



Full wwPDB EM Validation Report ⓘ

Nov 26, 2025 – 05:39 PM EST

PDB ID : 9OFP / pdb_00009ofp
EMDB ID : EMD-70441
Title : HCoV-229E S2P bound by two DH1533 Fabs
Authors : Wrapp, D.
Deposited on : 2025-04-30
Resolution : 3.16 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:

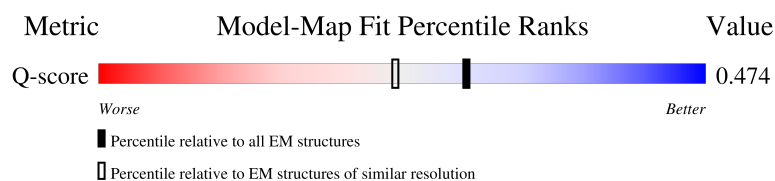
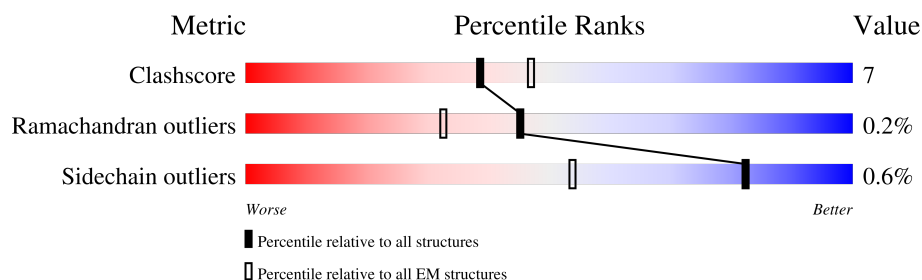
EMDB validation analysis : 0.0.1.dev129
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.46

1 Overall quality at a glance

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



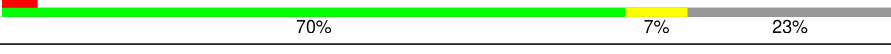

The reported resolution of this entry is 3.16 Å.

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.






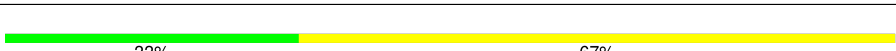
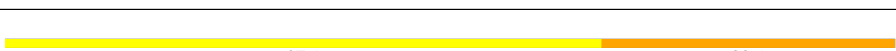
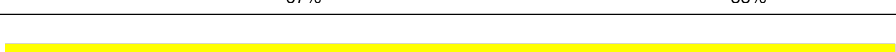
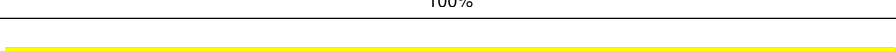
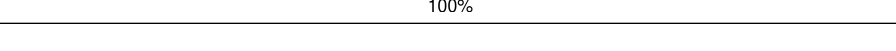

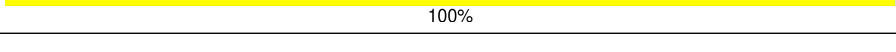
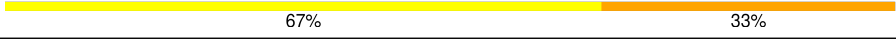
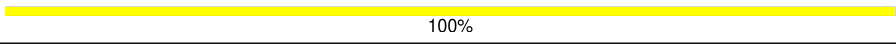
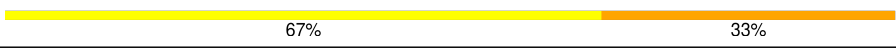
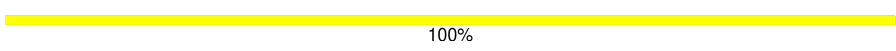
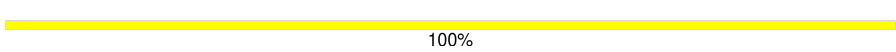

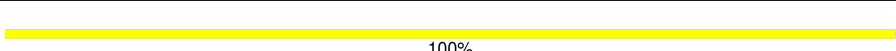
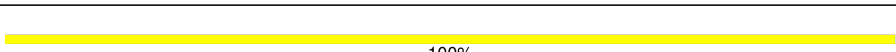
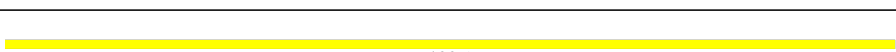
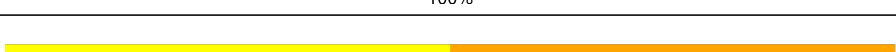

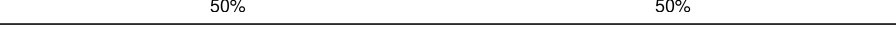
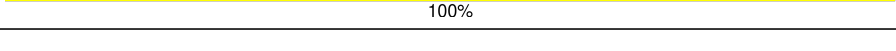
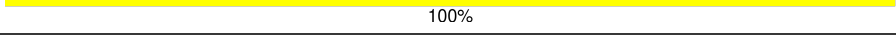
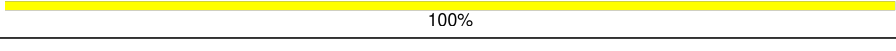
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	210492	15764	-
Ramachandran outliers	207382	16835	-
Sidechain outliers	206894	16415	-
Q-score	-	25397	14474 (2.66 - 3.66)

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	1134	
1	B	1134	
1	C	1134	
2	H	239	

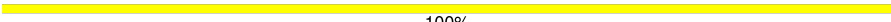
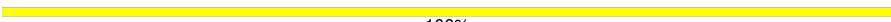
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Mol	Chain	Length	Quality of chain
2	M	239	
3	L	214	
3	N	214	
4	D	3	
4	E	3	
4	G	3	
4	P	3	
4	Q	3	
4	R	3	
4	T	3	
4	U	3	
4	V	3	
4	a	3	
4	c	3	
5	F	2	
5	I	2	
5	J	2	
5	K	2	
5	O	2	
5	S	2	
5	W	2	
5	X	2	
5	Y	2	
5	Z	2	
5	b	2	

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Mol	Chain	Length	Quality of chain
5	d	2	 100%
5	e	2	 100%

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
4	NAG	E	1	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 27163 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 Surface glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1012	Total	C	N	O	S	0	0
			7768	4930	1302	1496	40		
1	B	1012	Total	C	N	O	S	0	0
			7768	4930	1302	1496	40		
1	C	873	Total	C	N	O	S	0	0
			6694	4242	1116	1304	32		

There are 246 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	869	PRO	ILE	conflict	UNP A0A1B3TN99
A	870	PRO	ILE	conflict	UNP A0A1B3TN99
A	1055	GLY	-	expression tag	UNP A0A1B3TN99
A	1056	SER	-	expression tag	UNP A0A1B3TN99
A	1057	GLY	-	expression tag	UNP A0A1B3TN99
A	1058	TYR	-	expression tag	UNP A0A1B3TN99
A	1059	ILE	-	expression tag	UNP A0A1B3TN99
A	1060	PRO	-	expression tag	UNP A0A1B3TN99
A	1061	GLU	-	expression tag	UNP A0A1B3TN99
A	1062	ALA	-	expression tag	UNP A0A1B3TN99
A	1063	PRO	-	expression tag	UNP A0A1B3TN99
A	1064	ARG	-	expression tag	UNP A0A1B3TN99
A	1065	ASP	-	expression tag	UNP A0A1B3TN99
A	1066	GLY	-	expression tag	UNP A0A1B3TN99
A	1067	GLN	-	expression tag	UNP A0A1B3TN99
A	1068	ALA	-	expression tag	UNP A0A1B3TN99
A	1069	TYR	-	expression tag	UNP A0A1B3TN99
A	1070	VAL	-	expression tag	UNP A0A1B3TN99
A	1071	ARG	-	expression tag	UNP A0A1B3TN99
A	1072	LYS	-	expression tag	UNP A0A1B3TN99
A	1073	ASP	-	expression tag	UNP A0A1B3TN99
A	1074	GLY	-	expression tag	UNP A0A1B3TN99
A	1075	GLU	-	expression tag	UNP A0A1B3TN99
A	1076	TRP	-	expression tag	UNP A0A1B3TN99

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1077	VAL	-	expression tag	UNP A0A1B3TN99
A	1078	LEU	-	expression tag	UNP A0A1B3TN99
A	1079	LEU	-	expression tag	UNP A0A1B3TN99
A	1080	SER	-	expression tag	UNP A0A1B3TN99
A	1081	THR	-	expression tag	UNP A0A1B3TN99
A	1082	PHE	-	expression tag	UNP A0A1B3TN99
A	1083	LEU	-	expression tag	UNP A0A1B3TN99
A	1084	GLY	-	expression tag	UNP A0A1B3TN99
A	1085	ARG	-	expression tag	UNP A0A1B3TN99
A	1086	SER	-	expression tag	UNP A0A1B3TN99
A	1087	LEU	-	expression tag	UNP A0A1B3TN99
A	1088	GLU	-	expression tag	UNP A0A1B3TN99
A	1089	VAL	-	expression tag	UNP A0A1B3TN99
A	1090	LEU	-	expression tag	UNP A0A1B3TN99
A	1091	PHE	-	expression tag	UNP A0A1B3TN99
A	1092	GLN	-	expression tag	UNP A0A1B3TN99
A	1093	GLY	-	expression tag	UNP A0A1B3TN99
A	1094	PRO	-	expression tag	UNP A0A1B3TN99
A	1095	GLY	-	expression tag	UNP A0A1B3TN99
A	1096	HIS	-	expression tag	UNP A0A1B3TN99
A	1097	HIS	-	expression tag	UNP A0A1B3TN99
A	1098	HIS	-	expression tag	UNP A0A1B3TN99
A	1099	HIS	-	expression tag	UNP A0A1B3TN99
A	1100	HIS	-	expression tag	UNP A0A1B3TN99
A	1101	HIS	-	expression tag	UNP A0A1B3TN99
A	1102	HIS	-	expression tag	UNP A0A1B3TN99
A	1103	HIS	-	expression tag	UNP A0A1B3TN99
A	1104	SER	-	expression tag	UNP A0A1B3TN99
A	1105	ALA	-	expression tag	UNP A0A1B3TN99
A	1106	TRP	-	expression tag	UNP A0A1B3TN99
A	1107	SER	-	expression tag	UNP A0A1B3TN99
A	1108	HIS	-	expression tag	UNP A0A1B3TN99
A	1109	PRO	-	expression tag	UNP A0A1B3TN99
A	1110	GLN	-	expression tag	UNP A0A1B3TN99
A	1111	PHE	-	expression tag	UNP A0A1B3TN99
A	1112	GLU	-	expression tag	UNP A0A1B3TN99
A	1113	LYS	-	expression tag	UNP A0A1B3TN99
A	1114	GLY	-	expression tag	UNP A0A1B3TN99
A	1115	GLY	-	expression tag	UNP A0A1B3TN99
A	1116	GLY	-	expression tag	UNP A0A1B3TN99
A	1117	SER	-	expression tag	UNP A0A1B3TN99
A	1118	GLY	-	expression tag	UNP A0A1B3TN99

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1119	GLY	-	expression tag	UNP A0A1B3TN99
A	1120	GLY	-	expression tag	UNP A0A1B3TN99
A	1121	GLY	-	expression tag	UNP A0A1B3TN99
A	1122	SER	-	expression tag	UNP A0A1B3TN99
A	1123	GLY	-	expression tag	UNP A0A1B3TN99
A	1124	GLY	-	expression tag	UNP A0A1B3TN99
A	1125	SER	-	expression tag	UNP A0A1B3TN99
A	1126	ALA	-	expression tag	UNP A0A1B3TN99
A	1127	TRP	-	expression tag	UNP A0A1B3TN99
A	1128	SER	-	expression tag	UNP A0A1B3TN99
A	1129	HIS	-	expression tag	UNP A0A1B3TN99
A	1130	PRO	-	expression tag	UNP A0A1B3TN99
A	1131	GLN	-	expression tag	UNP A0A1B3TN99
A	1132	PHE	-	expression tag	UNP A0A1B3TN99
A	1133	GLU	-	expression tag	UNP A0A1B3TN99
A	1134	LYS	-	expression tag	UNP A0A1B3TN99
B	869	PRO	ILE	conflict	UNP A0A1B3TN99
B	870	PRO	ILE	conflict	UNP A0A1B3TN99
B	1055	GLY	-	expression tag	UNP A0A1B3TN99
B	1056	SER	-	expression tag	UNP A0A1B3TN99
B	1057	GLY	-	expression tag	UNP A0A1B3TN99
B	1058	TYR	-	expression tag	UNP A0A1B3TN99
B	1059	ILE	-	expression tag	UNP A0A1B3TN99
B	1060	PRO	-	expression tag	UNP A0A1B3TN99
B	1061	GLU	-	expression tag	UNP A0A1B3TN99
B	1062	ALA	-	expression tag	UNP A0A1B3TN99
B	1063	PRO	-	expression tag	UNP A0A1B3TN99
B	1064	ARG	-	expression tag	UNP A0A1B3TN99
B	1065	ASP	-	expression tag	UNP A0A1B3TN99
B	1066	GLY	-	expression tag	UNP A0A1B3TN99
B	1067	GLN	-	expression tag	UNP A0A1B3TN99
B	1068	ALA	-	expression tag	UNP A0A1B3TN99
B	1069	TYR	-	expression tag	UNP A0A1B3TN99
B	1070	VAL	-	expression tag	UNP A0A1B3TN99
B	1071	ARG	-	expression tag	UNP A0A1B3TN99
B	1072	LYS	-	expression tag	UNP A0A1B3TN99
B	1073	ASP	-	expression tag	UNP A0A1B3TN99
B	1074	GLY	-	expression tag	UNP A0A1B3TN99
B	1075	GLU	-	expression tag	UNP A0A1B3TN99
B	1076	TRP	-	expression tag	UNP A0A1B3TN99
B	1077	VAL	-	expression tag	UNP A0A1B3TN99
B	1078	LEU	-	expression tag	UNP A0A1B3TN99

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1079	LEU	-	expression tag	UNP A0A1B3TN99
B	1080	SER	-	expression tag	UNP A0A1B3TN99
B	1081	THR	-	expression tag	UNP A0A1B3TN99
B	1082	PHE	-	expression tag	UNP A0A1B3TN99
B	1083	LEU	-	expression tag	UNP A0A1B3TN99
B	1084	GLY	-	expression tag	UNP A0A1B3TN99
B	1085	ARG	-	expression tag	UNP A0A1B3TN99
B	1086	SER	-	expression tag	UNP A0A1B3TN99
B	1087	LEU	-	expression tag	UNP A0A1B3TN99
B	1088	GLU	-	expression tag	UNP A0A1B3TN99
B	1089	VAL	-	expression tag	UNP A0A1B3TN99
B	1090	LEU	-	expression tag	UNP A0A1B3TN99
B	1091	PHE	-	expression tag	UNP A0A1B3TN99
B	1092	GLN	-	expression tag	UNP A0A1B3TN99
B	1093	GLY	-	expression tag	UNP A0A1B3TN99
B	1094	PRO	-	expression tag	UNP A0A1B3TN99
B	1095	GLY	-	expression tag	UNP A0A1B3TN99
B	1096	HIS	-	expression tag	UNP A0A1B3TN99
B	1097	HIS	-	expression tag	UNP A0A1B3TN99
B	1098	HIS	-	expression tag	UNP A0A1B3TN99
B	1099	HIS	-	expression tag	UNP A0A1B3TN99
B	1100	HIS	-	expression tag	UNP A0A1B3TN99
B	1101	HIS	-	expression tag	UNP A0A1B3TN99
B	1102	HIS	-	expression tag	UNP A0A1B3TN99
B	1103	HIS	-	expression tag	UNP A0A1B3TN99
B	1104	SER	-	expression tag	UNP A0A1B3TN99
B	1105	ALA	-	expression tag	UNP A0A1B3TN99
B	1106	TRP	-	expression tag	UNP A0A1B3TN99
B	1107	SER	-	expression tag	UNP A0A1B3TN99
B	1108	HIS	-	expression tag	UNP A0A1B3TN99
B	1109	PRO	-	expression tag	UNP A0A1B3TN99
B	1110	GLN	-	expression tag	UNP A0A1B3TN99
B	1111	PHE	-	expression tag	UNP A0A1B3TN99
B	1112	GLU	-	expression tag	UNP A0A1B3TN99
B	1113	LYS	-	expression tag	UNP A0A1B3TN99
B	1114	GLY	-	expression tag	UNP A0A1B3TN99
B	1115	GLY	-	expression tag	UNP A0A1B3TN99
B	1116	GLY	-	expression tag	UNP A0A1B3TN99
B	1117	SER	-	expression tag	UNP A0A1B3TN99
B	1118	GLY	-	expression tag	UNP A0A1B3TN99
B	1119	GLY	-	expression tag	UNP A0A1B3TN99
B	1120	GLY	-	expression tag	UNP A0A1B3TN99

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1121	GLY	-	expression tag	UNP A0A1B3TN99
B	1122	SER	-	expression tag	UNP A0A1B3TN99
B	1123	GLY	-	expression tag	UNP A0A1B3TN99
B	1124	GLY	-	expression tag	UNP A0A1B3TN99
B	1125	SER	-	expression tag	UNP A0A1B3TN99
B	1126	ALA	-	expression tag	UNP A0A1B3TN99
B	1127	TRP	-	expression tag	UNP A0A1B3TN99
B	1128	SER	-	expression tag	UNP A0A1B3TN99
B	1129	HIS	-	expression tag	UNP A0A1B3TN99
B	1130	PRO	-	expression tag	UNP A0A1B3TN99
B	1131	GLN	-	expression tag	UNP A0A1B3TN99
B	1132	PHE	-	expression tag	UNP A0A1B3TN99
B	1133	GLU	-	expression tag	UNP A0A1B3TN99
B	1134	LYS	-	expression tag	UNP A0A1B3TN99
C	869	PRO	ILE	conflict	UNP A0A1B3TN99
C	870	PRO	ILE	conflict	UNP A0A1B3TN99
C	1055	GLY	-	expression tag	UNP A0A1B3TN99
C	1056	SER	-	expression tag	UNP A0A1B3TN99
C	1057	GLY	-	expression tag	UNP A0A1B3TN99
C	1058	TYR	-	expression tag	UNP A0A1B3TN99
C	1059	ILE	-	expression tag	UNP A0A1B3TN99
C	1060	PRO	-	expression tag	UNP A0A1B3TN99
C	1061	GLU	-	expression tag	UNP A0A1B3TN99
C	1062	ALA	-	expression tag	UNP A0A1B3TN99
C	1063	PRO	-	expression tag	UNP A0A1B3TN99
C	1064	ARG	-	expression tag	UNP A0A1B3TN99
C	1065	ASP	-	expression tag	UNP A0A1B3TN99
C	1066	GLY	-	expression tag	UNP A0A1B3TN99
C	1067	GLN	-	expression tag	UNP A0A1B3TN99
C	1068	ALA	-	expression tag	UNP A0A1B3TN99
C	1069	TYR	-	expression tag	UNP A0A1B3TN99
C	1070	VAL	-	expression tag	UNP A0A1B3TN99
C	1071	ARG	-	expression tag	UNP A0A1B3TN99
C	1072	LYS	-	expression tag	UNP A0A1B3TN99
C	1073	ASP	-	expression tag	UNP A0A1B3TN99
C	1074	GLY	-	expression tag	UNP A0A1B3TN99
C	1075	GLU	-	expression tag	UNP A0A1B3TN99
C	1076	TRP	-	expression tag	UNP A0A1B3TN99
C	1077	VAL	-	expression tag	UNP A0A1B3TN99
C	1078	LEU	-	expression tag	UNP A0A1B3TN99
C	1079	LEU	-	expression tag	UNP A0A1B3TN99
C	1080	SER	-	expression tag	UNP A0A1B3TN99

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1081	THR	-	expression tag	UNP A0A1B3TN99
C	1082	PHE	-	expression tag	UNP A0A1B3TN99
C	1083	LEU	-	expression tag	UNP A0A1B3TN99
C	1084	GLY	-	expression tag	UNP A0A1B3TN99
C	1085	ARG	-	expression tag	UNP A0A1B3TN99
C	1086	SER	-	expression tag	UNP A0A1B3TN99
C	1087	LEU	-	expression tag	UNP A0A1B3TN99
C	1088	GLU	-	expression tag	UNP A0A1B3TN99
C	1089	VAL	-	expression tag	UNP A0A1B3TN99
C	1090	LEU	-	expression tag	UNP A0A1B3TN99
C	1091	PHE	-	expression tag	UNP A0A1B3TN99
C	1092	GLN	-	expression tag	UNP A0A1B3TN99
C	1093	GLY	-	expression tag	UNP A0A1B3TN99
C	1094	PRO	-	expression tag	UNP A0A1B3TN99
C	1095	GLY	-	expression tag	UNP A0A1B3TN99
C	1096	HIS	-	expression tag	UNP A0A1B3TN99
C	1097	HIS	-	expression tag	UNP A0A1B3TN99
C	1098	HIS	-	expression tag	UNP A0A1B3TN99
C	1099	HIS	-	expression tag	UNP A0A1B3TN99
C	1100	HIS	-	expression tag	UNP A0A1B3TN99
C	1101	HIS	-	expression tag	UNP A0A1B3TN99
C	1102	HIS	-	expression tag	UNP A0A1B3TN99
C	1103	HIS	-	expression tag	UNP A0A1B3TN99
C	1104	SER	-	expression tag	UNP A0A1B3TN99
C	1105	ALA	-	expression tag	UNP A0A1B3TN99
C	1106	TRP	-	expression tag	UNP A0A1B3TN99
C	1107	SER	-	expression tag	UNP A0A1B3TN99
C	1108	HIS	-	expression tag	UNP A0A1B3TN99
C	1109	PRO	-	expression tag	UNP A0A1B3TN99
C	1110	GLN	-	expression tag	UNP A0A1B3TN99
C	1111	PHE	-	expression tag	UNP A0A1B3TN99
C	1112	GLU	-	expression tag	UNP A0A1B3TN99
C	1113	LYS	-	expression tag	UNP A0A1B3TN99
C	1114	GLY	-	expression tag	UNP A0A1B3TN99
C	1115	GLY	-	expression tag	UNP A0A1B3TN99
C	1116	GLY	-	expression tag	UNP A0A1B3TN99
C	1117	SER	-	expression tag	UNP A0A1B3TN99
C	1118	GLY	-	expression tag	UNP A0A1B3TN99
C	1119	GLY	-	expression tag	UNP A0A1B3TN99
C	1120	GLY	-	expression tag	UNP A0A1B3TN99
C	1121	GLY	-	expression tag	UNP A0A1B3TN99
C	1122	SER	-	expression tag	UNP A0A1B3TN99

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1123	GLY	-	expression tag	UNP A0A1B3TN99
C	1124	GLY	-	expression tag	UNP A0A1B3TN99
C	1125	SER	-	expression tag	UNP A0A1B3TN99
C	1126	ALA	-	expression tag	UNP A0A1B3TN99
C	1127	TRP	-	expression tag	UNP A0A1B3TN99
C	1128	SER	-	expression tag	UNP A0A1B3TN99
C	1129	HIS	-	expression tag	UNP A0A1B3TN99
C	1130	PRO	-	expression tag	UNP A0A1B3TN99
C	1131	GLN	-	expression tag	UNP A0A1B3TN99
C	1132	PHE	-	expression tag	UNP A0A1B3TN99
C	1133	GLU	-	expression tag	UNP A0A1B3TN99
C	1134	LYS	-	expression tag	UNP A0A1B3TN99

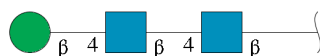
- Molecule 2 is a protein called DH1533 heavy chain.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	H	124	Total	C	N	O	S	
			944	598	159	183	4	
2	M	124	Total	C	N	O	S	
			944	598	159	183	4	

- Molecule 3 is a protein called DH1533 light chain.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	L	106	Total	C	N	O	S	
			811	511	130	167	3	
3	N	106	Total	C	N	O	S	
			811	511	130	167	3	

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



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

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Mol	Chain	Residues	Atoms				AltConf	Trace
4	G	3	Total	C	N	O	0	0
			39	22	2	15		
4	P	3	Total	C	N	O	0	0
			39	22	2	15		
4	Q	3	Total	C	N	O	0	0
			39	22	2	15		
4	R	3	Total	C	N	O	0	0
			39	22	2	15		
4	T	3	Total	C	N	O	0	0
			39	22	2	15		
4	U	3	Total	C	N	O	0	0
			39	22	2	15		
4	V	3	Total	C	N	O	0	0
			39	22	2	15		
4	a	3	Total	C	N	O	0	0
			39	22	2	15		
4	c	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



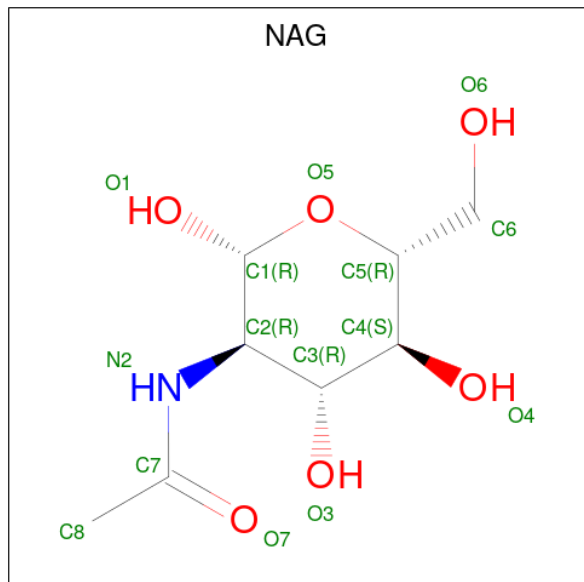
Mol	Chain	Residues	Atoms				AltConf	Trace
5	F	2	Total	C	N	O	0	0
			28	16	2	10		
5	I	2	Total	C	N	O	0	0
			28	16	2	10		
5	J	2	Total	C	N	O	0	0
			28	16	2	10		
5	K	2	Total	C	N	O	0	0
			28	16	2	10		
5	O	2	Total	C	N	O	0	0
			28	16	2	10		
5	S	2	Total	C	N	O	0	0
			28	16	2	10		
5	W	2	Total	C	N	O	0	0
			28	16	2	10		
5	X	2	Total	C	N	O	0	0
			28	16	2	10		

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Mol	Chain	Residues	Atoms				AltConf	Trace
5	Y	2	Total	C	N	O	0	0
			28	16	2	10		
5	Z	2	Total	C	N	O	0	0
			28	16	2	10		
5	b	2	Total	C	N	O	0	0
			28	16	2	10		
5	d	2	Total	C	N	O	0	0
			28	16	2	10		
5	e	2	Total	C	N	O	0	0
			28	16	2	10		

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



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

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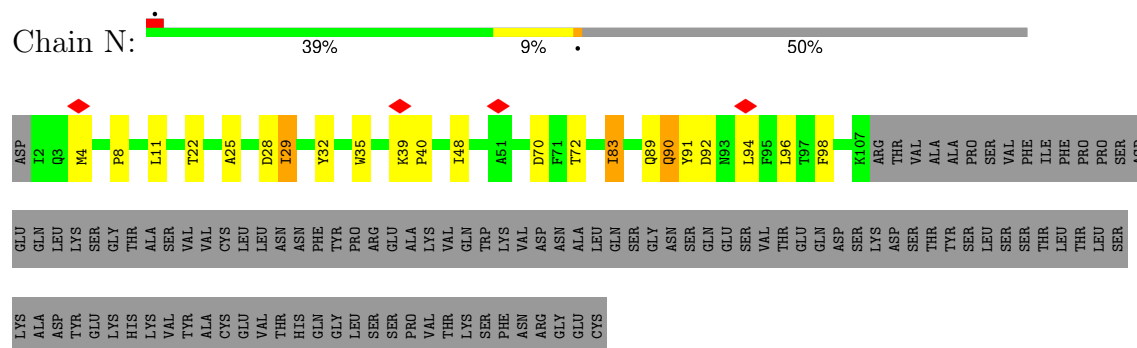
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Mol	Chain	Residues	Atoms				AltConf
6	A	1	Total 14	C 8	N 1	O 5	0
6	A	1	Total 14	C 8	N 1	O 5	0
6	A	1	Total 14	C 8	N 1	O 5	0
6	A	1	Total 14	C 8	N 1	O 5	0
6	A	1	Total 14	C 8	N 1	O 5	0
6	A	1	Total 14	C 8	N 1	O 5	0
6	A	1	Total 14	C 8	N 1	O 5	0
6	A	1	Total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

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



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



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



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



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



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



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

Chain R:  100%

MAG1
MAG2
BMA3

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

Chain T:  67% 33%

MAG1
MAG2
BMA3

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

Chain U:  100%

MAG1
MAG2
BMA3

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

Chain V:  67% 33%

MAG1
MAG2
BMA3

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

Chain a:  100%

MAG1
MAG2
BMA3

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

Chain c:  100%

MAG1
MAG2
BMA3

- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  50% 50%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  100%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  100%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  100%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  50% 50%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S:  50% 50%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain W:  100%




- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain X:  100%

MAG1
MAG2

- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Y:  100%

MAG1
MAG2

- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z:  100%

MAG1
MAG2

- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain b:  50%  50%

MAG1
MAG2

- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain d:  100%

MAG1
MAG2

- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain e:  100%

MAG1
MAG2

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	122657	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION; Patch CTF Estimation	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	49.8	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.447	Depositor
Minimum map value	-0.005	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.036	Depositor
Recommended contour level	0.34	Depositor
Map size (Å)	466.56003, 466.56003, 466.56003	wwPDB
Map dimensions	432, 432, 432	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.16	0/7928	0.45	1/10805 (0.0%)
1	B	0.17	0/7928	0.43	2/10805 (0.0%)
1	C	0.16	0/6822	0.45	5/9299 (0.1%)
2	H	0.16	0/969	0.48	0/1317
2	M	0.17	0/969	0.46	0/1317
3	L	0.15	0/827	0.50	2/1124 (0.2%)
3	N	0.24	0/827	0.63	2/1124 (0.2%)
All	All	0.17	0/26270	0.46	12/35791 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	6
1	C	0	4
2	M	0	3
All	All	0	15

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	38	ASN	N-CA-C	-9.35	100.97	112.38
1	C	36	SER	N-CA-C	-6.75	105.49	113.38
1	A	98	ASN	N-CA-C	-5.99	105.83	113.02
1	C	37	GLU	N-CA-C	-5.99	105.89	113.43
3	N	29	ILE	CA-C-N	5.61	132.25	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	N	29	ILE	C-N-CA	5.61	132.25	121.54
1	B	462	ASN	N-CA-C	-5.42	106.14	114.16
1	B	462	ASN	CA-CB-CG	5.41	118.01	112.60
1	C	38	ASN	CB-CA-C	5.29	121.48	110.32
3	L	29	ILE	CA-C-N	5.26	131.59	121.54
3	L	29	ILE	C-N-CA	5.26	131.59	121.54
1	C	39	VAL	N-CA-C	-5.19	98.55	109.34

There are no chirality outliers.

All (15) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	129	ARG	Sidechain
1	A	878	ARG	Sidechain
1	B	128	ARG	Sidechain
1	B	260	ARG	Sidechain
1	B	262	ARG	Sidechain
1	B	565	ARG	Sidechain
1	B	790	ARG	Sidechain
1	B	866	ARG	Sidechain
1	C	129	ARG	Sidechain
1	C	459	ARG	Sidechain
1	C	878	ARG	Sidechain
1	C	905	ARG	Sidechain
2	M	38	ARG	Sidechain
2	M	66	ARG	Sidechain
2	M	71	ARG	Sidechain

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	7768	0	7601	97	0
1	B	7768	0	7602	94	0
1	C	6694	0	6553	77	0
2	H	944	0	898	54	0
2	M	944	0	898	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	811	0	787	31	0
3	N	811	0	787	17	0
4	D	39	0	34	0	0
4	E	39	0	34	7	0
4	G	39	0	34	0	0
4	P	39	0	34	0	0
4	Q	39	0	34	0	0
4	R	39	0	34	0	0
4	T	39	0	34	1	0
4	U	39	0	34	0	0
4	V	39	0	34	1	0
4	a	39	0	34	0	0
4	c	39	0	34	0	0
5	F	28	0	25	0	0
5	I	28	0	25	0	0
5	J	28	0	25	0	0
5	K	28	0	25	0	0
5	O	28	0	25	1	0
5	S	28	0	25	0	0
5	W	28	0	25	0	0
5	X	28	0	25	0	0
5	Y	28	0	25	0	0
5	Z	28	0	25	0	0
5	b	28	0	25	1	0
5	d	28	0	25	0	0
5	e	28	0	25	0	0
6	A	224	0	208	1	0
6	B	196	0	182	4	0
6	C	210	0	195	4	0
All	All	27163	0	26410	379	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 7.

All (379) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:486:ASN:ND2	6:B:1206:NAG:C1	1.73	1.52
1:B:486:ASN:HD22	6:B:1206:NAG:C1	1.21	1.46
2:H:20:LEU:HD23	2:H:80:LEU:CD2	1.50	1.39
2:H:20:LEU:CD2	2:H:80:LEU:HD23	1.68	1.23

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:39:GLN:HE22	3:L:38:GLN:NE2	1.36	1.23
2:H:39:GLN:HG2	2:H:45:LEU:CD1	1.66	1.22
2:H:39:GLN:HG2	2:H:45:LEU:HD11	1.19	1.16
2:H:39:GLN:NE2	3:L:38:GLN:NE2	1.98	1.10
1:A:186:LEU:HD12	1:A:187:PRO:CD	1.83	1.09
2:H:39:GLN:CG	2:H:45:LEU:HD11	1.81	1.09
2:H:20:LEU:HD21	2:H:90:TYR:CD2	1.89	1.07
2:H:39:GLN:CG	2:H:45:LEU:CD1	2.34	1.06
1:B:759:LEU:HD11	1:B:765:ILE:HG23	1.39	1.02
1:B:1034:VAL:HG11	1:C:1034:VAL:HG11	1.42	0.99
1:A:186:LEU:HD12	1:A:187:PRO:HD3	1.46	0.94
1:C:134:PHE:CD1	1:C:217:VAL:HG12	2.02	0.94
1:A:186:LEU:HD21	1:A:190:VAL:HG11	1.50	0.92
2:H:20:LEU:HD21	2:H:90:TYR:HD2	1.26	0.92
2:M:52:ILE:HD13	2:M:100(D):SER:HA	1.52	0.91
1:A:100:THR:OG1	4:E:1:NAG:C7	2.19	0.90
1:A:186:LEU:HD12	1:A:187:PRO:HD2	1.52	0.89
1:A:100:THR:OG1	4:E:1:NAG:H83	1.71	0.89
1:A:186:LEU:HD11	1:A:203:ILE:HG21	1.54	0.88
1:B:486:ASN:HD21	6:B:1206:NAG:C1	1.86	0.88
2:H:39:GLN:HG2	2:H:45:LEU:HD13	1.55	0.87
1:A:186:LEU:CD1	1:A:187:PRO:HD2	2.05	0.85
1:C:39:VAL:HG21	1:C:84:SER:HB2	1.61	0.83
2:M:52:ILE:CD1	2:M:100(D):SER:HA	2.08	0.83
1:A:100:THR:OG1	4:E:1:NAG:C8	2.29	0.81
2:H:20:LEU:CD2	2:H:90:TYR:CD2	2.65	0.80
1:B:814:ILE:HG22	1:B:814:ILE:O	1.79	0.80
1:B:133:LEU:HD23	1:B:142:VAL:HG23	1.64	0.80
1:B:339:CYS:HB3	1:B:384:CYS:HA	1.62	0.79
1:B:759:LEU:CD1	1:B:765:ILE:HG23	2.11	0.79
2:H:20:LEU:CD2	2:H:90:TYR:CE2	2.66	0.78
1:B:759:LEU:HD11	1:B:765:ILE:CG2	2.12	0.77
2:H:20:LEU:HD23	2:H:80:LEU:HD23	0.80	0.77
1:A:186:LEU:CD1	1:A:203:ILE:HG21	2.13	0.77
1:B:765:ILE:HD12	1:B:769:LEU:HB2	1.66	0.77
1:B:207:ARG:HG2	1:B:207:ARG:HH11	1.48	0.76
1:A:128:ARG:HB2	1:A:129:ARG:HD3	1.69	0.75
1:C:565:ARG:NH2	1:C:819:GLN:OE1	2.20	0.75
1:A:1010:GLU:OE2	1:B:989:ARG:NE	2.17	0.73
1:C:772:GLN:HE22	1:C:786:GLN:HB3	1.52	0.73
2:H:39:GLN:HE22	3:L:38:GLN:CD	1.98	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:20:LEU:HD23	2:H:80:LEU:HD21	1.66	0.72
1:C:147:ASN:HB3	1:C:166:PHE:HA	1.72	0.71
2:M:52:ILE:HD12	2:M:100(E):GLY:H	1.54	0.71
1:A:100:THR:OG1	4:E:1:NAG:N2	2.24	0.70
1:B:173:THR:HG22	1:B:175:GLY:H	1.55	0.70
1:B:616:ARG:NH2	1:B:620:LEU:HD21	2.06	0.70
3:L:5:THR:OG1	3:L:24:GLN:HB2	1.91	0.70
1:B:1034:VAL:HG11	1:C:1034:VAL:CG1	2.21	0.70
2:H:20:LEU:CD2	2:H:80:LEU:CD2	2.46	0.70
2:H:39:GLN:NE2	3:L:38:GLN:HE21	1.88	0.69
1:B:324:ILE:HG13	1:B:324:ILE:O	1.93	0.69
2:M:38:ARG:HH21	2:M:88:ALA:HB2	1.58	0.68
1:A:186:LEU:HD13	1:A:203:ILE:HD13	1.75	0.68
1:A:577:THR:HG23	1:A:577:THR:O	1.93	0.68
1:B:76:PRO:HB2	1:B:116:VAL:HG11	1.75	0.68
2:H:39:GLN:HG3	2:H:45:LEU:CD1	2.23	0.68
1:B:564:PRO:HG3	1:C:742:ARG:HG2	1.74	0.67
2:M:52:ILE:CD1	2:M:100(E):GLY:H	2.08	0.67
3:L:83:ILE:O	3:L:83:ILE:HG13	1.93	0.66
3:L:89:GLN:HE21	3:L:96:LEU:HD13	1.60	0.66
1:A:39:VAL:HG22	1:A:106:LYS:HZ1	1.59	0.65
1:C:32:CYS:N	1:C:38:ASN:HB3	2.11	0.65
2:H:39:GLN:CD	3:L:38:GLN:NE2	2.54	0.65
1:C:134:PHE:HD1	1:C:217:VAL:HG12	1.61	0.65
1:B:133:LEU:HB2	1:B:218:ASN:HB3	1.79	0.64
1:B:485:THR:OG1	6:B:1206:NAG:H83	1.97	0.64
1:A:130:GLY:HA3	1:A:221:VAL:HA	1.79	0.64
1:A:186:LEU:HD21	1:A:190:VAL:CG1	2.25	0.64
2:H:20:LEU:HD21	2:H:90:TYR:CE2	2.31	0.64
1:B:738:ALA:HB1	1:B:743:MET:HE2	1.80	0.64
2:M:35:SER:HB2	2:M:100(J):PHE:CE2	2.34	0.63
1:B:422:VAL:HG13	1:B:423:ILE:HG23	1.81	0.63
2:H:20:LEU:HD22	2:H:90:TYR:HE2	1.65	0.62
2:H:20:LEU:HD22	2:H:90:TYR:CE2	2.33	0.62
1:A:186:LEU:CD1	1:A:187:PRO:CD	2.62	0.62
1:B:133:LEU:CD2	1:B:142:VAL:HG23	2.28	0.62
2:M:35:SER:HB2	2:M:100(J):PHE:HE2	1.64	0.62
1:A:964:LEU:HD23	1:A:1017:VAL:HB	1.82	0.62
2:H:20:LEU:CD2	2:H:90:TYR:HE2	2.13	0.61
2:M:100(F):TRP:CD1	2:M:100(G):TYR:H	2.19	0.61
1:B:437:MET:CE	1:B:483:ASP:HB2	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:15:VAL:HA	3:L:78:LEU:HD22	1.82	0.61
3:N:83:ILE:O	3:N:83:ILE:HG13	2.00	0.60
3:N:39:LYS:HD3	3:N:40:PRO:HD2	1.83	0.60
2:M:52(A):GLY:HA2	2:M:71:ARG:HH12	1.66	0.60
1:A:39:VAL:HG11	1:A:84:SER:HB2	1.84	0.60
2:M:59:TYR:HE1	2:M:69:ILE:HG22	1.66	0.60
1:C:126:ASN:HA	1:C:190:VAL:HG22	1.84	0.60
1:B:814:ILE:O	1:B:814:ILE:CG2	2.50	0.60
1:C:94:PHE:CD1	1:C:218:ASN:HB3	2.37	0.60
1:B:302:VAL:HG22	1:B:348:GLN:HE21	1.67	0.59
1:A:97:PHE:CE2	1:A:197:ARG:HA	2.37	0.59
1:C:134:PHE:CD1	1:C:217:VAL:CG1	2.83	0.59
1:A:597:ILE:HG12	1:A:942:GLY:HA2	1.85	0.59
1:C:37:GLU:HA	1:C:40:PHE:HA	1.84	0.59
2:M:15:GLY:HA2	2:M:82(B):SER:HA	1.84	0.59
1:B:528:MET:HE3	1:B:529:PRO:HD2	1.85	0.58
2:M:58:TYR:HB3	3:N:94:LEU:HD21	1.84	0.58
1:C:39:VAL:HG22	1:C:106:LYS:HE2	1.85	0.58
2:M:18:LEU:HD11	2:M:82(C):LEU:HD21	1.85	0.58
1:A:297:LYS:HB2	1:A:342:THR:HB	1.86	0.58
1:B:207:ARG:HH11	1:B:207:ARG:CG	2.16	0.58
2:H:46:GLU:HA	3:L:98:PHE:CE1	2.39	0.58
1:A:97:PHE:CZ	1:A:197:ARG:HA	2.38	0.57
1:B:850:ARG:HG3	1:B:850:ARG:HH11	1.69	0.57
1:C:37:GLU:C	1:C:39:VAL:N	2.56	0.57
3:N:22:THR:HG22	3:N:72:THR:HG22	1.84	0.57
1:B:188:LYS:HG3	1:B:189:THR:N	2.19	0.57
1:B:304:TYR:HB2	1:B:327:THR:OG1	2.04	0.57
1:B:1009:ILE:HD12	1:C:784:VAL:CG2	2.35	0.57
2:H:75:LYS:O	2:H:75:LYS:HG2	2.05	0.57
3:L:33:LEU:HD12	3:L:34:ASN:H	1.69	0.57
1:A:127:LEU:HD22	1:A:219:PHE:CZ	2.40	0.57
1:A:218:ASN:OD1	1:A:220:ASN:HB2	2.05	0.57
1:A:1010:GLU:HG2	1:B:1000:ILE:HD11	1.86	0.56
1:A:97:PHE:HE2	1:A:214:VAL:O	1.88	0.56
1:C:220:ASN:HB3	6:C:1204:NAG:O5	2.05	0.56
1:A:96:TYR:CD1	4:E:1:NAG:H81	2.41	0.56
1:B:528:MET:HE3	1:B:529:PRO:CD	2.36	0.56
3:N:96:LEU:HD23	3:N:96:LEU:H	1.70	0.56
1:A:39:VAL:HG22	1:A:106:LYS:NZ	2.19	0.56
2:H:36:TRP:CZ3	2:H:92:CYS:HB3	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:48:VAL:HG13	2:M:63:VAL:HG21	1.87	0.56
2:M:61:ASP:OD1	2:M:62:SER:N	2.37	0.56
1:B:324:ILE:HD12	1:B:410:ILE:O	2.06	0.56
1:C:786:GLN:O	1:C:789:GLN:HG3	2.06	0.56
2:H:20:LEU:CG	2:H:80:LEU:HD23	2.35	0.56
1:C:179:THR:HG22	1:C:180:SER:N	2.21	0.55
1:B:460:ILE:HD11	1:B:488:THR:HG21	1.87	0.55
1:B:1002:THR:HG22	1:B:1004:ALA:H	1.71	0.55
1:B:59:LEU:HB2	1:B:248:ILE:HD11	1.87	0.54
1:A:1036:LYS:HA	1:A:1036:LYS:HE2	1.89	0.54
1:B:786:GLN:HB3	1:B:789:GLN:HE22	1.72	0.54
1:C:111:ASN:HD22	6:C:1201:NAG:C7	2.21	0.54
2:H:39:GLN:HE21	2:H:45:LEU:HD11	1.72	0.54
1:B:1009:ILE:CD1	1:C:784:VAL:HG21	2.38	0.54
1:C:224:ALA:HA	1:C:227:THR:HG22	1.91	0.53
1:A:295:TYR:CE2	1:A:297:LYS:HD3	2.43	0.53
2:H:36:TRP:CD1	2:H:69:ILE:HD12	2.42	0.53
1:A:134:PHE:HD2	1:A:214:VAL:HG11	1.74	0.53
3:L:16:GLY:H	3:L:78:LEU:HB3	1.72	0.53
1:B:133:LEU:HD23	1:B:142:VAL:CG2	2.38	0.53
1:A:297:LYS:HA	1:A:297:LYS:HE3	1.91	0.53
1:B:188:LYS:HG3	1:B:189:THR:HG23	1.91	0.53
1:A:100:THR:CB	4:E:1:NAG:H83	2.39	0.52
1:B:156:HIS:CD2	1:B:157:ILE:HG12	2.43	0.52
1:B:145:CYS:HA	1:B:168:CYS:HA	1.91	0.52
1:A:953:THR:HG22	1:A:954:GLN:HG3	1.90	0.52
1:B:125:GLU:HG2	1:B:227:THR:CG2	2.39	0.52
1:B:296:HIS:O	1:C:156:HIS:HE1	1.92	0.52
1:A:1014:VAL:HG13	5:O:1:NAG:H3	1.91	0.52
1:B:188:LYS:HG3	1:B:189:THR:H	1.73	0.52
1:B:461:SER:OG	1:B:489:ILE:HB	2.10	0.52
1:B:125:GLU:HG2	1:B:227:THR:HG21	1.91	0.51
2:H:6:GLU:HA	2:H:22:CYS:HA	1.93	0.51
3:N:29:ILE:HG23	3:N:92:ASP:HB2	1.91	0.51
1:B:902:ARG:HG3	1:B:902:ARG:HH11	1.76	0.51
1:A:102:ARG:HD2	1:A:105:CYS:SG	2.51	0.51
1:A:381:GLY:HA2	1:A:422:VAL:HG12	1.93	0.51
2:H:46:GLU:HA	3:L:98:PHE:CZ	2.45	0.50
1:A:98:ASN:C	1:A:100:THR:H	2.18	0.50
3:L:33:LEU:HD11	3:L:88:CYS:HB2	1.94	0.50
1:B:464:THR:O	1:B:465:PHE:C	2.54	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:THR:O	4:E:1:NAG:H83	2.11	0.50
1:B:1009:ILE:HD12	1:C:784:VAL:HG21	1.94	0.50
1:C:39:VAL:O	1:C:106:LYS:CE	2.60	0.50
2:M:20:LEU:HD12	2:M:21:SER:N	2.27	0.50
1:C:953:THR:HG22	1:C:954:GLN:HG3	1.94	0.50
2:H:22:CYS:SG	2:H:78:LEU:HD12	2.52	0.50
1:A:169:PHE:CE1	1:A:182:PHE:HD1	2.29	0.49
1:B:301:ILE:HB	1:B:347:THR:HG22	1.94	0.49
1:C:669:ASN:OD1	5:b:1:NAG:C7	2.60	0.49
1:B:339:CYS:HB3	1:B:384:CYS:CA	2.32	0.49
1:C:155:ALA:O	1:C:158:PRO:HD3	2.13	0.49
1:C:819:GLN:O	1:C:823:THR:HG23	2.12	0.49
1:A:52:PHE:HD1	1:A:53:ALA:N	2.10	0.49
3:N:8:PRO:HG2	3:N:11:LEU:HD23	1.95	0.49
1:A:1013:ASN:O	1:A:1014:VAL:HB	2.12	0.49
1:B:585:ASN:HB3	1:B:953:THR:HG23	1.94	0.49
1:A:242:ASN:HB2	1:A:268:PHE:CE1	2.47	0.49
2:H:20:LEU:HD22	2:H:82:MET:HE2	1.94	0.49
2:H:4:LEU:HB3	2:H:104:GLY:HA2	1.94	0.49
3:L:20:THR:HG23	3:L:74:THR:HG22	1.95	0.49
1:A:485:THR:O	1:A:486:ASN:HB2	2.12	0.49
1:B:751:ILE:HD12	1:B:935:LEU:HD22	1.95	0.49
1:C:96:TYR:CD2	1:C:216:ALA:HB2	2.47	0.48
2:M:52:ILE:HG12	2:M:52(A):GLY:N	2.28	0.48
1:A:523:SER:HB3	6:A:1210:NAG:H61	1.95	0.48
1:B:430:VAL:HG11	1:B:433:VAL:HB	1.95	0.48
1:C:156:HIS:O	1:C:156:HIS:ND1	2.47	0.48
3:L:85:THR:HG23	3:L:103:LYS:HD3	1.95	0.48
2:H:39:GLN:NE2	2:H:45:LEU:HD11	2.28	0.47
1:A:303:LEU:HD13	1:A:328:LEU:HD13	1.95	0.47
2:M:43:LYS:HE3	2:M:44:GLY:H	1.79	0.47
1:A:145:CYS:HA	1:A:168:CYS:HA	1.97	0.47
1:A:390:ILE:O	1:A:390:ILE:HG23	2.15	0.47
1:B:1009:ILE:CD1	1:C:784:VAL:CG2	2.93	0.47
2:H:24:ALA:HB1	2:H:27:PHE:HE1	1.79	0.47
1:A:867:LEU:HB2	1:A:872:ALA:HB2	1.96	0.47
1:B:158:PRO:HB2	1:B:161:THR:HG21	1.95	0.47
1:B:437:MET:HE3	1:B:483:ASP:HB2	1.97	0.47
1:C:35:HIS:O	1:C:37:GLU:HG2	2.13	0.47
1:C:565:ARG:CZ	1:C:819:GLN:OE1	2.61	0.47
2:H:20:LEU:CD2	2:H:90:TYR:HD2	2.10	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:ARG:HH12	1:C:188:LYS:HD2	1.79	0.47
1:A:98:ASN:C	1:A:100:THR:N	2.73	0.47
1:A:195:ILE:HG21	1:A:217:VAL:HG11	1.97	0.47
1:A:74:PHE:CE1	1:A:238:ASP:HB3	2.49	0.47
1:A:186:LEU:HD11	1:A:203:ILE:CG2	2.37	0.47
1:A:403:VAL:HG22	1:A:404:ASN:OD1	2.15	0.47
1:C:133:LEU:HD23	1:C:220:ASN:OD1	2.15	0.47
1:C:784:VAL:O	1:C:784:VAL:HG12	2.14	0.47
1:C:516:ASN:OD1	6:C:1208:NAG:O5	2.31	0.47
1:A:698:LEU:HG	1:A:831:ILE:HG22	1.97	0.47
1:B:379:LYS:HB3	1:B:420:GLY:O	2.15	0.47
3:L:3:GLN:C	3:L:4:MET:HE2	2.39	0.46
1:A:737:VAL:HG23	1:C:549:SER:HA	1.97	0.46
1:B:799:ALA:O	1:B:803:ILE:HG12	2.15	0.46
3:L:45:LYS:HD2	3:L:46:LEU:N	2.31	0.46
1:C:131:THR:HG23	1:C:220:ASN:HB2	1.97	0.46
1:A:161:THR:HG23	1:A:167:TYR:CE2	2.50	0.46
1:B:307:PHE:CD2	1:B:321:PRO:HG3	2.51	0.46
1:A:169:PHE:HE1	1:A:182:PHE:HD1	1.61	0.46
1:B:76:PRO:CB	1:B:116:VAL:HG11	2.42	0.46
2:M:36:TRP:CH2	2:M:92:CYS:HB3	2.50	0.46
1:C:128:ARG:HH21	1:C:166:PHE:HB2	1.81	0.46
1:A:390:ILE:HD13	1:A:416:SER:OG	2.15	0.46
1:C:738:ALA:HB1	1:C:743:MET:HE3	1.97	0.46
1:C:35:HIS:C	1:C:37:GLU:N	2.71	0.46
1:C:124:GLU:H	1:C:191:ARG:HH12	1.64	0.46
2:H:99:ARG:HH12	2:H:100(F):TRP:N	2.14	0.46
3:L:59:PRO:HG2	3:L:61:ARG:HG2	1.98	0.46
2:M:36:TRP:CD1	2:M:80:LEU:HG	2.51	0.46
1:A:126:ASN:HA	1:A:190:VAL:HG22	1.97	0.45
1:C:145:CYS:HA	1:C:168:CYS:HA	1.98	0.45
1:C:37:GLU:HA	1:C:40:PHE:CA	2.46	0.45
2:M:66:ARG:HD2	2:M:82(B):SER:O	2.16	0.45
3:N:35:TRP:HB2	3:N:48:ILE:HB	1.98	0.45
1:A:52:PHE:CD1	1:A:53:ALA:N	2.84	0.45
2:H:38:ARG:HD2	2:H:90:TYR:HE1	1.82	0.45
1:A:127:LEU:HD22	1:A:219:PHE:CE1	2.51	0.45
1:A:270:VAL:HG23	1:A:270:VAL:O	2.16	0.45
1:A:288:SER:HB3	1:A:439:VAL:HG12	1.97	0.45
1:B:354:LYS:HE3	2:M:100(F):TRP:CE2	2.51	0.45
2:H:6:GLU:HB2	2:H:107:THR:OG1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:577:THR:O	1:A:577:THR:CG2	2.63	0.45
1:B:699:VAL:HG23	1:B:832:GLN:HE22	1.82	0.45
1:B:683:ARG:HG2	1:B:741:GLU:OE1	2.16	0.45
1:B:586:TRP:CZ2	1:B:995:MET:HG2	2.52	0.44
1:C:90:PHE:CE2	1:C:92:THR:HA	2.52	0.44
2:H:82:MET:HB3	2:H:82(C):LEU:HD11	1.98	0.44
3:L:35:TRP:O	3:L:46:LEU:HD12	2.17	0.44
1:B:616:ARG:HH22	1:B:620:LEU:HD21	1.81	0.44
1:C:35:HIS:C	1:C:37:GLU:H	2.26	0.44
1:C:39:VAL:O	1:C:39:VAL:HG13	2.17	0.44
2:H:34:MET:HE3	2:H:78:LEU:HD21	1.99	0.44
3:N:32:TYR:HB3	3:N:91:TYR:CD1	2.52	0.44
1:C:518:THR:HB	1:C:522:PHE:O	2.17	0.44
1:C:649:MET:HE1	1:C:905:ARG:HG3	1.99	0.44
1:A:708:ALA:HB2	1:A:723:ALA:HB1	1.98	0.44
1:B:133:LEU:CD2	1:B:142:VAL:CG2	2.96	0.44
1:A:790:ARG:HG3	1:A:790:ARG:HH11	1.82	0.44
1:B:161:THR:HB	1:B:167:TYR:CE2	2.53	0.44
2:H:36:TRP:HB3	2:H:48:VAL:HB	1.99	0.44
2:M:59:TYR:CE1	2:M:69:ILE:HG22	2.50	0.44
1:B:58:PHE:HE2	1:B:191:ARG:HB2	1.82	0.44
3:L:28:ASP:OD1	3:L:30:THR:HG22	2.17	0.44
2:M:51:ILE:HG23	2:M:71:ARG:NH2	2.33	0.44
2:H:34:MET:HB2	2:H:34:MET:HE2	1.69	0.44
3:L:94:LEU:HD22	3:L:94:LEU:H	1.83	0.44
1:C:221:VAL:HG13	1:C:227:THR:HG21	1.99	0.44
3:N:89:GLN:HB2	3:N:98:PHE:CD2	2.53	0.44
1:A:344:HIS:ND1	1:A:426:VAL:HG11	2.33	0.43
1:A:1024:LEU:HD23	1:A:1031:TYR:CD2	2.53	0.43
3:L:28:ASP:OD1	3:L:68:GLY:HA2	2.18	0.43
2:M:52:ILE:CG1	2:M:52(A):GLY:N	2.81	0.43
1:A:186:LEU:HD13	1:A:203:ILE:HG21	1.99	0.43
1:C:108:PHE:HE1	1:C:110:SER:HB2	1.82	0.43
3:L:11:LEU:HD21	3:L:13:ALA:HB2	1.99	0.43
1:B:296:HIS:H	1:C:156:HIS:CE1	2.36	0.43
1:C:179:THR:CG2	1:C:180:SER:N	2.81	0.43
1:A:800:MET:HE1	1:A:946:LEU:HD13	2.01	0.43
1:B:375:ASN:HB3	1:B:427:PRO:HA	2.00	0.43
1:C:124:GLU:H	1:C:191:ARG:NH1	2.16	0.43
1:C:134:PHE:CE1	1:C:217:VAL:CG1	3.02	0.43
2:H:60:ALA:HB3	2:H:63:VAL:HG23	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:ASN:O	1:A:100:THR:N	2.52	0.43
2:M:96:THR:HG23	2:M:97:THR:HG23	2.01	0.43
1:C:473:SER:HB3	1:C:479:LEU:HD11	2.00	0.43
1:C:856:ILE:HD11	1:C:866:ARG:HD2	1.99	0.43
2:H:18:LEU:HD13	2:H:19:ARG:N	2.33	0.43
3:N:28:ASP:C	3:N:29:ILE:HD13	2.44	0.43
1:A:242:ASN:HB2	1:A:268:PHE:CZ	2.54	0.42
1:B:580:LEU:HD23	1:C:759:LEU:HD23	2.01	0.42
1:A:132:ILE:HD12	1:A:193:PHE:CE2	2.54	0.42
1:A:586:TRP:CZ2	1:A:995:MET:HG2	2.54	0.42
1:B:161:THR:HB	1:B:167:TYR:HE2	1.84	0.42
3:L:35:TRP:CD1	3:L:88:CYS:HB3	2.54	0.42
1:B:616:ARG:HD2	1:B:616:ARG:HA	1.83	0.42
1:B:1034:VAL:HG12	1:B:1034:VAL:O	2.18	0.42
1:B:642:GLU:HG3	1:B:898:TYR:CE1	2.53	0.42
1:C:44:SER:HA	1:C:108:PHE:HD2	1.84	0.42
3:L:13:ALA:HB1	3:L:17:ASP:OD2	2.18	0.42
1:A:96:TYR:CD2	1:A:216:ALA:HB2	2.54	0.42
1:C:82:LEU:HD13	1:C:232:ALA:HB2	2.02	0.42
1:A:337:PRO:HB3	1:A:386:SER:HB2	2.00	0.42
1:A:371:PHE:HB2	1:A:396:MET:SD	2.60	0.42
1:C:677:LEU:HG	6:C:1214:NAG:H83	2.02	0.42
1:C:900:GLU:HA	1:C:900:GLU:OE2	2.20	0.42
1:A:193:PHE:CD1	1:A:193:PHE:C	2.98	0.42
1:A:568:SER:HA	1:B:653:ASP:OD2	2.19	0.42
1:B:90:PHE:CE2	1:B:92:THR:HA	2.55	0.42
1:B:449:ILE:HD13	1:B:481:PHE:CE2	2.54	0.42
1:B:485:THR:OG1	1:B:486:ASN:N	2.52	0.42
1:C:108:PHE:CD1	1:C:108:PHE:C	2.98	0.42
2:H:95:ASP:OD1	2:H:96:THR:N	2.53	0.42
3:L:6:GLN:HG3	3:L:23:CYS:SG	2.60	0.42
3:N:32:TYR:HD2	3:N:92:ASP:HB2	1.84	0.42
1:C:40:PHE:CE2	1:C:42:VAL:HG12	2.54	0.42
1:A:127:LEU:HD21	1:A:132:ILE:HD11	2.02	0.42
1:C:44:SER:HA	1:C:108:PHE:CD2	2.55	0.42
2:H:38:ARG:HB3	2:H:90:TYR:CE1	2.55	0.42
2:M:45:LEU:O	3:N:98:PHE:HB2	2.19	0.42
1:B:188:LYS:CG	1:B:189:THR:N	2.83	0.41
1:A:799:ALA:O	1:A:803:ILE:HG12	2.20	0.41
2:H:39:GLN:CD	3:L:38:GLN:HE21	2.24	0.41
1:A:127:LEU:CD2	1:A:132:ILE:HD11	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:TYR:CZ	1:A:297:LYS:HD3	2.55	0.41
1:B:157:ILE:O	1:B:157:ILE:HG13	2.20	0.41
1:C:37:GLU:O	1:C:40:PHE:N	2.53	0.41
1:C:814:ILE:O	1:C:814:ILE:HG22	2.20	0.41
2:M:20:LEU:HD11	2:M:36:TRP:CH2	2.56	0.41
2:M:89:VAL:HA	2:M:108:PRO:HA	2.01	0.41
1:A:648:GLU:OE2	1:A:648:GLU:HA	2.20	0.41
1:C:90:PHE:C	1:C:92:THR:H	2.28	0.41
2:H:71:ARG:HG2	2:H:72:ASP:N	2.35	0.41
3:N:90:GLN:OE1	3:N:96:LEU:HA	2.20	0.41
1:A:52:PHE:HD1	1:A:53:ALA:H	1.68	0.41
1:A:602:ILE:HD13	1:A:842:LEU:HD21	2.02	0.41
1:C:157:ILE:O	1:C:157:ILE:HD12	2.21	0.41
1:C:224:ALA:C	1:C:226:THR:H	2.29	0.41
2:H:3:GLN:OE1	2:H:5:VAL:HG13	2.21	0.41
2:H:36:TRP:CE2	2:H:80:LEU:HD22	2.56	0.41
2:M:52:ILE:CG1	2:M:52(A):GLY:H	2.33	0.41
2:M:72:ASP:HB3	2:M:77:THR:OG1	2.20	0.41
1:A:186:LEU:HD13	1:A:203:ILE:CD1	2.46	0.41
1:B:674:ILE:HD12	1:B:674:ILE:O	2.21	0.41
2:H:22:CYS:HB3	2:H:78:LEU:HB2	2.02	0.41
3:L:39:LYS:HB3	3:L:42:LYS:HB2	2.03	0.41
2:M:3:GLN:C	2:M:4:LEU:HD12	2.46	0.41
1:A:356:ALA:O	1:A:357:ARG:HB2	2.21	0.41
1:A:857:SER:H	1:A:863:ILE:HD11	1.85	0.41
3:L:8:PRO:HD2	3:L:21:ILE:HG12	2.02	0.41
1:B:327:THR:HG21	4:T:1:NAG:H5	2.02	0.41
1:C:37:GLU:C	1:C:39:VAL:H	2.28	0.41
2:M:43:LYS:HD2	2:M:43:LYS:HA	1.99	0.41
3:N:4:MET:HB3	3:N:25:ALA:HA	2.03	0.41
1:A:280:ILE:C	1:A:280:ILE:HD12	2.46	0.41
1:B:819:GLN:HG2	4:V:1:NAG:H61	2.03	0.41
1:B:830:LYS:HB2	1:B:830:LYS:HE2	1.76	0.40
1:B:1009:ILE:HD12	1:C:784:VAL:HG23	2.02	0.40
1:C:655:LYS:O	1:C:658:THR:HG22	2.21	0.40
1:A:964:LEU:HB2	1:A:972:TYR:HB2	2.04	0.40
1:B:1020:SER:HB3	1:B:1023:GLU:CD	2.46	0.40
2:M:58:TYR:CD1	3:N:94:LEU:HD11	2.56	0.40
1:B:128:ARG:HH21	1:B:187:PRO:C	2.29	0.40
1:B:142:VAL:HG12	1:B:171:ASN:HB2	2.04	0.40
1:A:142:VAL:CG1	1:A:171:ASN:HB2	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:659:LEU:C	1:C:661:ASN:H	2.30	0.40
1:A:280:ILE:HD12	1:A:280:ILE:O	2.22	0.40
3:L:35:TRP:HH2	3:L:71:PHE:HB3	1.86	0.40
2:M:29:PHE:CE2	2:M:71:ARG:HD3	2.56	0.40
2:M:103:TRP:CD1	2:M:103:TRP:N	2.89	0.40
3:N:70:ASP:OD1	3:N:70:ASP:C	2.65	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1010/1134 (89%)	955 (95%)	53 (5%)	2 (0%)	44	72
1	B	1010/1134 (89%)	962 (95%)	46 (5%)	2 (0%)	44	72
1	C	869/1134 (77%)	841 (97%)	27 (3%)	1 (0%)	48	77
2	H	122/239 (51%)	117 (96%)	5 (4%)	0	100	100
2	M	122/239 (51%)	113 (93%)	9 (7%)	0	100	100
3	L	104/214 (49%)	96 (92%)	8 (8%)	0	100	100
3	N	104/214 (49%)	93 (89%)	10 (10%)	1 (1%)	13	43
All	All	3341/4308 (78%)	3177 (95%)	158 (5%)	6 (0%)	45	72

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	33	VAL
1	A	486	ASN
3	N	83	ILE
1	B	484	VAL
1	B	42	VAL

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Mol	Chain	Res	Type
1	C	33	VAL

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	873/969 (90%)	869 (100%)	4 (0%)	86	92
1	B	873/969 (90%)	869 (100%)	4 (0%)	86	92
1	C	752/969 (78%)	749 (100%)	3 (0%)	89	94
2	H	99/199 (50%)	96 (97%)	3 (3%)	36	63
2	M	99/199 (50%)	98 (99%)	1 (1%)	73	85
3	L	92/189 (49%)	92 (100%)	0	100	100
3	N	92/189 (49%)	91 (99%)	1 (1%)	70	84
All	All	2880/3683 (78%)	2864 (99%)	16 (1%)	82	91

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	PHE
1	A	98	ASN
1	A	193	PHE
1	A	711	LYS
1	B	207	ARG
1	B	250	ASN
1	B	517	PHE
1	B	654	LYS
1	C	39	VAL
1	C	218	ASN
1	C	229	CYS
2	H	19	ARG
2	H	32	TYR
2	H	46	GLU
2	M	86	ASP
3	N	90	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	200	HIS
1	A	245	GLN
1	A	309	HIS
1	A	348	GLN
1	A	352	ASN
1	A	729	ASN
1	A	919	GLN
1	A	937	ASN
1	B	148	ASN
1	B	200	HIS
1	B	348	GLN
1	B	486	ASN
1	B	729	ASN
1	B	871	GLN
1	B	919	GLN
1	B	978	ASN
1	C	200	HIS
1	C	281	GLN
1	C	591	GLN
1	C	781	GLN
1	C	832	GLN
1	C	909	GLN
1	C	913	ASN
1	C	919	GLN
1	C	978	ASN
2	H	39	GLN
3	L	37	GLN
3	L	38	GLN
3	L	89	GLN
2	M	13	GLN
2	M	56	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

59 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	D	1	4,1	14,14,15	0.77	0	17,19,21	1.50	2 (11%)
4	NAG	D	2	4	14,14,15	0.67	0	17,19,21	1.83	3 (17%)
4	BMA	D	3	4	11,11,12	0.82	0	15,15,17	0.61	0
4	NAG	E	1	4,1	14,14,15	0.75	0	17,19,21	1.89	2 (11%)
4	NAG	E	2	4	14,14,15	0.79	0	17,19,21	1.30	2 (11%)
4	BMA	E	3	4	11,11,12	0.95	0	15,15,17	1.84	1 (6%)
5	NAG	F	1	5,1	14,14,15	0.70	0	17,19,21	1.24	3 (17%)
5	NAG	F	2	5	14,14,15	0.82	0	17,19,21	0.79	0
4	NAG	G	1	4,1	14,14,15	0.73	0	17,19,21	1.08	1 (5%)
4	NAG	G	2	4	14,14,15	0.67	0	17,19,21	1.63	3 (17%)
4	BMA	G	3	4	11,11,12	0.67	0	15,15,17	2.19	1 (6%)
5	NAG	I	1	5,1	14,14,15	0.75	0	17,19,21	0.80	1 (5%)
5	NAG	I	2	5	14,14,15	0.68	0	17,19,21	1.07	1 (5%)
5	NAG	J	1	5,1	14,14,15	0.81	1 (7%)	17,19,21	1.30	3 (17%)
5	NAG	J	2	5	14,14,15	0.72	0	17,19,21	1.40	4 (23%)
5	NAG	K	1	5,1	14,14,15	0.68	0	17,19,21	1.92	4 (23%)
5	NAG	K	2	5	14,14,15	0.80	1 (7%)	17,19,21	1.90	4 (23%)
5	NAG	O	1	5,1	14,14,15	0.66	0	17,19,21	0.85	1 (5%)
5	NAG	O	2	5	14,14,15	0.66	0	17,19,21	1.02	1 (5%)
4	NAG	P	1	4,1	14,14,15	0.69	0	17,19,21	1.42	3 (17%)
4	NAG	P	2	4	14,14,15	0.73	0	17,19,21	1.41	1 (5%)
4	BMA	P	3	4	11,11,12	0.81	0	15,15,17	1.22	1 (6%)
4	NAG	Q	1	4,1	14,14,15	0.71	0	17,19,21	1.22	3 (17%)
4	NAG	Q	2	4	14,14,15	0.63	0	17,19,21	0.82	0
4	BMA	Q	3	4	11,11,12	0.92	0	15,15,17	1.01	0
4	NAG	R	1	4,1	14,14,15	0.76	0	17,19,21	1.16	2 (11%)
4	NAG	R	2	4	14,14,15	0.71	0	17,19,21	1.41	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	BMA	R	3	4	11,11,12	0.75	0	15,15,17	1.11	1 (6%)
5	NAG	S	1	5,1	14,14,15	0.71	0	17,19,21	0.88	0
5	NAG	S	2	5	14,14,15	0.65	0	17,19,21	1.30	1 (5%)
4	NAG	T	1	4,1	14,14,15	0.89	1 (7%)	17,19,21	1.87	2 (11%)
4	NAG	T	2	4	14,14,15	0.71	0	17,19,21	1.11	1 (5%)
4	BMA	T	3	4	11,11,12	0.76	0	15,15,17	1.28	1 (6%)
4	NAG	U	1	4,1	14,14,15	0.71	0	17,19,21	0.93	1 (5%)
4	NAG	U	2	4	14,14,15	0.65	0	17,19,21	1.08	2 (11%)
4	BMA	U	3	4	11,11,12	0.70	0	15,15,17	1.14	1 (6%)
4	NAG	V	1	4,1	14,14,15	0.78	0	17,19,21	1.04	1 (5%)
4	NAG	V	2	4	14,14,15	0.67	0	17,19,21	1.13	1 (5%)
4	BMA	V	3	4	11,11,12	0.75	0	15,15,17	1.05	1 (6%)
5	NAG	W	1	5,1	14,14,15	0.69	0	17,19,21	1.43	1 (5%)
5	NAG	W	2	5	14,14,15	0.66	0	17,19,21	0.99	1 (5%)
5	NAG	X	1	5,1	14,14,15	0.71	0	17,19,21	1.05	1 (5%)
5	NAG	X	2	5	14,14,15	0.66	0	17,19,21	0.78	1 (5%)
5	NAG	Y	1	5,1	14,14,15	0.67	0	17,19,21	1.44	4 (23%)
5	NAG	Y	2	5	14,14,15	0.68	0	17,19,21	2.14	5 (29%)
5	NAG	Z	1	5,1	14,14,15	0.75	0	17,19,21	1.05	1 (5%)
5	NAG	Z	2	5	14,14,15	0.83	0	17,19,21	1.02	1 (5%)
4	NAG	a	1	4,1	14,14,15	0.79	0	17,19,21	1.61	3 (17%)
4	NAG	a	2	4	14,14,15	0.70	0	17,19,21	1.01	1 (5%)
4	BMA	a	3	4	11,11,12	0.73	0	15,15,17	1.51	1 (6%)
5	NAG	b	1	5,1	14,14,15	1.18	1 (7%)	17,19,21	1.36	2 (11%)
5	NAG	b	2	5	14,14,15	0.65	0	17,19,21	1.31	1 (5%)
4	NAG	c	1	4,1	14,14,15	0.72	0	17,19,21	1.17	2 (11%)
4	NAG	c	2	4	14,14,15	0.70	1 (7%)	17,19,21	2.54	5 (29%)
4	BMA	c	3	4	11,11,12	0.80	0	15,15,17	1.28	1 (6%)
5	NAG	d	1	5,1	14,14,15	0.82	0	17,19,21	1.30	3 (17%)
5	NAG	d	2	5	14,14,15	0.66	0	17,19,21	1.08	1 (5%)
5	NAG	e	1	5,1	14,14,15	0.70	0	17,19,21	1.35	1 (5%)
5	NAG	e	2	5	14,14,15	0.68	0	17,19,21	1.10	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	D	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	0/6/23/26	0/1/1/1
4	BMA	D	3	4	-	1/2/19/22	0/1/1/1
4	NAG	E	1	4,1	-	1/6/23/26	0/1/1/1
4	NAG	E	2	4	-	2/6/23/26	0/1/1/1
4	BMA	E	3	4	-	1/2/19/22	0/1/1/1
5	NAG	F	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	F	2	5	-	0/6/23/26	0/1/1/1
4	NAG	G	1	4,1	-	3/6/23/26	0/1/1/1
4	NAG	G	2	4	-	1/6/23/26	0/1/1/1
4	BMA	G	3	4	-	2/2/19/22	0/1/1/1
5	NAG	I	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	0/6/23/26	0/1/1/1
5	NAG	J	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	J	2	5	-	1/6/23/26	0/1/1/1
5	NAG	K	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	K	2	5	-	0/6/23/26	0/1/1/1
5	NAG	O	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	O	2	5	-	0/6/23/26	0/1/1/1
4	NAG	P	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	P	2	4	-	2/6/23/26	0/1/1/1
4	BMA	P	3	4	-	0/2/19/22	0/1/1/1
4	NAG	Q	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	Q	2	4	-	1/6/23/26	0/1/1/1
4	BMA	Q	3	4	-	2/2/19/22	0/1/1/1
4	NAG	R	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	R	2	4	-	3/6/23/26	0/1/1/1
4	BMA	R	3	4	-	0/2/19/22	0/1/1/1
5	NAG	S	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	S	2	5	-	0/6/23/26	0/1/1/1
4	NAG	T	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	T	2	4	-	1/6/23/26	0/1/1/1
4	BMA	T	3	4	-	0/2/19/22	0/1/1/1
4	NAG	U	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	U	2	4	-	0/6/23/26	0/1/1/1
4	BMA	U	3	4	-	0/2/19/22	0/1/1/1
4	NAG	V	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	V	2	4	-	0/6/23/26	0/1/1/1
4	BMA	V	3	4	-	0/2/19/22	0/1/1/1
5	NAG	W	1	5,1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	W	2	5	-	1/6/23/26	0/1/1/1
5	NAG	X	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	X	2	5	-	1/6/23/26	0/1/1/1
5	NAG	Y	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	Y	2	5	-	1/6/23/26	0/1/1/1
5	NAG	Z	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	Z	2	5	-	1/6/23/26	0/1/1/1
4	NAG	a	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	a	2	4	-	0/6/23/26	0/1/1/1
4	BMA	a	3	4	-	1/2/19/22	0/1/1/1
5	NAG	b	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	b	2	5	-	0/6/23/26	0/1/1/1
4	NAG	c	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	c	2	4	-	1/6/23/26	0/1/1/1
4	BMA	c	3	4	-	0/2/19/22	0/1/1/1
5	NAG	d	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	d	2	5	-	0/6/23/26	0/1/1/1
5	NAG	e	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	e	2	5	-	0/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	b	1	NAG	C1-C2	4.01	1.57	1.52
5	K	2	NAG	C1-C2	2.25	1.55	1.52
4	T	1	NAG	O5-C1	-2.10	1.40	1.43
5	J	1	NAG	C1-C2	2.08	1.55	1.52
4	c	2	NAG	C1-C2	2.06	1.55	1.52

All (100) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	3	BMA	C1-O5-C5	8.17	123.14	112.19
4	c	2	NAG	C2-N2-C7	7.16	132.50	122.90
4	E	3	BMA	C1-O5-C5	6.37	120.73	112.19
4	D	2	NAG	C1-O5-C5	6.16	120.44	112.19
4	T	1	NAG	C2-N2-C7	6.15	131.14	122.90
4	E	1	NAG	C1-C2-N2	5.88	119.70	110.43
5	K	2	NAG	C1-O5-C5	5.87	120.06	112.19
5	K	1	NAG	C1-C2-N2	5.84	119.64	110.43
4	a	3	BMA	C1-O5-C5	5.32	119.32	112.19
5	Y	2	NAG	C2-N2-C7	5.16	129.82	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	c	2	NAG	C1-C2-N2	4.74	117.90	110.43
5	Y	2	NAG	C1-C2-N2	4.53	117.57	110.43
4	D	1	NAG	C1-O5-C5	4.49	118.21	112.19
5	S	2	NAG	C1-O5-C5	4.44	118.14	112.19
4	T	3	BMA	C1-O5-C5	4.44	118.14	112.19
5	b	1	NAG	C2-N2-C7	4.34	128.71	122.90
5	e	1	NAG	C1-O5-C5	4.20	117.82	112.19
4	P	2	NAG	C2-N2-C7	4.18	128.50	122.90
5	b	2	NAG	C1-O5-C5	4.16	117.76	112.19
4	a	1	NAG	C1-O5-C5	4.09	117.66	112.19
4	G	2	NAG	C4-C3-C2	-3.93	105.26	111.02
4	c	3	BMA	C1-O5-C5	3.84	117.34	112.19
5	K	1	NAG	C2-N2-C7	3.73	127.90	122.90
4	U	3	BMA	C1-O5-C5	3.73	117.18	112.19
4	c	2	NAG	C1-O5-C5	3.70	117.15	112.19
5	W	1	NAG	C2-N2-C7	3.69	127.85	122.90
5	Y	1	NAG	C2-N2-C7	3.62	127.75	122.90
4	E	1	NAG	O5-C1-C2	-3.59	105.73	111.29
4	G	2	NAG	C1-O5-C5	3.59	116.99	112.19
5	Y	2	NAG	C1-O5-C5	3.58	116.98	112.19
4	P	1	NAG	C1-O5-C5	3.44	116.80	112.19
5	e	2	NAG	C1-O5-C5	3.38	116.71	112.19
4	T	2	NAG	C1-O5-C5	3.33	116.65	112.19
4	P	3	BMA	C1-O5-C5	3.29	116.60	112.19
4	E	2	NAG	C2-N2-C7	3.29	127.30	122.90
4	V	2	NAG	C1-O5-C5	3.20	116.47	112.19
5	F	1	NAG	C1-O5-C5	3.18	116.44	112.19
5	I	2	NAG	C1-O5-C5	3.08	116.31	112.19
5	J	2	NAG	C2-N2-C7	3.05	126.99	122.90
4	R	1	NAG	C2-N2-C7	3.03	126.96	122.90
4	c	2	NAG	C4-C3-C2	-3.03	106.58	111.02
4	R	2	NAG	C2-N2-C7	3.00	126.92	122.90
5	J	2	NAG	C1-O5-C5	2.98	116.18	112.19
4	D	1	NAG	C2-N2-C7	2.97	126.89	122.90
4	a	1	NAG	C2-N2-C7	2.96	126.86	122.90
4	G	1	NAG	C2-N2-C7	2.92	126.81	122.90
5	X	1	NAG	C2-N2-C7	2.86	126.74	122.90
4	R	3	BMA	C1-O5-C5	2.84	116.00	112.19
4	V	3	BMA	C1-O5-C5	2.79	115.92	112.19
5	J	1	NAG	C2-N2-C7	2.75	126.59	122.90
5	d	2	NAG	C4-C3-C2	-2.72	107.03	111.02
4	Q	1	NAG	C2-N2-C7	2.68	126.49	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	V	1	NAG	C2-N2-C7	2.68	126.49	122.90
4	c	1	NAG	O4-C4-C3	-2.66	104.11	110.38
5	W	2	NAG	C1-O5-C5	2.62	115.70	112.19
4	U	2	NAG	C1-O5-C5	2.62	115.70	112.19
4	Q	1	NAG	O4-C4-C3	-2.60	104.25	110.38
5	K	2	NAG	C2-N2-C7	2.57	126.34	122.90
5	O	1	NAG	C2-N2-C7	2.54	126.30	122.90
4	c	2	NAG	O7-C7-N2	2.52	126.43	121.98
4	P	1	NAG	C4-C3-C2	-2.50	107.36	111.02
5	d	1	NAG	C1-O5-C5	2.47	115.50	112.19
5	O	2	NAG	C1-O5-C5	2.47	115.50	112.19
4	a	1	NAG	C4-C3-C2	-2.47	107.40	111.02
5	b	1	NAG	O4-C4-C3	-2.45	104.61	110.38
5	Y	1	NAG	C4-C3-C2	-2.45	107.43	111.02
4	Q	1	NAG	C1-O5-C5	2.37	115.36	112.19
5	K	2	NAG	O5-C1-C2	2.34	114.92	111.29
4	E	2	NAG	O4-C4-C3	2.34	115.90	110.38
5	d	1	NAG	C2-N2-C7	2.33	126.02	122.90
4	R	2	NAG	C1-O5-C5	2.32	115.30	112.19
5	F	1	NAG	O5-C1-C2	2.31	114.87	111.29
5	J	1	NAG	C1-O5-C5	2.31	115.29	112.19
5	J	1	NAG	O4-C4-C3	-2.31	104.94	110.38
5	d	1	NAG	O5-C1-C2	2.30	114.85	111.29
4	G	2	NAG	C3-C4-C5	-2.28	106.09	110.23
5	J	2	NAG	C1-C2-N2	2.28	114.02	110.43
4	R	2	NAG	C1-C2-N2	2.25	113.98	110.43
5	Z	1	NAG	C2-N2-C7	2.22	125.87	122.90
4	R	1	NAG	C1-O5-C5	2.20	115.13	112.19
5	J	2	NAG	O4-C4-C3	-2.19	105.20	110.38
4	U	1	NAG	C1-O5-C5	2.17	115.09	112.19
4	D	2	NAG	O4-C4-C3	-2.14	105.33	110.38
4	D	2	NAG	C2-N2-C7	2.14	125.77	122.90
4	U	2	NAG	O4-C4-C3	-2.13	105.35	110.38
4	P	1	NAG	O4-C4-C3	-2.13	105.36	110.38
4	a	2	NAG	C2-N2-C7	2.12	125.73	122.90
5	K	1	NAG	O3-C3-C4	2.11	115.35	110.38
5	I	1	NAG	C4-C3-C2	-2.11	107.93	111.02
5	K	2	NAG	O4-C4-C3	-2.09	105.44	110.38
5	Y	1	NAG	C1-C2-N2	2.09	113.73	110.43
5	X	2	NAG	C1-O5-C5	2.07	114.97	112.19
5	K	1	NAG	C4-C3-C2	-2.06	108.00	111.02
5	Y	2	NAG	O5-C1-C2	-2.06	108.10	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	Y	2	NAG	O7-C7-N2	2.05	125.61	121.98
5	F	1	NAG	O5-C5-C4	-2.05	105.84	110.83
5	Y	1	NAG	O5-C1-C2	-2.02	108.16	111.29
5	Z	2	NAG	C2-N2-C7	2.01	125.60	122.90
4	T	1	NAG	C1-C2-N2	2.01	113.60	110.43
4	c	1	NAG	C3-C4-C5	2.00	113.86	110.23

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	1	NAG	C1-C2-N2-C7
4	R	2	NAG	C1-C2-N2-C7
4	c	2	NAG	C1-C2-N2-C7
5	Y	2	NAG	C1-C2-N2-C7
4	c	1	NAG	O5-C5-C6-O6
4	c	1	NAG	C4-C5-C6-O6
4	Q	3	BMA	O5-C5-C6-O6
4	R	1	NAG	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
4	D	3	BMA	O5-C5-C6-O6
4	E	3	BMA	O5-C5-C6-O6
4	T	2	NAG	O5-C5-C6-O6
4	R	2	NAG	O5-C5-C6-O6
5	J	2	NAG	O5-C5-C6-O6
4	G	3	BMA	O5-C5-C6-O6
4	a	3	BMA	O5-C5-C6-O6
5	W	2	NAG	O5-C5-C6-O6
4	Q	2	NAG	O5-C5-C6-O6
5	Z	2	NAG	O5-C5-C6-O6
5	X	2	NAG	O5-C5-C6-O6
4	E	2	NAG	C1-C2-N2-C7
4	G	1	NAG	C1-C2-N2-C7
4	P	2	NAG	C1-C2-N2-C7
5	Y	1	NAG	C1-C2-N2-C7
4	G	1	NAG	C3-C2-N2-C7
5	J	1	NAG	C3-C2-N2-C7
4	Q	3	BMA	C4-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	E	2	NAG	C3-C2-N2-C7
4	P	2	NAG	C3-C2-N2-C7
4	R	2	NAG	C3-C2-N2-C7

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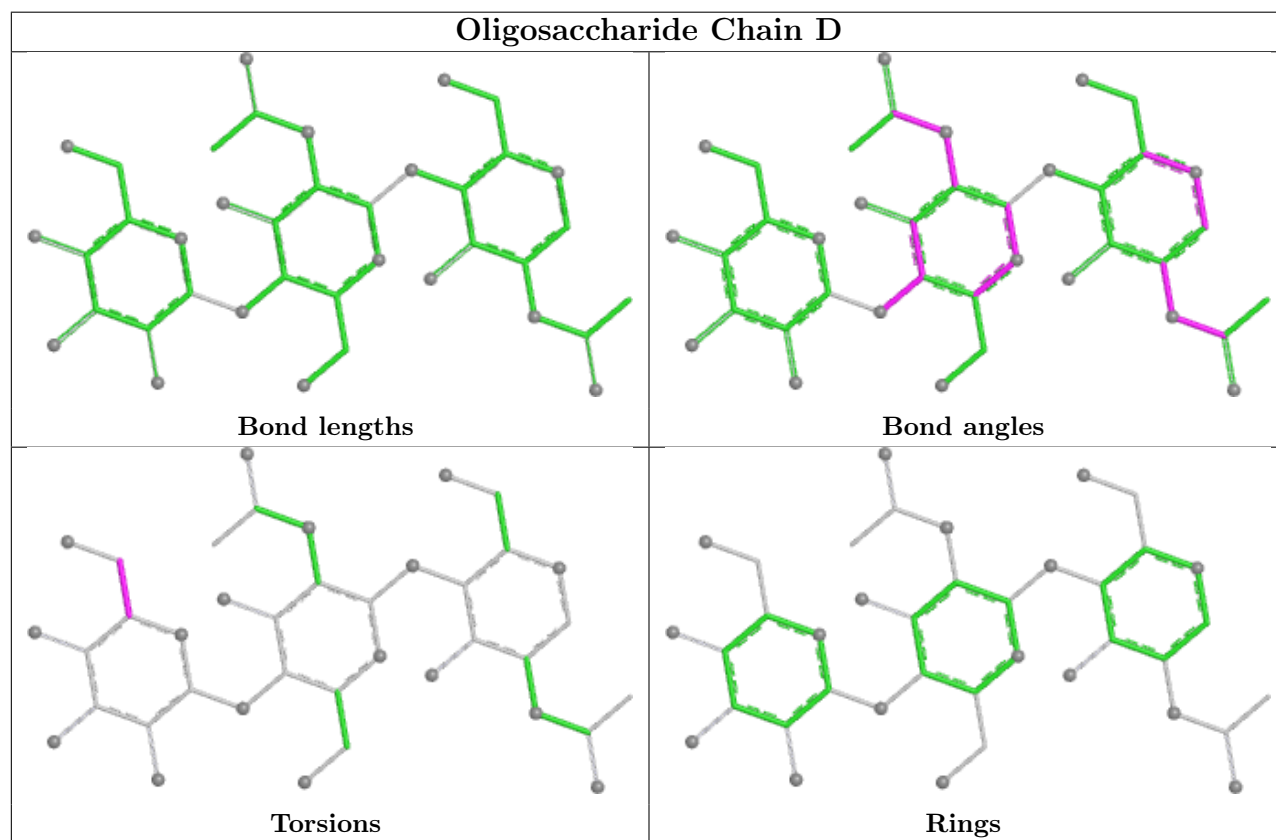
Mol	Chain	Res	Type	Atoms
5	Y	1	NAG	C3-C2-N2-C7
5	X	1	NAG	O5-C5-C6-O6
4	G	3	BMA	C4-C5-C6-O6
4	R	1	NAG	C4-C5-C6-O6

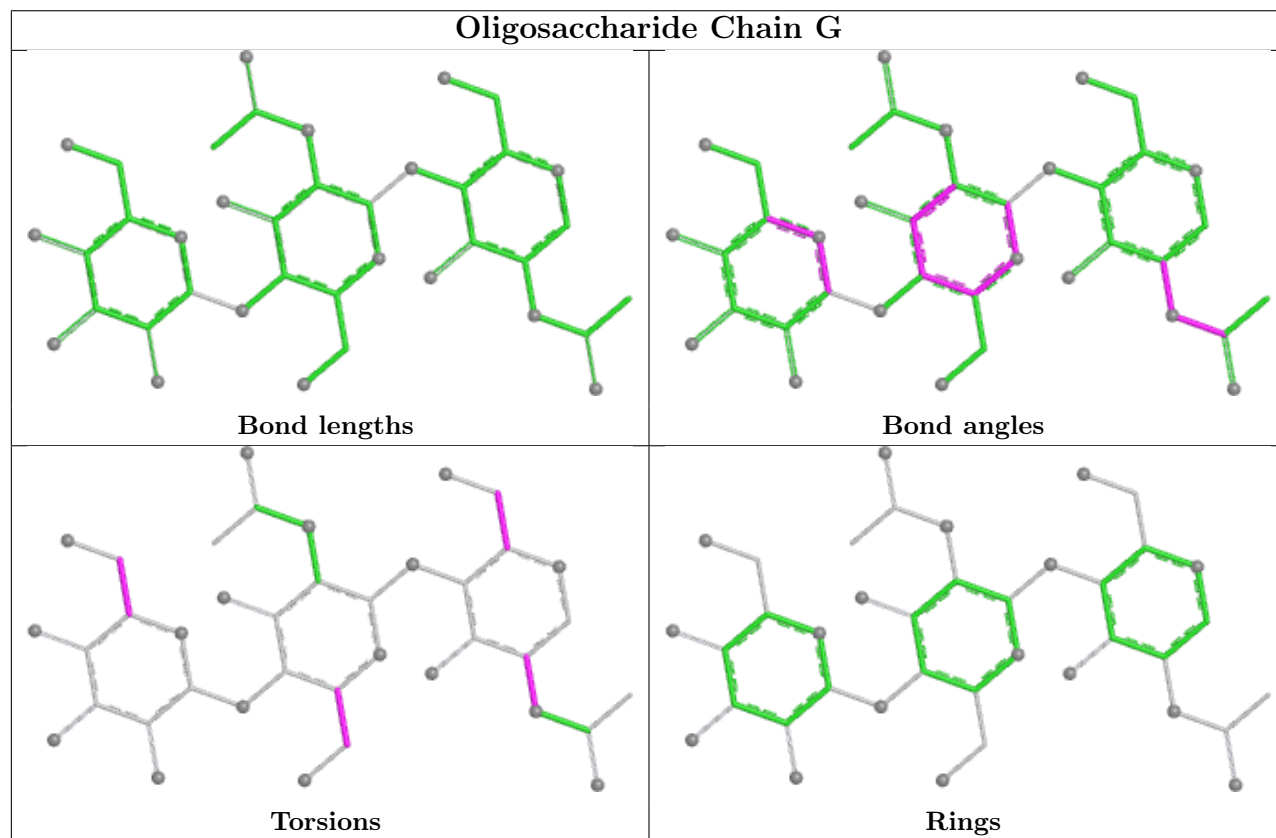
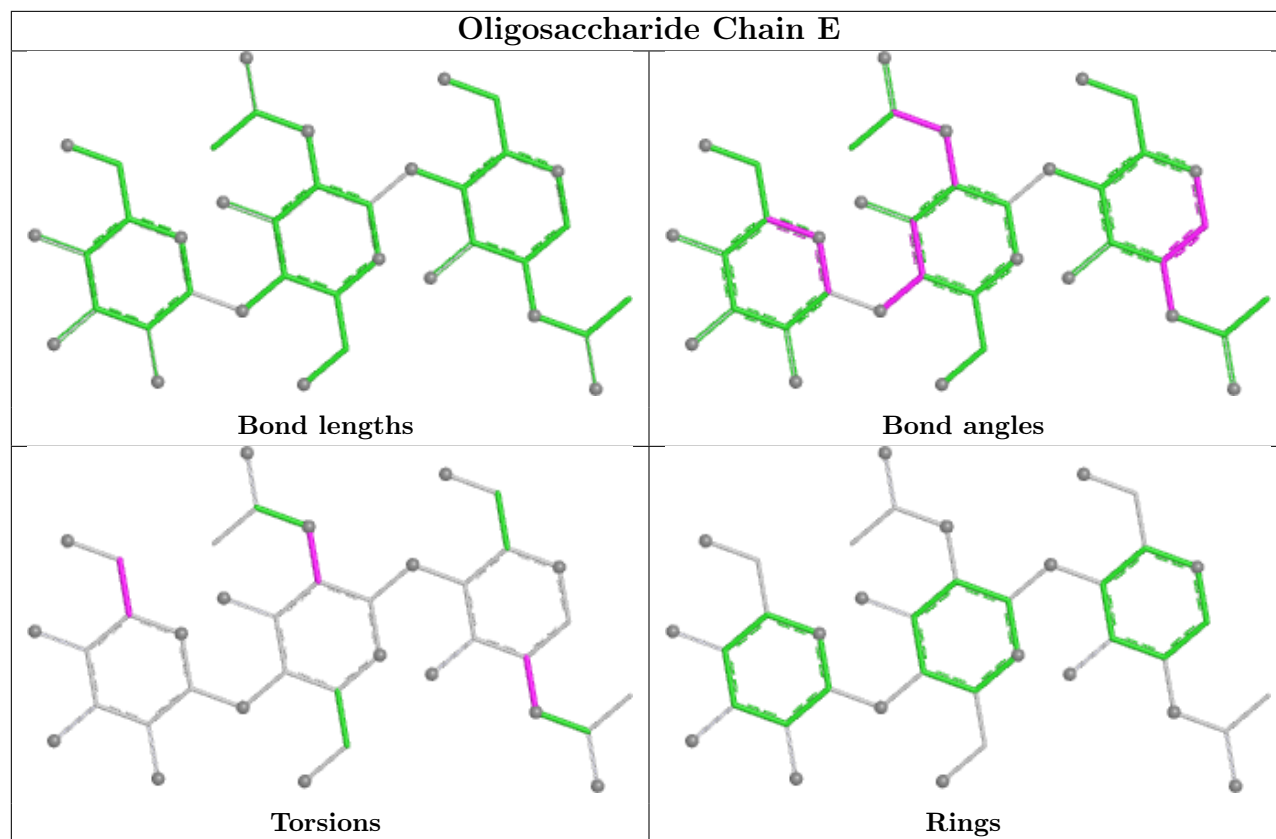
There are no ring outliers.

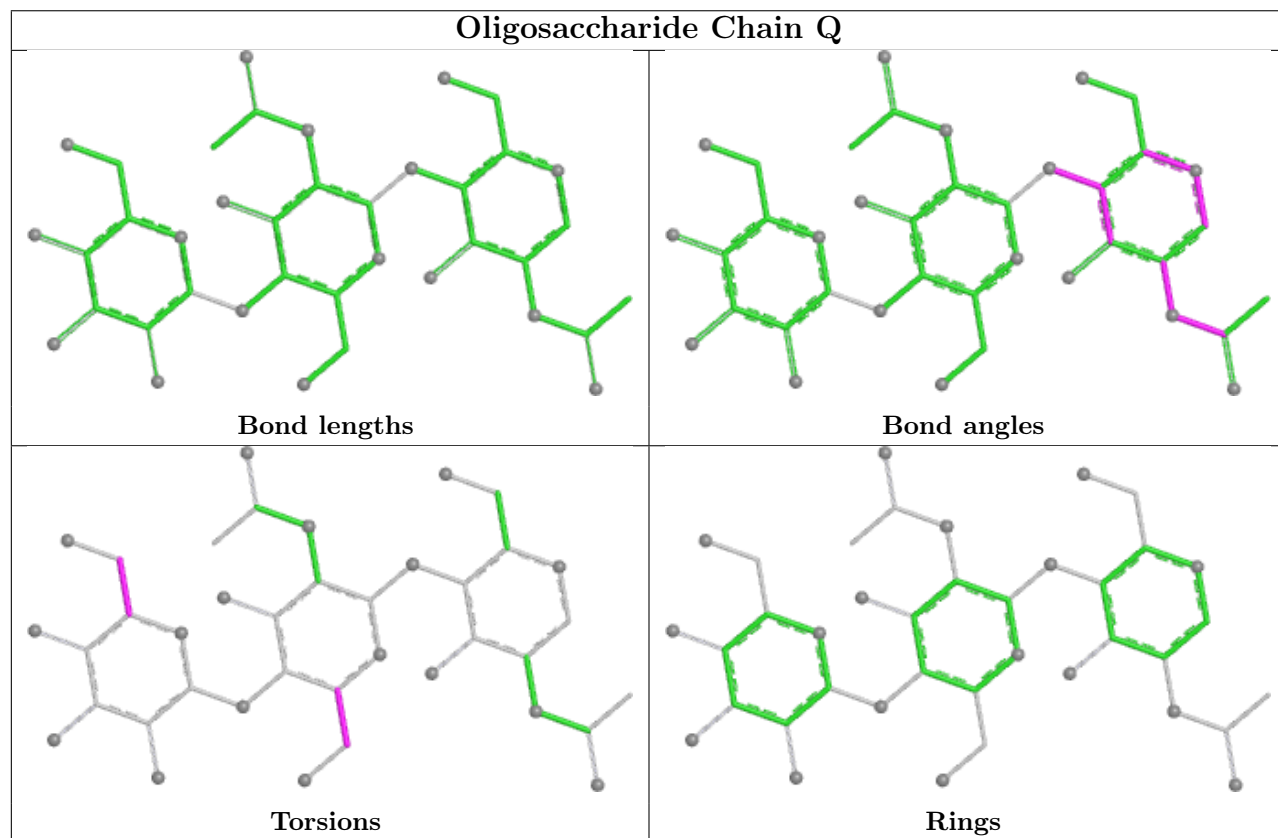
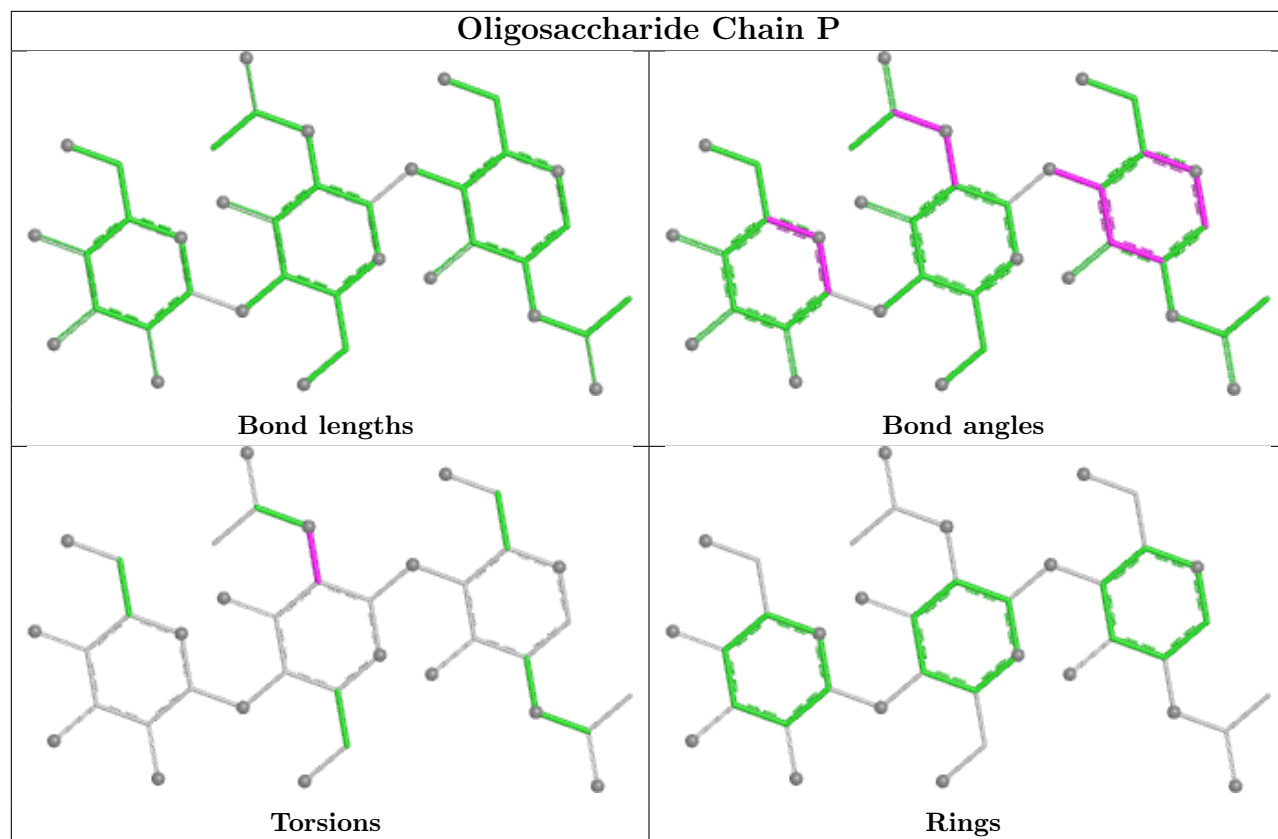
5 monomers are involved in 11 short contacts:

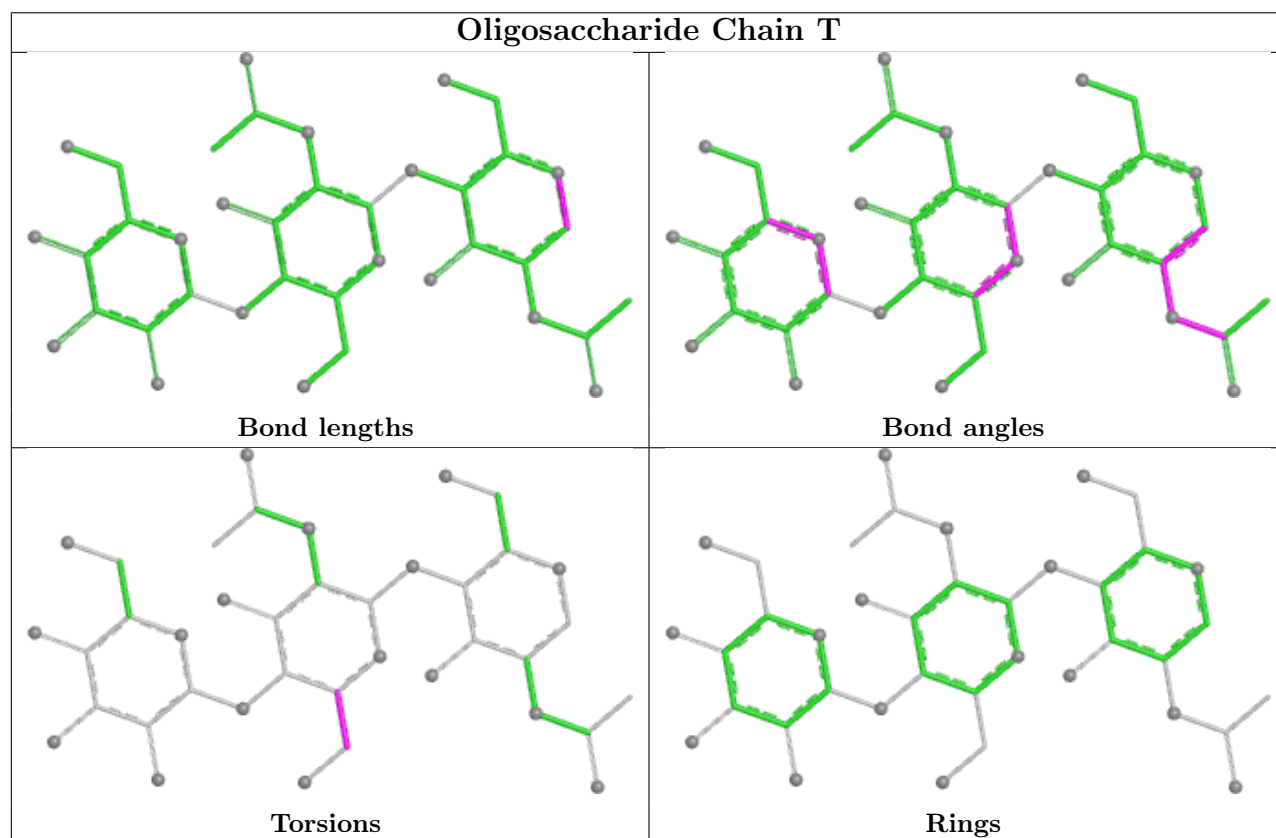
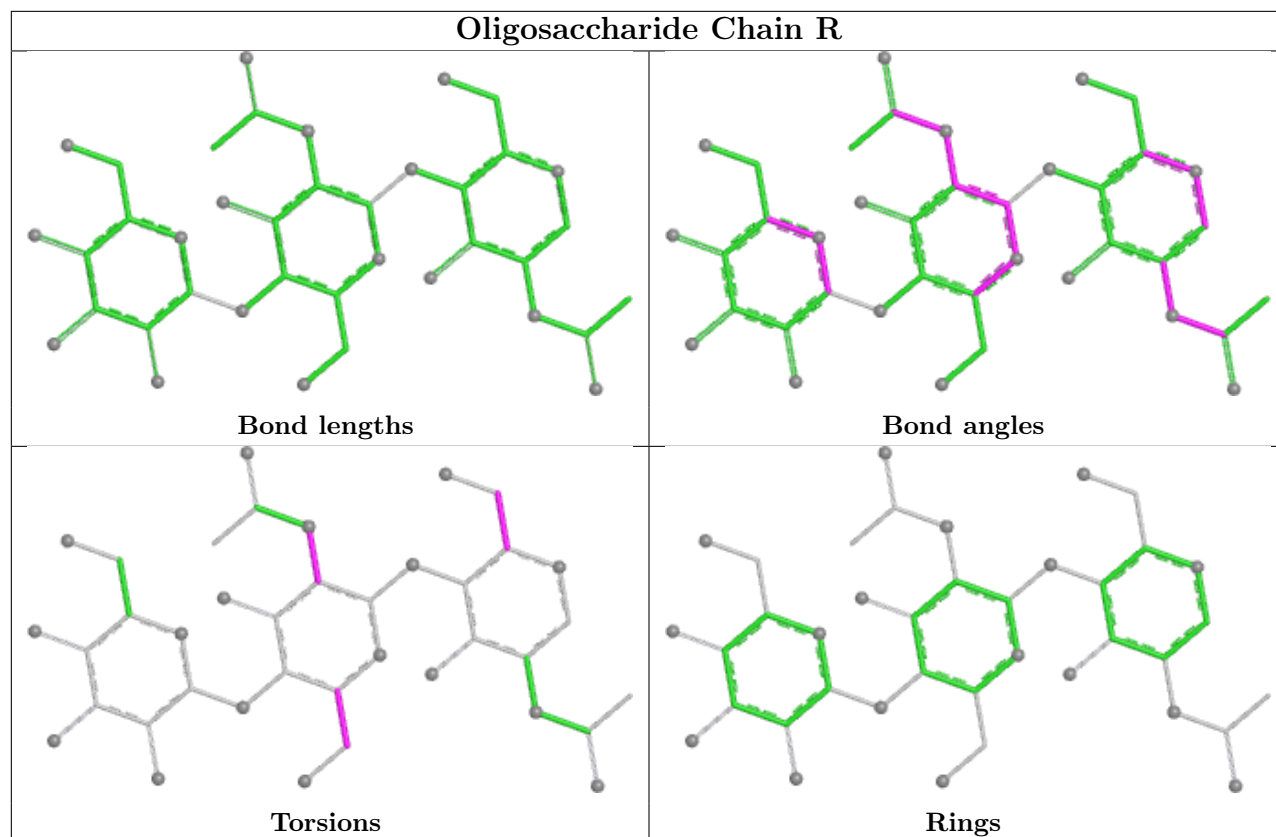
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	b	1	NAG	1	0
4	T	1	NAG	1	0
4	E	1	NAG	7	0
4	V	1	NAG	1	0
5	O	1	NAG	1	0

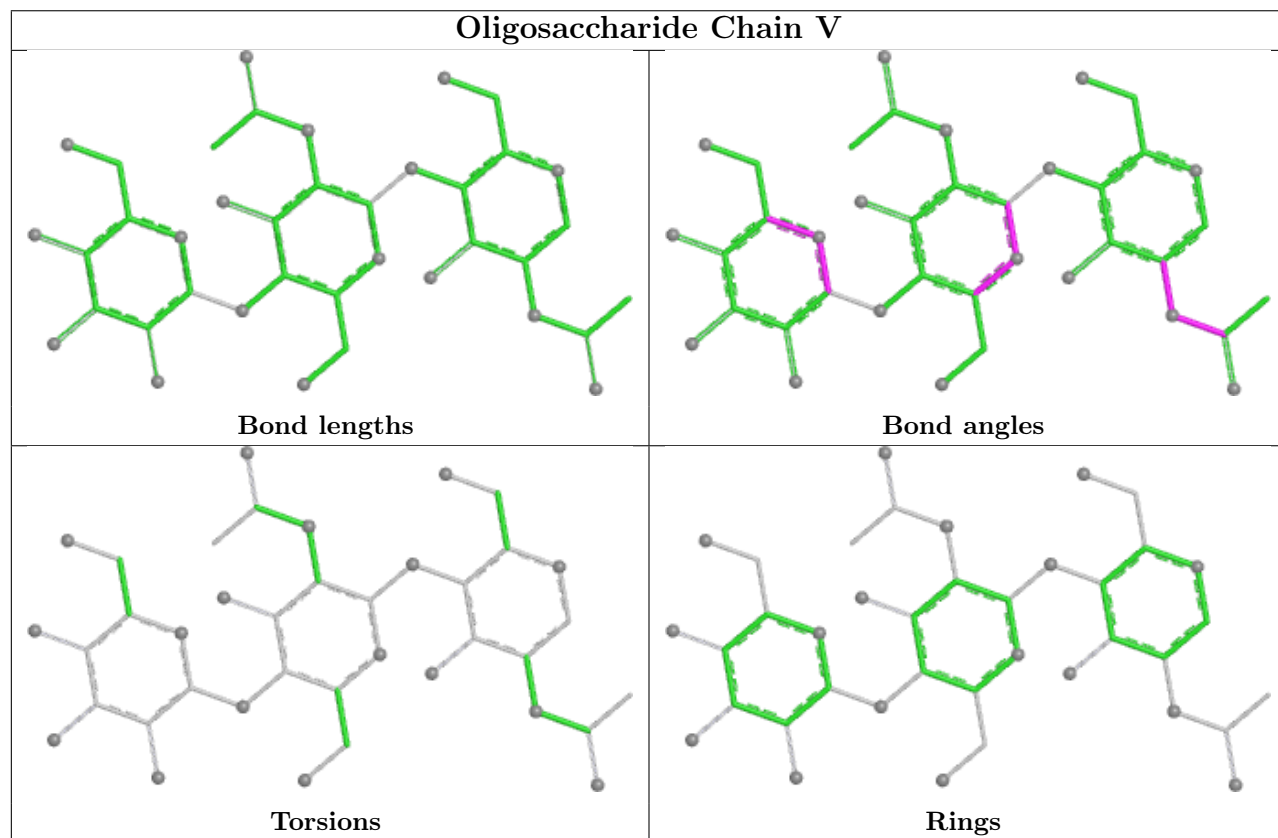
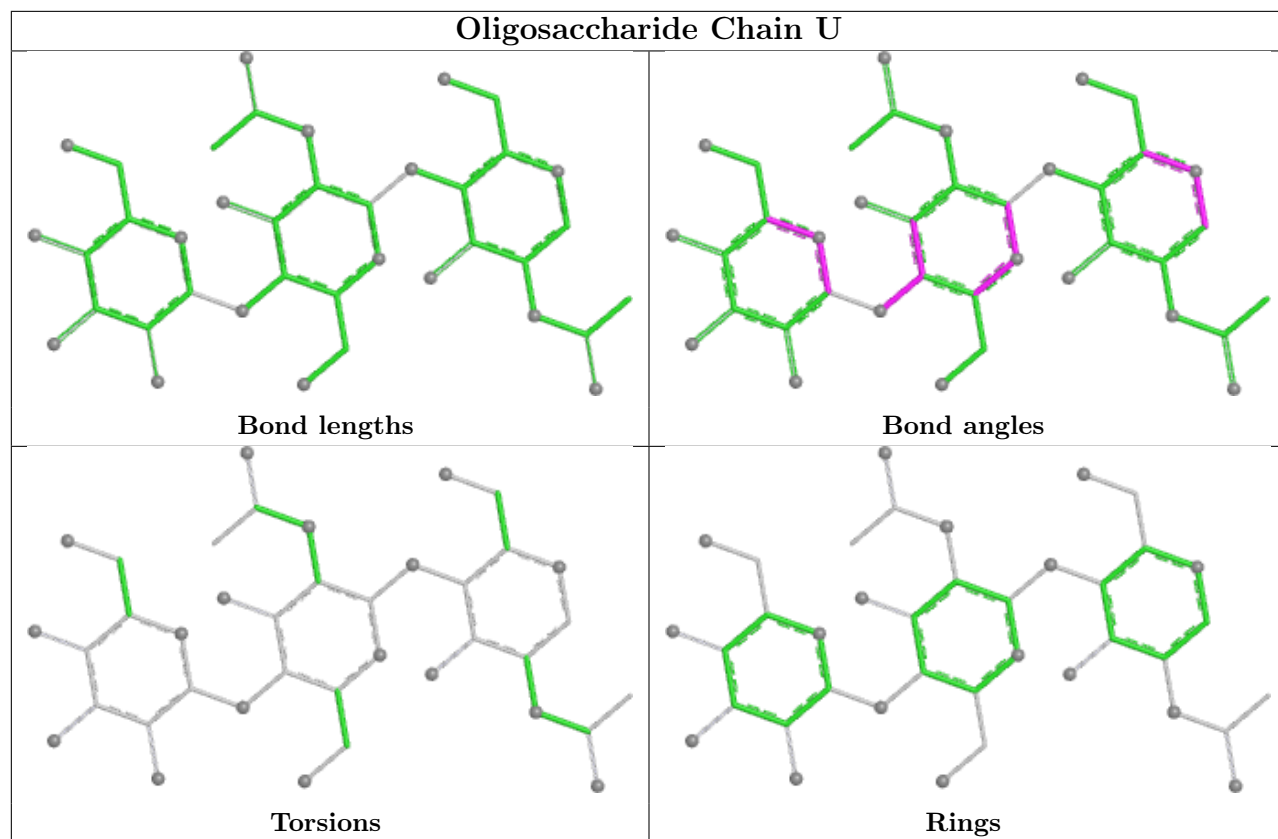
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

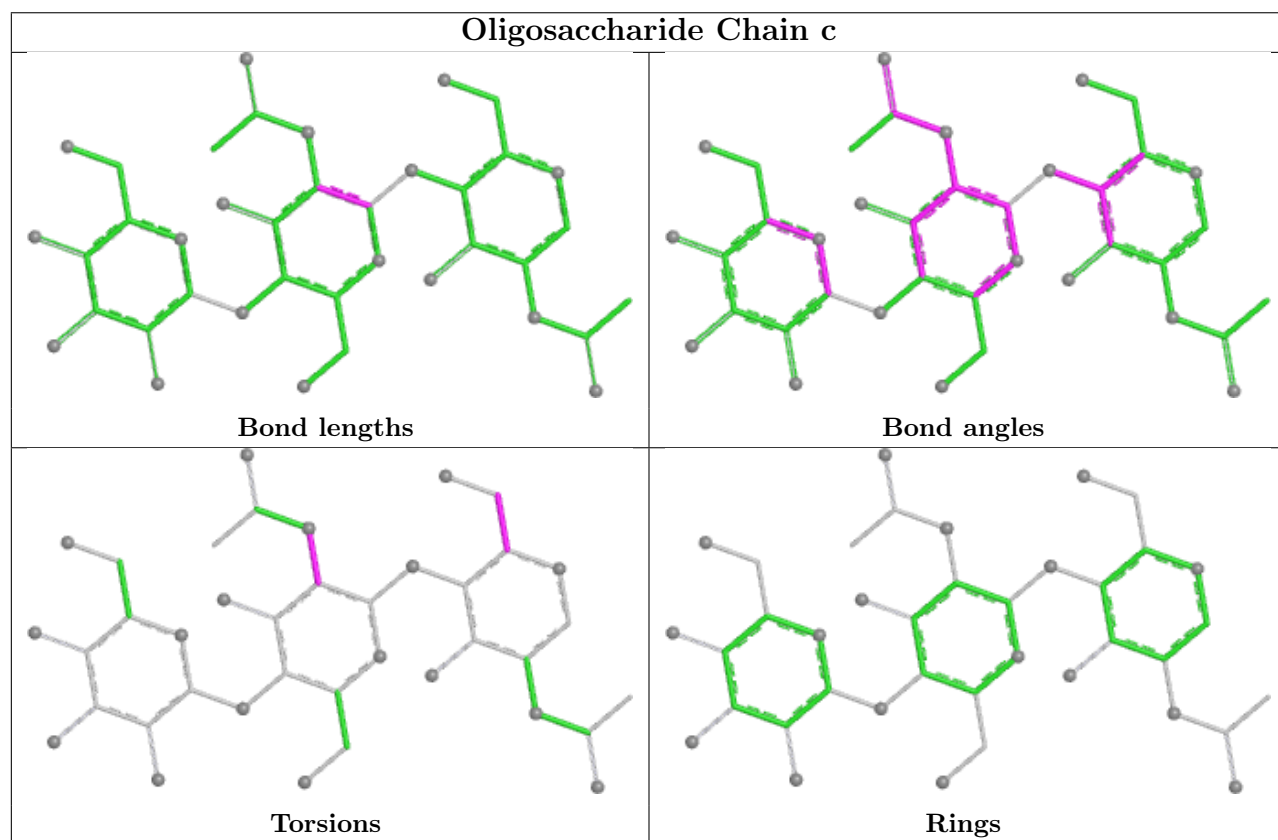
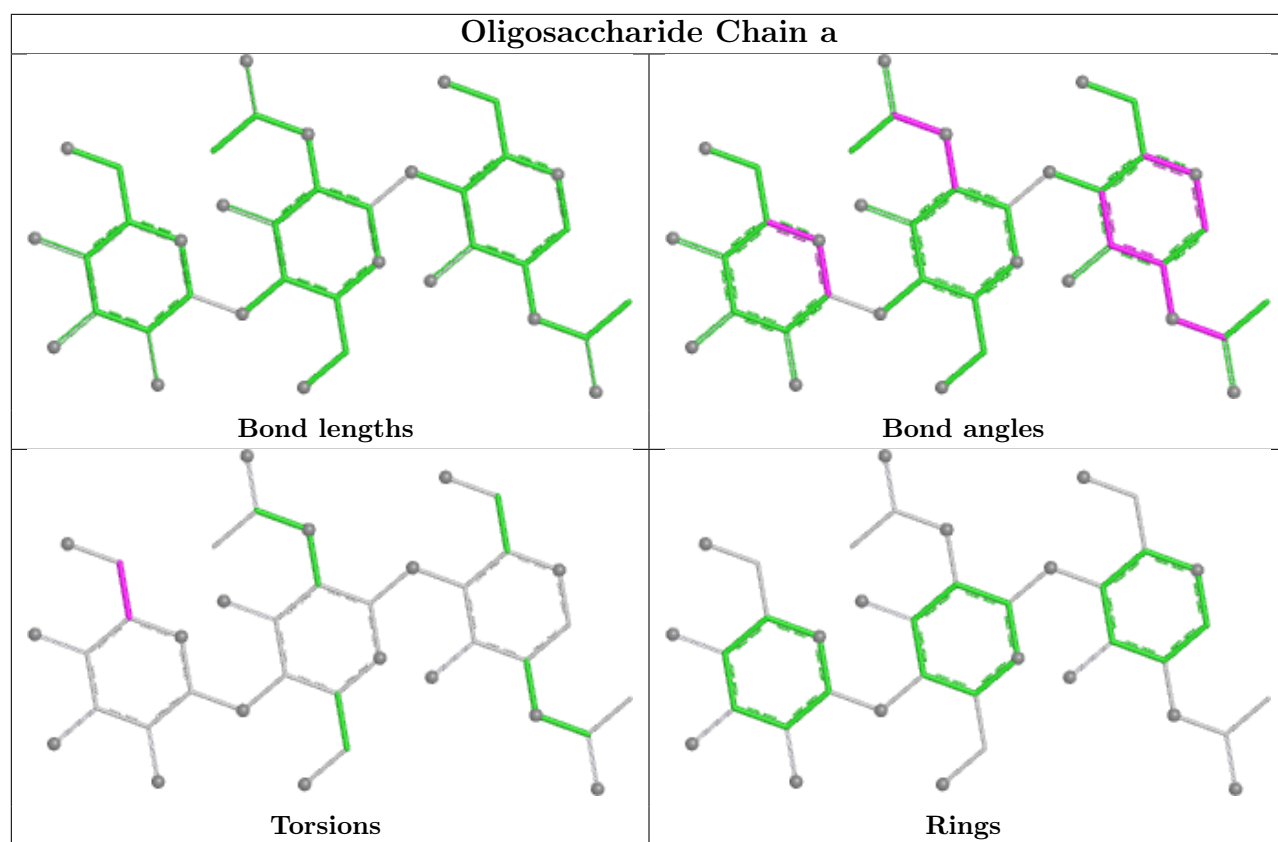


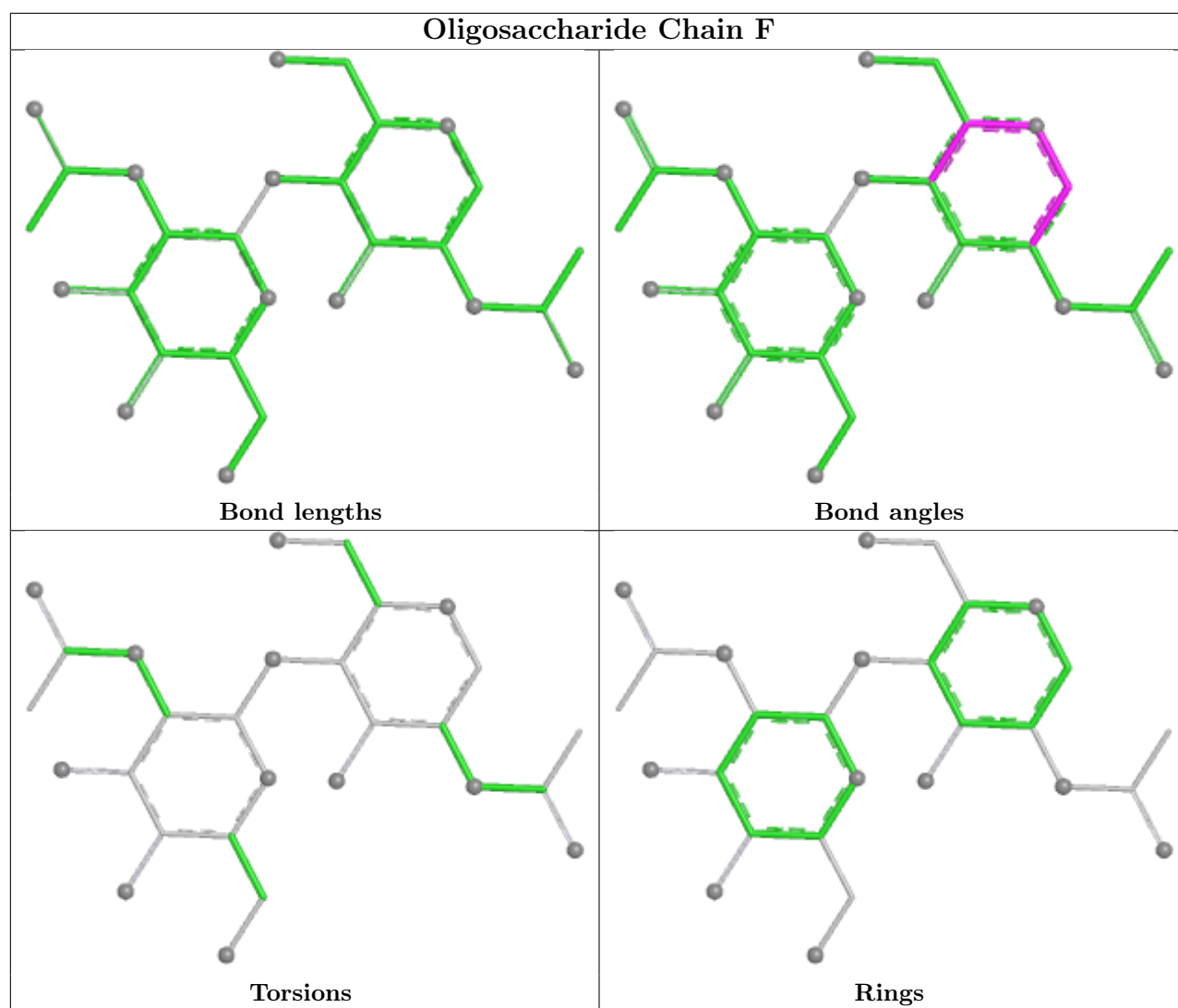


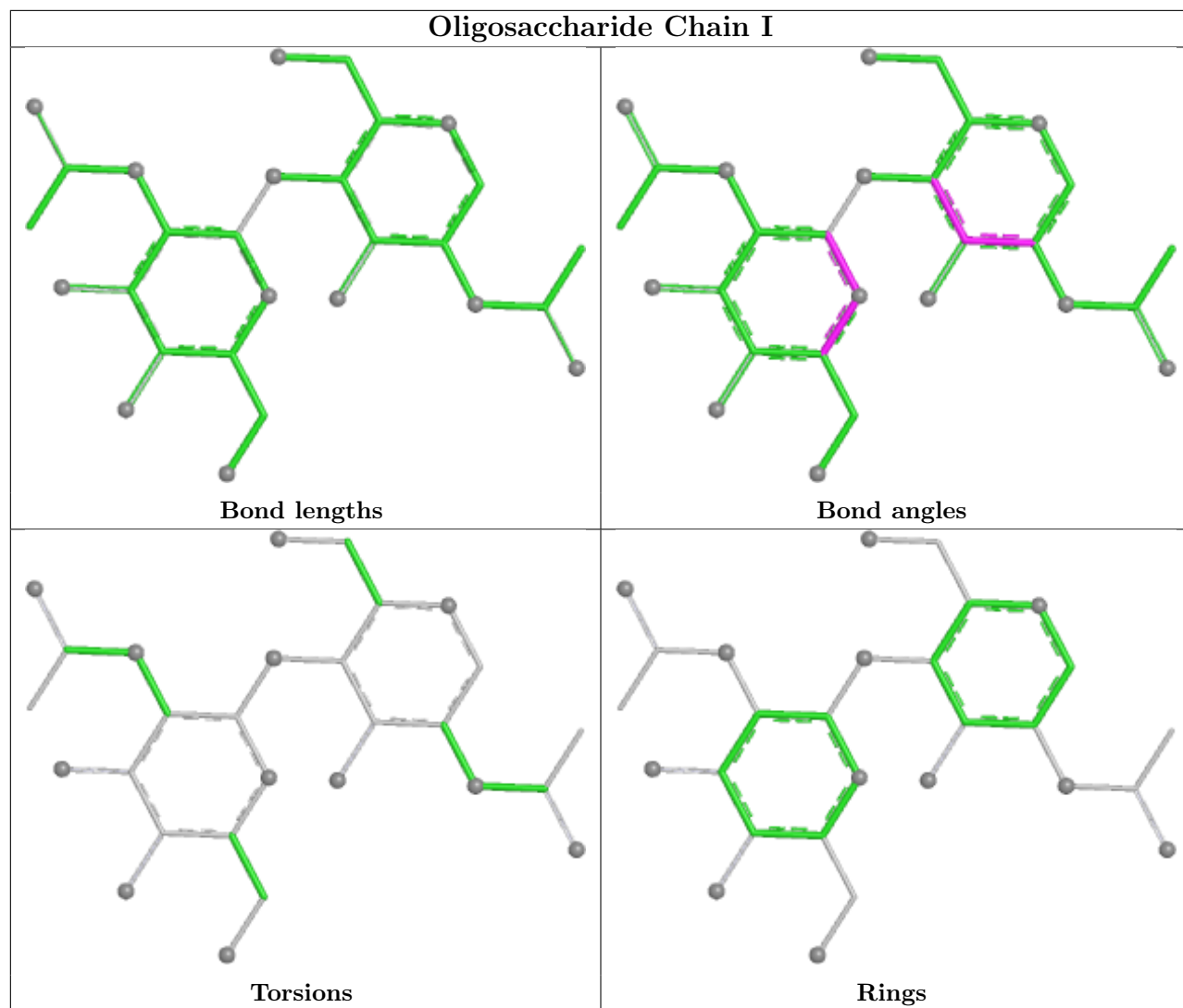


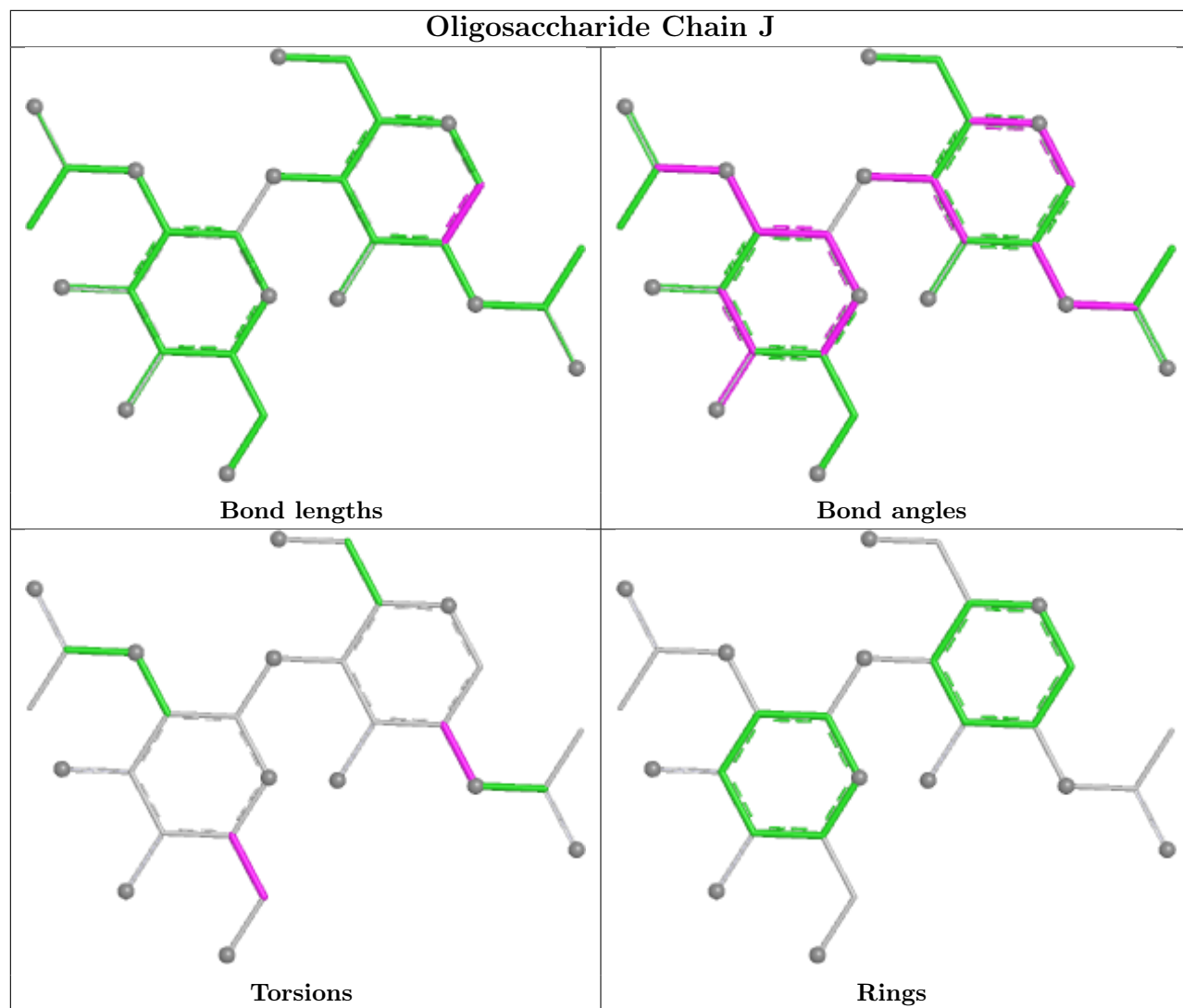


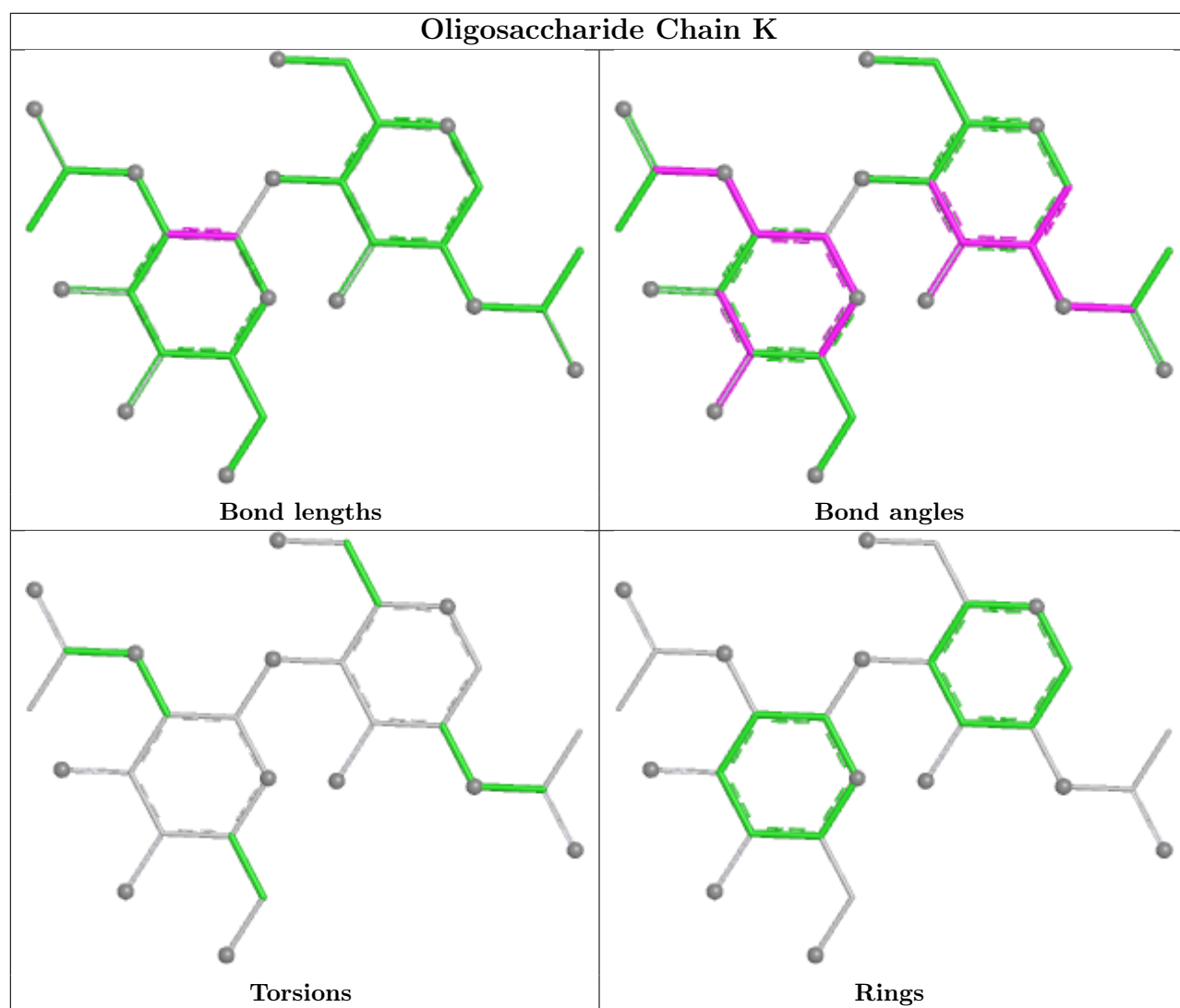


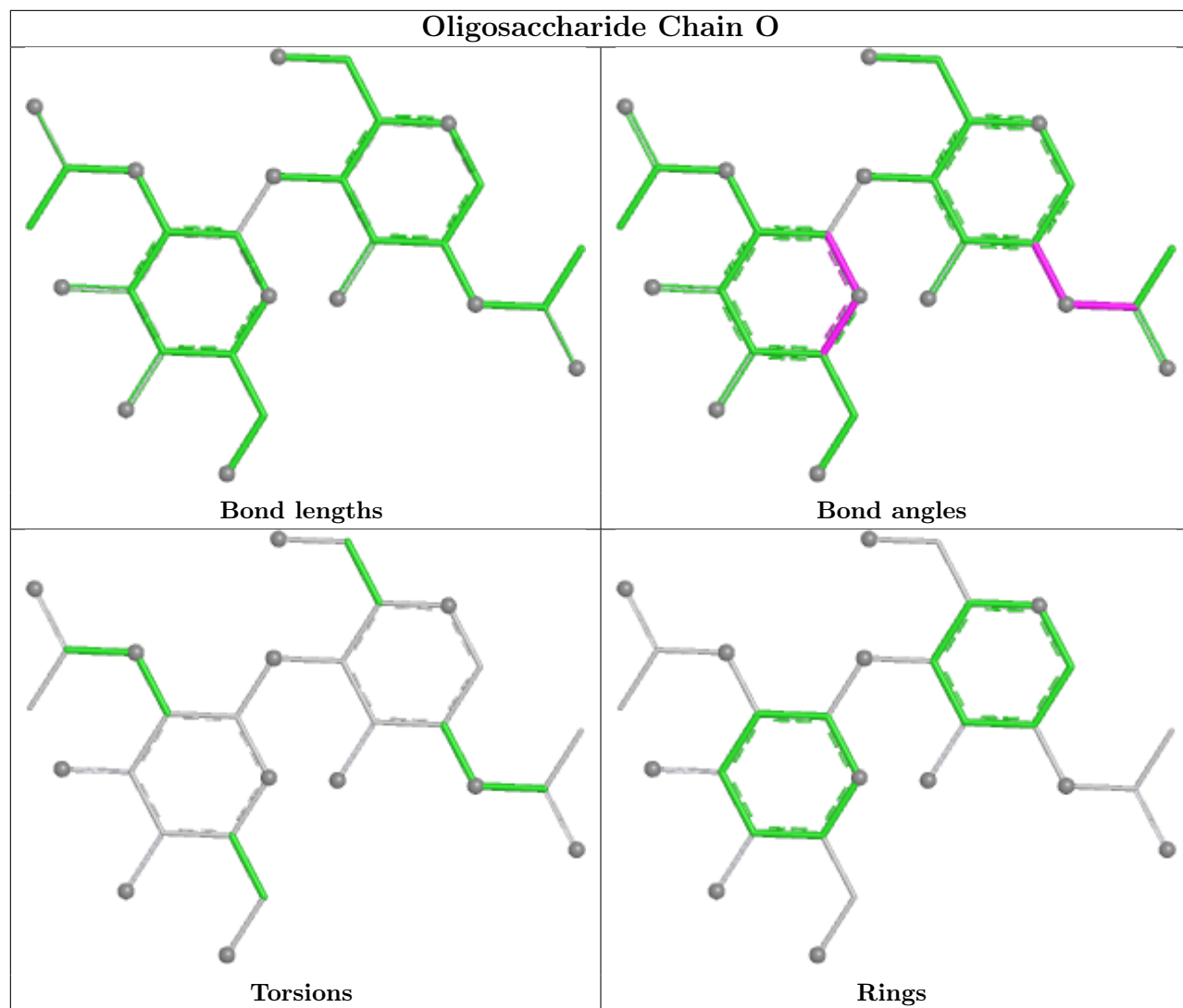


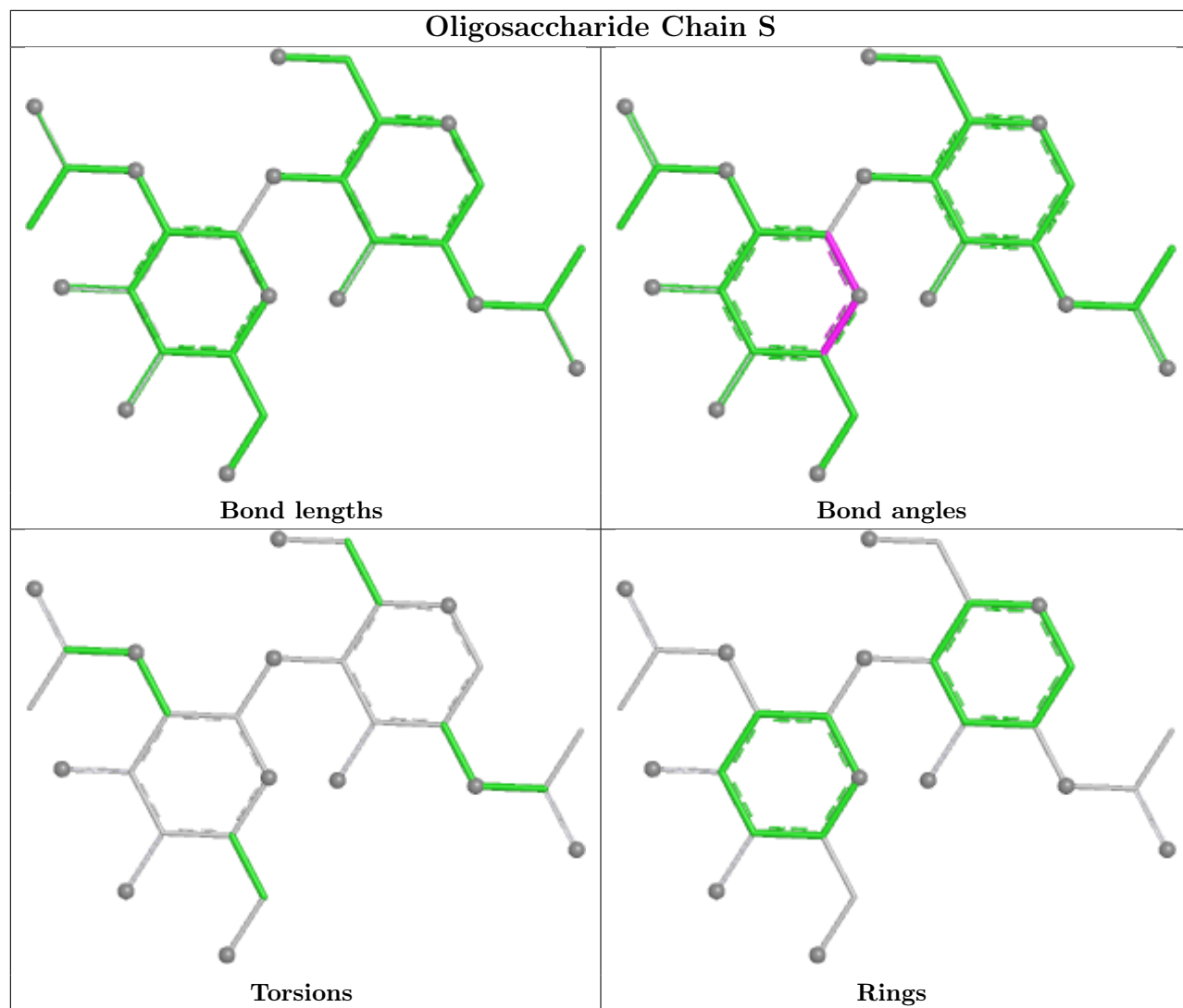


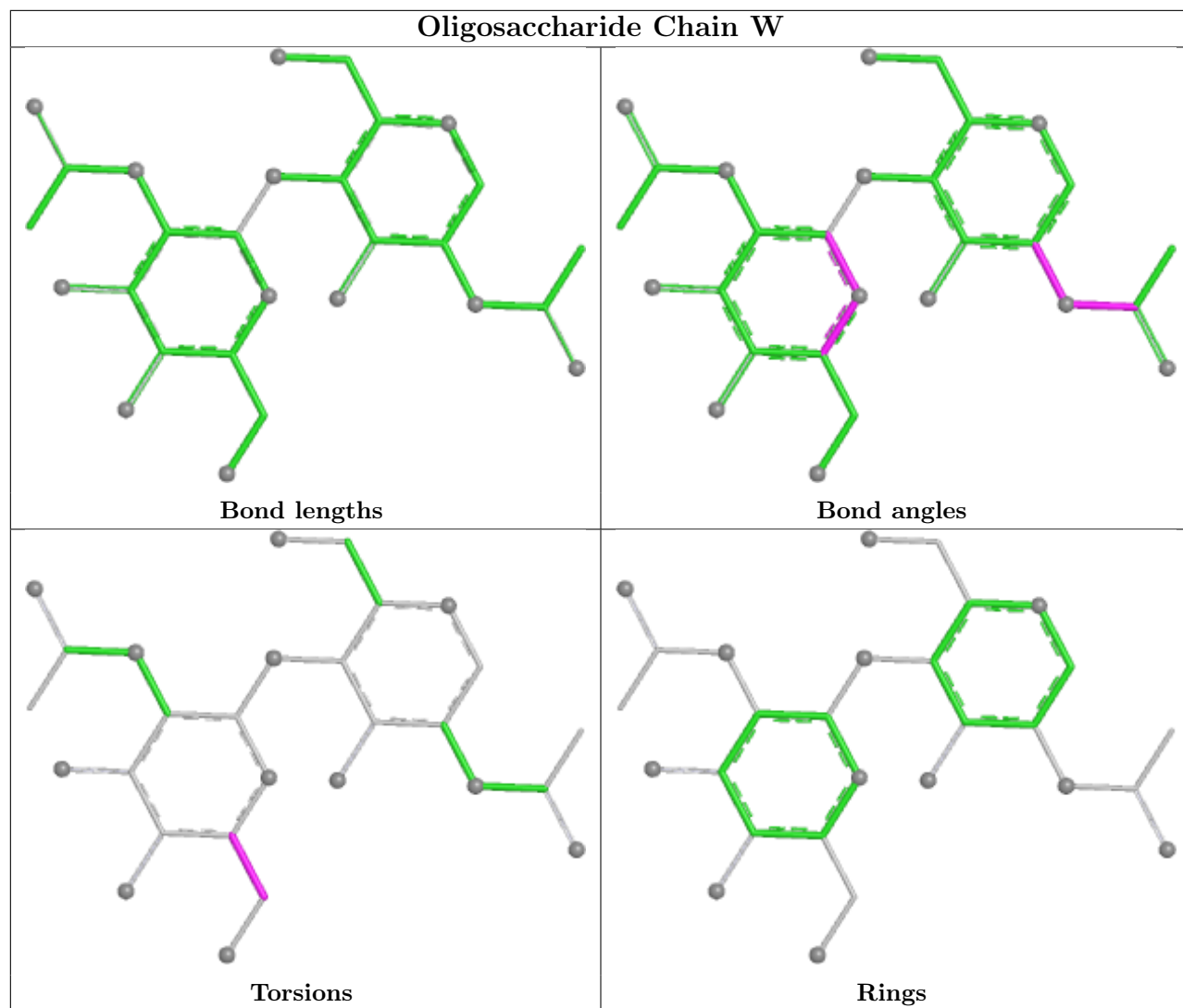


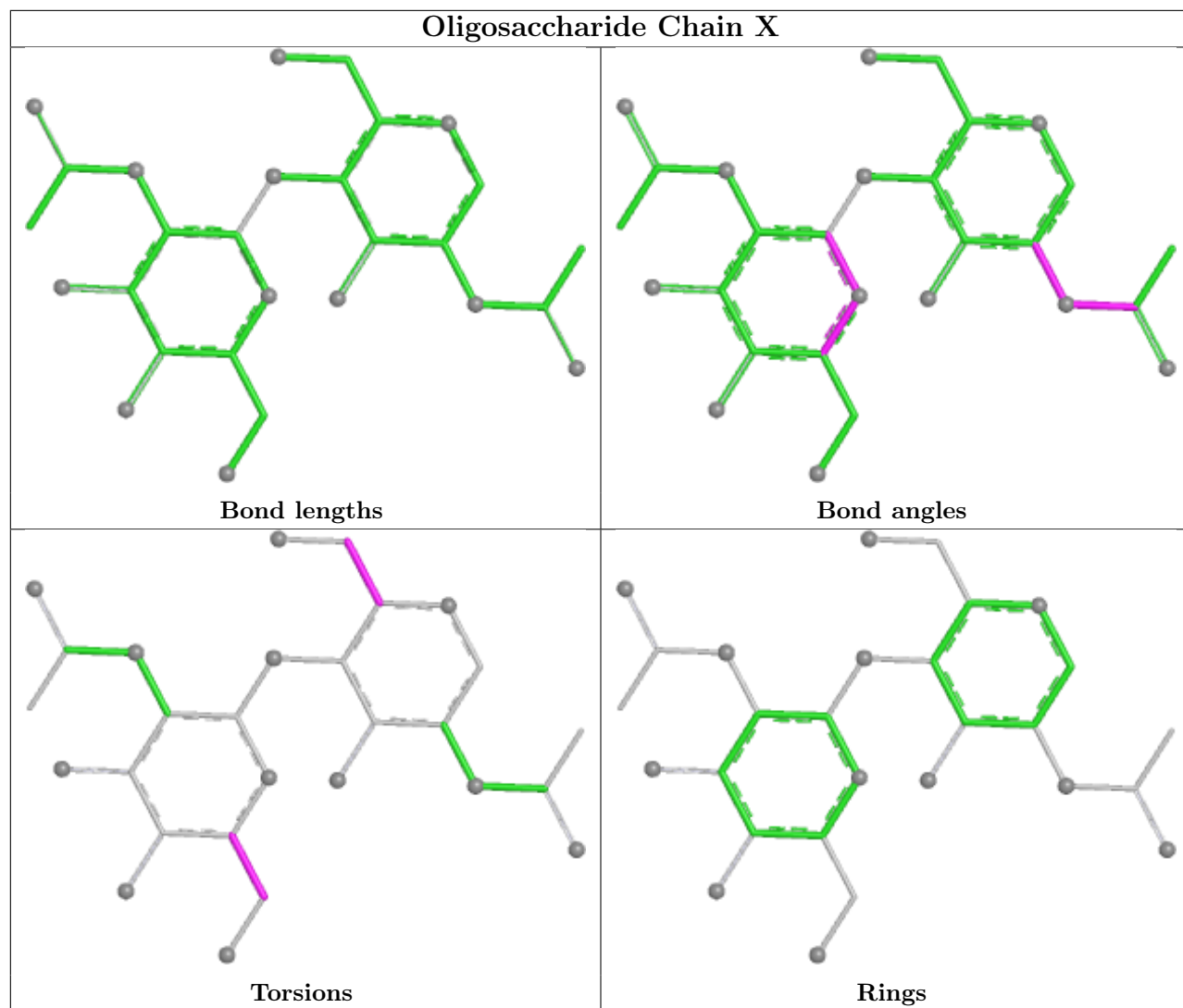


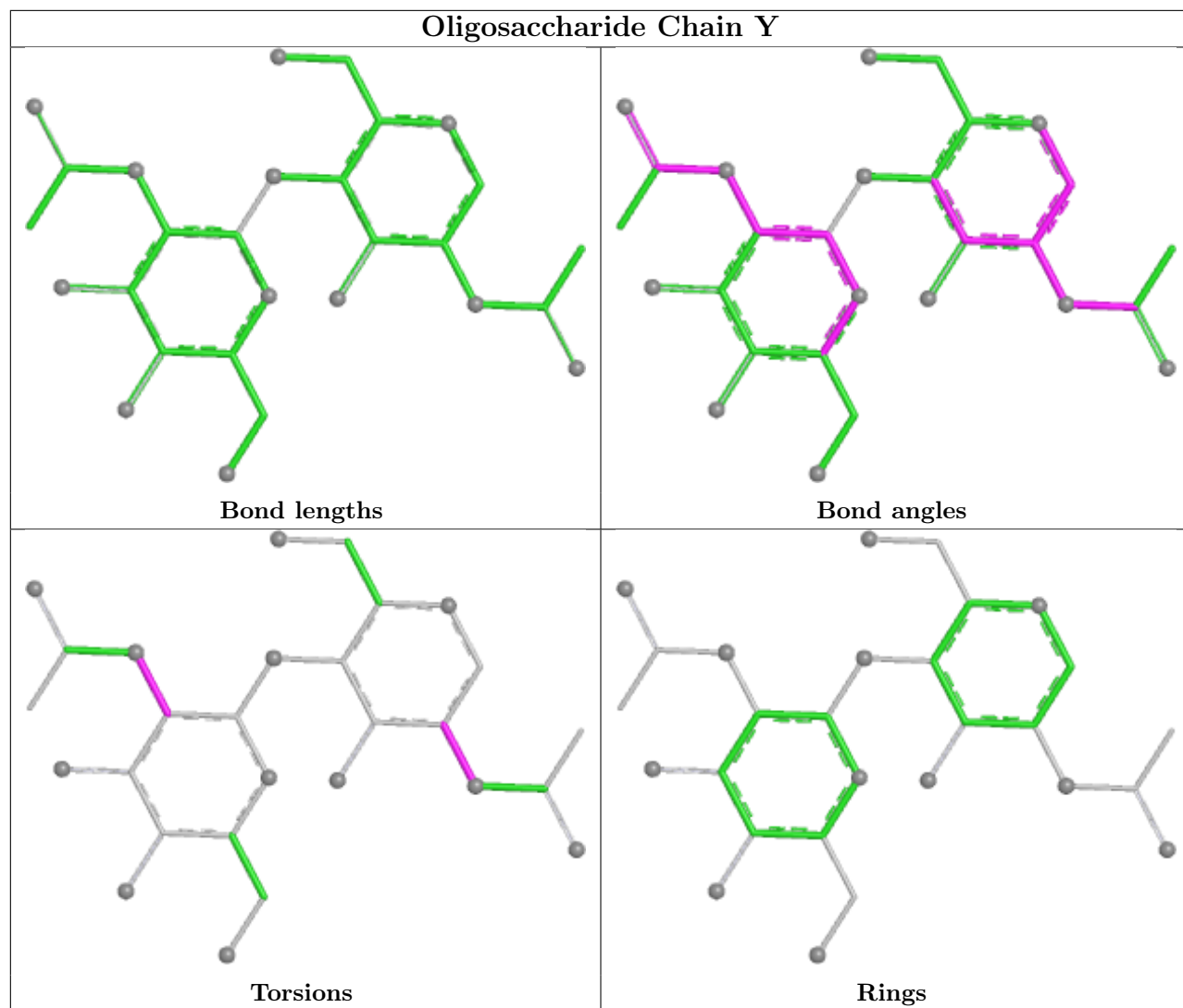


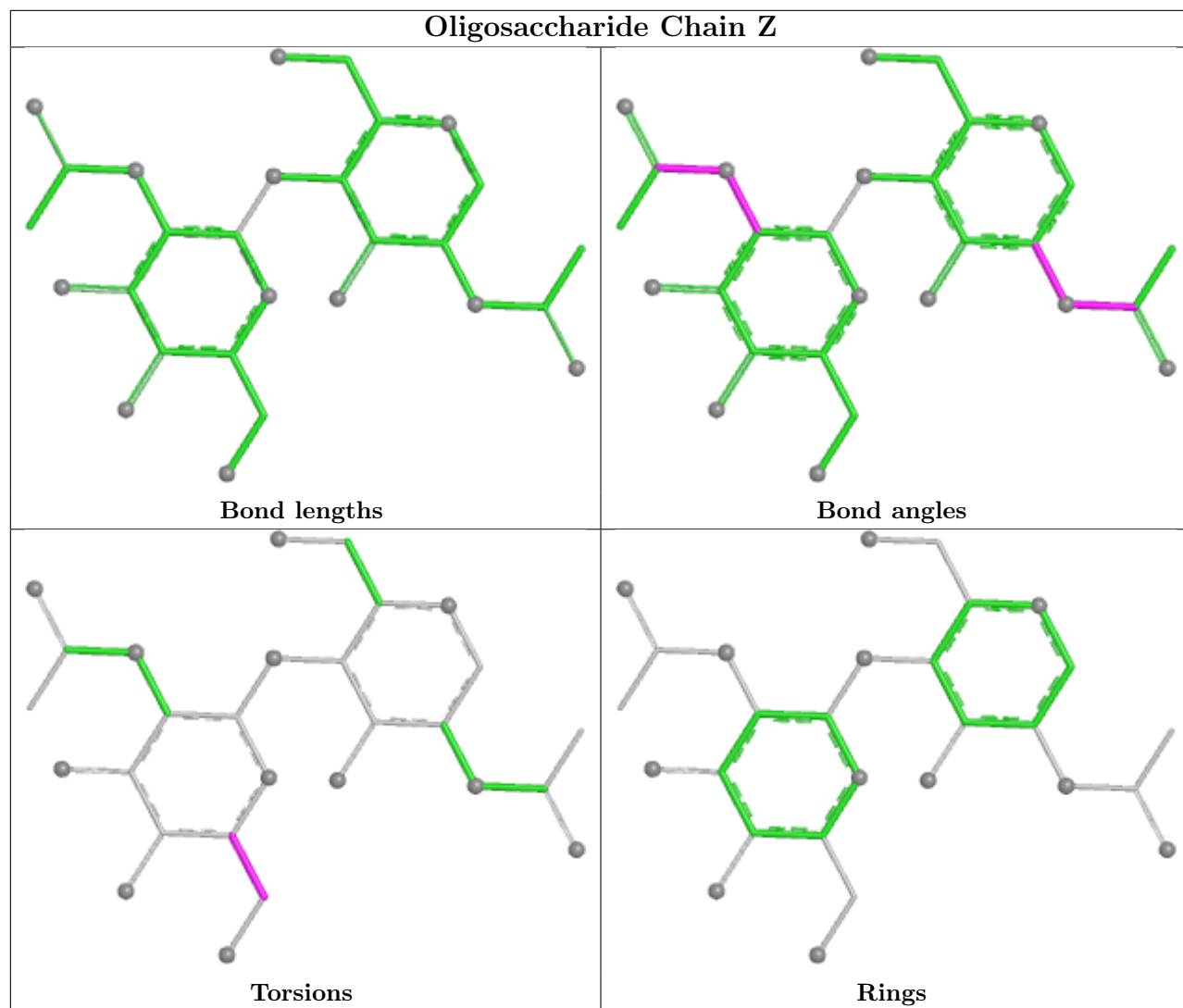


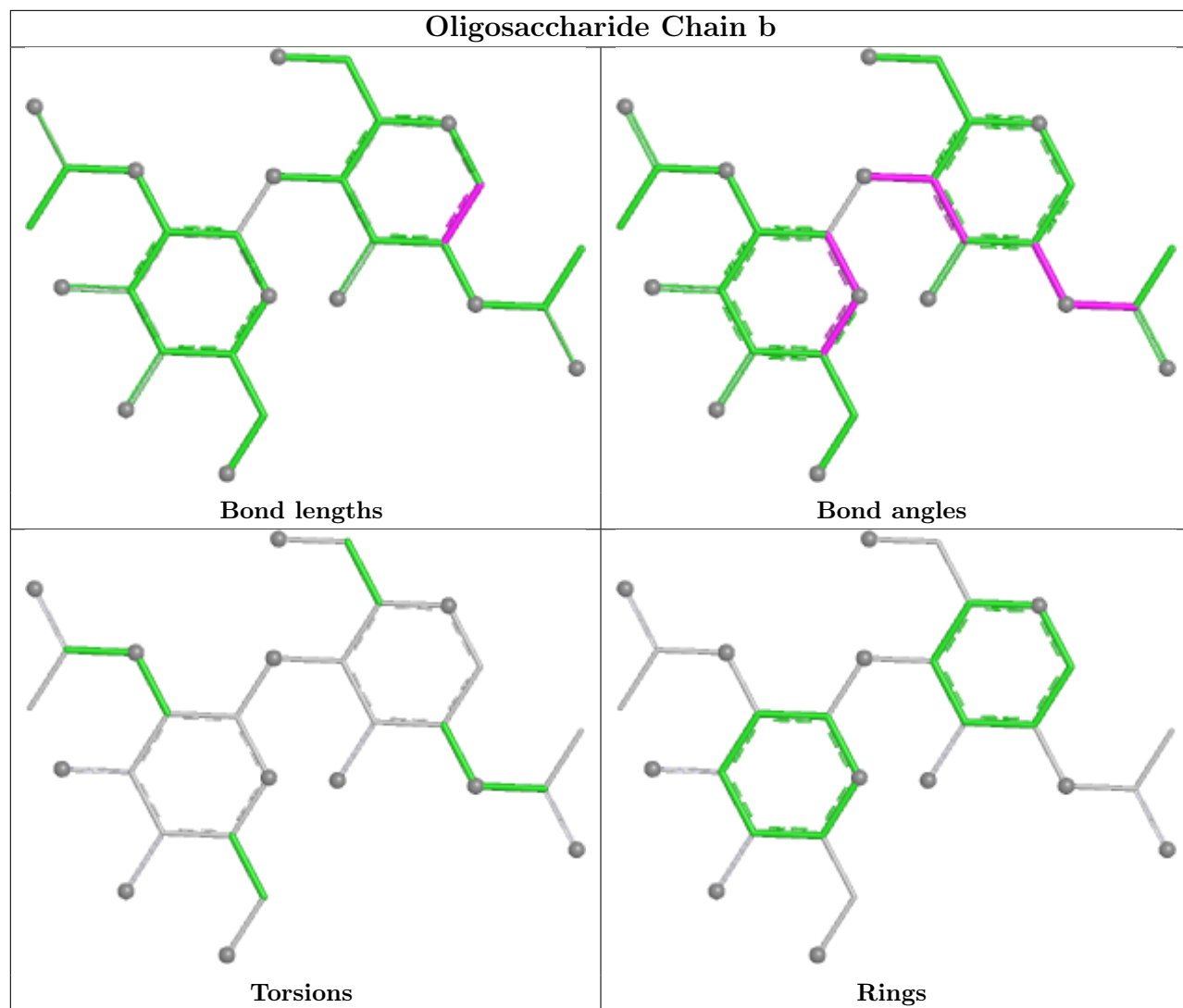


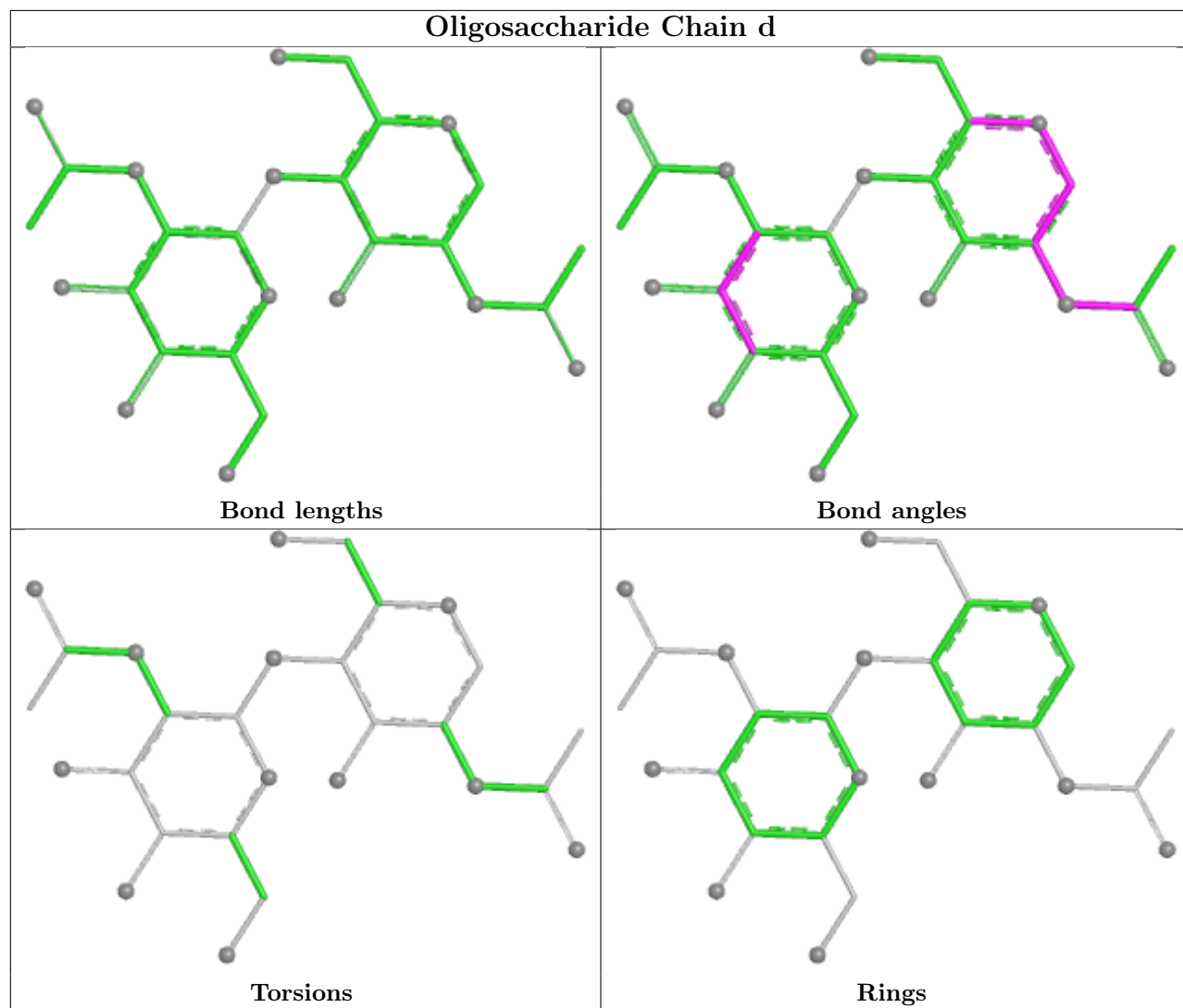


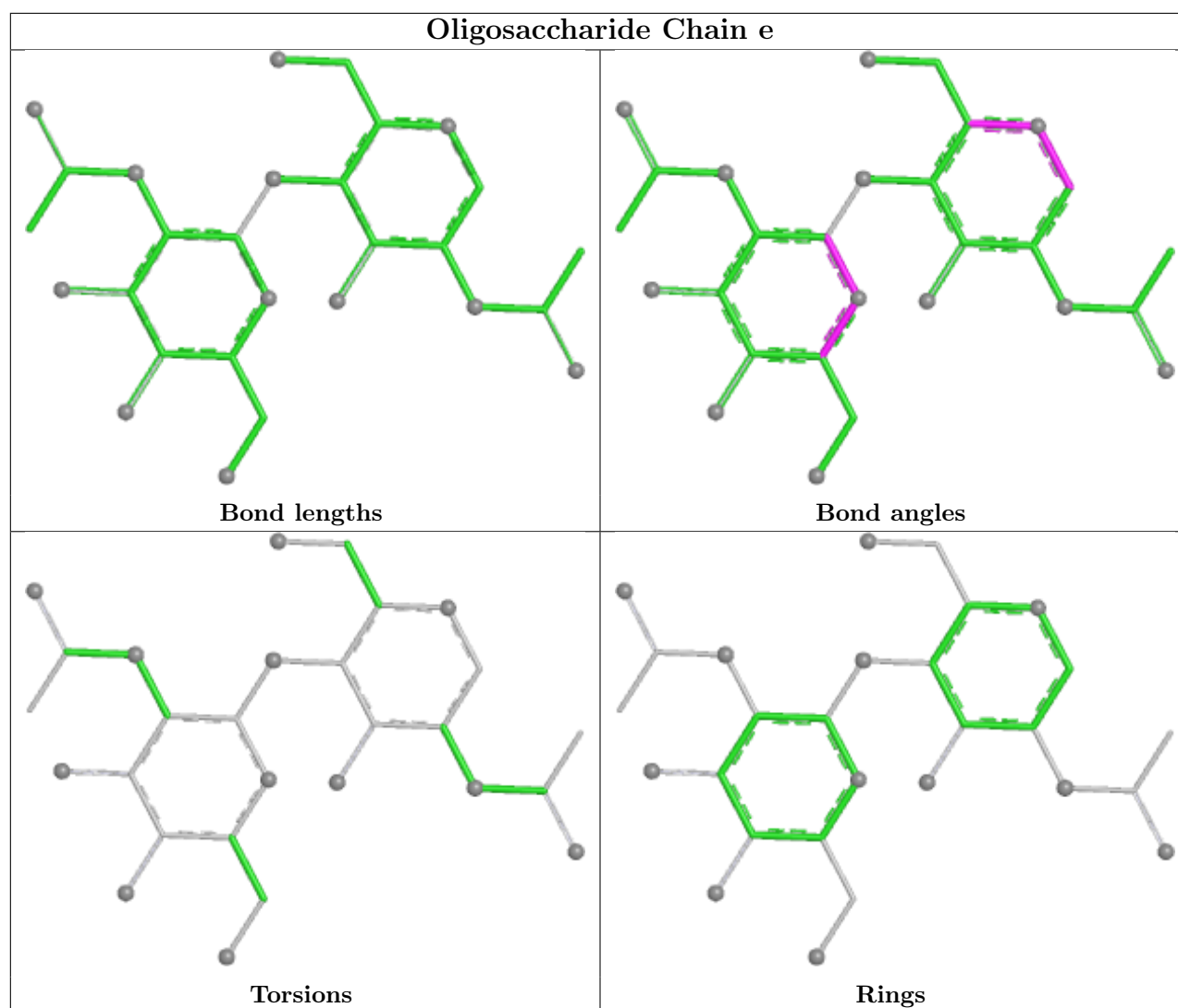












5.6 Ligand geometry [i](#)

45 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	NAG	A	1208	1	14,14,15	1.46	3 (21%)	17,19,21	1.83	2 (11%)
6	NAG	B	1214	1	14,14,15	0.95	0	17,19,21	0.83	1 (5%)
6	NAG	A	1210	1	14,14,15	0.60	0	17,19,21	1.55	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	1214	1	14,14,15	0.63	0	17,19,21	1.77	4 (23%)
6	NAG	A	1215	1	14,14,15	0.77	0	17,19,21	0.74	0
6	NAG	B	1209	1	14,14,15	0.71	0	17,19,21	1.41	2 (11%)
6	NAG	C	1202	1	14,14,15	0.77	0	17,19,21	1.46	2 (11%)
6	NAG	A	1201	1	14,14,15	0.69	0	17,19,21	1.12	2 (11%)
6	NAG	A	1216	1	14,14,15	0.66	0	17,19,21	1.36	3 (17%)
6	NAG	A	1207	1	14,14,15	0.68	0	17,19,21	1.03	2 (11%)
6	NAG	B	1208	1	14,14,15	0.70	0	17,19,21	1.13	1 (5%)
6	NAG	A	1206	1	14,14,15	0.68	0	17,19,21	1.16	2 (11%)
6	NAG	B	1205	1	14,14,15	0.86	0	17,19,21	1.41	2 (11%)
6	NAG	B	1212	1	14,14,15	0.76	0	17,19,21	1.31	2 (11%)
6	NAG	A	1205	1	14,14,15	0.66	0	17,19,21	1.33	2 (11%)
6	NAG	C	1215	1	14,14,15	0.72	0	17,19,21	2.67	4 (23%)
6	NAG	A	1203	1	14,14,15	0.82	0	17,19,21	1.51	1 (5%)
6	NAG	A	1211	1	14,14,15	0.60	0	17,19,21	1.82	3 (17%)
6	NAG	A	1212	1	14,14,15	0.72	0	17,19,21	0.98	0
6	NAG	A	1209	1	14,14,15	0.67	0	17,19,21	1.34	4 (23%)
6	NAG	B	1213	1	14,14,15	0.70	0	17,19,21	1.37	1 (5%)
6	NAG	A	1213	1	14,14,15	0.76	0	17,19,21	1.75	3 (17%)
6	NAG	C	1210	1	14,14,15	0.76	1 (7%)	17,19,21	1.66	2 (11%)
6	NAG	B	1202	1	14,14,15	0.75	0	17,19,21	0.99	1 (5%)
6	NAG	B	1204	1	14,14,15	0.71	0	17,19,21	1.05	2 (11%)
6	NAG	A	1204	1	14,14,15	0.89	1 (7%)	17,19,21	1.31	2 (11%)
6	NAG	B	1203	1	14,14,15	0.80	0	17,19,21	1.41	1 (5%)
6	NAG	B	1207	1	14,14,15	0.74	0	17,19,21	0.97	0
6	NAG	C	1201	1	14,14,15	0.52	0	17,19,21	1.42	2 (11%)
6	NAG	C	1207	1	14,14,15	0.68	0	17,19,21	1.67	4 (23%)
6	NAG	C	1205	1	14,14,15	0.75	0	17,19,21	0.95	1 (5%)
6	NAG	C	1203	1	14,14,15	0.73	0	17,19,21	1.04	2 (11%)
6	NAG	C	1212	1	14,14,15	0.90	0	17,19,21	1.48	3 (17%)
6	NAG	B	1210	1	14,14,15	0.70	0	17,19,21	0.78	0
6	NAG	C	1211	1	14,14,15	0.75	0	17,19,21	1.44	1 (5%)
6	NAG	C	1213	1	14,14,15	0.73	0	17,19,21	0.97	1 (5%)
6	NAG	C	1206	1	14,14,15	0.70	0	17,19,21	1.12	1 (5%)
6	NAG	A	1202	1	14,14,15	0.70	0	17,19,21	1.32	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	C	1209	1	14,14,15	0.74	0	17,19,21	1.34	3 (17%)
6	NAG	C	1214	1	14,14,15	0.73	0	17,19,21	1.31	2 (11%)
6	NAG	B	1211	1	14,14,15	0.78	0	17,19,21	0.89	1 (5%)
6	NAG	C	1204	1	14,14,15	0.70	0	17,19,21	1.48	1 (5%)
6	NAG	B	1201	1	14,14,15	0.70	0	17,19,21	0.91	0
6	NAG	C	1208	1	14,14,15	0.89	0	17,19,21	1.79	5 (29%)
6	NAG	B	1206	-	14,14,15	0.77	0	17,19,21	1.71	4 (23%)

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	A	1208	1	-	0/6/23/26	0/1/1/1
6	NAG	B	1214	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1210	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1214	1	-	3/6/23/26	0/1/1/1
6	NAG	A	1215	1	-	0/6/23/26	0/1/1/1
6	NAG	B	1209	1	-	2/6/23/26	0/1/1/1
6	NAG	C	1202	1	-	2/6/23/26	0/1/1/1
6	NAG	A	1201	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1216	1	-	0/6/23/26	0/1/1/1
6	NAG	A	1207	1	-	1/6/23/26	0/1/1/1
6	NAG	B	1208	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1206	1	-	0/6/23/26	0/1/1/1
6	NAG	B	1205	1	-	0/6/23/26	0/1/1/1
6	NAG	B	1212	1	-	0/6/23/26	0/1/1/1
6	NAG	A	1205	1	-	0/6/23/26	0/1/1/1
6	NAG	C	1215	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1203	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1211	1	-	0/6/23/26	0/1/1/1
6	NAG	A	1212	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1209	1	-	1/6/23/26	0/1/1/1
6	NAG	B	1213	1	-	0/6/23/26	0/1/1/1
6	NAG	A	1213	1	-	1/6/23/26	0/1/1/1
6	NAG	C	1210	1	-	2/6/23/26	0/1/1/1
6	NAG	B	1202	1	-	1/6/23/26	0/1/1/1
6	NAG	B	1204	1	-	0/6/23/26	0/1/1/1
6	NAG	A	1204	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	B	1203	1	-	0/6/23/26	0/1/1/1
6	NAG	B	1207	1	-	2/6/23/26	0/1/1/1
6	NAG	C	1201	1	-	1/6/23/26	0/1/1/1
6	NAG	C	1207	1	-	1/6/23/26	0/1/1/1
6	NAG	C	1205	1	-	1/6/23/26	0/1/1/1
6	NAG	C	1203	1	-	0/6/23/26	0/1/1/1
6	NAG	C	1212	1	-	3/6/23/26	0/1/1/1
6	NAG	B	1210	1	-	0/6/23/26	0/1/1/1
6	NAG	C	1211	1	-	1/6/23/26	0/1/1/1
6	NAG	C	1213	1	-	0/6/23/26	0/1/1/1
6	NAG	C	1206	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1202	1	-	2/6/23/26	0/1/1/1
6	NAG	C	1209	1	-	3/6/23/26	0/1/1/1
6	NAG	C	1214	1	-	1/6/23/26	0/1/1/1
6	NAG	B	1211	1	-	0/6/23/26	0/1/1/1
6	NAG	C	1204	1	-	0/6/23/26	0/1/1/1
6	NAG	B	1201	1	-	1/6/23/26	0/1/1/1
6	NAG	C	1208	1	-	1/6/23/26	0/1/1/1
6	NAG	B	1206	-	-	3/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1208	NAG	C1-C2	3.75	1.57	1.52
6	A	1204	NAG	C1-C2	2.33	1.55	1.52
6	A	1208	NAG	C3-C2	-2.32	1.47	1.52
6	A	1208	NAG	O4-C4	-2.08	1.37	1.43
6	C	1210	NAG	C1-C2	2.07	1.55	1.52

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	1215	NAG	C2-N2-C7	8.20	133.88	122.90
6	A	1211	NAG	C1-O5-C5	5.61	119.71	112.19
6	A	1208	NAG	C1-O5-C5	5.57	119.66	112.19
6	C	1204	NAG	C1-O5-C5	5.04	118.94	112.19
6	A	1210	NAG	C1-O5-C5	4.98	118.86	112.19
6	C	1215	NAG	C1-O5-C5	4.94	118.80	112.19
6	A	1213	NAG	O5-C1-C2	-4.79	103.88	111.29
6	A	1203	NAG	C2-N2-C7	4.79	129.31	122.90
6	C	1208	NAG	C2-N2-C7	4.74	129.25	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	1210	NAG	C1-O5-C5	4.38	118.05	112.19
6	A	1214	NAG	C1-O5-C5	4.07	117.64	112.19
6	C	1211	NAG	C2-N2-C7	4.07	128.35	122.90
6	B	1203	NAG	C2-N2-C7	4.01	128.27	122.90
6	A	1205	NAG	C1-O5-C5	3.88	117.38	112.19
6	B	1205	NAG	C2-N2-C7	3.84	128.04	122.90
6	B	1206	NAG	C1-O5-C5	3.73	117.18	112.19
6	C	1210	NAG	C2-N2-C7	3.72	127.89	122.90
6	C	1202	NAG	C1-O5-C5	3.72	117.17	112.19
6	A	1204	NAG	C1-O5-C5	3.72	117.17	112.19
6	B	1212	NAG	C1-O5-C5	3.70	117.14	112.19
6	C	1201	NAG	C4-C3-C2	-3.61	105.73	111.02
6	B	1213	NAG	C1-O5-C5	3.60	117.01	112.19
6	C	1212	NAG	C1-O5-C5	3.50	116.87	112.19
6	A	1214	NAG	C2-N2-C7	3.46	127.54	122.90
6	C	1202	NAG	C2-N2-C7	3.34	127.37	122.90
6	A	1211	NAG	C2-N2-C7	3.32	127.35	122.90
6	C	1207	NAG	C2-N2-C7	3.28	127.29	122.90
6	C	1201	NAG	O5-C1-C2	-3.26	106.25	111.29
6	B	1209	NAG	C2-N2-C7	3.26	127.26	122.90
6	B	1206	NAG	C1-C2-N2	3.24	115.54	110.43
6	A	1214	NAG	C4-C3-C2	-3.22	106.29	111.02
6	A	1216	NAG	C1-O5-C5	3.12	116.37	112.19
6	A	1213	NAG	C1-C2-N2	3.11	115.34	110.43
6	A	1202	NAG	C2-N2-C7	3.10	127.06	122.90
6	A	1208	NAG	C4-C3-C2	-2.95	106.70	111.02
6	A	1210	NAG	C4-C3-C2	-2.94	106.71	111.02
6	A	1213	NAG	C2-N2-C7	2.94	126.84	122.90
6	C	1206	NAG	C1-O5-C5	2.94	116.12	112.19
6	C	1207	NAG	C4-C3-C2	-2.92	106.74	111.02
6	B	1209	NAG	C1-O5-C5	2.92	116.10	112.19
6	B	1208	NAG	C1-O5-C5	2.90	116.08	112.19
6	C	1207	NAG	C1-O5-C5	2.88	116.05	112.19
6	C	1214	NAG	C1-O5-C5	2.88	116.04	112.19
6	A	1202	NAG	C4-C3-C2	-2.87	106.81	111.02
6	C	1208	NAG	C1-O5-C5	2.84	115.99	112.19
6	A	1201	NAG	C1-O5-C5	2.84	115.99	112.19
6	C	1215	NAG	C1-C2-N2	2.81	114.86	110.43
6	A	1206	NAG	C1-O5-C5	2.80	115.94	112.19
6	C	1214	NAG	C2-N2-C7	2.76	126.59	122.90
6	C	1215	NAG	O7-C7-N2	2.74	126.82	121.98
6	A	1209	NAG	C1-O5-C5	2.72	115.83	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	1203	NAG	O5-C1-C2	-2.72	107.09	111.29
6	B	1204	NAG	C1-O5-C5	2.71	115.82	112.19
6	A	1206	NAG	C4-C3-C2	-2.66	107.12	111.02
6	B	1206	NAG	C2-N2-C7	2.63	126.43	122.90
6	C	1209	NAG	O5-C1-C2	-2.63	107.22	111.29
6	C	1212	NAG	O5-C1-C2	-2.55	107.34	111.29
6	A	1205	NAG	C4-C3-C2	-2.54	107.30	111.02
6	C	1208	NAG	O5-C1-C2	-2.53	107.37	111.29
6	A	1204	NAG	C2-N2-C7	2.48	126.23	122.90
6	C	1209	NAG	C2-N2-C7	2.48	126.22	122.90
6	B	1202	NAG	C2-N2-C7	2.46	126.20	122.90
6	C	1208	NAG	C8-C7-N2	-2.44	112.07	116.12
6	C	1203	NAG	C2-N2-C7	2.40	126.12	122.90
6	A	1216	NAG	C2-N2-C7	2.40	126.12	122.90
6	A	1209	NAG	C4-C3-C2	-2.35	107.58	111.02
6	A	1209	NAG	O5-C1-C2	-2.34	107.66	111.29
6	A	1209	NAG	C1-C2-N2	2.31	114.07	110.43
6	C	1209	NAG	C1-C2-N2	2.29	114.05	110.43
6	B	1205	NAG	C1-O5-C5	2.27	115.23	112.19
6	C	1213	NAG	C4-C3-C2	-2.25	107.73	111.02
6	C	1208	NAG	O7-C7-N2	2.23	125.92	121.98
6	A	1211	NAG	O4-C4-C3	-2.22	105.15	110.38
6	B	1212	NAG	C4-C3-C2	-2.19	107.81	111.02
6	B	1214	NAG	C3-C4-C5	2.17	114.17	110.23
6	A	1216	NAG	C1-C2-N2	2.15	113.82	110.43
6	C	1205	NAG	C3-C4-C5	2.12	114.08	110.23
6	A	1207	NAG	C2-N2-C7	2.09	125.70	122.90
6	B	1204	NAG	C2-N2-C7	2.08	125.69	122.90
6	C	1212	NAG	C4-C3-C2	-2.08	107.97	111.02
6	A	1214	NAG	C1-C2-N2	2.08	113.71	110.43
6	B	1206	NAG	C4-C3-C2	-2.08	107.98	111.02
6	A	1207	NAG	C1-O5-C5	2.06	114.95	112.19
6	B	1211	NAG	O5-C5-C4	-2.02	105.92	110.83
6	A	1201	NAG	O5-C1-C2	-2.01	108.19	111.29
6	C	1207	NAG	O5-C1-C2	-2.01	108.19	111.29

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1214	NAG	C1-C2-N2-C7
6	C	1202	NAG	C1-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
6	C	1207	NAG	C3-C2-N2-C7
6	C	1210	NAG	C1-C2-N2-C7
6	C	1215	NAG	C1-C2-N2-C7
6	C	1212	NAG	O5-C5-C6-O6
6	C	1212	NAG	C4-C5-C6-O6
6	A	1203	NAG	O5-C5-C6-O6
6	B	1202	NAG	O5-C5-C6-O6
6	B	1207	NAG	O5-C5-C6-O6
6	B	1207	NAG	C4-C5-C6-O6
6	A	1201	NAG	O5-C5-C6-O6
6	C	1201	NAG	O5-C5-C6-O6
6	C	1206	NAG	O5-C5-C6-O6
6	A	1207	NAG	O5-C5-C6-O6
6	A	1212	NAG	O5-C5-C6-O6
6	B	1201	NAG	O5-C5-C6-O6
6	A	1214	NAG	O5-C5-C6-O6
6	C	1214	NAG	O5-C5-C6-O6
6	C	1209	NAG	O5-C5-C6-O6
6	B	1208	NAG	O5-C5-C6-O6
6	B	1214	NAG	O5-C5-C6-O6
6	A	1210	NAG	O5-C5-C6-O6
6	C	1202	NAG	O5-C5-C6-O6
6	C	1205	NAG	O5-C5-C6-O6
6	A	1209	NAG	O5-C5-C6-O6
6	A	1202	NAG	C1-C2-N2-C7
6	A	1204	NAG	C1-C2-N2-C7
6	B	1209	NAG	C1-C2-N2-C7
6	C	1209	NAG	C1-C2-N2-C7
6	A	1202	NAG	C3-C2-N2-C7
6	A	1213	NAG	C3-C2-N2-C7
6	A	1214	NAG	C3-C2-N2-C7
6	B	1206	NAG	C3-C2-N2-C7
6	B	1206	NAG	O5-C5-C6-O6
6	B	1206	NAG	C1-C2-N2-C7
6	C	1208	NAG	C1-C2-N2-C7
6	C	1212	NAG	C1-C2-N2-C7
6	A	1204	NAG	C3-C2-N2-C7
6	B	1209	NAG	C3-C2-N2-C7
6	C	1209	NAG	C3-C2-N2-C7
6	C	1210	NAG	C3-C2-N2-C7
6	C	1211	NAG	O5-C5-C6-O6

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1210	NAG	1	0
6	C	1201	NAG	1	0
6	C	1214	NAG	1	0
6	C	1204	NAG	1	0
6	C	1208	NAG	1	0
6	B	1206	NAG	4	0

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.

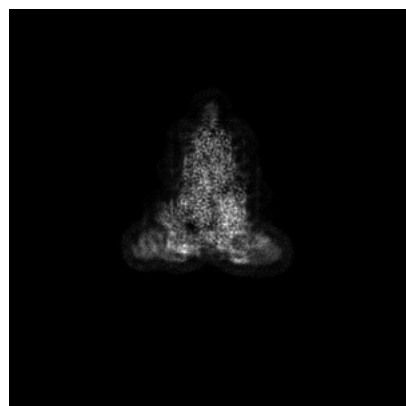
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-70441. These allow visual inspection of the internal detail of the map and identification of artifacts.

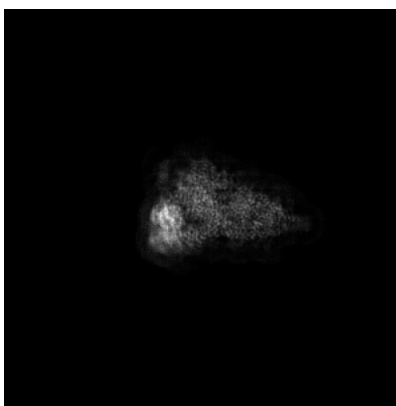
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

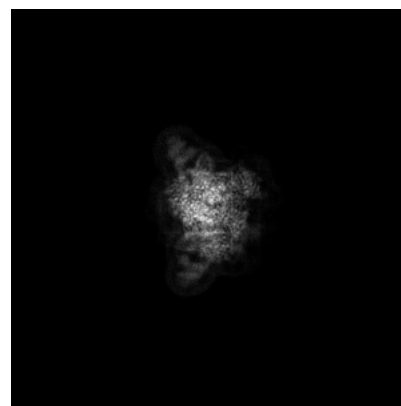
6.1.1 Primary map



X

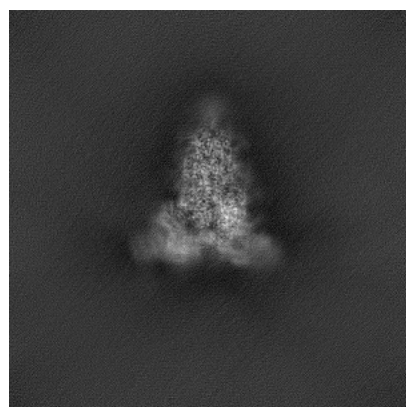


Y

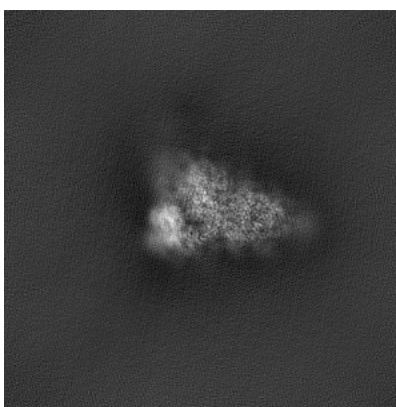


Z

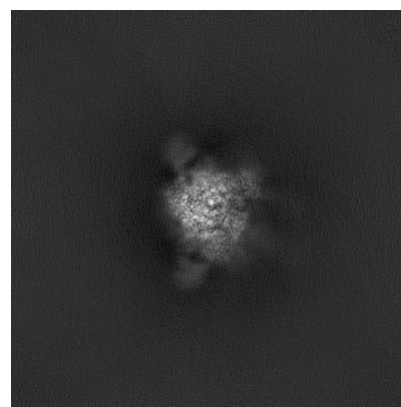
6.1.2 Raw map



X



Y

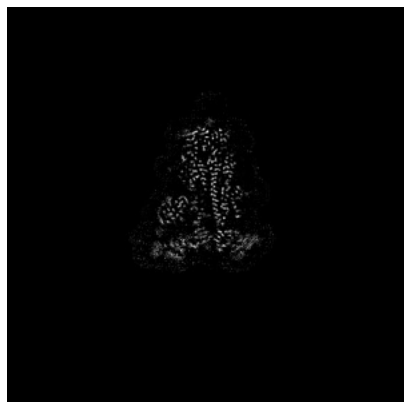


Z

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

6.2 Central slices [i](#)

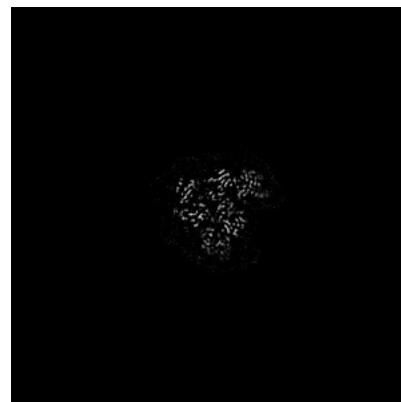
6.2.1 Primary map



X Index: 216

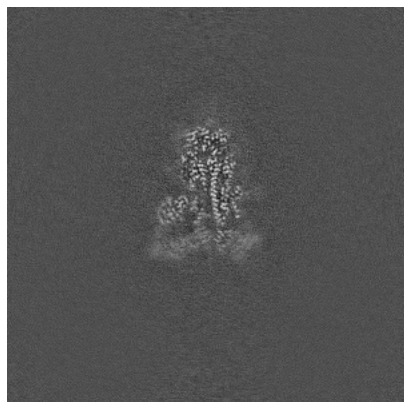


Y Index: 216

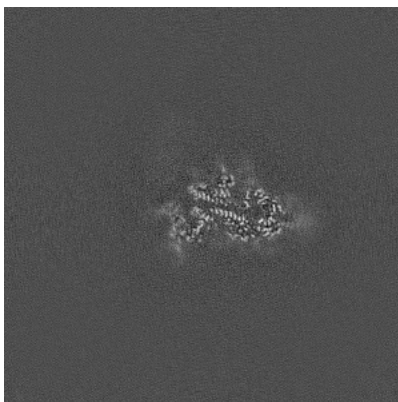


Z Index: 216

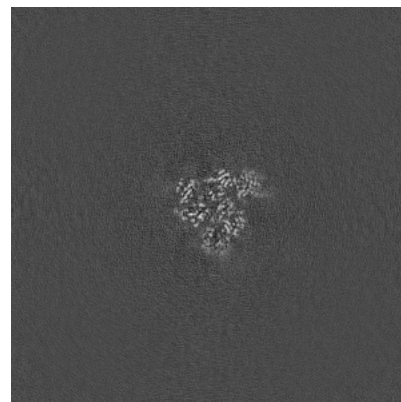
6.2.2 Raw map



X Index: 216



Y Index: 216

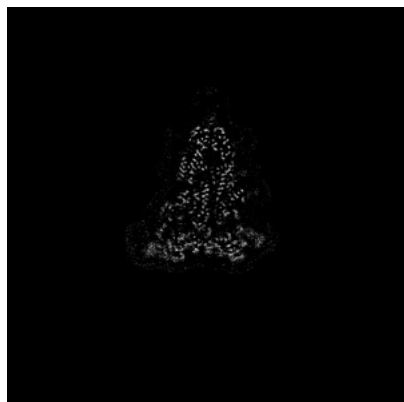


Z Index: 216

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

6.3 Largest variance slices [i](#)

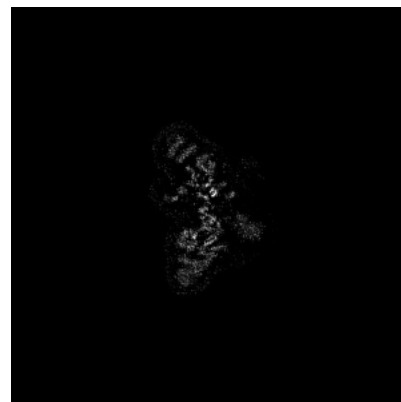
6.3.1 Primary map



X Index: 211

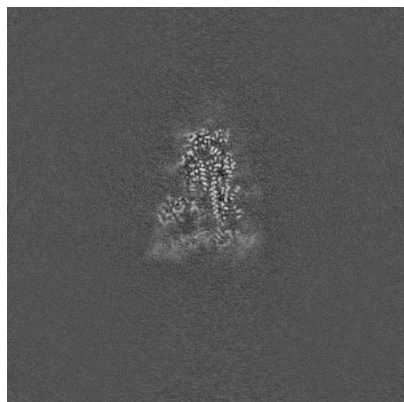


Y Index: 230

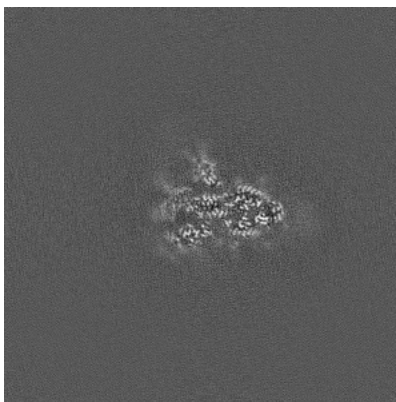


Z Index: 174

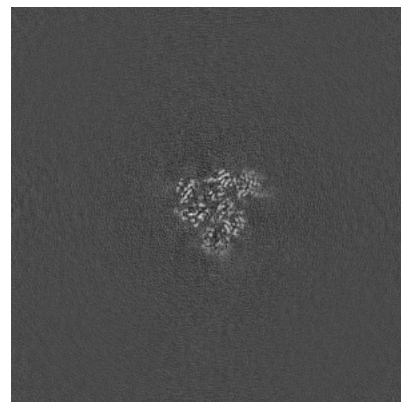
6.3.2 Raw map



X Index: 217



Y Index: 229

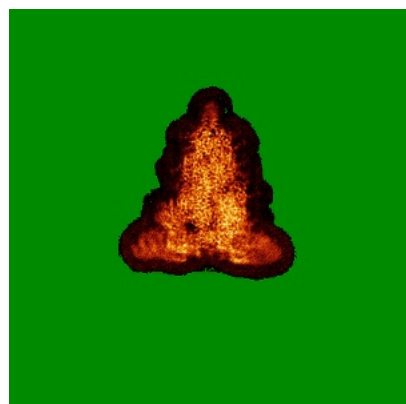


Z Index: 216

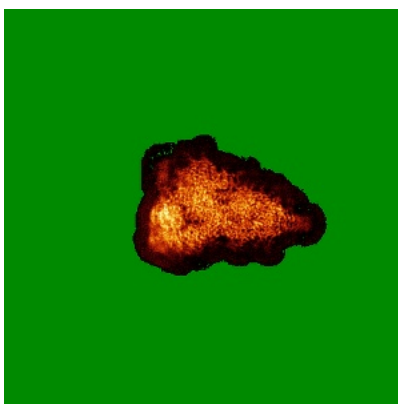
The images above show the largest variance slices of the map in three orthogonal directions.

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

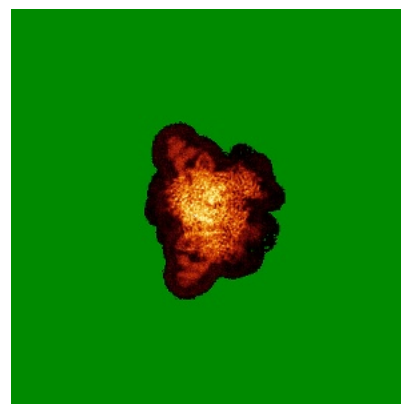
6.4.1 Primary map



X

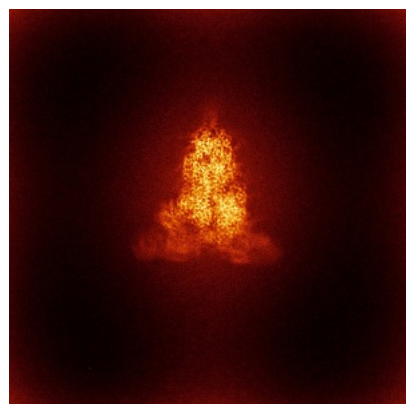


Y

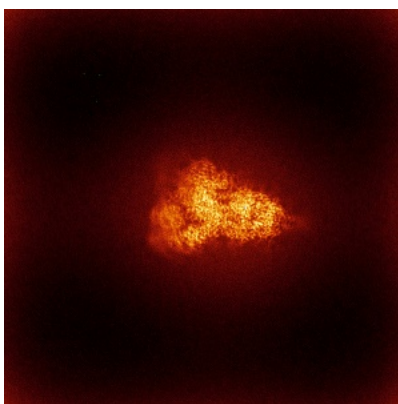


Z

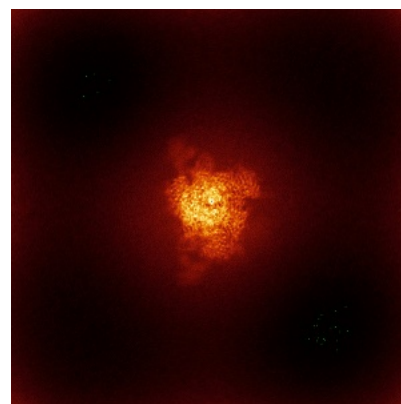
6.4.2 Raw map



X



Y

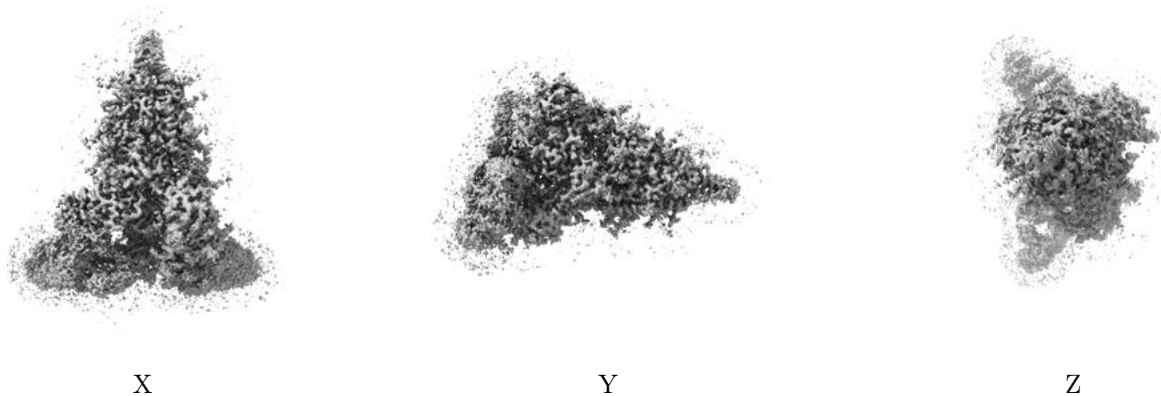


Z

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.

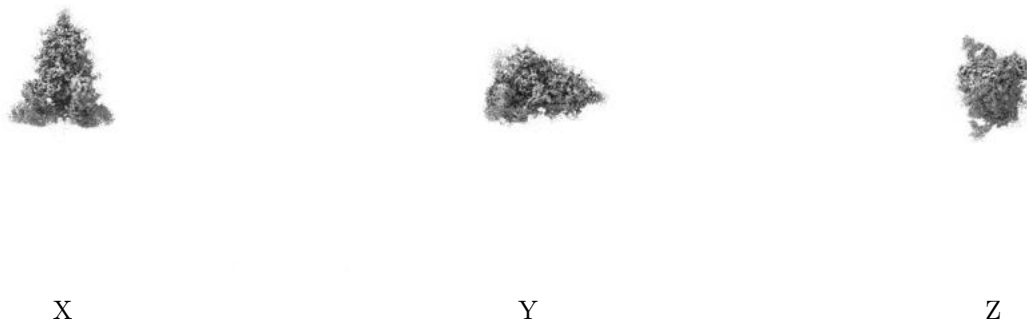
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.34. 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.

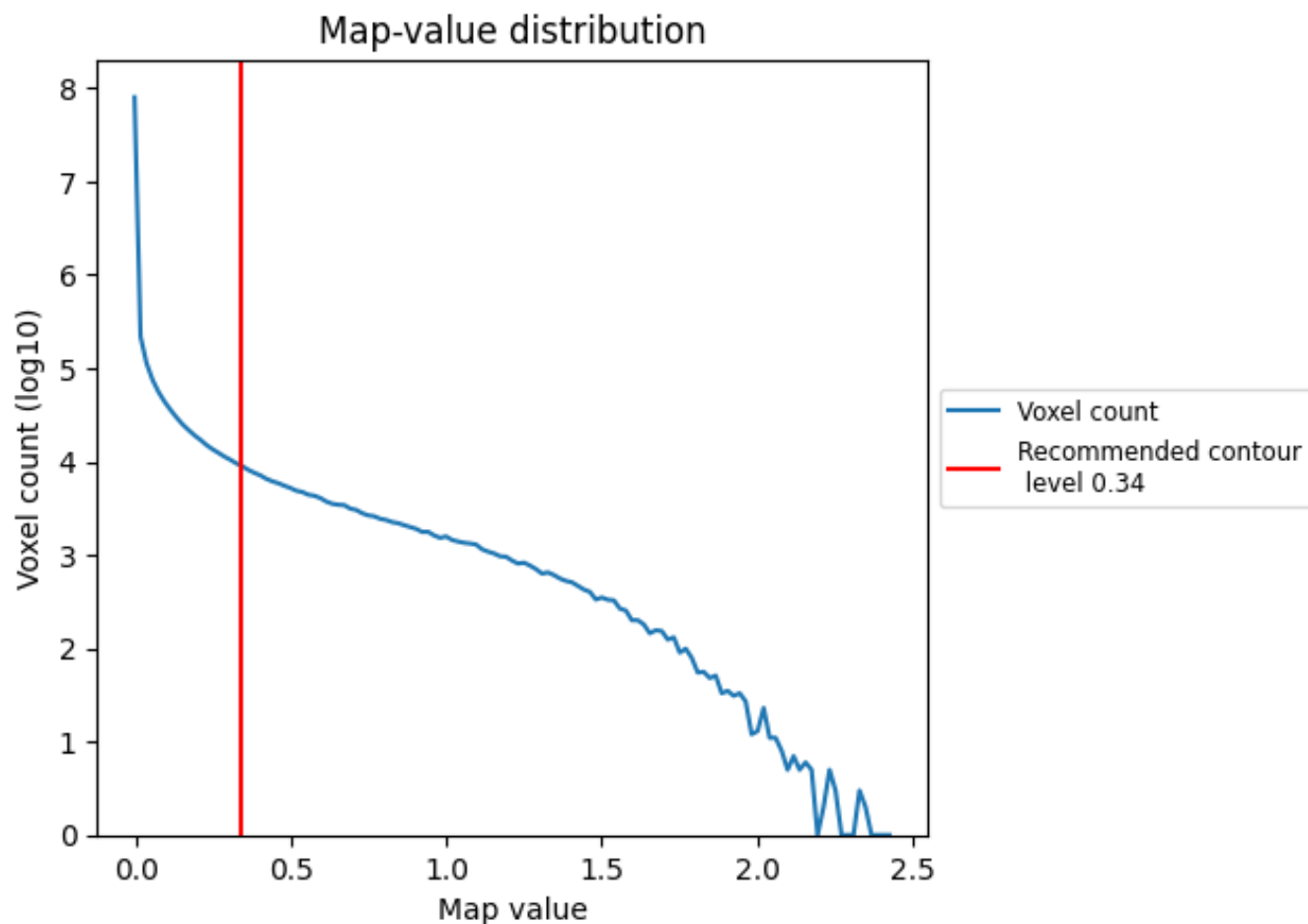
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

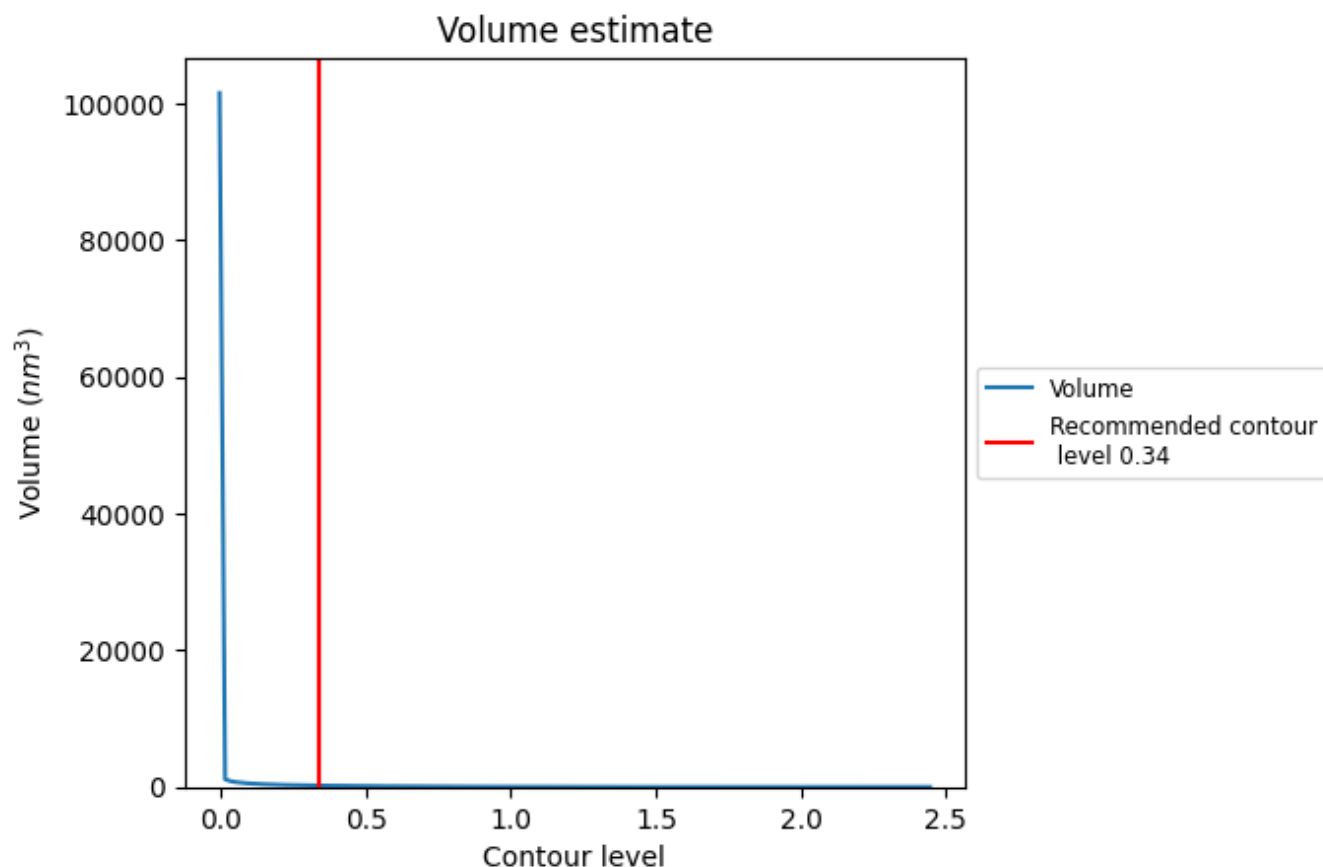
This section contains the results of statistical analysis of the map.

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.

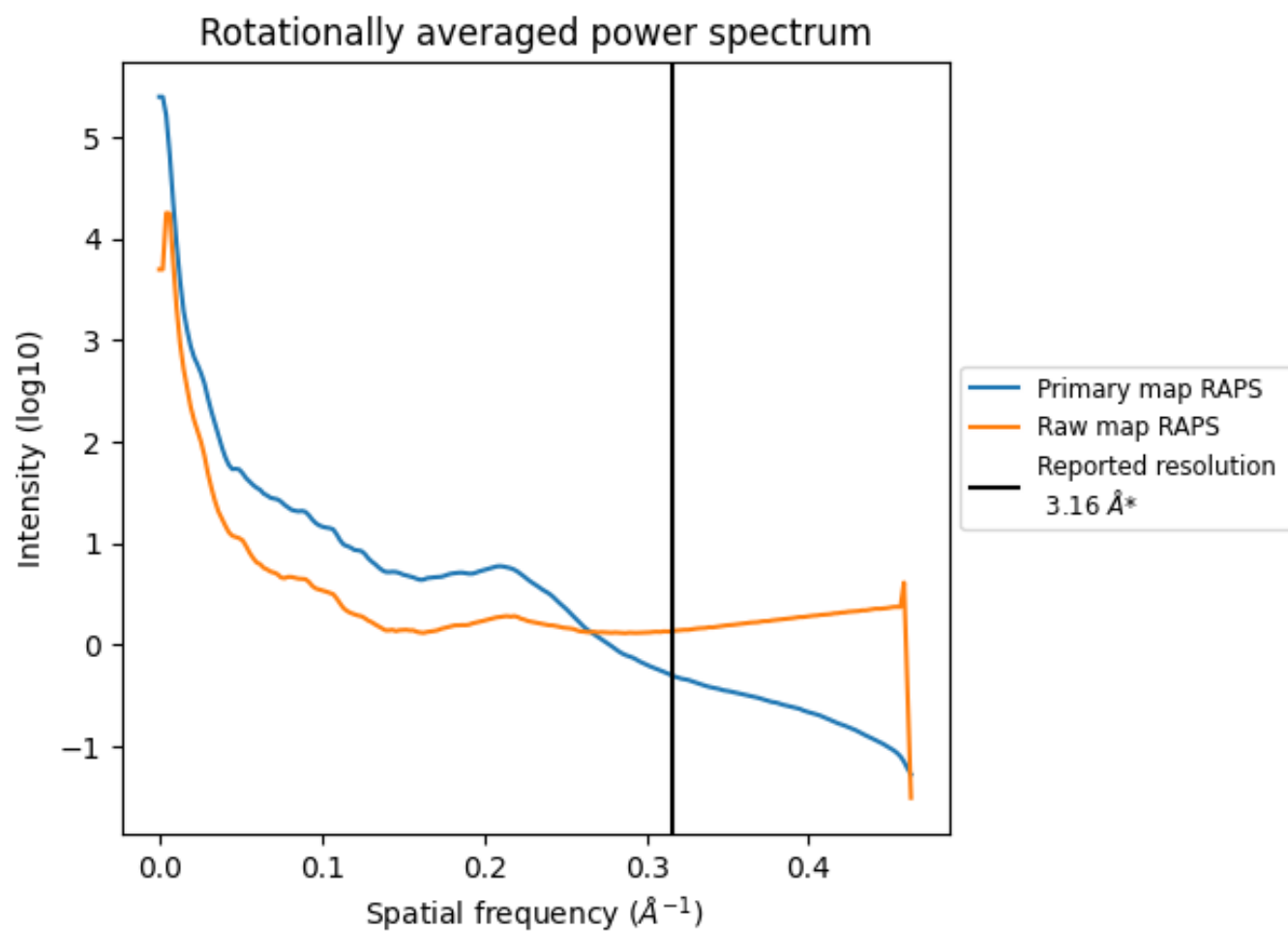
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 203 nm^3 ; this corresponds to an approximate mass of 183 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 ⓘ

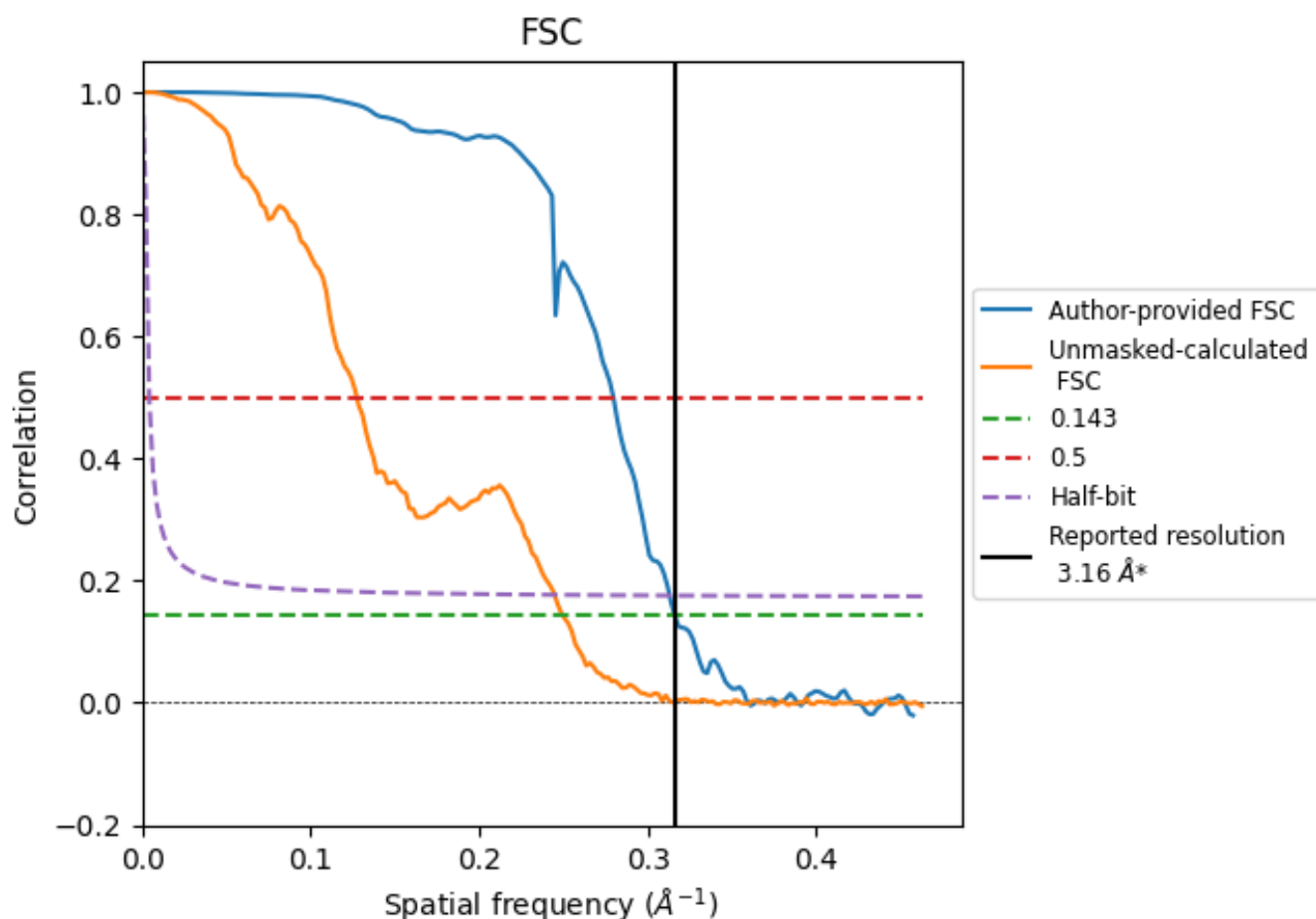


*Reported resolution corresponds to spatial frequency of 0.316 \AA^{-1}

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.316 \AA^{-1}

8.2 Resolution estimates [i](#)

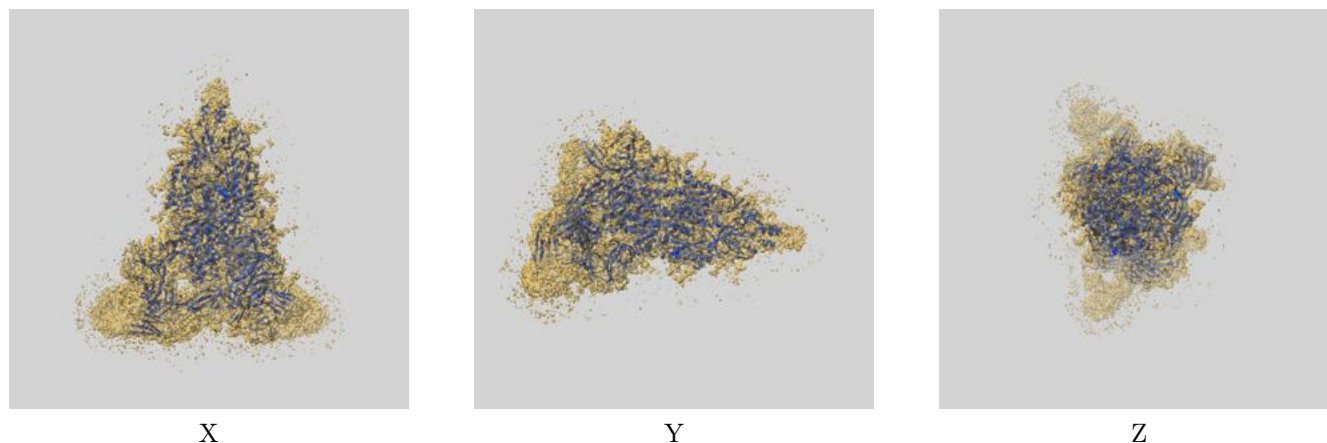
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.16	-	-
Author-provided FSC curve	3.16	3.58	3.19
Unmasked-calculated*	4.00	7.85	4.08

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.00 differs from the reported value 3.16 by more than 10 %

9 Map-model fit [i](#)

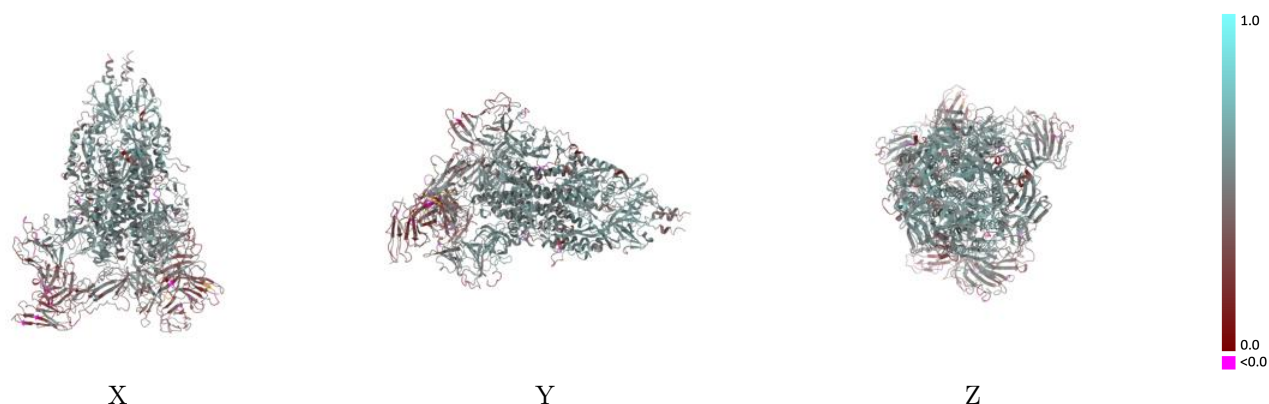
This section contains information regarding the fit between EMDB map EMD-70441 and PDB model 9OFP. Per-residue inclusion information can be found in section 3 on page 16.

9.1 Map-model overlay [i](#)



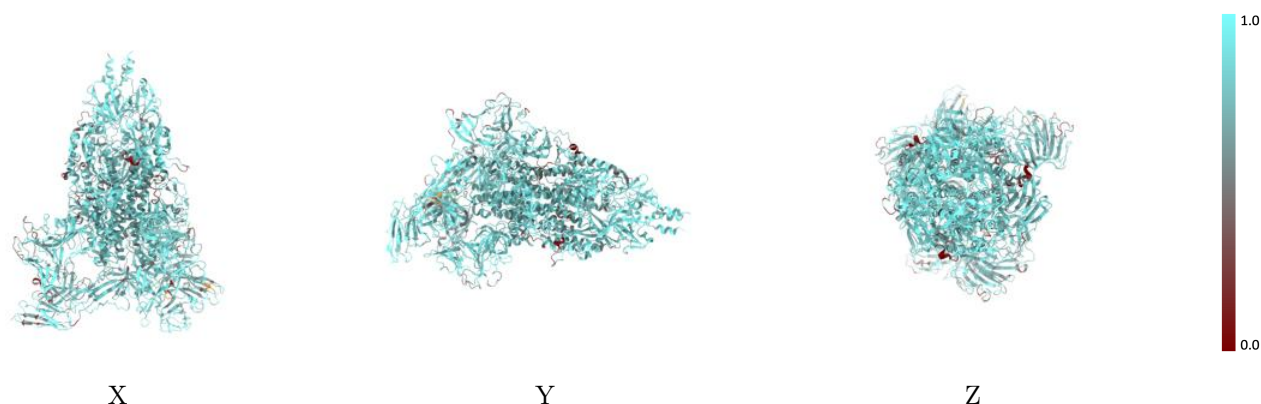
The images above show the 3D surface view of the map at the recommended contour level 0.34 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



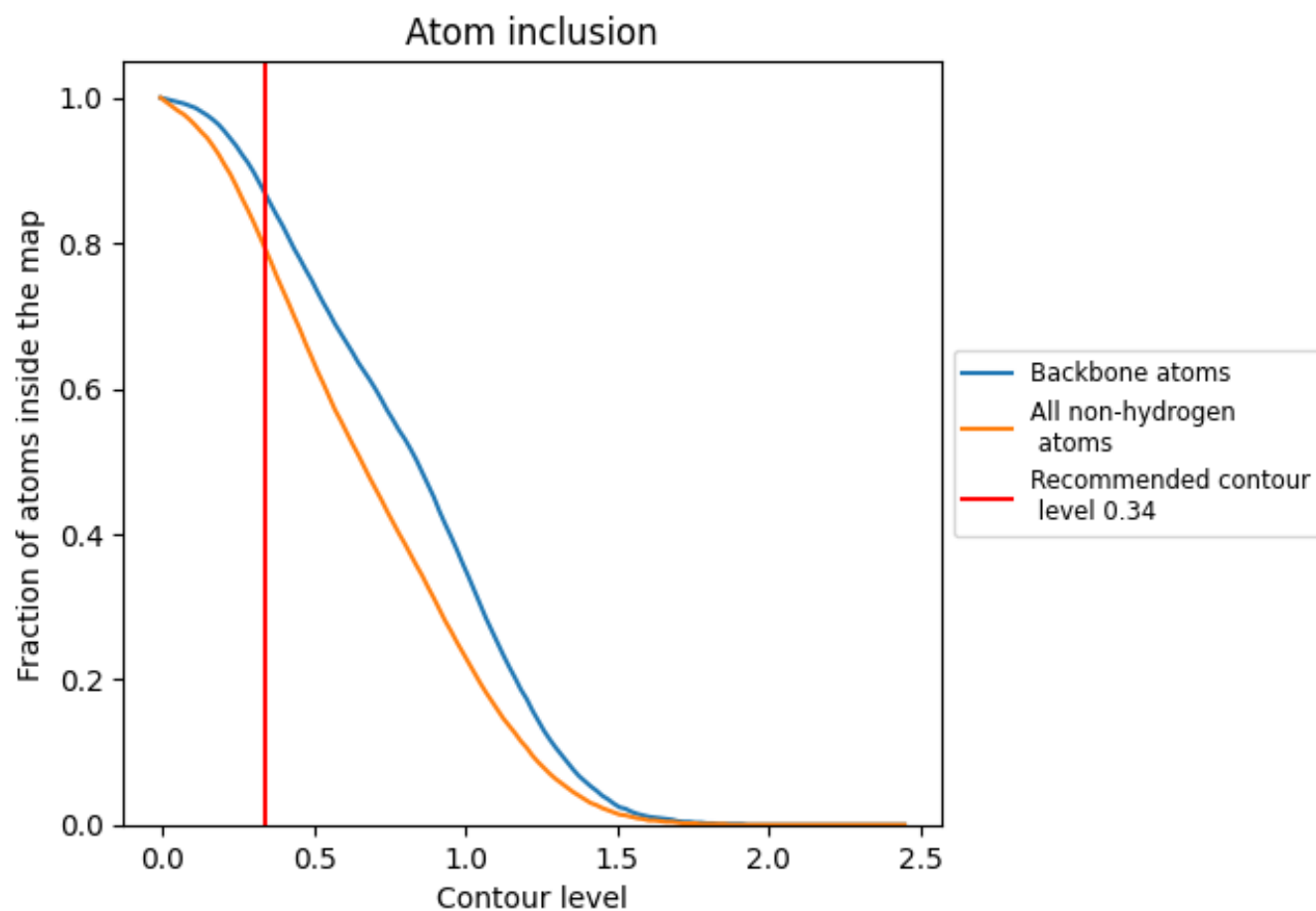
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.34).

































































9.4 Atom inclusion [i](#)



At the recommended contour level, 87% of all backbone atoms, 79% 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 (0.34) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7920	 0.4740
A	 0.7890	 0.4870
B	 0.8070	 0.5090
C	 0.7830	 0.5030
D	 0.7440	 0.4670
E	 0.7950	 0.3880
F	 0.6790	 0.4320
G	 0.9740	 0.4550
H	 0.7630	 0.3170
I	 0.7860	 0.4840
J	 0.6790	 0.4690
K	 0.5000	 0.3400
L	 0.7700	 0.2850
M	 0.7750	 0.3410
N	 0.8380	 0.3020
O	 0.7140	 0.5110
P	 0.7440	 0.4900
Q	 0.7690	 0.4080
R	 0.8970	 0.4400
S	 0.8570	 0.5140
T	 0.8970	 0.3880
U	 0.6920	 0.4330
V	 0.7950	 0.5000
W	 0.7140	 0.4710
X	 0.7860	 0.4920
Y	 0.7860	 0.4680
Z	 0.6070	 0.3930
a	 0.7690	 0.4760
b	 0.7860	 0.5090
c	 0.8460	 0.5240
d	 0.7140	 0.5020
e	 0.8570	 0.5530

