



Full wwPDB EM Validation Report ⓘ

Apr 12, 2026 – 11:32 AM UTC

PDB ID : 9RPB / pdb_00009rpb
EMDB ID : EMD-54151
Title : a5b3 GABAAR bound to GABA, and Mb25 in a desensitized state in saposin nanodiscs, topiramate-free
Authors : Cowgill, J.; Lindahl, E.; Howard, R.J.
Deposited on : 2025-06-24
Resolution : 3.10 Å(reported)
Based on initial model : 9HNS

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

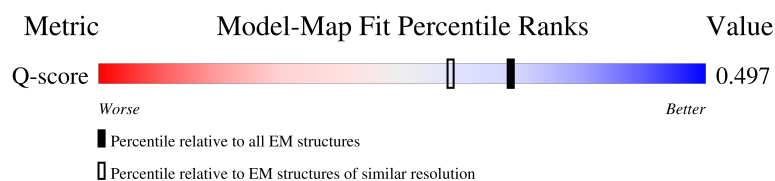
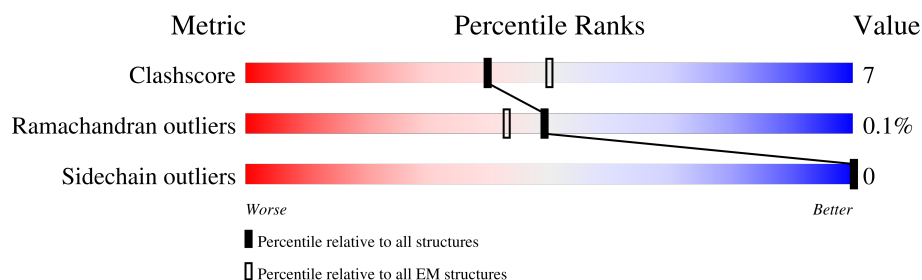
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

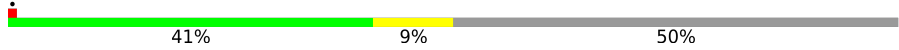
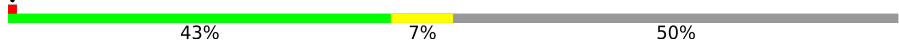
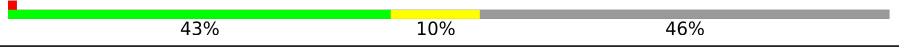
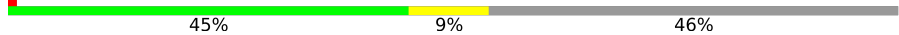
The reported resolution of this entry is 3.10 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14724 (2.60 - 3.60)

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	679	
1	D	679	
2	B	623	
2	C	623	

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Mol	Chain	Length	Quality of chain
2	E	623	
3	P	541	
4	F	2	
5	G	3	
5	I	3	
5	M	3	
6	H	6	
6	N	6	
7	J	3	
8	K	6	
9	L	2	

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 15127 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 Green fluorescent protein, Gamma-aminobutyric acid receptor subunit alpha-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	342	Total	C	N	O	S	0	0
			2745	1769	454	506	16		
1	D	339	Total	C	N	O	S	0	0
			2719	1754	451	498	16		

There are 194 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-315	MET	-	initiating methionine	UNP P42212
A	-314	ARG	-	expression tag	UNP P42212
A	-313	LYS	-	expression tag	UNP P42212
A	-312	SER	-	expression tag	UNP P42212
A	-311	PRO	-	expression tag	UNP P42212
A	-310	GLY	-	expression tag	UNP P42212
A	-309	LEU	-	expression tag	UNP P42212
A	-308	SER	-	expression tag	UNP P42212
A	-307	ASP	-	expression tag	UNP P42212
A	-306	CYS	-	expression tag	UNP P42212
A	-305	LEU	-	expression tag	UNP P42212
A	-304	TRP	-	expression tag	UNP P42212
A	-303	ALA	-	expression tag	UNP P42212
A	-302	TRP	-	expression tag	UNP P42212
A	-301	ILE	-	expression tag	UNP P42212
A	-300	LEU	-	expression tag	UNP P42212
A	-299	LEU	-	expression tag	UNP P42212
A	-298	LEU	-	expression tag	UNP P42212
A	-297	SER	-	expression tag	UNP P42212
A	-296	THR	-	expression tag	UNP P42212
A	-295	LEU	-	expression tag	UNP P42212
A	-294	THR	-	expression tag	UNP P42212
A	-293	GLY	-	expression tag	UNP P42212
A	-292	ARG	-	expression tag	UNP P42212
A	-291	SER	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-290	TYR	-	expression tag	UNP P42212
A	-289	GLY	-	expression tag	UNP P42212
A	-288	GLN	-	expression tag	UNP P42212
A	-258	ARG	SER	conflict	UNP P42212
A	-249	ASN	TYR	conflict	UNP P42212
A	-224	LEU	PHE	conflict	UNP P42212
A	-223	THR	SER	conflict	UNP P42212
A	-208	ARG	GLN	conflict	UNP P42212
A	-189	SER	PHE	conflict	UNP P42212
A	-183	THR	ASN	conflict	UNP P42212
A	-143	PHE	TYR	conflict	UNP P42212
A	-135	THR	MET	conflict	UNP P42212
A	-125	ALA	VAL	conflict	UNP P42212
A	-117	VAL	ILE	conflict	UNP P42212
A	-82	VAL	ALA	conflict	UNP P42212
A	-49	ALA	-	linker	UNP P42212
A	-48	ALA	-	linker	UNP P42212
A	-47	ASN	-	linker	UNP P42212
A	-46	ALA	-	linker	UNP P42212
A	-45	LEU	-	linker	UNP P42212
A	-44	ALA	-	linker	UNP P42212
A	-43	ALA	-	linker	UNP P42212
A	-42	TRP	-	linker	UNP P42212
A	-41	SER	-	linker	UNP P42212
A	-40	HIS	-	linker	UNP P42212
A	-39	PRO	-	linker	UNP P42212
A	-38	GLN	-	linker	UNP P42212
A	-37	PHE	-	linker	UNP P42212
A	-36	GLU	-	linker	UNP P42212
A	-35	LYS	-	linker	UNP P42212
A	-34	GLY	-	linker	UNP P42212
A	-33	GLY	-	linker	UNP P42212
A	-32	GLY	-	linker	UNP P42212
A	-31	SER	-	linker	UNP P42212
A	-30	GLY	-	linker	UNP P42212
A	-29	GLY	-	linker	UNP P42212
A	-28	GLY	-	linker	UNP P42212
A	-27	SER	-	linker	UNP P42212
A	-26	GLY	-	linker	UNP P42212
A	-25	GLY	-	linker	UNP P42212
A	-24	GLY	-	linker	UNP P42212
A	-23	SER	-	linker	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	TRP	-	linker	UNP P42212
A	-21	SER	-	linker	UNP P42212
A	-20	HIS	-	linker	UNP P42212
A	-19	PRO	-	linker	UNP P42212
A	-18	GLN	-	linker	UNP P42212
A	-17	PHE	-	linker	UNP P42212
A	-16	GLU	-	linker	UNP P42212
A	-15	LYS	-	linker	UNP P42212
A	-14	SER	-	linker	UNP P42212
A	-13	SER	-	linker	UNP P42212
A	-12	SER	-	linker	UNP P42212
A	-11	ASN	-	linker	UNP P42212
A	-10	ASN	-	linker	UNP P42212
A	-9	ASN	-	linker	UNP P42212
A	-8	ASN	-	linker	UNP P42212
A	-7	ASN	-	linker	UNP P42212
A	-6	GLU	-	linker	UNP P42212
A	-5	ASN	-	linker	UNP P42212
A	-4	LEU	-	linker	UNP P42212
A	-3	TYR	-	linker	UNP P42212
A	-2	PHE	-	linker	UNP P42212
A	-1	GLN	-	linker	UNP P42212
A	0	ALA	-	linker	UNP P42212
A	315	SER	-	linker	UNP P31644
A	316	GLN	-	linker	UNP P31644
A	317	PRO	-	linker	UNP P31644
A	318	ALA	-	linker	UNP P31644
A	319	ARG	-	linker	UNP P31644
A	320	ALA	-	linker	UNP P31644
A	321	ALA	-	linker	UNP P31644
D	-315	MET	-	initiating methionine	UNP P42212
D	-314	ARG	-	expression tag	UNP P42212
D	-313	LYS	-	expression tag	UNP P42212
D	-312	SER	-	expression tag	UNP P42212
D	-311	PRO	-	expression tag	UNP P42212
D	-310	GLY	-	expression tag	UNP P42212
D	-309	LEU	-	expression tag	UNP P42212
D	-308	SER	-	expression tag	UNP P42212
D	-307	ASP	-	expression tag	UNP P42212
D	-306	CYS	-	expression tag	UNP P42212
D	-305	LEU	-	expression tag	UNP P42212
D	-304	TRP	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-303	ALA	-	expression tag	UNP P42212
D	-302	TRP	-	expression tag	UNP P42212
D	-301	ILE	-	expression tag	UNP P42212
D	-300	LEU	-	expression tag	UNP P42212
D	-299	LEU	-	expression tag	UNP P42212
D	-298	LEU	-	expression tag	UNP P42212
D	-297	SER	-	expression tag	UNP P42212
D	-296	THR	-	expression tag	UNP P42212
D	-295	LEU	-	expression tag	UNP P42212
D	-294	THR	-	expression tag	UNP P42212
D	-293	GLY	-	expression tag	UNP P42212
D	-292	ARG	-	expression tag	UNP P42212
D	-291	SER	-	expression tag	UNP P42212
D	-290	TYR	-	expression tag	UNP P42212
D	-289	GLY	-	expression tag	UNP P42212
D	-288	GLN	-	expression tag	UNP P42212
D	-258	ARG	SER	conflict	UNP P42212
D	-249	ASN	TYR	conflict	UNP P42212
D	-224	LEU	PHE	conflict	UNP P42212
D	-223	THR	SER	conflict	UNP P42212
D	-208	ARG	GLN	conflict	UNP P42212
D	-189	SER	PHE	conflict	UNP P42212
D	-183	THR	ASN	conflict	UNP P42212
D	-143	PHE	TYR	conflict	UNP P42212
D	-135	THR	MET	conflict	UNP P42212
D	-125	ALA	VAL	conflict	UNP P42212
D	-117	VAL	ILE	conflict	UNP P42212
D	-82	VAL	ALA	conflict	UNP P42212
D	-49	ALA	-	linker	UNP P42212
D	-48	ALA	-	linker	UNP P42212
D	-47	ASN	-	linker	UNP P42212
D	-46	ALA	-	linker	UNP P42212
D	-45	LEU	-	linker	UNP P42212
D	-44	ALA	-	linker	UNP P42212
D	-43	ALA	-	linker	UNP P42212
D	-42	TRP	-	linker	UNP P42212
D	-41	SER	-	linker	UNP P42212
D	-40	HIS	-	linker	UNP P42212
D	-39	PRO	-	linker	UNP P42212
D	-38	GLN	-	linker	UNP P42212
D	-37	PHE	-	linker	UNP P42212
D	-36	GLU	-	linker	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-35	LYS	-	linker	UNP P42212
D	-34	GLY	-	linker	UNP P42212
D	-33	GLY	-	linker	UNP P42212
D	-32	GLY	-	linker	UNP P42212
D	-31	SER	-	linker	UNP P42212
D	-30	GLY	-	linker	UNP P42212
D	-29	GLY	-	linker	UNP P42212
D	-28	GLY	-	linker	UNP P42212
D	-27	SER	-	linker	UNP P42212
D	-26	GLY	-	linker	UNP P42212
D	-25	GLY	-	linker	UNP P42212
D	-24	GLY	-	linker	UNP P42212
D	-23	SER	-	linker	UNP P42212
D	-22	TRP	-	linker	UNP P42212
D	-21	SER	-	linker	UNP P42212
D	-20	HIS	-	linker	UNP P42212
D	-19	PRO	-	linker	UNP P42212
D	-18	GLN	-	linker	UNP P42212
D	-17	PHE	-	linker	UNP P42212
D	-16	GLU	-	linker	UNP P42212
D	-15	LYS	-	linker	UNP P42212
D	-14	SER	-	linker	UNP P42212
D	-13	SER	-	linker	UNP P42212
D	-12	SER	-	linker	UNP P42212
D	-11	ASN	-	linker	UNP P42212
D	-10	ASN	-	linker	UNP P42212
D	-9	ASN	-	linker	UNP P42212
D	-8	ASN	-	linker	UNP P42212
D	-7	ASN	-	linker	UNP P42212
D	-6	GLU	-	linker	UNP P42212
D	-5	ASN	-	linker	UNP P42212
D	-4	LEU	-	linker	UNP P42212
D	-3	TYR	-	linker	UNP P42212
D	-2	PHE	-	linker	UNP P42212
D	-1	GLN	-	linker	UNP P42212
D	0	ALA	-	linker	UNP P42212
D	315	SER	-	linker	UNP P31644
D	316	GLN	-	linker	UNP P31644
D	317	PRO	-	linker	UNP P31644
D	318	ALA	-	linker	UNP P31644
D	319	ARG	-	linker	UNP P31644
D	320	ALA	-	linker	UNP P31644

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Chain	Residue	Modelled	Actual	Comment	Reference
D	321	ALA	-	linker	UNP P31644

- Molecule 2 is a protein called Gamma-aminobutyric acid receptor subunit beta-3, Green fluorescent protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	336	Total	C	N	O	S	0	0
			2747	1796	450	485	16		
2	C	338	Total	C	N	O	S	0	0
			2767	1812	452	487	16		
2	E	336	Total	C	N	O	S	0	0
			2747	1797	450	484	16		

There are 153 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	308	SER	-	linker	UNP P28472
B	309	GLN	-	linker	UNP P28472
B	310	PRO	-	linker	UNP P28472
B	311	GLY	-	linker	UNP P28472
B	312	ARG	-	linker	UNP P28472
B	313	ALA	-	linker	UNP P28472
B	314	MET	-	linker	UNP P28472
B	315	VAL	-	linker	UNP P28472
B	332	MET	LEU	conflict	UNP P42212
B	339	ARG	HIS	conflict	UNP P42212
B	344	ARG	SER	conflict	UNP P42212
B	346	VAL	GLU	conflict	UNP P42212
B	353	HIS	TYR	conflict	UNP P42212
B	364	SER	THR	conflict	UNP P42212
B	378	LEU	PHE	conflict	UNP P42212
B	419	SER	ASN	conflict	UNP P42212
B	438	VAL	GLU	conflict	UNP P42212
B	442	THR	ILE	conflict	UNP P42212
B	459	MET	TYR	conflict	UNP P42212
B	461	VAL	SER	conflict	UNP P42212
B	462	GLY	HIS	conflict	UNP P42212
B	467	THR	MET	conflict	UNP P42212
B	477	ALA	VAL	conflict	UNP P42212
B	480	GLU	LYS	conflict	UNP P42212
B	485	VAL	ILE	conflict	UNP P42212
B	489	GLY	SER	conflict	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
B	506	SER	PRO	conflict	UNP P42212
B	517	VAL	THR	conflict	UNP P42212
B	519	VAL	SER	conflict	UNP P42212
B	520	LYS	ALA	conflict	UNP P42212
B	538	ARG	VAL	conflict	UNP P42212
B	545	PRO	HIS	conflict	UNP P42212
B	553	GLY	-	linker	UNP P42212
B	554	ARG	-	linker	UNP P42212
B	555	ALA	-	linker	UNP P42212
B	556	ALA	-	linker	UNP P42212
B	584	SER	-	expression tag	UNP P28472
B	585	ARG	-	expression tag	UNP P28472
B	586	GLU	-	expression tag	UNP P28472
B	587	ASN	-	expression tag	UNP P28472
B	588	LEU	-	expression tag	UNP P28472
B	589	TYR	-	expression tag	UNP P28472
B	590	PHE	-	expression tag	UNP P28472
B	591	GLN	-	expression tag	UNP P28472
B	592	ALA	-	expression tag	UNP P28472
B	593	HIS	-	expression tag	UNP P28472
B	594	HIS	-	expression tag	UNP P28472
B	595	HIS	-	expression tag	UNP P28472
B	596	HIS	-	expression tag	UNP P28472
B	597	HIS	-	expression tag	UNP P28472
B	598	HIS	-	expression tag	UNP P28472
C	308	SER	-	linker	UNP P28472
C	309	GLN	-	linker	UNP P28472
C	310	PRO	-	linker	UNP P28472
C	311	GLY	-	linker	UNP P28472
C	312	ARG	-	linker	UNP P28472
C	313	ALA	-	linker	UNP P28472
C	314	MET	-	linker	UNP P28472
C	315	VAL	-	linker	UNP P28472
C	332	MET	LEU	conflict	UNP P42212
C	339	ARG	HIS	conflict	UNP P42212
C	344	ARG	SER	conflict	UNP P42212
C	346	VAL	GLU	conflict	UNP P42212
C	353	HIS	TYR	conflict	UNP P42212
C	364	SER	THR	conflict	UNP P42212
C	378	LEU	PHE	conflict	UNP P42212
C	419	SER	ASN	conflict	UNP P42212
C	438	VAL	GLU	conflict	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
C	442	THR	ILE	conflict	UNP P42212
C	459	MET	TYR	conflict	UNP P42212
C	461	VAL	SER	conflict	UNP P42212
C	462	GLY	HIS	conflict	UNP P42212
C	467	THR	MET	conflict	UNP P42212
C	477	ALA	VAL	conflict	UNP P42212
C	480	GLU	LYS	conflict	UNP P42212
C	485	VAL	ILE	conflict	UNP P42212
C	489	GLY	SER	conflict	UNP P42212
C	506	SER	PRO	conflict	UNP P42212
C	517	VAL	THR	conflict	UNP P42212
C	519	VAL	SER	conflict	UNP P42212
C	520	LYS	ALA	conflict	UNP P42212
C	538	ARG	VAL	conflict	UNP P42212
C	545	PRO	HIS	conflict	UNP P42212
C	553	GLY	-	linker	UNP P42212
C	554	ARG	-	linker	UNP P42212
C	555	ALA	-	linker	UNP P42212
C	556	ALA	-	linker	UNP P42212
C	584	SER	-	expression tag	UNP P28472
C	585	ARG	-	expression tag	UNP P28472
C	586	GLU	-	expression tag	UNP P28472
C	587	ASN	-	expression tag	UNP P28472
C	588	LEU	-	expression tag	UNP P28472
C	589	TYR	-	expression tag	UNP P28472
C	590	PHE	-	expression tag	UNP P28472
C	591	GLN	-	expression tag	UNP P28472
C	592	ALA	-	expression tag	UNP P28472
C	593	HIS	-	expression tag	UNP P28472
C	594	HIS	-	expression tag	UNP P28472
C	595	HIS	-	expression tag	UNP P28472
C	596	HIS	-	expression tag	UNP P28472
C	597	HIS	-	expression tag	UNP P28472
C	598	HIS	-	expression tag	UNP P28472
E	308	SER	-	linker	UNP P28472
E	309	GLN	-	linker	UNP P28472
E	310	PRO	-	linker	UNP P28472
E	311	GLY	-	linker	UNP P28472
E	312	ARG	-	linker	UNP P28472
E	313	ALA	-	linker	UNP P28472
E	314	MET	-	linker	UNP P28472
E	315	VAL	-	linker	UNP P28472

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Chain	Residue	Modelled	Actual	Comment	Reference
E	332	MET	LEU	conflict	UNP P42212
E	339	ARG	HIS	conflict	UNP P42212
E	344	ARG	SER	conflict	UNP P42212
E	346	VAL	GLU	conflict	UNP P42212
E	353	HIS	TYR	conflict	UNP P42212
E	364	SER	THR	conflict	UNP P42212
E	378	LEU	PHE	conflict	UNP P42212
E	419	SER	ASN	conflict	UNP P42212
E	438	VAL	GLU	conflict	UNP P42212
E	442	THR	ILE	conflict	UNP P42212
E	459	MET	TYR	conflict	UNP P42212
E	461	VAL	SER	conflict	UNP P42212
E	462	GLY	HIS	conflict	UNP P42212
E	467	THR	MET	conflict	UNP P42212
E	477	ALA	VAL	conflict	UNP P42212
E	480	GLU	LYS	conflict	UNP P42212
E	485	VAL	ILE	conflict	UNP P42212
E	489	GLY	SER	conflict	UNP P42212
E	506	SER	PRO	conflict	UNP P42212
E	517	VAL	THR	conflict	UNP P42212
E	519	VAL	SER	conflict	UNP P42212
E	520	LYS	ALA	conflict	UNP P42212
E	538	ARG	VAL	conflict	UNP P42212
E	545	PRO	HIS	conflict	UNP P42212
E	553	GLY	-	linker	UNP P42212
E	554	ARG	-	linker	UNP P42212
E	555	ALA	-	linker	UNP P42212
E	556	ALA	-	linker	UNP P42212
E	584	SER	-	expression tag	UNP P28472
E	585	ARG	-	expression tag	UNP P28472
E	586	GLU	-	expression tag	UNP P28472
E	587	ASN	-	expression tag	UNP P28472
E	588	LEU	-	expression tag	UNP P28472
E	589	TYR	-	expression tag	UNP P28472
E	590	PHE	-	expression tag	UNP P28472
E	591	GLN	-	expression tag	UNP P28472
E	592	ALA	-	expression tag	UNP P28472
E	593	HIS	-	expression tag	UNP P28472
E	594	HIS	-	expression tag	UNP P28472
E	595	HIS	-	expression tag	UNP P28472
E	596	HIS	-	expression tag	UNP P28472
E	597	HIS	-	expression tag	UNP P28472

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Chain	Residue	Modelled	Actual	Comment	Reference
E	598	HIS	-	expression tag	UNP P28472

- Molecule 3 is a protein called Outer membrane protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	P	120	Total	C	N	O	S	0	0
			931	588	159	180	4		

There are 320 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	-18	MET	-	initiating methionine	UNP B5Z8H1
P	-17	LYS	-	expression tag	UNP B5Z8H1
P	-16	LYS	-	expression tag	UNP B5Z8H1
P	-15	ILE	-	expression tag	UNP B5Z8H1
P	-14	TRP	-	expression tag	UNP B5Z8H1
P	-13	LEU	-	expression tag	UNP B5Z8H1
P	-12	ALA	-	expression tag	UNP B5Z8H1
P	-11	LEU	-	expression tag	UNP B5Z8H1
P	-10	ALA	-	expression tag	UNP B5Z8H1
P	-9	GLY	-	expression tag	UNP B5Z8H1
P	-8	LEU	-	expression tag	UNP B5Z8H1
P	-7	VAL	-	expression tag	UNP B5Z8H1
P	-6	LEU	-	expression tag	UNP B5Z8H1
P	-5	ALA	-	expression tag	UNP B5Z8H1
P	-4	PHE	-	expression tag	UNP B5Z8H1
P	-3	SER	-	expression tag	UNP B5Z8H1
P	-2	ALA	-	expression tag	UNP B5Z8H1
P	-1	SER	-	expression tag	UNP B5Z8H1
P	0	ALA	-	expression tag	UNP B5Z8H1
P	1	GLN	-	expression tag	UNP B5Z8H1
P	2	VAL	-	expression tag	UNP B5Z8H1
P	3	GLN	-	expression tag	UNP B5Z8H1
P	4	LEU	-	expression tag	UNP B5Z8H1
P	5	VAL	-	expression tag	UNP B5Z8H1
P	6	GLU	-	expression tag	UNP B5Z8H1
P	7	SER	-	expression tag	UNP B5Z8H1
P	8	GLY	-	expression tag	UNP B5Z8H1
P	9	GLY	-	expression tag	UNP B5Z8H1
P	10	GLY	-	expression tag	UNP B5Z8H1
P	11	LEU	-	expression tag	UNP B5Z8H1
P	12	VAL	-	expression tag	UNP B5Z8H1

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Chain	Residue	Modelled	Actual	Comment	Reference
P	234	THR	PRO	conflict	UNP B5Z8H1
P	235	TYR	PHE	conflict	UNP B5Z8H1
P	236	GLU	ARG	conflict	UNP B5Z8H1
P	237	GLN	ASN	conflict	UNP B5Z8H1
P	238	LEU	MET	conflict	UNP B5Z8H1
P	239	SER	GLY	conflict	UNP B5Z8H1
P	240	ARG	MET	conflict	UNP B5Z8H1
P	241	LEU	ILE	conflict	UNP B5Z8H1
P	242	LEU	-	expression tag	UNP B5Z8H1
P	243	THR	-	expression tag	UNP B5Z8H1
P	244	ASN	-	expression tag	UNP B5Z8H1
P	245	ASP	-	expression tag	UNP B5Z8H1
P	246	ASN	-	expression tag	UNP B5Z8H1
P	247	GLY	-	expression tag	UNP B5Z8H1
P	248	THR	-	expression tag	UNP B5Z8H1
P	249	ASN	-	expression tag	UNP B5Z8H1
P	250	SER	-	expression tag	UNP B5Z8H1
P	251	LYS	-	expression tag	UNP B5Z8H1
P	252	THR	-	expression tag	UNP B5Z8H1
P	253	SER	-	expression tag	UNP B5Z8H1
P	254	ALA	-	expression tag	UNP B5Z8H1
P	255	GLN	-	expression tag	UNP B5Z8H1
P	256	ALA	-	expression tag	UNP B5Z8H1
P	257	ILE	-	expression tag	UNP B5Z8H1
P	258	ASN	-	expression tag	UNP B5Z8H1
P	259	GLN	-	expression tag	UNP B5Z8H1
P	260	ALA	-	expression tag	UNP B5Z8H1
P	261	VAL	-	expression tag	UNP B5Z8H1
P	262	ASN	-	expression tag	UNP B5Z8H1
P	263	ASN	-	expression tag	UNP B5Z8H1
P	264	LEU	-	expression tag	UNP B5Z8H1
P	265	ASN	-	expression tag	UNP B5Z8H1
P	266	GLU	-	expression tag	UNP B5Z8H1
P	267	ARG	-	expression tag	UNP B5Z8H1
P	268	ALA	-	expression tag	UNP B5Z8H1
P	269	LYS	-	expression tag	UNP B5Z8H1
P	270	THR	-	expression tag	UNP B5Z8H1
P	271	LEU	-	expression tag	UNP B5Z8H1
P	272	ALA	-	expression tag	UNP B5Z8H1
P	273	GLY	-	expression tag	UNP B5Z8H1
P	274	GLY	-	expression tag	UNP B5Z8H1
P	275	THR	-	expression tag	UNP B5Z8H1

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Chain	Residue	Modelled	Actual	Comment	Reference
P	276	THR	-	expression tag	UNP B5Z8H1
P	277	ASN	-	expression tag	UNP B5Z8H1
P	278	SER	-	expression tag	UNP B5Z8H1
P	279	PRO	-	expression tag	UNP B5Z8H1
P	280	ALA	-	expression tag	UNP B5Z8H1
P	281	TYR	-	expression tag	UNP B5Z8H1
P	282	GLN	-	expression tag	UNP B5Z8H1
P	283	ALA	-	expression tag	UNP B5Z8H1
P	284	THR	-	expression tag	UNP B5Z8H1
P	285	LEU	-	expression tag	UNP B5Z8H1
P	286	LEU	-	expression tag	UNP B5Z8H1
P	287	ALA	-	expression tag	UNP B5Z8H1
P	288	LEU	-	expression tag	UNP B5Z8H1
P	289	ARG	-	expression tag	UNP B5Z8H1
P	290	SER	-	expression tag	UNP B5Z8H1
P	291	VAL	-	expression tag	UNP B5Z8H1
P	292	LEU	-	expression tag	UNP B5Z8H1
P	293	GLY	-	expression tag	UNP B5Z8H1
P	294	LEU	-	expression tag	UNP B5Z8H1
P	295	TRP	-	expression tag	UNP B5Z8H1
P	296	ASN	-	expression tag	UNP B5Z8H1
P	297	SER	-	expression tag	UNP B5Z8H1
P	298	MET	-	expression tag	UNP B5Z8H1
P	299	GLY	-	expression tag	UNP B5Z8H1
P	300	TYR	-	expression tag	UNP B5Z8H1
P	301	ALA	-	expression tag	UNP B5Z8H1
P	302	VAL	-	expression tag	UNP B5Z8H1
P	303	ILE	-	expression tag	UNP B5Z8H1
P	304	CYS	-	expression tag	UNP B5Z8H1
P	305	GLY	-	expression tag	UNP B5Z8H1
P	306	GLY	-	expression tag	UNP B5Z8H1
P	307	TYR	-	expression tag	UNP B5Z8H1
P	308	THR	-	expression tag	UNP B5Z8H1
P	309	LYS	-	expression tag	UNP B5Z8H1
P	310	SER	-	expression tag	UNP B5Z8H1
P	311	PRO	-	expression tag	UNP B5Z8H1
P	312	GLY	-	expression tag	UNP B5Z8H1
P	313	GLU	-	expression tag	UNP B5Z8H1
P	314	ASN	-	expression tag	UNP B5Z8H1
P	315	ASN	-	expression tag	UNP B5Z8H1
P	316	GLN	-	expression tag	UNP B5Z8H1
P	317	LYS	-	expression tag	UNP B5Z8H1

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Chain	Residue	Modelled	Actual	Comment	Reference
P	318	ASP	-	expression tag	UNP B5Z8H1
P	319	PHE	-	expression tag	UNP B5Z8H1
P	320	HIS	-	expression tag	UNP B5Z8H1
P	321	TYR	-	expression tag	UNP B5Z8H1
P	322	THR	-	expression tag	UNP B5Z8H1
P	323	ASP	-	expression tag	UNP B5Z8H1
P	324	GLU	-	expression tag	UNP B5Z8H1
P	325	ASN	-	expression tag	UNP B5Z8H1
P	326	GLY	-	expression tag	UNP B5Z8H1
P	327	ASN	-	expression tag	UNP B5Z8H1
P	328	GLY	-	expression tag	UNP B5Z8H1
P	329	THR	-	expression tag	UNP B5Z8H1
P	330	THR	-	expression tag	UNP B5Z8H1
P	331	ILE	-	expression tag	UNP B5Z8H1
P	332	ASN	-	expression tag	UNP B5Z8H1
P	333	CYS	-	expression tag	UNP B5Z8H1
P	334	GLY	-	expression tag	UNP B5Z8H1
P	335	GLY	-	expression tag	UNP B5Z8H1
P	336	SER	-	expression tag	UNP B5Z8H1
P	337	THR	-	expression tag	UNP B5Z8H1
P	338	ASN	-	expression tag	UNP B5Z8H1
P	339	SER	-	expression tag	UNP B5Z8H1
P	340	ASN	-	expression tag	UNP B5Z8H1
P	341	GLY	-	expression tag	UNP B5Z8H1
P	342	THR	-	expression tag	UNP B5Z8H1
P	343	HIS	-	expression tag	UNP B5Z8H1
P	344	SER	-	expression tag	UNP B5Z8H1
P	345	TYR	-	expression tag	UNP B5Z8H1
P	346	ASN	-	expression tag	UNP B5Z8H1
P	347	GLY	-	expression tag	UNP B5Z8H1
P	348	THR	-	expression tag	UNP B5Z8H1
P	349	ASN	-	expression tag	UNP B5Z8H1
P	350	THR	-	expression tag	UNP B5Z8H1
P	351	LEU	-	expression tag	UNP B5Z8H1
P	352	LYS	-	expression tag	UNP B5Z8H1
P	353	ALA	-	expression tag	UNP B5Z8H1
P	354	ASP	-	expression tag	UNP B5Z8H1
P	355	LYS	-	expression tag	UNP B5Z8H1
P	356	ASN	-	expression tag	UNP B5Z8H1
P	357	VAL	-	expression tag	UNP B5Z8H1
P	358	SER	-	expression tag	UNP B5Z8H1
P	359	LEU	-	expression tag	UNP B5Z8H1

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Chain	Residue	Modelled	Actual	Comment	Reference
P	360	SER	-	expression tag	UNP B5Z8H1
P	361	ILE	-	expression tag	UNP B5Z8H1
P	362	GLU	-	expression tag	UNP B5Z8H1
P	363	GLN	-	expression tag	UNP B5Z8H1
P	364	TYR	-	expression tag	UNP B5Z8H1
P	365	GLU	-	expression tag	UNP B5Z8H1
P	366	LYS	-	expression tag	UNP B5Z8H1
P	367	ILE	-	expression tag	UNP B5Z8H1
P	368	HIS	-	expression tag	UNP B5Z8H1
P	369	GLU	-	expression tag	UNP B5Z8H1
P	370	ALA	-	expression tag	UNP B5Z8H1
P	371	TYR	-	expression tag	UNP B5Z8H1
P	372	GLN	-	expression tag	UNP B5Z8H1
P	373	ILE	-	expression tag	UNP B5Z8H1
P	374	LEU	-	expression tag	UNP B5Z8H1
P	375	SER	-	expression tag	UNP B5Z8H1
P	376	LYS	-	expression tag	UNP B5Z8H1
P	377	ALA	-	expression tag	UNP B5Z8H1
P	378	LEU	-	expression tag	UNP B5Z8H1
P	379	LYS	-	expression tag	UNP B5Z8H1
P	380	GLN	-	expression tag	UNP B5Z8H1
P	381	ALA	-	expression tag	UNP B5Z8H1
P	382	GLY	-	expression tag	UNP B5Z8H1
P	383	LEU	-	expression tag	UNP B5Z8H1
P	384	ALA	-	expression tag	UNP B5Z8H1
P	385	PRO	-	expression tag	UNP B5Z8H1
P	386	LEU	-	expression tag	UNP B5Z8H1
P	387	ASN	-	expression tag	UNP B5Z8H1
P	388	SER	-	expression tag	UNP B5Z8H1
P	389	LYS	-	expression tag	UNP B5Z8H1
P	390	GLY	-	expression tag	UNP B5Z8H1
P	391	GLU	-	expression tag	UNP B5Z8H1
P	392	LYS	-	expression tag	UNP B5Z8H1
P	393	LEU	-	expression tag	UNP B5Z8H1
P	394	GLU	-	expression tag	UNP B5Z8H1
P	395	ALA	-	expression tag	UNP B5Z8H1
P	396	HIS	-	expression tag	UNP B5Z8H1
P	397	VAL	-	expression tag	UNP B5Z8H1
P	398	THR	-	expression tag	UNP B5Z8H1
P	399	THR	-	expression tag	UNP B5Z8H1
P	400	SER	-	expression tag	UNP B5Z8H1
P	401	LYS	-	expression tag	UNP B5Z8H1

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Chain	Residue	Modelled	Actual	Comment	Reference
P	402	TYR	-	expression tag	UNP B5Z8H1
P	403	GLY	-	expression tag	UNP B5Z8H1
P	404	SER	-	expression tag	UNP B5Z8H1
P	405	LEU	-	expression tag	UNP B5Z8H1
P	406	ARG	-	expression tag	UNP B5Z8H1
P	407	LEU	-	expression tag	UNP B5Z8H1
P	408	SER	-	expression tag	UNP B5Z8H1
P	409	CYS	-	expression tag	UNP B5Z8H1
P	410	ALA	-	expression tag	UNP B5Z8H1
P	411	ALA	-	expression tag	UNP B5Z8H1
P	412	SER	-	expression tag	UNP B5Z8H1
P	413	GLY	-	expression tag	UNP B5Z8H1
P	414	HIS	-	expression tag	UNP B5Z8H1
P	415	THR	-	expression tag	UNP B5Z8H1
P	416	PHE	-	expression tag	UNP B5Z8H1
P	417	ASN	-	expression tag	UNP B5Z8H1
P	418	TYR	-	expression tag	UNP B5Z8H1
P	419	PRO	-	expression tag	UNP B5Z8H1
P	420	ILE	-	expression tag	UNP B5Z8H1
P	421	MET	-	expression tag	UNP B5Z8H1
P	422	GLY	-	expression tag	UNP B5Z8H1
P	423	TRP	-	expression tag	UNP B5Z8H1
P	424	PHE	-	expression tag	UNP B5Z8H1
P	425	ARG	-	expression tag	UNP B5Z8H1
P	426	GLN	-	expression tag	UNP B5Z8H1
P	427	ALA	-	expression tag	UNP B5Z8H1
P	428	PRO	-	expression tag	UNP B5Z8H1
P	429	GLY	-	expression tag	UNP B5Z8H1
P	430	LYS	-	expression tag	UNP B5Z8H1
P	431	GLU	-	expression tag	UNP B5Z8H1
P	432	ARG	-	expression tag	UNP B5Z8H1
P	433	GLU	-	expression tag	UNP B5Z8H1
P	434	PHE	-	expression tag	UNP B5Z8H1
P	435	VAL	-	expression tag	UNP B5Z8H1
P	436	GLY	-	expression tag	UNP B5Z8H1
P	437	ALA	-	expression tag	UNP B5Z8H1
P	438	ILE	-	expression tag	UNP B5Z8H1
P	439	SER	-	expression tag	UNP B5Z8H1
P	440	TRP	-	expression tag	UNP B5Z8H1
P	441	SER	-	expression tag	UNP B5Z8H1
P	442	GLY	-	expression tag	UNP B5Z8H1
P	443	GLY	-	expression tag	UNP B5Z8H1

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Chain	Residue	Modelled	Actual	Comment	Reference
P	444	SER	-	expression tag	UNP B5Z8H1
P	445	THR	-	expression tag	UNP B5Z8H1
P	446	SER	-	expression tag	UNP B5Z8H1
P	447	TYR	-	expression tag	UNP B5Z8H1
P	448	ALA	-	expression tag	UNP B5Z8H1
P	449	ASP	-	expression tag	UNP B5Z8H1
P	450	SER	-	expression tag	UNP B5Z8H1
P	451	VAL	-	expression tag	UNP B5Z8H1
P	452	LYS	-	expression tag	UNP B5Z8H1
P	453	ASP	-	expression tag	UNP B5Z8H1
P	454	ARG	-	expression tag	UNP B5Z8H1
P	455	PHE	-	expression tag	UNP B5Z8H1
P	456	THR	-	expression tag	UNP B5Z8H1
P	457	ILE	-	expression tag	UNP B5Z8H1
P	458	SER	-	expression tag	UNP B5Z8H1
P	459	ARG	-	expression tag	UNP B5Z8H1
P	460	ASP	-	expression tag	UNP B5Z8H1
P	461	ASN	-	expression tag	UNP B5Z8H1
P	462	ALA	-	expression tag	UNP B5Z8H1
P	463	LYS	-	expression tag	UNP B5Z8H1
P	464	ASN	-	expression tag	UNP B5Z8H1
P	465	THR	-	expression tag	UNP B5Z8H1
P	466	VAL	-	expression tag	UNP B5Z8H1
P	467	TYR	-	expression tag	UNP B5Z8H1
P	468	LEU	-	expression tag	UNP B5Z8H1
P	469	GLU	-	expression tag	UNP B5Z8H1
P	470	MET	-	expression tag	UNP B5Z8H1
P	471	ASN	-	expression tag	UNP B5Z8H1
P	472	ASN	-	expression tag	UNP B5Z8H1
P	473	LEU	-	expression tag	UNP B5Z8H1
P	474	LYS	-	expression tag	UNP B5Z8H1
P	475	PRO	-	expression tag	UNP B5Z8H1
P	476	GLU	-	expression tag	UNP B5Z8H1
P	477	ASP	-	expression tag	UNP B5Z8H1
P	478	THR	-	expression tag	UNP B5Z8H1
P	479	ALA	-	expression tag	UNP B5Z8H1
P	480	VAL	-	expression tag	UNP B5Z8H1
P	481	TYR	-	expression tag	UNP B5Z8H1
P	482	TYR	-	expression tag	UNP B5Z8H1
P	483	CYS	-	expression tag	UNP B5Z8H1
P	484	ALA	-	expression tag	UNP B5Z8H1
P	485	ALA	-	expression tag	UNP B5Z8H1

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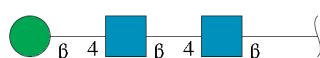
Chain	Residue	Modelled	Actual	Comment	Reference
P	486	LYS	-	expression tag	UNP B5Z8H1
P	487	GLY	-	expression tag	UNP B5Z8H1
P	488	ARG	-	expression tag	UNP B5Z8H1
P	489	TYR	-	expression tag	UNP B5Z8H1
P	490	SER	-	expression tag	UNP B5Z8H1
P	491	GLY	-	expression tag	UNP B5Z8H1
P	492	GLY	-	expression tag	UNP B5Z8H1
P	493	LEU	-	expression tag	UNP B5Z8H1
P	494	TYR	-	expression tag	UNP B5Z8H1
P	495	TYR	-	expression tag	UNP B5Z8H1
P	496	PRO	-	expression tag	UNP B5Z8H1
P	497	THR	-	expression tag	UNP B5Z8H1
P	498	ASN	-	expression tag	UNP B5Z8H1
P	499	TYR	-	expression tag	UNP B5Z8H1
P	500	ASP	-	expression tag	UNP B5Z8H1
P	501	TYR	-	expression tag	UNP B5Z8H1
P	502	TRP	-	expression tag	UNP B5Z8H1
P	503	GLY	-	expression tag	UNP B5Z8H1
P	504	GLN	-	expression tag	UNP B5Z8H1
P	505	GLY	-	expression tag	UNP B5Z8H1
P	506	THR	-	expression tag	UNP B5Z8H1
P	507	GLN	-	expression tag	UNP B5Z8H1
P	508	VAL	-	expression tag	UNP B5Z8H1
P	509	THR	-	expression tag	UNP B5Z8H1
P	510	VAL	-	expression tag	UNP B5Z8H1
P	511	SER	-	expression tag	UNP B5Z8H1
P	512	SER	-	expression tag	UNP B5Z8H1
P	513	HIS	-	expression tag	UNP B5Z8H1
P	514	HIS	-	expression tag	UNP B5Z8H1
P	515	HIS	-	expression tag	UNP B5Z8H1
P	516	HIS	-	expression tag	UNP B5Z8H1
P	517	HIS	-	expression tag	UNP B5Z8H1
P	518	HIS	-	expression tag	UNP B5Z8H1
P	519	GLU	-	expression tag	UNP B5Z8H1
P	520	PRO	-	expression tag	UNP B5Z8H1
P	521	GLU	-	expression tag	UNP B5Z8H1
P	522	ALA	-	expression tag	UNP B5Z8H1

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



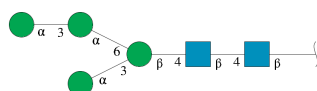
Mol	Chain	Residues	Atoms				AltConf	Trace
4	F	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 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.



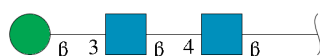
Mol	Chain	Residues	Atoms				AltConf	Trace
5	G	3	Total	C	N	O	0	0
			39	22	2	15		
5	I	3	Total	C	N	O	0	0
			39	22	2	15		
5	M	3	Total	C	N	O	0	0
			39	22	2	15		

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



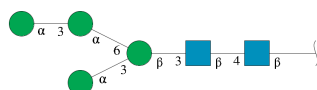
Mol	Chain	Residues	Atoms				AltConf	Trace
6	H	6	Total	C	N	O	0	0
			72	40	2	30		
6	N	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 7 is an oligosaccharide called beta-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
7	J	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 8 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
8	K	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 9 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.



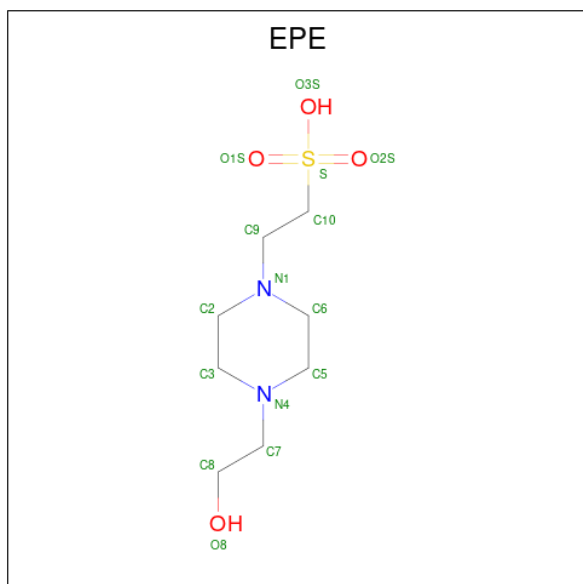
Mol	Chain	Residues	Atoms				AltConf	Trace
9	L	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 10 is GAMMA-AMINO-BUTANOIC ACID (CCD ID: ABU) (formula: C₄H₉NO₂) (labeled as "Ligand of Interest" by depositor).



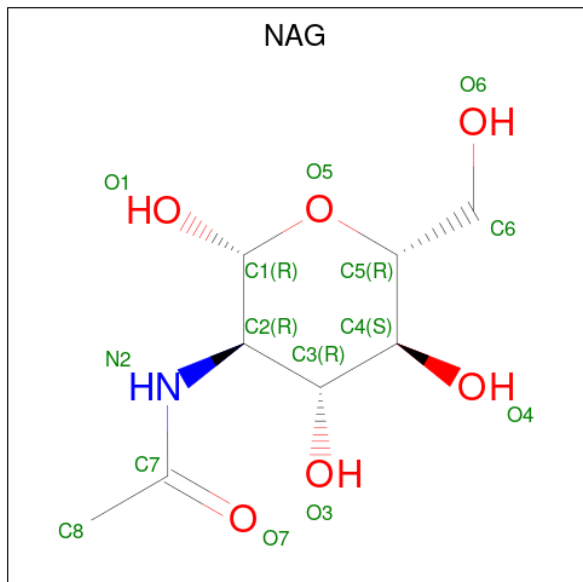
Mol	Chain	Residues	Atoms				AltConf
10	B	1	Total	C	N	O	0
			7	4	1	2	
10	E	1	Total	C	N	O	0
			7	4	1	2	

- Molecule 11 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).

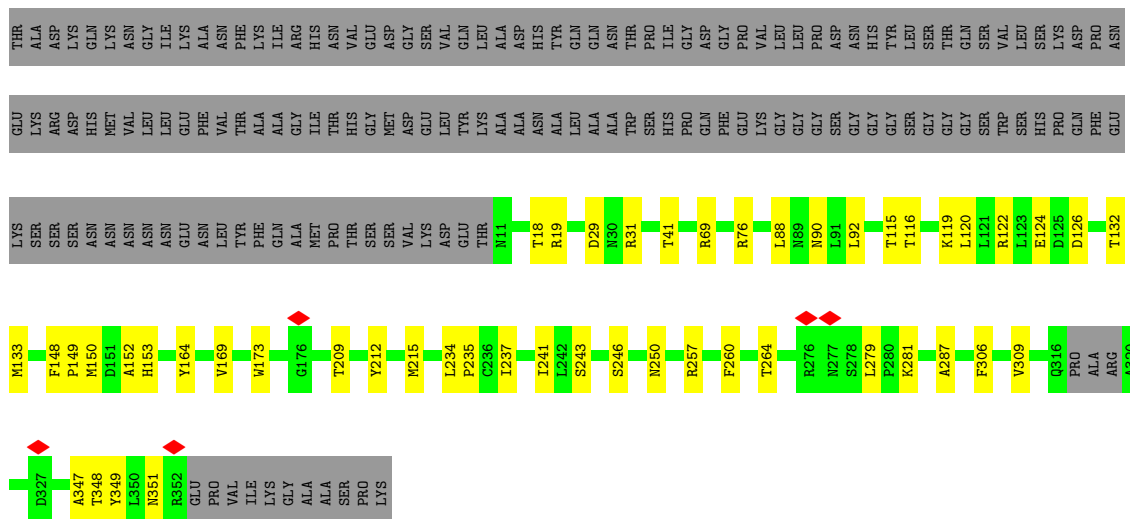


Mol	Chain	Residues	Atoms					AltConf
11	C	1	Total	C	N	O	S	0
			15	8	2	4	1	

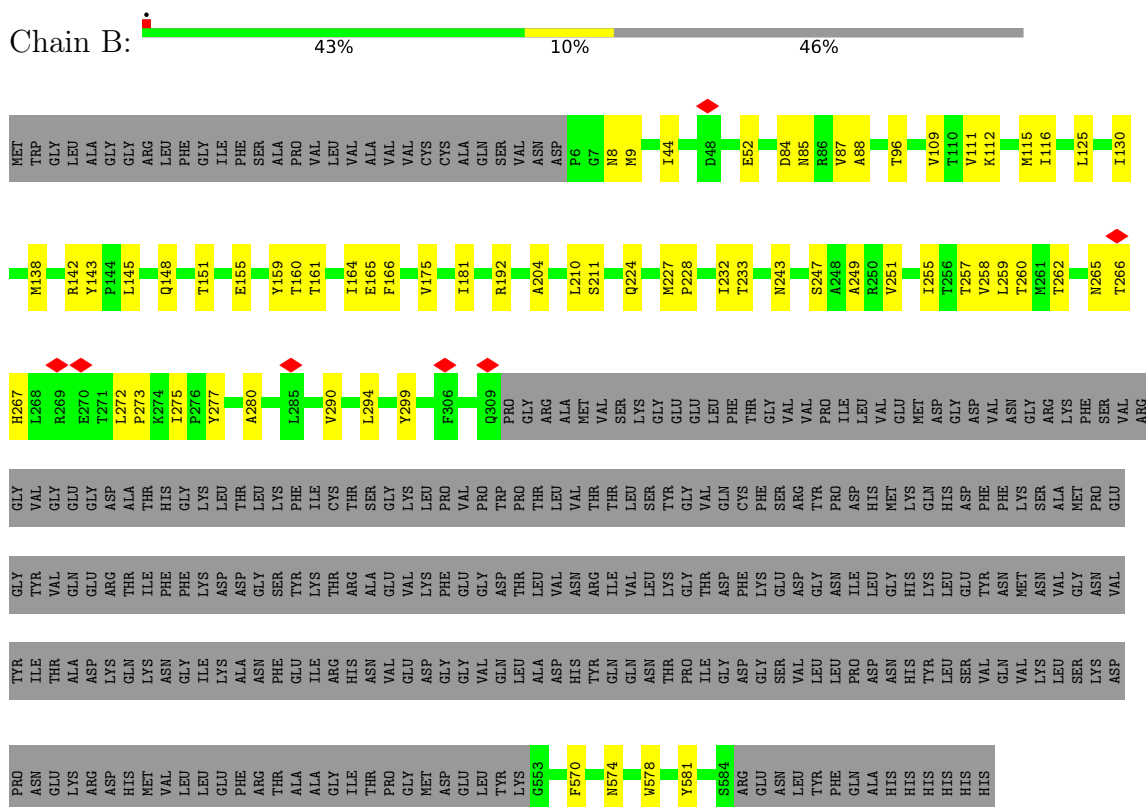
- Molecule 12 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



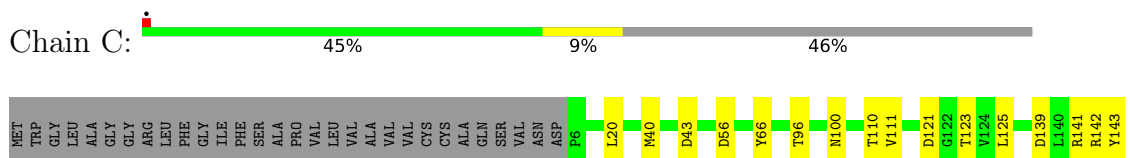
Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
12	D	1	14	8	1	5	0



- Molecule 2: Gamma-aminobutyric acid receptor subunit beta-3, Green fluorescent protein

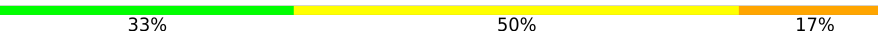


- Molecule 2: Gamma-aminobutyric acid receptor subunit beta-3, Green fluorescent protein





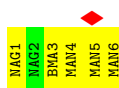
- Molecule 6: α -D-mannopyranose-(1-3)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain H:  33% 50% 17%



- Molecule 6: α -D-mannopyranose-(1-3)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain N:  17% 17% 83%



- Molecule 7: β -D-mannopyranose-(1-3)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain J:  33% 67%



- Molecule 8: α -D-mannopyranose-(1-3)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] β -D-mannopyranose-(1-3)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain K:  100%



- Molecule 9: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-6)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain L:  50% 50%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	96876	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOCONTINUUM (6k x 4k)	Depositor
Maximum map value	36.189	Depositor
Minimum map value	-23.614	Depositor
Average map value	0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	3	Depositor
Map size (Å)	286.0, 286.0, 286.0	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.65, 0.65, 0.65	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EPE, MAN, ABU, 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	0.15	0/2813	0.45	0/3826
1	D	0.14	0/2786	0.43	0/3788
2	B	0.15	0/2820	0.44	0/3835
2	C	0.15	0/2841	0.46	0/3863
2	E	0.14	0/2820	0.44	0/3834
3	P	0.16	0/954	0.44	0/1292
All	All	0.15	0/15034	0.44	0/20438

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2745	0	2719	51	0
1	D	2719	0	2700	35	0
2	B	2747	0	2737	48	0
2	C	2767	0	2762	42	0
2	E	2747	0	2742	44	0
3	P	931	0	879	14	0
4	F	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	G	39	0	34	1	0
5	I	39	0	34	0	0
5	M	39	0	34	0	0
6	H	72	0	61	2	0
6	N	72	0	61	1	0
7	J	39	0	34	0	0
8	K	72	0	61	0	0
9	L	28	0	25	0	0
10	B	7	0	0	1	0
10	E	7	0	0	0	0
11	C	15	0	18	0	0
12	D	14	0	13	0	0
All	All	15127	0	14939	213	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 (213) 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)
2:E:263:THR:HG23	2:E:267:HIS:CE1	1.85	1.11
1:A:133:MET:HE1	2:E:110:THR:HG23	1.57	0.86
2:E:263:THR:CG2	2:E:267:HIS:CE1	2.58	0.86
1:D:115:THR:O	1:D:116:THR:HG23	1.81	0.80
1:D:124:GLU:N	1:D:124:GLU:OE1	2.17	0.76
2:B:112:LYS:NZ	5:G:2:NAG:O7	2.22	0.71
1:A:347:ALA:O	1:A:351:ASN:ND2	2.25	0.68
3:P:405:LEU:HD23	3:P:406:ARG:N	2.10	0.66
2:E:144:PRO:HD3	2:E:281:ILE:HG23	1.78	0.66
1:A:205:ILE:HD13	1:A:214:ILE:HD11	1.78	0.65
1:A:251:ARG:NH1	1:A:313:THR:OG1	2.30	0.65
1:D:348:THR:HG23	1:D:349:TYR:HD1	1.60	0.65
2:B:233:THR:HG21	2:B:570:PHE:CE2	2.32	0.63
1:A:256:ALA:HB1	2:B:251:VAL:HG21	1.80	0.63
2:E:574:ASN:OD1	2:E:578:TRP:NE1	2.31	0.62
2:B:142:ARG:O	2:B:145:LEU:N	2.31	0.62
1:A:109:SER:HB2	2:E:109:VAL:HG22	1.81	0.62
2:B:280:ALA:HB2	2:B:581:TYR:CE2	2.35	0.61
2:C:139:ASP:OD1	2:C:141:ARG:NH2	2.33	0.61
2:E:280:ALA:HB2	2:E:581:TYR:HE2	1.66	0.61
2:C:66:TYR:CZ	2:C:125:LEU:HD13	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:246:SER:O	1:D:257:ARG:NE	2.34	0.61
2:C:293:PHE:CE1	2:C:297:LEU:HD11	2.36	0.60
1:D:90:ASN:ND2	1:D:119:LYS:O	2.35	0.60
2:B:44:ILE:HG21	2:B:181:ILE:HD11	1.84	0.60
2:B:574:ASN:OD1	2:B:578:TRP:NE1	2.35	0.59
2:C:143:TYR:HB3	2:C:144:PRO:HD3	1.83	0.59
2:E:271:THR:C	2:E:272:LEU:HD12	2.27	0.59
2:C:224:GLN:O	2:C:268:LEU:HD21	2.02	0.58
1:A:8:ASP:O	1:A:13:ASN:ND2	2.34	0.58
2:E:142:ARG:NH2	2:E:145:LEU:O	2.37	0.58
2:E:242:ILE:O	2:E:250:ARG:NH2	2.34	0.58
2:B:142:ARG:NH1	2:B:581:TYR:O	2.37	0.58
1:D:120:LEU:HD13	2:E:158:GLY:HA2	1.85	0.58
2:E:66:TYR:CZ	2:E:125:LEU:HD13	2.38	0.58
1:A:215:MET:HE2	1:A:215:MET:HA	1.85	0.58
1:D:120:LEU:HD11	1:D:122:ARG:HE	1.67	0.58
2:B:262:THR:O	2:B:265:ASN:OD1	2.21	0.57
1:A:314:LYS:NZ	2:E:243:ASN:HB2	2.20	0.57
2:B:233:THR:HG21	2:B:570:PHE:CD2	2.40	0.57
2:E:280:ALA:HB2	2:E:581:TYR:CE2	2.41	0.56
1:A:133:MET:CE	2:E:110:THR:HG23	2.34	0.55
1:A:15:THR:O	1:A:18:THR:HG22	2.07	0.55
1:A:115:THR:HG23	2:B:96:THR:O	2.06	0.54
1:D:306:PHE:HA	1:D:309:VAL:HG12	1.89	0.54
3:P:447:TYR:OH	3:P:457:ILE:N	2.39	0.54
2:E:40:MET:HG3	2:E:208:LEU:HD12	1.88	0.54
2:B:159:TYR:HB2	2:B:164:ILE:HD12	1.88	0.54
2:B:267:HIS:CD2	2:B:267:HIS:O	2.60	0.54
1:A:336:VAL:O	1:A:340:THR:HG23	2.08	0.54
1:A:310:ASN:O	1:A:314:LYS:HB2	2.08	0.53
2:E:143:TYR:O	2:E:280:ALA:HB3	2.08	0.53
2:B:255:ILE:O	2:B:259:LEU:HD13	2.07	0.53
2:C:283:MET:HA	2:C:283:MET:HE2	1.89	0.53
1:A:250:ASN:OD1	1:A:253:SER:N	2.39	0.53
2:B:143:TYR:O	2:B:280:ALA:HB3	2.10	0.52
2:C:179:GLU:OE2	2:C:191:HIS:ND1	2.42	0.52
2:C:142:ARG:NH1	2:C:581:TYR:O	2.43	0.52
2:B:138:MET:HB3	2:B:148:GLN:OE1	2.09	0.52
2:E:290:VAL:O	2:E:294:LEU:HD23	2.09	0.52
3:P:472:ASN:OD1	3:P:474:LYS:NZ	2.42	0.51
1:A:46:ASP:OD2	1:A:69:ARG:NH1	2.38	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:287:ALA:HB2	1:D:349:TYR:HE2	1.75	0.51
2:B:224:GLN:OE1	2:C:269:ARG:NH2	2.44	0.51
1:D:260:PHE:CE1	2:E:255:ILE:HD13	2.46	0.51
1:A:69:ARG:HG2	1:A:132:THR:HG23	1.93	0.50
1:D:348:THR:HG23	1:D:349:TYR:CD1	2.44	0.50
1:A:92:LEU:HB3	1:A:96:ILE:HD12	1.93	0.50
2:B:211:SER:HB2	6:H:1:NAG:H82	1.93	0.50
2:C:302:VAL:HG12	2:C:558:ILE:HG21	1.93	0.50
1:A:75:GLU:O	1:A:78:ARG:NH2	2.44	0.50
2:B:115:MET:HE3	2:B:125:LEU:CD2	2.42	0.50
2:C:144:PRO:HD2	2:C:280:ALA:HB3	1.94	0.50
2:E:263:THR:CG2	2:E:267:HIS:ND1	2.75	0.50
2:E:213:ARG:HG3	6:N:1:NAG:H81	1.92	0.50
2:C:66:TYR:CE1	2:C:125:LEU:HD13	2.47	0.50
1:A:314:LYS:HZ1	2:E:243:ASN:HB2	1.77	0.49
2:C:110:THR:O	2:C:111:VAL:HG13	2.13	0.49
2:E:243:ASN:ND2	2:E:245:ASP:OD2	2.43	0.49
2:B:52:GLU:OE1	2:B:52:GLU:N	2.46	0.49
1:D:243:SER:OG	1:D:264:THR:HG21	2.13	0.49
2:C:286:MET:SD	2:C:286:MET:C	2.96	0.48
2:C:96:THR:HG22	2:C:166:PHE:HZ	1.77	0.48
2:C:251:VAL:HG23	2:C:299:TYR:CE2	2.49	0.48
1:A:265:VAL:HG11	2:E:235:LEU:HD12	1.96	0.48
1:D:88:LEU:HD13	1:D:92:LEU:CD2	2.44	0.48
2:B:84:ASP:O	2:B:87:VAL:HG12	2.14	0.48
2:B:272:LEU:HB2	2:B:273:PRO:HD2	1.96	0.48
1:A:33:ARG:HG3	1:A:34:PRO:HD2	1.95	0.48
2:B:160:THR:HG22	2:B:161:THR:N	2.29	0.48
2:B:265:ASN:OD1	2:B:266:THR:N	2.47	0.48
2:C:110:THR:O	2:C:111:VAL:CG1	2.62	0.48
1:A:17:PHE:CZ	1:A:83:MET:HG2	2.50	0.47
2:B:249:ALA:HB1	2:C:251:VAL:HG21	1.96	0.47
1:A:133:MET:HE1	2:E:110:THR:CG2	2.39	0.47
1:A:310:ASN:OD1	1:A:314:LYS:HE3	2.13	0.47
1:D:41:THR:HG23	1:D:41:THR:O	2.14	0.47
3:P:469:GLU:HA	3:P:469:GLU:OE1	2.15	0.47
1:A:109:SER:CB	2:E:109:VAL:HG13	2.43	0.47
1:D:281:LYS:HG3	1:D:281:LYS:O	2.13	0.47
1:D:69:ARG:HG2	1:D:132:THR:HG23	1.96	0.47
2:E:146:ASP:OD2	2:E:148:GLN:NE2	2.38	0.47
2:C:290:VAL:O	2:C:294:LEU:HD23	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:85:ASN:ND2	2:B:112:LYS:O	2.48	0.46
2:B:257:THR:O	2:B:260:THR:HG22	2.14	0.46
2:E:257:THR:HG22	2:E:292:VAL:HG11	1.97	0.46
2:C:227:MET:HB3	2:C:228:PRO:HD3	1.98	0.46
1:D:164:TYR:HB3	1:D:169:VAL:HG23	1.97	0.46
1:A:196:MET:HE3	1:A:196:MET:HA	1.97	0.46
2:C:261:MET:HE1	2:C:288:CYS:HB3	1.97	0.46
2:E:101:ASP:O	2:E:101:ASP:OD1	2.34	0.46
2:E:144:PRO:CD	2:E:281:ILE:HG23	2.46	0.46
3:P:480:VAL:HG23	3:P:480:VAL:O	2.16	0.46
2:C:580:TYR:HD1	2:C:581:TYR:HD1	1.65	0.45
2:E:146:ASP:OD1	2:E:146:ASP:N	2.49	0.45
2:C:166:PHE:HB3	2:C:208:LEU:HD11	1.98	0.45
3:P:431:GLU:H	3:P:431:GLU:CD	2.25	0.45
1:A:114:MET:HE1	2:B:130:ILE:HG12	1.99	0.45
2:B:88:ALA:HB2	2:B:116:ILE:HD11	1.99	0.45
2:B:255:ILE:HA	2:B:258:VAL:HG22	1.99	0.45
2:C:40:MET:HG3	2:C:208:LEU:HD12	1.99	0.45
3:P:440:TRP:O	3:P:441:SER:OG	2.25	0.45
2:B:164:ILE:HG21	2:B:166:PHE:CE1	2.52	0.45
3:P:2:VAL:HG21	3:P:418:TYR:HD2	1.82	0.45
2:C:56:ASP:OD1	2:C:56:ASP:C	2.60	0.44
2:B:262:THR:O	2:B:266:THR:HG23	2.17	0.44
2:B:109:VAL:O	2:B:111:VAL:N	2.50	0.44
1:A:256:ALA:O	1:A:260:PHE:CD2	2.71	0.44
1:A:300:PHE:CD2	1:A:304:ILE:HD11	2.53	0.44
1:A:294:VAL:HG21	1:A:341:PHE:CE1	2.53	0.44
1:D:250:ASN:ND2	2:E:307:PHE:CG	2.86	0.44
2:C:190:GLU:OE2	2:C:192:ARG:NH1	2.50	0.44
1:D:148:PHE:HB3	1:D:149:PRO:HD3	1.99	0.44
1:A:78:ARG:HH21	1:A:78:ARG:HG3	1.83	0.44
2:C:142:ARG:O	2:C:145:LEU:N	2.51	0.44
1:D:150:MET:SD	1:D:349:TYR:HB3	2.57	0.44
1:D:237:ILE:O	1:D:241:ILE:HG12	2.18	0.44
2:C:43:ASP:OD1	2:C:176:THR:OG1	2.35	0.43
1:D:18:THR:HG23	1:D:19:ARG:N	2.33	0.43
2:B:155:GLU:OE2	10:B:601:ABU:N	2.51	0.43
2:B:228:PRO:O	2:B:232:ILE:HG12	2.18	0.43
2:B:247:SER:HA	2:B:299:TYR:OH	2.18	0.43
1:D:115:THR:O	1:D:116:THR:CG2	2.61	0.43
2:E:554:ARG:O	2:E:558:ILE:HG12	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:488:ARG:HH21	3:P:488:ARG:HG3	1.83	0.43
2:E:247:SER:HA	2:E:299:TYR:OH	2.19	0.43
2:B:164:ILE:HG22	2:B:165:GLU:N	2.34	0.43
1:A:13:ASN:O	1:A:16:ILE:HG22	2.19	0.43
1:A:256:ALA:O	1:A:259:VAL:HG22	2.19	0.43
2:C:96:THR:HG22	2:C:166:PHE:CZ	2.53	0.43
2:E:301:PHE:O	2:E:305:ILE:HG12	2.18	0.43
1:A:115:THR:C	1:A:116:THR:HG23	2.43	0.43
2:C:20:LEU:HD12	1:D:31:ARG:NH2	2.33	0.43
3:P:406:ARG:HD2	3:P:406:ARG:C	2.44	0.43
2:B:151:THR:HG22	2:B:211:SER:CB	2.48	0.43
2:C:281:ILE:HG23	2:C:282:ASP:N	2.34	0.43
1:D:173:TRP:CE2	1:D:215:MET:HE3	2.54	0.43
2:B:290:VAL:O	2:B:294:LEU:HD13	2.18	0.43
2:C:179:GLU:HA	2:C:179:GLU:OE1	2.19	0.43
2:E:190:GLU:OE1	2:E:213:ARG:NH2	2.38	0.43
1:A:49:VAL:HA	1:A:66:VAL:HG12	2.01	0.43
1:D:29:ASP:O	1:D:76:ARG:NH2	2.51	0.43
1:D:148:PHE:CE2	1:D:279:LEU:HD11	2.53	0.43
1:D:347:ALA:O	1:D:351:ASN:ND2	2.48	0.43
1:A:306:PHE:O	1:A:309:VAL:HG12	2.18	0.42
1:A:310:ASN:O	1:A:313:THR:HG22	2.19	0.42
2:C:121:ASP:OD2	2:C:123:THR:OG1	2.29	0.42
2:B:175:VAL:HG21	2:B:210:LEU:HD13	2.00	0.42
2:B:192:ARG:NH2	6:H:1:NAG:H83	2.34	0.42
3:P:5:VAL:HG13	3:P:5:VAL:O	2.19	0.42
1:A:101:THR:HB	1:A:171:TYR:OH	2.18	0.42
2:E:180:ARG:HD2	2:E:180:ARG:O	2.19	0.42
1:A:13:ASN:OD1	1:A:17:PHE:CE2	2.72	0.42
1:A:292:ILE:HG22	1:A:296:TYR:CE2	2.55	0.42
1:D:152:ALA:O	1:D:153:HIS:ND1	2.52	0.42
1:A:148:PHE:CD1	1:A:279:LEU:HD21	2.54	0.42
2:C:251:VAL:CG2	2:C:299:TYR:HE2	2.33	0.42
2:E:66:TYR:CE2	2:E:125:LEU:HD13	2.54	0.42
2:B:161:THR:HG23	2:B:204:ALA:O	2.20	0.42
1:A:174:THR:HG23	1:A:175:ASN:N	2.33	0.42
1:A:109:SER:HB3	2:E:109:VAL:HG13	2.01	0.41
1:A:16:ILE:O	1:A:20:ILE:HG13	2.20	0.41
1:A:314:LYS:HZ1	2:E:243:ASN:CB	2.33	0.41
2:C:228:PRO:O	2:C:232:ILE:HG12	2.20	0.41
1:A:148:PHE:HB3	1:A:149:PRO:HD3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:265:ASN:OD1	2:B:266:THR:HG23	2.20	0.41
1:D:149:PRO:HG2	1:D:150:MET:HE3	2.01	0.41
2:B:8:ASN:OD1	2:B:9:MET:N	2.54	0.41
2:B:243:ASN:HB2	2:C:307:PHE:HD2	1.85	0.41
2:B:275:ILE:HG13	2:B:277:TYR:CE1	2.56	0.41
2:C:275:ILE:HG21	2:C:277:TYR:CZ	2.55	0.41
1:D:126:ASP:OD1	1:D:126:ASP:C	2.62	0.41
2:E:263:THR:HG22	2:E:267:HIS:ND1	2.34	0.41
2:B:258:VAL:HG23	2:B:259:LEU:HD12	2.02	0.41
2:C:306:PHE:CE2	2:C:307:PHE:CE1	3.09	0.41
2:E:85:ASN:OD1	2:E:86:ARG:N	2.53	0.41
1:A:45:THR:HG22	1:A:46:ASP:N	2.35	0.41
1:D:260:PHE:CD1	2:E:255:ILE:HD13	2.56	0.41
2:C:100:ASN:HD22	2:C:153:GLU:HG2	1.86	0.41
2:C:561:TRP:O	2:C:565:VAL:HG23	2.21	0.41
3:P:418:TYR:CE2	3:P:501:TYR:HE2	2.39	0.41
2:C:299:TYR:HA	2:C:302:VAL:HG22	2.02	0.41
1:D:234:LEU:N	1:D:235:PRO:HD2	2.36	0.41
3:P:507:GLN:OE1	3:P:509:THR:OG1	2.30	0.41
1:A:115:THR:O	1:A:116:THR:OG1	2.31	0.40
1:A:259:VAL:O	1:A:263:THR:HG23	2.21	0.40
1:D:132:THR:HG22	1:D:133:MET:N	2.36	0.40
2:B:227:MET:HB3	2:B:228:PRO:HD3	2.04	0.40
3:P:455:PHE:N	3:P:455:PHE:CD1	2.90	0.40
2:E:261:MET:HE2	2:E:264:ILE:HD11	2.03	0.40
1:A:301:SER:O	1:A:305:GLU:HG2	2.21	0.40
2:C:185:GLN:HA	2:C:185:GLN:OE1	2.22	0.40
1:D:209:THR:OG1	1:D:212:TYR:OH	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	338/679 (50%)	331 (98%)	6 (2%)	1 (0%)	36	67
1	D	335/679 (49%)	324 (97%)	11 (3%)	0	100	100
2	B	332/623 (53%)	327 (98%)	5 (2%)	0	100	100
2	C	334/623 (54%)	324 (97%)	10 (3%)	0	100	100
2	E	332/623 (53%)	326 (98%)	6 (2%)	0	100	100
3	P	116/541 (21%)	113 (97%)	3 (3%)	0	100	100
All	All	1787/3768 (47%)	1745 (98%)	41 (2%)	1 (0%)	49	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	THR

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	306/589 (52%)	306 (100%)	0	100	100
1	D	302/589 (51%)	302 (100%)	0	100	100
2	B	300/544 (55%)	300 (100%)	0	100	100
2	C	302/544 (56%)	302 (100%)	0	100	100
2	E	300/544 (55%)	300 (100%)	0	100	100
3	P	96/443 (22%)	96 (100%)	0	100	100
All	All	1606/3253 (49%)	1606 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	218	HIS
2	B	64	GLN

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Mol	Chain	Res	Type
2	B	191	HIS
2	B	267	HIS
2	C	185	GLN
1	D	42	GLN
1	D	218	HIS
1	D	220	HIS
2	E	303	ASN
3	P	464	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

34 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	F	1	1,4	14,14,15	0.74	0	17,19,21	0.88	0
4	NAG	F	2	4	14,14,15	0.71	0	17,19,21	0.86	0
5	NAG	G	1	1,5	14,14,15	0.62	0	17,19,21	1.47	3 (17%)
5	NAG	G	2	5	14,14,15	0.76	0	17,19,21	0.88	0
5	BMA	G	3	5	11,11,12	0.78	0	15,15,17	1.97	1 (6%)
6	NAG	H	1	6,2	14,14,15	0.77	0	17,19,21	1.05	1 (5%)
6	NAG	H	2	6	14,14,15	0.67	0	17,19,21	0.98	0
6	BMA	H	3	6	11,11,12	0.80	0	15,15,17	1.84	3 (20%)
6	MAN	H	4	6	11,11,12	0.66	0	15,15,17	1.15	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	MAN	H	5	6	11,11,12	0.67	0	15,15,17	1.16	1 (6%)
6	MAN	H	6	6	11,11,12	0.73	0	15,15,17	0.88	0
5	NAG	I	1	2,5	14,14,15	0.74	0	17,19,21	1.02	1 (5%)
5	NAG	I	2	5	14,14,15	0.71	0	17,19,21	0.84	0
5	BMA	I	3	5	11,11,12	0.80	0	15,15,17	1.68	1 (6%)
7	NAG	J	1	2,7	14,14,15	0.79	0	17,19,21	0.98	1 (5%)
7	NAG	J	2	7	14,14,15	0.75	0	17,19,21	0.83	0
7	BMA	J	3	7	11,11,12	0.81	0	15,15,17	1.73	1 (6%)
8	NAG	K	1	2,8	14,14,15	0.66	0	17,19,21	2.60	5 (29%)
8	NAG	K	2	8	14,14,15	0.79	0	17,19,21	0.88	1 (5%)
8	BMA	K	3	8	11,11,12	0.75	0	15,15,17	2.13	3 (20%)
8	MAN	K	4	8	11,11,12	0.65	0	15,15,17	1.48	2 (13%)
8	MAN	K	5	8	11,11,12	0.69	0	15,15,17	0.98	1 (6%)
8	MAN	K	6	8	11,11,12	0.69	0	15,15,17	1.00	1 (6%)
9	NAG	L	1	1,9	14,14,15	0.67	0	17,19,21	0.95	1 (5%)
9	NAG	L	2	9	14,14,15	0.69	0	17,19,21	0.80	0
5	NAG	M	1	2,5	14,14,15	0.74	0	17,19,21	0.96	1 (5%)
5	NAG	M	2	5	14,14,15	0.73	0	17,19,21	1.01	1 (5%)
5	BMA	M	3	5	11,11,12	0.76	0	15,15,17	2.00	1 (6%)
6	NAG	N	1	6,2	14,14,15	0.76	0	17,19,21	1.02	0
6	NAG	N	2	6	14,14,15	0.70	0	17,19,21	0.91	0
6	BMA	N	3	6	11,11,12	0.76	0	15,15,17	1.82	2 (13%)
6	MAN	N	4	6	11,11,12	0.66	0	15,15,17	1.15	1 (6%)
6	MAN	N	5	6	11,11,12	0.69	0	15,15,17	1.03	1 (6%)
6	MAN	N	6	6	11,11,12	0.67	0	15,15,17	1.11	1 (6%)

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	F	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1
5	NAG	G	1	1,5	-	1/6/23/26	0/1/1/1
5	NAG	G	2	5	-	0/6/23/26	0/1/1/1
5	BMA	G	3	5	-	0/2/19/22	0/1/1/1
6	NAG	H	1	6,2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	H	2	6	-	0/6/23/26	0/1/1/1
6	BMA	H	3	6	-	0/2/19/22	0/1/1/1
6	MAN	H	4	6	-	0/2/19/22	0/1/1/1
6	MAN	H	5	6	-	1/2/19/22	0/1/1/1
6	MAN	H	6	6	-	1/2/19/22	0/1/1/1
5	NAG	I	1	2,5	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	2/6/23/26	0/1/1/1
5	BMA	I	3	5	-	0/2/19/22	0/1/1/1
7	NAG	J	1	2,7	-	0/6/23/26	0/1/1/1
7	NAG	J	2	7	-	0/6/23/26	0/1/1/1
7	BMA	J	3	7	-	1/2/19/22	0/1/1/1
8	NAG	K	1	2,8	-	0/6/23/26	0/1/1/1
8	NAG	K	2	8	-	0/6/23/26	0/1/1/1
8	BMA	K	3	8	-	1/2/19/22	0/1/1/1
8	MAN	K	4	8	-	2/2/19/22	0/1/1/1
8	MAN	K	5	8	-	0/2/19/22	0/1/1/1
8	MAN	K	6	8	-	0/2/19/22	0/1/1/1
9	NAG	L	1	1,9	-	1/6/23/26	0/1/1/1
9	NAG	L	2	9	-	1/6/23/26	0/1/1/1
5	NAG	M	1	2,5	-	1/6/23/26	0/1/1/1
5	NAG	M	2	5	-	2/6/23/26	0/1/1/1
5	BMA	M	3	5	-	0/2/19/22	0/1/1/1
6	NAG	N	1	6,2	-	0/6/23/26	0/1/1/1
6	NAG	N	2	6	-	0/6/23/26	0/1/1/1
6	BMA	N	3	6	-	0/2/19/22	0/1/1/1
6	MAN	N	4	6	-	0/2/19/22	0/1/1/1
6	MAN	N	5	6	-	0/2/19/22	0/1/1/1
6	MAN	N	6	6	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	K	1	NAG	C1-O5-C5	8.69	123.83	112.19
5	M	3	BMA	C1-O5-C5	6.67	121.12	112.19
5	G	3	BMA	C1-O5-C5	6.61	121.04	112.19
8	K	3	BMA	C1-O5-C5	6.27	120.59	112.19
7	J	3	BMA	C1-O5-C5	5.24	119.21	112.19
6	H	3	BMA	C1-O5-C5	5.17	119.12	112.19
6	N	3	BMA	C1-O5-C5	5.14	119.08	112.19
5	I	3	BMA	C1-O5-C5	4.94	118.81	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	K	3	BMA	O3-C3-C4	3.84	119.43	110.38
8	K	4	MAN	C1-O5-C5	3.62	117.04	112.19
5	G	1	NAG	C4-C3-C2	-3.57	105.79	111.02
6	N	4	MAN	C1-O5-C5	3.42	116.77	112.19
6	H	5	MAN	C1-O5-C5	3.41	116.76	112.19
6	N	6	MAN	C1-O5-C5	3.24	116.53	112.19
6	H	4	MAN	C1-O5-C5	3.22	116.50	112.19
8	K	1	NAG	O5-C1-C2	3.21	116.25	111.29
8	K	1	NAG	C4-C3-C2	-3.03	106.58	111.02
8	K	4	MAN	O3-C3-C4	2.88	117.16	110.38
8	K	1	NAG	O4-C4-C5	2.87	116.39	109.32
6	N	3	BMA	O3-C3-C4	2.81	117.01	110.38
6	H	3	BMA	C3-C4-C5	2.77	115.26	110.23
6	N	5	MAN	C1-O5-C5	2.68	115.78	112.19
5	G	1	NAG	C2-N2-C7	2.63	126.42	122.90
8	K	6	MAN	C1-O5-C5	2.58	115.65	112.19
5	I	1	NAG	O5-C1-C2	-2.52	107.39	111.29
8	K	5	MAN	C1-O5-C5	2.49	115.52	112.19
5	M	1	NAG	C1-O5-C5	2.37	115.36	112.19
5	G	1	NAG	O4-C4-C5	2.31	115.02	109.32
8	K	1	NAG	C3-C4-C5	-2.28	106.09	110.23
8	K	2	NAG	C1-O5-C5	2.26	115.22	112.19
6	H	3	BMA	O4-C4-C3	-2.21	105.16	110.38
9	L	1	NAG	C1-O5-C5	2.20	115.14	112.19
7	J	1	NAG	O5-C1-C2	-2.09	108.06	111.29
6	H	1	NAG	C4-C3-C2	2.06	114.04	111.02
8	K	3	BMA	O3-C3-C2	-2.02	105.92	110.05
5	M	2	NAG	C2-N2-C7	2.00	125.58	122.90

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	G	1	NAG	C1-C2-N2-C7
6	H	1	NAG	O5-C5-C6-O6
6	H	1	NAG	C4-C5-C6-O6
8	K	4	MAN	O5-C5-C6-O6
6	H	5	MAN	O5-C5-C6-O6
9	L	1	NAG	O5-C5-C6-O6
5	M	1	NAG	O5-C5-C6-O6
6	H	6	MAN	O5-C5-C6-O6
7	J	3	BMA	O5-C5-C6-O6

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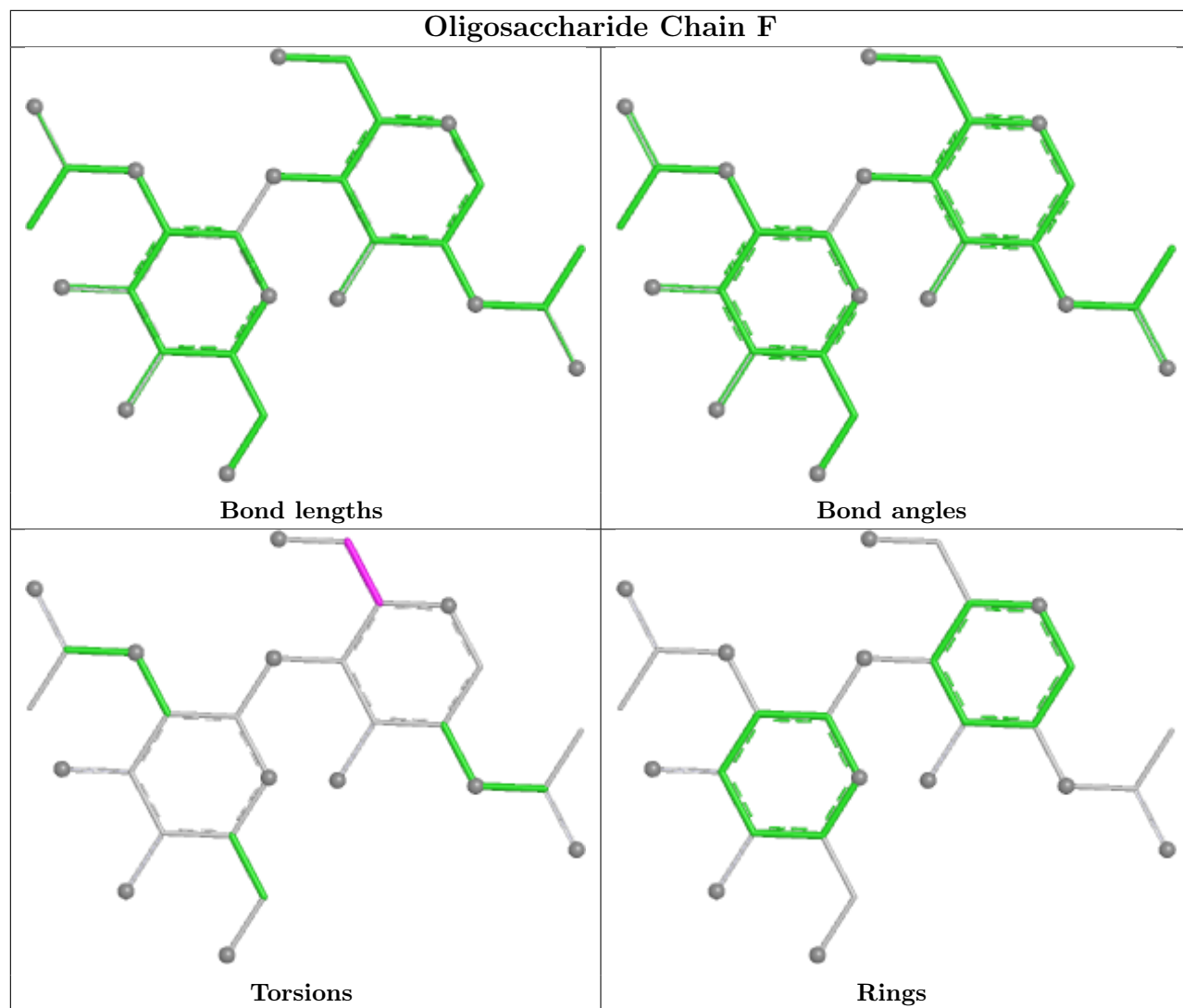
Mol	Chain	Res	Type	Atoms
5	I	2	NAG	C3-C2-N2-C7
5	M	2	NAG	C3-C2-N2-C7
4	F	1	NAG	C4-C5-C6-O6
9	L	2	NAG	C4-C5-C6-O6
5	I	2	NAG	C1-C2-N2-C7
5	M	2	NAG	C4-C5-C6-O6
8	K	3	BMA	O5-C5-C6-O6
8	K	4	MAN	C4-C5-C6-O6

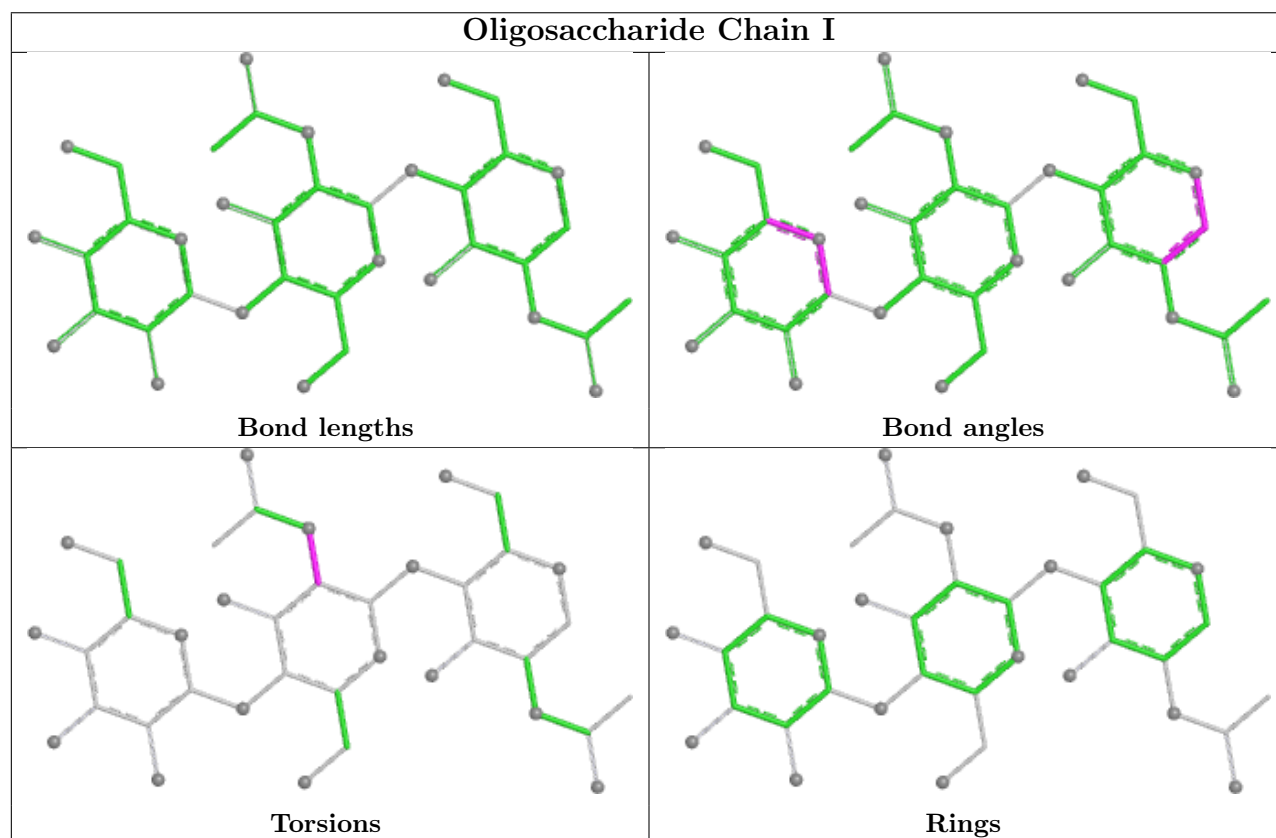
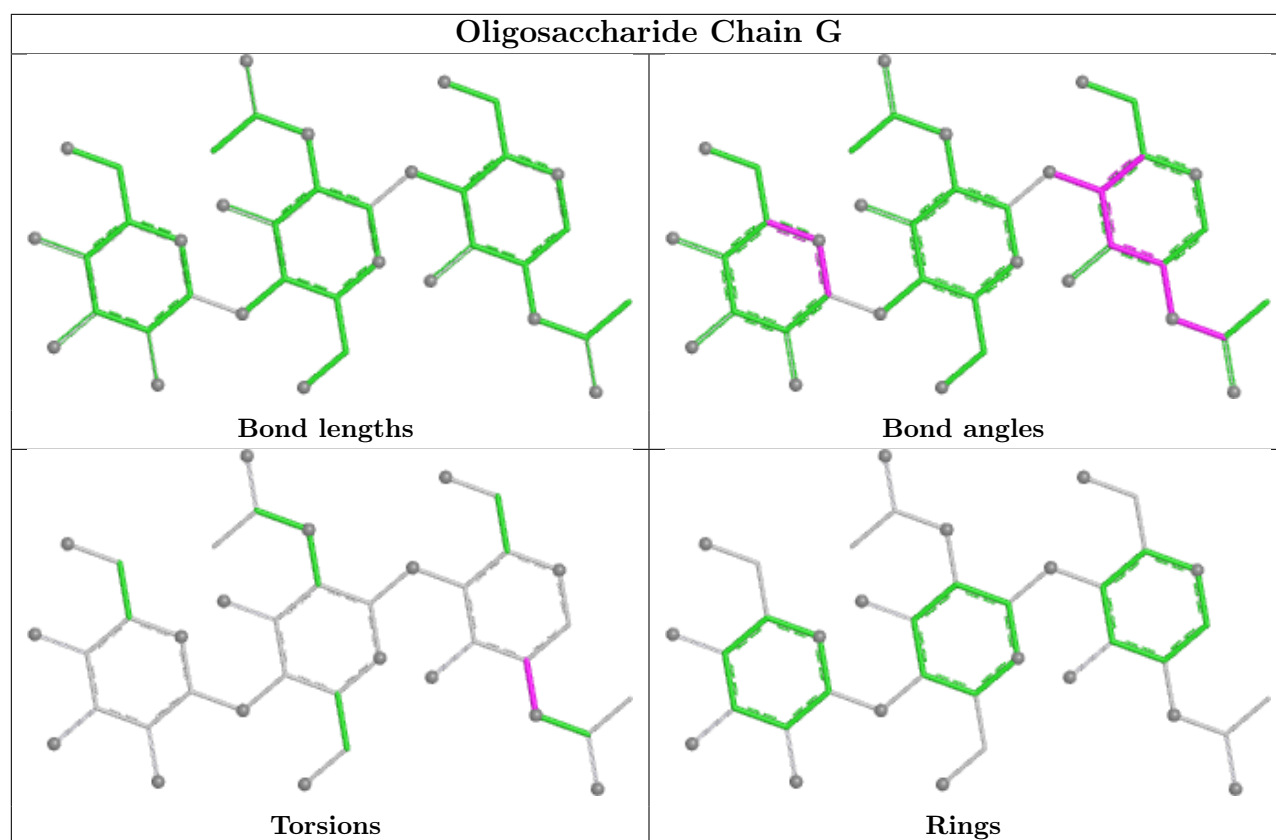
There are no ring outliers.

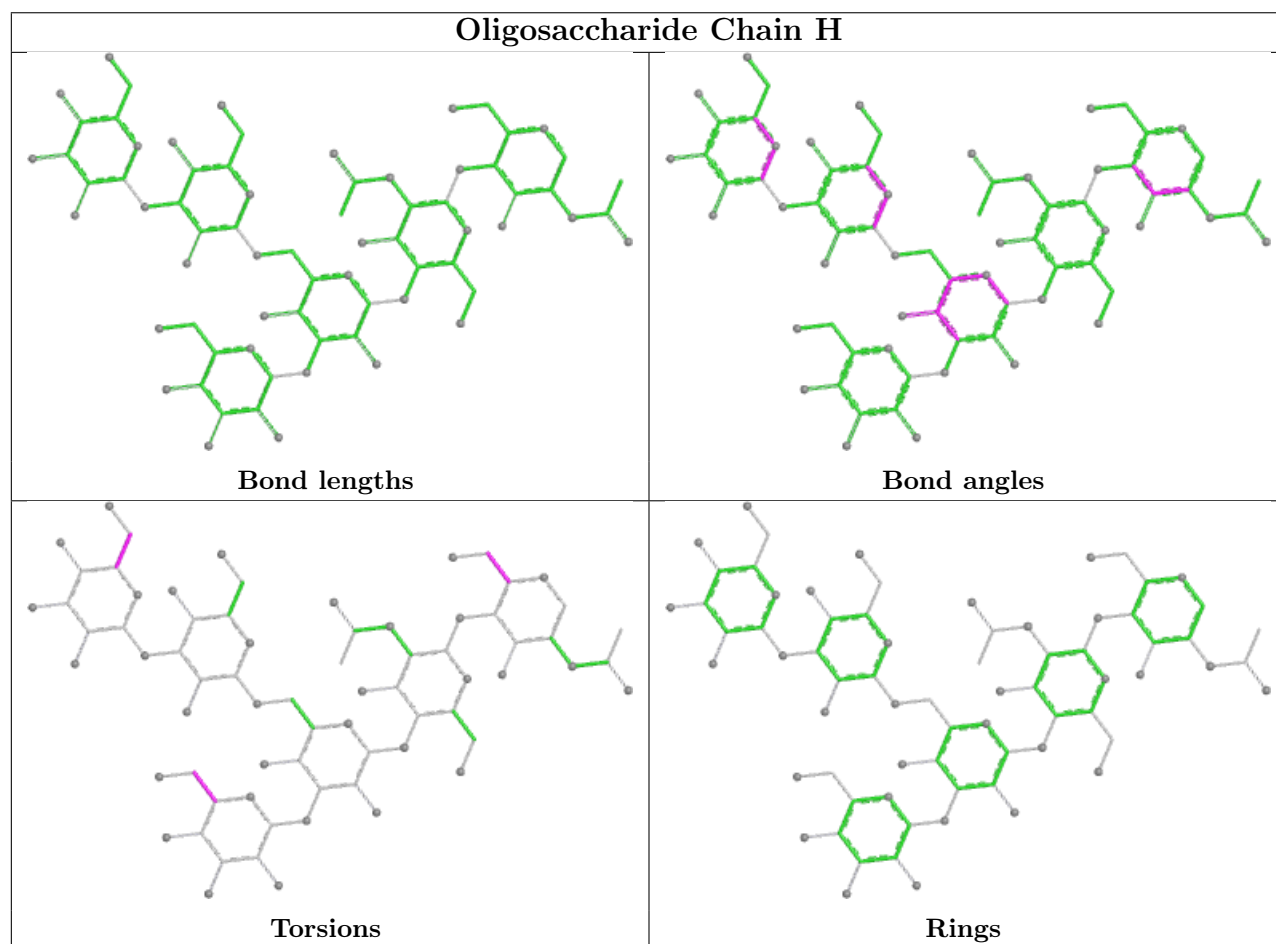
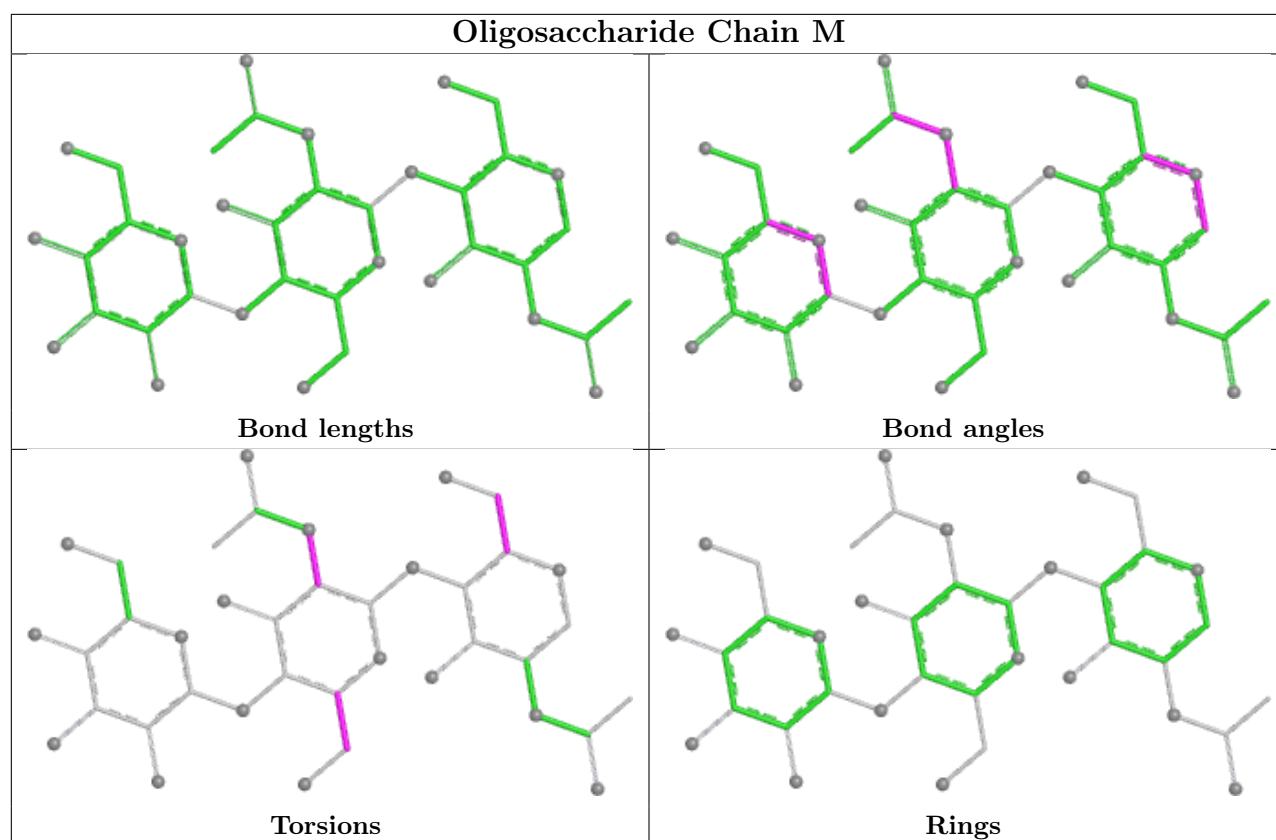
3 monomers are involved in 4 short contacts:

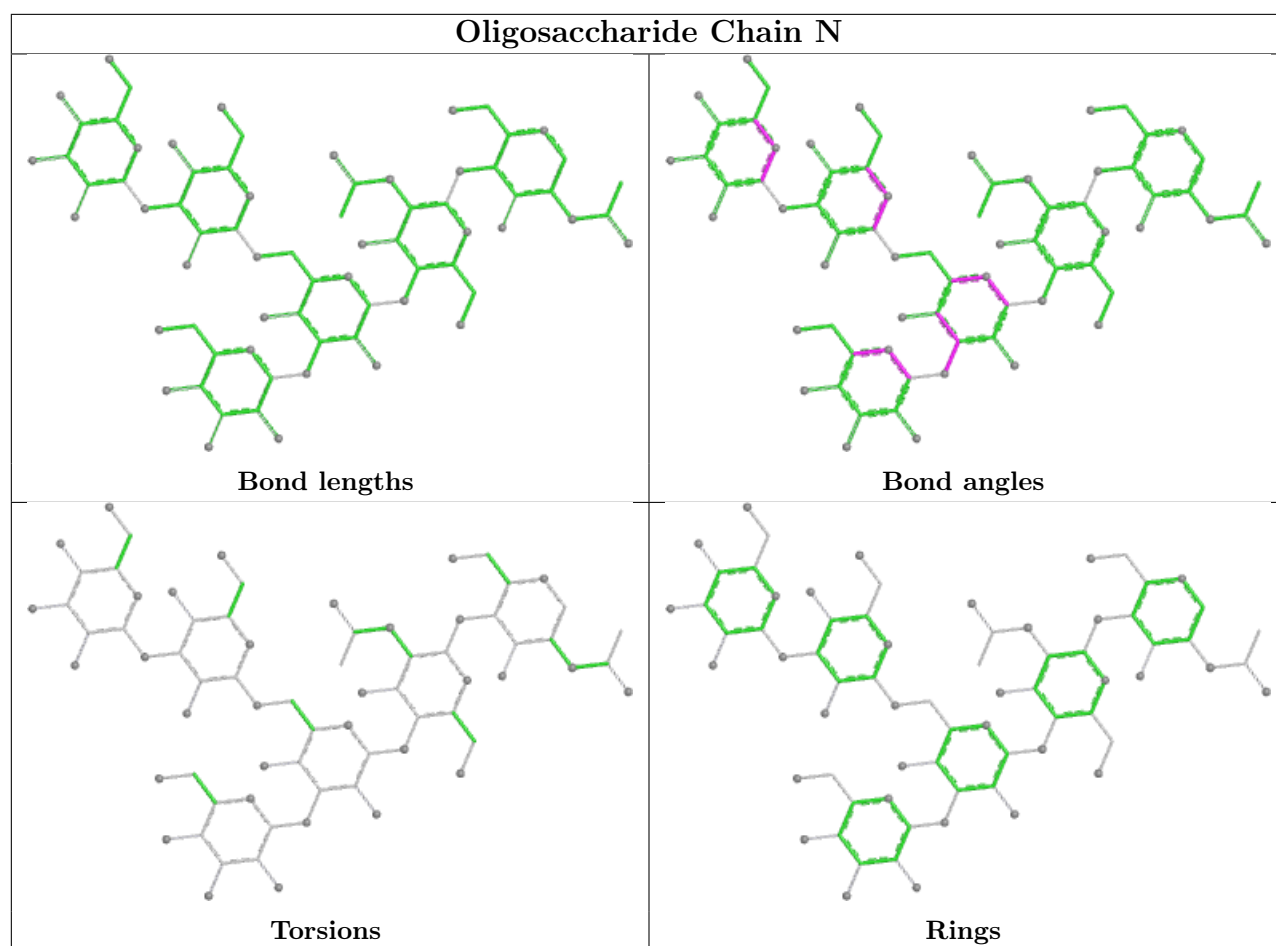
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	G	2	NAG	1	0
6	H	1	NAG	2	0
6	N	1	NAG	1	0

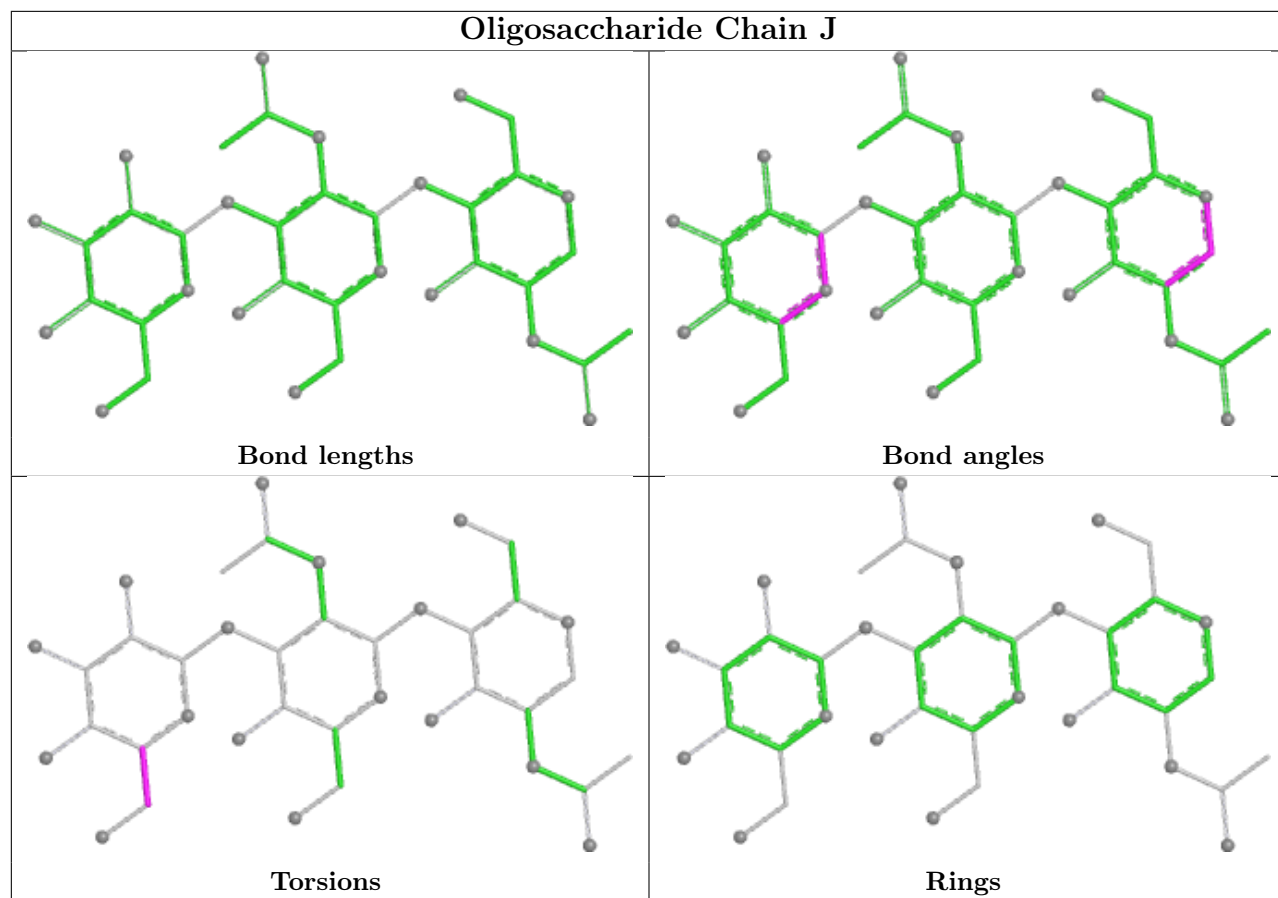
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

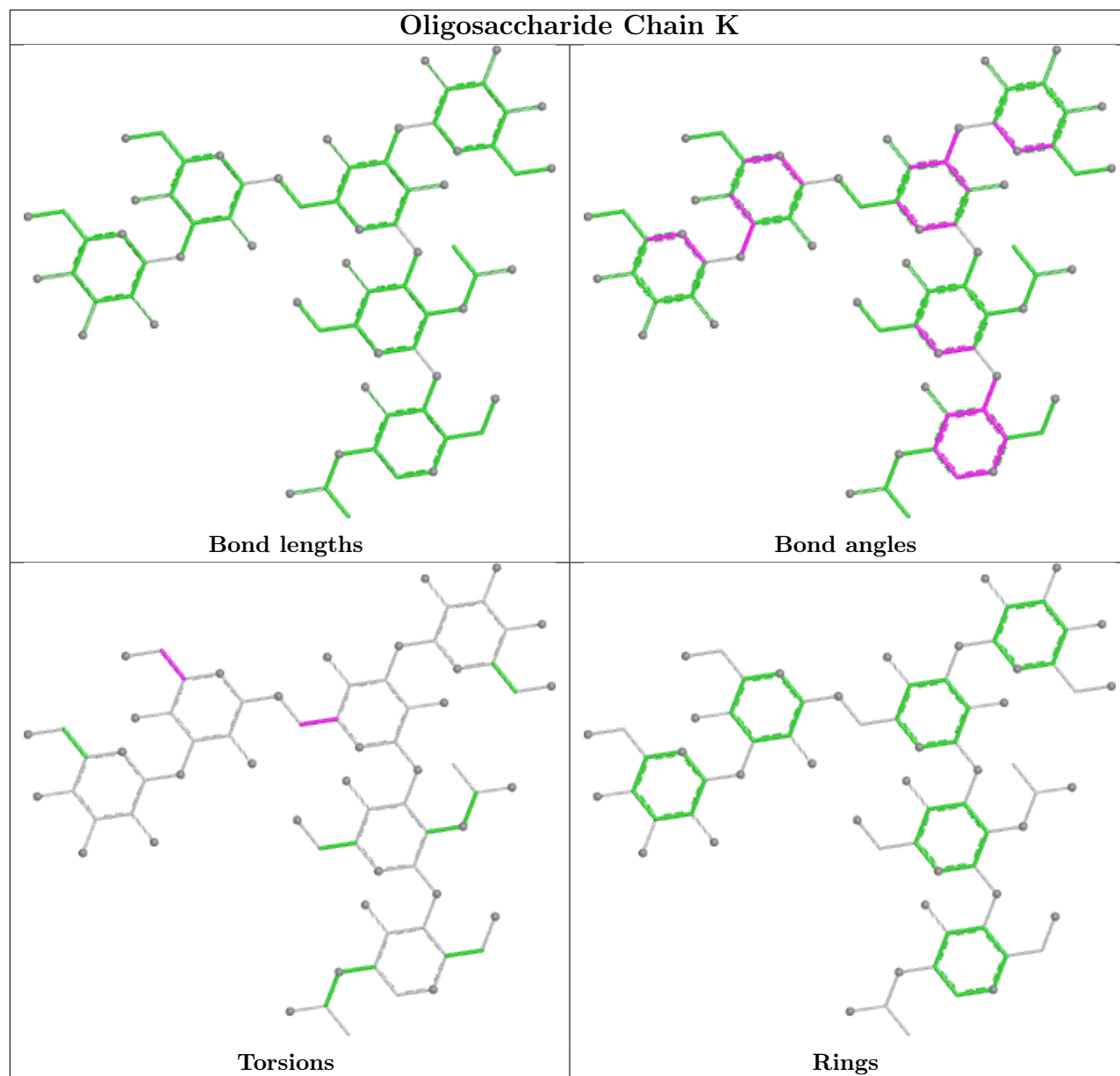


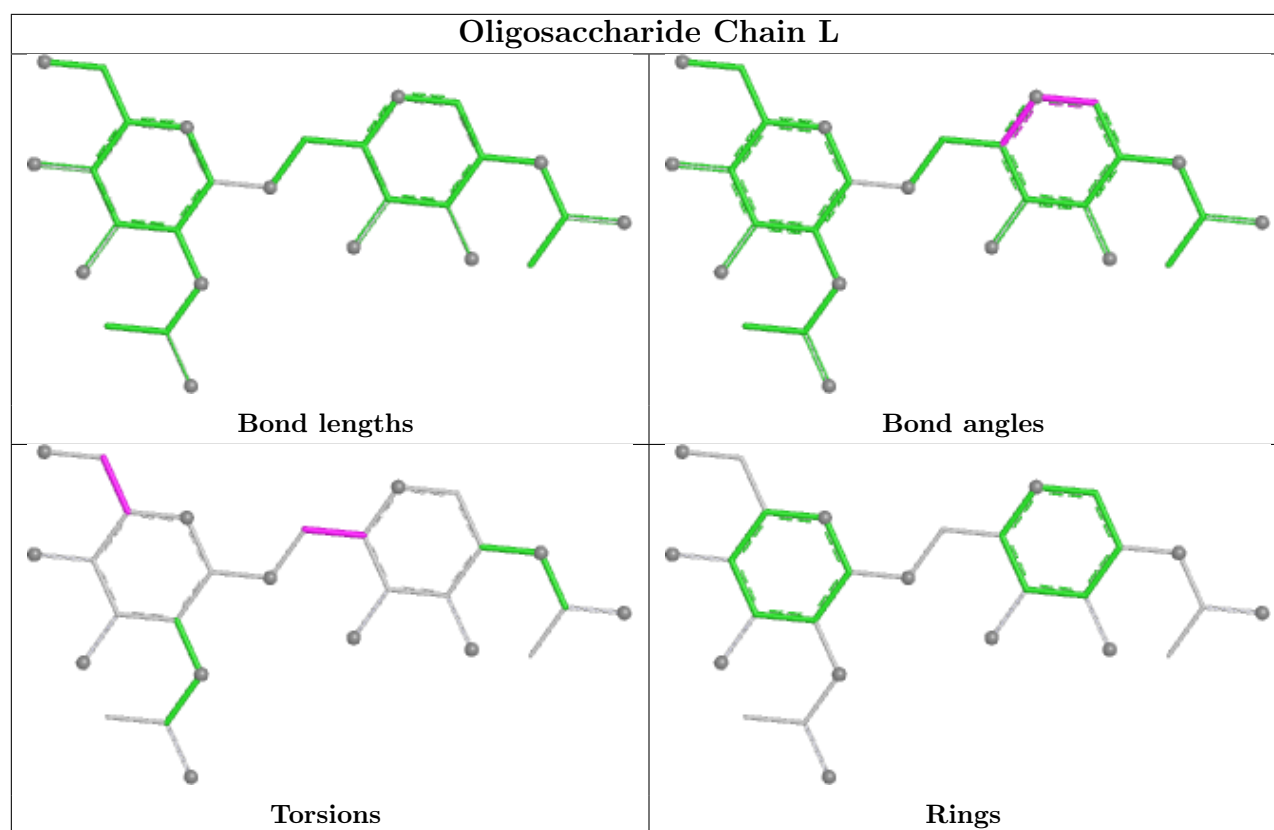












5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
10	ABU	B	601	-	6,6,6	1.05	0	6,6,6	1.17	1 (16%)
12	NAG	D	401	1	14,14,15	0.74	0	17,19,21	1.04	1 (5%)
11	EPE	C	601	-	15,15,15	0.81	1 (6%)	19,20,20	0.68	0
10	ABU	E	601	-	6,6,6	1.03	0	6,6,6	1.19	1 (16%)

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
10	ABU	B	601	-	-	2/4/4/4	-
12	NAG	D	401	1	-	1/6/23/26	0/1/1/1
11	EPE	C	601	-	-	1/9/19/19	0/1/1/1
10	ABU	E	601	-	-	0/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	C	601	EPE	C10-S	2.20	1.80	1.77

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	D	401	NAG	O5-C1-C2	-2.69	107.12	111.29
10	E	601	ABU	O-C-CG	-2.30	115.81	123.09
10	B	601	ABU	O-C-CG	-2.25	115.95	123.09

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	C	601	EPE	N4-C7-C8-O8
12	D	401	NAG	O5-C5-C6-O6
10	B	601	ABU	OXT-C-CG-CB
10	B	601	ABU	O-C-CG-CB

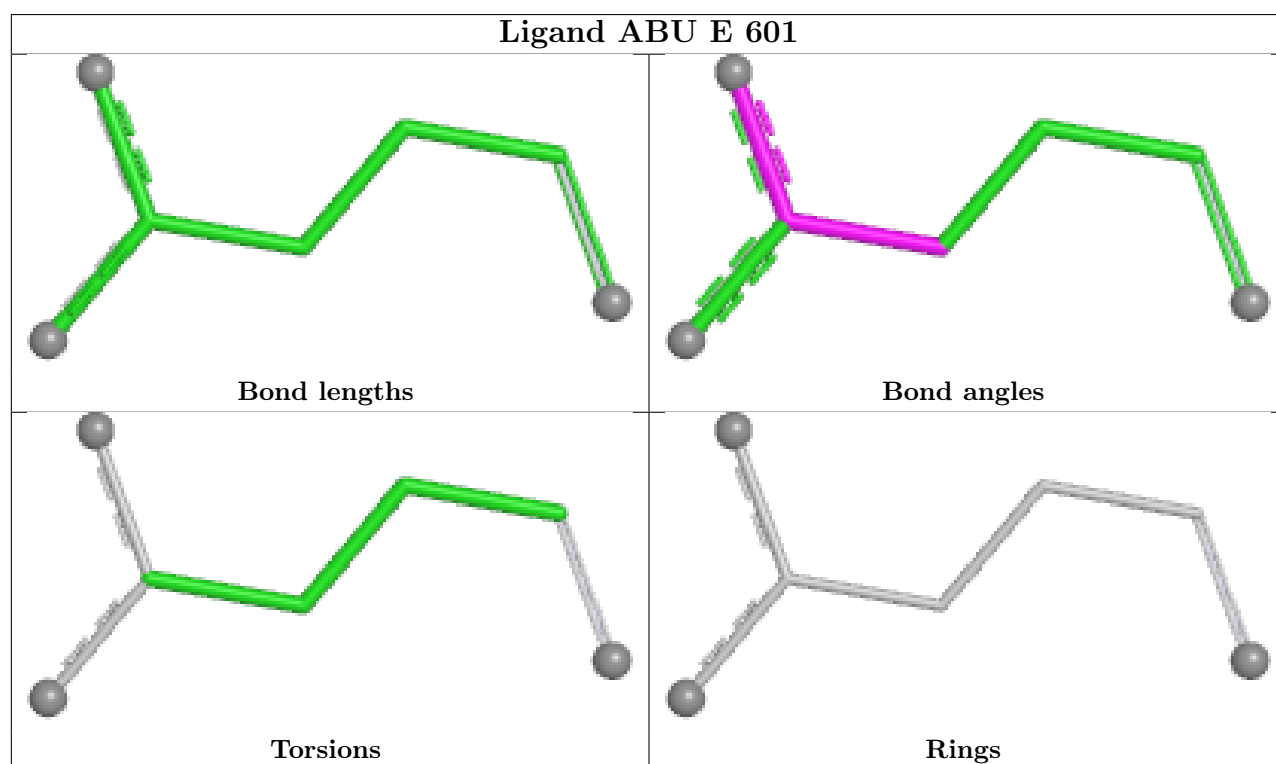
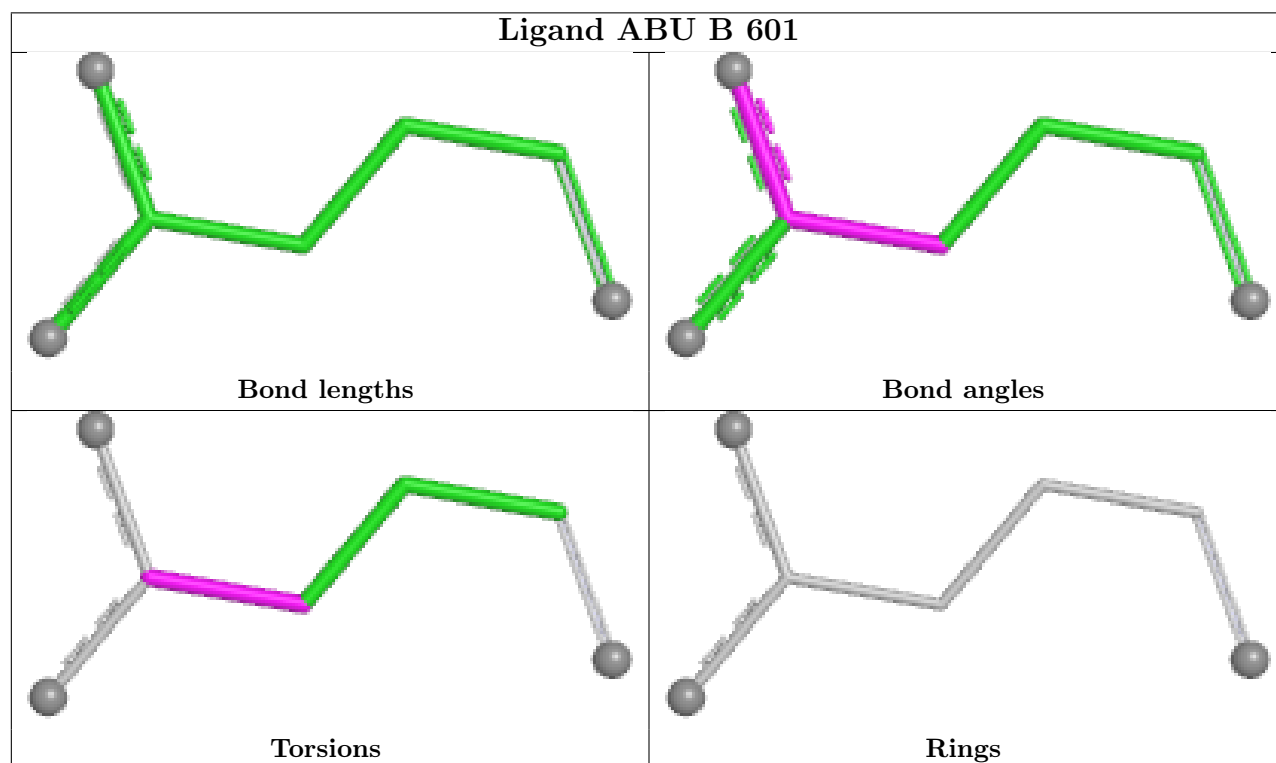
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	B	601	ABU	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

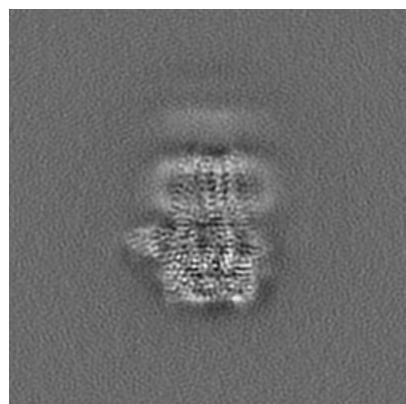
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-54151. These allow visual inspection of the internal detail of the map and identification of artifacts.

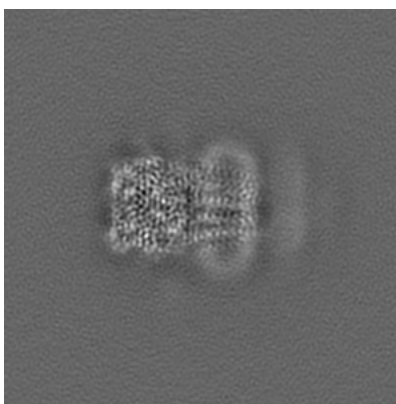
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

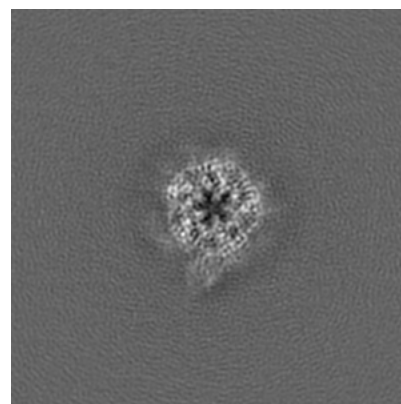
6.1.1 Primary map



X

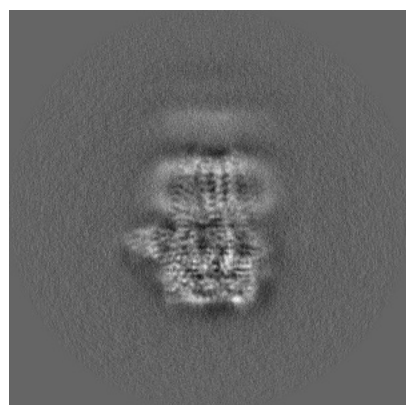


Y

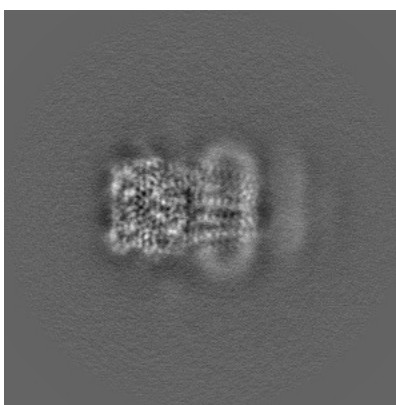


Z

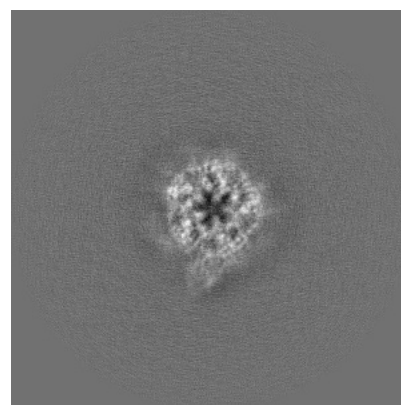
6.1.2 Raw map



X



Y

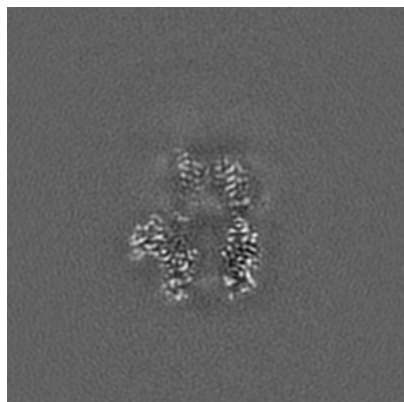


Z

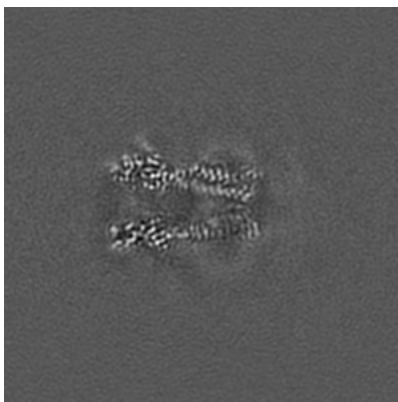
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

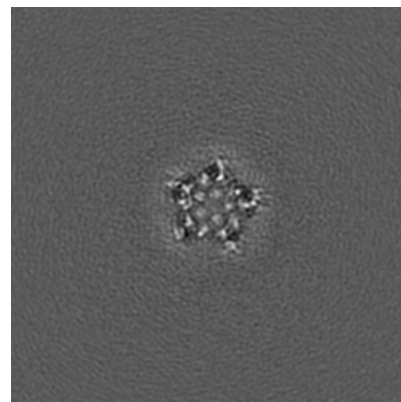
6.2.1 Primary map



X Index: 220

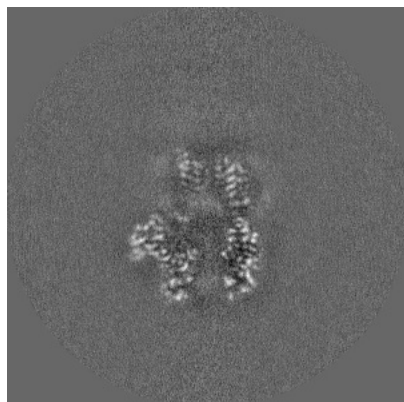


Y Index: 220

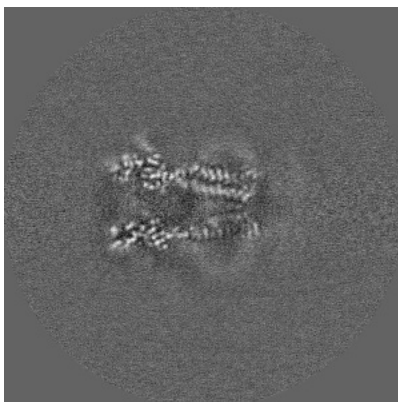


Z Index: 220

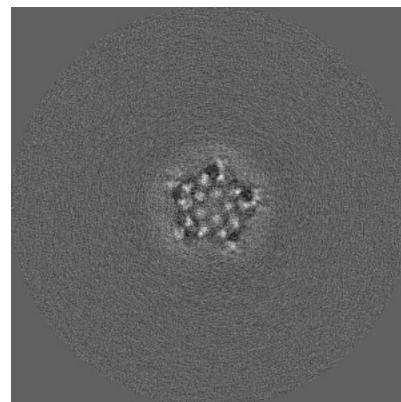
6.2.2 Raw map



X Index: 220



Y Index: 220

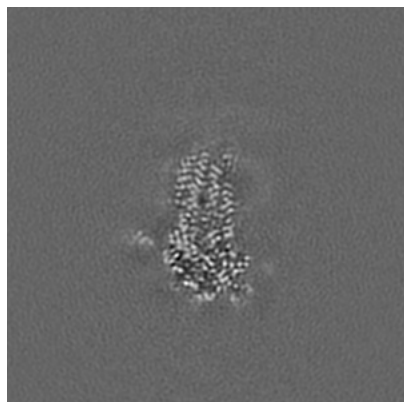


Z Index: 220

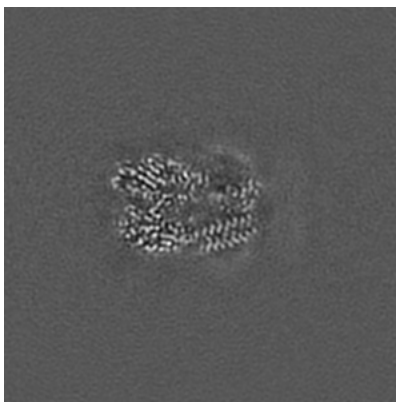
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

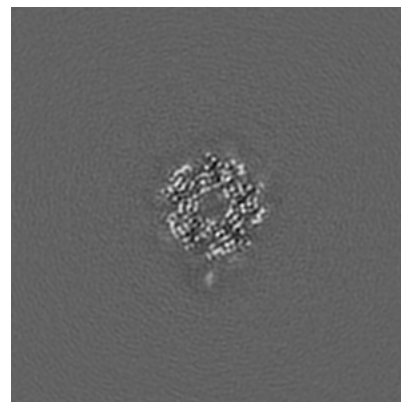
6.3.1 Primary map



X Index: 194

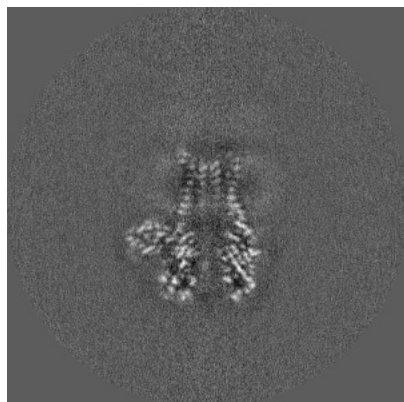


Y Index: 240

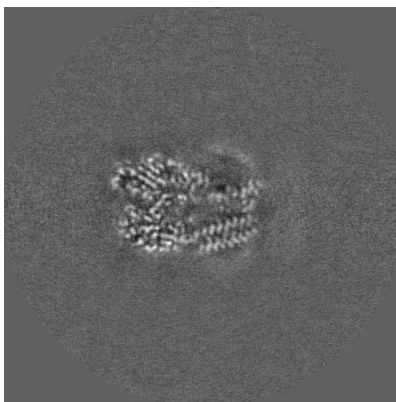


Z Index: 164

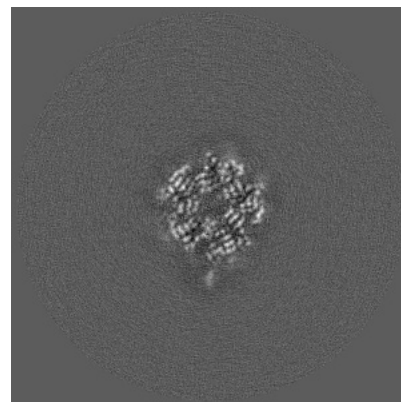
6.3.2 Raw map



X Index: 213



Y Index: 240

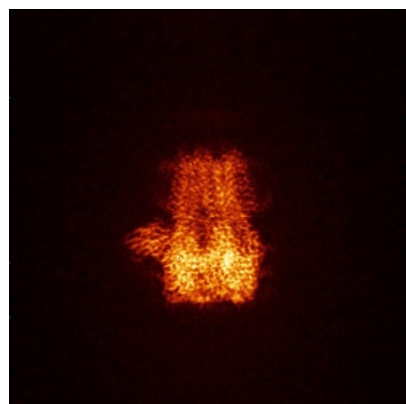


Z Index: 165

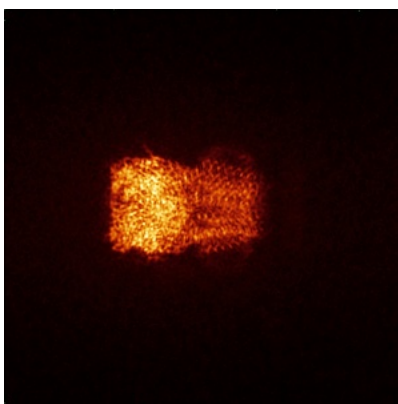
The images above show the largest variance slices of the map in three orthogonal directions.

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

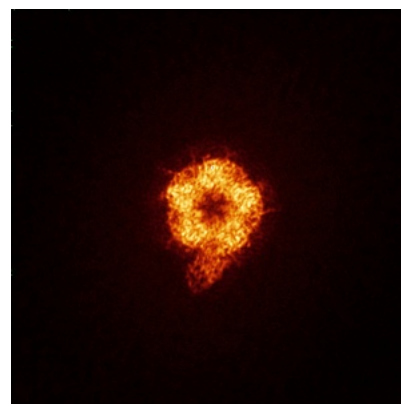
6.4.1 Primary map



X

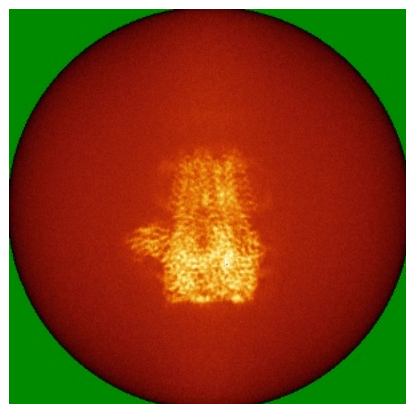


Y

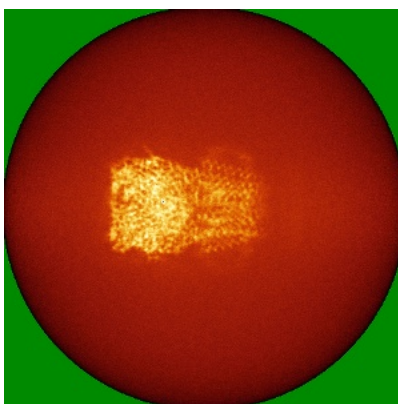


Z

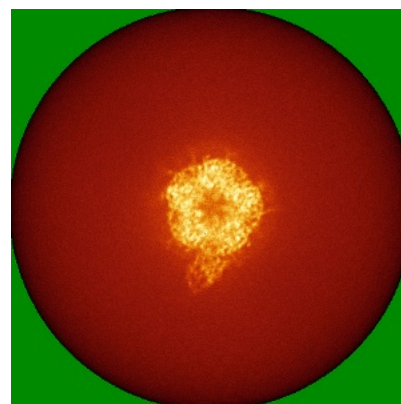
6.4.2 Raw map



X



Y

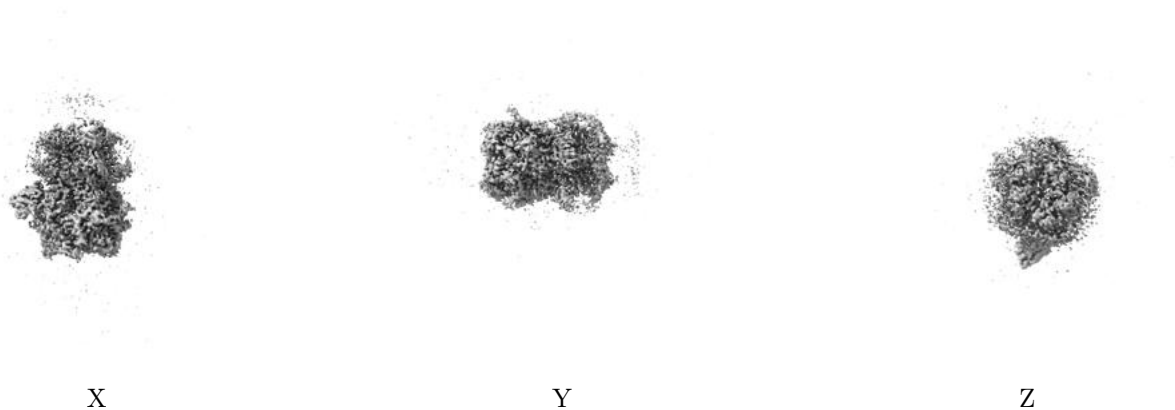


Z

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.

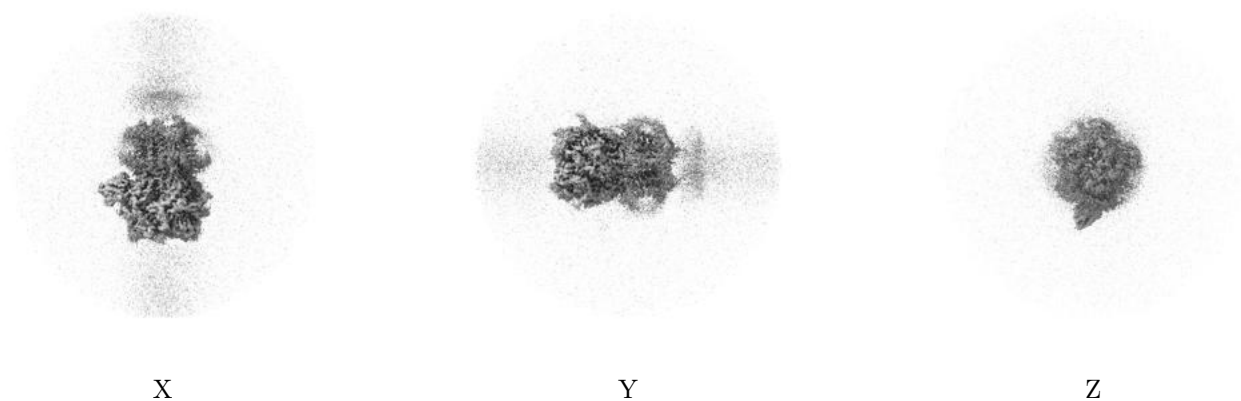
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 3.0. 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.

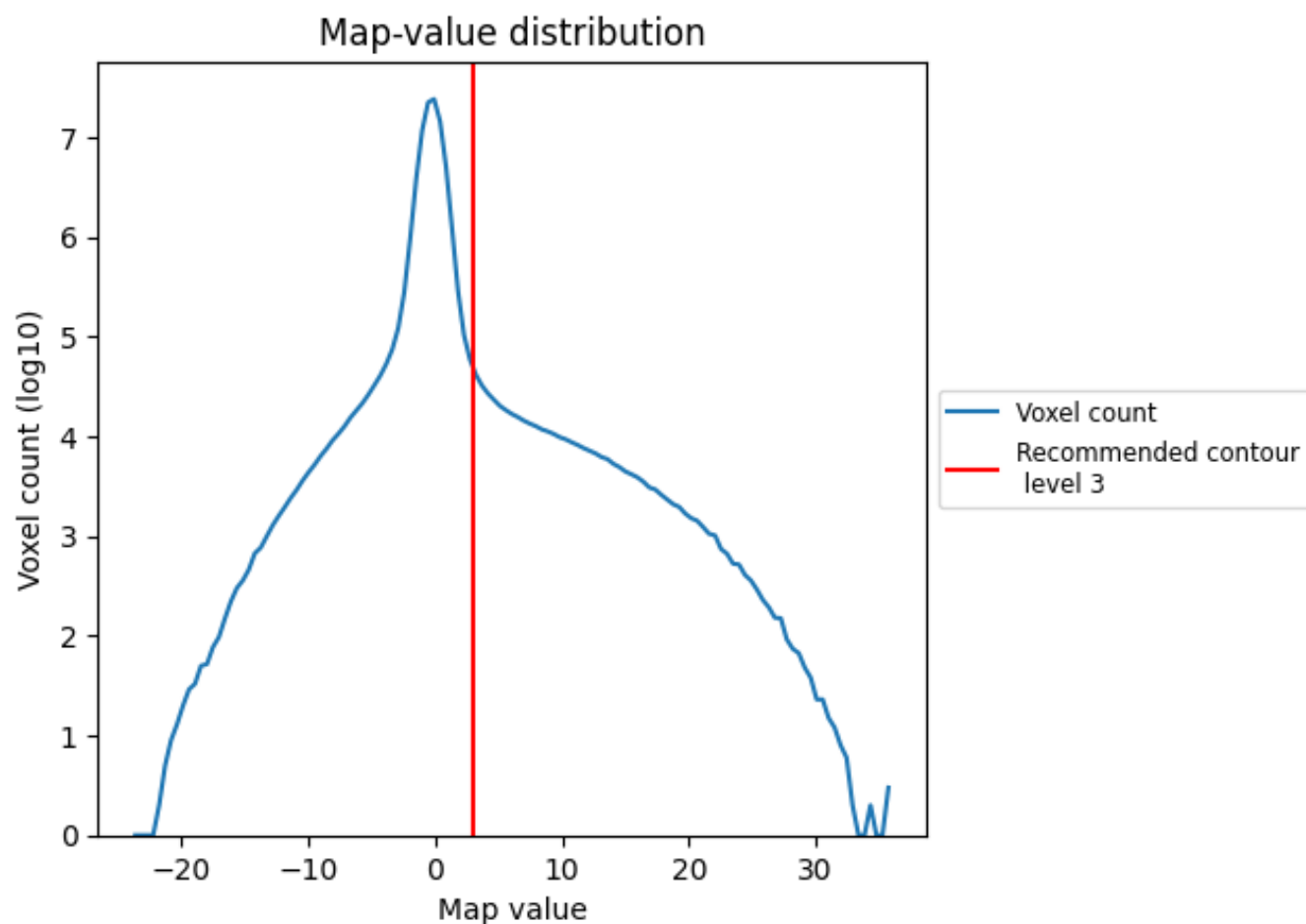
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

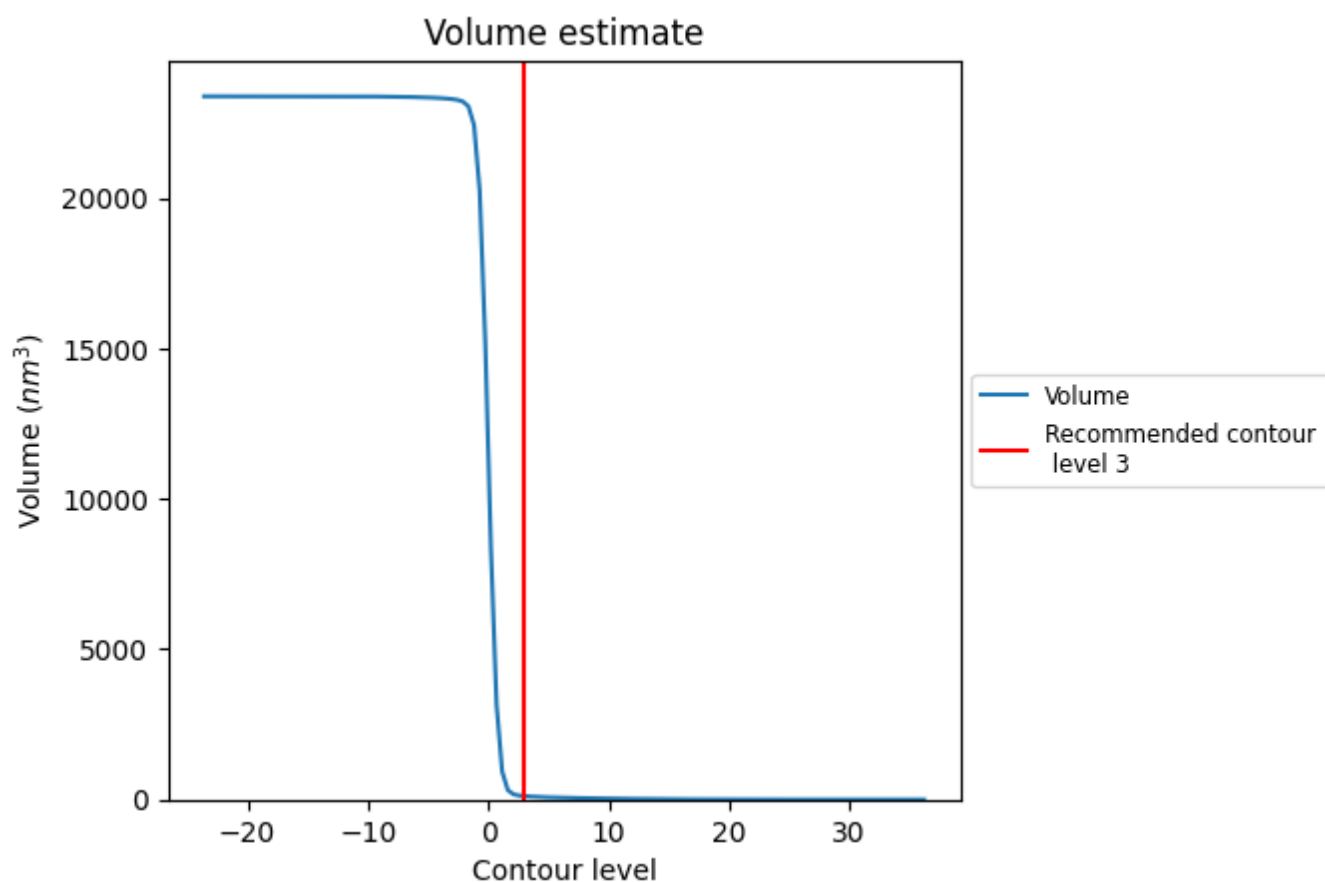
This section contains the results of statistical analysis of the map.

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.

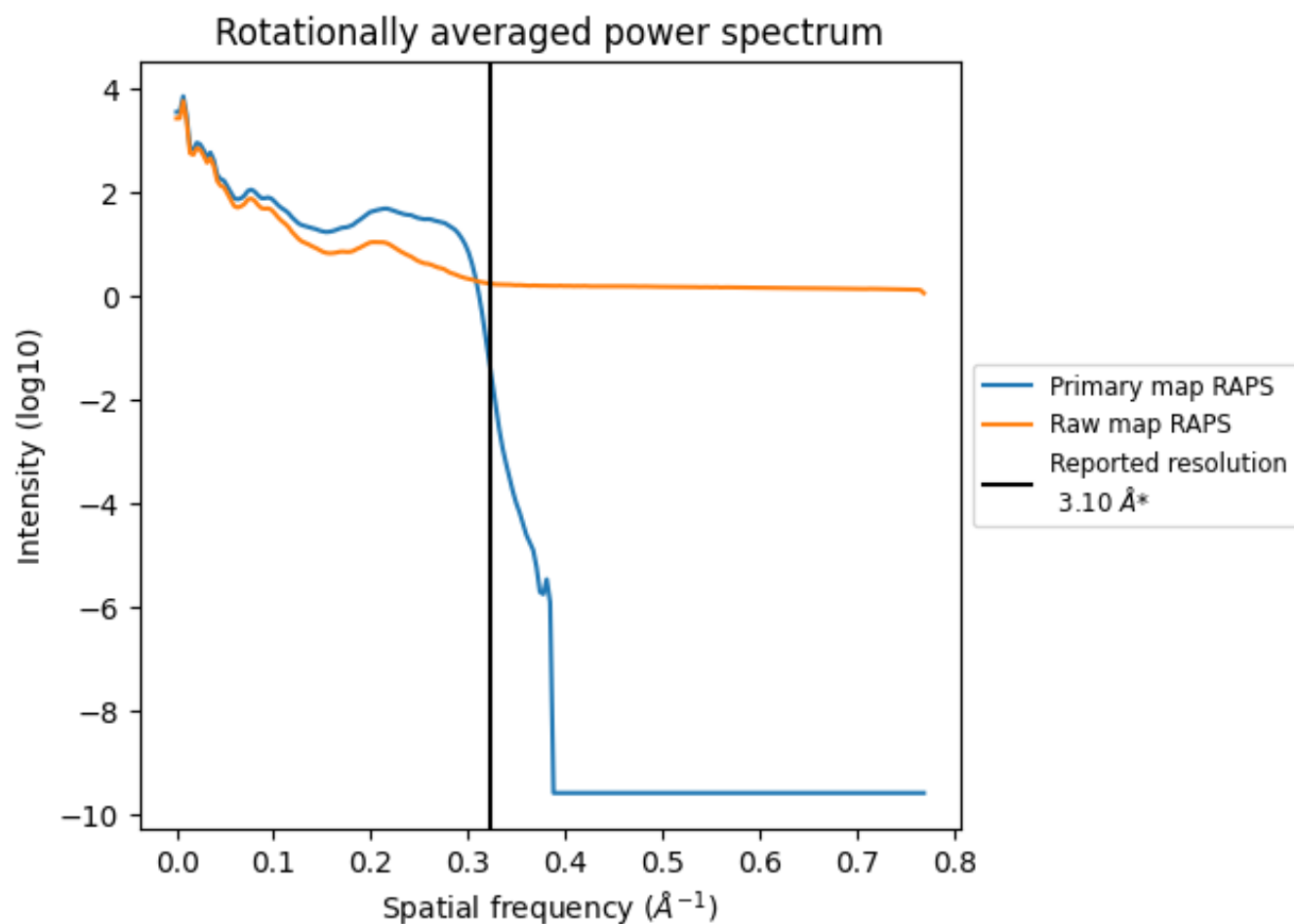
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 115 nm³; this corresponds to an approximate mass of 104 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 ⓘ

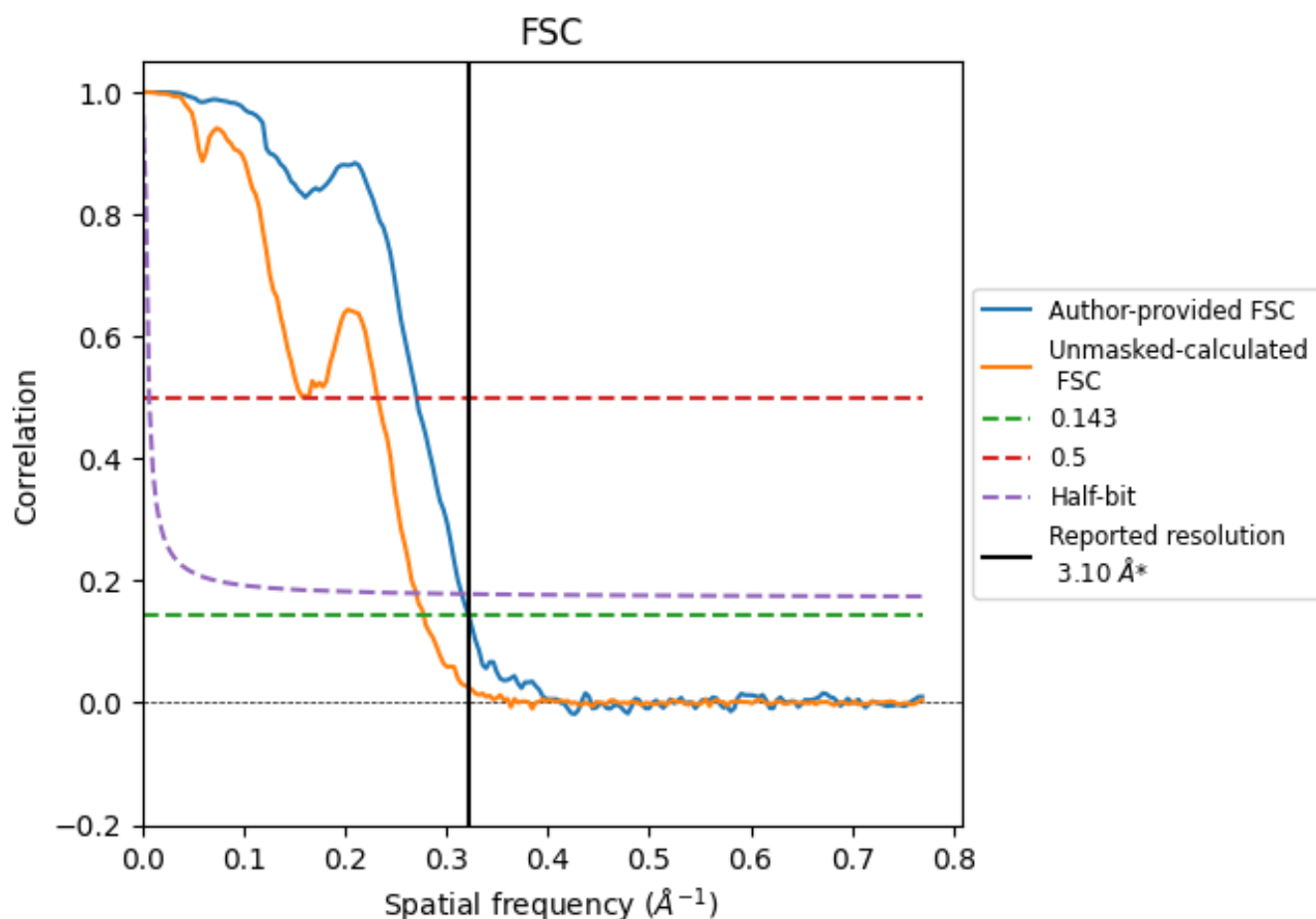


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.323 \AA^{-1}

8.2 Resolution estimates [i](#)

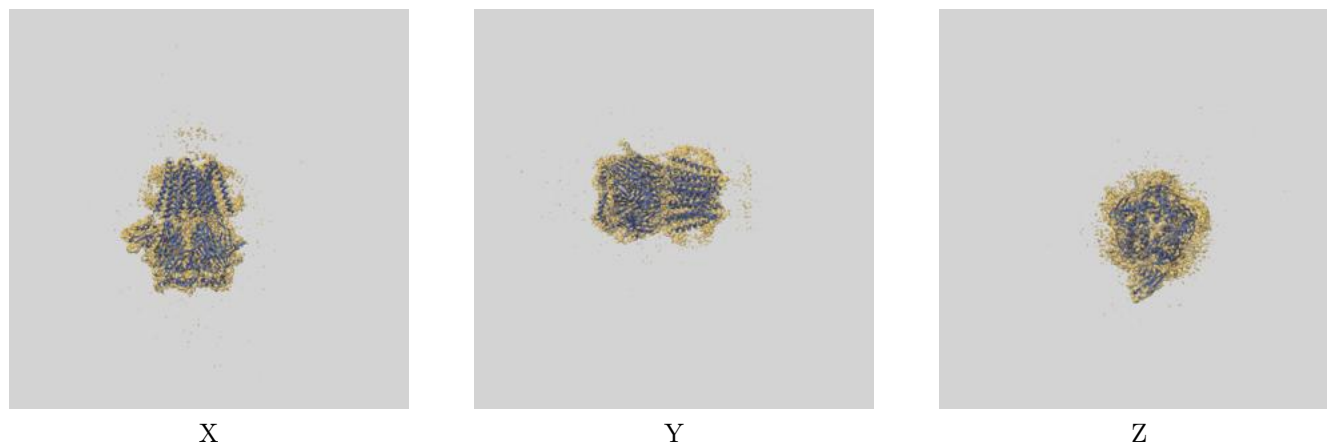
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.11	3.70	3.17
Unmasked-calculated*	3.60	4.31	3.70

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.60 differs from the reported value 3.1 by more than 10 %

9 Map-model fit [i](#)

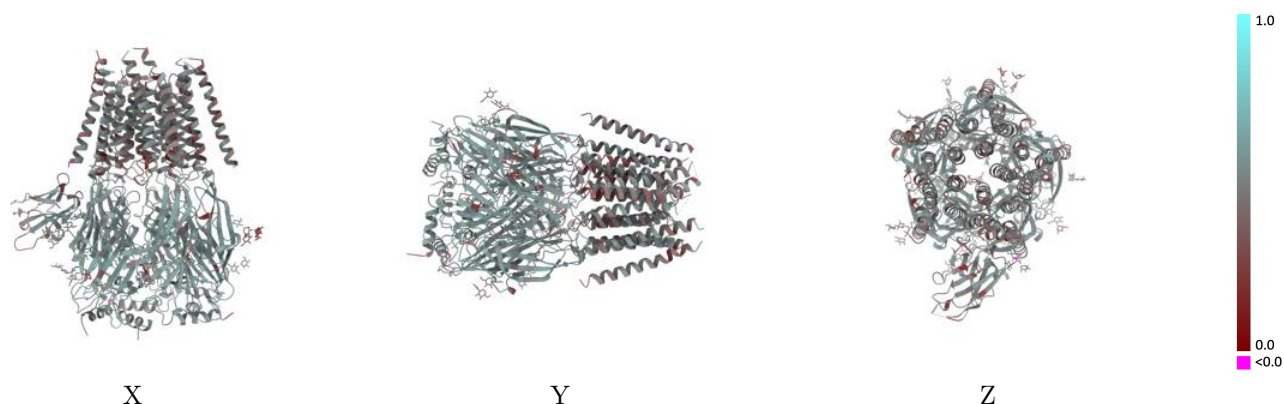
This section contains information regarding the fit between EMDB map EMD-54151 and PDB model 9RPB. Per-residue inclusion information can be found in section 3 on page 25.

9.1 Map-model overlay [i](#)



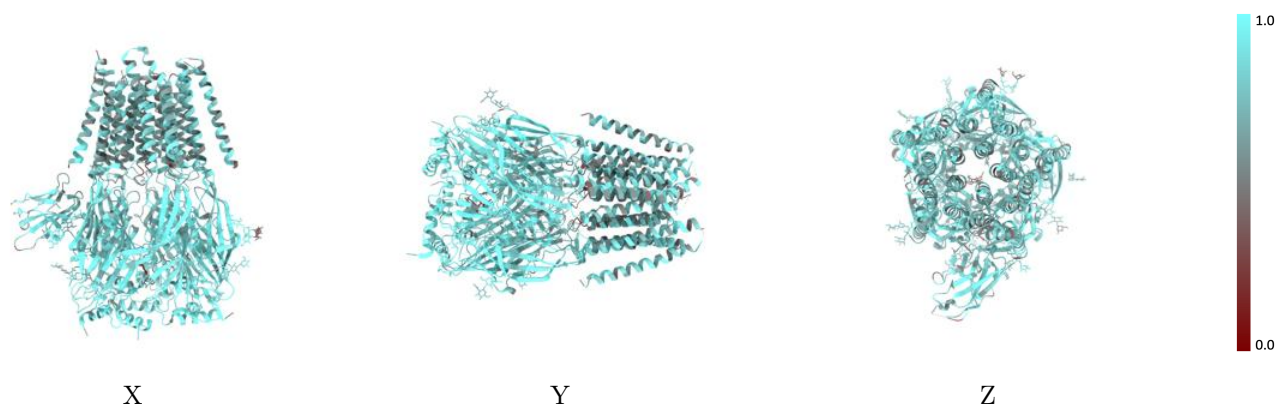
The images above show the 3D surface view of the map at the recommended contour level 3.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



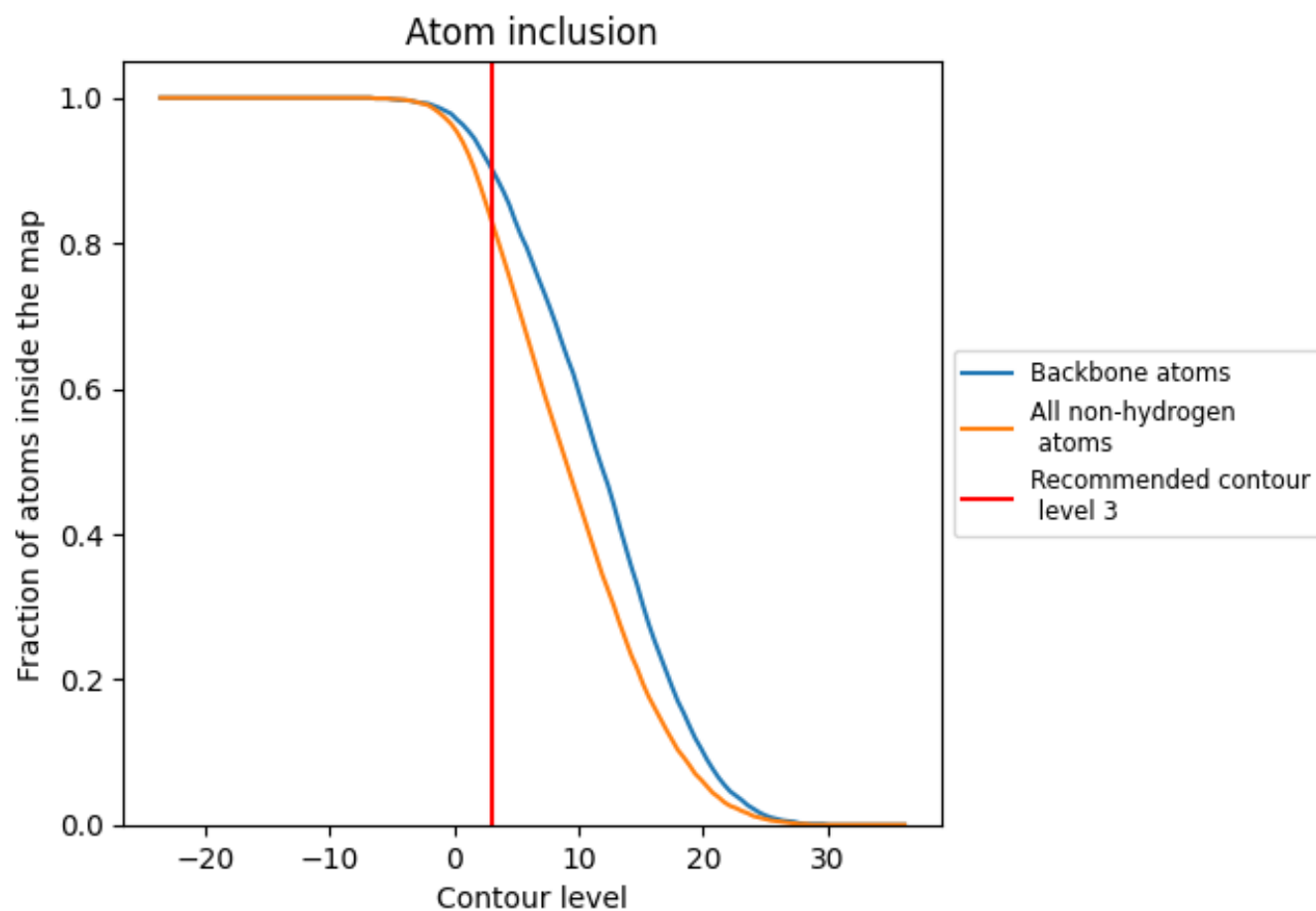
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (3).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 83% 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 (3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8310	<div></div> 0.4970
A	<div></div> 0.8360	<div></div> 0.5030
B	<div></div> 0.8290	<div></div> 0.5020
C	<div></div> 0.8290	<div></div> 0.5070
D	<div></div> 0.8340	<div></div> 0.4980
E	<div></div> 0.8510	<div></div> 0.5040
F	<div></div> 0.7860	<div></div> 0.4290
G	<div></div> 0.2820	<div></div> 0.2360
H	<div></div> 0.7920	<div></div> 0.4580
I	<div></div> 0.7690	<div></div> 0.4250
J	<div></div> 0.6920	<div></div> 0.3970
K	<div></div> 0.9030	<div></div> 0.4740
L	<div></div> 0.6790	<div></div> 0.4050
M	<div></div> 0.8460	<div></div> 0.4500
N	<div></div> 0.7640	<div></div> 0.3790
P	<div></div> 0.7990	<div></div> 0.4470

