



Full wwPDB EM Validation Report ⓘ

Apr 28, 2026 – 10:47 pm BST

PDB ID : 9TBJ / pdb_00009tbj
EMDB ID : EMD-55773
Title : Leishmania mexicana secreted acid phosphatase
Authors : Bose, P.; Grossman-Haham, I.
Deposited on : 2025-11-20
Resolution : 3.00 Å(reported)
Based on initial model : .

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

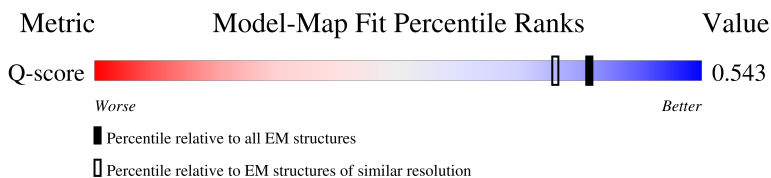
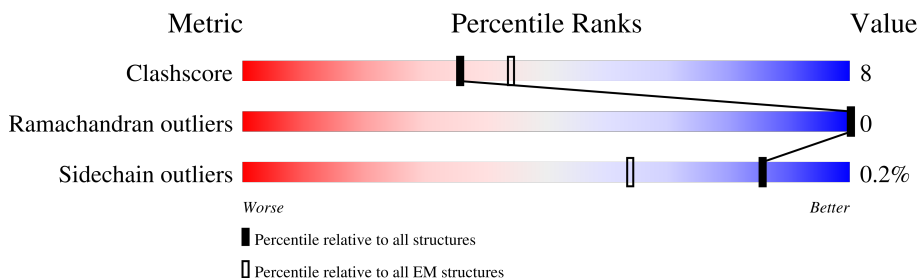
EMDB validation analysis : 0.0.1.dev132
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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



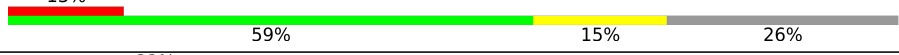

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14081 (2.50 - 3.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	565	
1	B	565	
1	C	565	
2	D	3	

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Mol	Chain	Length	Quality of chain
2	I	3	 33% 67%
3	F	2	 100%
3	G	2	 50% 50% 50%
3	K	2	 50% 50%
3	L	2	 50% 50%
3	M	2	 50% 100%
3	P	2	 50% 100%
3	Q	2	 50% 50%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	PO4	C	604	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10259 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 Secreted acid phosphatase 1 (SAP1).

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	423	Total	C	N	O	S	0	0
			3292	2067	573	631	21		
1	B	424	Total	C	N	O	S	0	0
			3300	2071	574	634	21		
1	C	420	Total	C	N	O	S	0	0
			3275	2057	570	627	21		

There are 252 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	HIS	ARG	conflict	UNP Q25332
A	126	ASN	TYR	conflict	UNP Q25332
A	152	ALA	GLY	conflict	UNP Q25332
A	180	CYS	SER	conflict	UNP Q25332
A	199	SER	ASN	conflict	UNP Q25332
A	374	ASN	SER	conflict	UNP Q25332
A	379	ILE	VAL	conflict	UNP Q25332
A	392	THR	ALA	conflict	UNP Q25332
A	405	PRO	ARG	conflict	UNP Q25332
A	412	SER	PRO	conflict	UNP Q25332
A	422	TYR	HIS	conflict	UNP Q25332
A	446	SER	PRO	conflict	UNP Q25332
A	451	LEU	PRO	conflict	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	GLU	deletion	UNP Q25332
A	?	-	GLY	deletion	UNP Q25332
A	?	-	THR	deletion	UNP Q25332
A	?	-	THR	deletion	UNP Q25332
A	?	-	THR	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	SER	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	GLU	deletion	UNP Q25332
A	?	-	GLY	deletion	UNP Q25332
A	?	-	THR	deletion	UNP Q25332
A	?	-	THR	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	?	-	SER	deletion	UNP Q25332
A	493	GLY	ARG	conflict	UNP Q25332
A	517	SER	-	expression tag	UNP Q25332
A	518	ARG	-	expression tag	UNP Q25332
A	519	MET	-	expression tag	UNP Q25332
A	520	ASP	-	expression tag	UNP Q25332
A	521	GLU	-	expression tag	UNP Q25332
A	522	LYS	-	expression tag	UNP Q25332
A	523	THR	-	expression tag	UNP Q25332
A	524	THR	-	expression tag	UNP Q25332
A	525	GLY	-	expression tag	UNP Q25332
A	526	TRP	-	expression tag	UNP Q25332
A	527	ARG	-	expression tag	UNP Q25332
A	528	GLY	-	expression tag	UNP Q25332
A	529	GLY	-	expression tag	UNP Q25332
A	530	HIS	-	expression tag	UNP Q25332
A	531	VAL	-	expression tag	UNP Q25332
A	532	VAL	-	expression tag	UNP Q25332
A	533	GLU	-	expression tag	UNP Q25332
A	534	GLY	-	expression tag	UNP Q25332
A	535	LEU	-	expression tag	UNP Q25332
A	536	ALA	-	expression tag	UNP Q25332
A	537	GLY	-	expression tag	UNP Q25332
A	538	GLU	-	expression tag	UNP Q25332
A	539	LEU	-	expression tag	UNP Q25332
A	540	GLU	-	expression tag	UNP Q25332
A	541	GLN	-	expression tag	UNP Q25332
A	542	LEU	-	expression tag	UNP Q25332
A	543	ARG	-	expression tag	UNP Q25332
A	544	ALA	-	expression tag	UNP Q25332
A	545	ARG	-	expression tag	UNP Q25332
A	546	LEU	-	expression tag	UNP Q25332
A	547	GLU	-	expression tag	UNP Q25332

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Chain	Residue	Modelled	Actual	Comment	Reference
A	548	HIS	-	expression tag	UNP Q25332
A	549	HIS	-	expression tag	UNP Q25332
A	550	PRO	-	expression tag	UNP Q25332
A	551	GLN	-	expression tag	UNP Q25332
A	552	GLY	-	expression tag	UNP Q25332
A	553	GLN	-	expression tag	UNP Q25332
A	554	ARG	-	expression tag	UNP Q25332
A	555	GLU	-	expression tag	UNP Q25332
A	556	PRO	-	expression tag	UNP Q25332
A	557	SER	-	expression tag	UNP Q25332
A	558	GLY	-	expression tag	UNP Q25332
A	559	GLY	-	expression tag	UNP Q25332
A	560	CYS	-	expression tag	UNP Q25332
A	561	LYS	-	expression tag	UNP Q25332
A	562	LEU	-	expression tag	UNP Q25332
A	563	GLY	-	expression tag	UNP Q25332
A	564	LEU	-	expression tag	UNP Q25332
A	565	TYR	-	expression tag	UNP Q25332
B	24	HIS	ARG	conflict	UNP Q25332
B	126	ASN	TYR	conflict	UNP Q25332
B	152	ALA	GLY	conflict	UNP Q25332
B	180	CYS	SER	conflict	UNP Q25332
B	199	SER	ASN	conflict	UNP Q25332
B	374	ASN	SER	conflict	UNP Q25332
B	379	ILE	VAL	conflict	UNP Q25332
B	392	THR	ALA	conflict	UNP Q25332
B	405	PRO	ARG	conflict	UNP Q25332
B	412	SER	PRO	conflict	UNP Q25332
B	422	TYR	HIS	conflict	UNP Q25332
B	446	SER	PRO	conflict	UNP Q25332
B	451	LEU	PRO	conflict	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	GLU	deletion	UNP Q25332
B	?	-	GLY	deletion	UNP Q25332
B	?	-	THR	deletion	UNP Q25332
B	?	-	THR	deletion	UNP Q25332
B	?	-	THR	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	SER	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	GLU	deletion	UNP Q25332
B	?	-	GLY	deletion	UNP Q25332
B	?	-	THR	deletion	UNP Q25332
B	?	-	THR	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	?	-	SER	deletion	UNP Q25332
B	493	GLY	ARG	conflict	UNP Q25332
B	517	SER	-	expression tag	UNP Q25332
B	518	ARG	-	expression tag	UNP Q25332
B	519	MET	-	expression tag	UNP Q25332
B	520	ASP	-	expression tag	UNP Q25332
B	521	GLU	-	expression tag	UNP Q25332
B	522	LYS	-	expression tag	UNP Q25332
B	523	THR	-	expression tag	UNP Q25332
B	524	THR	-	expression tag	UNP Q25332
B	525	GLY	-	expression tag	UNP Q25332
B	526	TRP	-	expression tag	UNP Q25332
B	527	ARG	-	expression tag	UNP Q25332
B	528	GLY	-	expression tag	UNP Q25332
B	529	GLY	-	expression tag	UNP Q25332
B	530	HIS	-	expression tag	UNP Q25332
B	531	VAL	-	expression tag	UNP Q25332
B	532	VAL	-	expression tag	UNP Q25332
B	533	GLU	-	expression tag	UNP Q25332
B	534	GLY	-	expression tag	UNP Q25332
B	535	LEU	-	expression tag	UNP Q25332
B	536	ALA	-	expression tag	UNP Q25332
B	537	GLY	-	expression tag	UNP Q25332
B	538	GLU	-	expression tag	UNP Q25332
B	539	LEU	-	expression tag	UNP Q25332
B	540	GLU	-	expression tag	UNP Q25332
B	541	GLN	-	expression tag	UNP Q25332
B	542	LEU	-	expression tag	UNP Q25332
B	543	ARG	-	expression tag	UNP Q25332
B	544	ALA	-	expression tag	UNP Q25332
B	545	ARG	-	expression tag	UNP Q25332
B	546	LEU	-	expression tag	UNP Q25332
B	547	GLU	-	expression tag	UNP Q25332

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Chain	Residue	Modelled	Actual	Comment	Reference
B	548	HIS	-	expression tag	UNP Q25332
B	549	HIS	-	expression tag	UNP Q25332
B	550	PRO	-	expression tag	UNP Q25332
B	551	GLN	-	expression tag	UNP Q25332
B	552	GLY	-	expression tag	UNP Q25332
B	553	GLN	-	expression tag	UNP Q25332
B	554	ARG	-	expression tag	UNP Q25332
B	555	GLU	-	expression tag	UNP Q25332
B	556	PRO	-	expression tag	UNP Q25332
B	557	SER	-	expression tag	UNP Q25332
B	558	GLY	-	expression tag	UNP Q25332
B	559	GLY	-	expression tag	UNP Q25332
B	560	CYS	-	expression tag	UNP Q25332
B	561	LYS	-	expression tag	UNP Q25332
B	562	LEU	-	expression tag	UNP Q25332
B	563	GLY	-	expression tag	UNP Q25332
B	564	LEU	-	expression tag	UNP Q25332
B	565	TYR	-	expression tag	UNP Q25332
C	24	HIS	ARG	conflict	UNP Q25332
C	126	ASN	TYR	conflict	UNP Q25332
C	152	ALA	GLY	conflict	UNP Q25332
C	180	CYS	SER	conflict	UNP Q25332
C	199	SER	ASN	conflict	UNP Q25332
C	374	ASN	SER	conflict	UNP Q25332
C	379	ILE	VAL	conflict	UNP Q25332
C	392	THR	ALA	conflict	UNP Q25332
C	405	PRO	ARG	conflict	UNP Q25332
C	412	SER	PRO	conflict	UNP Q25332
C	422	TYR	HIS	conflict	UNP Q25332
C	446	SER	PRO	conflict	UNP Q25332
C	451	LEU	PRO	conflict	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	GLU	deletion	UNP Q25332
C	?	-	GLY	deletion	UNP Q25332
C	?	-	THR	deletion	UNP Q25332
C	?	-	THR	deletion	UNP Q25332
C	?	-	THR	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332

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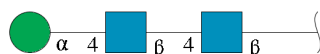
Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	SER	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	GLU	deletion	UNP Q25332
C	?	-	GLY	deletion	UNP Q25332
C	?	-	THR	deletion	UNP Q25332
C	?	-	THR	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	?	-	SER	deletion	UNP Q25332
C	493	GLY	ARG	conflict	UNP Q25332
C	517	SER	-	expression tag	UNP Q25332
C	518	ARG	-	expression tag	UNP Q25332
C	519	MET	-	expression tag	UNP Q25332
C	520	ASP	-	expression tag	UNP Q25332
C	521	GLU	-	expression tag	UNP Q25332
C	522	LYS	-	expression tag	UNP Q25332
C	523	THR	-	expression tag	UNP Q25332
C	524	THR	-	expression tag	UNP Q25332
C	525	GLY	-	expression tag	UNP Q25332
C	526	TRP	-	expression tag	UNP Q25332
C	527	ARG	-	expression tag	UNP Q25332
C	528	GLY	-	expression tag	UNP Q25332
C	529	GLY	-	expression tag	UNP Q25332
C	530	HIS	-	expression tag	UNP Q25332
C	531	VAL	-	expression tag	UNP Q25332
C	532	VAL	-	expression tag	UNP Q25332
C	533	GLU	-	expression tag	UNP Q25332
C	534	GLY	-	expression tag	UNP Q25332
C	535	LEU	-	expression tag	UNP Q25332
C	536	ALA	-	expression tag	UNP Q25332
C	537	GLY	-	expression tag	UNP Q25332
C	538	GLU	-	expression tag	UNP Q25332
C	539	LEU	-	expression tag	UNP Q25332
C	540	GLU	-	expression tag	UNP Q25332
C	541	GLN	-	expression tag	UNP Q25332
C	542	LEU	-	expression tag	UNP Q25332
C	543	ARG	-	expression tag	UNP Q25332
C	544	ALA	-	expression tag	UNP Q25332
C	545	ARG	-	expression tag	UNP Q25332
C	546	LEU	-	expression tag	UNP Q25332
C	547	GLU	-	expression tag	UNP Q25332

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Chain	Residue	Modelled	Actual	Comment	Reference
C	548	HIS	-	expression tag	UNP Q25332
C	549	HIS	-	expression tag	UNP Q25332
C	550	PRO	-	expression tag	UNP Q25332
C	551	GLN	-	expression tag	UNP Q25332
C	552	GLY	-	expression tag	UNP Q25332
C	553	GLN	-	expression tag	UNP Q25332
C	554	ARG	-	expression tag	UNP Q25332
C	555	GLU	-	expression tag	UNP Q25332
C	556	PRO	-	expression tag	UNP Q25332
C	557	SER	-	expression tag	UNP Q25332
C	558	GLY	-	expression tag	UNP Q25332
C	559	GLY	-	expression tag	UNP Q25332
C	560	CYS	-	expression tag	UNP Q25332
C	561	LYS	-	expression tag	UNP Q25332
C	562	LEU	-	expression tag	UNP Q25332
C	563	GLY	-	expression tag	UNP Q25332
C	564	LEU	-	expression tag	UNP Q25332
C	565	TYR	-	expression tag	UNP Q25332

- Molecule 2 is an oligosaccharide called alpha-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
2	D	3	Total	C	N	O	0	0
			39	22	2	15		
2	I	3	Total	C	N	O	0	0
			39	22	2	15		

- 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
3	F	2	Total	C	N	O	0	0
			28	16	2	10		

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Mol	Chain	Residues	Atoms				AltConf	Trace
3	G	2	Total	C	N	O	0	0
			28	16	2	10		
3	K	2	Total	C	N	O	0	0
			28	16	2	10		
3	L	2	Total	C	N	O	0	0
			28	16	2	10		
3	M	2	Total	C	N	O	0	0
			28	16	2	10		
3	P	2	Total	C	N	O	0	0
			28	16	2	10		
3	Q	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).



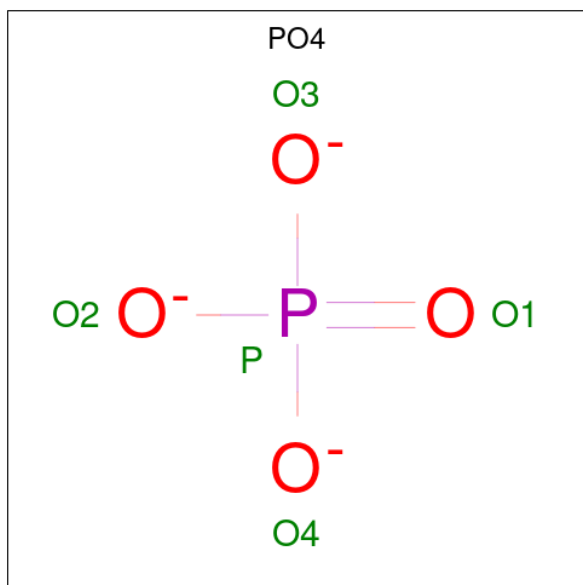
Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	

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

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P) (labeled as "Ligand of Interest" by depositor).

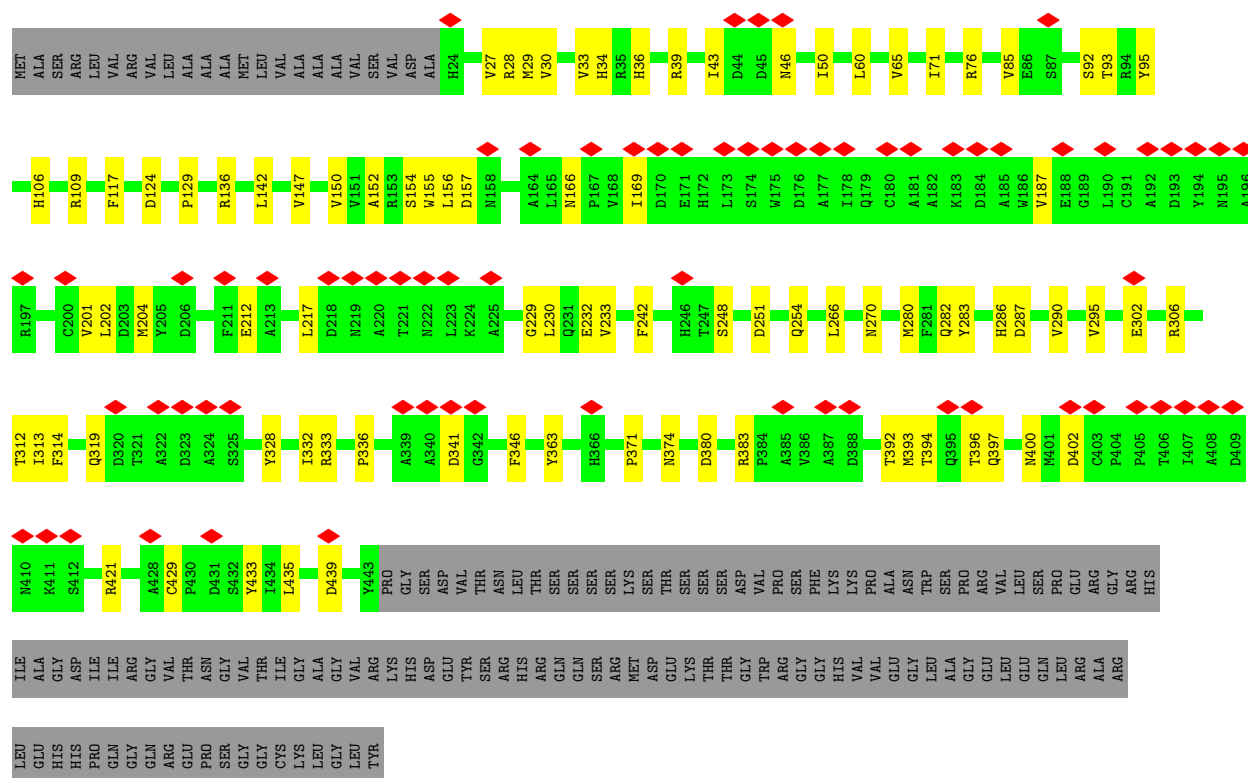


Mol	Chain	Residues	Atoms			AltConf
5	A	1	Total	O	P	0
			5	4	1	
5	B	1	Total	O	P	0
			5	4	1	
5	C	1	Total	O	P	0
			5	4	1	

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		AltConf
6	A	8	Total	O	0
			8	8	
6	B	8	Total	O	0
			8	8	
6	C	3	Total	O	0
			3	3	

- Molecule 1: Secreted acid phosphatase 1 (SAP1)



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



- Molecule 2: alpha-D-mannopyranose-(1-4)-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 F:  100%



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

Chain G:  50% 50%



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

Chain K:  50% 50%



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

Chain L:  50% 50%



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

Chain M:  50% 100%



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

Chain P:  50% 100%



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

Chain Q:  50% 50%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	524666	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS GLACIOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	30	Depositor
Minimum defocus (nm)	750	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	130000	Depositor
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	2.519	Depositor
Minimum map value	-1.855	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.036	Depositor
Recommended contour level	0.3	Depositor
Map size (Å)	341.76, 341.76, 341.76	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.89000005, 0.89000005, 0.89000005	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: MAN, PO4, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.15	0/3375	0.34	0/4612
1	B	0.17	0/3383	0.36	0/4623
1	C	0.13	0/3357	0.31	0/4587
All	All	0.15	0/10115	0.34	0/13822

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	3292	0	3136	51	0
1	B	3300	0	3140	52	0
1	C	3275	0	3121	56	0
2	D	39	0	34	1	0
2	I	39	0	34	4	0
3	F	28	0	25	0	0
3	G	28	0	25	0	0
3	K	28	0	25	0	0
3	L	28	0	25	0	0
3	M	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	P	28	0	25	0	0
3	Q	28	0	25	0	0
4	A	28	0	26	0	0
4	B	14	0	13	0	0
4	C	42	0	39	0	0
5	A	5	0	0	1	0
5	B	5	0	0	1	0
5	C	5	0	0	3	0
6	A	8	0	0	1	0
6	B	8	0	0	1	0
6	C	3	0	0	0	0
All	All	10259	0	9718	155	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 (155) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:ALA:HA	1:A:156:LEU:HB2	1.57	0.86
1:A:431:ASP:OD2	1:A:432:SER:N	2.16	0.78
1:C:152:ALA:HA	1:C:156:LEU:HB2	1.69	0.74
1:A:190:LEU:HG	2:D:1:NAG:H81	1.72	0.71
1:B:152:ALA:HB2	1:B:393:MET:HE2	1.71	0.71
1:B:358:GLU:OE1	1:B:358:GLU:N	2.21	0.67
1:A:444:PRO:HB2	1:B:446:SER:HB3	1.75	0.66
1:B:109:ARG:NH2	5:B:602:PO4:O1	2.27	0.66
1:B:106:HIS:HB3	1:B:136:ARG:HB2	1.79	0.64
1:C:109:ARG:NH2	5:C:604:PO4:O3	2.31	0.64
1:C:117:PHE:CE1	1:C:280:MET:HE1	2.33	0.63
1:A:248:SER:HB3	1:A:251:ASP:HB2	1.81	0.62
1:B:103:ARG:NH2	6:B:701:HOH:O	2.32	0.62
1:A:103:ARG:NH1	6:A:701:HOH:O	2.28	0.61
1:B:232:GLU:HG2	1:B:306:ARG:HH22	1.67	0.59
1:A:436:SER:HB2	1:A:443:TYR:HE1	1.67	0.59
1:B:258:SER:HB2	1:B:292:PRO:HB2	1.85	0.59
1:C:30:VAL:HG13	1:C:280:MET:HG3	1.84	0.59
1:B:147:VAL:HB	1:B:150:VAL:HG22	1.84	0.58
1:A:30:VAL:HG13	1:A:280:MET:HG3	1.84	0.58
1:C:34:HIS:HB3	1:C:312:THR:HG23	1.85	0.58
1:B:30:VAL:HG13	1:B:280:MET:HG3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:394:THR:HB	1:C:397:GLN:HG3	1.86	0.58
1:A:147:VAL:HB	1:A:150:VAL:HG22	1.86	0.57
1:A:34:HIS:HB3	1:A:312:THR:HG23	1.85	0.57
1:C:154:SER:HG	1:C:155:TRP:CD1	2.23	0.57
1:B:287:ASP:N	1:B:287:ASP:OD1	2.38	0.57
1:A:336:PRO:HA	1:A:346:PHE:HA	1.87	0.56
1:A:103:ARG:HH11	1:A:103:ARG:HG2	1.70	0.56
1:B:106:HIS:HB2	1:B:134:THR:O	2.05	0.56
1:C:287:ASP:N	1:C:287:ASP:OD1	2.38	0.56
1:B:64:GLY:HA2	1:B:67:MET:HG2	1.87	0.56
1:A:100:VAL:HG22	1:A:280:MET:HB3	1.87	0.56
1:C:232:GLU:HG3	1:C:306:ARG:HH22	1.71	0.56
1:A:287:ASP:OD2	1:A:287:ASP:N	2.39	0.55
1:A:394:THR:HB	1:A:397:GLN:HG3	1.89	0.54
1:C:106:HIS:HB3	1:C:136:ARG:HB2	1.90	0.54
1:A:154:SER:HG	1:A:155:TRP:CD1	2.26	0.54
1:A:436:SER:HB2	1:A:443:TYR:CE1	2.42	0.54
1:C:157:ASP:HB2	1:C:397:GLN:HG2	1.90	0.53
1:C:363:TYR:OH	1:C:374:ASN:ND2	2.39	0.53
1:A:254:GLN:NE2	1:A:392:THR:O	2.37	0.53
1:C:314:PHE:HB2	1:C:332:ILE:HB	1.91	0.53
1:B:352:LYS:HG2	1:B:364:LEU:HD11	1.91	0.53
1:B:313:ILE:HG12	1:B:333:ARG:HG3	1.91	0.53
1:C:95:TYR:CE2	1:C:129:PRO:HA	2.44	0.52
1:C:147:VAL:HB	1:C:150:VAL:HG22	1.90	0.52
1:B:254:GLN:NE2	1:B:392:THR:O	2.42	0.52
1:A:396:THR:O	1:A:400:ASN:ND2	2.41	0.52
1:A:185:ALA:HB3	1:A:187:VAL:HG12	1.91	0.52
1:B:190:LEU:HG	2:I:1:NAG:H81	1.91	0.51
1:C:46:ASN:O	1:C:50:ILE:HG13	2.09	0.51
1:C:28:ARG:HE	1:C:319:GLN:HB3	1.75	0.51
1:C:229:GLY:O	1:C:233:VAL:HG23	2.11	0.51
1:A:333:ARG:O	1:A:348:GLU:HG2	2.11	0.51
1:C:60:LEU:HD23	1:C:65:VAL:HG12	1.92	0.51
1:A:421:ARG:HD3	1:A:435:LEU:HD11	1.93	0.50
1:A:258:SER:HB2	1:A:292:PRO:HB2	1.93	0.50
1:A:165:LEU:O	1:A:168:VAL:HG12	2.12	0.50
1:A:109:ARG:NH2	5:A:603:PO4:O1	2.45	0.49
1:B:212:GLU:HG3	1:B:217:LEU:HD22	1.94	0.49
1:C:36:HIS:NE2	5:C:604:PO4:O1	2.45	0.49
1:C:286:HIS:O	1:C:290:VAL:HG23	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:242:PHE:CD1	1:B:292:PRO:HG3	2.48	0.49
1:B:189:GLY:HA2	2:I:2:NAG:H62	1.93	0.49
1:B:156:LEU:HD13	1:B:397:GLN:O	2.13	0.49
1:B:380:ASP:HA	1:B:383:ARG:HG3	1.94	0.49
1:C:254:GLN:NE2	1:C:392:THR:O	2.42	0.49
1:B:64:GLY:O	1:B:68:VAL:HG23	2.13	0.49
1:A:363:TYR:OH	1:A:374:ASN:ND2	2.42	0.48
1:A:95:TYR:CE2	1:A:129:PRO:HA	2.48	0.48
1:A:106:HIS:HB3	1:A:136:ARG:HB2	1.95	0.48
1:A:242:PHE:CD1	1:A:292:PRO:HG3	2.49	0.48
1:A:187:VAL:HG23	1:A:190:LEU:HD12	1.96	0.48
1:B:363:TYR:OH	1:B:374:ASN:ND2	2.42	0.48
1:A:404:PRO:O	1:A:421:ARG:NH2	2.41	0.48
1:B:40:SER:HB2	1:B:57:CYS:HA	1.96	0.48
1:B:401:MET:HE1	1:B:420:TYR:HB2	1.95	0.48
1:C:380:ASP:HA	1:C:383:ARG:HG3	1.94	0.48
1:C:421:ARG:HD3	1:C:435:LEU:HD11	1.95	0.48
1:A:152:ALA:HB1	1:A:397:GLN:HB3	1.95	0.47
1:B:235:ALA:HB2	1:B:306:ARG:HG2	1.96	0.47
1:C:439:ASP:OD1	1:C:439:ASP:N	2.43	0.47
1:C:39:ARG:NH2	5:C:604:PO4:O2	2.36	0.47
1:C:302:GLU:O	1:C:306:ARG:HG3	2.15	0.47
1:A:44:ASP:OD1	1:A:44:ASP:N	2.46	0.47
1:B:65:VAL:O	1:B:69:ARG:HG3	2.16	0.46
1:B:200:CYS:O	1:B:204:MET:HG3	2.15	0.46
1:B:439:ASP:OD1	1:B:439:ASP:N	2.43	0.46
1:C:71:ILE:HG22	1:C:117:PHE:HE2	1.79	0.46
1:C:27:VAL:HB	1:C:85:VAL:HG12	1.97	0.46
1:A:187:VAL:HG11	1:A:204:MET:HG2	1.98	0.46
1:C:242:PHE:O	1:C:295:VAL:HG21	2.16	0.46
1:C:280:MET:HE2	1:C:282:GLN:HB2	1.98	0.46
1:C:43:ILE:O	1:C:43:ILE:HD12	2.16	0.46
1:C:152:ALA:HB2	1:C:393:MET:HE3	1.98	0.46
1:B:336:PRO:HA	1:B:346:PHE:HA	1.97	0.45
1:A:286:HIS:O	1:A:290:VAL:HG23	2.16	0.45
1:C:76:ARG:NH2	1:C:92:SER:O	2.50	0.45
1:B:67:MET:O	1:B:336:PRO:HG2	2.17	0.45
1:B:74:PHE:CZ	1:B:78:ARG:HD2	2.52	0.45
1:C:166:ASN:O	1:C:169:ILE:HG13	2.17	0.44
1:C:429:CYS:HB3	1:C:433:TYR:HB3	1.98	0.44
1:A:157:ASP:HB2	1:A:397:GLN:HG2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:336:PRO:HA	1:C:346:PHE:HA	2.00	0.44
1:A:76:ARG:NH2	1:A:92:SER:O	2.51	0.44
1:B:174:SER:OG	1:B:175:TRP:N	2.51	0.44
1:B:248:SER:HB3	1:B:251:ASP:HB2	1.98	0.44
1:C:341:ASP:OD2	1:C:341:ASP:N	2.48	0.44
1:A:165:LEU:O	1:A:169:ILE:HG13	2.18	0.44
1:C:380:ASP:OD1	1:C:383:ARG:NH1	2.50	0.44
1:B:76:ARG:NH2	1:B:92:SER:O	2.51	0.43
1:A:235:ALA:HB2	1:A:306:ARG:HG2	1.99	0.43
1:C:33:VAL:HG22	1:C:283:TYR:HB2	2.01	0.43
1:A:119:ARG:HA	1:A:119:ARG:HD2	1.67	0.43
1:A:434:ILE:HB	1:B:434:ILE:HD11	1.99	0.43
1:C:229:GLY:HA2	1:C:232:GLU:OE1	2.19	0.43
1:C:29:MET:HE2	1:C:29:MET:HB3	1.75	0.43
1:A:64:GLY:O	1:A:68:VAL:HG23	2.20	0.42
1:A:64:GLY:HA2	1:A:67:MET:HG2	2.02	0.42
1:C:396:THR:O	1:C:400:ASN:ND2	2.52	0.42
1:C:201:VAL:HG11	1:C:233:VAL:HG11	2.01	0.42
1:C:212:GLU:HG3	1:C:217:LEU:HD22	2.01	0.42
1:A:401:MET:HE3	1:A:403:CYS:SG	2.60	0.42
1:C:93:THR:O	1:C:124:ASP:HB2	2.19	0.42
1:A:436:SER:O	1:B:437:ALA:HB2	2.20	0.42
1:C:328:TYR:HD1	1:C:371:PRO:HA	1.85	0.42
1:B:29:MET:HE1	1:B:269:ILE:HA	2.02	0.41
1:C:43:ILE:HG12	1:C:50:ILE:HD12	2.02	0.41
1:C:142:LEU:HD21	1:C:283:TYR:HB3	2.02	0.41
1:B:67:MET:HE2	1:B:344:TYR:CE2	2.54	0.41
1:B:333:ARG:O	1:B:348:GLU:HG2	2.20	0.41
1:C:156:LEU:HD23	1:C:156:LEU:HA	1.82	0.41
1:B:71:ILE:HD13	1:B:312:THR:OG1	2.21	0.41
1:C:201:VAL:HG13	1:C:230:LEU:CD2	2.50	0.41
1:A:193:ASP:OD1	1:A:193:ASP:N	2.44	0.41
1:B:156:LEU:HD23	1:B:156:LEU:HA	1.85	0.41
1:B:224:LYS:HE3	1:B:224:LYS:HB3	1.91	0.41
1:B:286:HIS:O	1:B:290:VAL:HG23	2.20	0.41
1:C:266:LEU:O	1:C:270:ASN:ND2	2.54	0.41
1:A:129:PRO:O	1:B:115:THR:HG21	2.20	0.41
1:B:45:ASP:HB3	2:I:1:NAG:H82	2.03	0.41
1:A:33:VAL:HG22	1:A:283:TYR:HB2	2.02	0.41
1:A:78:ARG:NH1	1:A:84:LEU:HD13	2.35	0.41
1:B:202:LEU:HD23	1:B:202:LEU:HA	1.92	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:248:SER:HB3	1:C:251:ASP:HB2	2.01	0.41
1:A:416:ARG:HG3	1:A:416:ARG:NH1	2.36	0.41
1:C:187:VAL:HG11	1:C:204:MET:HG3	2.03	0.41
1:C:313:ILE:HG12	1:C:333:ARG:HG3	2.02	0.41
1:B:201:VAL:HA	1:B:204:MET:HG3	2.02	0.41
1:A:266:LEU:O	1:A:270:ASN:ND2	2.54	0.40
1:C:202:LEU:HD23	1:C:202:LEU:HA	1.95	0.40
1:B:29:MET:HE2	1:B:29:MET:HB3	1.78	0.40
1:B:45:ASP:CB	2:I:1:NAG:H82	2.51	0.40
1:B:60:LEU:HD23	1:B:65:VAL:HG12	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	421/565 (74%)	405 (96%)	16 (4%)	0	100	100
1	B	422/565 (75%)	415 (98%)	7 (2%)	0	100	100
1	C	418/565 (74%)	410 (98%)	8 (2%)	0	100	100
All	All	1261/1695 (74%)	1230 (98%)	31 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	357/471 (76%)	356 (100%)	1 (0%)	86	91
1	B	358/471 (76%)	358 (100%)	0	100	100
1	C	355/471 (75%)	354 (100%)	1 (0%)	86	91
All	All	1070/1413 (76%)	1068 (100%)	2 (0%)	85	92

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	347	GLN
1	C	402	ASP

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

Mol	Chain	Res	Type
1	A	31	GLN
1	A	101	HIS
1	A	166	ASN
1	A	172	HIS
1	A	268	ASN
1	A	272	HIS
1	A	282	GLN
1	A	286	HIS
1	A	300	GLN
1	A	335	ASN
1	B	106	HIS
1	B	166	ASN
1	B	231	GLN
1	B	272	HIS
1	B	373	ASN
1	C	31	GLN
1	C	101	HIS
1	C	166	ASN
1	C	179	GLN
1	C	231	GLN
1	C	268	ASN
1	C	272	HIS
1	C	279	ASN
1	C	282	GLN
1	C	335	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

20 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	$\# Z > 2$
2	NAG	D	1	1,2	14,14,15	0.81	0	17,19,21	1.28	2 (11%)
2	NAG	D	2	2	14,14,15	0.82	1 (7%)	17,19,21	1.56	3 (17%)
2	MAN	D	3	2	11,11,12	0.67	0	15,15,17	1.43	1 (6%)
3	NAG	F	1	3,1	14,14,15	0.97	0	17,19,21	1.03	0
3	NAG	F	2	3	14,14,15	0.70	0	17,19,21	1.13	0
3	NAG	G	1	3,1	14,14,15	0.80	0	17,19,21	0.97	0
3	NAG	G	2	3	14,14,15	0.81	0	17,19,21	1.11	1 (5%)
2	NAG	I	1	1,2	14,14,15	0.72	0	17,19,21	1.44	3 (17%)
2	NAG	I	2	2	14,14,15	0.81	0	17,19,21	1.30	2 (11%)
2	MAN	I	3	2	11,11,12	0.65	0	15,15,17	1.44	1 (6%)
3	NAG	K	1	3,1	14,14,15	0.87	0	17,19,21	1.02	0
3	NAG	K	2	3	14,14,15	0.74	0	17,19,21	1.26	1 (5%)
3	NAG	L	1	3,1	14,14,15	0.86	0	17,19,21	0.92	0
3	NAG	L	2	3	14,14,15	0.79	0	17,19,21	1.31	1 (5%)
3	NAG	M	1	3,1	14,14,15	0.80	0	17,19,21	1.07	1 (5%)
3	NAG	M	2	3	14,14,15	0.82	0	17,19,21	1.07	1 (5%)
3	NAG	P	1	3,1	14,14,15	0.82	0	17,19,21	1.09	2 (11%)
3	NAG	P	2	3	14,14,15	0.85	0	17,19,21	1.12	1 (5%)
3	NAG	Q	1	3,1	14,14,15	0.79	0	17,19,21	0.85	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	Q	2	3	14,14,15	0.85	0	17,19,21	1.22	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1
2	MAN	D	3	2	-	2/2/19/22	1/1/1/1
3	NAG	F	1	3,1	-	3/6/23/26	0/1/1/1
3	NAG	F	2	3	-	1/6/23/26	0/1/1/1
3	NAG	G	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
2	NAG	I	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	I	2	2	-	2/6/23/26	0/1/1/1
2	MAN	I	3	2	-	2/2/19/22	1/1/1/1
3	NAG	K	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	K	2	3	-	2/6/23/26	0/1/1/1
3	NAG	L	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	L	2	3	-	2/6/23/26	0/1/1/1
3	NAG	M	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	M	2	3	-	2/6/23/26	0/1/1/1
3	NAG	P	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	P	2	3	-	2/6/23/26	0/1/1/1
3	NAG	Q	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	Q	2	3	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	NAG	O5-C1	-2.07	1.40	1.43

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	NAG	O5-C1-C2	-4.74	103.81	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3	MAN	C1-O5-C5	4.08	117.72	112.19
2	I	3	MAN	C1-O5-C5	4.08	117.72	112.19
3	L	2	NAG	C1-O5-C5	4.01	117.62	112.19
3	K	2	NAG	O5-C5-C6	3.13	112.11	107.20
2	D	1	NAG	O4-C4-C3	-3.11	103.15	110.35
3	Q	2	NAG	O5-C5-C6	3.03	111.95	107.20
2	I	1	NAG	C1-O5-C5	3.02	116.29	112.19
2	I	1	NAG	O5-C1-C2	-2.99	106.57	111.29
3	M	2	NAG	O5-C5-C6	2.89	111.73	107.20
3	G	2	NAG	C1-O5-C5	2.82	116.02	112.19
2	I	2	NAG	O3-C3-C2	-2.60	104.09	109.47
2	D	2	NAG	O3-C3-C2	-2.49	104.31	109.47
3	P	1	NAG	C1-O5-C5	2.47	115.53	112.19
3	P	2	NAG	O5-C5-C6	2.44	111.03	107.20
2	I	1	NAG	C1-C2-N2	2.33	114.47	110.49
3	P	1	NAG	C2-N2-C7	2.25	126.11	122.90
2	I	2	NAG	O4-C4-C3	2.22	115.48	110.35
3	M	1	NAG	C2-N2-C7	2.19	126.02	122.90
2	D	1	NAG	C4-C3-C2	2.11	114.12	111.02
2	D	2	NAG	C6-C5-C4	-2.00	108.31	113.00

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	I	3	MAN	C4-C5-C6-O6
3	Q	2	NAG	O5-C5-C6-O6
3	K	2	NAG	C4-C5-C6-O6
3	M	2	NAG	O5-C5-C6-O6
3	M	2	NAG	C4-C5-C6-O6
2	D	2	NAG	C8-C7-N2-C2
2	D	2	NAG	O7-C7-N2-C2
2	I	2	NAG	C8-C7-N2-C2
2	I	2	NAG	O7-C7-N2-C2
3	F	1	NAG	C8-C7-N2-C2
3	F	1	NAG	O7-C7-N2-C2
3	G	1	NAG	C8-C7-N2-C2
3	G	1	NAG	O7-C7-N2-C2
3	K	1	NAG	C8-C7-N2-C2
3	K	1	NAG	O7-C7-N2-C2
3	L	1	NAG	C8-C7-N2-C2
3	L	1	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	M	1	NAG	C8-C7-N2-C2
3	M	1	NAG	O7-C7-N2-C2
3	P	1	NAG	C8-C7-N2-C2
3	P	1	NAG	O7-C7-N2-C2
3	Q	1	NAG	C8-C7-N2-C2
3	Q	1	NAG	O7-C7-N2-C2
2	I	3	MAN	O5-C5-C6-O6
3	Q	2	NAG	C4-C5-C6-O6
2	D	3	MAN	C4-C5-C6-O6
3	K	2	NAG	O5-C5-C6-O6
2	D	3	MAN	O5-C5-C6-O6
3	P	2	NAG	C4-C5-C6-O6
3	P	2	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	L	2	NAG	C4-C5-C6-O6
3	L	1	NAG	C4-C5-C6-O6
3	L	2	NAG	O5-C5-C6-O6
3	K	1	NAG	C4-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
3	L	1	NAG	O5-C5-C6-O6
2	I	1	NAG	C4-C5-C6-O6
3	K	1	NAG	O5-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
2	I	1	NAG	O5-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6

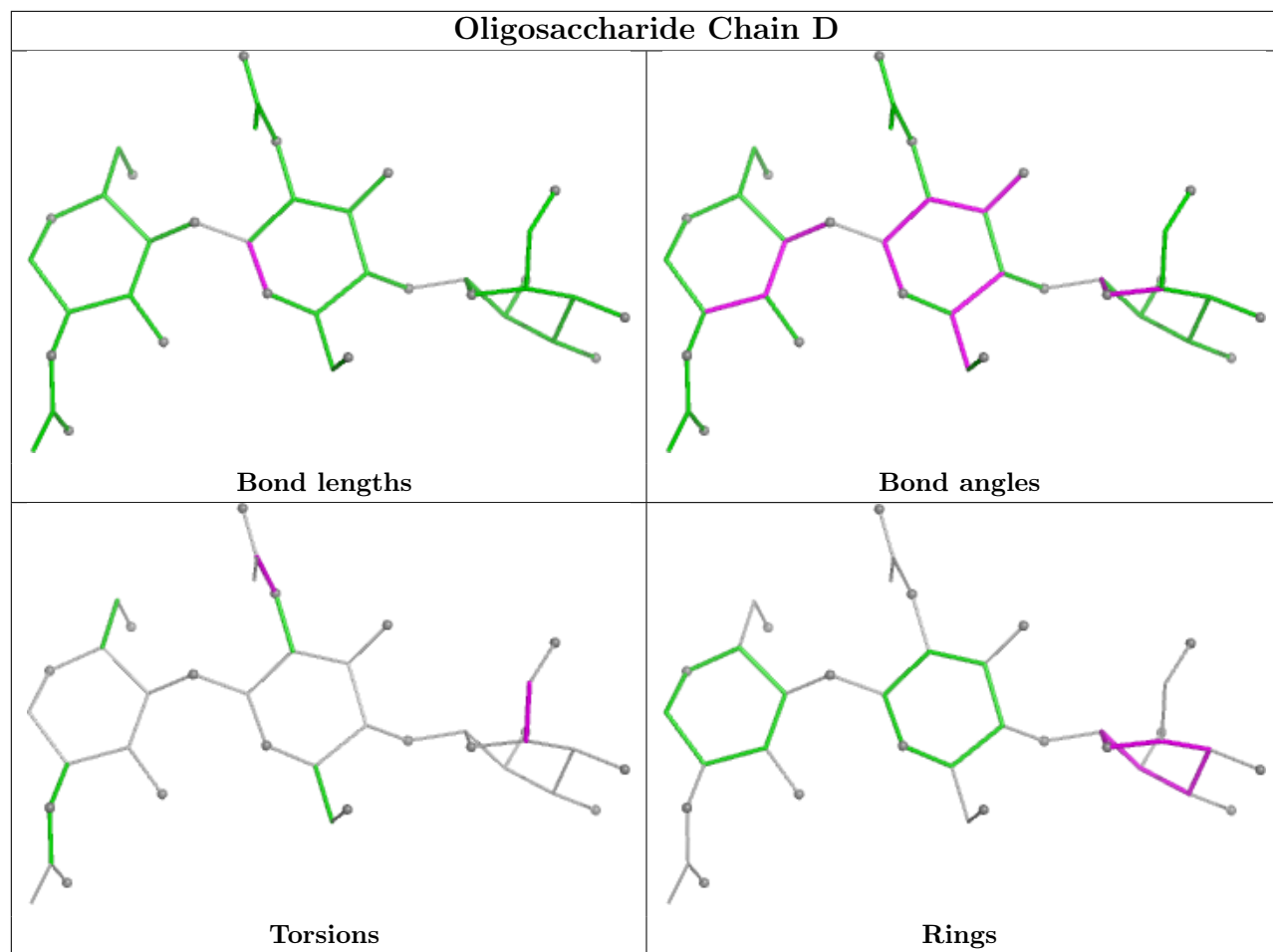
All (2) ring outliers are listed below:

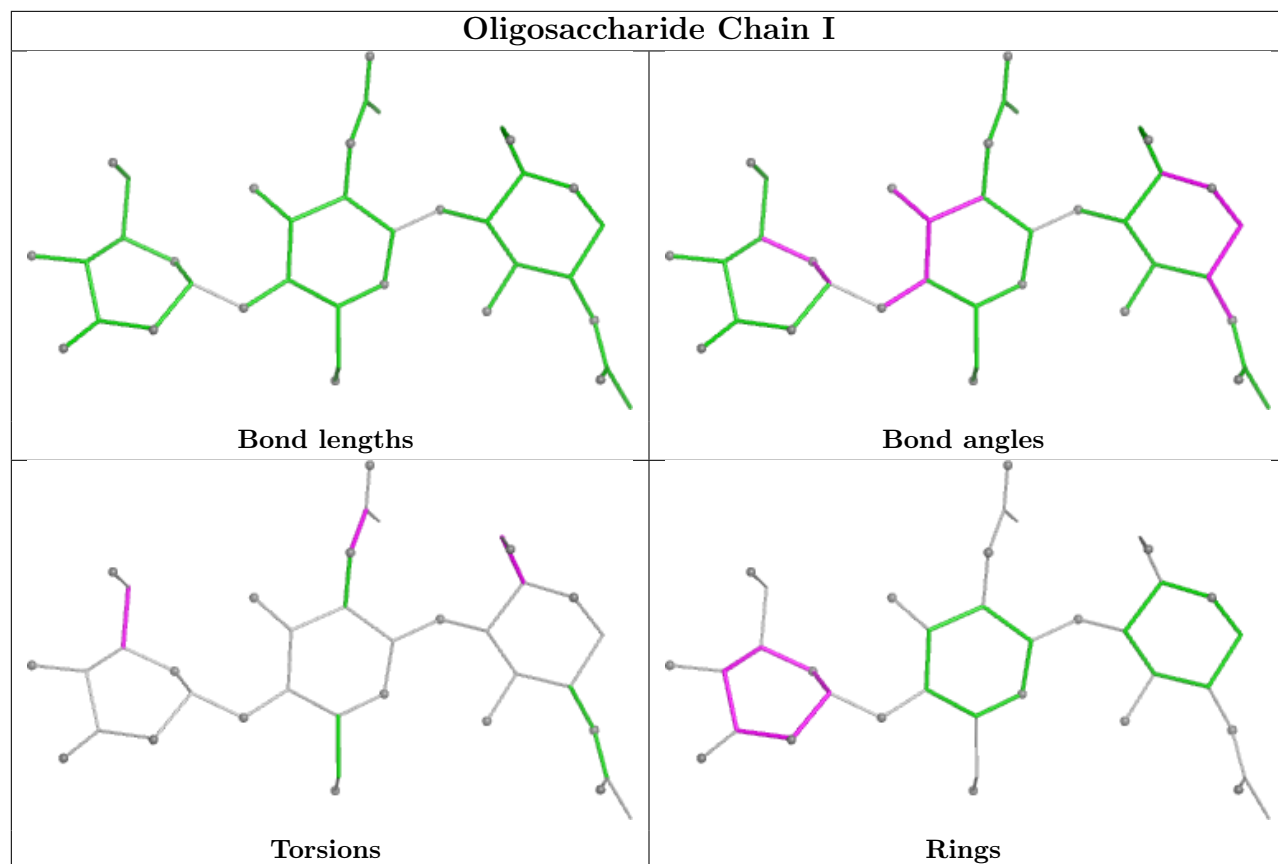
Mol	Chain	Res	Type	Atoms
2	I	3	MAN	C1-C2-C3-C4-C5-O5
2	D	3	MAN	C1-C2-C3-C4-C5-O5

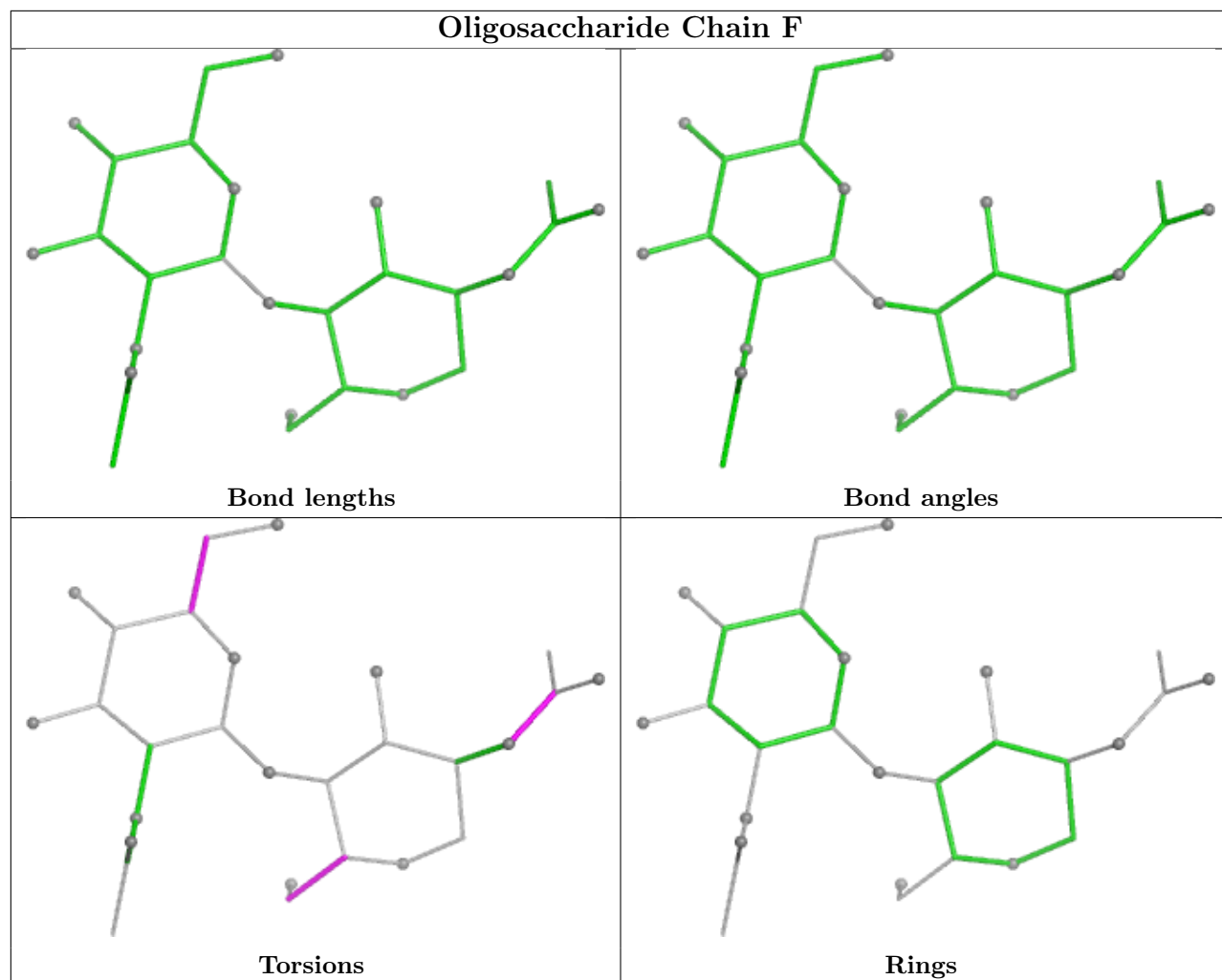
3 monomers are involved in 5 short contacts:

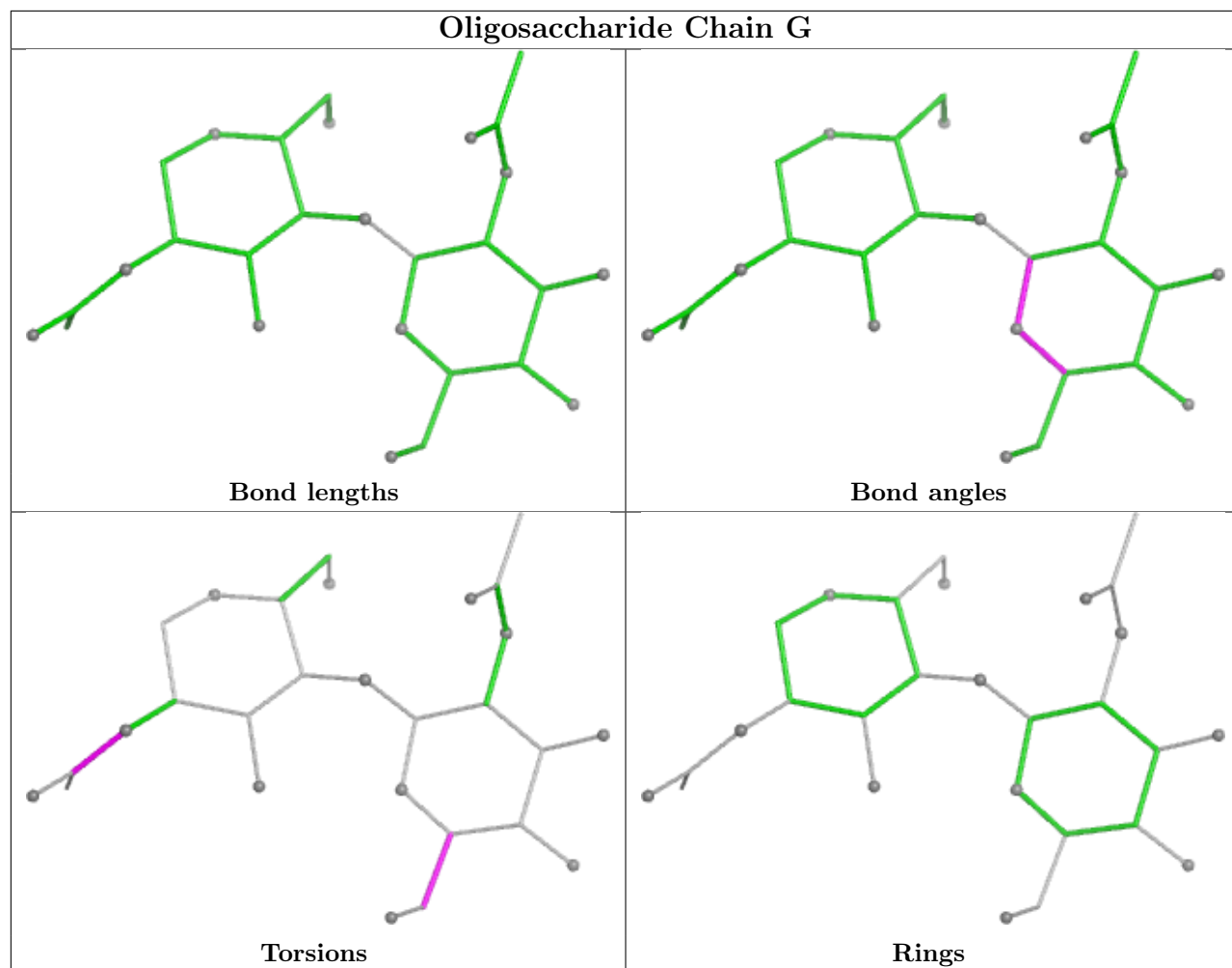
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	NAG	1	0
2	I	2	NAG	1	0
2	I	1	NAG	3	0

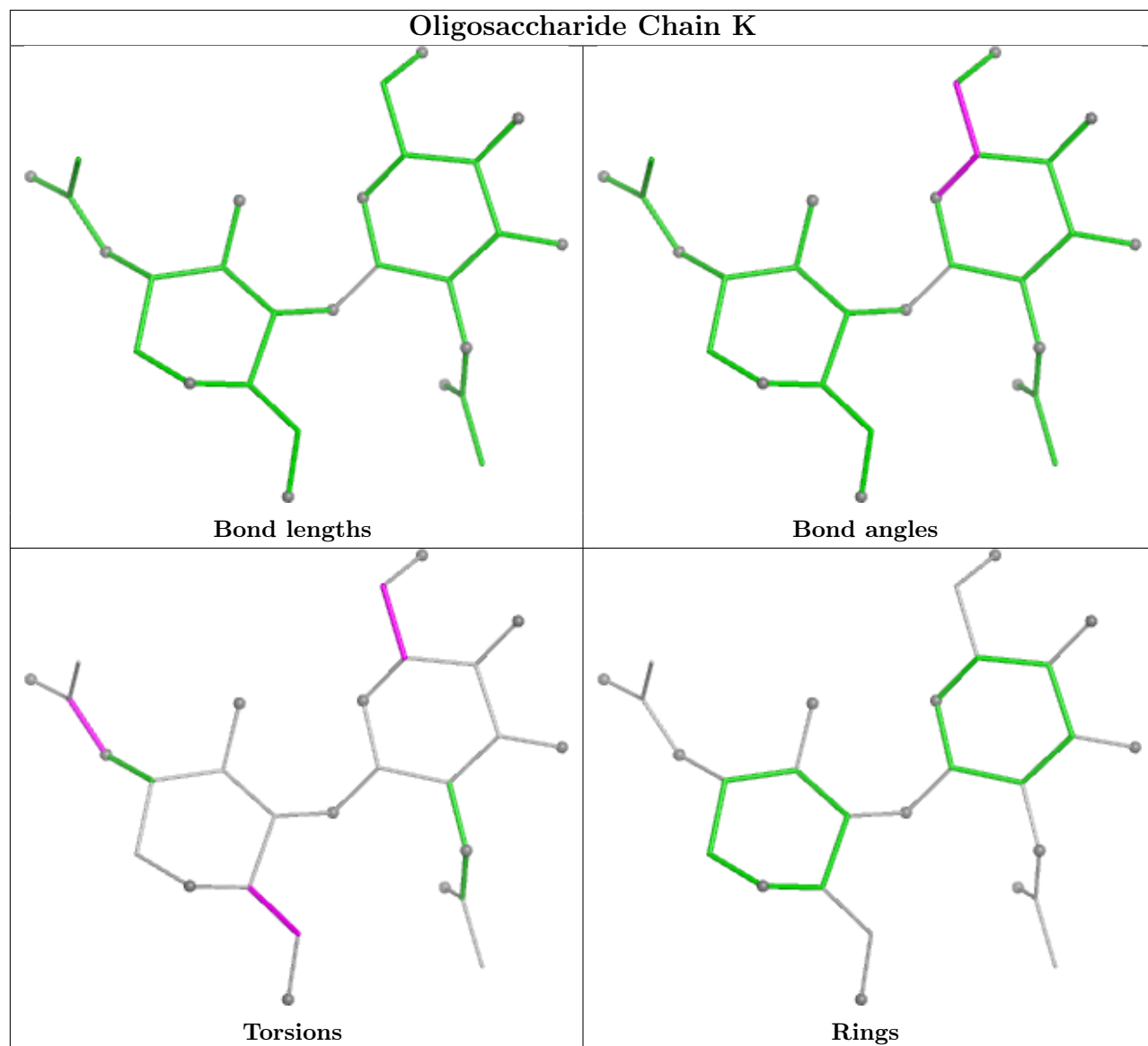
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

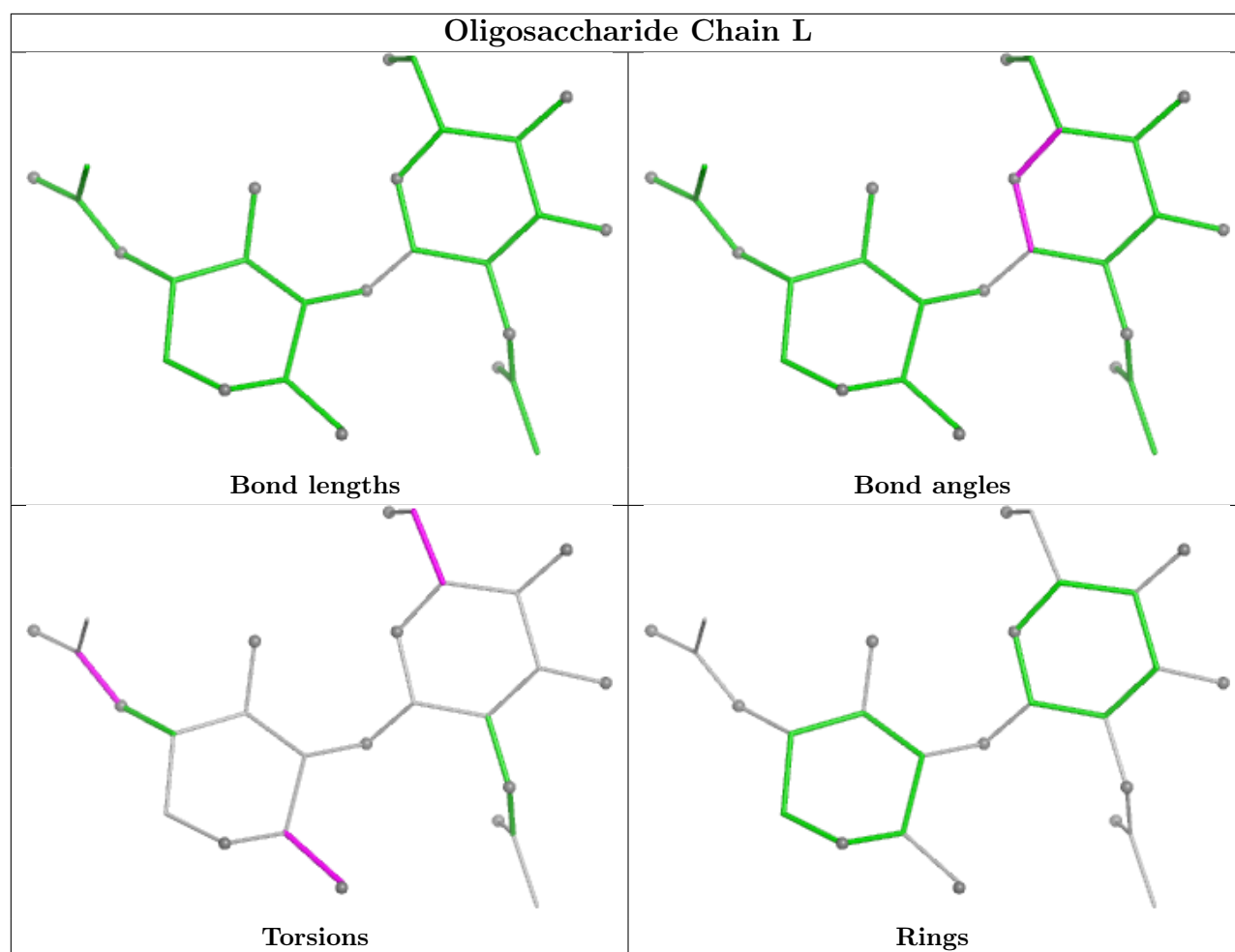


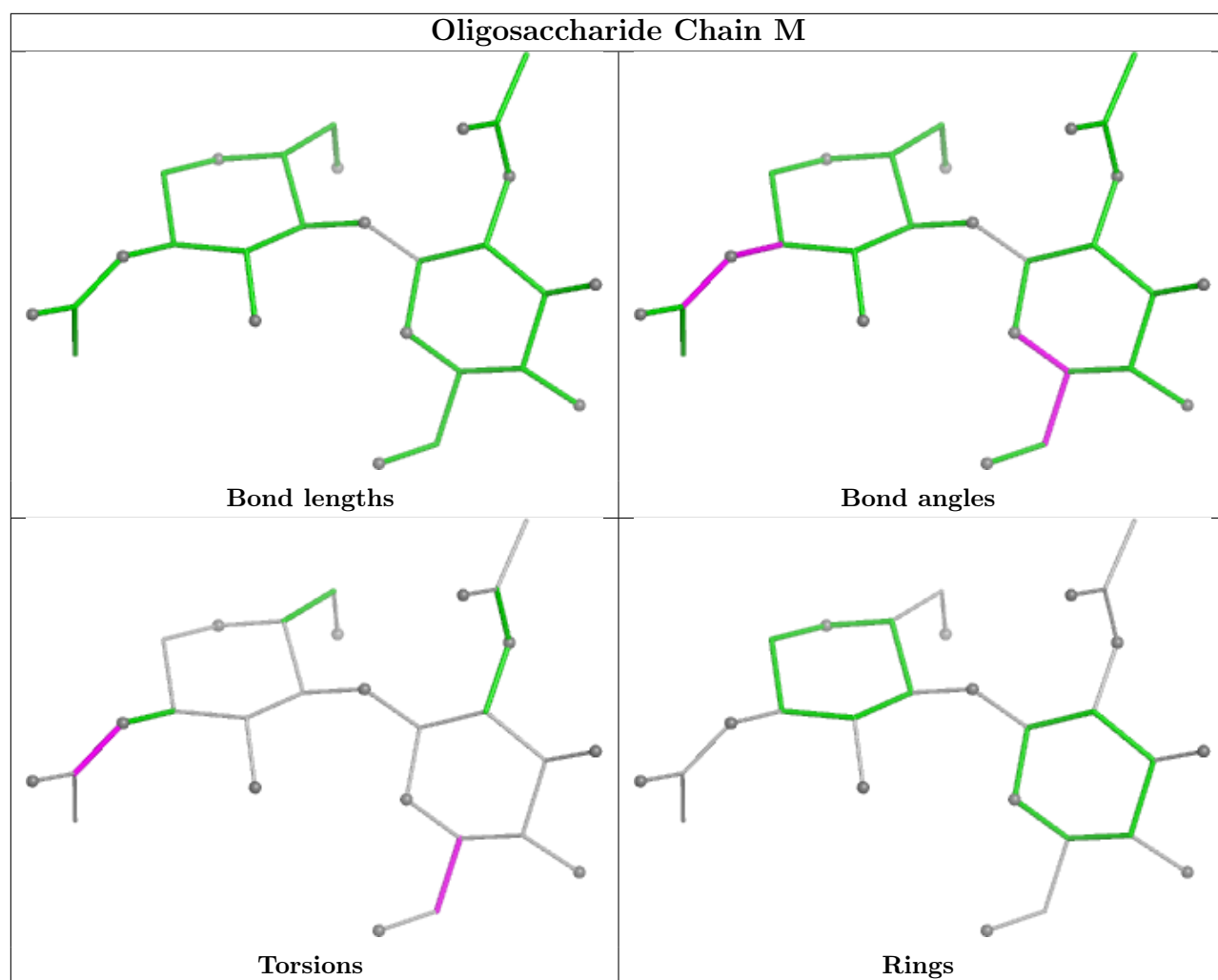


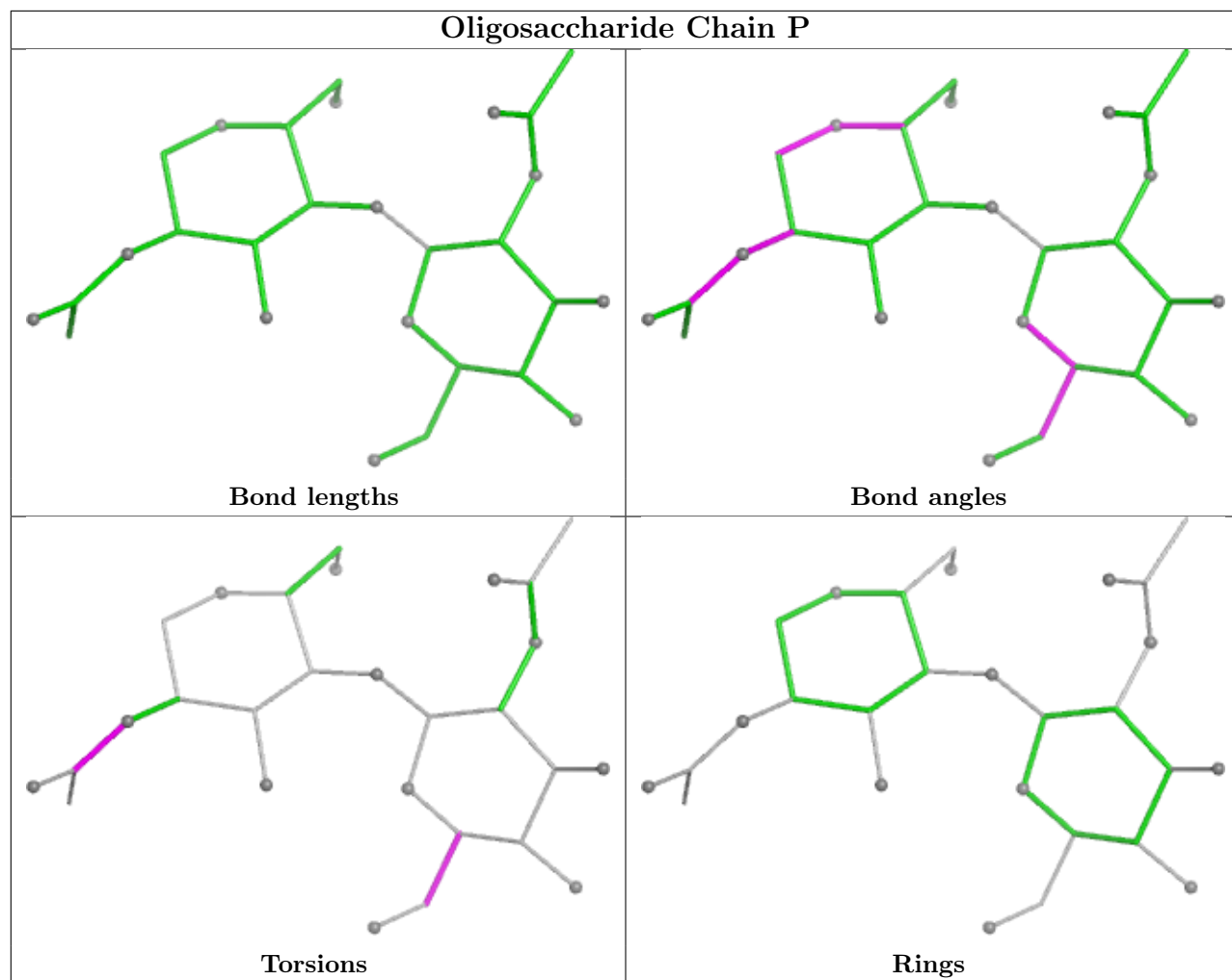


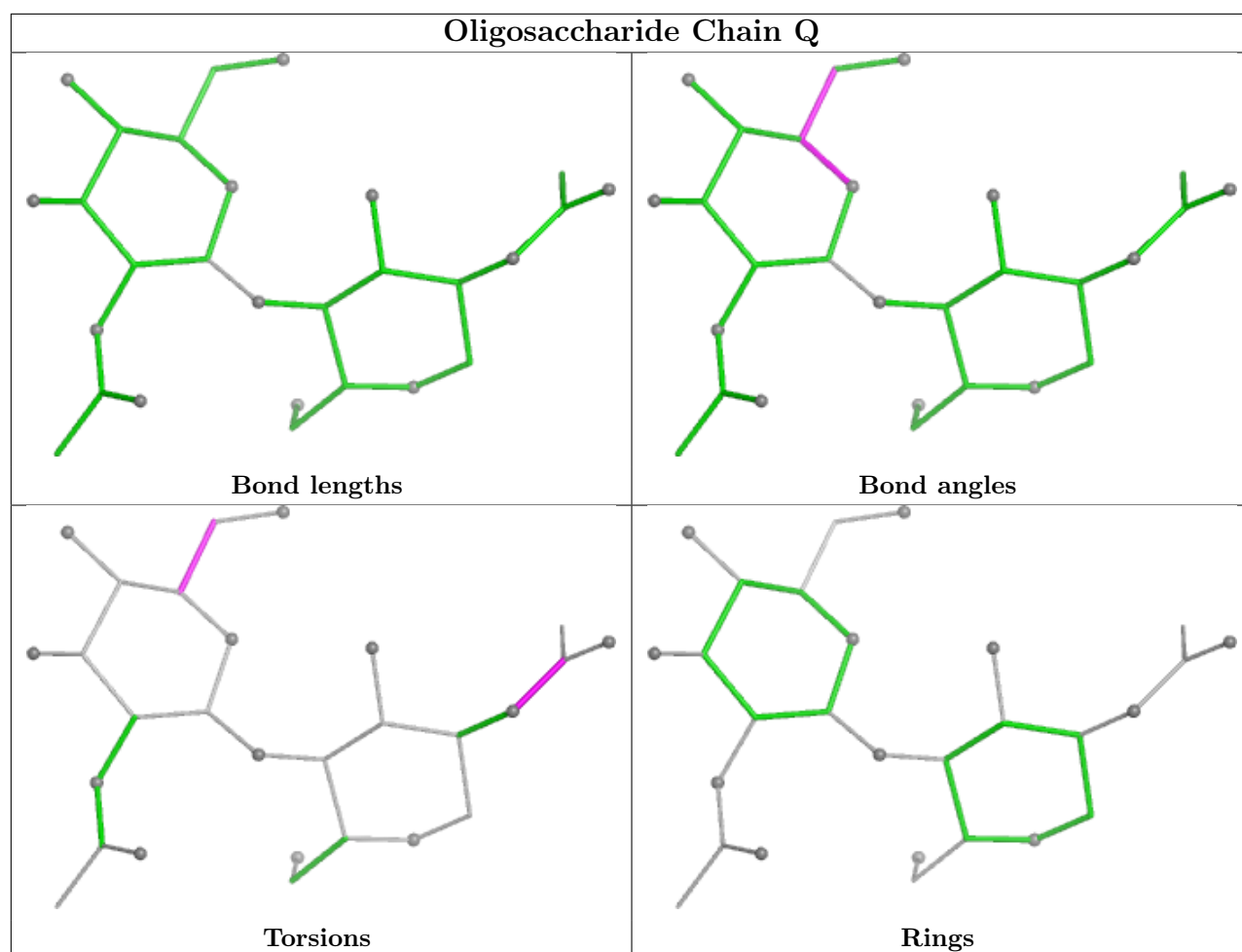












5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	PO4	A	603	-	4,4,4	1.42	1 (25%)	6,6,6	0.45	0
5	PO4	C	604	-	4,4,4	1.44	1 (25%)	6,6,6	0.48	0
4	NAG	B	601	1	14,14,15	0.74	0	17,19,21	1.26	1 (5%)
4	NAG	C	602	1	14,14,15	0.74	0	17,19,21	0.97	1 (5%)
4	NAG	C	601	1	14,14,15	0.73	0	17,19,21	0.95	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PO4	B	602	-	4,4,4	1.40	1 (25%)	6,6,6	0.49	0
4	NAG	A	602	1	14,14,15	0.80	0	17,19,21	1.07	1 (5%)
4	NAG	A	601	1	14,14,15	0.75	0	17,19,21	1.02	1 (5%)
4	NAG	C	603	1	14,14,15	0.73	0	17,19,21	0.94	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	601	1	-	3/6/23/26	0/1/1/1
4	NAG	C	602	1	-	3/6/23/26	0/1/1/1
4	NAG	C	601	1	-	3/6/23/26	0/1/1/1
4	NAG	A	602	1	-	3/6/23/26	0/1/1/1
4	NAG	A	601	1	-	3/6/23/26	0/1/1/1
4	NAG	C	603	1	-	3/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	604	PO4	P-O1	2.51	1.56	1.50
5	A	603	PO4	P-O1	2.50	1.56	1.50
5	B	602	PO4	P-O1	2.44	1.56	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	601	NAG	C1-O5-C5	3.71	117.22	112.19
4	A	601	NAG	C1-O5-C5	2.55	115.65	112.19
4	C	602	NAG	C1-O5-C5	2.44	115.50	112.19
4	A	602	NAG	C1-O5-C5	2.43	115.48	112.19
4	C	601	NAG	C1-O5-C5	2.15	115.11	112.19
4	C	603	NAG	C1-O5-C5	2.14	115.09	112.19

There are no chirality outliers.

All (18) torsion outliers are listed below:

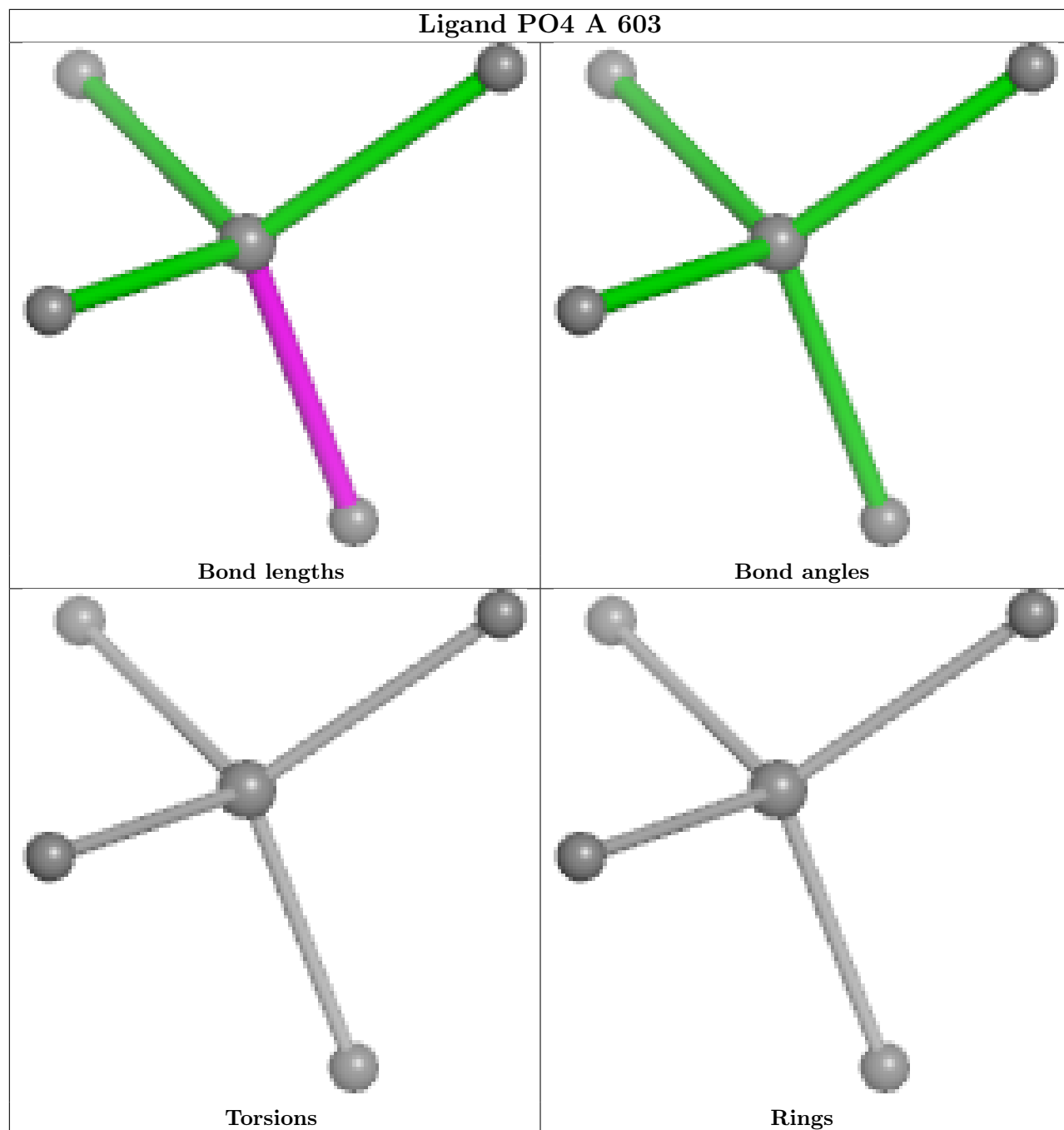
Mol	Chain	Res	Type	Atoms
4	A	601	NAG	C8-C7-N2-C2
4	A	601	NAG	O7-C7-N2-C2
4	A	602	NAG	C8-C7-N2-C2
4	A	602	NAG	O7-C7-N2-C2
4	B	601	NAG	C8-C7-N2-C2
4	B	601	NAG	O7-C7-N2-C2
4	C	601	NAG	C8-C7-N2-C2
4	C	601	NAG	O7-C7-N2-C2
4	C	602	NAG	C8-C7-N2-C2
4	C	602	NAG	O7-C7-N2-C2
4	C	603	NAG	C8-C7-N2-C2
4	C	603	NAG	O7-C7-N2-C2
4	B	601	NAG	O5-C5-C6-O6
4	C	603	NAG	O5-C5-C6-O6
4	C	601	NAG	O5-C5-C6-O6
4	A	602	NAG	O5-C5-C6-O6
4	A	601	NAG	O5-C5-C6-O6
4	C	602	NAG	O5-C5-C6-O6

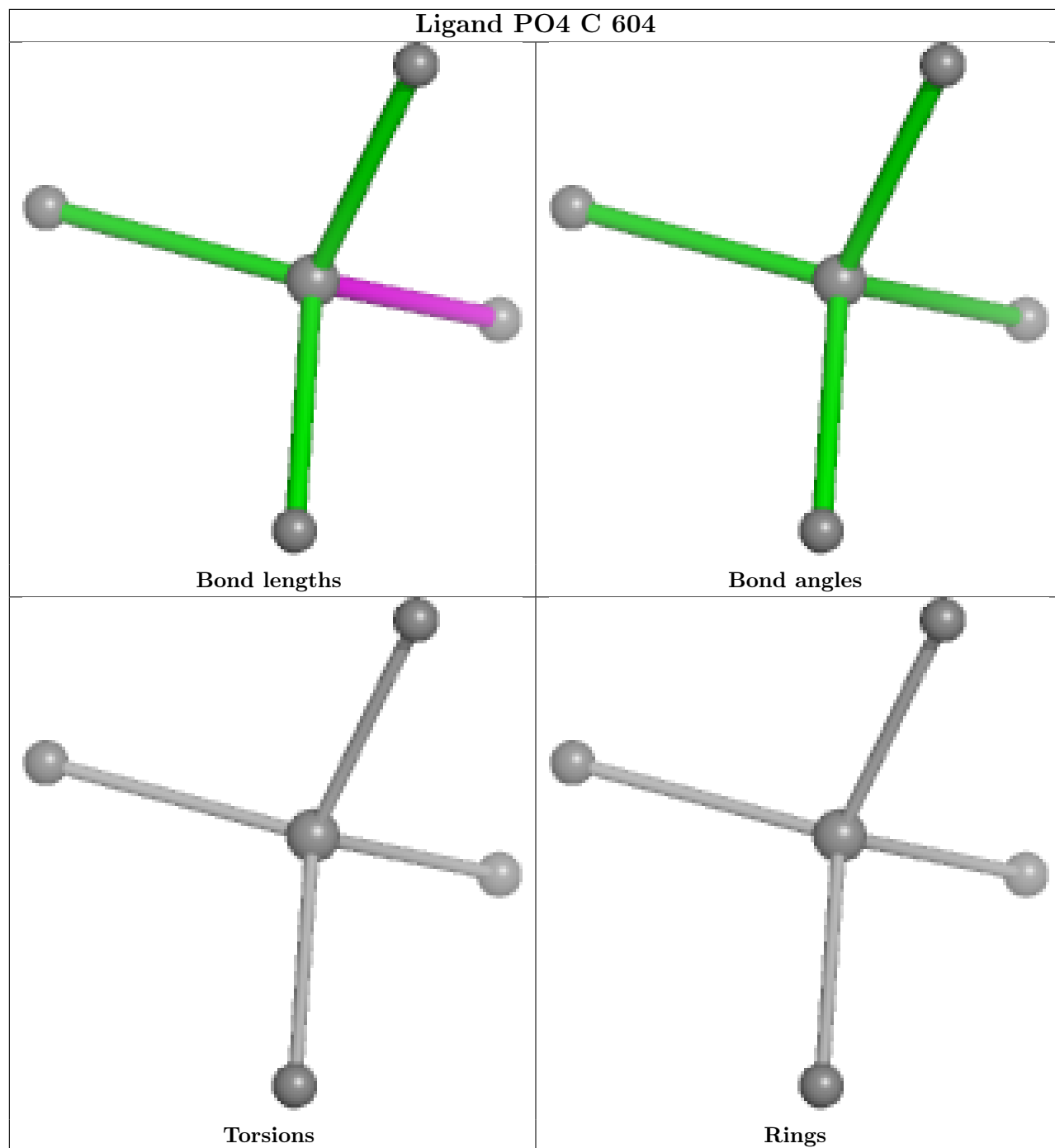
There are no ring outliers.

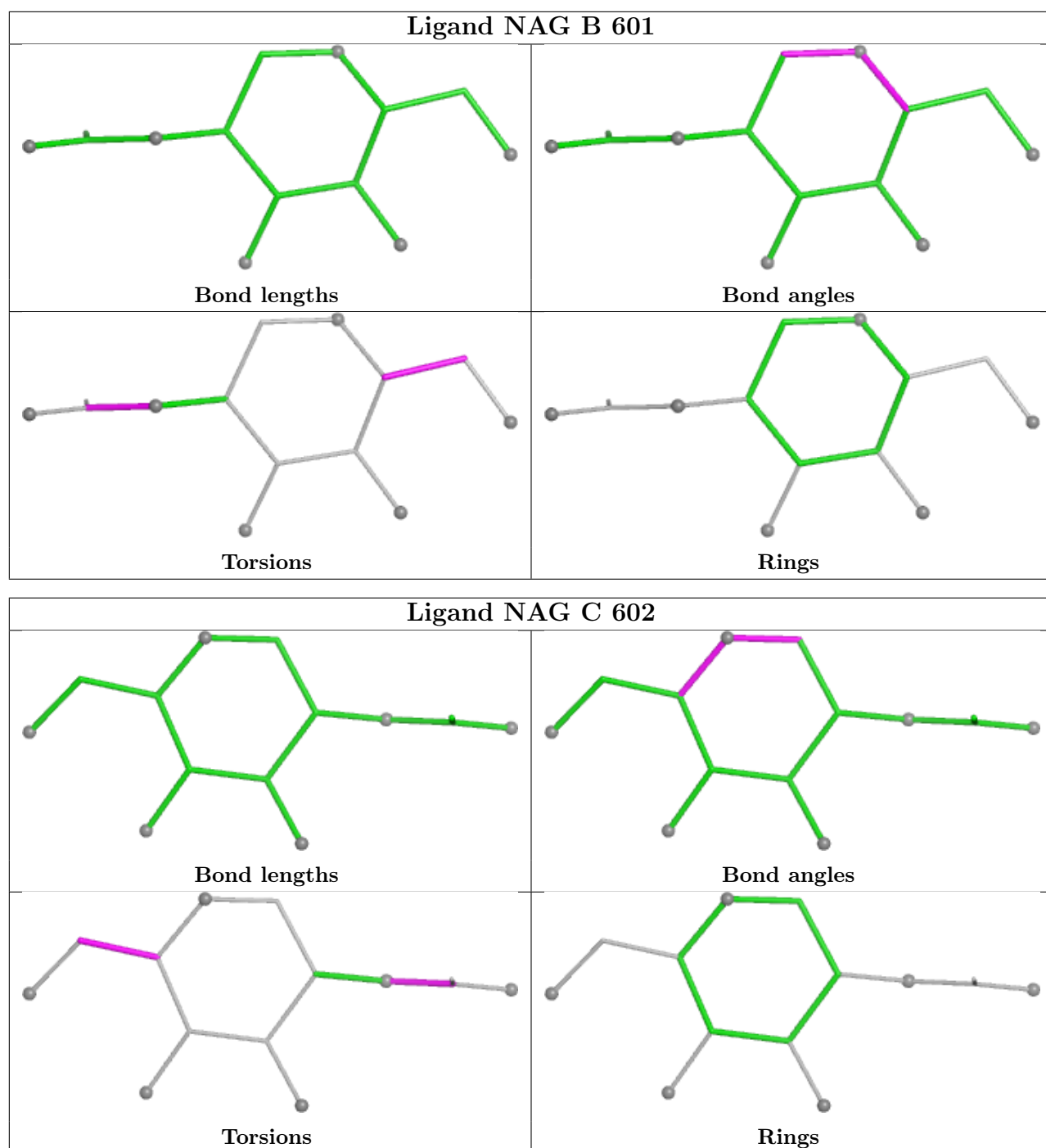
3 monomers are involved in 5 short contacts:

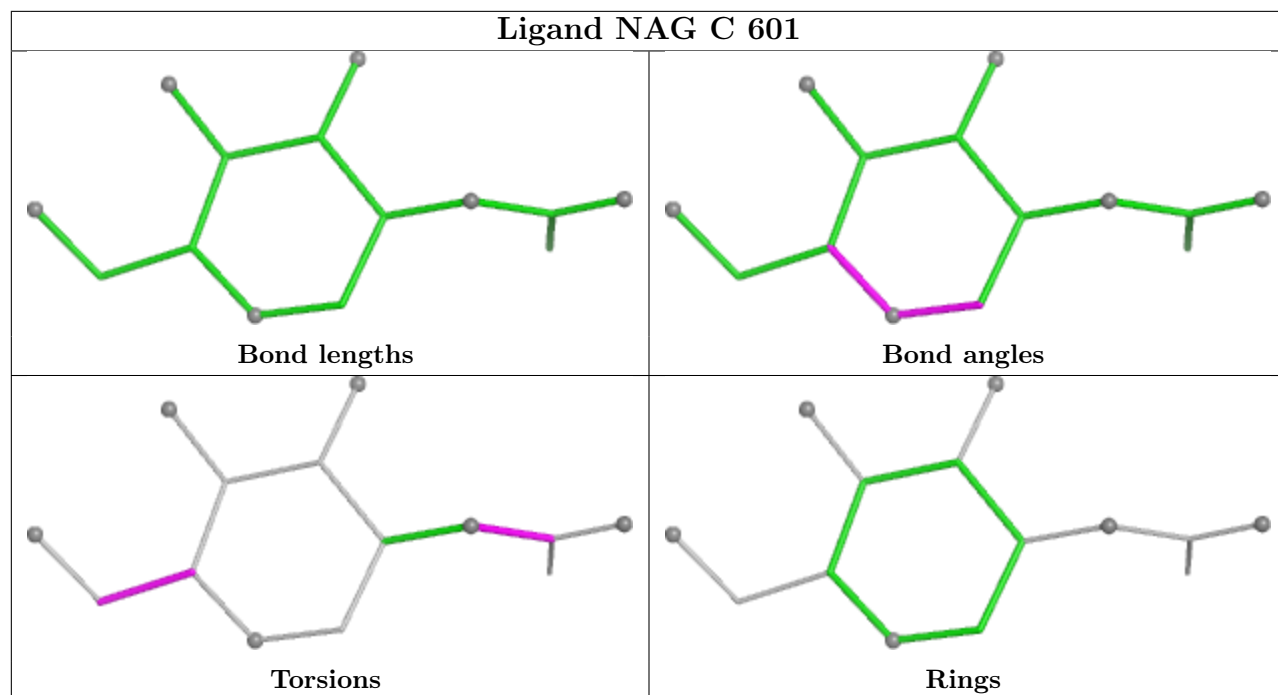
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	603	PO4	1	0
5	C	604	PO4	3	0
5	B	602	PO4	1	0

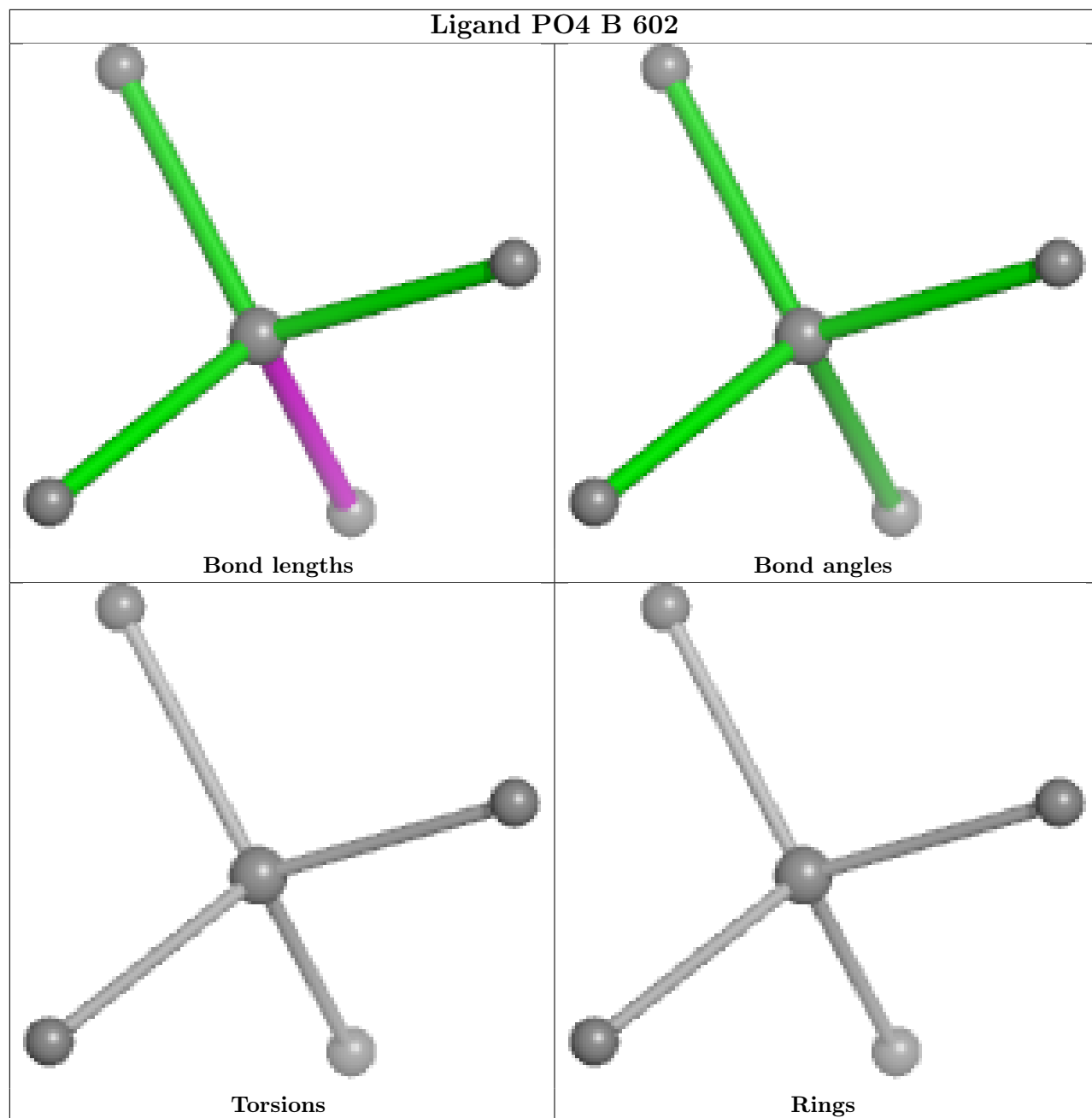
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

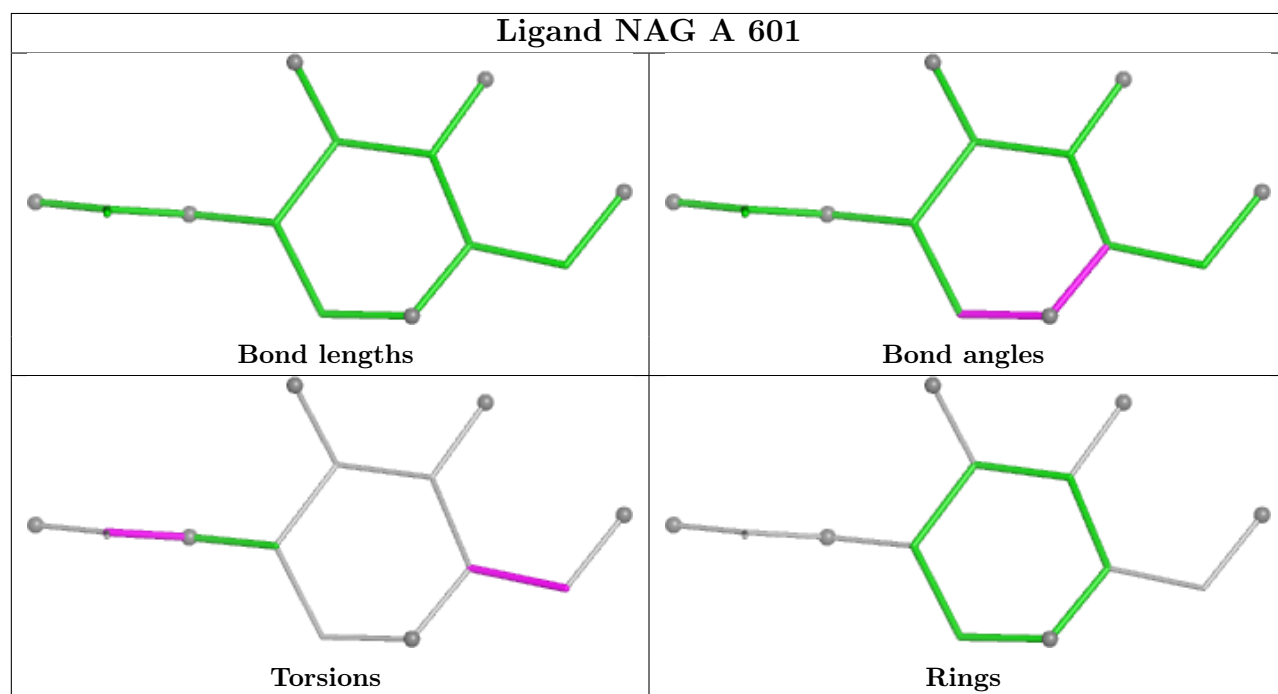
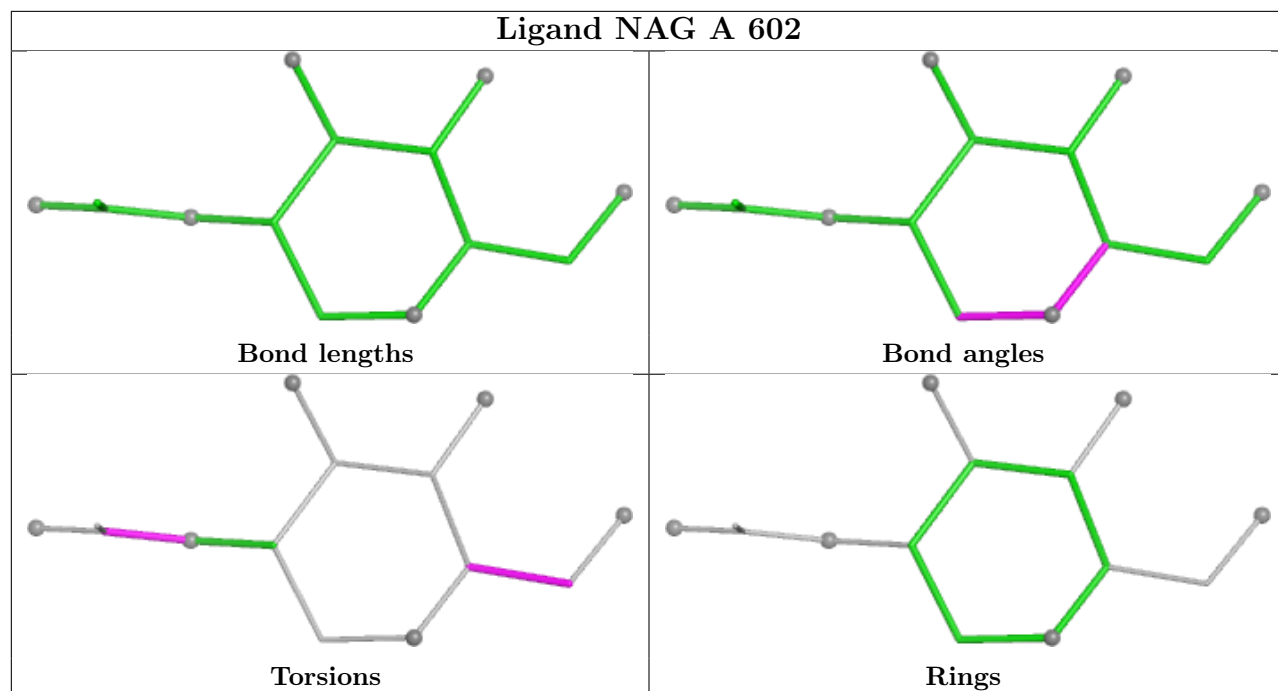


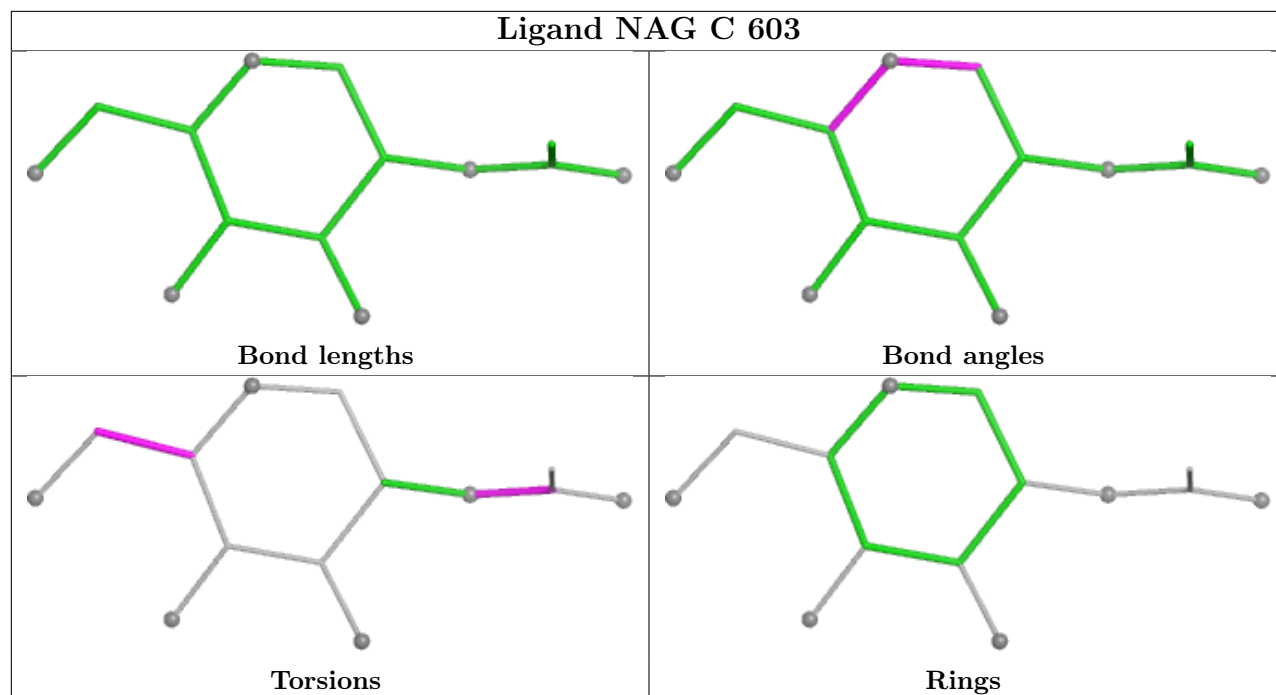












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

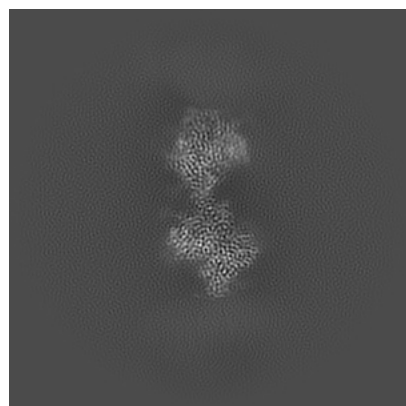
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-55773. These allow visual inspection of the internal detail of the map and identification of artifacts.

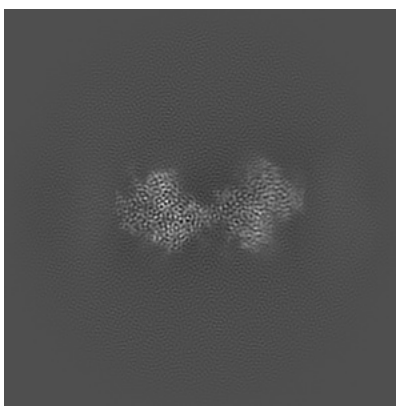
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

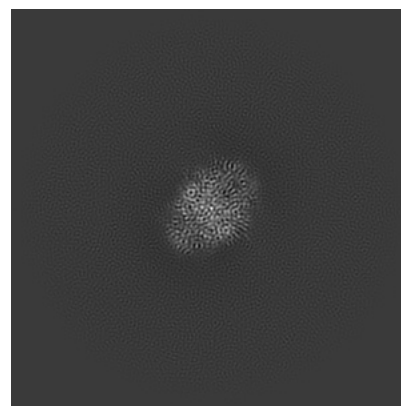
6.1.1 Primary map



X

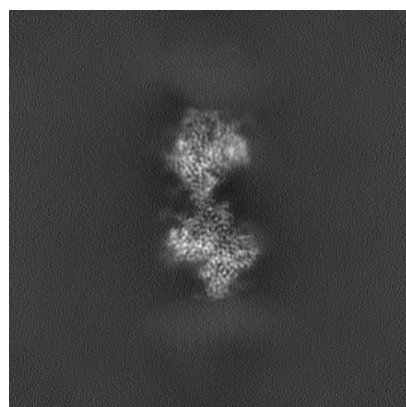


Y

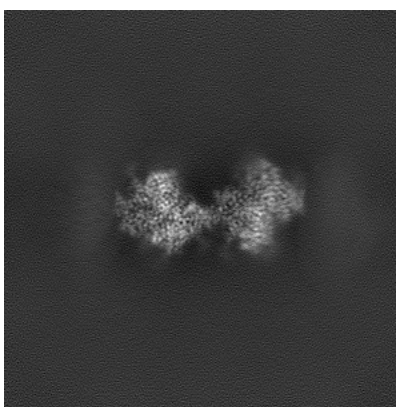


Z

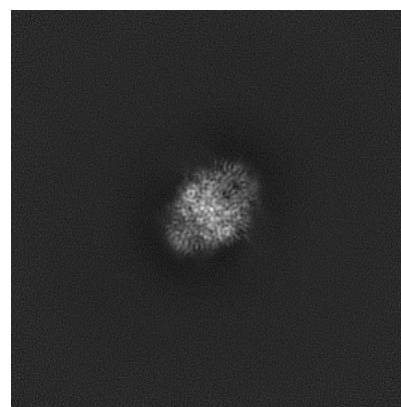
6.1.2 Raw map



X



Y

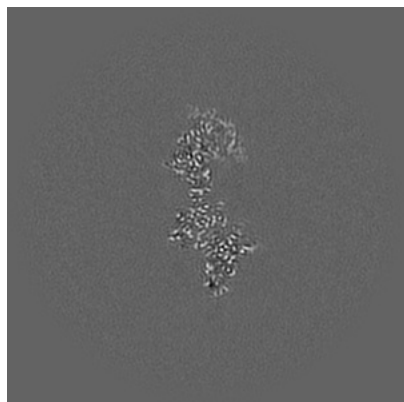


Z

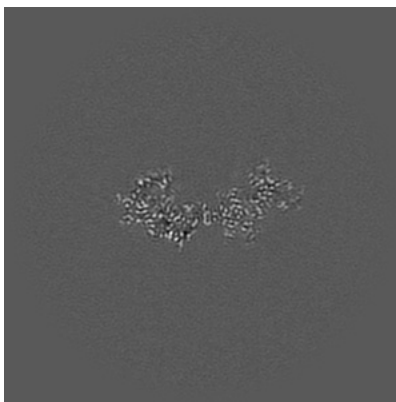
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

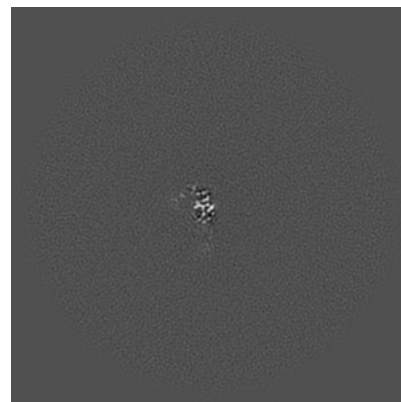
6.2.1 Primary map



X Index: 192

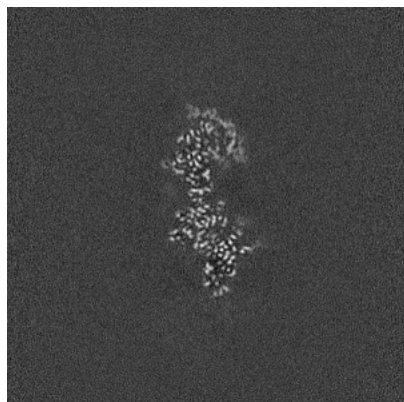


Y Index: 192

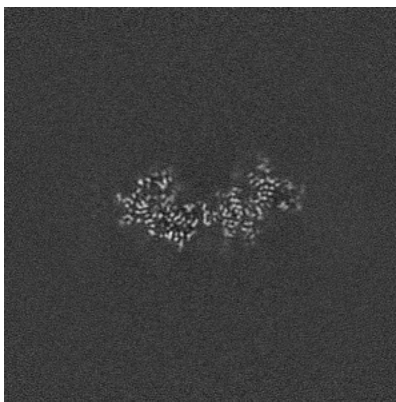


Z Index: 192

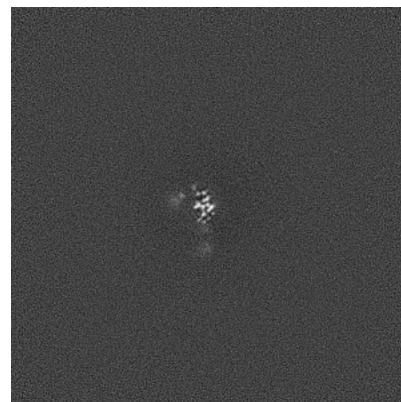
6.2.2 Raw map



X Index: 192



Y Index: 192

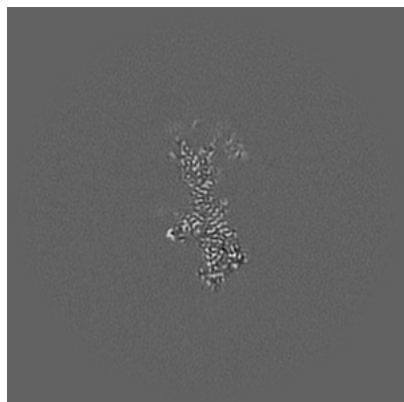


Z Index: 192

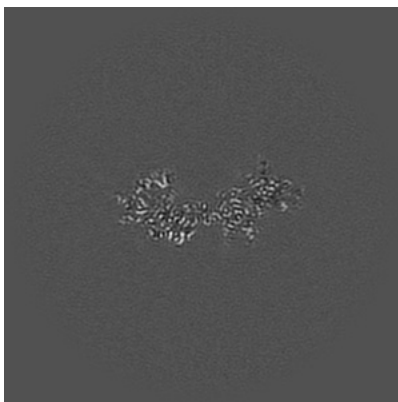
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

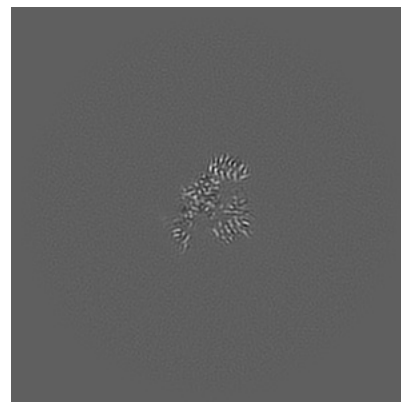
6.3.1 Primary map



X Index: 183

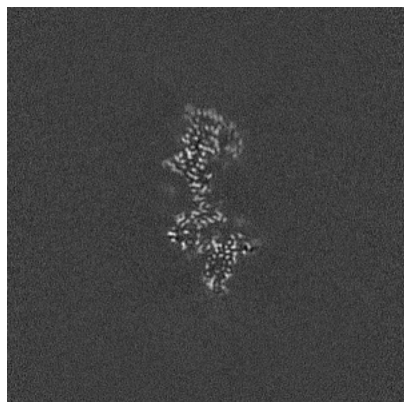


Y Index: 191

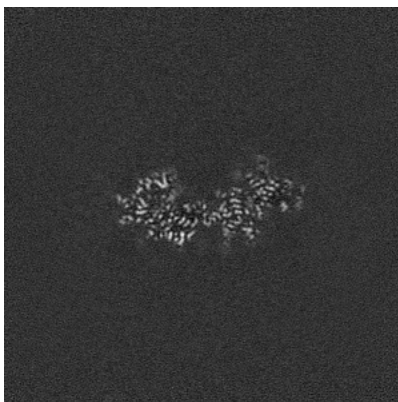


Z Index: 152

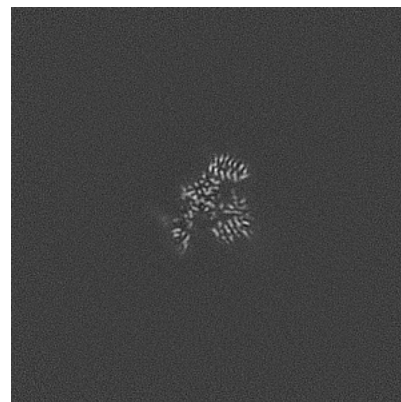
6.3.2 Raw map



X Index: 194



Y Index: 191

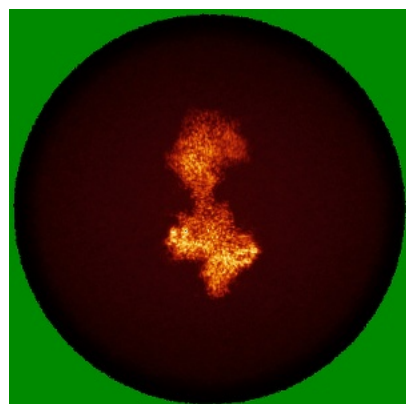


Z Index: 152

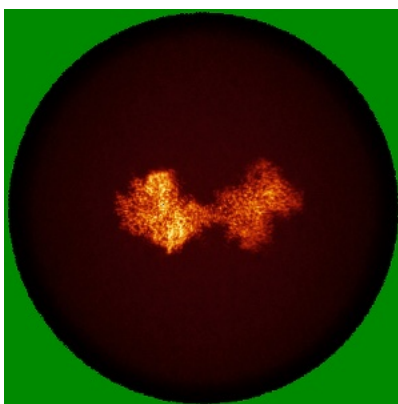
The images above show the largest variance slices of the map in three orthogonal directions.

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

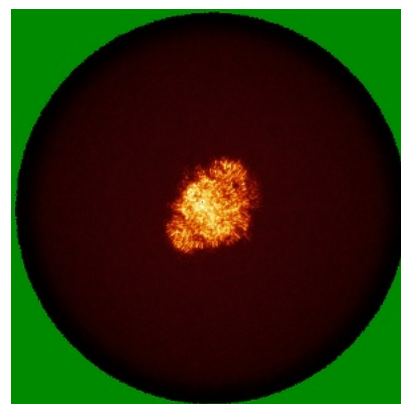
6.4.1 Primary map



X

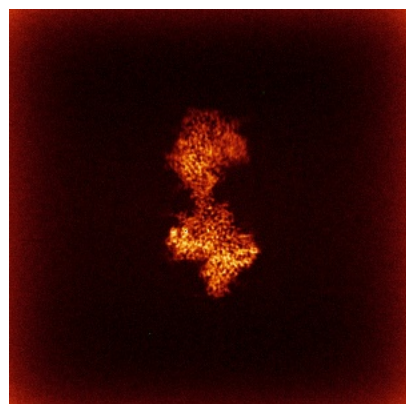


Y

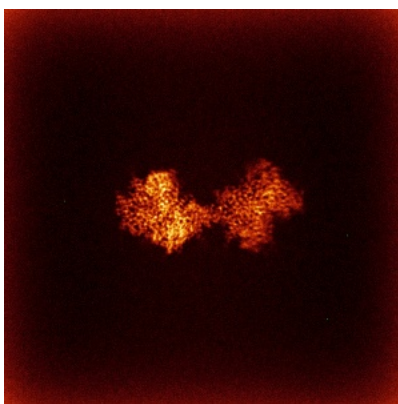


Z

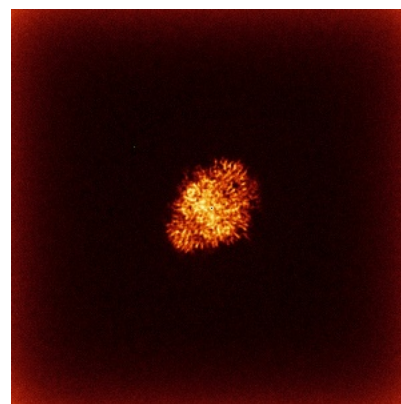
6.4.2 Raw map



X



Y

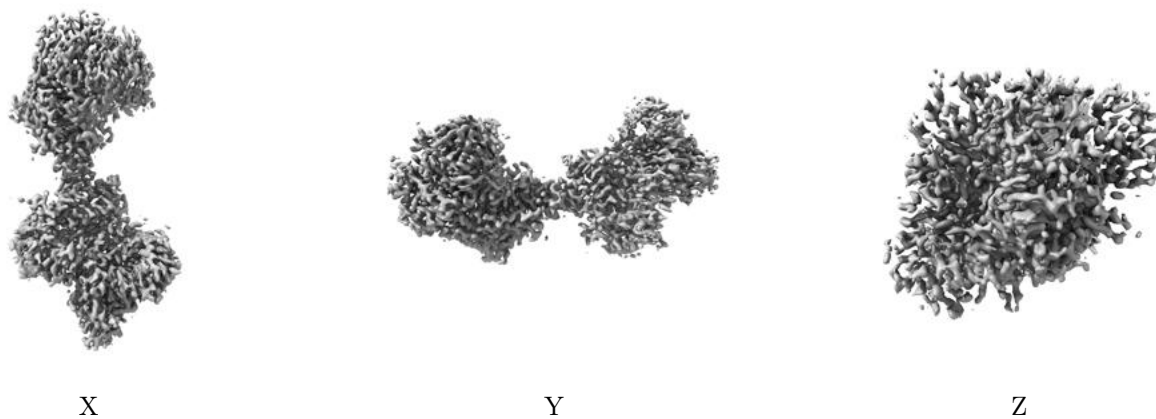


Z

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.

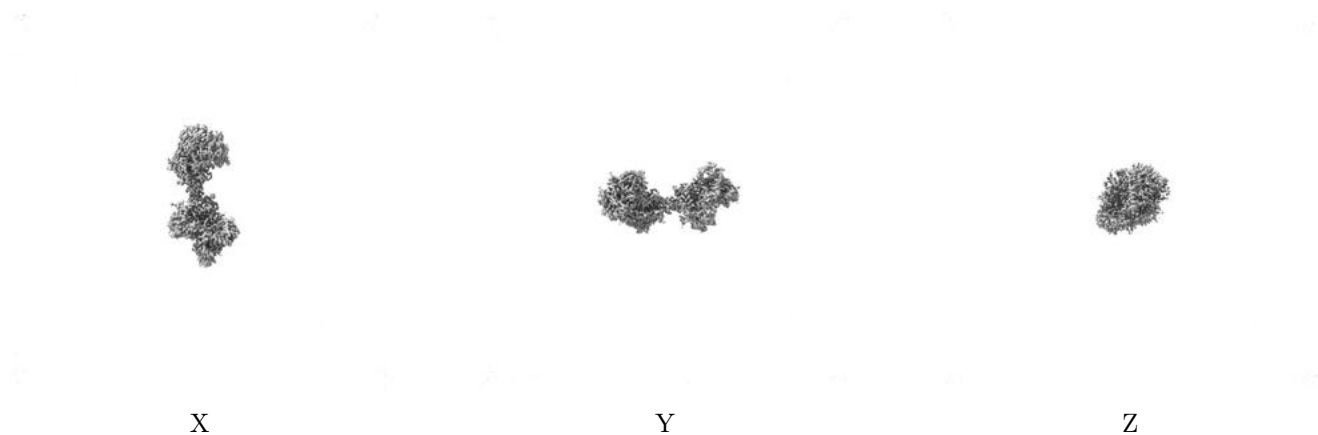
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.3. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

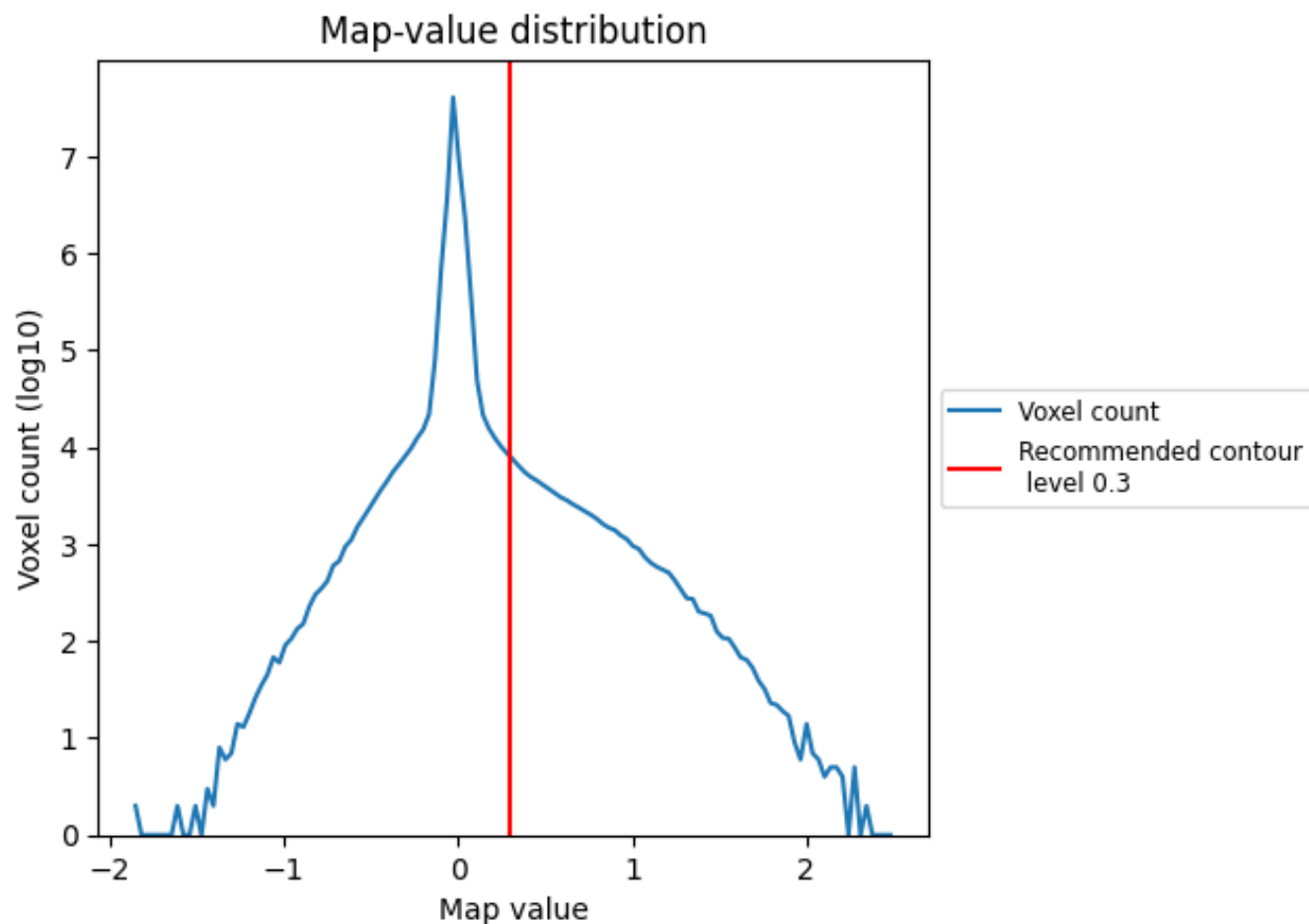
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

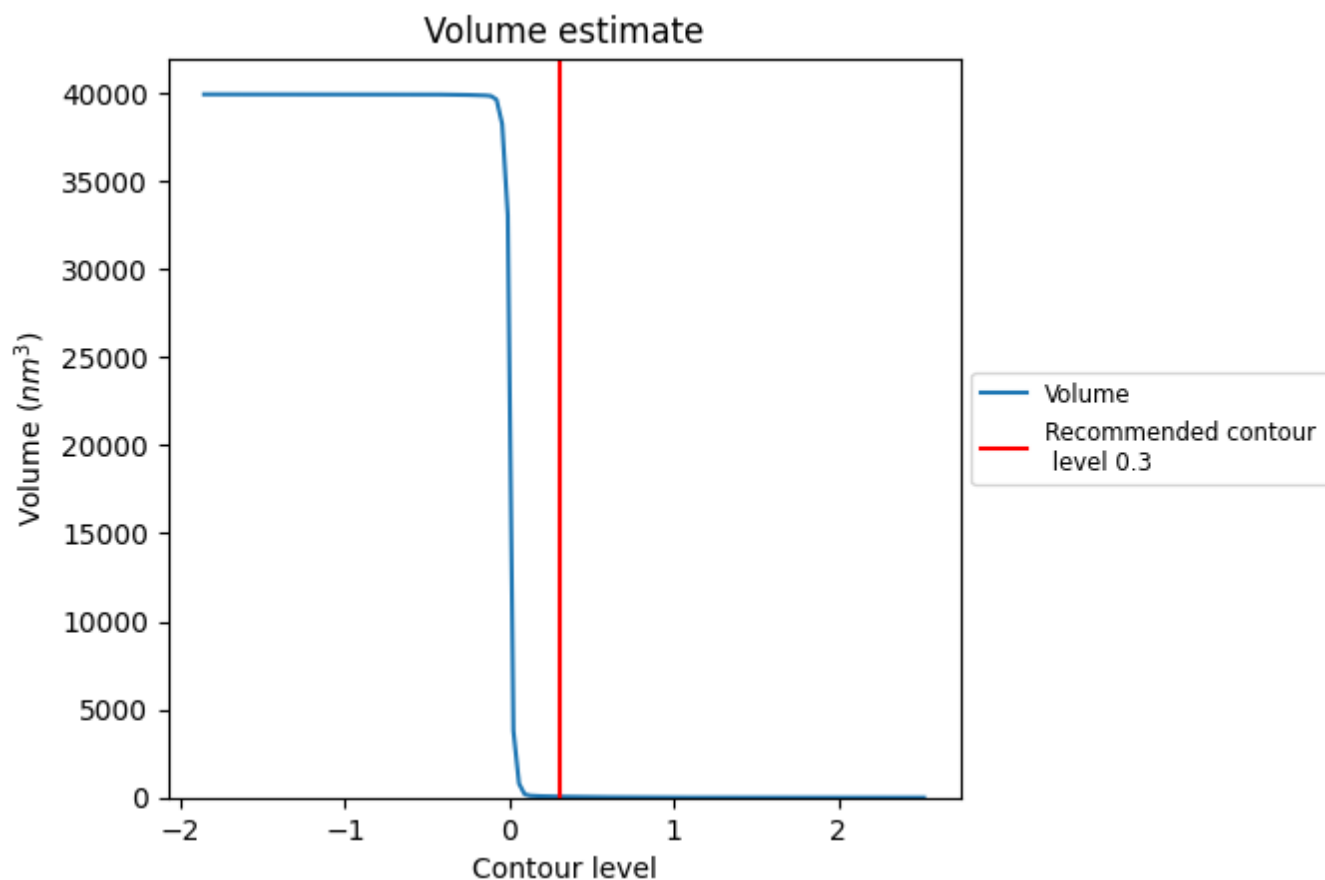
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

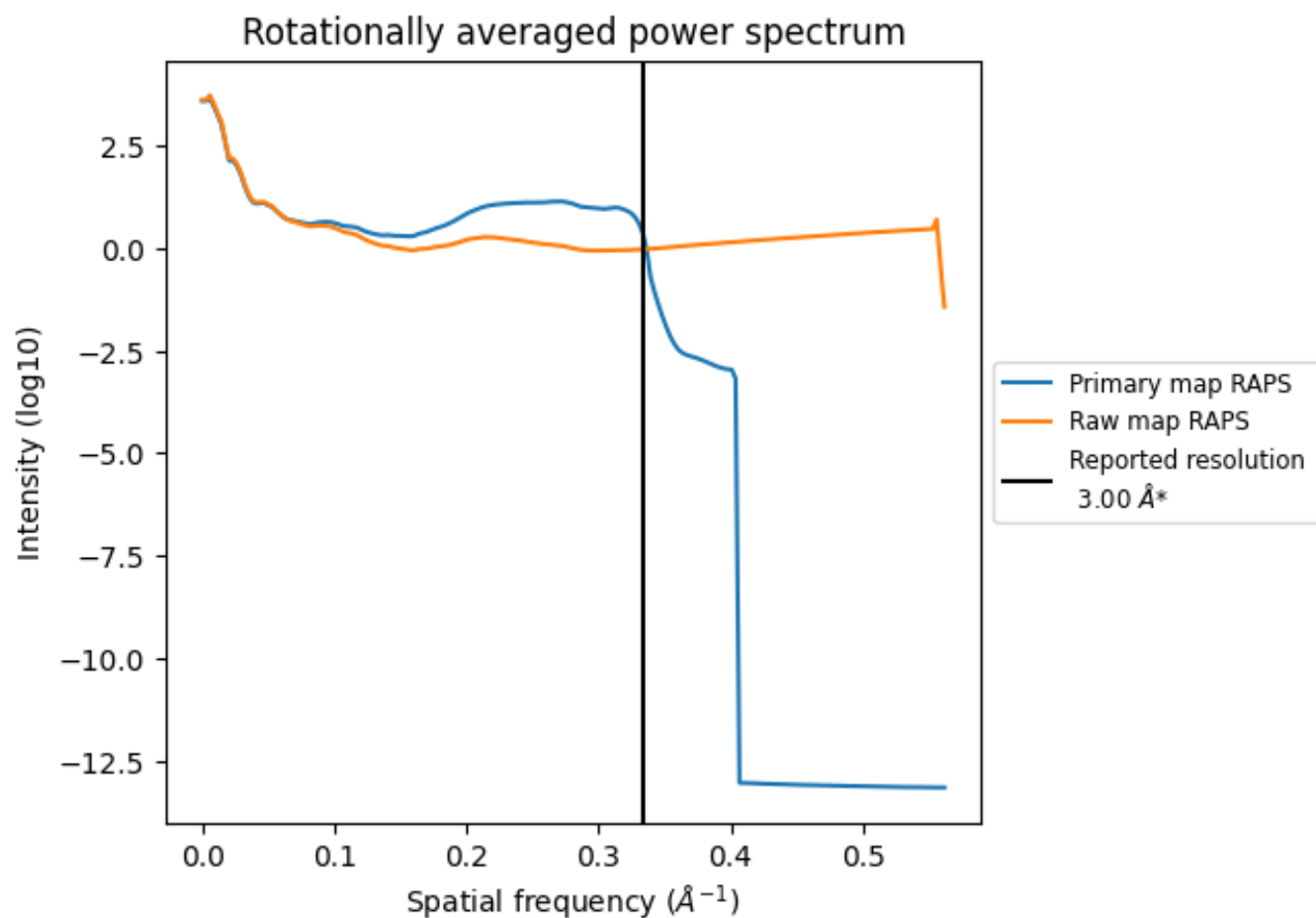
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 53 nm³; this corresponds to an approximate mass of 48 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

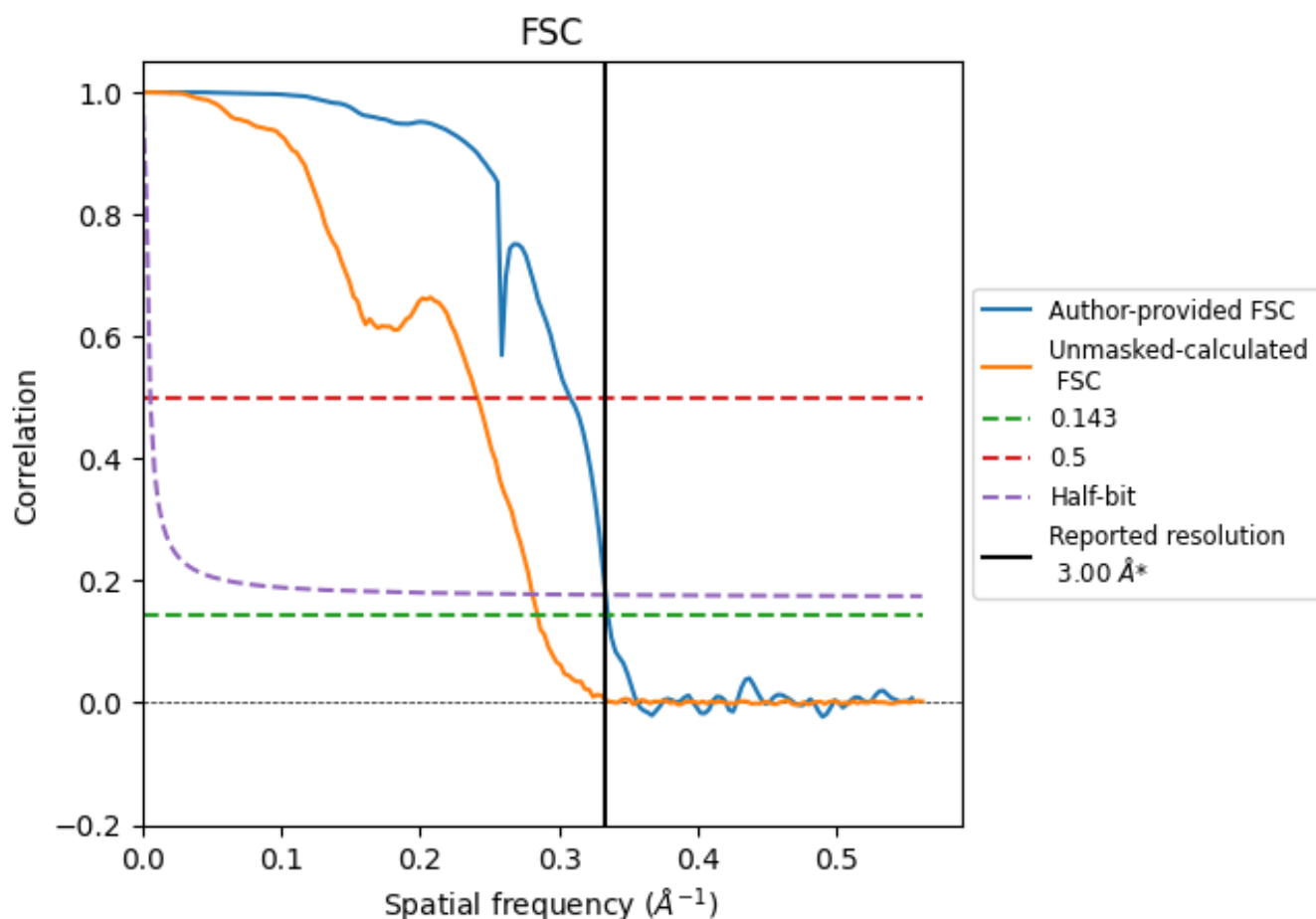


*Reported resolution corresponds to spatial frequency of 0.333 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.333 \AA^{-1}

8.2 Resolution estimates [i](#)

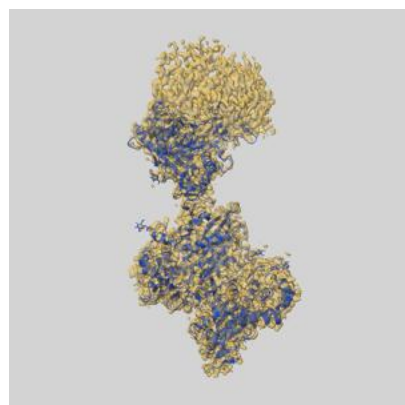
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	2.98	3.25	3.00
Unmasked-calculated*	3.51	4.14	3.56

*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 3.51 differs from the reported value 3.0 by more than 10 %

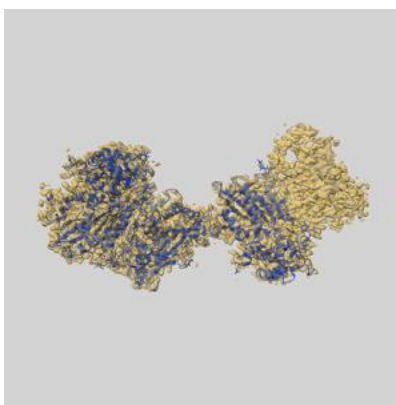
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-55773 and PDB model 9TBJ. Per-residue inclusion information can be found in section [3](#) on page [13](#).

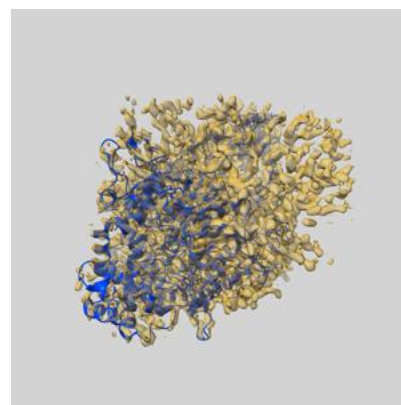
9.1 Map-model overlay [i](#)



X



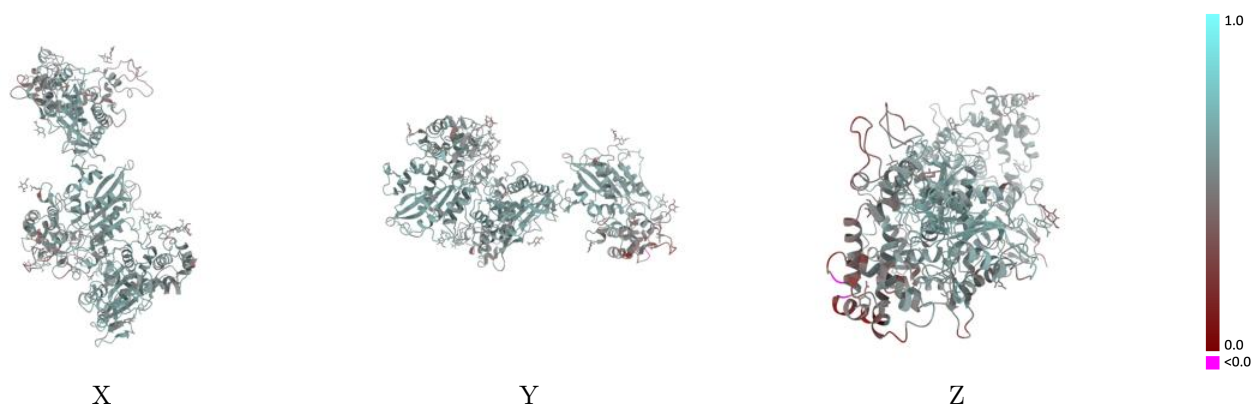
Y



Z

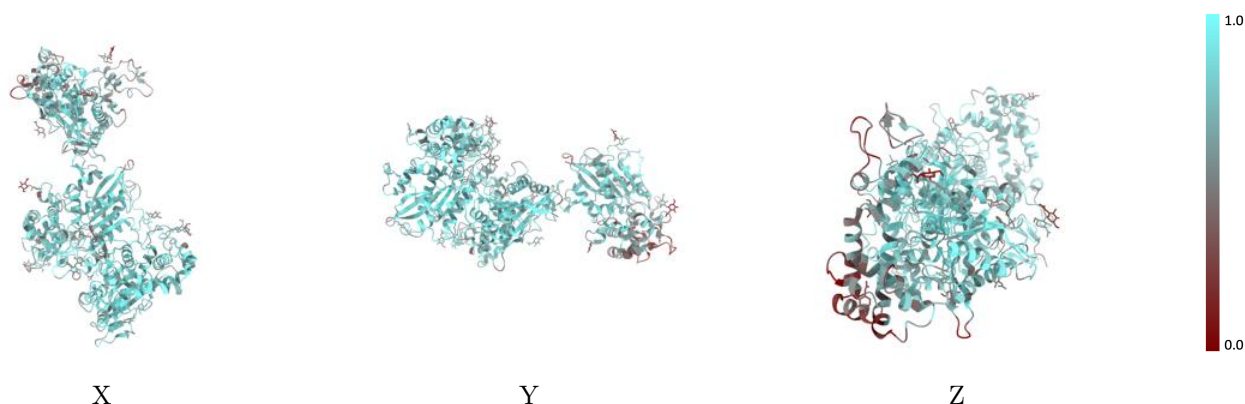
The images above show the 3D surface view of the map at the recommended contour level 0.3 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



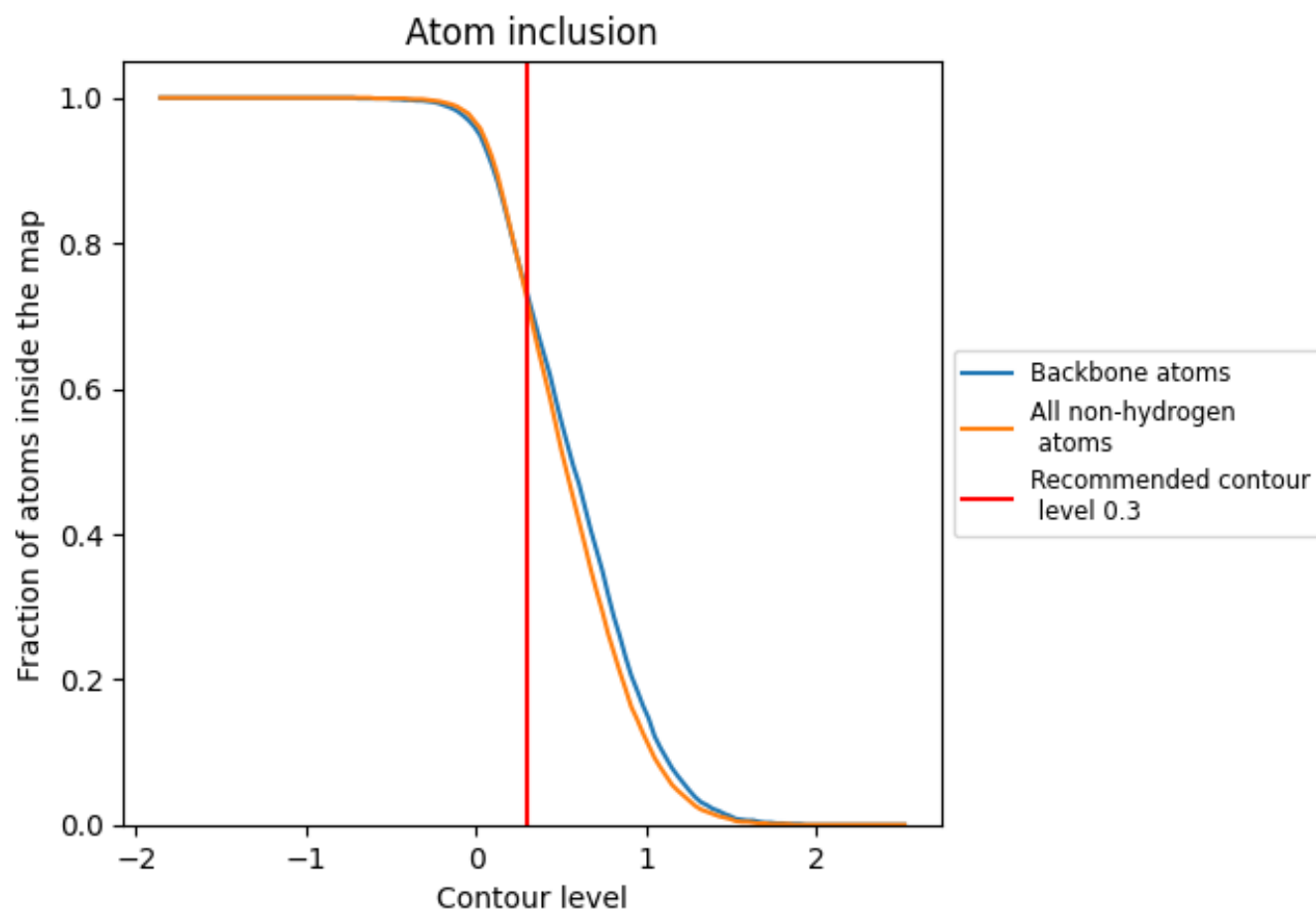
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.3).

9.4 Atom inclusion [i](#)



At the recommended contour level, 73% of all backbone atoms, 72% 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 (0.3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.7230	<div></div> 0.5430
A	<div></div> 0.7780	<div></div> 0.5650
B	<div></div> 0.7770	<div></div> 0.5550
C	<div></div> 0.6460	<div></div> 0.5140
D	<div></div> 0.5380	<div></div> 0.4650
F	<div></div> 0.7500	<div></div> 0.5460
G	<div></div> 0.3930	<div></div> 0.4710
I	<div></div> 0.6150	<div></div> 0.5250
K	<div></div> 0.6430	<div></div> 0.5110
L	<div></div> 0.4290	<div></div> 0.5180
M	<div></div> 0.3210	<div></div> 0.4760
P	<div></div> 0.3930	<div></div> 0.4700
Q	<div></div> 0.2860	<div></div> 0.4750

1.0

0.0

<0.0