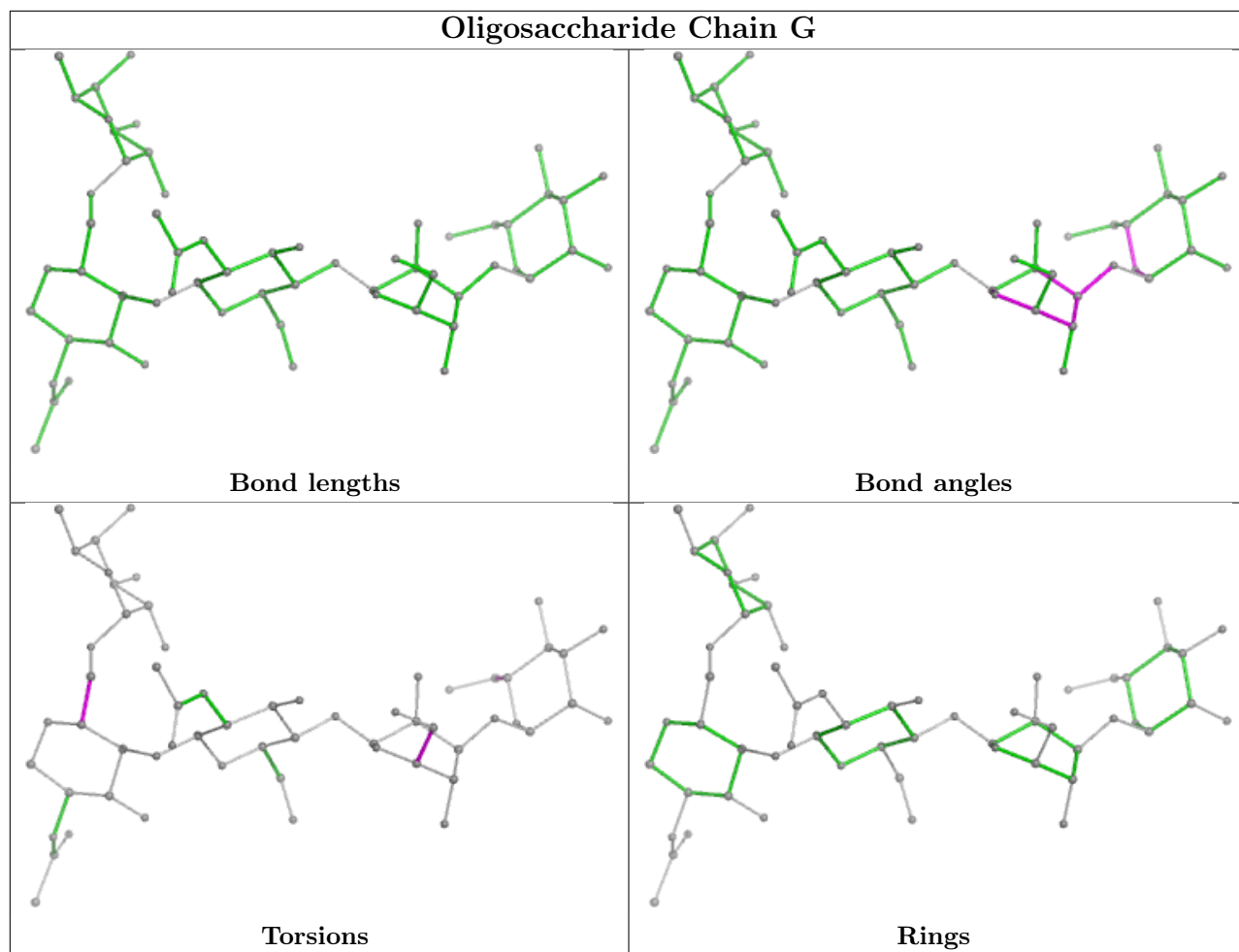
There are no ring outliers.

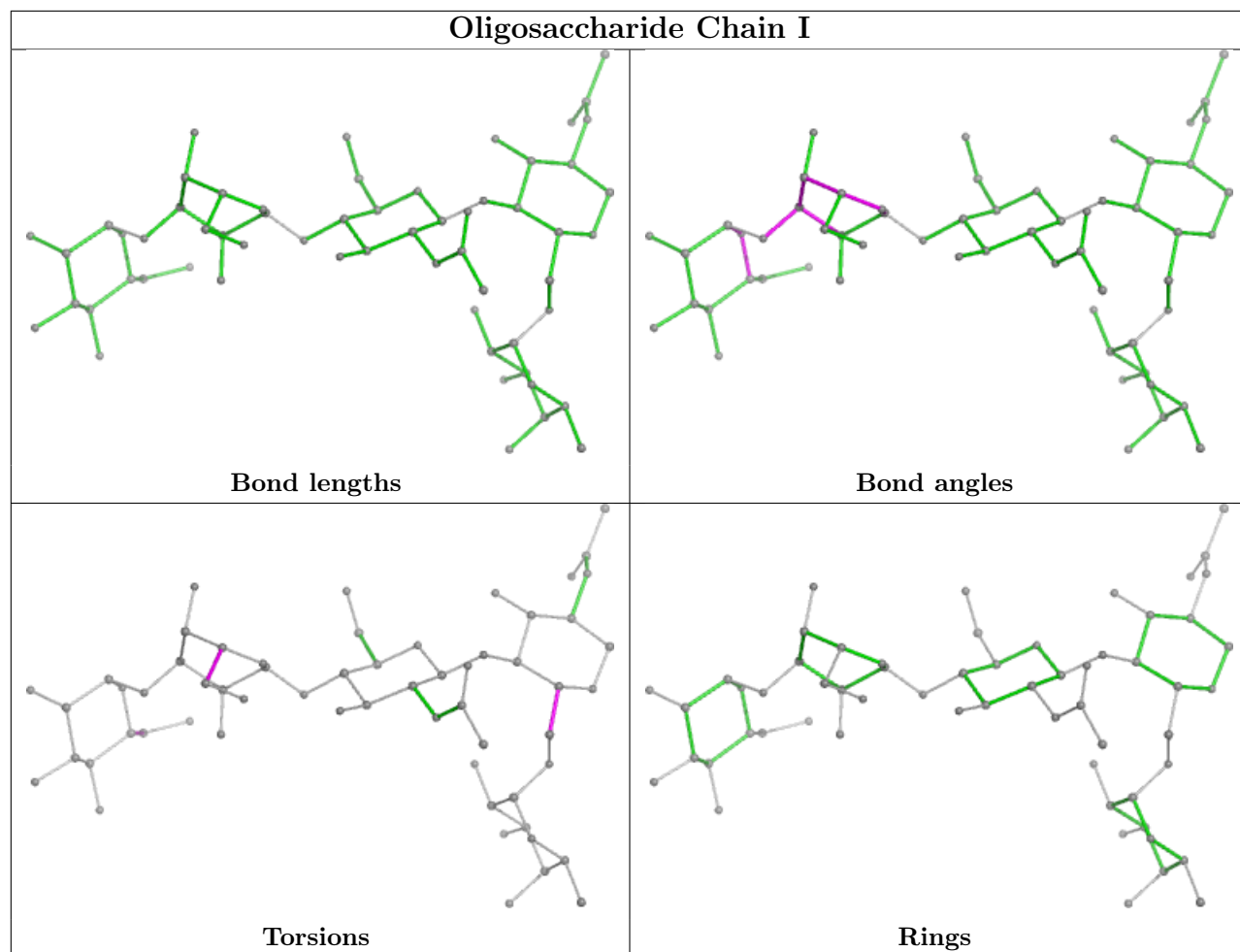
No monomer is involved in short contacts.

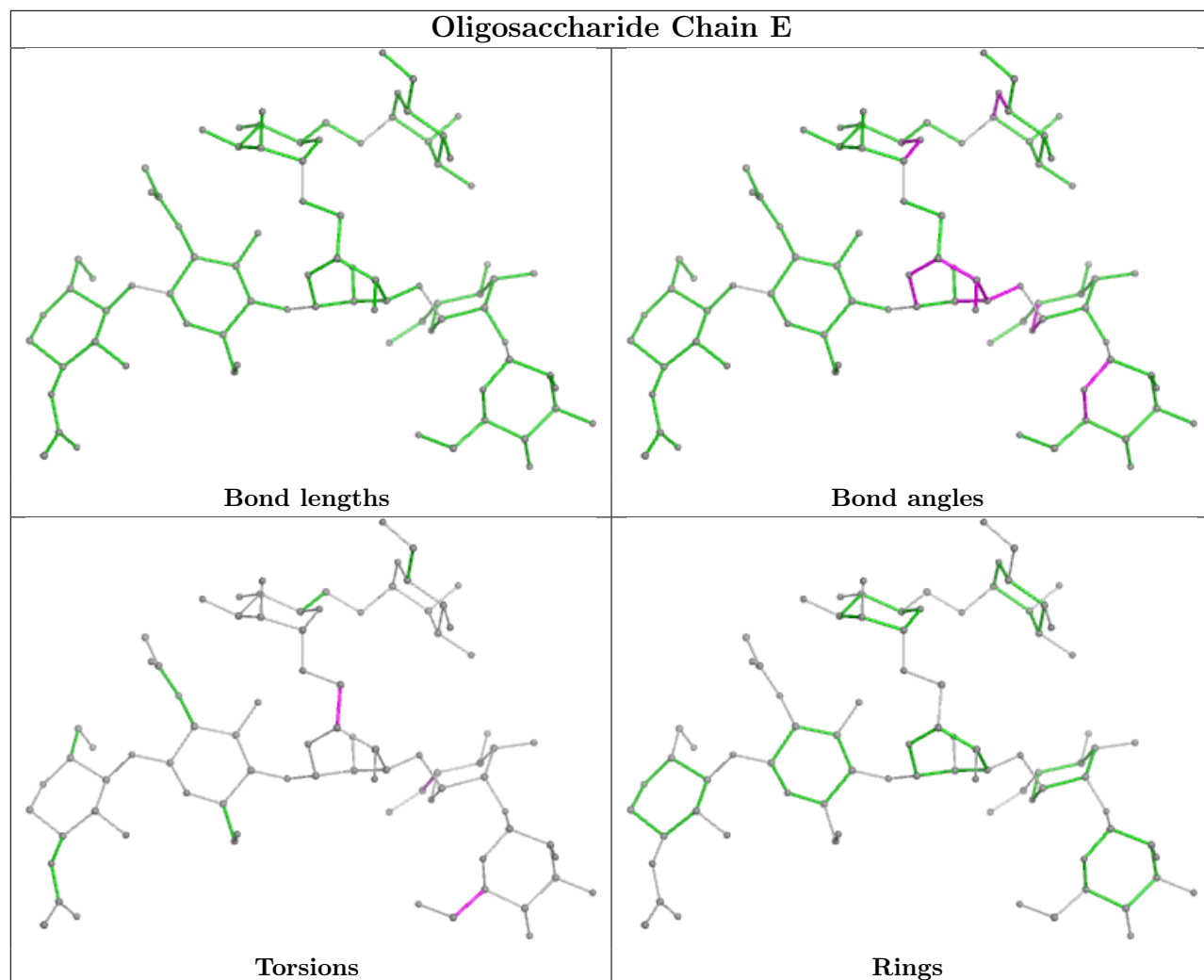
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

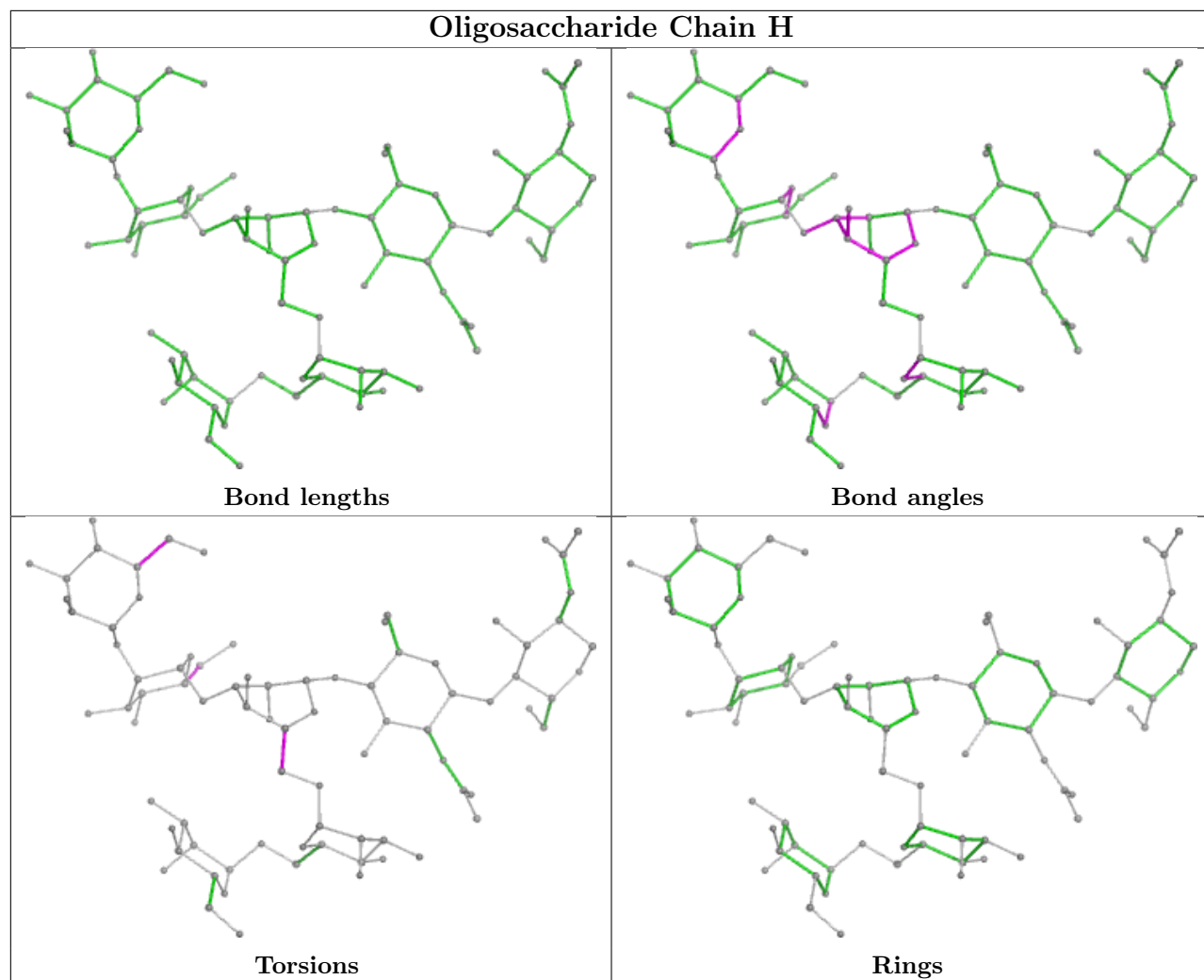


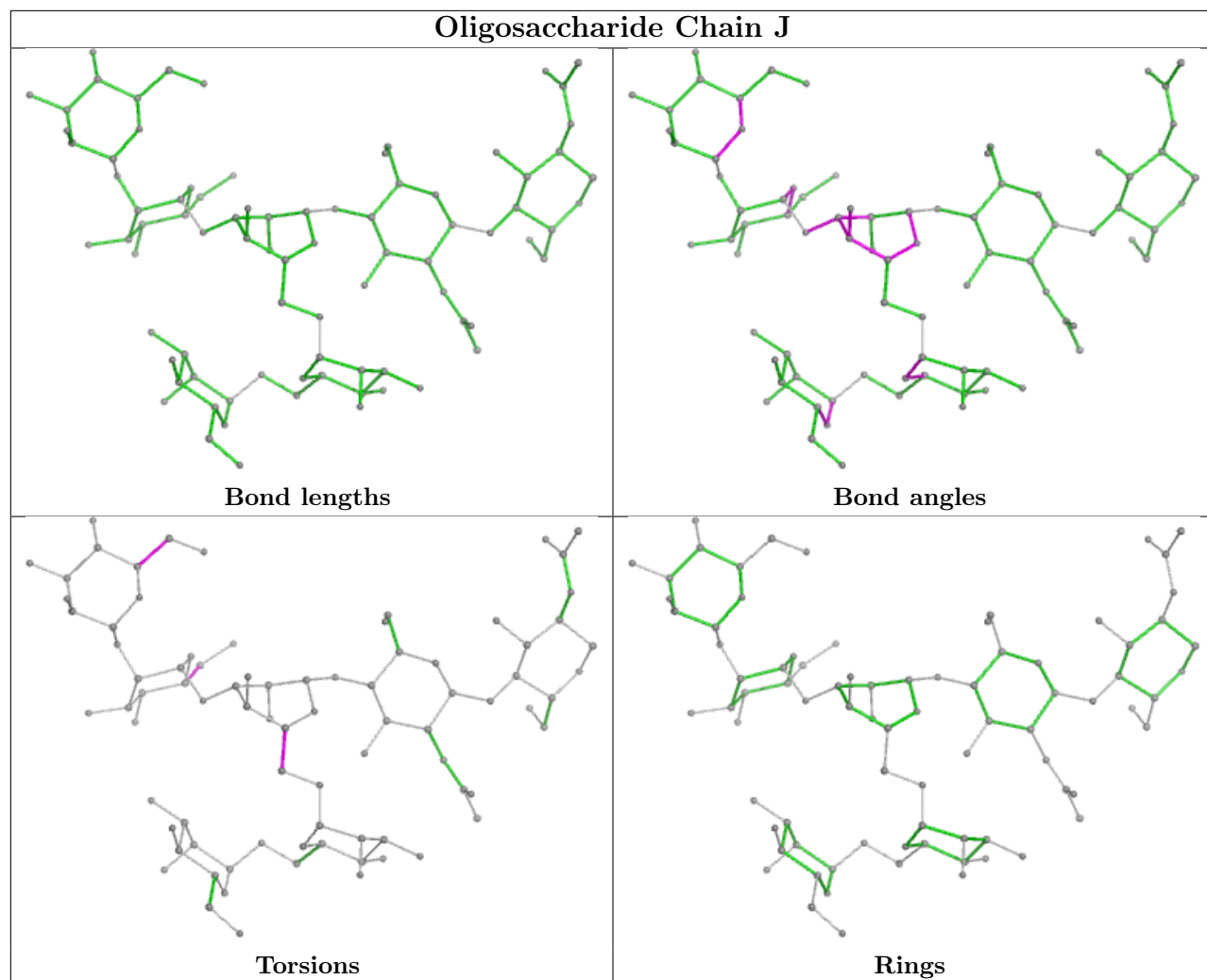


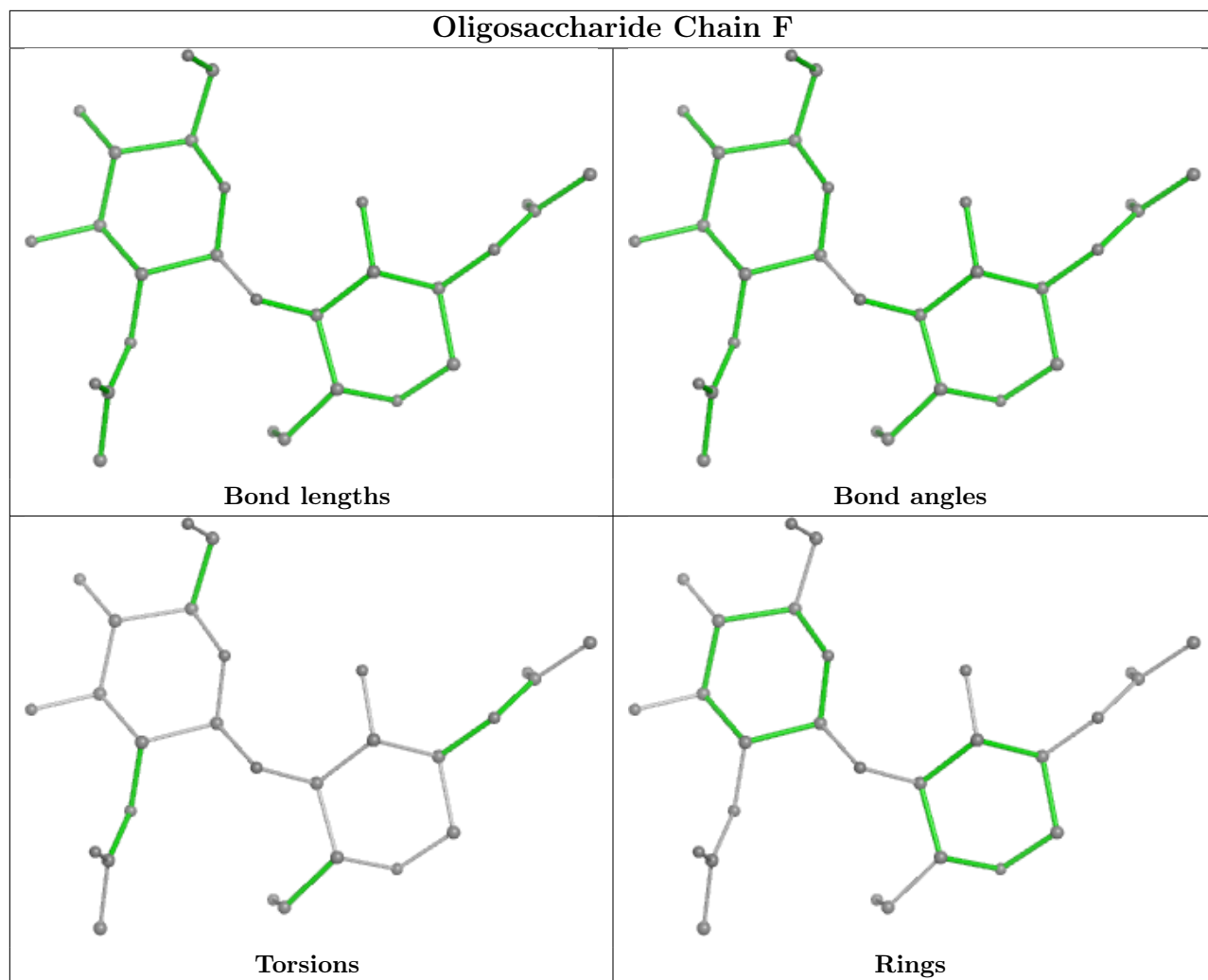


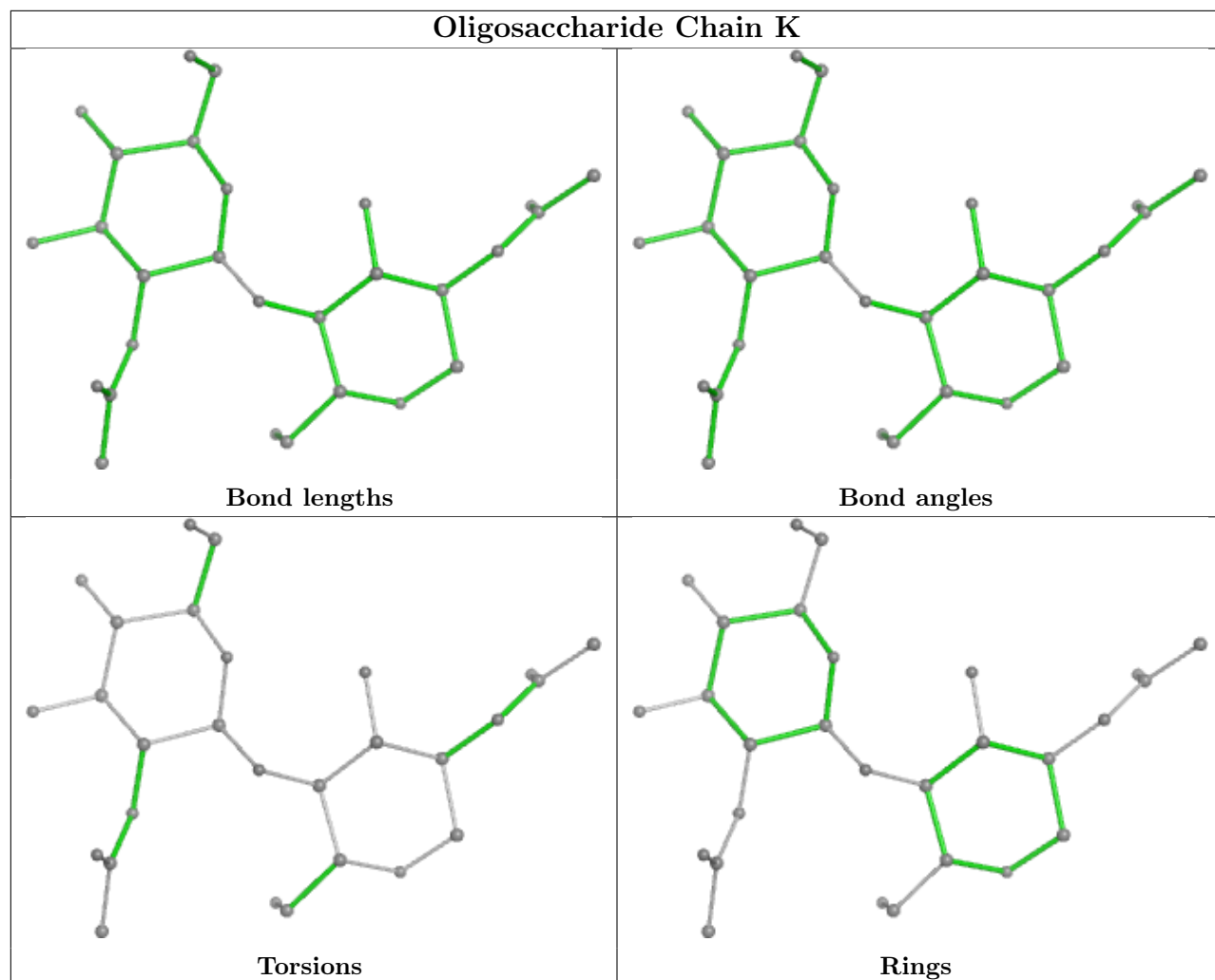


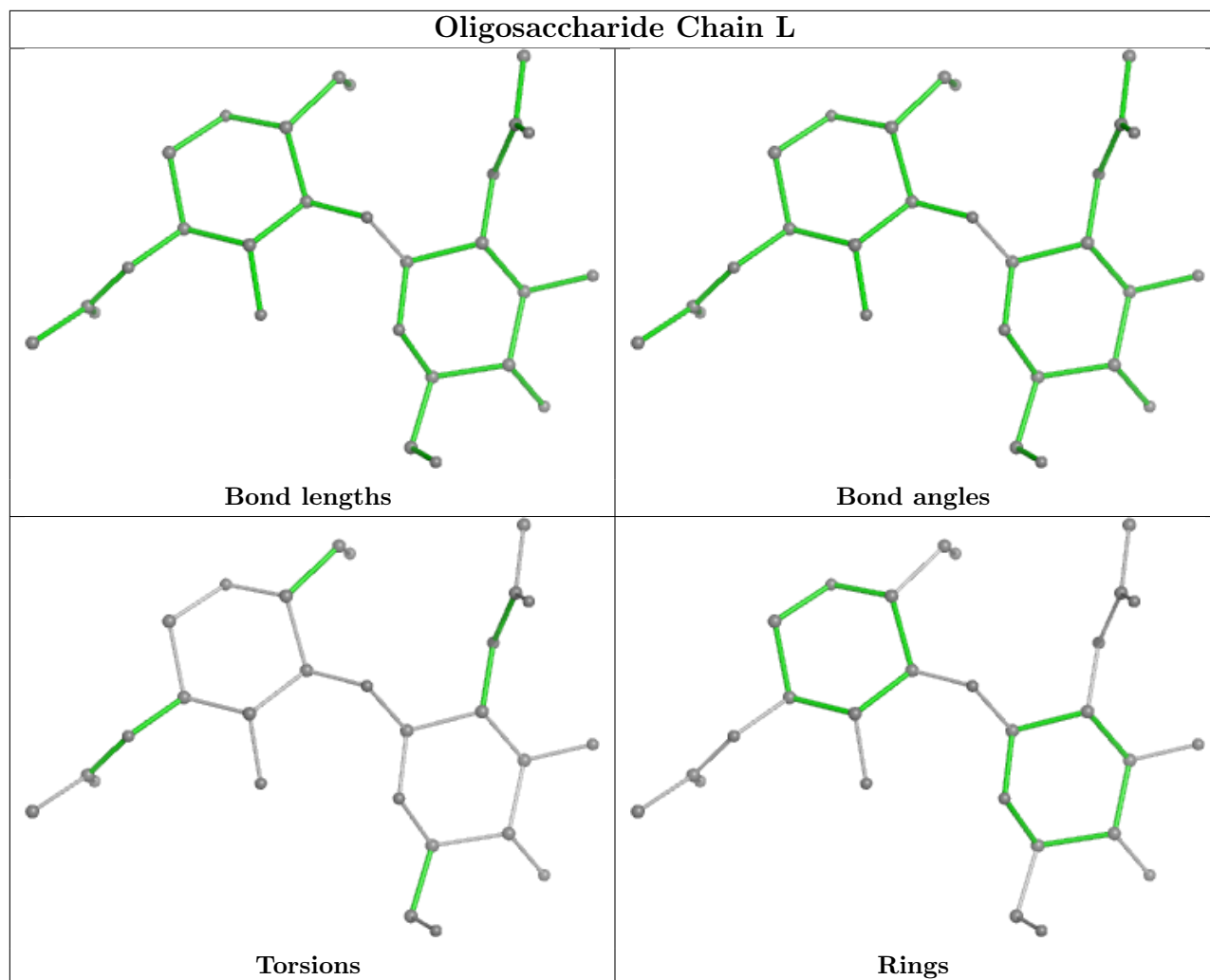












## 5.6 Ligand geometry [i](#)

Of 22 ligands modelled in this entry, 1 is monoatomic - leaving 21 for Mogul analysis.

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$
7	NAG	B	301	1	14,14,15	0.74	0	17,19,21	0.99	1 (5%)
7	NAG	C	301	1	14,14,15	0.74	0	17,19,21	0.99	1 (5%)
7	NAG	A	301	1	14,14,15	0.74	0	17,19,21	0.99	1 (5%)
7	NAG	a	502	2	14,14,15	0.70	0	17,19,21	1.25	3 (17%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
7	NAG	b	501	2	14,14,15	0.72	0	17,19,21	0.82	0
7	NAG	b	503	2	14,14,15	0.73	0	17,19,21	0.79	0
7	NAG	a	503	2	14,14,15	0.73	0	17,19,21	0.79	0
7	NAG	C	303	1	14,14,15	0.70	0	17,19,21	1.22	1 (5%)
7	NAG	c	503	2	14,14,15	0.73	0	17,19,21	0.79	0
7	NAG	a	501	2	14,14,15	0.72	0	17,19,21	0.82	0
7	NAG	A	303	1	14,14,15	0.70	0	17,19,21	1.21	1 (5%)
7	NAG	C	302	1	14,14,15	0.72	0	17,19,21	0.82	0
7	NAG	B	303	1	14,14,15	0.70	0	17,19,21	1.22	1 (5%)
7	NAG	c	502	2	14,14,15	0.70	0	17,19,21	1.25	3 (17%)
7	NAG	B	304	1	14,14,15	0.72	0	17,19,21	0.81	0
7	NAG	A	302	1	14,14,15	0.72	0	17,19,21	0.82	0
7	NAG	A	304	1	14,14,15	0.72	0	17,19,21	0.81	0
7	NAG	C	304	1	14,14,15	0.72	0	17,19,21	0.81	0
7	NAG	B	302	1	14,14,15	0.72	0	17,19,21	0.81	0
7	NAG	c	501	2	14,14,15	0.72	0	17,19,21	0.82	0
7	NAG	b	502	2	14,14,15	0.70	0	17,19,21	1.25	3 (17%)

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
7	NAG	B	301	1	-	0/6/23/26	0/1/1/1
7	NAG	C	301	1	-	0/6/23/26	0/1/1/1
7	NAG	A	301	1	-	0/6/23/26	0/1/1/1
7	NAG	a	502	2	-	1/6/23/26	0/1/1/1
7	NAG	b	501	2	-	0/6/23/26	0/1/1/1
7	NAG	b	503	2	-	0/6/23/26	0/1/1/1
7	NAG	a	503	2	-	0/6/23/26	0/1/1/1
7	NAG	C	303	1	-	2/6/23/26	0/1/1/1
7	NAG	c	503	2	-	0/6/23/26	0/1/1/1
7	NAG	a	501	2	-	0/6/23/26	0/1/1/1
7	NAG	A	303	1	-	2/6/23/26	0/1/1/1
7	NAG	C	302	1	-	0/6/23/26	0/1/1/1
7	NAG	B	303	1	-	2/6/23/26	0/1/1/1
7	NAG	c	502	2	-	1/6/23/26	0/1/1/1
7	NAG	B	304	1	-	0/6/23/26	0/1/1/1
7	NAG	A	302	1	-	0/6/23/26	0/1/1/1
7	NAG	A	304	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	C	304	1	-	0/6/23/26	0/1/1/1
7	NAG	B	302	1	-	0/6/23/26	0/1/1/1
7	NAG	c	501	2	-	0/6/23/26	0/1/1/1
7	NAG	b	502	2	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	303	NAG	C2-N2-C7	3.27	127.56	122.90
7	A	303	NAG	C2-N2-C7	3.27	127.56	122.90
7	C	303	NAG	C2-N2-C7	3.27	127.56	122.90
7	a	502	NAG	C1-O5-C5	2.70	115.86	112.19
7	b	502	NAG	C1-O5-C5	2.70	115.85	112.19
7	c	502	NAG	C1-O5-C5	2.70	115.85	112.19
7	c	502	NAG	C4-C3-C2	-2.66	107.12	111.02
7	b	502	NAG	C4-C3-C2	-2.66	107.12	111.02
7	a	502	NAG	C4-C3-C2	-2.66	107.12	111.02
7	a	502	NAG	O5-C1-C2	-2.57	107.23	111.29
7	b	502	NAG	O5-C1-C2	-2.57	107.23	111.29
7	c	502	NAG	O5-C1-C2	-2.57	107.23	111.29
7	A	301	NAG	C1-O5-C5	2.52	115.61	112.19
7	B	301	NAG	C1-O5-C5	2.52	115.60	112.19
7	C	301	NAG	C1-O5-C5	2.52	115.60	112.19

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	a	502	NAG	O5-C5-C6-O6
7	b	502	NAG	O5-C5-C6-O6
7	c	502	NAG	O5-C5-C6-O6
7	A	303	NAG	C3-C2-N2-C7
7	B	303	NAG	C3-C2-N2-C7
7	C	303	NAG	C3-C2-N2-C7
7	B	303	NAG	C1-C2-N2-C7
7	C	303	NAG	C1-C2-N2-C7
7	A	303	NAG	C1-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

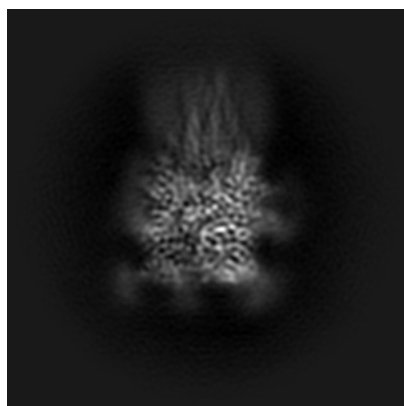
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17428. These allow visual inspection of the internal detail of the map and identification of artifacts.

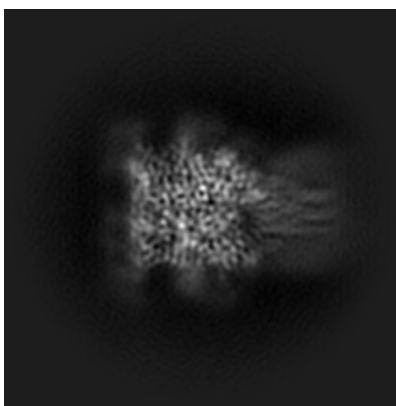
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

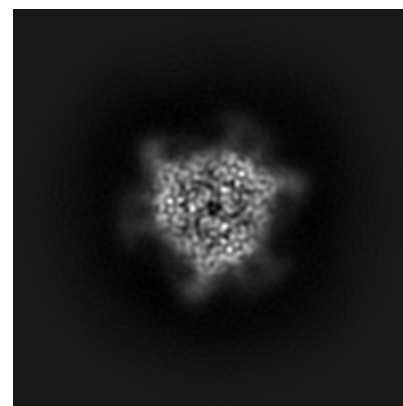
#### 6.1.1 Primary map



X

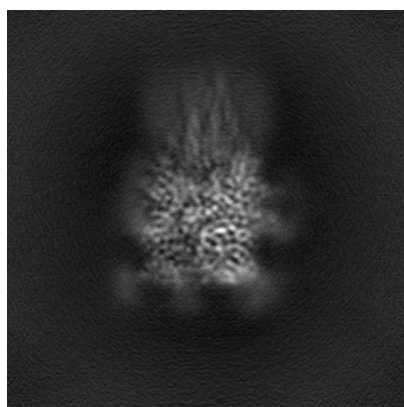


Y

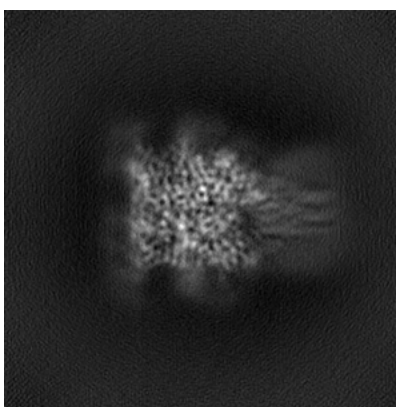


Z

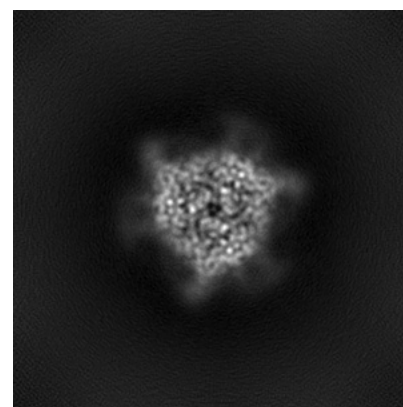
#### 6.1.2 Raw map



X



Y

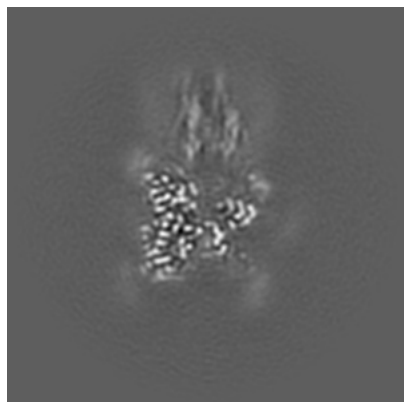


Z

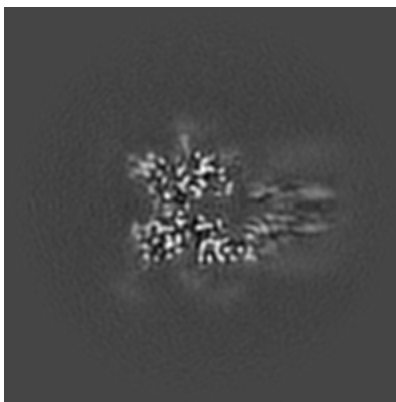
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

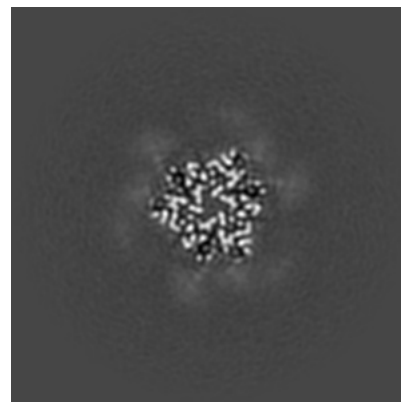
### 6.2.1 Primary map



X Index: 128

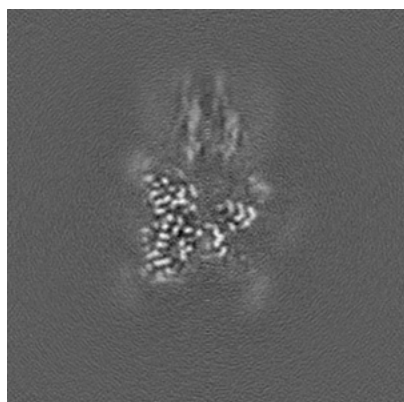


Y Index: 128

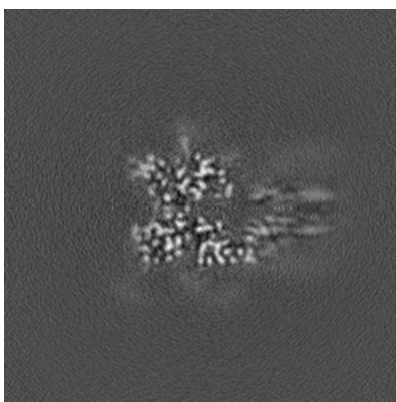


Z Index: 128

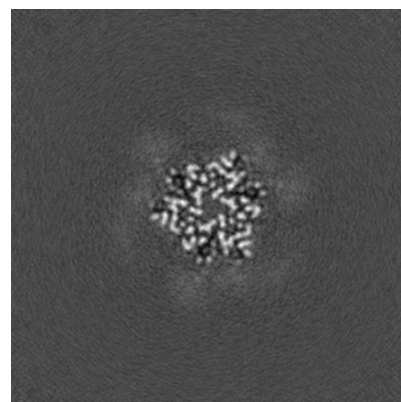
### 6.2.2 Raw map



X Index: 128



Y Index: 128

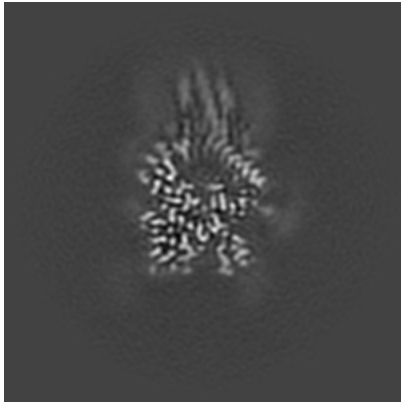


Z Index: 128

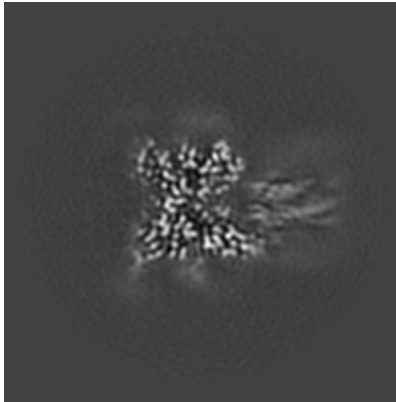
The images above show central slices of the map in three orthogonal directions.

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

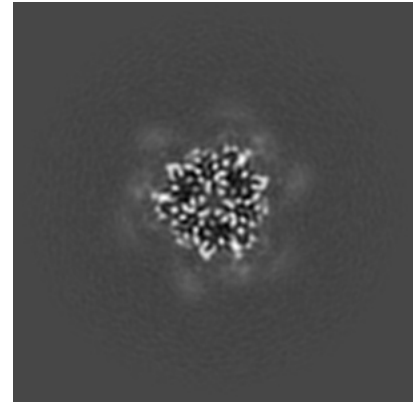
### 6.3.1 Primary map



X Index: 134

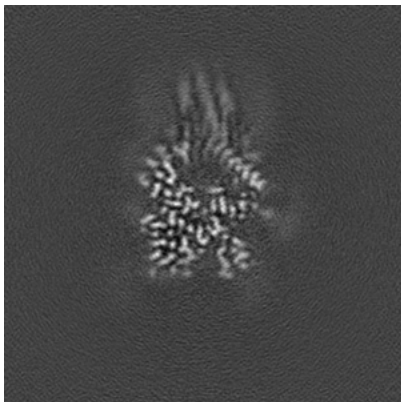


Y Index: 134

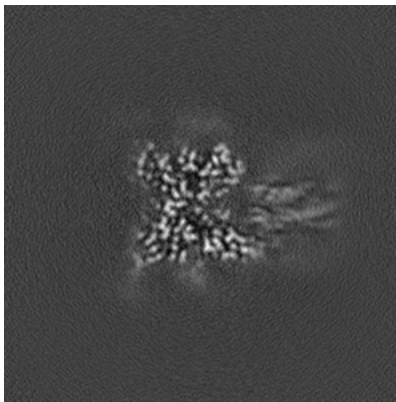


Z Index: 125

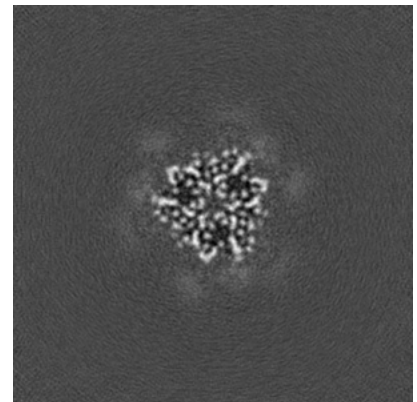
### 6.3.2 Raw map



X Index: 134



Y Index: 134

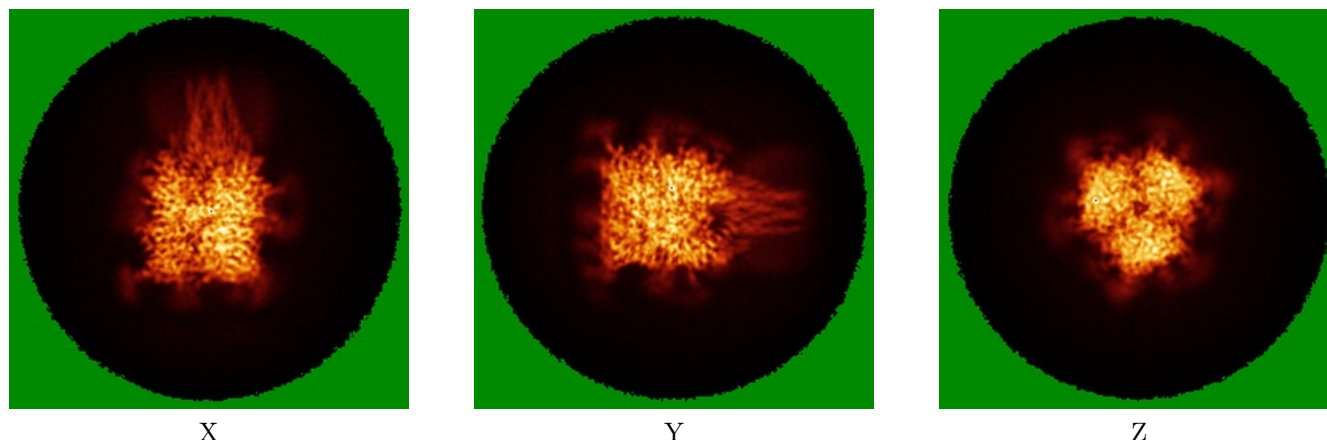


Z Index: 126

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

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

### 6.4.1 Primary map

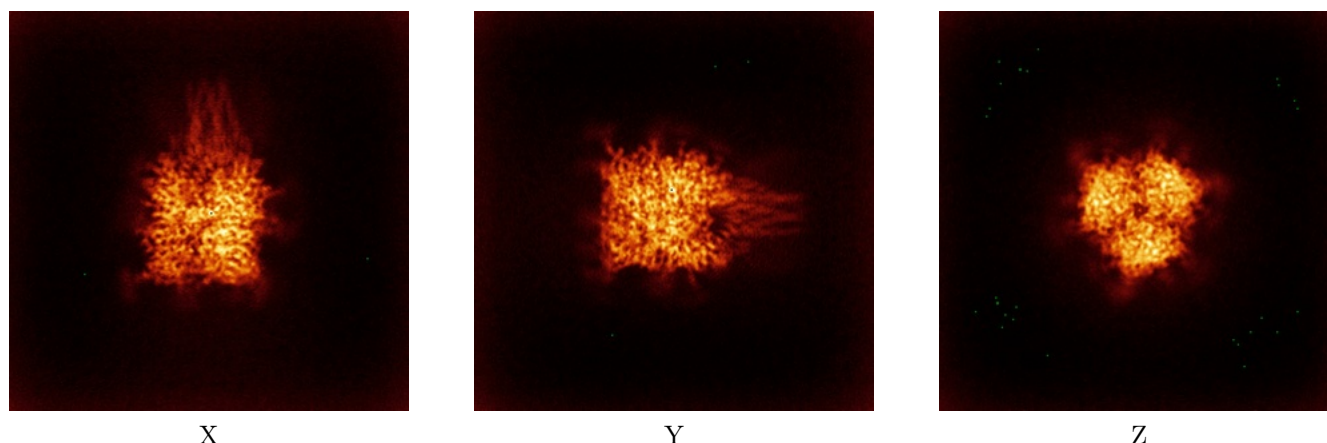


X

Y

Z

### 6.4.2 Raw map



X

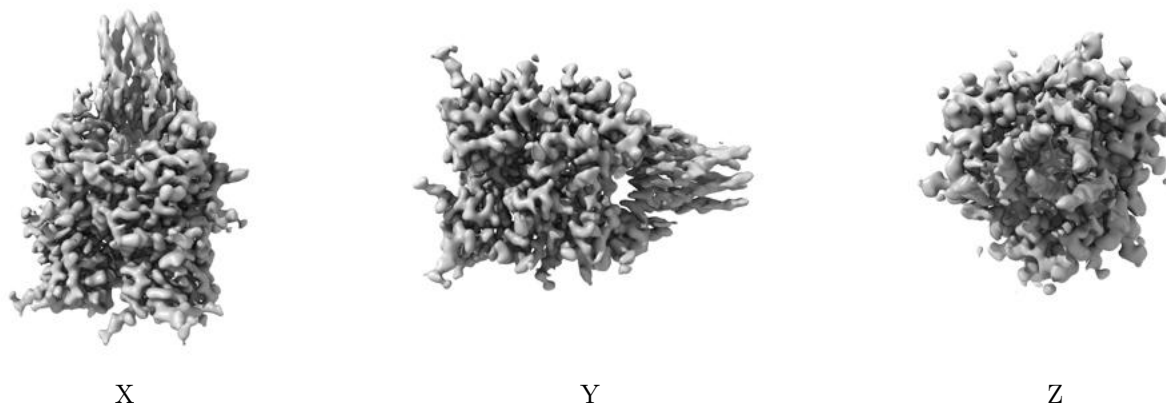
Y

Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

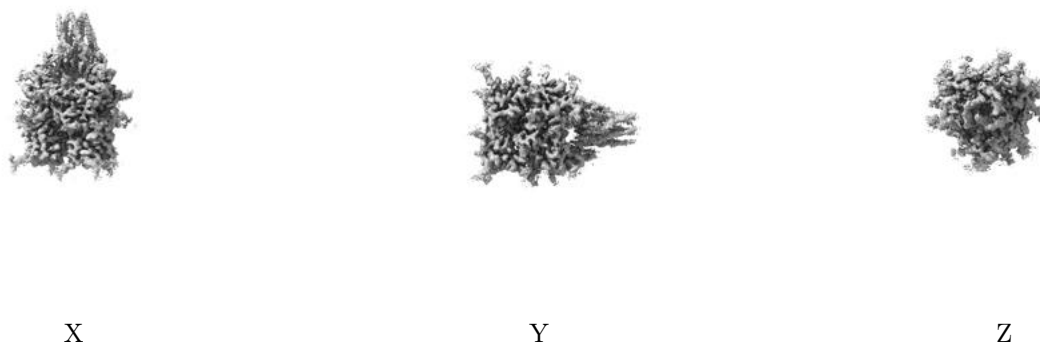
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.08. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.5.2 Raw map



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



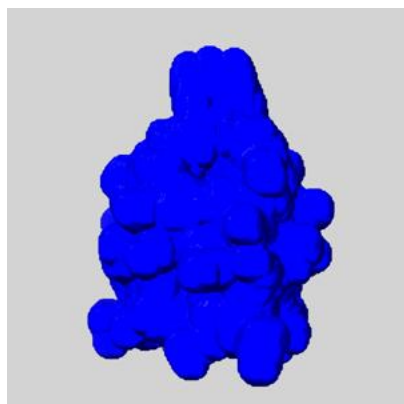
## 6.6 Mask visualisation [i](#)

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

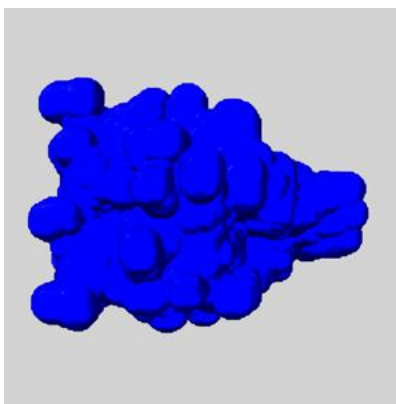
A mask typically either:

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

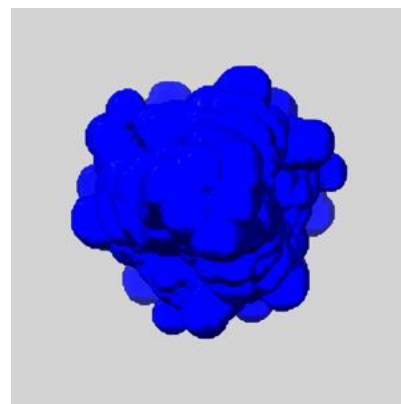
### 6.6.1 emd\_17428\_msk\_1.map [i](#)



X



Y

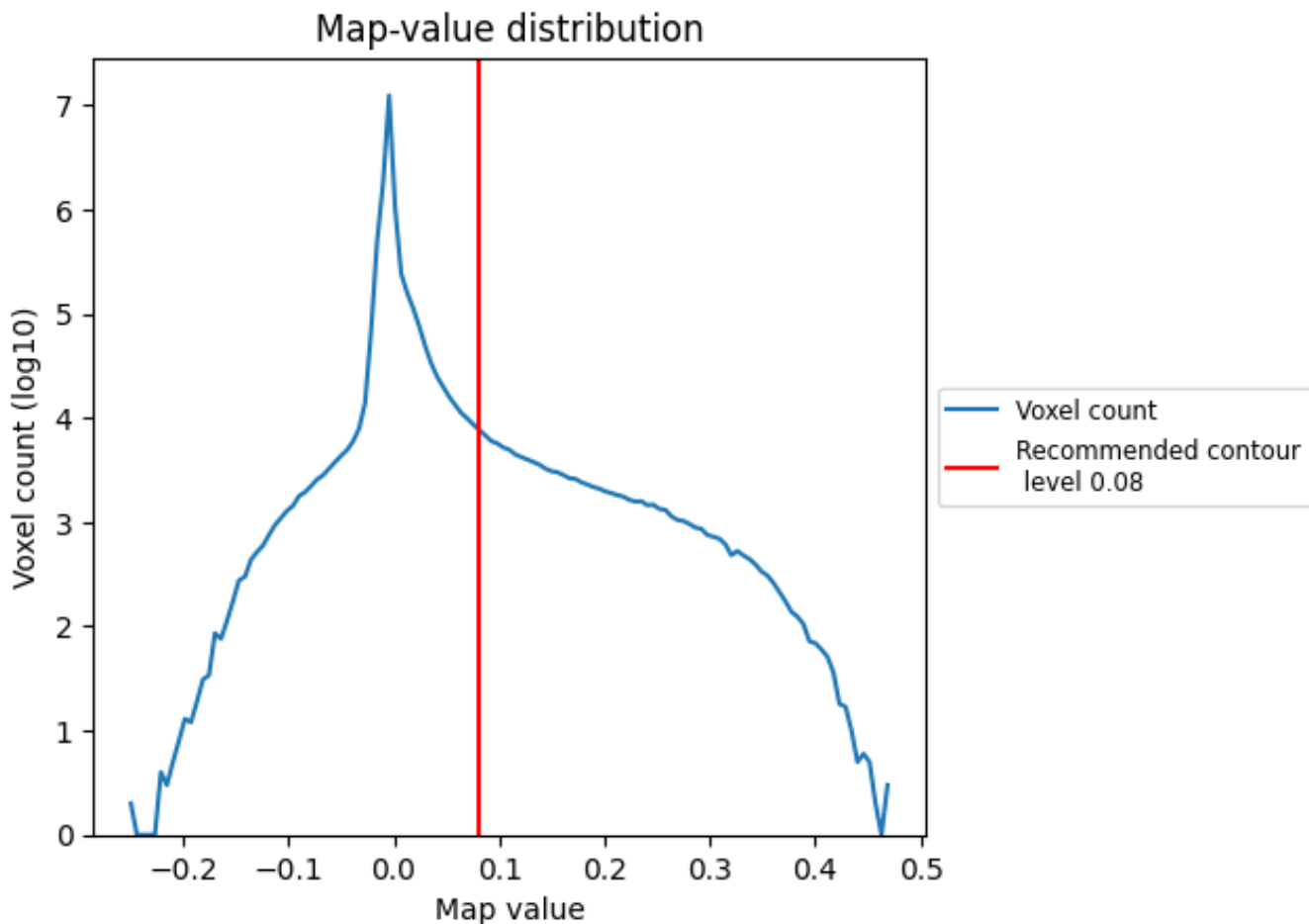


Z

## 7 Map analysis [i](#)

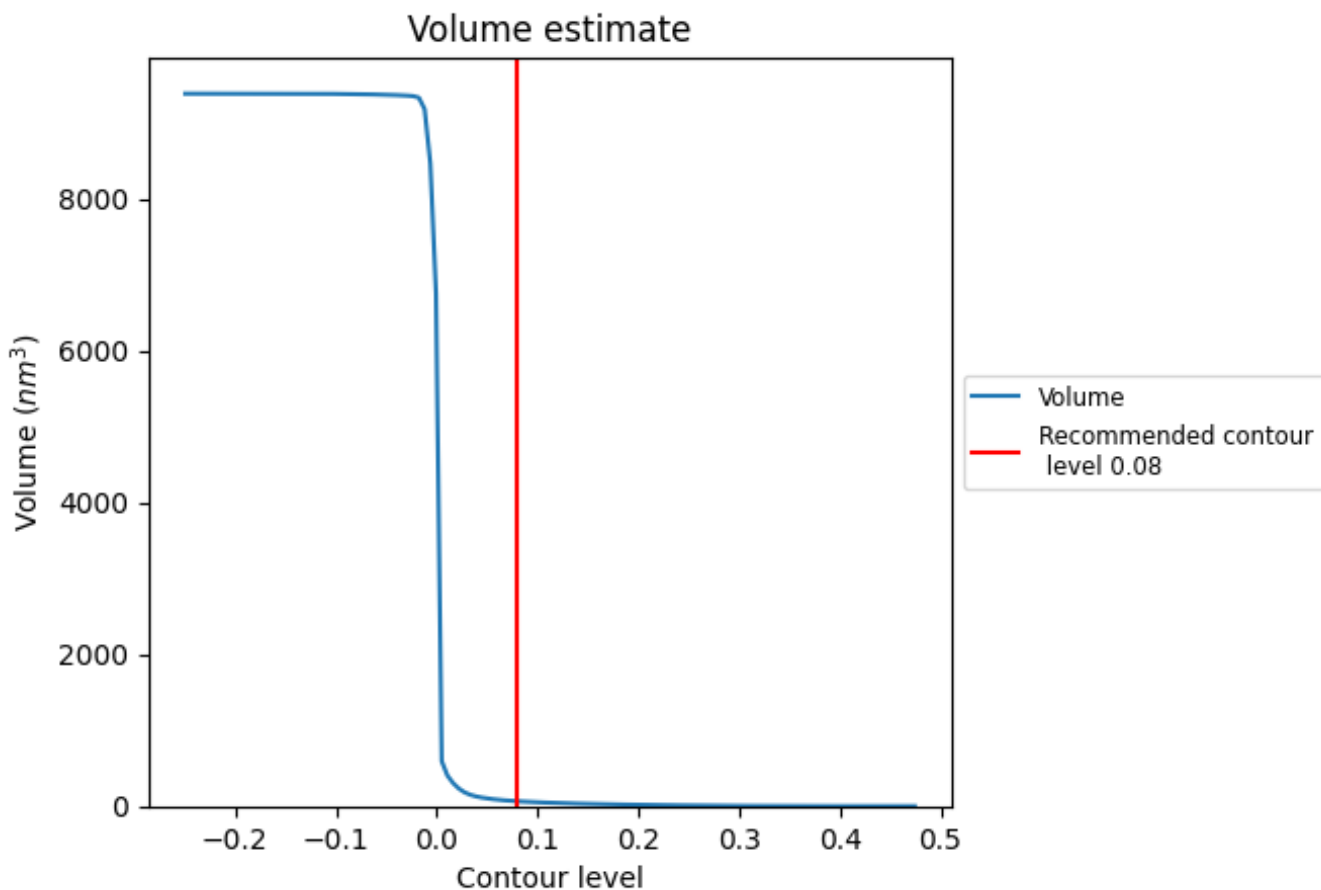
This section contains the results of statistical analysis of the map.

### 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.

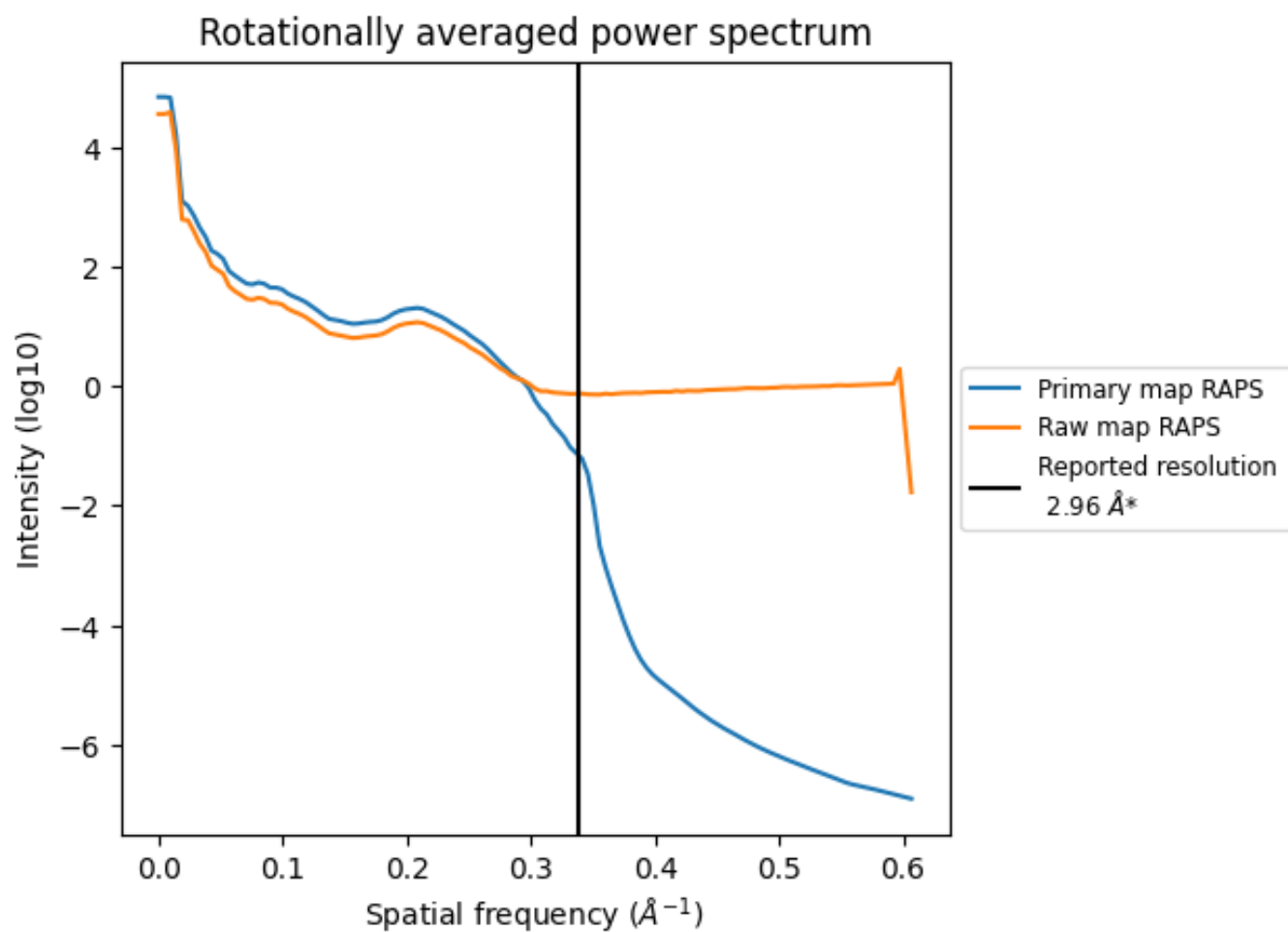
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is  $65 \text{ nm}^3$ ; this corresponds to an approximate mass of 58 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

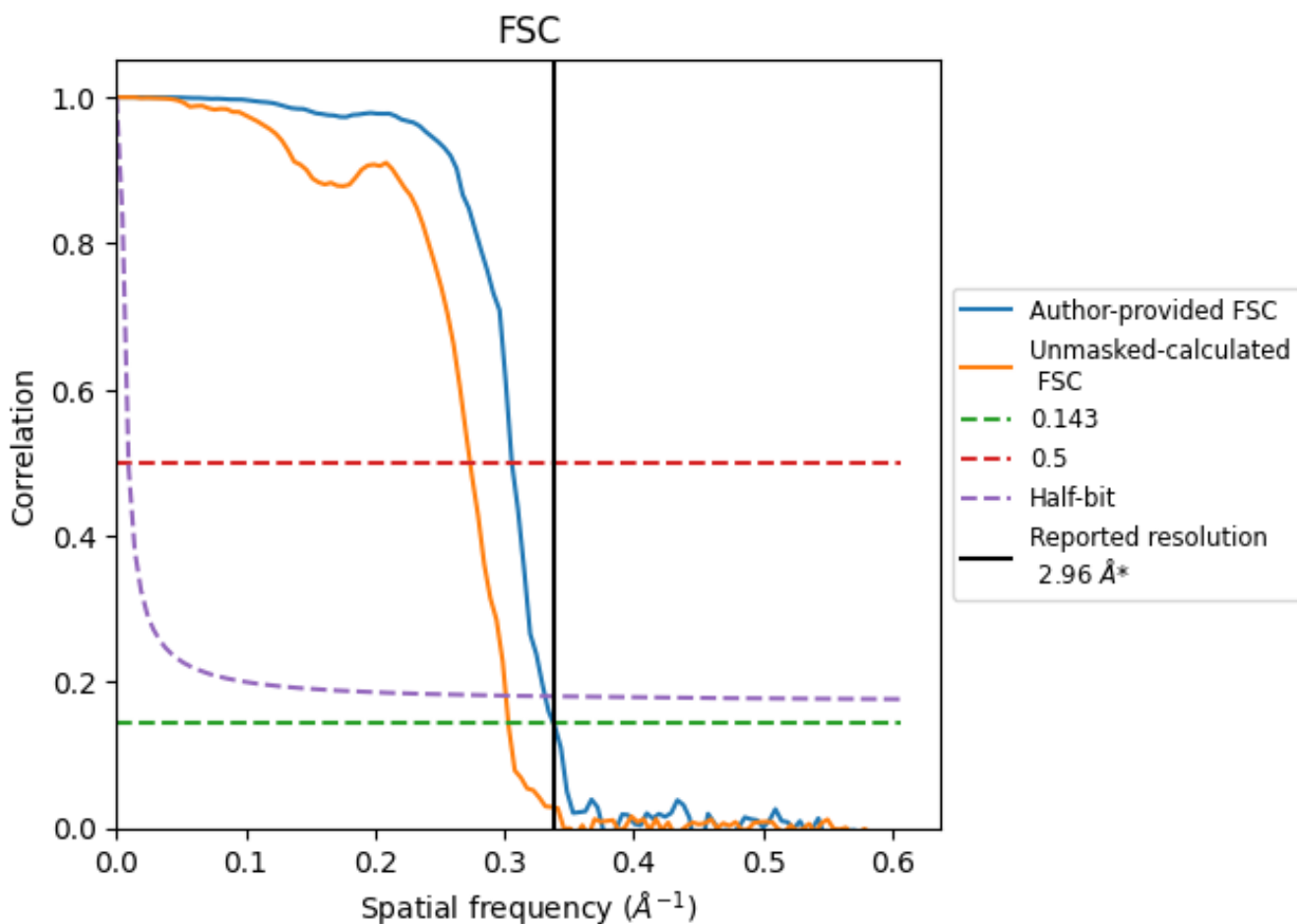


\*Reported resolution corresponds to spatial frequency of  $0.338 \text{ \AA}^{-1}$

## 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 0.338  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

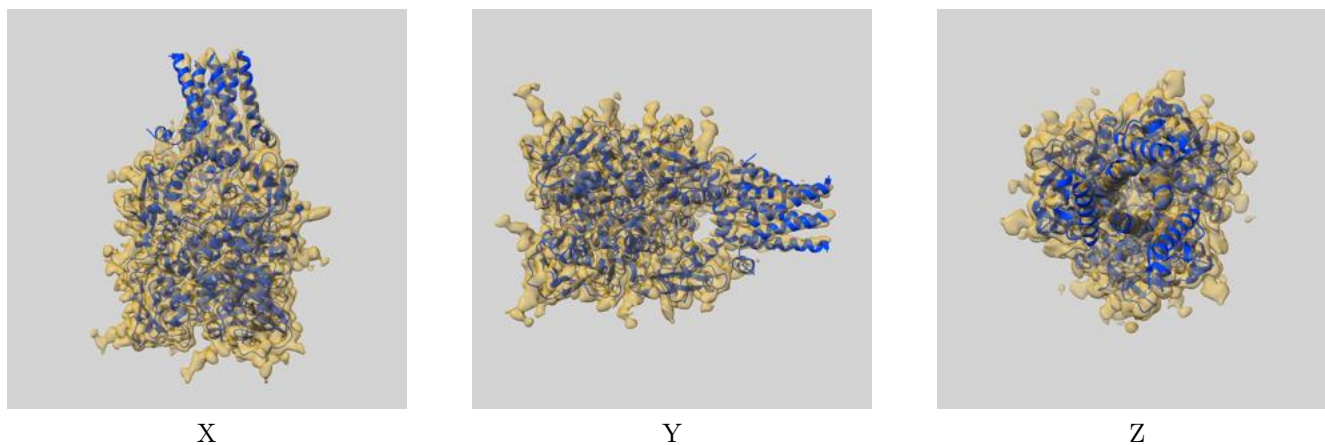
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.96	-	-
Author-provided FSC curve	2.96	3.27	3.02
Unmasked-calculated*	3.30	3.66	3.32

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.30 differs from the reported value 2.96 by more than 10 %

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

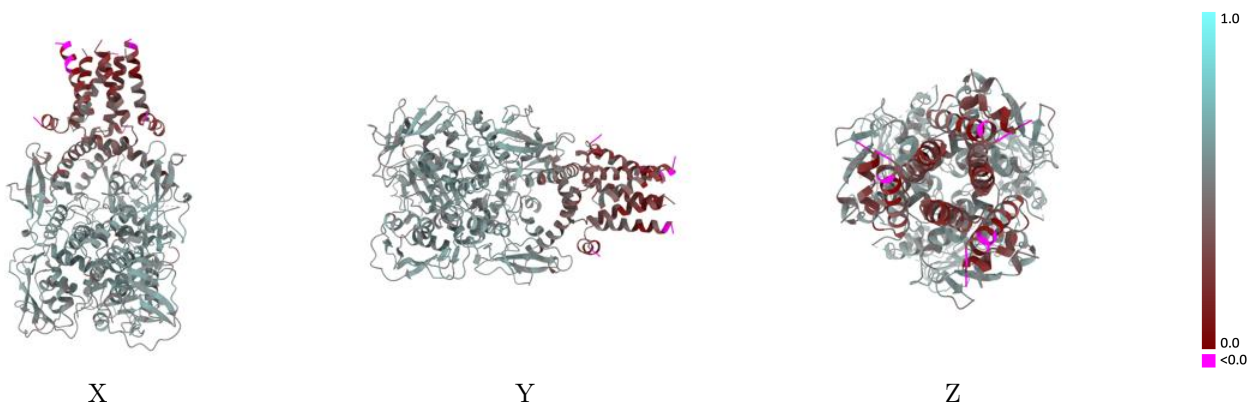
This section contains information regarding the fit between EMDB map EMD-17428 and PDB model 8P4T. Per-residue inclusion information can be found in section 3 on page 9.

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



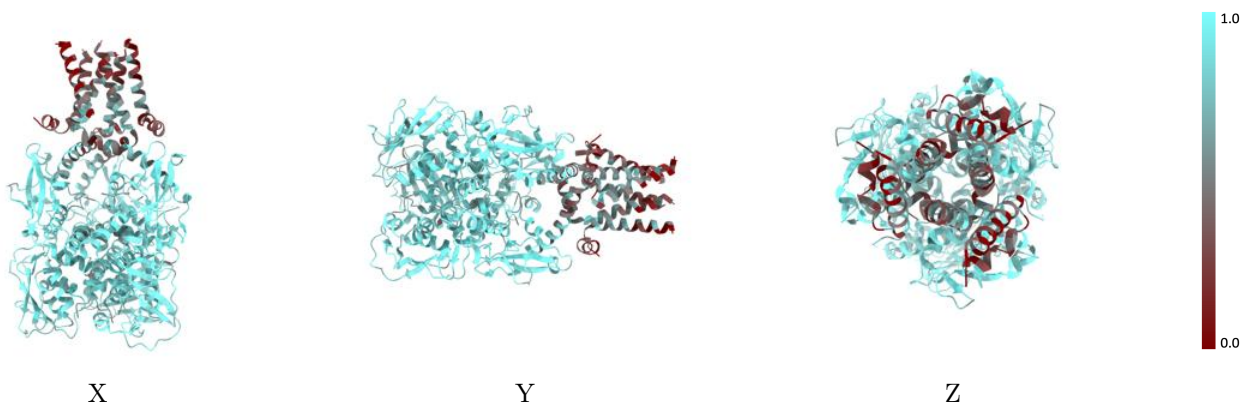
The images above show the 3D surface view of the map at the recommended contour level 0.08 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

## 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.

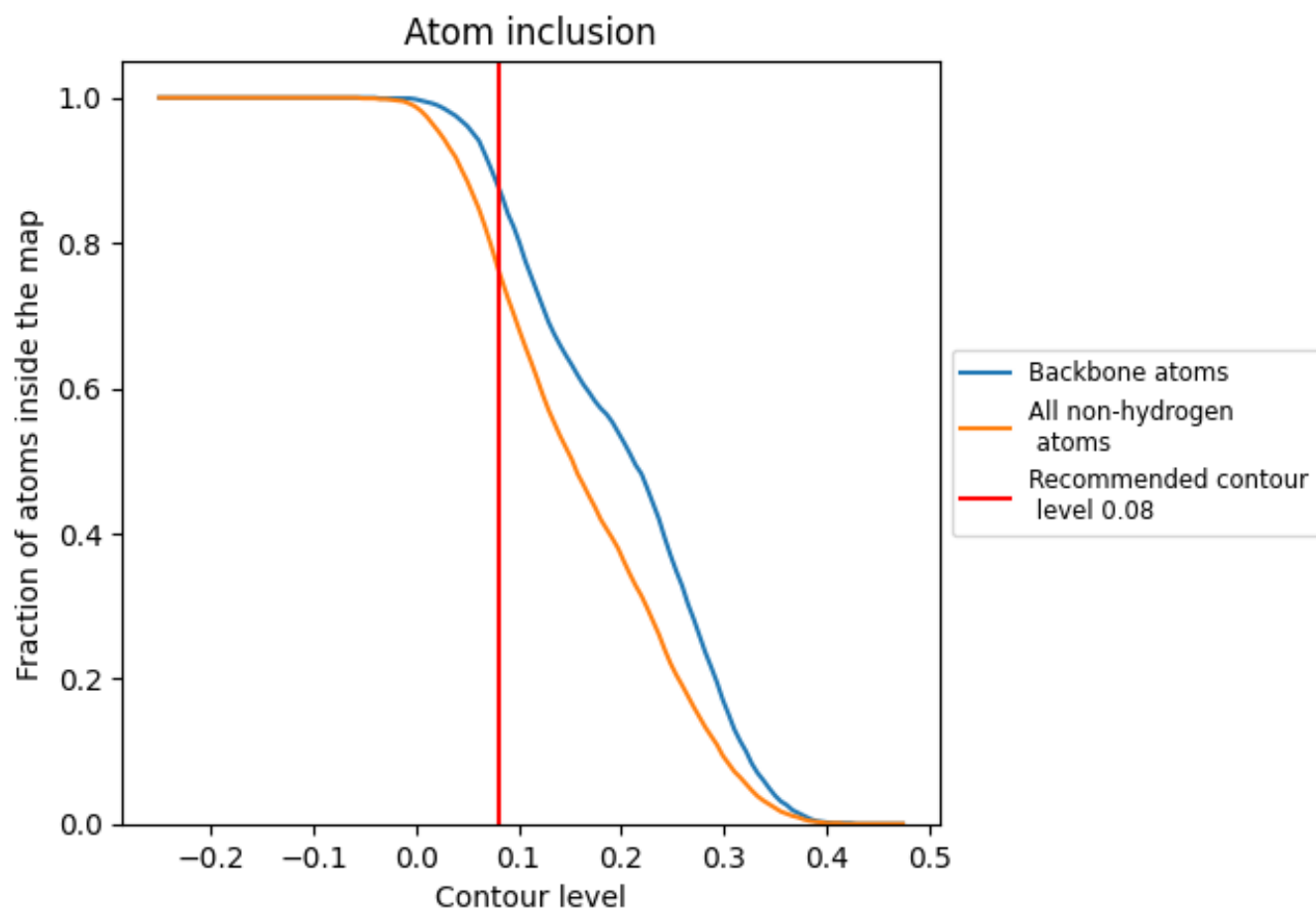
## 9.3 Atom inclusion mapped to coordinate model [\(i\)](#)



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 (0.08).









































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 76% 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 (0.08) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7650	 0.4810
A	 0.8590	 0.5340
B	 0.8590	 0.5340
C	 0.8520	 0.5320
D	 0.5170	 0.4120
E	 0.3130	 0.3290
F	 0.7140	 0.4340
G	 0.5500	 0.4170
H	 0.3490	 0.3410
I	 0.5170	 0.4140
J	 0.3250	 0.3310
K	 0.7500	 0.4230
L	 0.7140	 0.4280
SA	 0.3400	 0.2420
SB	 0.2940	 0.2150
SC	 0.3400	 0.2530
a	 0.7980	 0.4890
b	 0.7920	 0.4840
c	 0.7960	 0.4880

