



Full wwPDB EM Validation Report ⓘ

Jul 19, 2026 – 12:38 PM EDT

PDB ID : 9Y9D / pdb_00009y9d
EMDB ID : EMD-72694
Title : Cryo-EM structure of wild-type human CNNM4 tetramer with Magnesium and MgATP in outward-facing state
Authors : Bai, Z.; Zhou, E.; Lu, W.; Du, J.
Deposited on : 2025-09-14
Resolution : 3.48 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

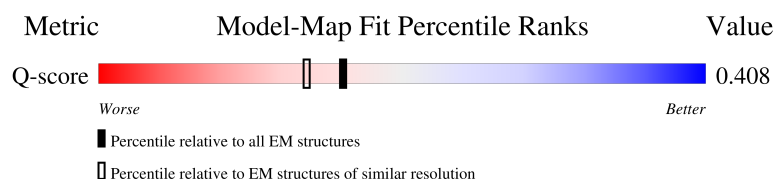
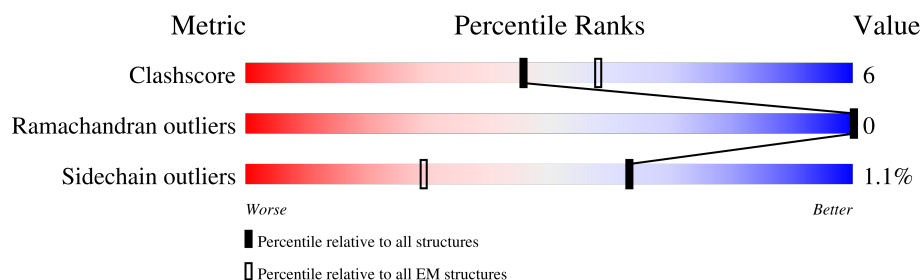
EMDB validation analysis : 0.0.1.dev133
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

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 3.48 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.



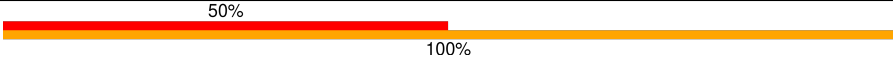
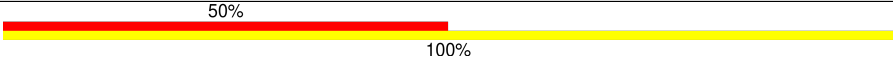
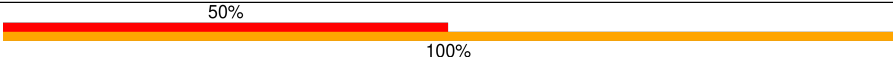
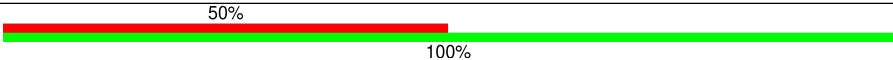
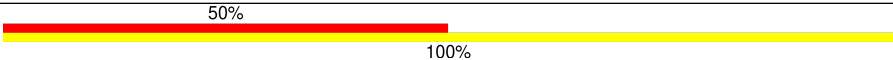
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13672 (2.98 - 3.98)

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	844	 10% 63% 12% 24%
1	B	844	 10% 63% 12% 25%
1	C	844	 10% 63% 12% 24%
1	D	844	 10% 63% 12% 25%

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Mol	Chain	Length	Quality of chain
2	E	2	
2	F	2	
2	G	2	
2	H	2	
2	I	2	
2	J	2	
2	K	2	
2	L	2	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 20658 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 Metal transporter CNNM4.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	639	Total	C	N	O	S	0	0
			5076	3259	838	951	28		
1	B	637	Total	C	N	O	S	0	0
			5040	3238	824	950	28		
1	C	639	Total	C	N	O	S	0	0
			5076	3259	838	951	28		
1	D	637	Total	C	N	O	S	0	0
			5040	3238	824	950	28		

There are 276 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	776	VAL	-	expression tag	UNP Q6P4Q7
A	777	ASP	-	expression tag	UNP Q6P4Q7
A	778	GLY	-	expression tag	UNP Q6P4Q7
A	779	SER	-	expression tag	UNP Q6P4Q7
A	780	LEU	-	expression tag	UNP Q6P4Q7
A	781	GLU	-	expression tag	UNP Q6P4Q7
A	782	VAL	-	expression tag	UNP Q6P4Q7
A	783	LEU	-	expression tag	UNP Q6P4Q7
A	784	PHE	-	expression tag	UNP Q6P4Q7
A	785	GLN	-	expression tag	UNP Q6P4Q7
A	786	GLY	-	expression tag	UNP Q6P4Q7
A	787	PRO	-	expression tag	UNP Q6P4Q7
A	788	SER	-	expression tag	UNP Q6P4Q7
A	789	GLY	-	expression tag	UNP Q6P4Q7
A	790	GLY	-	expression tag	UNP Q6P4Q7
A	791	SER	-	expression tag	UNP Q6P4Q7
A	792	ARG	-	expression tag	UNP Q6P4Q7
A	793	LEU	-	expression tag	UNP Q6P4Q7
A	794	GLU	-	expression tag	UNP Q6P4Q7
A	795	GLU	-	expression tag	UNP Q6P4Q7
A	796	GLU	-	expression tag	UNP Q6P4Q7
A	797	LEU	-	expression tag	UNP Q6P4Q7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	798	ARG	-	expression tag	UNP Q6P4Q7
A	799	ARG	-	expression tag	UNP Q6P4Q7
A	800	ARG	-	expression tag	UNP Q6P4Q7
A	801	LEU	-	expression tag	UNP Q6P4Q7
A	802	THR	-	expression tag	UNP Q6P4Q7
A	803	GLU	-	expression tag	UNP Q6P4Q7
A	804	GLY	-	expression tag	UNP Q6P4Q7
A	805	GLY	-	expression tag	UNP Q6P4Q7
A	806	SER	-	expression tag	UNP Q6P4Q7
A	807	TRP	-	expression tag	UNP Q6P4Q7
A	808	SER	-	expression tag	UNP Q6P4Q7
A	809	HIS	-	expression tag	UNP Q6P4Q7
A	810	PRO	-	expression tag	UNP Q6P4Q7
A	811	GLN	-	expression tag	UNP Q6P4Q7
A	812	PHE	-	expression tag	UNP Q6P4Q7
A	813	GLU	-	expression tag	UNP Q6P4Q7
A	814	LYS	-	expression tag	UNP Q6P4Q7
A	815	GLY	-	expression tag	UNP Q6P4Q7
A	816	GLY	-	expression tag	UNP Q6P4Q7
A	817	GLY	-	expression tag	UNP Q6P4Q7
A	818	SER	-	expression tag	UNP Q6P4Q7
A	819	GLY	-	expression tag	UNP Q6P4Q7
A	820	GLY	-	expression tag	UNP Q6P4Q7
A	821	GLY	-	expression tag	UNP Q6P4Q7
A	822	SER	-	expression tag	UNP Q6P4Q7
A	823	GLY	-	expression tag	UNP Q6P4Q7
A	824	GLY	-	expression tag	UNP Q6P4Q7
A	825	SER	-	expression tag	UNP Q6P4Q7
A	826	ALA	-	expression tag	UNP Q6P4Q7
A	827	TRP	-	expression tag	UNP Q6P4Q7
A	828	SER	-	expression tag	UNP Q6P4Q7
A	829	HIS	-	expression tag	UNP Q6P4Q7
A	830	PRO	-	expression tag	UNP Q6P4Q7
A	831	GLN	-	expression tag	UNP Q6P4Q7
A	832	PHE	-	expression tag	UNP Q6P4Q7
A	833	GLU	-	expression tag	UNP Q6P4Q7
A	834	LYS	-	expression tag	UNP Q6P4Q7
A	835	GLY	-	expression tag	UNP Q6P4Q7
A	836	SER	-	expression tag	UNP Q6P4Q7
A	837	HIS	-	expression tag	UNP Q6P4Q7
A	838	HIS	-	expression tag	UNP Q6P4Q7
A	839	HIS	-	expression tag	UNP Q6P4Q7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	840	HIS	-	expression tag	UNP Q6P4Q7
A	841	HIS	-	expression tag	UNP Q6P4Q7
A	842	HIS	-	expression tag	UNP Q6P4Q7
A	843	HIS	-	expression tag	UNP Q6P4Q7
A	844	HIS	-	expression tag	UNP Q6P4Q7
B	776	VAL	-	expression tag	UNP Q6P4Q7
B	777	ASP	-	expression tag	UNP Q6P4Q7
B	778	GLY	-	expression tag	UNP Q6P4Q7
B	779	SER	-	expression tag	UNP Q6P4Q7
B	780	LEU	-	expression tag	UNP Q6P4Q7
B	781	GLU	-	expression tag	UNP Q6P4Q7
B	782	VAL	-	expression tag	UNP Q6P4Q7
B	783	LEU	-	expression tag	UNP Q6P4Q7
B	784	PHE	-	expression tag	UNP Q6P4Q7
B	785	GLN	-	expression tag	UNP Q6P4Q7
B	786	GLY	-	expression tag	UNP Q6P4Q7
B	787	PRO	-	expression tag	UNP Q6P4Q7
B	788	SER	-	expression tag	UNP Q6P4Q7
B	789	GLY	-	expression tag	UNP Q6P4Q7
B	790	GLY	-	expression tag	UNP Q6P4Q7
B	791	SER	-	expression tag	UNP Q6P4Q7
B	792	ARG	-	expression tag	UNP Q6P4Q7
B	793	LEU	-	expression tag	UNP Q6P4Q7
B	794	GLU	-	expression tag	UNP Q6P4Q7
B	795	GLU	-	expression tag	UNP Q6P4Q7
B	796	GLU	-	expression tag	UNP Q6P4Q7
B	797	LEU	-	expression tag	UNP Q6P4Q7
B	798	ARG	-	expression tag	UNP Q6P4Q7
B	799	ARG	-	expression tag	UNP Q6P4Q7
B	800	ARG	-	expression tag	UNP Q6P4Q7
B	801	LEU	-	expression tag	UNP Q6P4Q7
B	802	THR	-	expression tag	UNP Q6P4Q7
B	803	GLU	-	expression tag	UNP Q6P4Q7
B	804	GLY	-	expression tag	UNP Q6P4Q7
B	805	GLY	-	expression tag	UNP Q6P4Q7
B	806	SER	-	expression tag	UNP Q6P4Q7
B	807	TRP	-	expression tag	UNP Q6P4Q7
B	808	SER	-	expression tag	UNP Q6P4Q7
B	809	HIS	-	expression tag	UNP Q6P4Q7
B	810	PRO	-	expression tag	UNP Q6P4Q7
B	811	GLN	-	expression tag	UNP Q6P4Q7
B	812	PHE	-	expression tag	UNP Q6P4Q7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	813	GLU	-	expression tag	UNP Q6P4Q7
B	814	LYS	-	expression tag	UNP Q6P4Q7
B	815	GLY	-	expression tag	UNP Q6P4Q7
B	816	GLY	-	expression tag	UNP Q6P4Q7
B	817	GLY	-	expression tag	UNP Q6P4Q7
B	818	SER	-	expression tag	UNP Q6P4Q7
B	819	GLY	-	expression tag	UNP Q6P4Q7
B	820	GLY	-	expression tag	UNP Q6P4Q7
B	821	GLY	-	expression tag	UNP Q6P4Q7
B	822	SER	-	expression tag	UNP Q6P4Q7
B	823	GLY	-	expression tag	UNP Q6P4Q7
B	824	GLY	-	expression tag	UNP Q6P4Q7
B	825	SER	-	expression tag	UNP Q6P4Q7
B	826	ALA	-	expression tag	UNP Q6P4Q7
B	827	TRP	-	expression tag	UNP Q6P4Q7
B	828	SER	-	expression tag	UNP Q6P4Q7
B	829	HIS	-	expression tag	UNP Q6P4Q7
B	830	PRO	-	expression tag	UNP Q6P4Q7
B	831	GLN	-	expression tag	UNP Q6P4Q7
B	832	PHE	-	expression tag	UNP Q6P4Q7
B	833	GLU	-	expression tag	UNP Q6P4Q7
B	834	LYS	-	expression tag	UNP Q6P4Q7
B	835	GLY	-	expression tag	UNP Q6P4Q7
B	836	SER	-	expression tag	UNP Q6P4Q7
B	837	HIS	-	expression tag	UNP Q6P4Q7
B	838	HIS	-	expression tag	UNP Q6P4Q7
B	839	HIS	-	expression tag	UNP Q6P4Q7
B	840	HIS	-	expression tag	UNP Q6P4Q7
B	841	HIS	-	expression tag	UNP Q6P4Q7
B	842	HIS	-	expression tag	UNP Q6P4Q7
B	843	HIS	-	expression tag	UNP Q6P4Q7
B	844	HIS	-	expression tag	UNP Q6P4Q7
C	776	VAL	-	expression tag	UNP Q6P4Q7
C	777	ASP	-	expression tag	UNP Q6P4Q7
C	778	GLY	-	expression tag	UNP Q6P4Q7
C	779	SER	-	expression tag	UNP Q6P4Q7
C	780	LEU	-	expression tag	UNP Q6P4Q7
C	781	GLU	-	expression tag	UNP Q6P4Q7
C	782	VAL	-	expression tag	UNP Q6P4Q7
C	783	LEU	-	expression tag	UNP Q6P4Q7
C	784	PHE	-	expression tag	UNP Q6P4Q7
C	785	GLN	-	expression tag	UNP Q6P4Q7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	786	GLY	-	expression tag	UNP Q6P4Q7
C	787	PRO	-	expression tag	UNP Q6P4Q7
C	788	SER	-	expression tag	UNP Q6P4Q7
C	789	GLY	-	expression tag	UNP Q6P4Q7
C	790	GLY	-	expression tag	UNP Q6P4Q7
C	791	SER	-	expression tag	UNP Q6P4Q7
C	792	ARG	-	expression tag	UNP Q6P4Q7
C	793	LEU	-	expression tag	UNP Q6P4Q7
C	794	GLU	-	expression tag	UNP Q6P4Q7
C	795	GLU	-	expression tag	UNP Q6P4Q7
C	796	GLU	-	expression tag	UNP Q6P4Q7
C	797	LEU	-	expression tag	UNP Q6P4Q7
C	798	ARG	-	expression tag	UNP Q6P4Q7
C	799	ARG	-	expression tag	UNP Q6P4Q7
C	800	ARG	-	expression tag	UNP Q6P4Q7
C	801	LEU	-	expression tag	UNP Q6P4Q7
C	802	THR	-	expression tag	UNP Q6P4Q7
C	803	GLU	-	expression tag	UNP Q6P4Q7
C	804	GLY	-	expression tag	UNP Q6P4Q7
C	805	GLY	-	expression tag	UNP Q6P4Q7
C	806	SER	-	expression tag	UNP Q6P4Q7
C	807	TRP	-	expression tag	UNP Q6P4Q7
C	808	SER	-	expression tag	UNP Q6P4Q7
C	809	HIS	-	expression tag	UNP Q6P4Q7
C	810	PRO	-	expression tag	UNP Q6P4Q7
C	811	GLN	-	expression tag	UNP Q6P4Q7
C	812	PHE	-	expression tag	UNP Q6P4Q7
C	813	GLU	-	expression tag	UNP Q6P4Q7
C	814	LYS	-	expression tag	UNP Q6P4Q7
C	815	GLY	-	expression tag	UNP Q6P4Q7
C	816	GLY	-	expression tag	UNP Q6P4Q7
C	817	GLY	-	expression tag	UNP Q6P4Q7
C	818	SER	-	expression tag	UNP Q6P4Q7
C	819	GLY	-	expression tag	UNP Q6P4Q7
C	820	GLY	-	expression tag	UNP Q6P4Q7
C	821	GLY	-	expression tag	UNP Q6P4Q7
C	822	SER	-	expression tag	UNP Q6P4Q7
C	823	GLY	-	expression tag	UNP Q6P4Q7
C	824	GLY	-	expression tag	UNP Q6P4Q7
C	825	SER	-	expression tag	UNP Q6P4Q7
C	826	ALA	-	expression tag	UNP Q6P4Q7
C	827	TRP	-	expression tag	UNP Q6P4Q7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	828	SER	-	expression tag	UNP Q6P4Q7
C	829	HIS	-	expression tag	UNP Q6P4Q7
C	830	PRO	-	expression tag	UNP Q6P4Q7
C	831	GLN	-	expression tag	UNP Q6P4Q7
C	832	PHE	-	expression tag	UNP Q6P4Q7
C	833	GLU	-	expression tag	UNP Q6P4Q7
C	834	LYS	-	expression tag	UNP Q6P4Q7
C	835	GLY	-	expression tag	UNP Q6P4Q7
C	836	SER	-	expression tag	UNP Q6P4Q7
C	837	HIS	-	expression tag	UNP Q6P4Q7
C	838	HIS	-	expression tag	UNP Q6P4Q7
C	839	HIS	-	expression tag	UNP Q6P4Q7
C	840	HIS	-	expression tag	UNP Q6P4Q7
C	841	HIS	-	expression tag	UNP Q6P4Q7
C	842	HIS	-	expression tag	UNP Q6P4Q7
C	843	HIS	-	expression tag	UNP Q6P4Q7
C	844	HIS	-	expression tag	UNP Q6P4Q7
D	776	VAL	-	expression tag	UNP Q6P4Q7
D	777	ASP	-	expression tag	UNP Q6P4Q7
D	778	GLY	-	expression tag	UNP Q6P4Q7
D	779	SER	-	expression tag	UNP Q6P4Q7
D	780	LEU	-	expression tag	UNP Q6P4Q7
D	781	GLU	-	expression tag	UNP Q6P4Q7
D	782	VAL	-	expression tag	UNP Q6P4Q7
D	783	LEU	-	expression tag	UNP Q6P4Q7
D	784	PHE	-	expression tag	UNP Q6P4Q7
D	785	GLN	-	expression tag	UNP Q6P4Q7
D	786	GLY	-	expression tag	UNP Q6P4Q7
D	787	PRO	-	expression tag	UNP Q6P4Q7
D	788	SER	-	expression tag	UNP Q6P4Q7
D	789	GLY	-	expression tag	UNP Q6P4Q7
D	790	GLY	-	expression tag	UNP Q6P4Q7
D	791	SER	-	expression tag	UNP Q6P4Q7
D	792	ARG	-	expression tag	UNP Q6P4Q7
D	793	LEU	-	expression tag	UNP Q6P4Q7
D	794	GLU	-	expression tag	UNP Q6P4Q7
D	795	GLU	-	expression tag	UNP Q6P4Q7
D	796	GLU	-	expression tag	UNP Q6P4Q7
D	797	LEU	-	expression tag	UNP Q6P4Q7
D	798	ARG	-	expression tag	UNP Q6P4Q7
D	799	ARG	-	expression tag	UNP Q6P4Q7
D	800	ARG	-	expression tag	UNP Q6P4Q7

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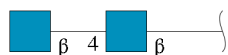
Chain	Residue	Modelled	Actual	Comment	Reference
D	801	LEU	-	expression tag	UNP Q6P4Q7
D	802	THR	-	expression tag	UNP Q6P4Q7
D	803	GLU	-	expression tag	UNP Q6P4Q7
D	804	GLY	-	expression tag	UNP Q6P4Q7
D	805	GLY	-	expression tag	UNP Q6P4Q7
D	806	SER	-	expression tag	UNP Q6P4Q7
D	807	TRP	-	expression tag	UNP Q6P4Q7
D	808	SER	-	expression tag	UNP Q6P4Q7
D	809	HIS	-	expression tag	UNP Q6P4Q7
D	810	PRO	-	expression tag	UNP Q6P4Q7
D	811	GLN	-	expression tag	UNP Q6P4Q7
D	812	PHE	-	expression tag	UNP Q6P4Q7
D	813	GLU	-	expression tag	UNP Q6P4Q7
D	814	LYS	-	expression tag	UNP Q6P4Q7
D	815	GLY	-	expression tag	UNP Q6P4Q7
D	816	GLY	-	expression tag	UNP Q6P4Q7
D	817	GLY	-	expression tag	UNP Q6P4Q7
D	818	SER	-	expression tag	UNP Q6P4Q7
D	819	GLY	-	expression tag	UNP Q6P4Q7
D	820	GLY	-	expression tag	UNP Q6P4Q7
D	821	GLY	-	expression tag	UNP Q6P4Q7
D	822	SER	-	expression tag	UNP Q6P4Q7
D	823	GLY	-	expression tag	UNP Q6P4Q7
D	824	GLY	-	expression tag	UNP Q6P4Q7
D	825	SER	-	expression tag	UNP Q6P4Q7
D	826	ALA	-	expression tag	UNP Q6P4Q7
D	827	TRP	-	expression tag	UNP Q6P4Q7
D	828	SER	-	expression tag	UNP Q6P4Q7
D	829	HIS	-	expression tag	UNP Q6P4Q7
D	830	PRO	-	expression tag	UNP Q6P4Q7
D	831	GLN	-	expression tag	UNP Q6P4Q7
D	832	PHE	-	expression tag	UNP Q6P4Q7
D	833	GLU	-	expression tag	UNP Q6P4Q7
D	834	LYS	-	expression tag	UNP Q6P4Q7
D	835	GLY	-	expression tag	UNP Q6P4Q7
D	836	SER	-	expression tag	UNP Q6P4Q7
D	837	HIS	-	expression tag	UNP Q6P4Q7
D	838	HIS	-	expression tag	UNP Q6P4Q7
D	839	HIS	-	expression tag	UNP Q6P4Q7
D	840	HIS	-	expression tag	UNP Q6P4Q7
D	841	HIS	-	expression tag	UNP Q6P4Q7
D	842	HIS	-	expression tag	UNP Q6P4Q7

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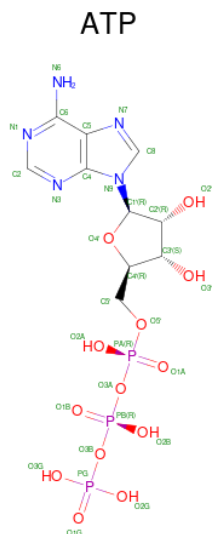
Chain	Residue	Modelled	Actual	Comment	Reference
D	843	HIS	-	expression tag	UNP Q6P4Q7
D	844	HIS	-	expression tag	UNP Q6P4Q7

- Molecule 2 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
2	E	2	Total	C	N	O	0	0
			28	16	2	10		
2	F	2	Total	C	N	O	0	0
			28	16	2	10		
2	G	2	Total	C	N	O	0	0
			28	16	2	10		
2	H	2	Total	C	N	O	0	0
			28	16	2	10		
2	I	2	Total	C	N	O	0	0
			28	16	2	10		
2	J	2	Total	C	N	O	0	0
			28	16	2	10		
2	K	2	Total	C	N	O	0	0
			28	16	2	10		
2	L	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
3	A	1	Total 31	C 10	N 5	O 13	P 3	0
3	B	1	Total 31	C 10	N 5	O 13	P 3	0
3	C	1	Total 31	C 10	N 5	O 13	P 3	0
3	D	1	Total 31	C 10	N 5	O 13	P 3	0

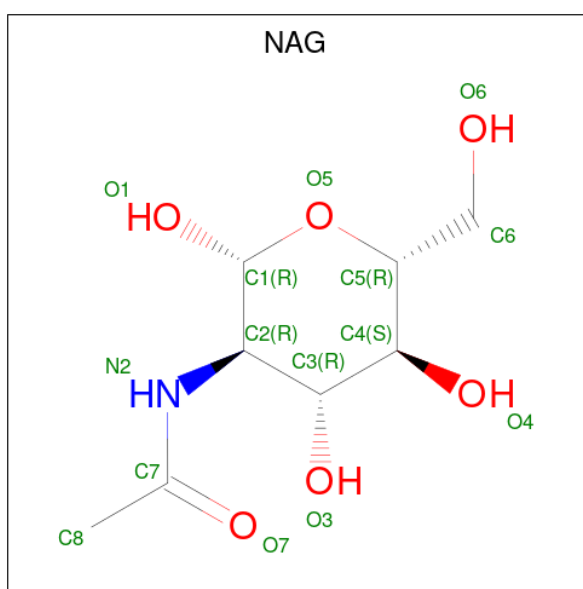
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	AltConf
4	A	4	Total Mg 4 4	0
4	B	4	Total Mg 4 4	0
4	C	4	Total Mg 4 4	0
4	D	4	Total Mg 4 4	0

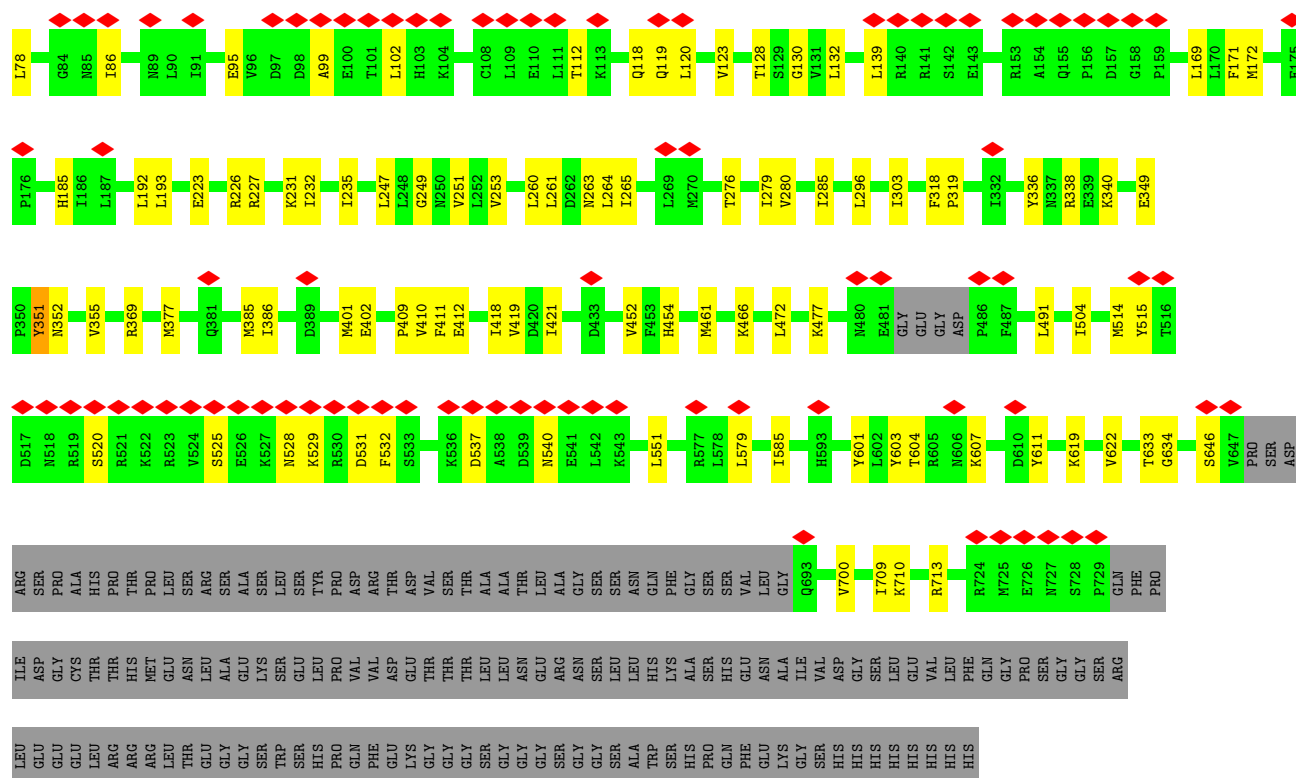
- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
5	A	2	Total	Na	0
			2	2	
5	B	1	Total	Na	0
			1	1	
5	C	2	Total	Na	0
			2	2	
5	D	1	Total	Na	0
			1	1	

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).

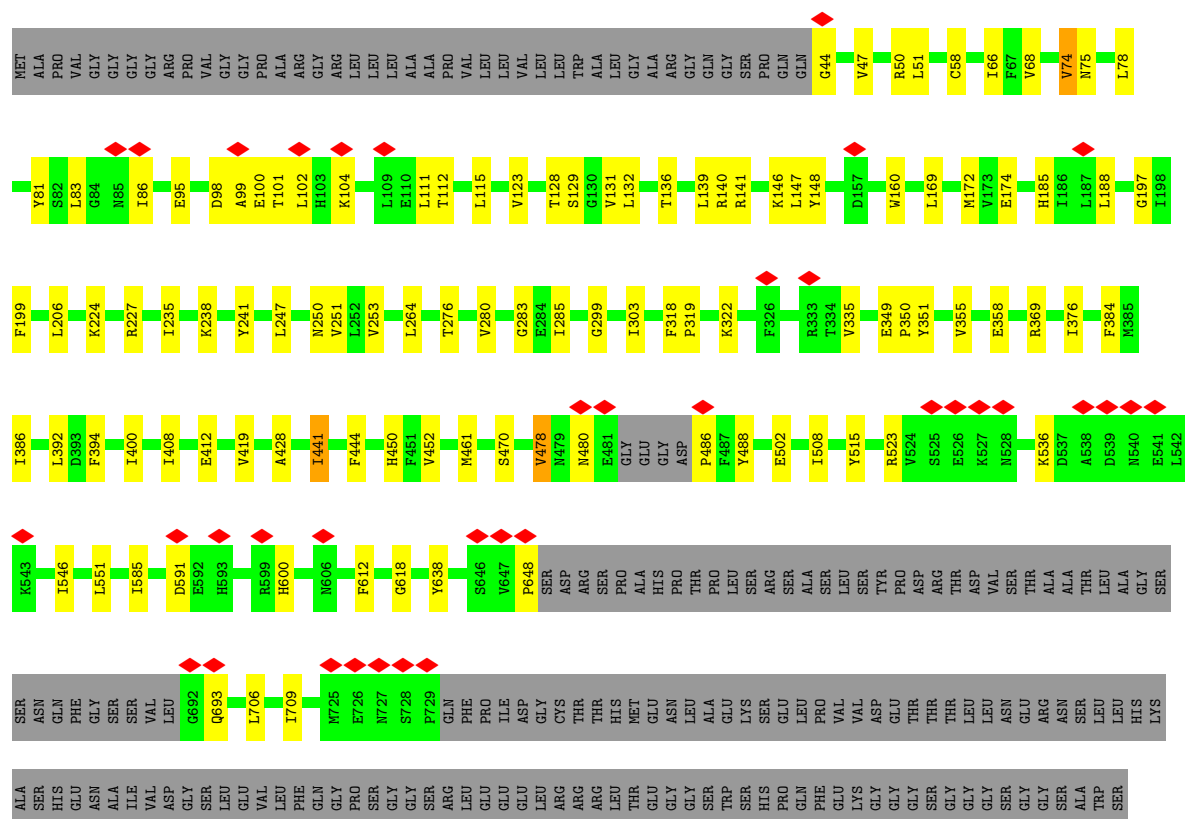


Mol	Chain	Residues	Atoms				AltConf
6	A	1	Total	C	N	O	0
			14	8	1	5	
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	

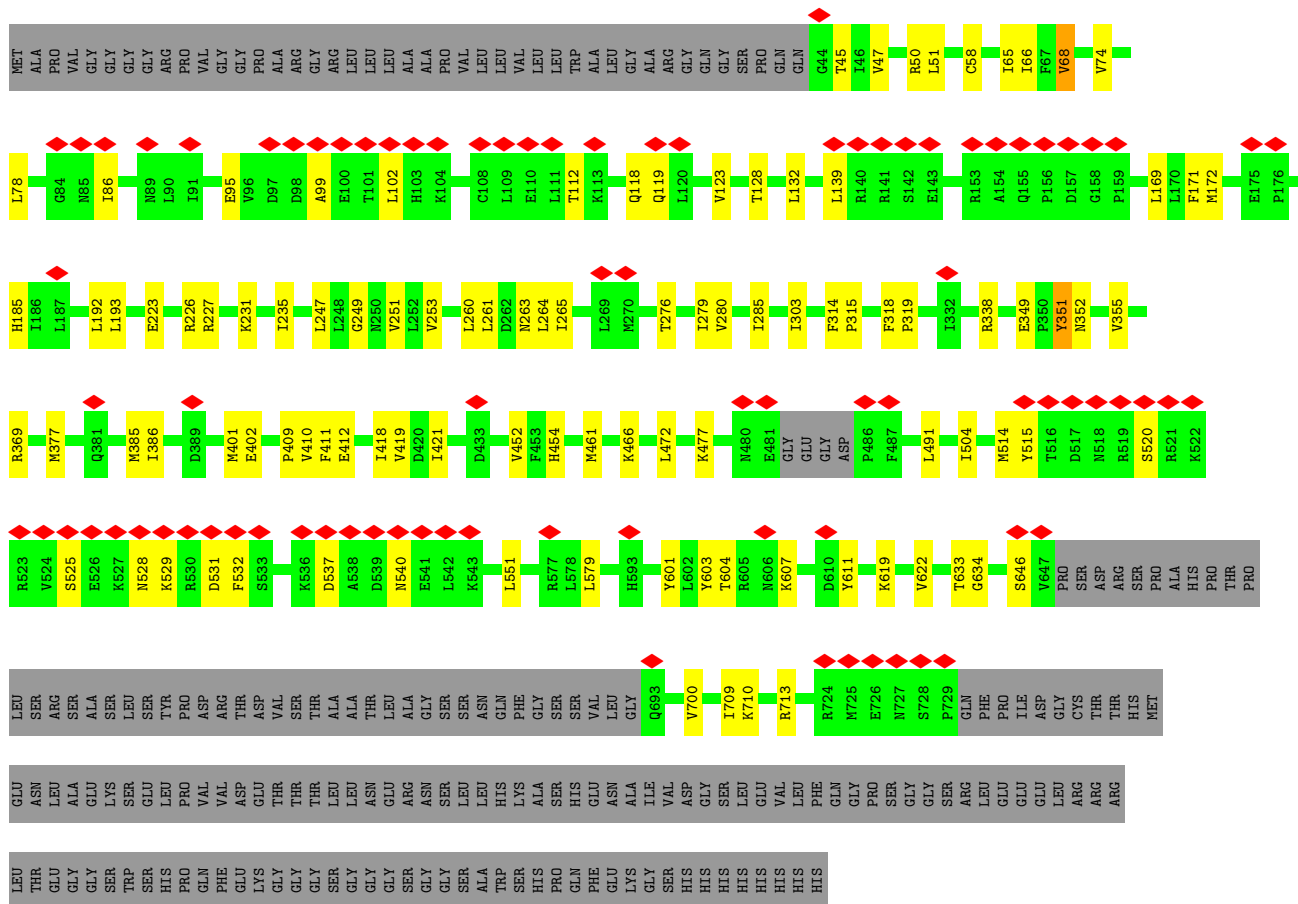


• Molecule 1: Metal transporter CNNM4

Chain C: 63% 12% 24%



- Molecule 1: Metal transporter CNNM4



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



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



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



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



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



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



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



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





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	946137	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 ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	1200	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	0.019	Depositor
Minimum map value	-0.004	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.000	Depositor
Recommended contour level	0.00376	Depositor
Map size (Å)	361.504, 361.504, 361.504	wwPDB
Map dimensions	416, 416, 416	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.869, 0.869, 0.869	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, ATP, NA, MG

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	0.24	0/5165	0.33	0/6984
1	B	0.17	0/5128	0.27	0/6939
1	C	0.24	0/5165	0.33	0/6984
1	D	0.17	0/5128	0.27	0/6939
All	All	0.21	0/20586	0.30	0/27846

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	5076	0	5157	74	0
1	B	5040	0	5096	69	0
1	C	5076	0	5157	82	0
1	D	5040	0	5096	67	0
2	E	28	0	25	0	0
2	F	28	0	25	1	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
2	I	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	J	28	0	25	1	0
2	K	28	0	25	0	0
2	L	28	0	25	0	0
3	A	31	0	12	0	0
3	B	31	0	12	0	0
3	C	31	0	12	0	0
3	D	31	0	12	0	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
4	C	4	0	0	0	0
4	D	4	0	0	0	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0
5	C	2	0	0	0	0
5	D	1	0	0	0	0
6	A	14	0	13	0	0
6	B	14	0	13	0	0
6	C	14	0	13	0	0
6	D	14	0	13	0	0
All	All	20658	0	20806	258	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 6.

All (258) 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:50:ARG:HG2	1:B:50:ARG:HG2	1.50	0.92
1:C:50:ARG:HG2	1:D:50:ARG:HG2	1.50	0.91
1:C:335:VAL:HG11	1:D:351:TYR:HB2	1.73	0.71
1:C:251:VAL:HG11	1:D:280:VAL:HG21	1.74	0.70
1:A:335:VAL:HG11	1:B:351:TYR:HB2	1.73	0.70
1:A:174:GLU:HG3	1:C:141:ARG:HH12	1.58	0.69
1:A:251:VAL:HG11	1:B:280:VAL:HG21	1.74	0.68
1:A:141:ARG:HH12	1:C:174:GLU:HG3	1.59	0.68
1:A:508:ILE:HG12	1:B:466:LYS:HG2	1.78	0.65
1:A:280:VAL:HG22	1:B:251:VAL:HG21	1.78	0.65
1:A:47:VAL:HG13	1:B:128:THR:HB	1.78	0.65
1:A:102:LEU:HD13	1:A:169:LEU:HD21	1.80	0.64
1:C:280:VAL:HG22	1:D:251:VAL:HG21	1.78	0.64
1:C:508:ILE:HG12	1:D:466:LYS:HG2	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:514:MET:HB3	1:D:525:SER:HB3	1.80	0.63
1:C:47:VAL:HG13	1:D:128:THR:HB	1.78	0.63
1:C:102:LEU:HD13	1:C:169:LEU:HD21	1.80	0.63
1:B:514:MET:HB3	1:B:525:SER:HB3	1.80	0.63
1:C:50:ARG:HD3	1:C:66:ILE:HD11	1.81	0.62
1:B:412:GLU:HB2	1:B:419:VAL:HG21	1.82	0.62
1:C:78:LEU:HD11	1:C:132:LEU:HD22	1.82	0.61
1:D:412:GLU:HB2	1:D:419:VAL:HG21	1.82	0.61
1:C:355:VAL:HG21	1:D:369:ARG:HH21	1.66	0.61
1:A:355:VAL:HG21	1:B:369:ARG:HH21	1.66	0.61
1:D:452:VAL:HG21	1:D:461:MET:HE1	1.83	0.60
1:A:50:ARG:HD3	1:A:66:ILE:HD11	1.81	0.60
1:A:78:LEU:HD11	1:A:132:LEU:HD22	1.83	0.60
1:C:44:GLY:HA2	1:C:81:TYR:O	2.01	0.60
1:B:452:VAL:HG21	1:B:461:MET:HE1	1.83	0.60
1:A:44:GLY:HA2	1:A:81:TYR:O	2.01	0.59
1:B:338:ARG:HH12	1:B:369:ARG:HG2	1.67	0.58
1:D:338:ARG:HH12	1:D:369:ARG:HG2	1.67	0.58
1:C:206:LEU:HD21	1:C:299:GLY:HA3	1.86	0.58
1:D:477:LYS:HE3	1:D:491:LEU:HD11	1.85	0.58
1:B:477:LYS:HE3	1:B:491:LEU:HD11	1.86	0.58
1:B:622:VAL:HG22	1:B:700:VAL:HG22	1.87	0.57
1:A:206:LEU:HD21	1:A:299:GLY:HA3	1.86	0.57
1:B:276:THR:O	1:B:280:VAL:HG23	2.05	0.57
1:B:185:HIS:HD2	1:B:264:LEU:HD13	1.71	0.56
1:D:622:VAL:HG22	1:D:700:VAL:HG22	1.87	0.56
1:A:115:LEU:HD23	1:A:136:THR:HG22	1.88	0.56
1:A:235:ILE:HG13	1:A:303:ILE:HD11	1.88	0.56
1:D:276:THR:O	1:D:280:VAL:HG23	2.05	0.56
1:A:141:ARG:HH21	1:C:139:LEU:HB2	1.71	0.55
1:A:140:ARG:HG2	1:C:140:ARG:HG2	1.88	0.55
1:C:95:GLU:HG3	1:C:112:THR:HB	1.89	0.55
1:C:115:LEU:HD23	1:C:136:THR:HG22	1.88	0.55
1:C:478:VAL:HG22	1:C:488:TYR:HE1	1.71	0.55
1:A:478:VAL:HG22	1:A:488:TYR:HE1	1.71	0.55
1:C:235:ILE:HG13	1:C:303:ILE:HD11	1.88	0.55
1:B:249:GLY:O	1:B:253:VAL:HG23	2.07	0.55
1:A:139:LEU:HB2	1:C:141:ARG:HH21	1.72	0.55
1:D:185:HIS:HD2	1:D:264:LEU:HD13	1.71	0.55
1:D:249:GLY:O	1:D:253:VAL:HG23	2.07	0.54
1:A:95:GLU:HG3	1:A:112:THR:HB	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:648:PRO:HG3	1:A:693:GLN:HA	1.90	0.54
1:B:95:GLU:HG3	1:B:112:THR:HB	1.90	0.54
1:D:95:GLU:HG3	1:D:112:THR:HB	1.90	0.54
1:C:648:PRO:HG3	1:C:693:GLN:HA	1.90	0.54
1:B:349:GLU:HA	1:B:352:ASN:HB2	1.91	0.53
1:D:102:LEU:HD22	1:D:169:LEU:HD21	1.91	0.53
1:A:68:VAL:HG11	1:A:74:VAL:HG22	1.91	0.53
1:A:376:ILE:HG21	1:A:502:GLU:HG3	1.92	0.52
1:D:385:MET:HE2	1:D:411:PHE:HD2	1.74	0.52
1:C:68:VAL:HG11	1:C:74:VAL:HG22	1.91	0.52
1:C:376:ILE:HG21	1:C:502:GLU:HG3	1.91	0.52
1:D:349:GLU:HA	1:D:352:ASN:HB2	1.91	0.52
1:B:102:LEU:HD22	1:B:169:LEU:HD21	1.91	0.52
1:B:385:MET:HE2	1:B:411:PHE:HD2	1.74	0.52
1:C:83:LEU:HD12	1:C:86:ILE:HD12	1.93	0.51
1:B:235:ILE:HG13	1:B:303:ILE:HD11	1.92	0.51
1:D:235:ILE:HG13	1:D:303:ILE:HD11	1.92	0.51
1:A:461:MET:HE3	1:A:461:MET:HA	1.93	0.51
1:A:355:VAL:HG21	1:B:369:ARG:NH2	2.25	0.51
1:A:111:LEU:HD11	1:A:160:TRP:CE2	2.46	0.50
1:C:238:LYS:HG3	1:C:241:TYR:HB3	1.93	0.50
1:A:238:LYS:HG3	1:A:241:TYR:HB3	1.93	0.50
1:C:618:GLY:HA3	1:C:706:LEU:HD23	1.93	0.50
1:D:421:ILE:HD13	1:D:472:LEU:HD23	1.93	0.50
1:C:461:MET:HE3	1:C:461:MET:HA	1.93	0.50
1:A:83:LEU:HD12	1:A:86:ILE:HD12	1.93	0.50
1:A:185:HIS:HD2	1:A:264:LEU:HD13	1.77	0.50
1:B:452:VAL:HG11	1:B:461:MET:HE1	1.94	0.49
1:C:111:LEU:HD11	1:C:160:TRP:CE2	2.46	0.49
1:A:318:PHE:CZ	1:A:322:LYS:HE3	2.47	0.49
1:D:514:MET:HG3	1:D:515:TYR:CD2	2.47	0.49
1:A:276:THR:O	1:A:280:VAL:HG23	2.13	0.49
1:D:452:VAL:HG11	1:D:461:MET:HE1	1.94	0.49
1:B:421:ILE:HD13	1:B:472:LEU:HD23	1.94	0.49
1:C:355:VAL:HG21	1:D:369:ARG:NH2	2.25	0.49
1:D:604:THR:HB	1:D:607:LYS:HD2	1.95	0.49
1:C:546:ILE:HD11	1:C:551:LEU:HD13	1.95	0.49
1:A:618:GLY:HA3	1:A:706:LEU:HD23	1.93	0.49
1:B:514:MET:HG3	1:B:515:TYR:CD2	2.47	0.49
1:A:452:VAL:HG11	1:A:461:MET:SD	2.54	0.48
1:C:276:THR:O	1:C:280:VAL:HG23	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:619:LYS:HG2	1:B:634:GLY:HA2	1.96	0.48
1:C:185:HIS:HD2	1:C:264:LEU:HD13	1.77	0.48
1:C:98:ASP:CG	1:C:99:ALA:H	2.22	0.48
1:C:318:PHE:CZ	1:C:322:LYS:HE3	2.47	0.48
1:C:452:VAL:HG11	1:C:461:MET:SD	2.54	0.48
1:D:402:GLU:HA	1:D:528:ASN:HD22	1.79	0.48
1:A:98:ASP:CG	1:A:99:ALA:H	2.22	0.48
1:B:402:GLU:HA	1:B:528:ASN:HD22	1.79	0.48
1:B:454:HIS:CE1	1:B:477:LYS:HG3	2.49	0.48
1:D:386:ILE:HG23	1:D:410:VAL:HG22	1.96	0.48
1:A:546:ILE:HD11	1:A:551:LEU:HD13	1.95	0.48
1:B:386:ILE:HG23	1:B:410:VAL:HG22	1.96	0.47
1:A:224:LYS:HA	1:A:227:ARG:HE	1.79	0.47
1:A:515:TYR:HE1	1:A:523:ARG:HG2	1.80	0.47
1:C:224:LYS:HA	1:C:227:ARG:HE	1.79	0.47
1:C:355:VAL:HG23	1:C:358:GLU:H	1.80	0.47
1:B:51:LEU:HD21	1:B:58:CYS:SG	2.54	0.47
1:B:604:THR:HB	1:B:607:LYS:HD2	1.95	0.47
1:D:51:LEU:HD21	1:D:58:CYS:SG	2.55	0.47
1:D:454:HIS:CE1	1:D:477:LYS:HG3	2.49	0.47
1:D:619:LYS:HG2	1:D:634:GLY:HA2	1.96	0.47
1:C:515:TYR:HE1	1:C:523:ARG:HG2	1.80	0.47
1:A:428:ALA:HB1	1:B:401:MET:HE3	1.96	0.47
1:C:185:HIS:CD2	1:C:264:LEU:HB2	2.50	0.47
1:D:118:GLN:HG2	1:D:119:GLN:HG2	1.97	0.47
1:A:238:LYS:CG	1:A:241:TYR:HB3	2.45	0.47
1:A:355:VAL:HG23	1:A:358:GLU:H	1.80	0.47
1:B:118:GLN:HG2	1:B:119:GLN:HG2	1.97	0.46
1:A:450:HIS:HB2	1:A:470:SER:HB3	1.97	0.46
1:B:646:SER:HB3	1:B:713:ARG:NH2	2.29	0.46
1:D:385:MET:HE2	1:D:411:PHE:CD2	2.50	0.46
1:B:385:MET:HE2	1:B:411:PHE:CD2	2.50	0.46
1:C:450:HIS:HB2	1:C:470:SER:HB3	1.98	0.46
1:D:338:ARG:NH1	1:D:369:ARG:HG2	2.31	0.46
1:D:646:SER:HB3	1:D:713:ARG:NH2	2.29	0.46
1:C:480:ASN:HA	1:C:486:PRO:HB3	1.98	0.46
1:A:480:ASN:HA	1:A:486:PRO:HB3	1.98	0.46
1:B:185:HIS:CE1	1:B:263:ASN:HD22	2.34	0.46
1:D:529:LYS:HA	1:D:532:PHE:CD2	2.51	0.46
1:A:185:HIS:CD2	1:A:264:LEU:HB2	2.50	0.46
1:D:185:HIS:CE1	1:D:263:ASN:HD22	2.34	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:238:LYS:CG	1:C:241:TYR:HB3	2.45	0.45
1:C:428:ALA:HB1	1:D:401:MET:HE3	1.96	0.45
1:A:247:LEU:HD12	1:B:285:ILE:HG12	1.97	0.45
1:C:536:LYS:HG3	1:C:536:LYS:O	2.16	0.45
1:D:318:PHE:HB3	1:D:319:PRO:HD3	1.98	0.45
1:B:227:ARG:HG2	1:B:231:LYS:HE3	1.99	0.45
1:C:247:LEU:HD12	1:D:285:ILE:HG12	1.97	0.45
1:B:318:PHE:HB3	1:B:319:PRO:HD3	1.98	0.45
1:B:529:LYS:HA	1:B:532:PHE:CD2	2.51	0.45
1:D:377:MET:HE1	1:D:491:LEU:HB3	1.99	0.45
1:A:536:LYS:O	1:A:536:LYS:HG3	2.16	0.45
1:B:99:ALA:O	1:B:102:LEU:HB3	2.17	0.45
1:C:199:PHE:HB3	1:C:250:ASN:HB2	1.99	0.45
1:A:99:ALA:O	1:A:102:LEU:HB3	2.17	0.44
1:A:128:THR:HB	1:B:47:VAL:HG13	1.99	0.44
1:A:51:LEU:HD21	1:A:58:CYS:SG	2.58	0.44
1:C:50:ARG:HG3	1:C:51:LEU:H	1.82	0.44
1:C:147:LEU:HD11	1:C:169:LEU:HB3	1.99	0.44
1:C:99:ALA:O	1:C:102:LEU:HB3	2.17	0.44
1:D:227:ARG:HG2	1:D:231:LYS:HE3	1.99	0.44
1:B:338:ARG:NH1	1:B:369:ARG:HG2	2.31	0.44
1:C:50:ARG:CG	1:D:50:ARG:HG2	2.34	0.44
1:D:45:THR:N	1:D:86:ILE:HD11	2.32	0.44
1:A:147:LEU:HD11	1:A:169:LEU:HB3	2.00	0.44
1:B:377:MET:HE1	1:B:491:LEU:HB3	1.99	0.44
1:C:51:LEU:HD21	1:C:58:CYS:SG	2.58	0.44
1:C:285:ILE:HG12	1:D:247:LEU:HD12	2.00	0.44
1:C:444:PHE:CZ	1:D:531:ASP:HA	2.53	0.44
1:A:444:PHE:CZ	1:B:531:ASP:HA	2.53	0.44
1:A:112:THR:HG21	1:A:148:TYR:CD2	2.53	0.43
1:C:78:LEU:O	1:C:129:SER:HA	2.18	0.43
1:C:128:THR:HB	1:D:47:VAL:HG13	2.00	0.43
1:D:515:TYR:HD1	1:D:520:SER:O	2.01	0.43
1:A:50:ARG:HG3	1:A:51:LEU:H	1.82	0.43
1:B:515:TYR:HD1	1:B:520:SER:O	2.01	0.43
1:B:45:THR:N	1:B:86:ILE:HD11	2.32	0.43
1:B:385:MET:HE3	1:B:409:PRO:HB2	2.00	0.43
1:C:386:ILE:HD12	1:C:400:ILE:HD11	2.00	0.43
1:A:199:PHE:HB3	1:A:250:ASN:HB2	1.99	0.43
1:B:601:TYR:HB3	1:B:603:TYR:O	2.19	0.43
1:C:112:THR:HG21	1:C:148:TYR:CD2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:579:LEU:HD22	1:D:709:ILE:HD13	2.01	0.43
1:A:318:PHE:HB3	1:A:319:PRO:HD3	2.00	0.43
1:C:585:ILE:HG12	1:C:709:ILE:HG12	2.01	0.43
1:A:78:LEU:O	1:A:129:SER:HA	2.18	0.43
1:B:65:ILE:HG21	1:B:171:PHE:HD2	1.83	0.43
1:D:99:ALA:O	1:D:102:LEU:HB3	2.18	0.43
1:A:285:ILE:HG12	1:B:247:LEU:HD12	2.00	0.43
1:C:197:GLY:HA2	1:C:283:GLY:O	2.18	0.43
1:A:75:ASN:HB3	1:A:131:VAL:HG13	2.01	0.43
1:A:197:GLY:HA2	1:A:283:GLY:O	2.18	0.43
1:A:280:VAL:HG21	1:B:251:VAL:HG11	2.01	0.43
1:A:386:ILE:HD12	1:A:400:ILE:HD11	2.00	0.43
1:A:585:ILE:HG12	1:A:709:ILE:HG12	2.01	0.43
1:C:394:PHE:CE2	1:C:536:LYS:HB3	2.54	0.43
1:D:65:ILE:HG21	1:D:171:PHE:HD2	1.83	0.43
1:D:223:GLU:HA	1:D:226:ARG:HE	1.84	0.43
1:A:394:PHE:CE2	1:A:536:LYS:HB3	2.54	0.43
1:B:579:LEU:HD22	1:B:709:ILE:HD13	2.01	0.43
1:B:223:GLU:HA	1:B:226:ARG:HE	1.84	0.42
1:C:75:ASN:HB3	1:C:131:VAL:HG13	2.01	0.42
1:C:98:ASP:CG	1:C:99:ALA:N	2.78	0.42
1:D:537:ASP:HA	1:D:540:ASN:HD22	1.84	0.42
1:A:98:ASP:CG	1:A:99:ALA:N	2.78	0.42
1:A:349:GLU:N	1:A:350:PRO:HD2	2.34	0.42
1:B:551:LEU:HD11	1:B:579:LEU:HB3	2.01	0.42
1:C:280:VAL:HG21	1:D:251:VAL:HG11	2.01	0.42
1:C:349:GLU:N	1:C:350:PRO:HD2	2.34	0.42
1:C:100:GLU:HB3	1:C:104:LYS:HE3	2.01	0.42
1:D:601:TYR:HB3	1:D:603:TYR:O	2.19	0.42
1:C:146:LYS:HE2	1:C:146:LYS:HB3	1.85	0.42
1:D:385:MET:HE3	1:D:409:PRO:HB2	2.00	0.42
1:C:172:MET:HE3	1:C:172:MET:HB2	1.96	0.42
1:D:551:LEU:HD11	1:D:579:LEU:HB3	2.01	0.42
1:C:318:PHE:HB3	1:C:319:PRO:HD3	2.01	0.42
1:D:50:ARG:HD3	1:D:66:ILE:HD11	2.02	0.42
1:A:100:GLU:HB3	1:A:104:LYS:HE3	2.01	0.42
1:A:392:LEU:HD12	1:A:441:ILE:HD13	2.02	0.41
1:A:612:PHE:CE2	1:A:638:TYR:HB2	2.55	0.41
1:C:369:ARG:HH21	1:D:355:VAL:HG21	1.85	0.41
1:C:612:PHE:CE2	1:C:638:TYR:HB2	2.55	0.41
1:D:192:LEU:HD12	1:D:260:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:LEU:HD11	1:B:172:MET:HG2	2.03	0.41
1:B:193:LEU:HD13	1:B:279:ILE:HD12	2.02	0.41
1:B:537:ASP:HA	1:B:540:ASN:HD22	1.84	0.41
1:C:384:PHE:HE2	1:C:408:ILE:HD13	1.84	0.41
2:F:1:NAG:H4	2:F:2:NAG:N2	2.36	0.41
1:A:384:PHE:HE2	1:A:408:ILE:HD13	1.85	0.41
1:C:185:HIS:CD2	1:C:264:LEU:HD13	2.55	0.41
1:C:392:LEU:HD12	1:C:441:ILE:HD13	2.02	0.41
1:C:478:VAL:HG22	1:C:488:TYR:CE1	2.53	0.41
1:D:78:LEU:HD11	1:D:132:LEU:HD22	2.02	0.41
2:J:1:NAG:H4	2:J:2:NAG:N2	2.36	0.41
1:B:68:VAL:HG11	1:B:74:VAL:HG21	2.03	0.41
1:A:591:ASP:H	1:A:600:HIS:CE1	2.39	0.41
1:B:232:ILE:HD12	1:B:296:LEU:HD12	2.03	0.41
1:A:199:PHE:HD1	1:A:199:PHE:HA	1.76	0.41
1:B:50:ARG:HD3	1:B:66:ILE:HD11	2.02	0.41
1:B:611:TYR:CG	1:B:710:LYS:HE3	2.56	0.41
1:D:139:LEU:HD11	1:D:172:MET:HG2	2.03	0.41
1:D:611:TYR:CG	1:D:710:LYS:HE3	2.56	0.41
1:B:336:TYR:HA	1:B:340:LYS:HD3	2.03	0.41
1:A:369:ARG:HH21	1:B:355:VAL:HG21	1.85	0.40
1:B:120:LEU:HD11	1:B:130:GLY:HA3	2.03	0.40
1:B:551:LEU:HD13	1:B:585:ILE:HD11	2.03	0.40
1:C:199:PHE:HD1	1:C:199:PHE:HA	1.76	0.40
1:C:247:LEU:HD23	1:C:247:LEU:HA	1.88	0.40
1:C:591:ASP:H	1:C:600:HIS:CE1	2.39	0.40
1:D:261:LEU:O	1:D:265:ILE:HG12	2.22	0.40
1:A:146:LYS:HE2	1:A:146:LYS:HB3	1.85	0.40
1:C:412:GLU:HB2	1:C:419:VAL:HG11	2.04	0.40
1:B:192:LEU:HD12	1:B:260:LEU:HD12	2.02	0.40
1:B:261:LEU:O	1:B:265:ILE:HG12	2.22	0.40
1:C:101:THR:HA	1:C:104:LYS:HB2	2.04	0.40
1:C:199:PHE:HE2	1:C:253:VAL:HG21	1.87	0.40
1:C:322:LYS:HD3	1:C:322:LYS:HA	1.89	0.40
1:D:193:LEU:HD13	1:D:279:ILE:HD12	2.03	0.40
1:D:314:PHE:N	1:D:315:PRO:HD2	2.37	0.40
1:A:412:GLU:HB2	1:A:419:VAL:HG11	2.04	0.40
1:B:78:LEU:HD11	1:B:132:LEU:HD22	2.02	0.40
1:D:68:VAL:HG11	1:D:74:VAL:HG21	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	633/844 (75%)	609 (96%)	24 (4%)	0	100	100
1	B	631/844 (75%)	616 (98%)	15 (2%)	0	100	100
1	C	633/844 (75%)	608 (96%)	25 (4%)	0	100	100
1	D	631/844 (75%)	616 (98%)	15 (2%)	0	100	100
All	All	2528/3376 (75%)	2449 (97%)	79 (3%)	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	572/737 (78%)	566 (99%)	6 (1%)	68	75
1	B	567/737 (77%)	561 (99%)	6 (1%)	65	74
1	C	572/737 (78%)	566 (99%)	6 (1%)	68	75
1	D	567/737 (77%)	561 (99%)	6 (1%)	65	74
All	All	2278/2948 (77%)	2254 (99%)	24 (1%)	63	74

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	VAL
1	A	123	VAL

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Mol	Chain	Res	Type
1	A	188	LEU
1	A	351	TYR
1	A	441	ILE
1	A	478	VAL
1	B	68	VAL
1	B	123	VAL
1	B	351	TYR
1	B	418	ILE
1	B	504	ILE
1	B	633	THR
1	C	74	VAL
1	C	123	VAL
1	C	188	LEU
1	C	351	TYR
1	C	441	ILE
1	C	478	VAL
1	D	68	VAL
1	D	123	VAL
1	D	351	TYR
1	D	418	ILE
1	D	504	ILE
1	D	633	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	HIS
1	A	219	ASN
1	A	250	ASN
1	A	254	ASN
1	A	479	ASN
1	A	518	ASN
1	B	185	HIS
1	B	203	ASN
1	B	250	ASN
1	B	254	ASN
1	B	263	ASN
1	B	454	HIS
1	C	185	HIS
1	C	219	ASN
1	C	250	ASN
1	C	254	ASN

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Mol	Chain	Res	Type
1	C	518	ASN
1	D	185	HIS
1	D	203	ASN
1	D	254	ASN
1	D	263	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 ⓘ

16 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	E	1	2,1	14,14,15	0.76	0	17,19,21	0.87	1 (5%)
2	NAG	E	2	2	14,14,15	0.75	0	17,19,21	0.87	0
2	NAG	F	1	2,1	14,14,15	0.73	0	17,19,21	1.54	2 (11%)
2	NAG	F	2	2	14,14,15	0.82	0	17,19,21	1.53	2 (11%)
2	NAG	G	1	2,1	14,14,15	0.71	0	17,19,21	0.99	0
2	NAG	G	2	2	14,14,15	0.75	0	17,19,21	0.81	0
2	NAG	H	1	2,1	14,14,15	0.78	0	17,19,21	1.30	3 (17%)
2	NAG	H	2	2	14,14,15	0.73	0	17,19,21	0.93	1 (5%)
2	NAG	I	1	2,1	14,14,15	0.74	0	17,19,21	0.87	1 (5%)
2	NAG	I	2	2	14,14,15	0.77	0	17,19,21	0.88	0
2	NAG	J	1	2,1	14,14,15	0.75	0	17,19,21	1.54	2 (11%)
2	NAG	J	2	2	14,14,15	0.84	1 (7%)	17,19,21	1.53	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	K	1	2,1	14,14,15	0.71	0	17,19,21	1.00	0
2	NAG	K	2	2	14,14,15	0.76	0	17,19,21	0.81	0
2	NAG	L	1	2,1	14,14,15	0.77	0	17,19,21	1.30	3 (17%)
2	NAG	L	2	2	14,14,15	0.72	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
2	NAG	E	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	NAG	F	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	F	2	2	-	4/6/23/26	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	1/6/23/26	0/1/1/1
2	NAG	H	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	1/6/23/26	0/1/1/1
2	NAG	I	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	NAG	J	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	J	2	2	-	4/6/23/26	0/1/1/1
2	NAG	K	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	K	2	2	-	1/6/23/26	0/1/1/1
2	NAG	L	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	L	2	2	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	2	NAG	C1-C2	2.02	1.55	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	2	NAG	C1-O5-C5	3.95	117.48	112.19
2	F	2	NAG	C1-O5-C5	3.94	117.47	112.19
2	F	1	NAG	C2-N2-C7	3.44	127.51	122.90
2	J	1	NAG	C2-N2-C7	3.44	127.51	122.90
2	J	1	NAG	O5-C1-C2	-3.04	106.58	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	NAG	O5-C1-C2	-3.03	106.60	111.29
2	J	2	NAG	C2-N2-C7	3.01	126.93	122.90
2	F	2	NAG	C2-N2-C7	2.98	126.89	122.90
2	H	1	NAG	O5-C1-C2	-2.65	107.20	111.29
2	L	1	NAG	O5-C1-C2	-2.63	107.22	111.29
2	H	1	NAG	C1-O5-C5	-2.45	108.90	112.19
2	L	1	NAG	C1-O5-C5	-2.44	108.91	112.19
2	L	2	NAG	O5-C1-C2	-2.20	107.89	111.29
2	H	2	NAG	O5-C1-C2	-2.15	107.96	111.29
2	L	1	NAG	C3-C4-C5	2.02	113.90	110.23
2	H	1	NAG	C3-C4-C5	2.02	113.90	110.23
2	E	1	NAG	O5-C1-C2	-2.01	108.19	111.29
2	I	1	NAG	O5-C1-C2	-2.01	108.19	111.29

There are no chirality outliers.

All (20) torsion outliers are listed below:

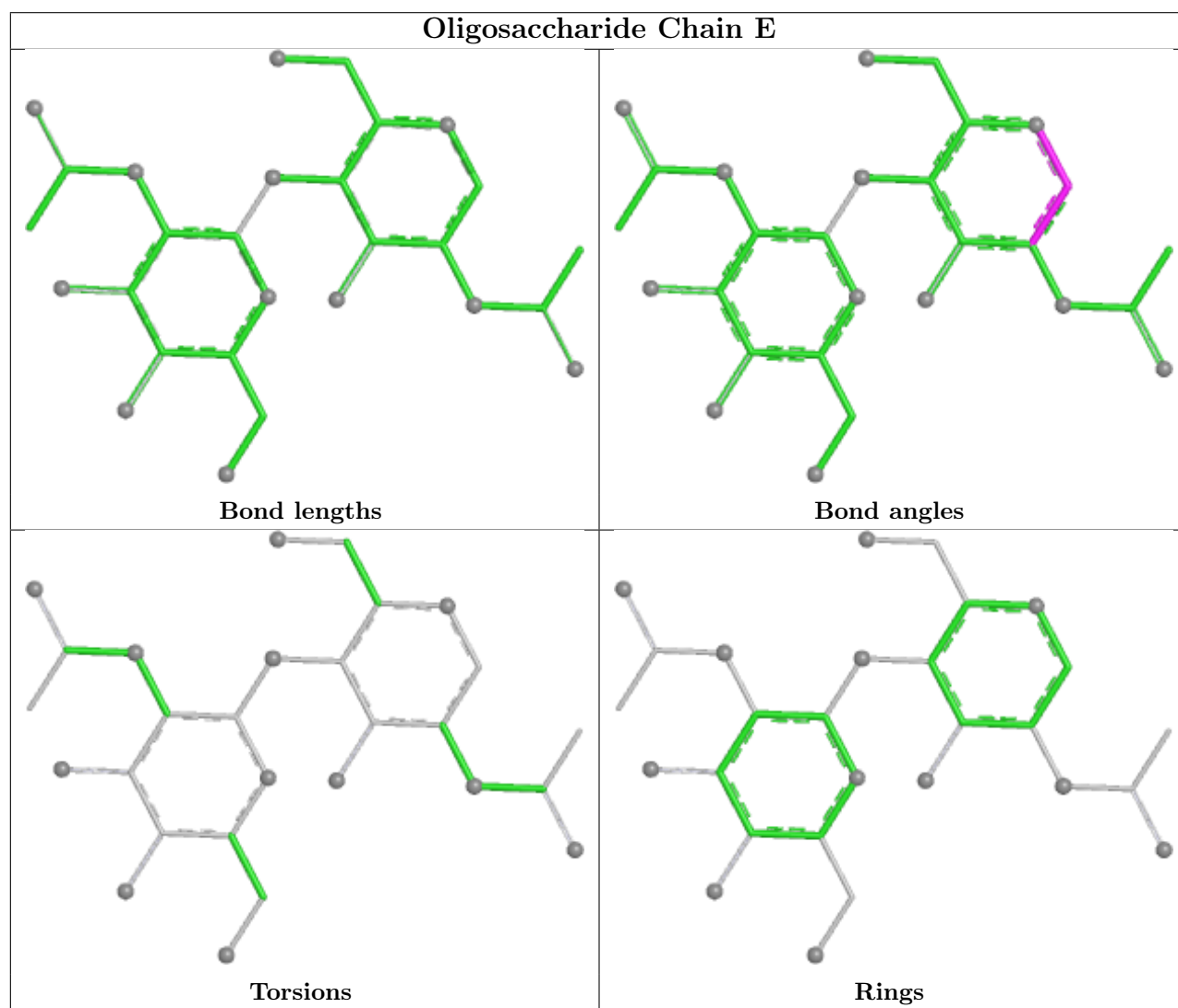
Mol	Chain	Res	Type	Atoms
2	F	2	NAG	O5-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
2	J	2	NAG	C4-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
2	J	1	NAG	O5-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
2	L	2	NAG	O5-C5-C6-O6
2	J	1	NAG	C4-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	F	1	NAG	C1-C2-N2-C7
2	F	2	NAG	C1-C2-N2-C7
2	J	1	NAG	C1-C2-N2-C7
2	J	2	NAG	C1-C2-N2-C7
2	F	1	NAG	C3-C2-N2-C7
2	F	2	NAG	C3-C2-N2-C7
2	J	1	NAG	C3-C2-N2-C7
2	J	2	NAG	C3-C2-N2-C7

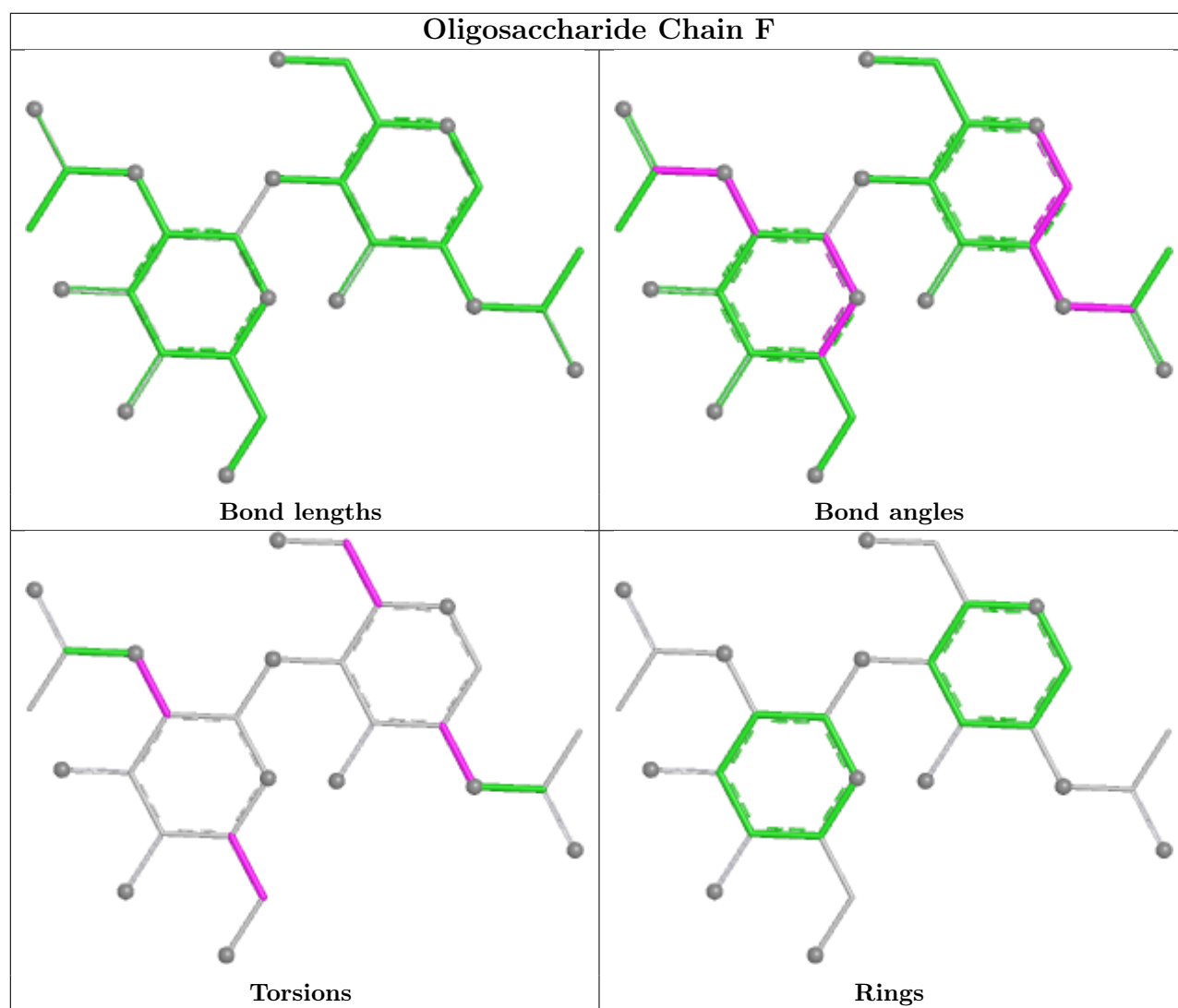
There are no ring outliers.

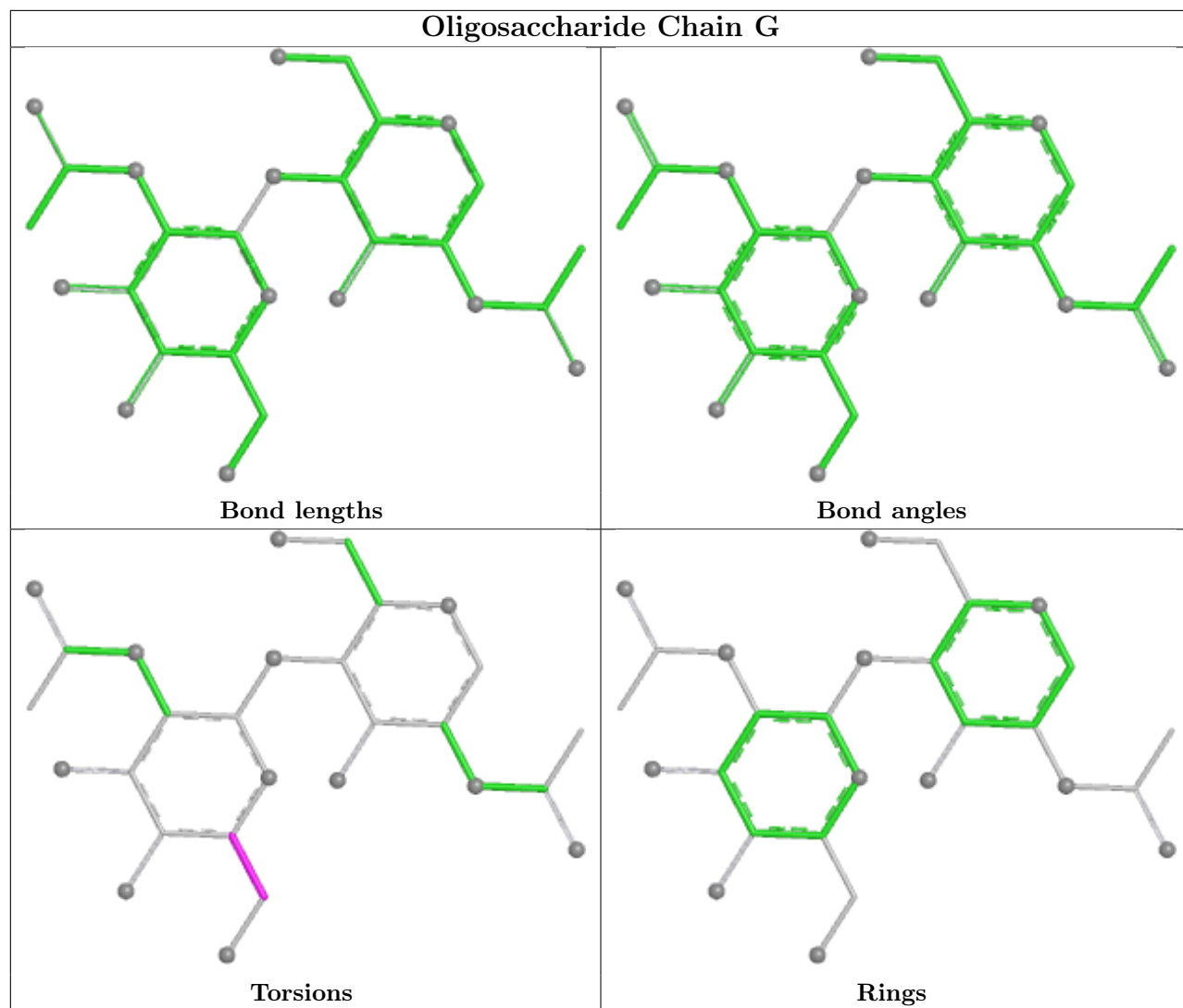
4 monomers are involved in 2 short contacts:

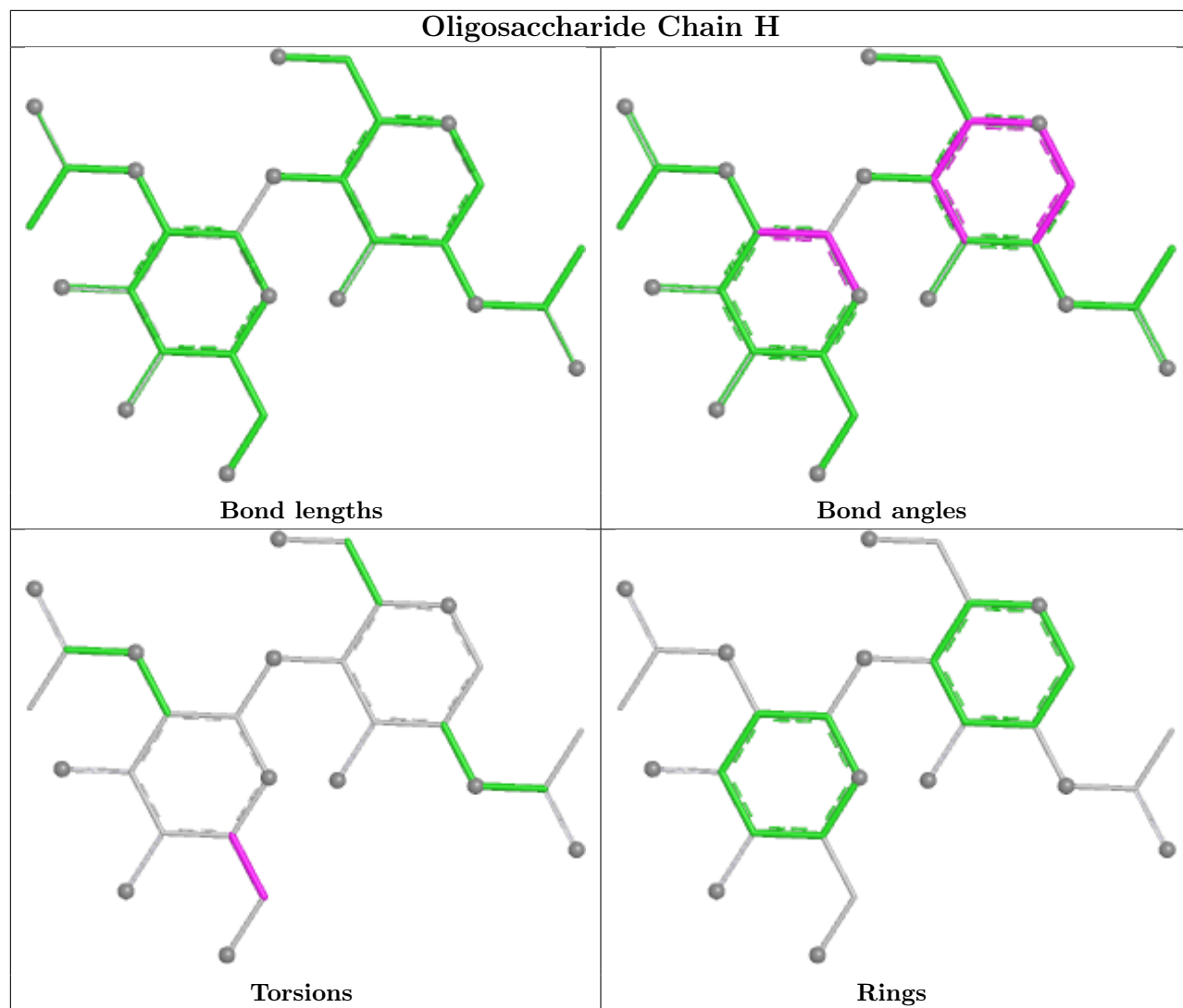
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	2	NAG	1	0
2	J	1	NAG	1	0
2	F	1	NAG	1	0
2	F	2	NAG	1	0

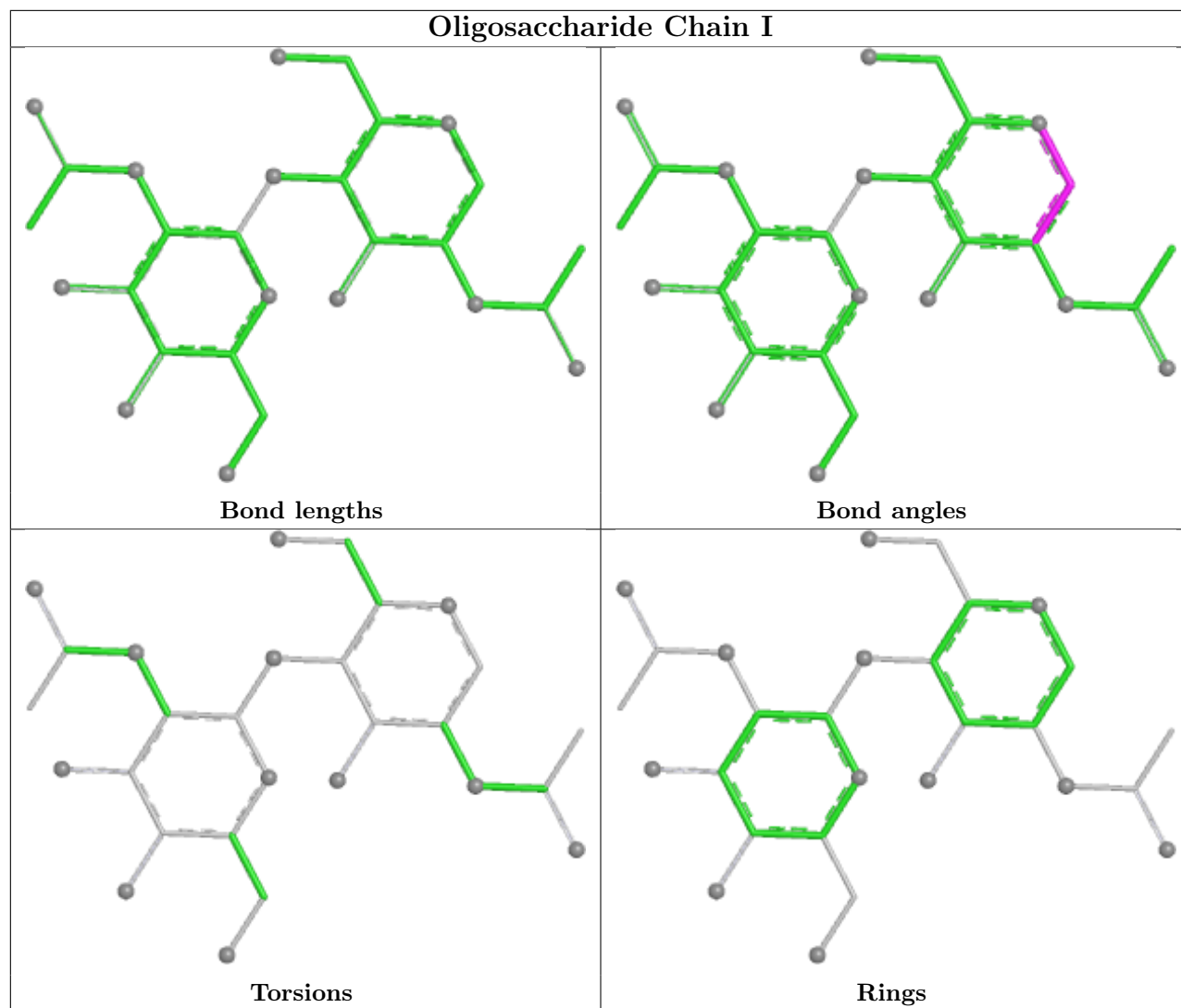
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

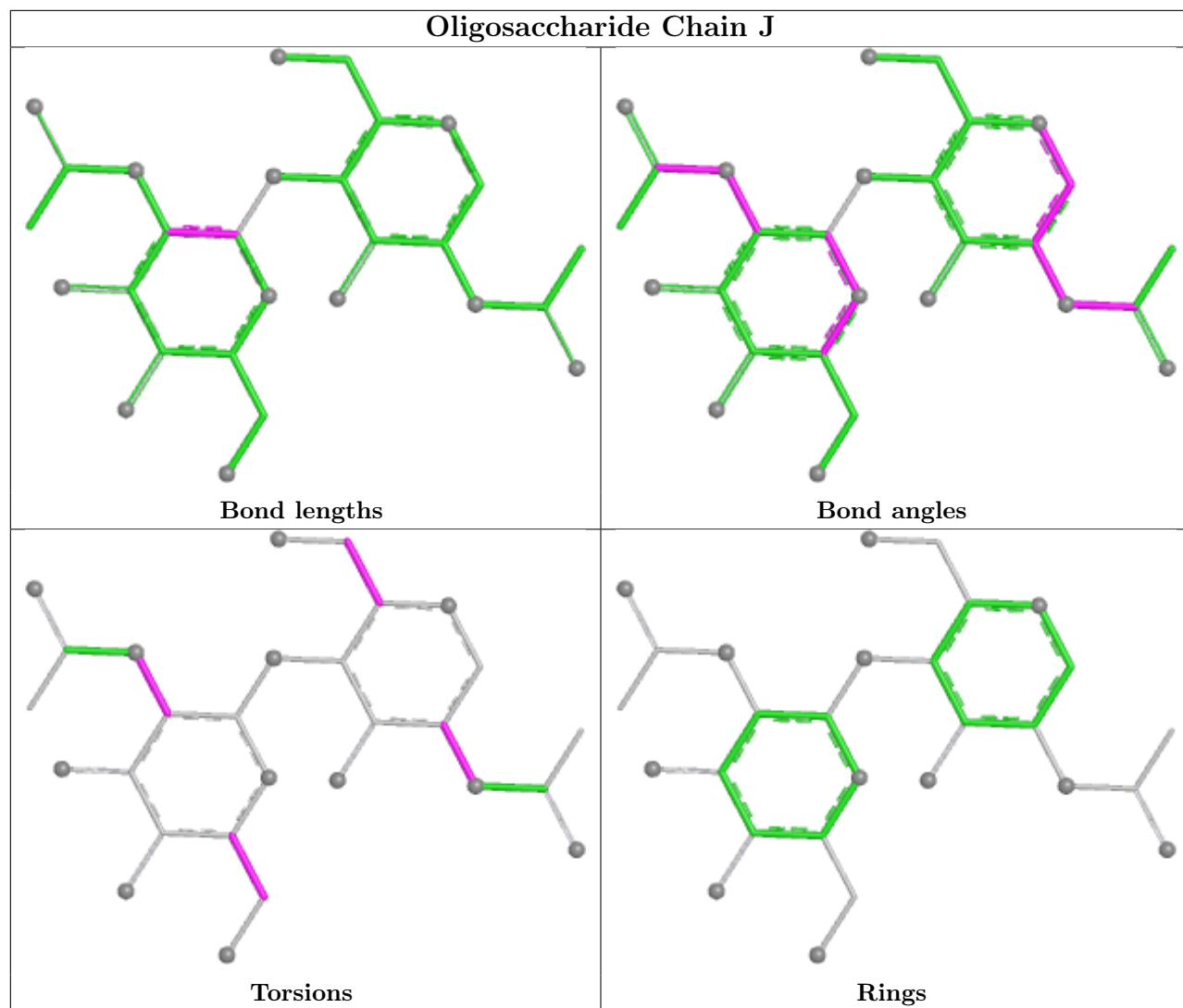


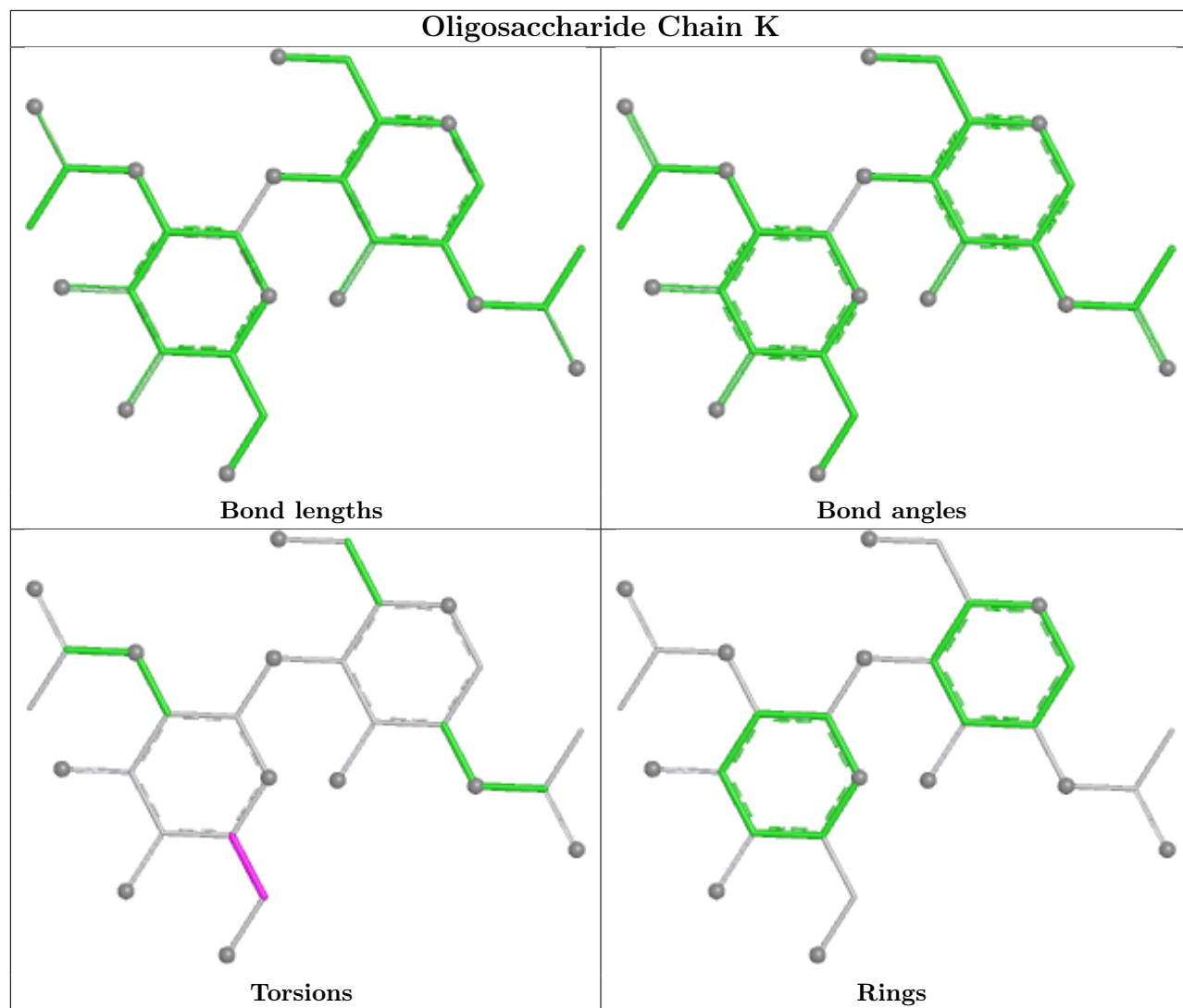


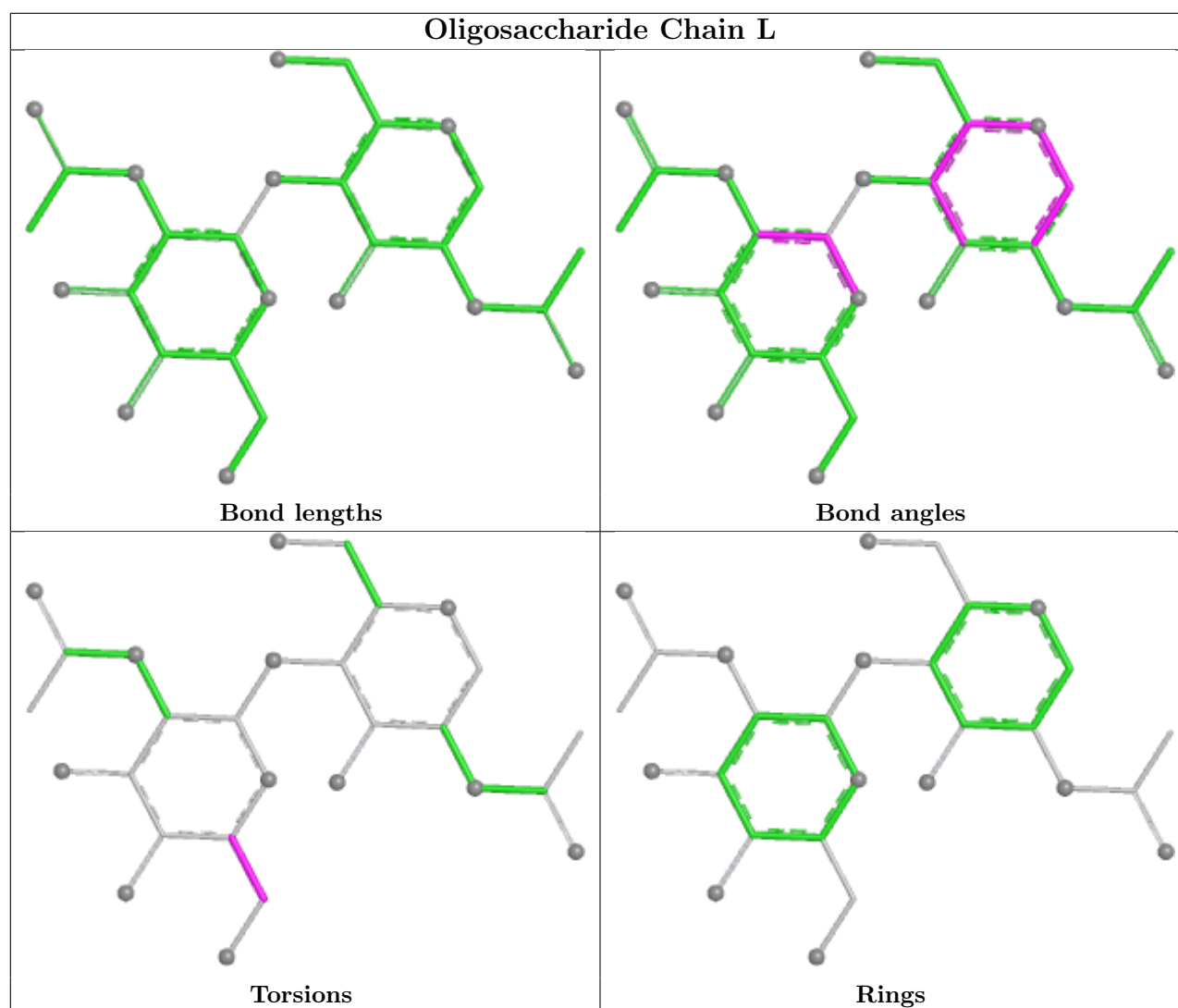












5.6 Ligand geometry [i](#)

Of 30 ligands modelled in this entry, 22 are monoatomic - leaving 8 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	$\# Z > 2$
6	NAG	C	908	1	14,14,15	0.74	0	17,19,21	1.31	1 (5%)
6	NAG	B	907	1	14,14,15	0.72	0	17,19,21	0.99	1 (5%)
3	ATP	D	901	4	32,33,33	0.43	0	48,52,52	0.40	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ATP	A	901	4	32,33,33	0.46	0	48,52,52	0.41	0
6	NAG	D	907	1	14,14,15	0.71	0	17,19,21	0.99	1 (5%)
3	ATP	C	901	4	32,33,33	0.45	0	48,52,52	0.41	0
3	ATP	B	901	4	32,33,33	0.43	0	48,52,52	0.40	0
6	NAG	A	908	1	14,14,15	0.77	0	17,19,21	1.31	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
6	NAG	C	908	1	-	0/6/23/26	0/1/1/1
6	NAG	B	907	1	-	1/6/23/26	0/1/1/1
3	ATP	D	901	4	-	4/22/38/38	0/3/3/3
3	ATP	A	901	4	-	4/22/38/38	0/3/3/3
6	NAG	D	907	1	-	1/6/23/26	0/1/1/1
3	ATP	C	901	4	-	4/22/38/38	0/3/3/3
3	ATP	B	901	4	-	4/22/38/38	0/3/3/3
6	NAG	A	908	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	908	NAG	C1-O5-C5	3.96	117.49	112.19
6	C	908	NAG	C1-O5-C5	3.94	117.46	112.19
6	B	907	NAG	C1-O5-C5	2.81	115.95	112.19
6	D	907	NAG	C1-O5-C5	2.79	115.93	112.19

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	901	ATP	PB-O3B-PG-O2G
3	D	901	ATP	PB-O3B-PG-O2G
6	B	907	NAG	O5-C5-C6-O6
6	D	907	NAG	O5-C5-C6-O6
3	B	901	ATP	PB-O3B-PG-O1G
3	D	901	ATP	PB-O3B-PG-O1G

Continued on next page...

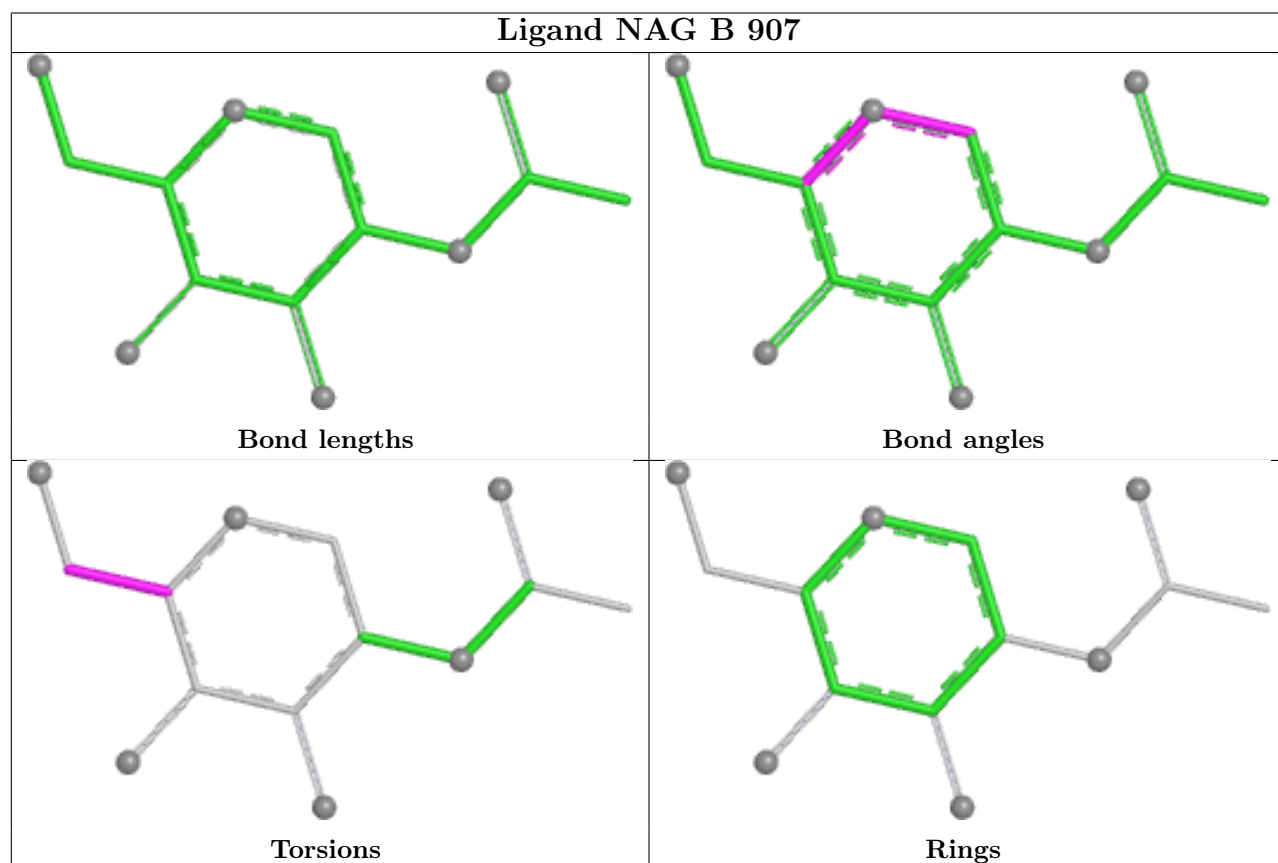
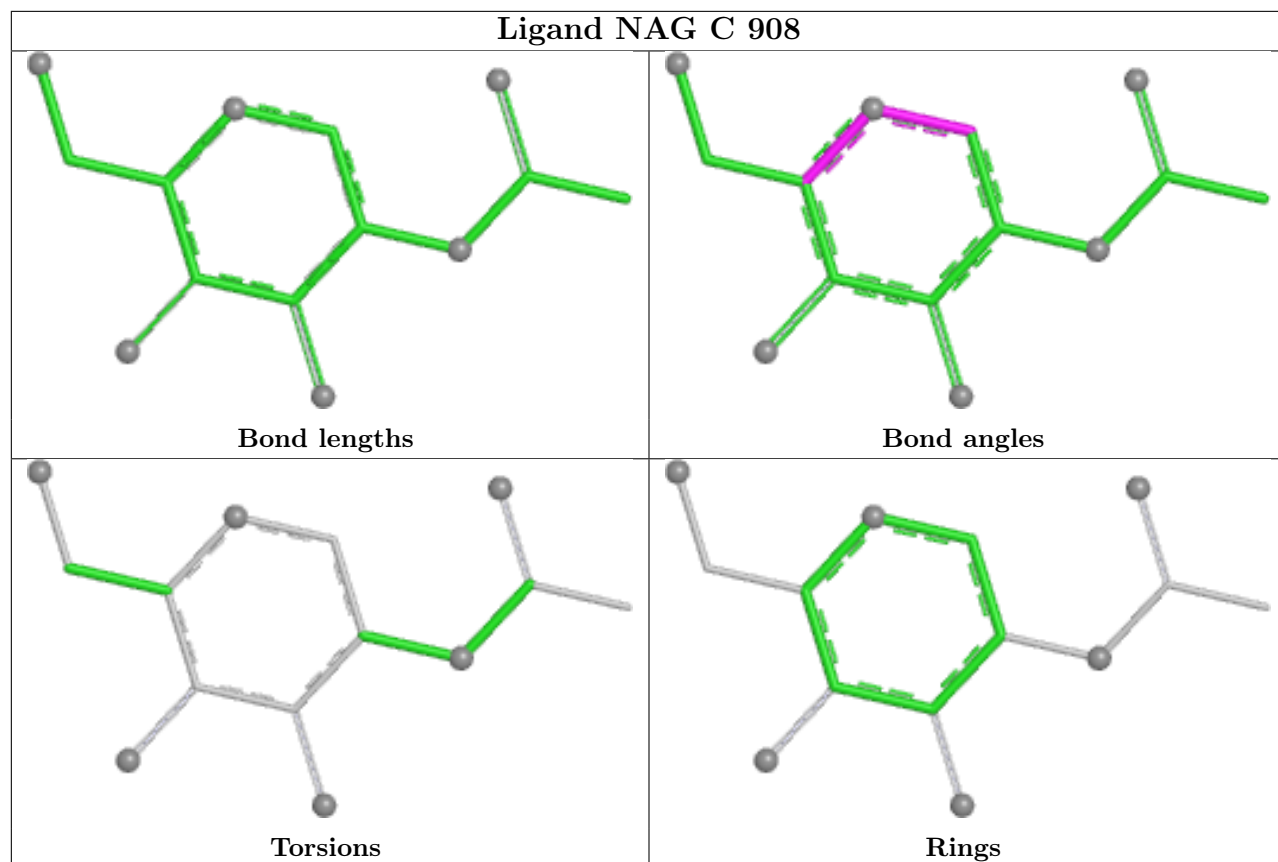
Continued from previous page...

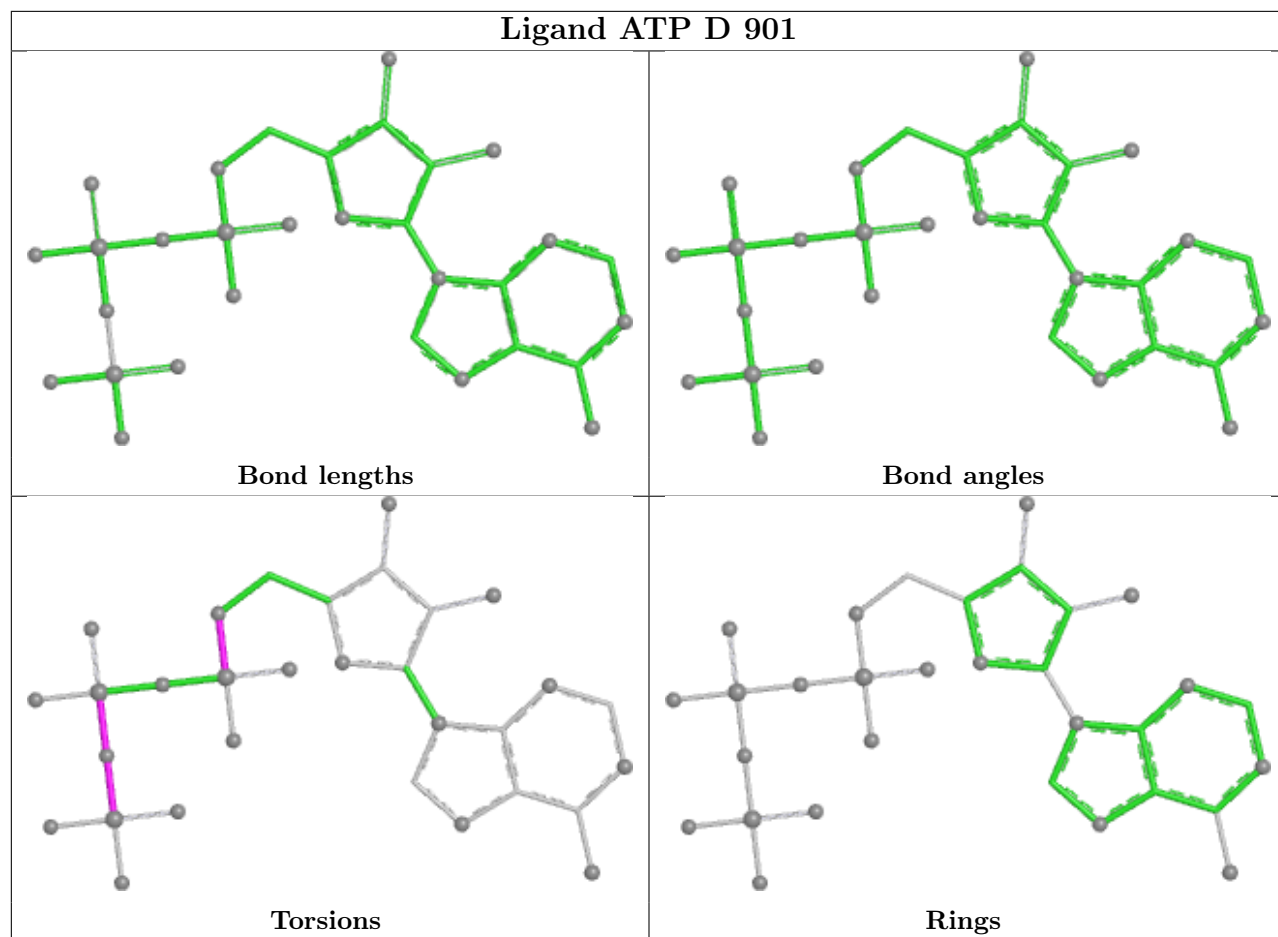
Mol	Chain	Res	Type	Atoms
3	A	901	ATP	PG-O3B-PB-O1B
3	C	901	ATP	PG-O3B-PB-O1B
3	A	901	ATP	C5'-O5'-PA-O1A
3	B	901	ATP	C5'-O5'-PA-O1A
3	C	901	ATP	C5'-O5'-PA-O1A
3	D	901	ATP	C5'-O5'-PA-O1A
3	B	901	ATP	PG-O3B-PB-O2B
3	D	901	ATP	PG-O3B-PB-O2B
3	A	901	ATP	PG-O3B-PB-O2B
3	A	901	ATP	PA-O3A-PB-O2B
3	C	901	ATP	PG-O3B-PB-O2B
3	C	901	ATP	PA-O3A-PB-O2B

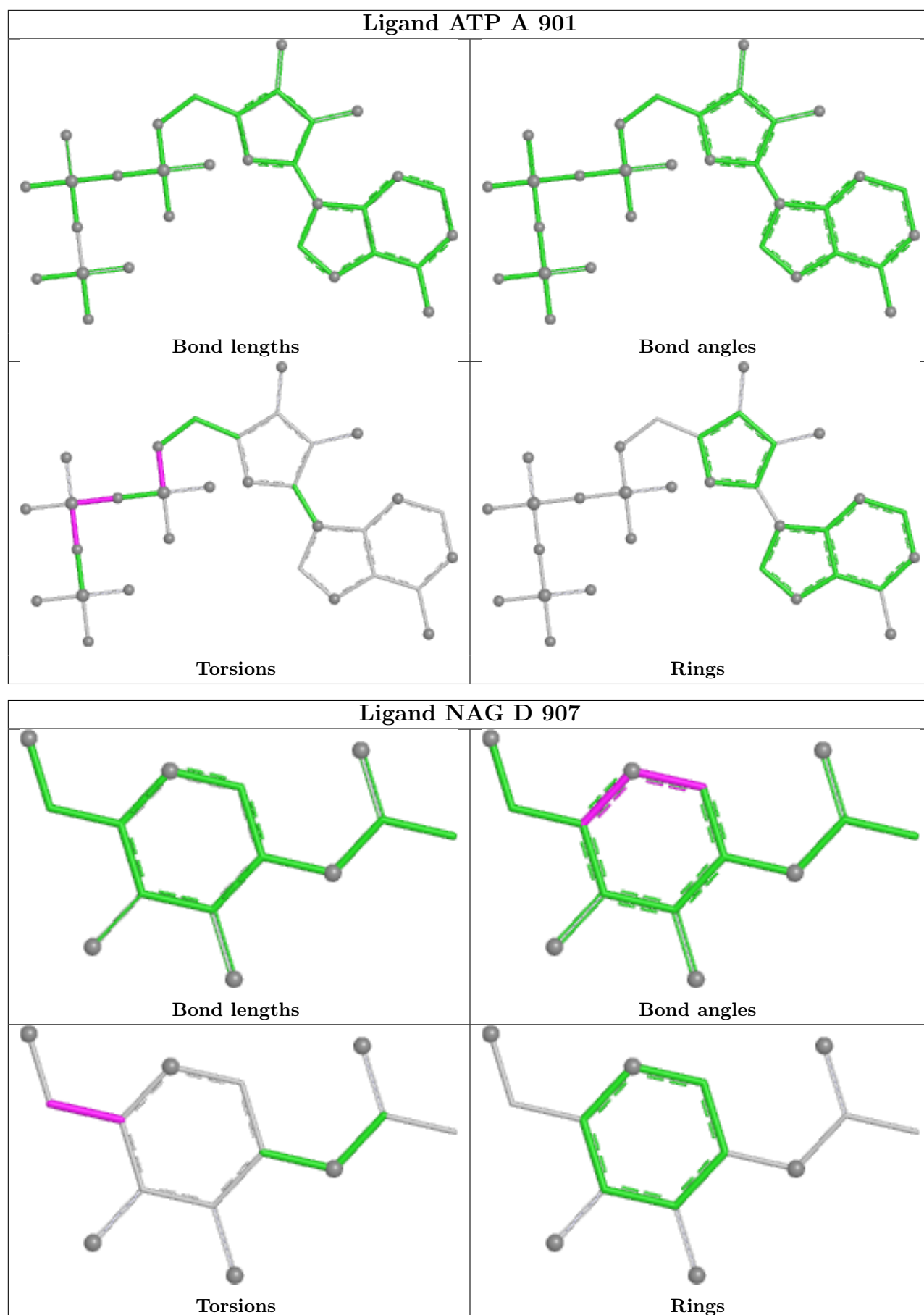
There are no ring outliers.

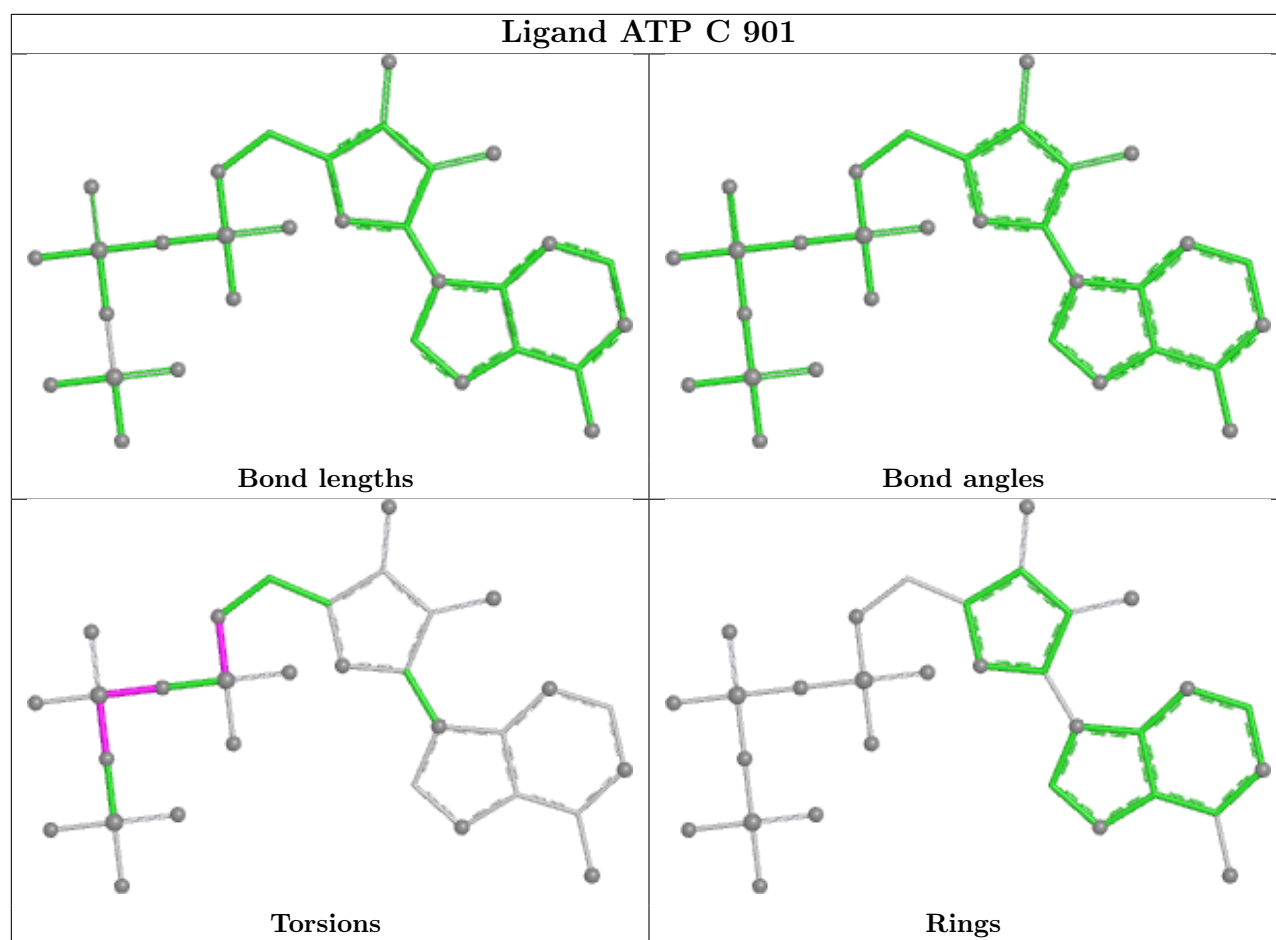
No monomer is involved in short contacts.

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

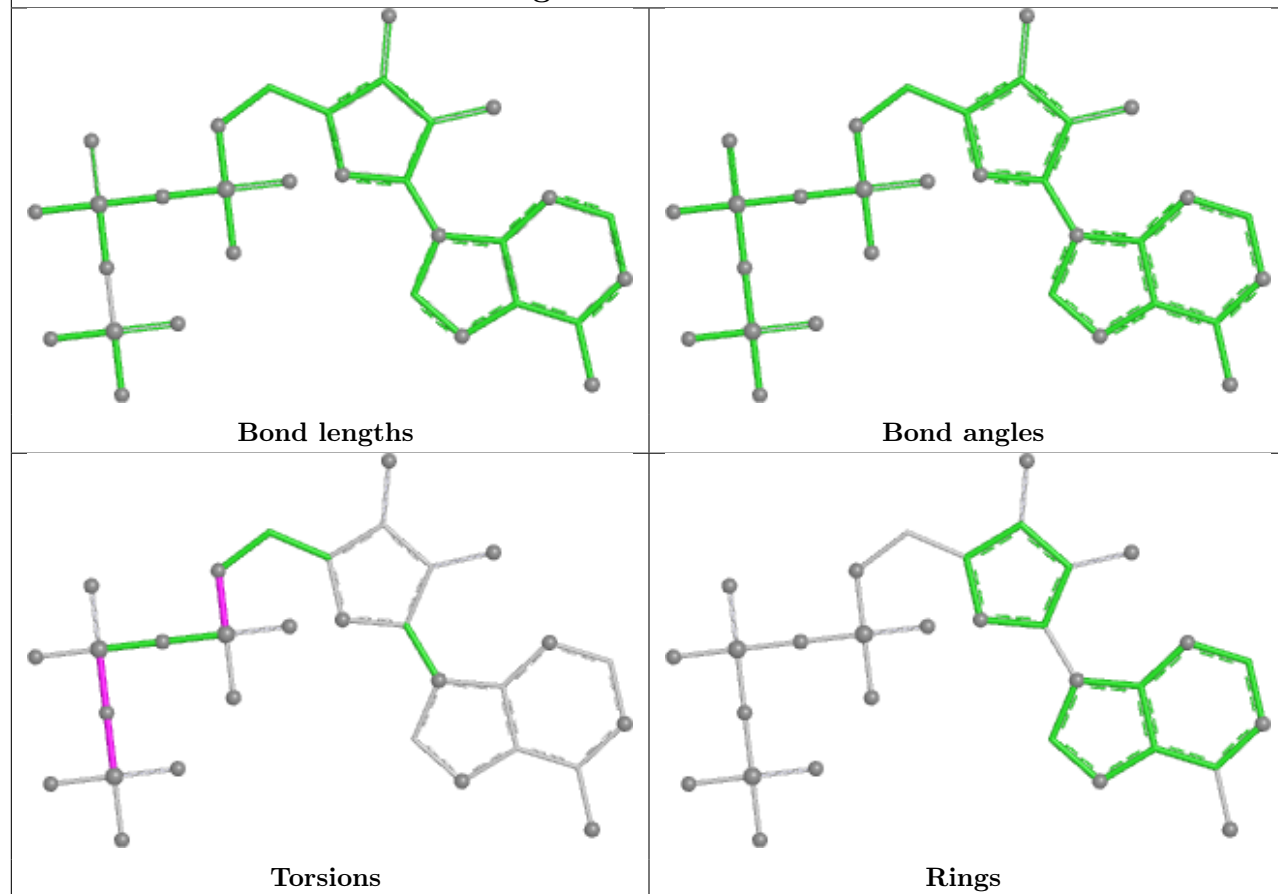




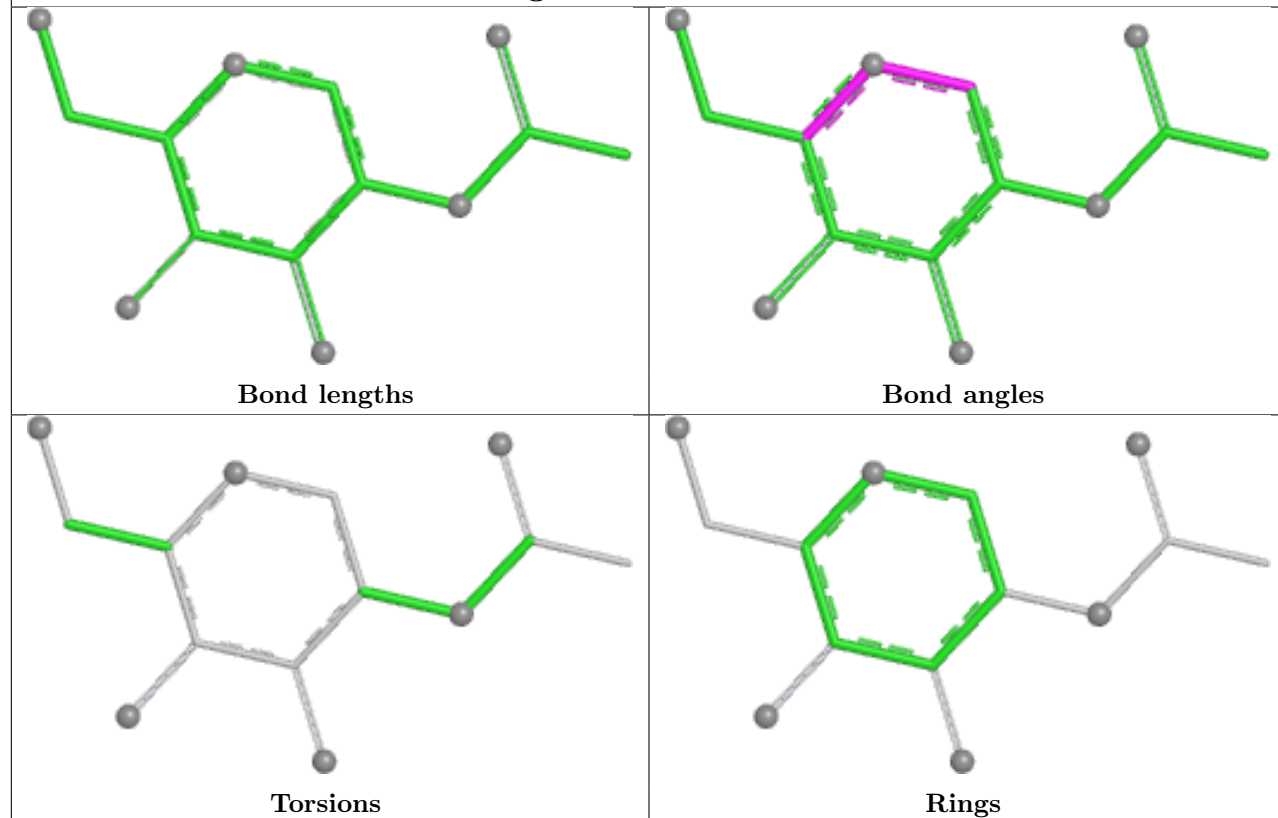




Ligand ATP B 901



Ligand NAG A 908



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.

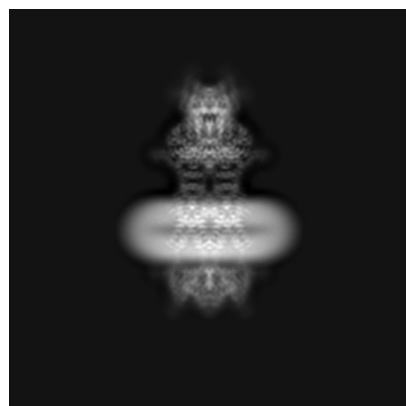
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-72694. These allow visual inspection of the internal detail of the map and identification of artifacts.

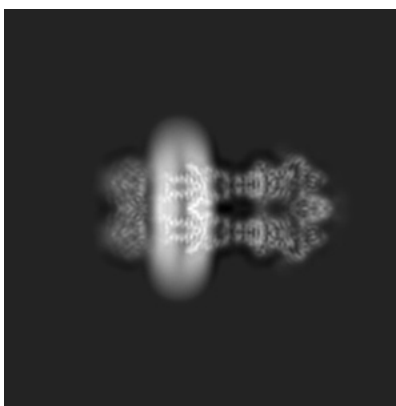
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

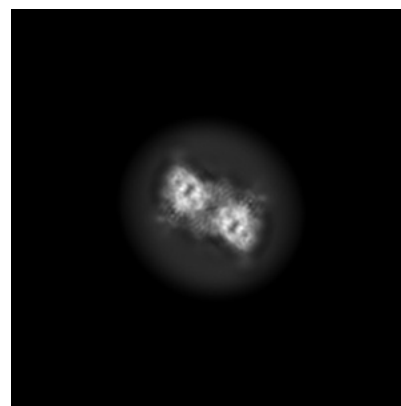
6.1.1 Primary map



X

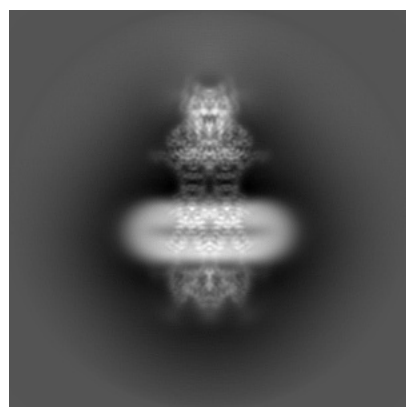


Y

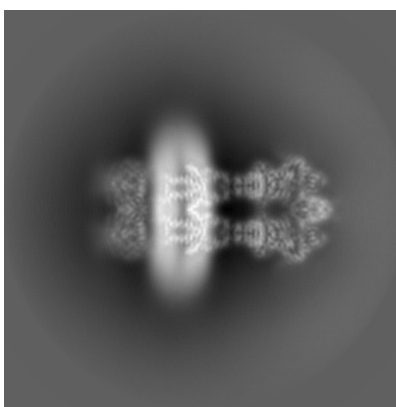


Z

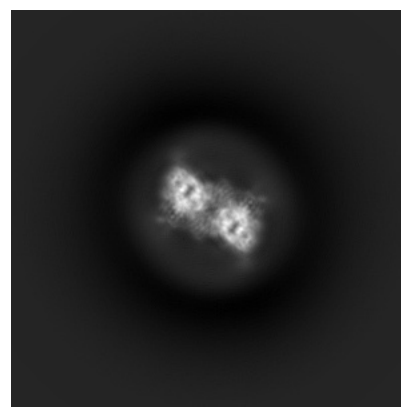
6.1.2 Raw map



X



Y

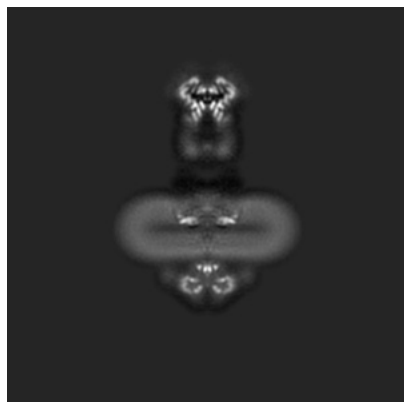


Z

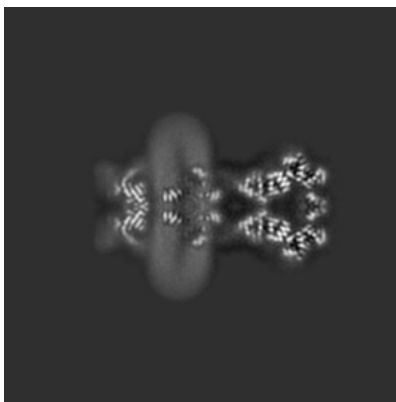
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

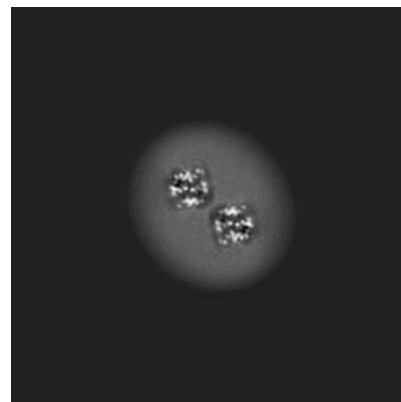
6.2.1 Primary map



X Index: 208

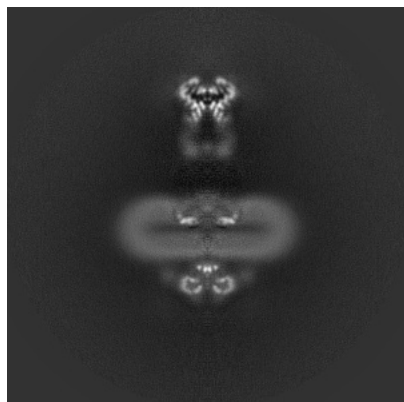


Y Index: 208

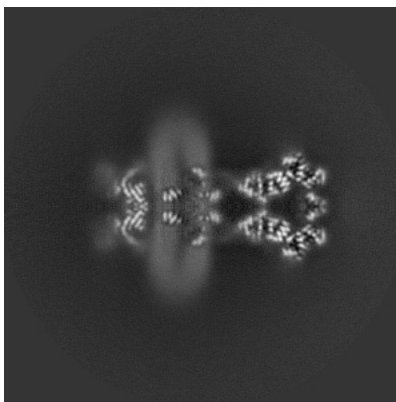


Z Index: 208

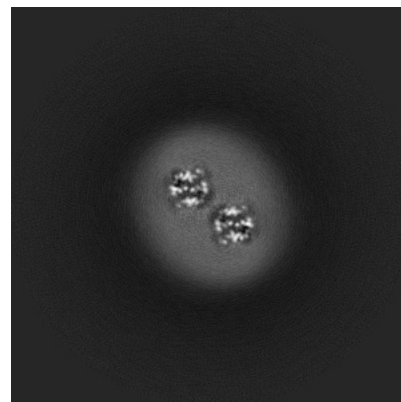
6.2.2 Raw map



X Index: 208



Y Index: 208

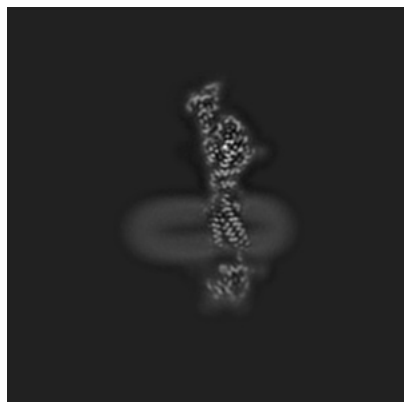


Z Index: 208

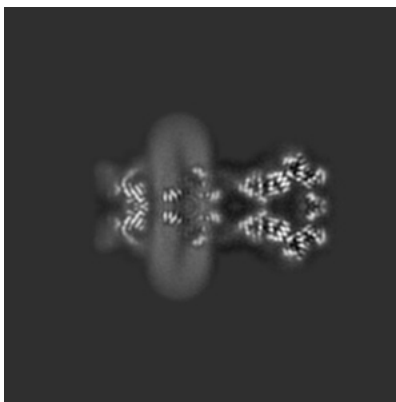
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

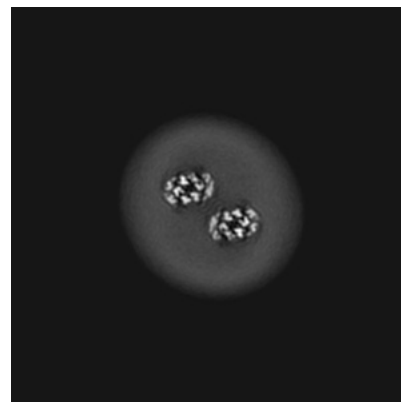
6.3.1 Primary map



X Index: 178

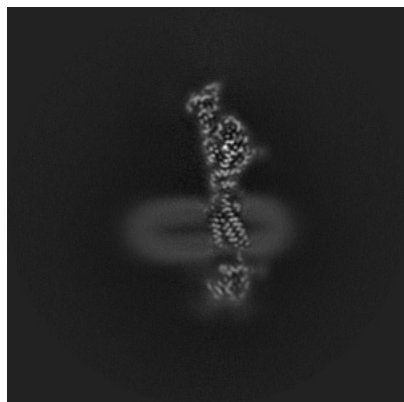


Y Index: 208

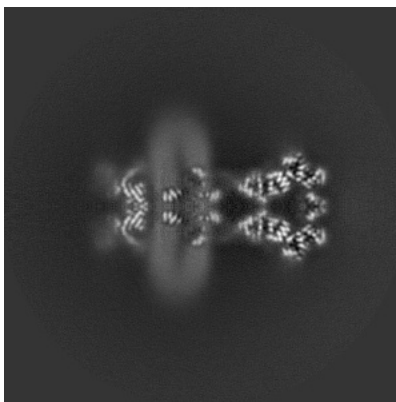


Z Index: 195

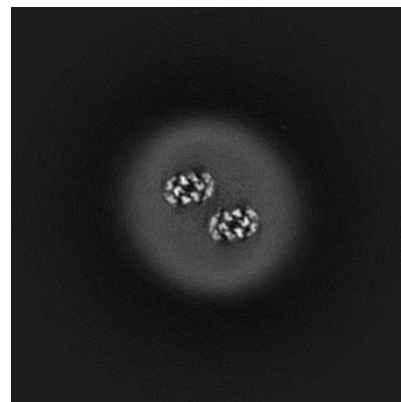
6.3.2 Raw map



X Index: 178



Y Index: 208

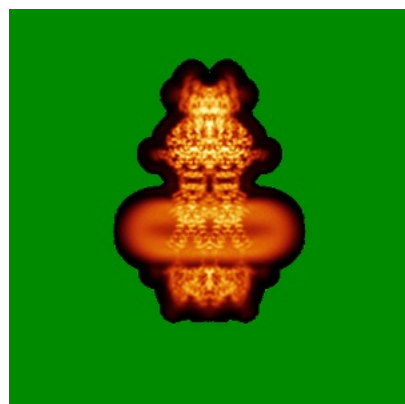


Z Index: 195

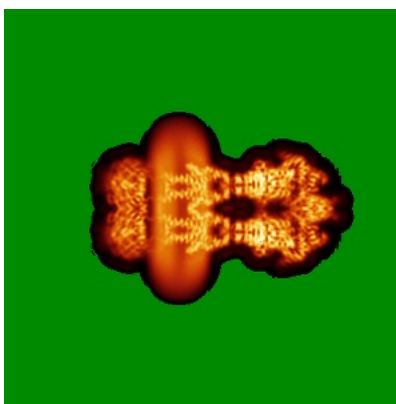
The images above show the largest variance slices of the map in three orthogonal directions.

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

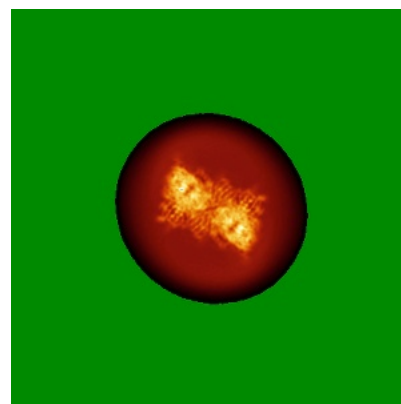
6.4.1 Primary map



X

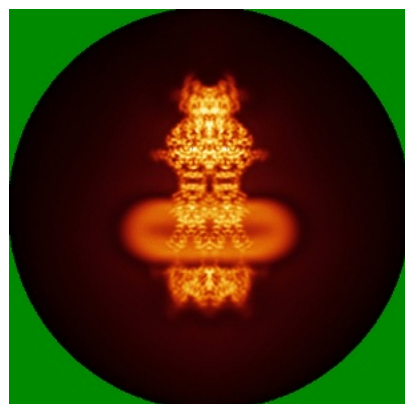


Y

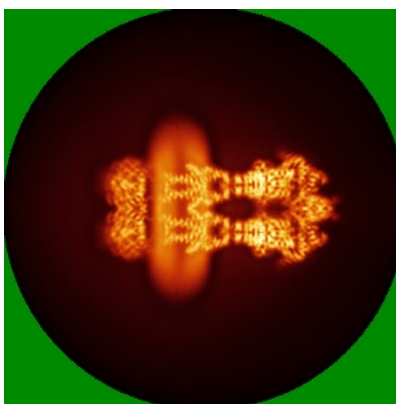


Z

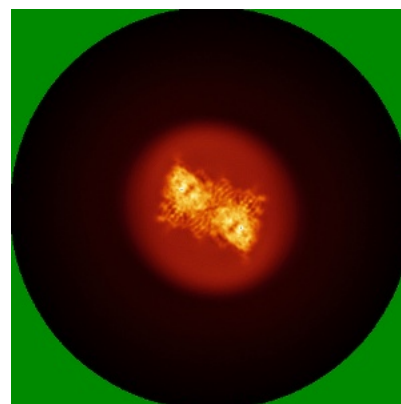
6.4.2 Raw map



X



Y

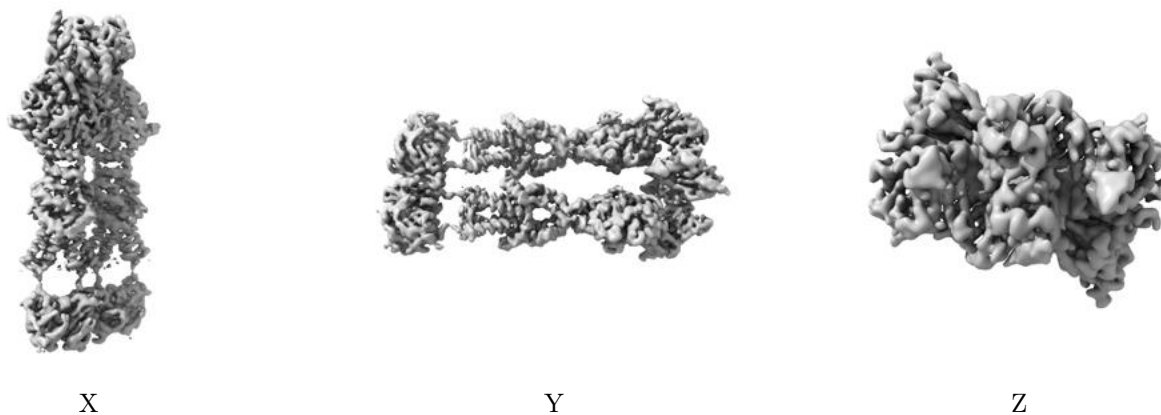


Z

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.

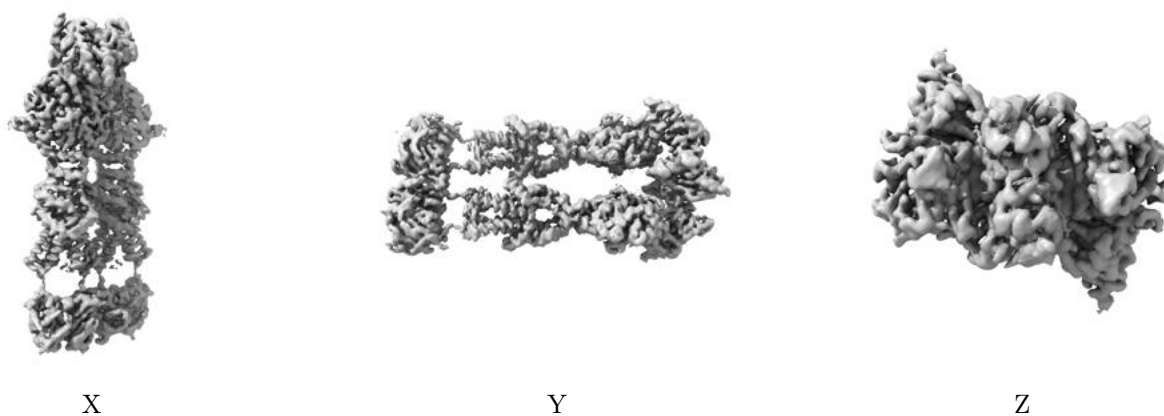
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.00376. 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.

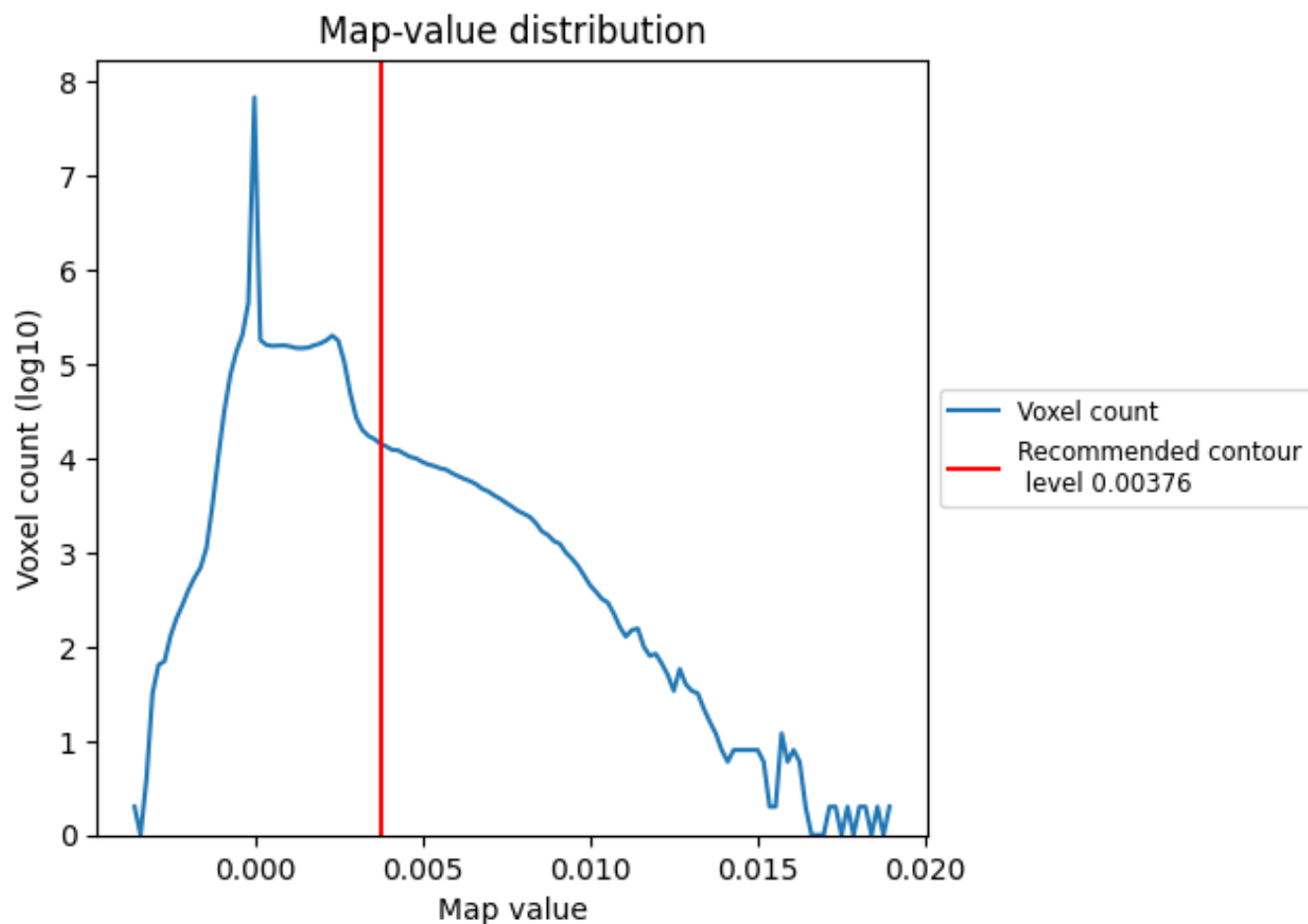
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

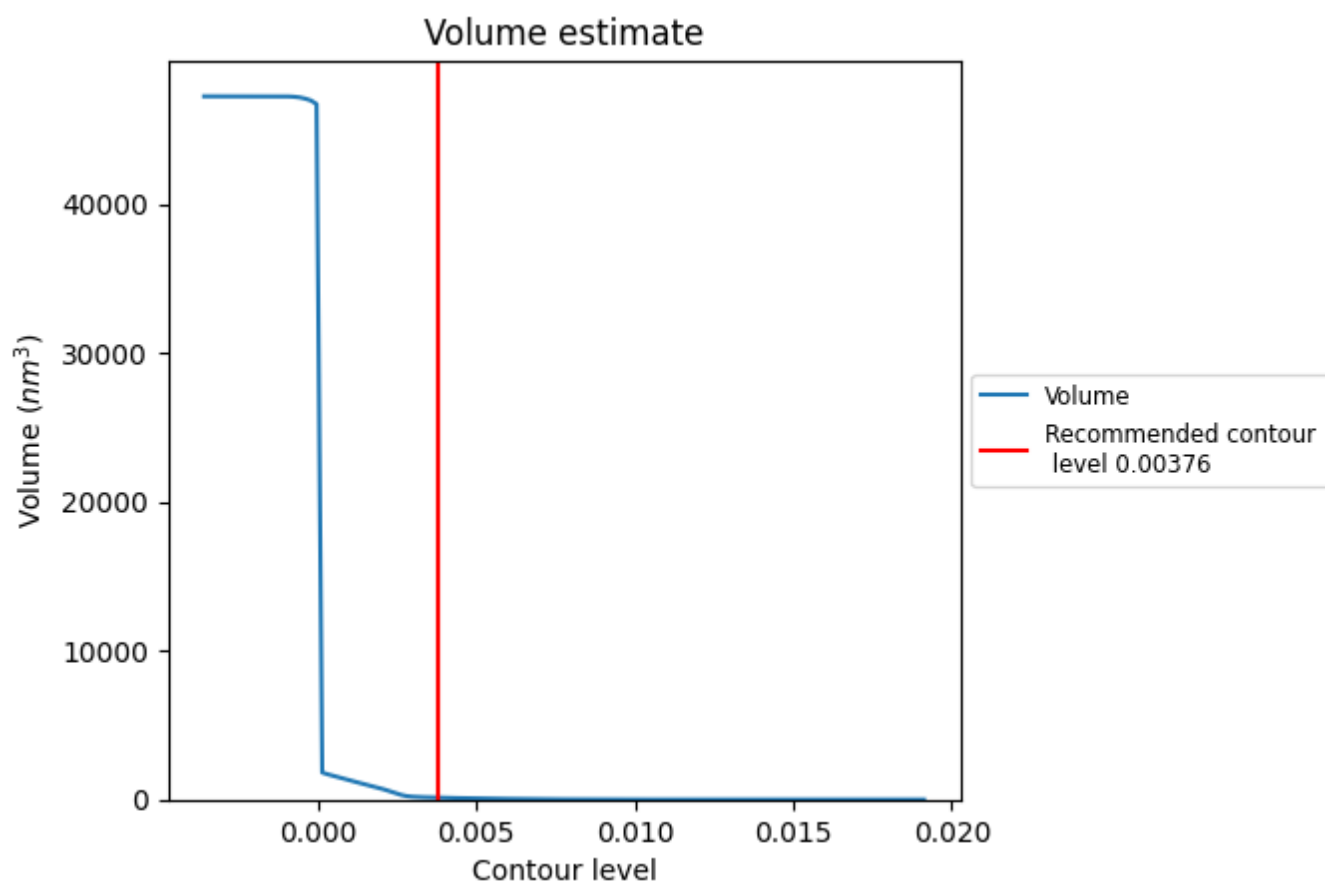
This section contains the results of statistical analysis of the map.

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.

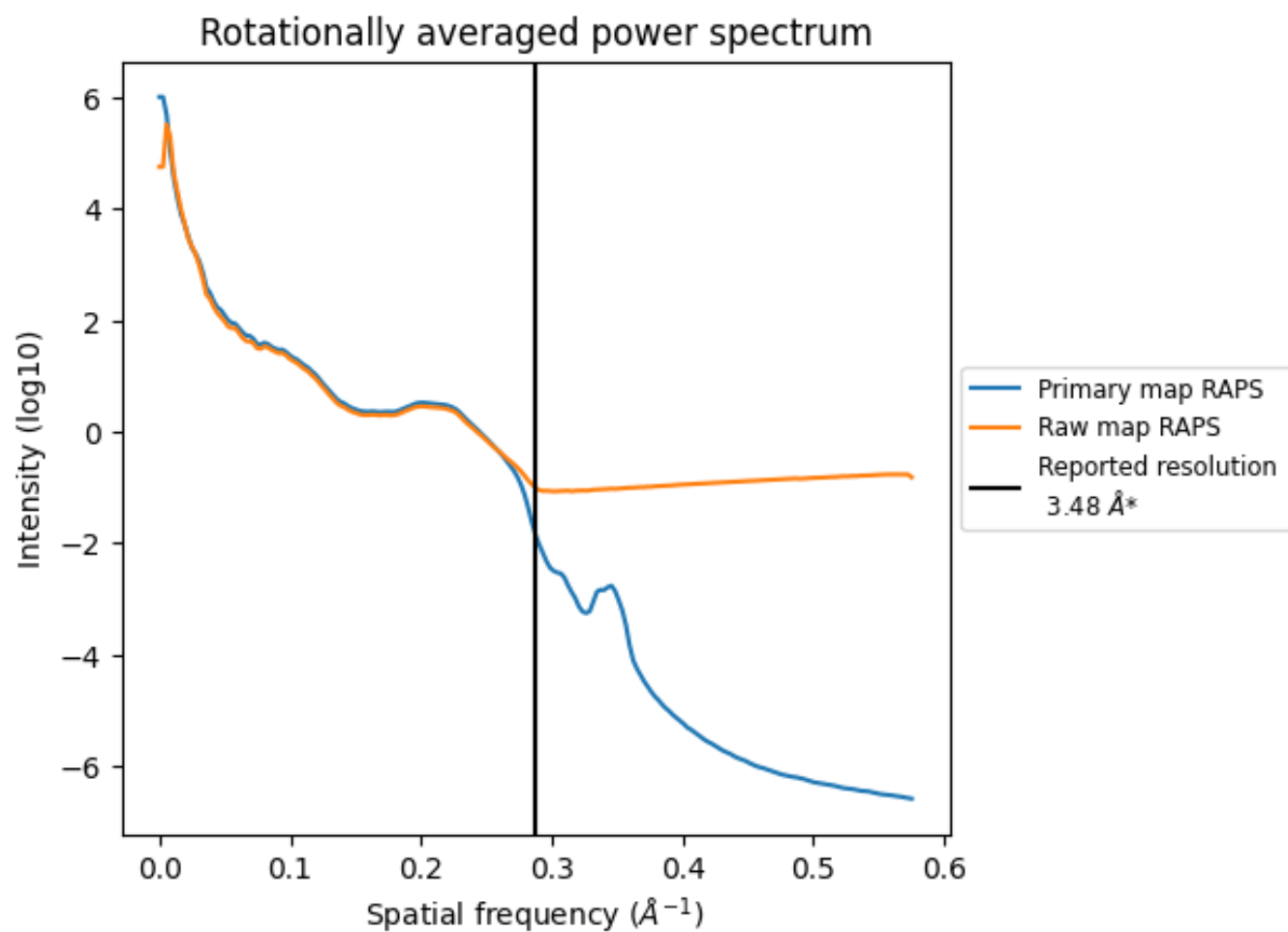
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 131 nm^3 ; this corresponds to an approximate mass of 118 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 ⓘ

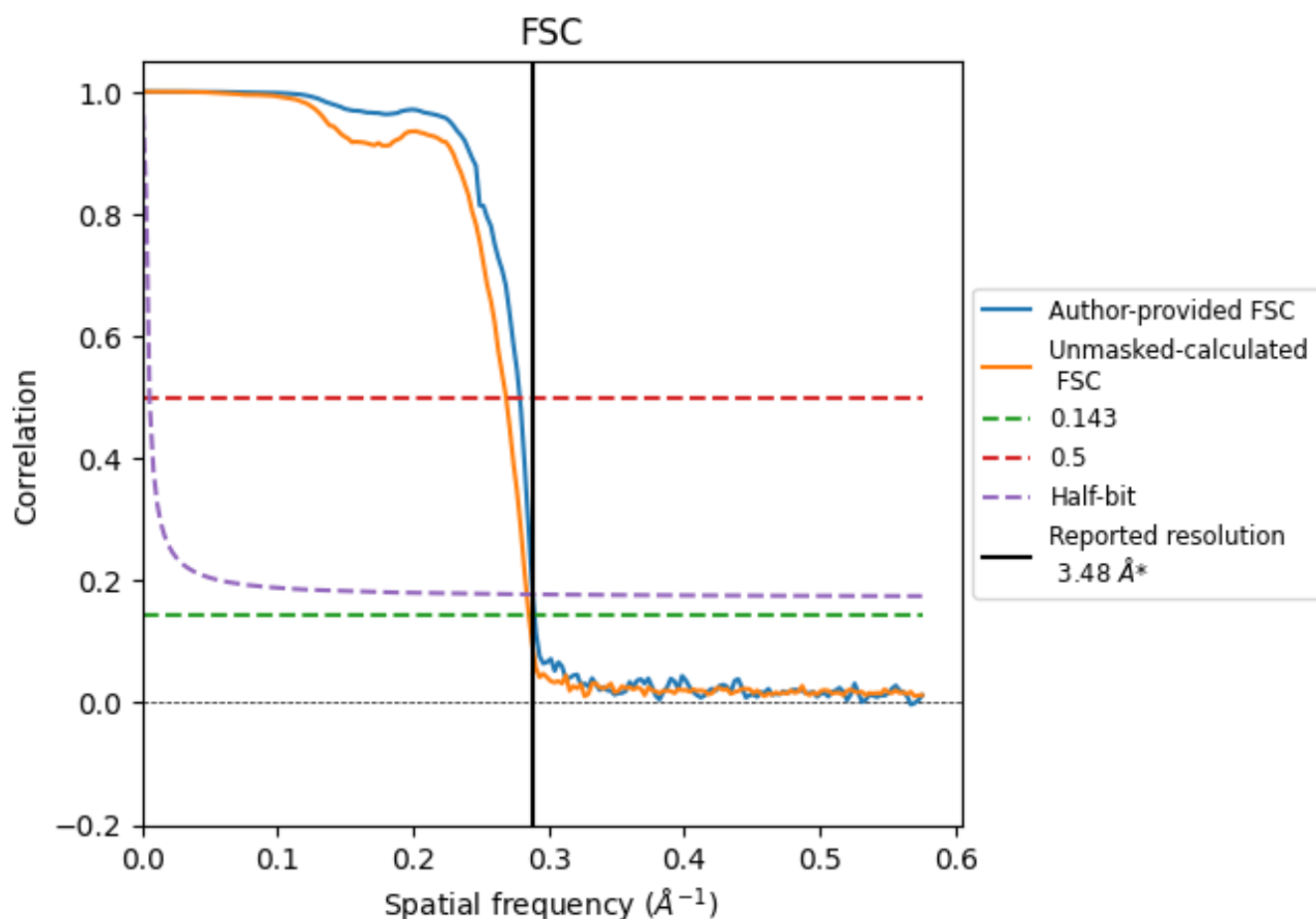


*Reported resolution corresponds to spatial frequency of 0.287 Å⁻¹

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 0.287 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

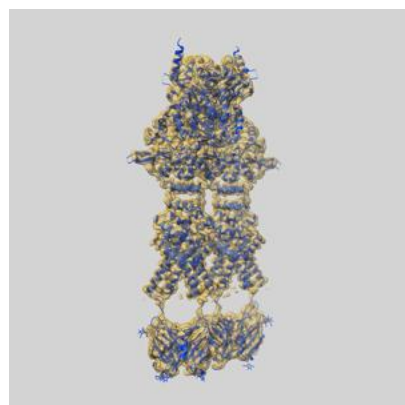
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.48	-	-
Author-provided FSC curve	3.46	3.59	3.47
Unmasked-calculated*	3.51	3.72	3.52

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

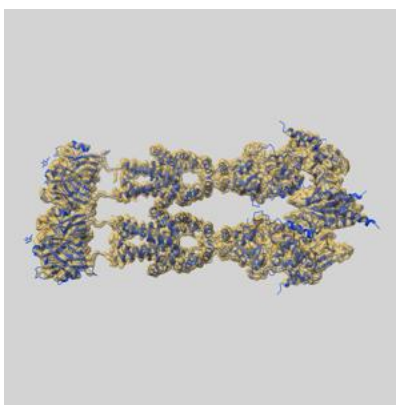
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-72694 and PDB model 9Y9D. Per-residue inclusion information can be found in [section 3](#) on [page 14](#).

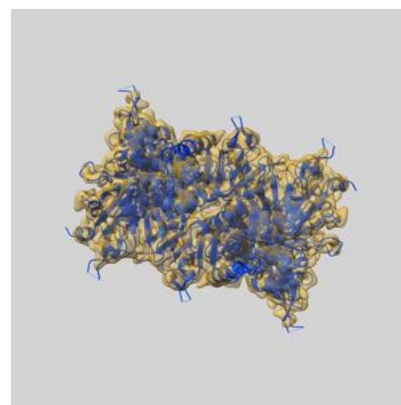
9.1 Map-model overlay [i](#)



X



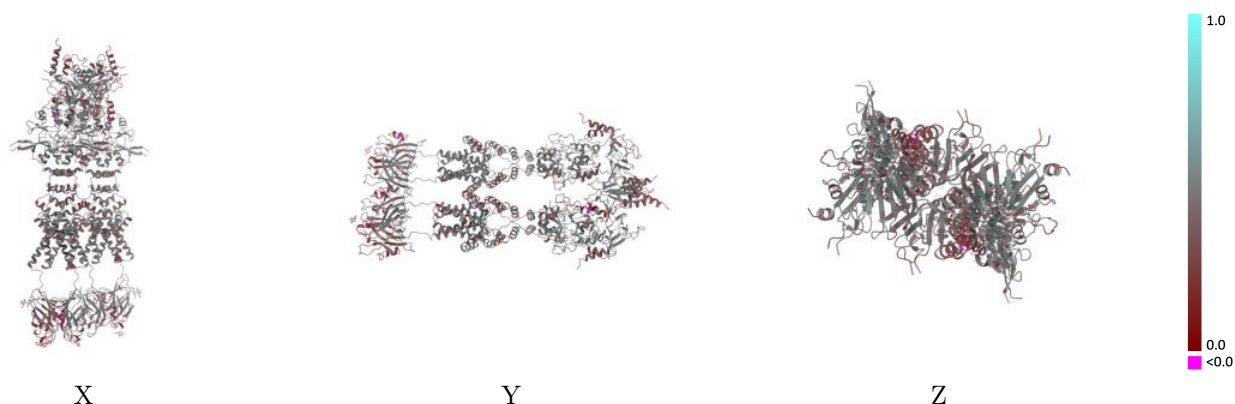
Y



Z

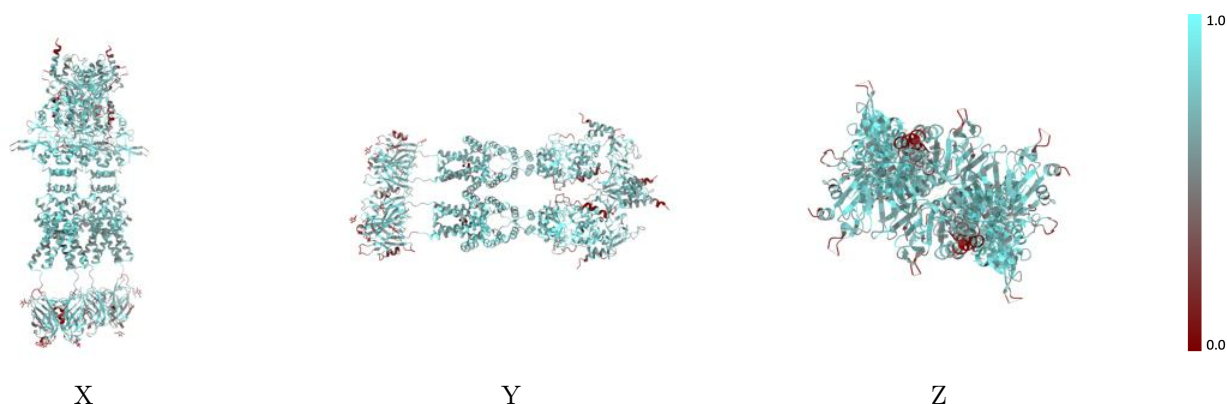
The images above show the 3D surface view of the map at the recommended contour level 0.00376 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



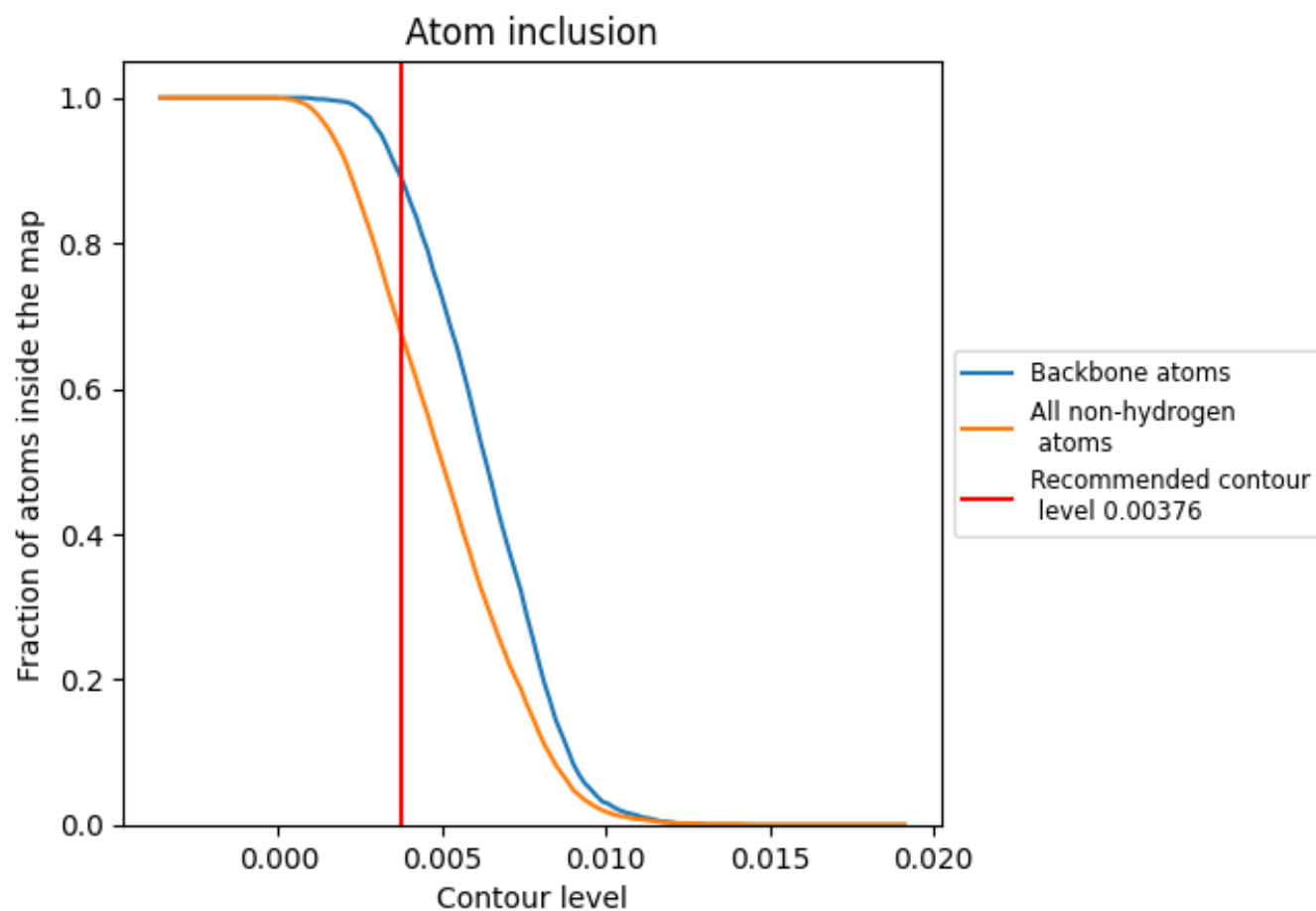
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.00376).

9.4 Atom inclusion [i](#)



At the recommended contour level, 89% of all backbone atoms, 68% 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 (0.00376) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6750	0.4080
A	0.7120	0.4280
B	0.6460	0.3900
C	0.7120	0.4280
D	0.6470	0.3920
E	0.4640	0.3340
F	0.2500	0.3670
G	0.2860	0.3010
H	0.3210	0.2790
I	0.4640	0.3540
J	0.2500	0.3530
K	0.2860	0.3100
L	0.3210	0.2780

