



wwPDB EM Validation Summary Report ⓘ

Oct 28, 2024 – 11:35 am GMT

PDB ID : 7QNE
EMDB ID : EMD-14076
Title : Cryo-EM structure of human full-length synaptic alpha1beta3gamma2 GABA(A)R in complex with Ro15-4513 and megabody Mb38
Authors : Sente, A.; Desai, R.; Naydenova, K.; Malinauskas, T.; Jounaidi, Y.; Miehl, J.; Zhou, X.; Masiulis, S.; Hardwick, S.W.; Chirgadze, D.Y.; Miller, K.W.; Aricescu, A.R.
Deposited on : 2021-12-20
Resolution : 2.70 Å(reported)

This is a wwPDB EM Validation Summary 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.dev113
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

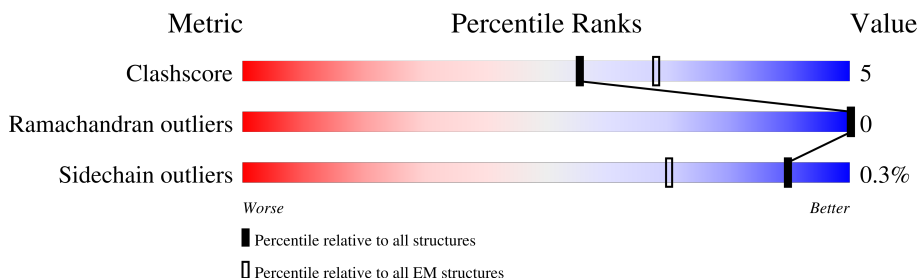
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.70 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	456	<div> <div>15%</div> <div>64%</div> <div>12%</div> <div>24%</div> </div>
1	D	456	<div> <div>22%</div> <div>63%</div> <div>12%</div> <div>25%</div> </div>
2	B	473	<div> <div>13%</div> <div>61%</div> <div>9%</div> <div>29%</div> </div>
2	E	473	<div> <div>16%</div> <div>60%</div> <div>10%</div> <div>29%</div> </div>
3	C	495	<div> <div>16%</div> <div>60%</div> <div>7%</div> <div>34%</div> </div>
4	G	123	<div> <div>49%</div> <div>85%</div> <div>15%</div> </div>
5	a	9	<div> <div>100%</div> </div>
6	b	2	<div> <div>100%</div> </div>

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Mol	Chain	Length	Quality of chain
6	e	2	100%
			100%
7	F	6	83%
			100%
7	H	6	67%
			100%
8	d	5	100%
			100%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 15302 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 GABA(A) receptor subunit alpha-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	345	Total	C	N	O	S	1	0
			2801	1812	473	500	16		
1	D	343	Total	C	N	O	S	2	0
			2795	1810	473	496	16		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	334	Total	C	N	O	S	1	0
			2751	1802	450	484	15		
2	E	334	Total	C	N	O	S	1	0
			2745	1796	449	484	16		

- Molecule 3 is a protein called GABA(A) receptor subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	329	Total	C	N	O	S	0	0
			2712	1770	447	480	15		

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	PRO	deletion	UNP A0A286YFI6
C	?	-	ILE	deletion	UNP A0A286YFI6
C	?	-	HIS	deletion	UNP A0A286YFI6
C	?	-	HIS	deletion	UNP A0A286YFI6
C	?	-	HIS	deletion	UNP A0A286YFI6
C	?	-	ILE	deletion	UNP A0A286YFI6
C	?	-	GLY	deletion	UNP A0A286YFI6
C	?	-	ASP	deletion	UNP A0A286YFI6
C	?	-	ILE	deletion	UNP A0A286YFI6
C	?	-	VAL	deletion	UNP A0A286YFI6
C	?	-	GLU	deletion	UNP A0A286YFI6

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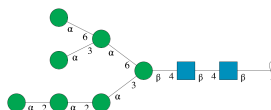
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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	LYS	deletion	UNP A0A286YFI6
C	?	-	GLN	deletion	UNP A0A286YFI6
C	?	-	PRO	deletion	UNP A0A286YFI6
C	?	-	ARG	deletion	UNP A0A286YFI6
C	?	-	ILE	deletion	UNP A0A286YFI6
C	?	-	SER	deletion	UNP A0A286YFI6
C	?	-	ARG	deletion	UNP A0A286YFI6
C	?	-	GLU	deletion	UNP A0A286YFI6
C	?	-	GLN	deletion	UNP A0A286YFI6
C	?	-	LEU	deletion	UNP A0A286YFI6
C	?	-	THR	deletion	UNP A0A286YFI6
C	?	-	ASP	deletion	UNP A0A286YFI6
C	338	LEU	SER	conflict	UNP A0A286YFI6
C	339	LEU	PRO	conflict	UNP A0A286YFI6
C	340	ARG	SER	conflict	UNP A0A286YFI6
C	341	MET	LEU	conflict	UNP A0A286YFI6
C	342	PHE	PRO	conflict	UNP A0A286YFI6
C	344	PHE	SER	conflict	UNP A0A286YFI6
C	345	LYS	GLN	conflict	UNP A0A286YFI6
C	437	GLY	-	expression tag	UNP A0A286YFI6
C	438	GLY	-	expression tag	UNP A0A286YFI6
C	439	SER	-	expression tag	UNP A0A286YFI6
C	440	GLY	-	expression tag	UNP A0A286YFI6
C	441	GLY	-	expression tag	UNP A0A286YFI6
C	442	SER	-	expression tag	UNP A0A286YFI6
C	443	GLY	-	expression tag	UNP A0A286YFI6
C	444	GLY	-	expression tag	UNP A0A286YFI6
C	445	SER	-	expression tag	UNP A0A286YFI6
C	446	GLY	-	expression tag	UNP A0A286YFI6
C	447	LYS	-	expression tag	UNP A0A286YFI6
C	448	THR	-	expression tag	UNP A0A286YFI6
C	449	GLU	-	expression tag	UNP A0A286YFI6
C	450	THR	-	expression tag	UNP A0A286YFI6
C	451	SER	-	expression tag	UNP A0A286YFI6
C	452	GLN	-	expression tag	UNP A0A286YFI6
C	453	VAL	-	expression tag	UNP A0A286YFI6
C	454	ALA	-	expression tag	UNP A0A286YFI6
C	455	PRO	-	expression tag	UNP A0A286YFI6
C	456	ALA	-	expression tag	UNP A0A286YFI6

- Molecule 4 is a protein called Megabody Mb38.

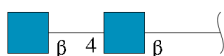
Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	123	Total	C	N	O	S	0	0
			949	593	171	181	4		

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



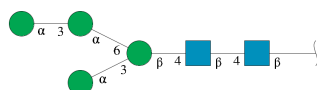
Mol	Chain	Residues	Atoms				AltConf	Trace
5	a	9	Total	C	N	O	0	0
			105	58	2	45		

- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



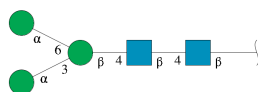
Mol	Chain	Residues	Atoms				AltConf	Trace
6	b	2	Total	C	N	O	0	0
			28	16	2	10		
6	e	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 7 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.



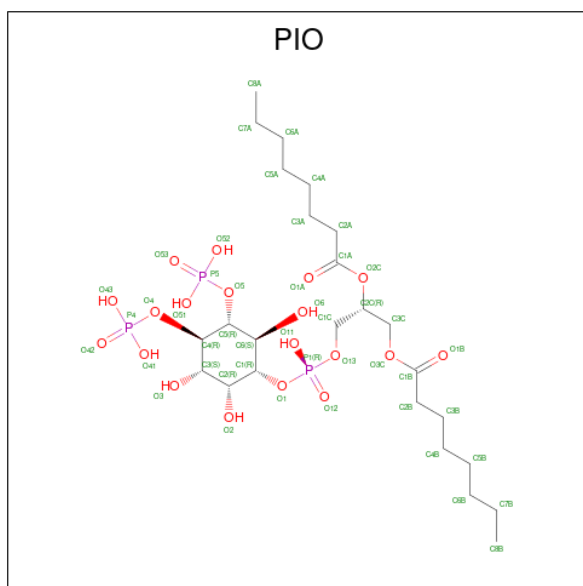
Mol	Chain	Residues	Atoms				AltConf	Trace
7	F	6	Total	C	N	O	0	0
			72	40	2	30		
7	H	6	Total	C	N	O	0	0
			72	40	2	30		

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



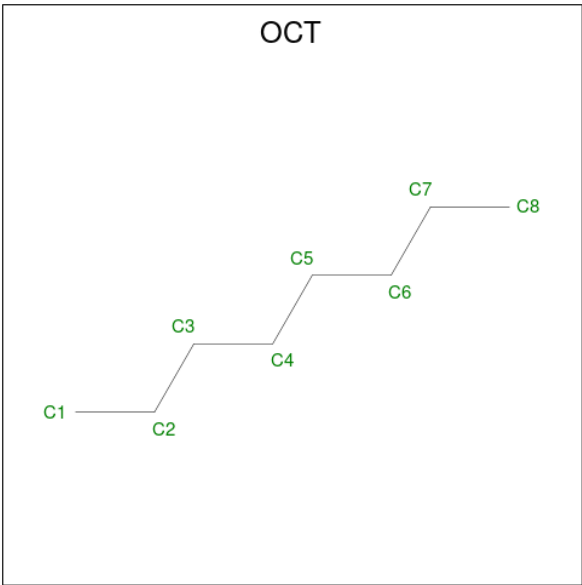
Mol	Chain	Residues	Atoms				AltConf	Trace
8	d	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 9 is [(2R)-2-octanoyloxy-3-[oxidanyl-[(1R,2R,3S,4R,5R,6S)-2,3,6-tris(oxidanyl)-4,5-diphosphonooxy-cyclohexyl]oxy-phosphoryl]oxy-propyl] octanoate (three-letter code: PIO) (formula: C₂₅H₄₉O₁₉P₃).



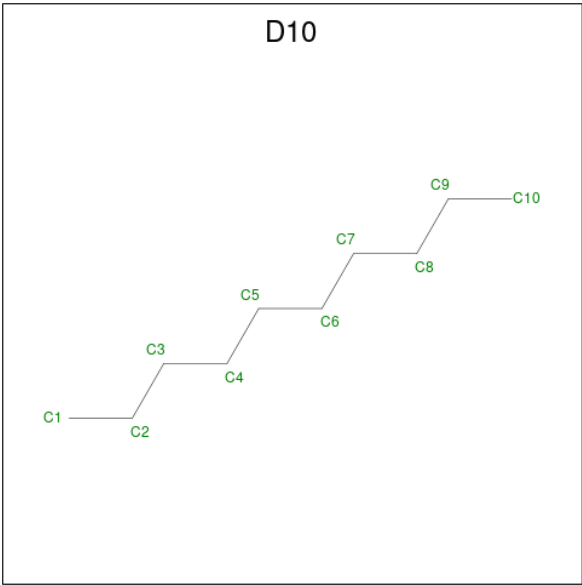
Mol	Chain	Residues	Atoms				AltConf
9	A	1	Total	C	O	P	0
			47	25	19	3	
9	D	1	Total	C	O	P	0
			47	25	19	3	

- Molecule 10 is N-OCTANE (three-letter code: OCT) (formula: C₈H₁₈).



Mol	Chain	Residues	Atoms		AltConf
10	A	1	Total	C	0

- Molecule 11 is DECANE (three-letter code: D10) (formula: C₁₀H₂₂).



Mol	Chain	Residues	Atoms		AltConf
11	A	1	Total 10	C 10	0
11	B	1	Total 10	C 10	0
11	C	1	Total 10	C 10	0

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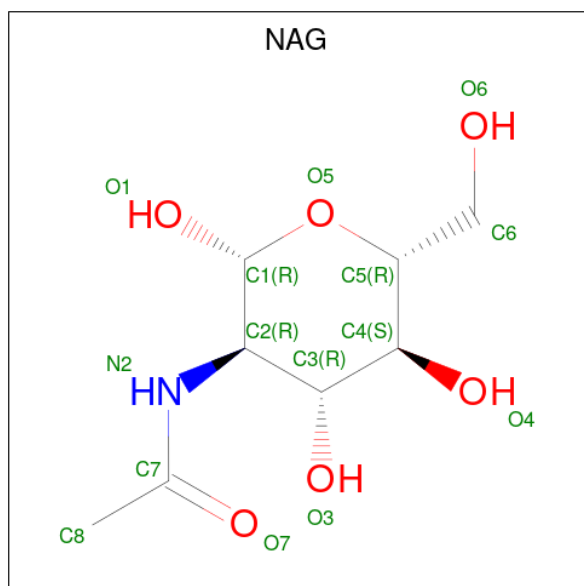
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Mol	Chain	Residues	Atoms	AltConf
11	E	1	Total C 10 10	0

- Molecule 12 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

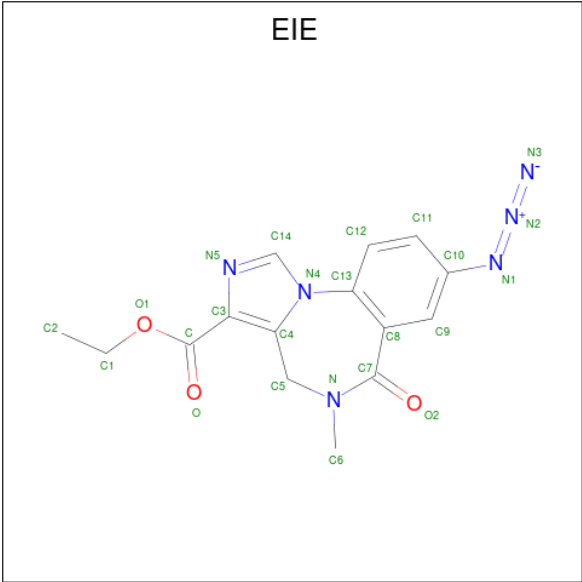
Mol	Chain	Residues	Atoms	AltConf
12	A	1	Total Cl 1 1	0
12	C	1	Total Cl 1 1	0
12	D	1	Total Cl 1 1	0

- Molecule 13 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 14 is ethyl 8-[(azanylidene- λ^4 -azanylidene)amino]-5-methyl-6-oxidanylidene-4 {H}-imidazo[1,5-a][1,4]benzodiazepine-3-carboxylate (three-letter code: EIE) (formula: C₁₅H₁₄N₆O₃) (labeled as "Ligand of Interest" by depositor).

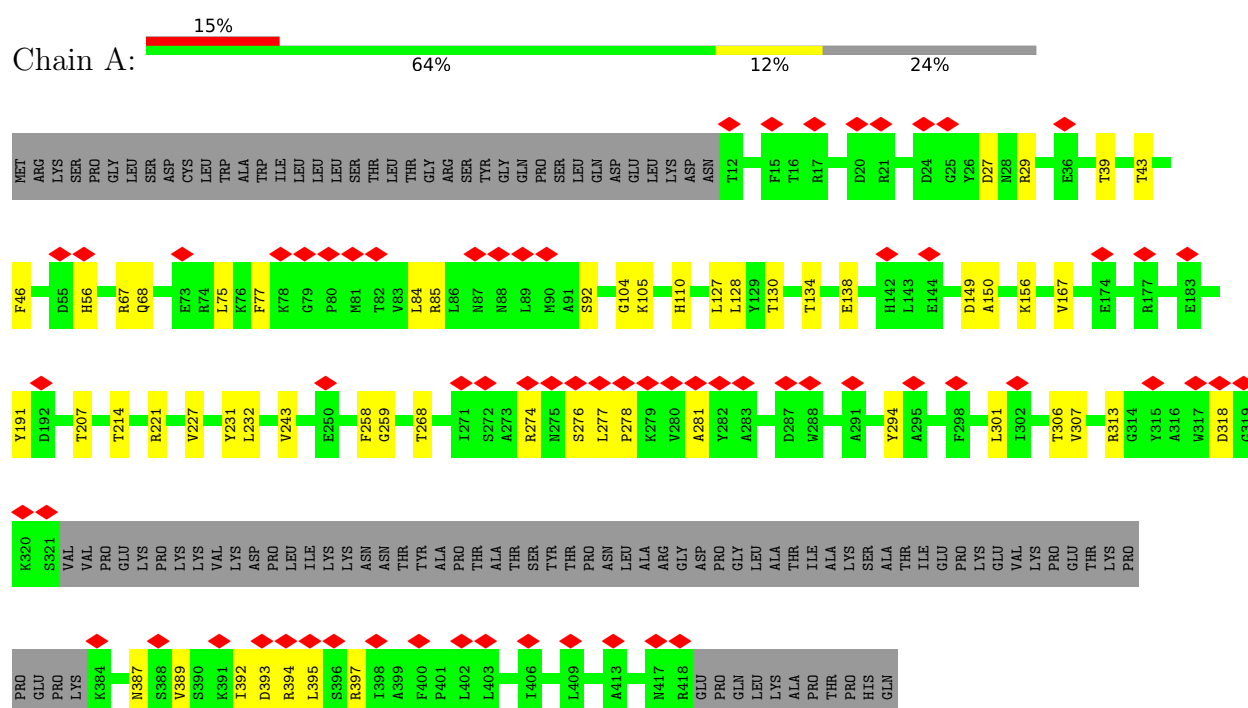


Mol	Chain	Residues	Atoms				AltConf
14	D	1	Total	C	N	O	0
			24	15	6	3	

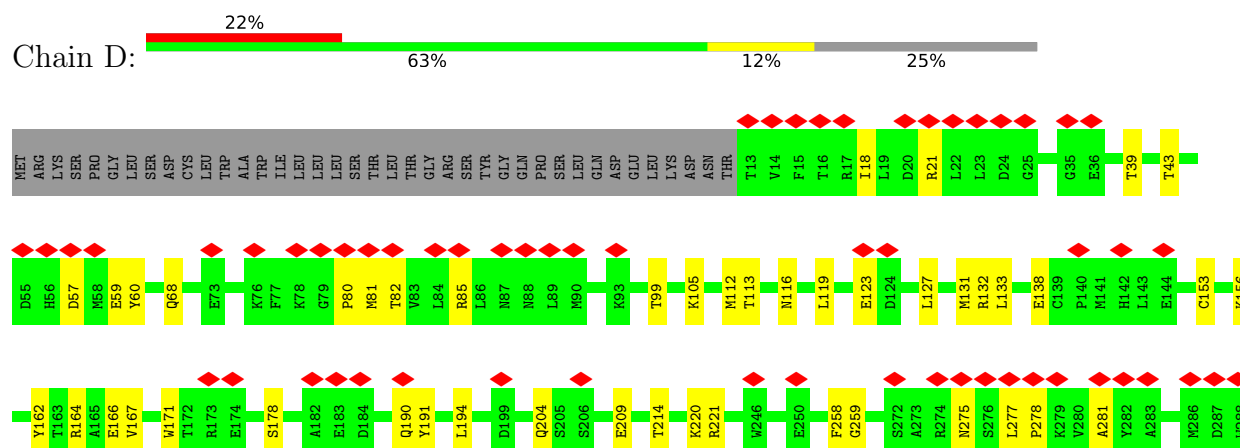
3 Residue-property plots

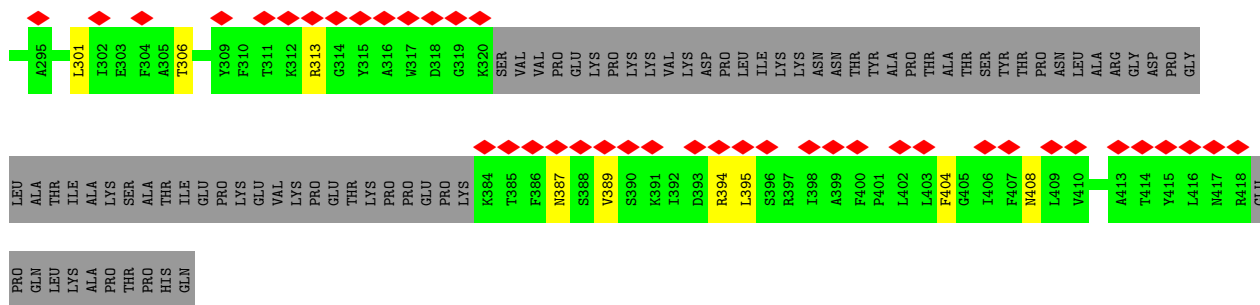
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: GABA(A) receptor subunit alpha-1

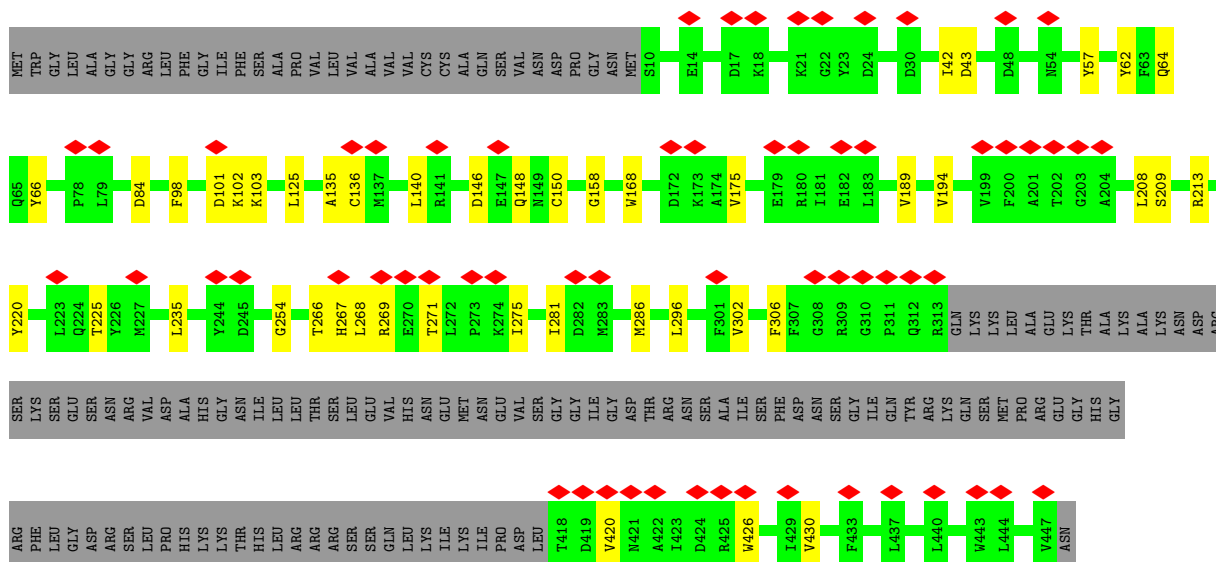


• Molecule 1: GABA(A) receptor subunit alpha-1

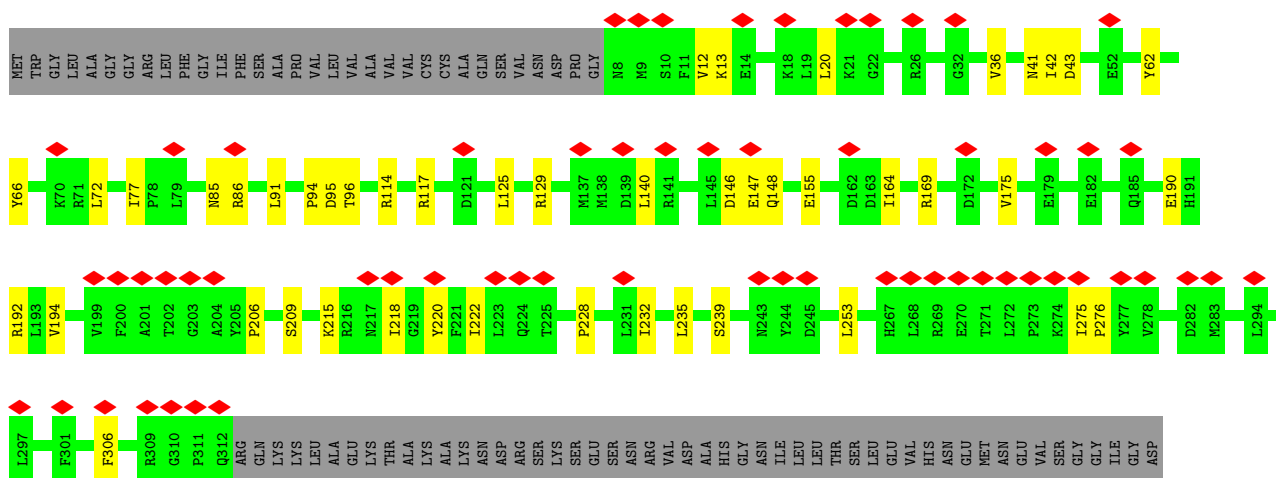


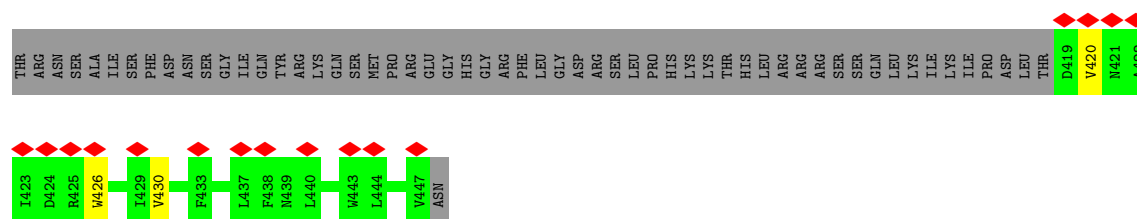


- Molecule 2: Gamma-aminobutyric acid receptor subunit beta-3

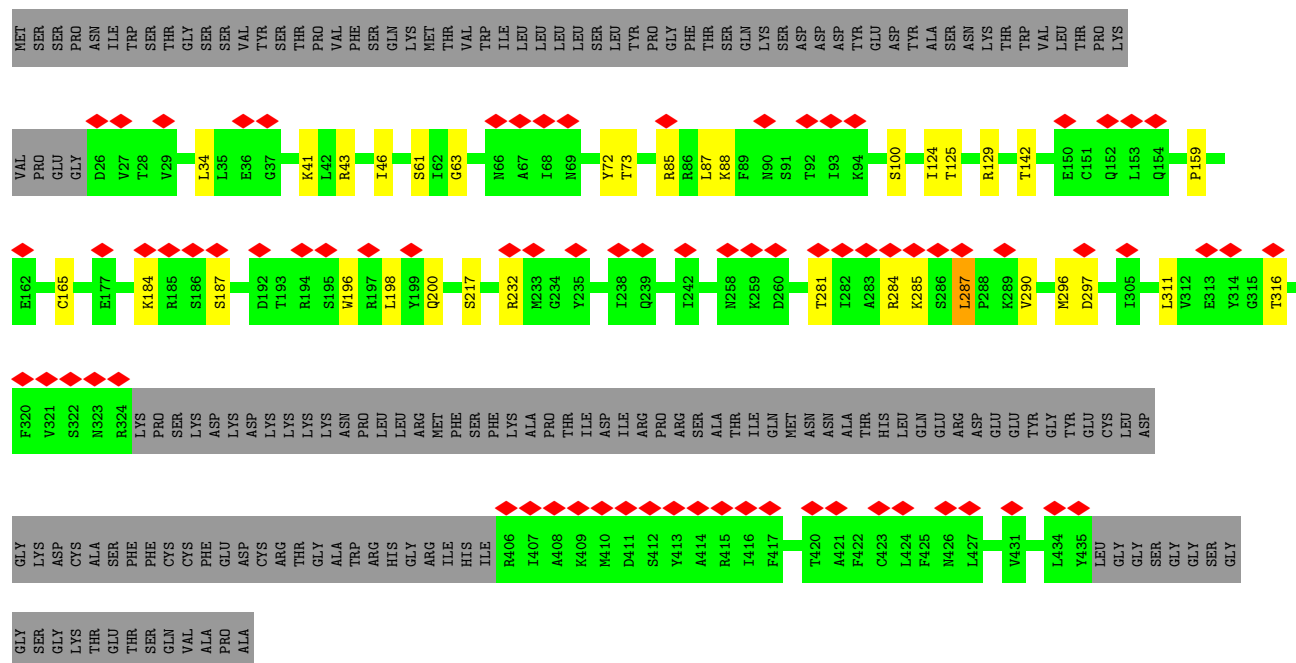


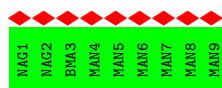
- Molecule 2: Gamma-aminobutyric acid receptor subunit beta-3





• Molecule 3: GABA(A) receptor subunit gamma-2





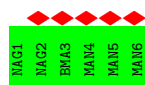
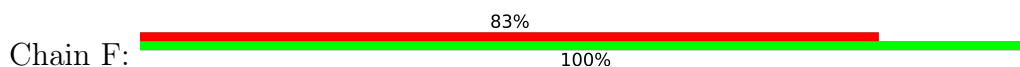
- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



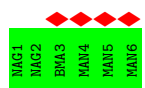
- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



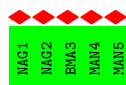
- Molecule 7: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 7: 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



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



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	119901	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2100	Depositor
Magnification	130000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.358	Depositor
Minimum map value	-0.224	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.08	Depositor
Map size (Å)	259.84, 259.84, 259.84	wwPDB
Map dimensions	290, 290, 290	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.89599997, 0.89599997, 0.89599997	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: CL, PIO, D10, MAN, EIE, OCT, BMA, NAG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.26	0/2875	0.47	0/3903
1	D	0.25	0/2873	0.48	0/3900
2	B	0.26	0/2828	0.48	0/3848
2	E	0.26	0/2821	0.47	0/3838
3	C	0.25	0/2786	0.47	0/3793
4	G	0.27	0/970	0.53	0/1309
All	All	0.26	0/15153	0.48	0/20591

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	2801	0	2797	38	0
1	D	2795	0	2792	34	0
2	B	2751	0	2743	27	0
2	E	2745	0	2737	30	0
3	C	2712	0	2699	25	0
4	G	949	0	901	12	0
5	a	105	0	88	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	b	28	0	25	0	0
6	e	28	0	25	0	0
7	F	72	0	61	0	0
7	H	72	0	61	0	0
8	d	61	0	52	0	0
9	A	47	0	44	4	0
9	D	47	0	44	2	0
10	A	8	0	18	0	0
11	A	10	0	22	1	0
11	B	10	0	22	1	0
11	C	10	0	22	0	0
11	E	10	0	22	0	0
12	A	1	0	0	0	0
12	C	1	0	0	0	0
12	D	1	0	0	0	0
13	C	14	0	13	0	0
14	D	24	0	0	1	0
All	All	15302	0	15188	147	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 5.

The worst 5 of 147 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:D:59:GLU:HB3	1:D:105:LYS:HZ2	1.50	0.77
1:D:277:LEU:HD12	1:D:278:PRO:HD2	1.74	0.70
2:E:41[A]:ASN:ND2	2:E:169:ARG:O	2.29	0.66
1:A:259:GLY:HA3	1:A:301:LEU:HD13	1.78	0.65
1:D:105:LYS:HE3	1:D:138:GLU:HG2	1.78	0.65

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	342/456 (75%)	337 (98%)	5 (2%)	0	100	100
1	D	341/456 (75%)	337 (99%)	4 (1%)	0	100	100
2	B	331/473 (70%)	325 (98%)	6 (2%)	0	100	100
2	E	331/473 (70%)	328 (99%)	3 (1%)	0	100	100
3	C	325/495 (66%)	320 (98%)	5 (2%)	0	100	100
4	G	121/123 (98%)	116 (96%)	5 (4%)	0	100	100
All	All	1791/2476 (72%)	1763 (98%)	28 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	308/405 (76%)	306 (99%)	2 (1%)	84	94
1	D	307/405 (76%)	306 (100%)	1 (0%)	91	97
2	B	301/417 (72%)	300 (100%)	1 (0%)	91	97
2	E	301/417 (72%)	301 (100%)	0	100	100
3	C	304/448 (68%)	303 (100%)	1 (0%)	91	97
4	G	93/93 (100%)	93 (100%)	0	100	100
All	All	1614/2185 (74%)	1609 (100%)	5 (0%)	90	97

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	258	PHE
1	A	294	TYR
2	B	136	CYS
3	C	287	LEU
1	D	258	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	229	GLN
2	E	119	HIS
2	E	312	GLN
3	C	426	ASN
1	D	68	GLN

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 ⓘ

30 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	$\# Z > 2$
7	NAG	F	1	2,7	14,14,15	0.25	0	17,19,21	0.41	0
7	NAG	F	2	7	14,14,15	0.23	0	17,19,21	0.42	0
7	BMA	F	3	7	11,11,12	0.60	0	15,15,17	0.77	0
7	MAN	F	4	7	11,11,12	0.20	0	15,15,17	0.34	0
7	MAN	F	5	7	11,11,12	0.19	0	15,15,17	0.24	0
7	MAN	F	6	7	11,11,12	0.21	0	15,15,17	0.30	0
7	NAG	H	1	2,7	14,14,15	0.23	0	17,19,21	0.43	0
7	NAG	H	2	7	14,14,15	0.24	0	17,19,21	0.42	0
7	BMA	H	3	7	11,11,12	0.61	0	15,15,17	0.73	0
7	MAN	H	4	7	11,11,12	0.19	0	15,15,17	0.23	0
7	MAN	H	5	7	11,11,12	0.20	0	15,15,17	0.24	0
7	MAN	H	6	7	11,11,12	0.21	0	15,15,17	0.26	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	a	1	1,5	14,14,15	0.24	0	17,19,21	0.50	0
5	NAG	a	2	5	14,14,15	0.23	0	17,19,21	0.38	0
5	BMA	a	3	5	11,11,12	0.60	0	15,15,17	0.92	0
5	MAN	a	4	5	11,11,12	0.20	0	15,15,17	0.48	0
5	MAN	a	5	5	11,11,12	0.19	0	15,15,17	0.25	0
5	MAN	a	6	5	11,11,12	0.19	0	15,15,17	0.25	0
5	MAN	a	7	5	11,11,12	0.19	0	15,15,17	0.36	0
5	MAN	a	8	5	11,11,12	0.21	0	15,15,17	0.24	0
5	MAN	a	9	5	11,11,12	0.19	0	15,15,17	0.25	0
6	NAG	b	1	2,6	14,14,15	0.25	0	17,19,21	0.46	0
6	NAG	b	2	6	14,14,15	0.22	0	17,19,21	0.41	0
8	NAG	d	1	8,1	14,14,15	0.25	0	17,19,21	0.36	0
8	NAG	d	2	8	14,14,15	0.22	0	17,19,21	0.56	0
8	BMA	d	3	8	11,11,12	0.74	0	15,15,17	0.75	0
8	MAN	d	4	8	11,11,12	0.23	0	15,15,17	0.31	0
8	MAN	d	5	8	11,11,12	0.20	0	15,15,17	0.24	0
6	NAG	e	1	2,6	14,14,15	0.23	0	17,19,21	0.54	0
6	NAG	e	2	6	14,14,15	0.23	0	17,19,21	0.40	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	F	1	2,7	-	0/6/23/26	0/1/1/1
7	NAG	F	2	7	-	0/6/23/26	0/1/1/1
7	BMA	F	3	7	-	2/2/19/22	0/1/1/1
7	MAN	F	4	7	-	1/2/19/22	0/1/1/1
7	MAN	F	5	7	-	0/2/19/22	0/1/1/1
7	MAN	F	6	7	-	1/2/19/22	0/1/1/1
7	NAG	H	1	2,7	-	0/6/23/26	0/1/1/1
7	NAG	H	2	7	-	0/6/23/26	0/1/1/1
7	BMA	H	3	7	-	0/2/19/22	0/1/1/1
7	MAN	H	4	7	-	0/2/19/22	0/1/1/1
7	MAN	H	5	7	-	1/2/19/22	0/1/1/1
7	MAN	H	6	7	-	0/2/19/22	0/1/1/1
5	NAG	a	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	a	2	5	-	0/6/23/26	0/1/1/1
5	BMA	a	3	5	-	0/2/19/22	0/1/1/1
5	MAN	a	4	5	-	1/2/19/22	0/1/1/1
5	MAN	a	5	5	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	MAN	a	6	5	-	0/2/19/22	0/1/1/1
5	MAN	a	7	5	-	0/2/19/22	0/1/1/1
5	MAN	a	8	5	-	0/2/19/22	0/1/1/1
5	MAN	a	9	5	-	0/2/19/22	0/1/1/1
6	NAG	b	1	2,6	-	0/6/23/26	0/1/1/1
6	NAG	b	2	6	-	0/6/23/26	0/1/1/1
8	NAG	d	1	8,1	-	0/6/23/26	0/1/1/1
8	NAG	d	2	8	-	0/6/23/26	0/1/1/1
8	BMA	d	3	8	-	0/2/19/22	0/1/1/1
8	MAN	d	4	8	-	0/2/19/22	0/1/1/1
8	MAN	d	5	8	-	0/2/19/22	0/1/1/1
6	NAG	e	1	2,6	-	0/6/23/26	0/1/1/1
6	NAG	e	2	6	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

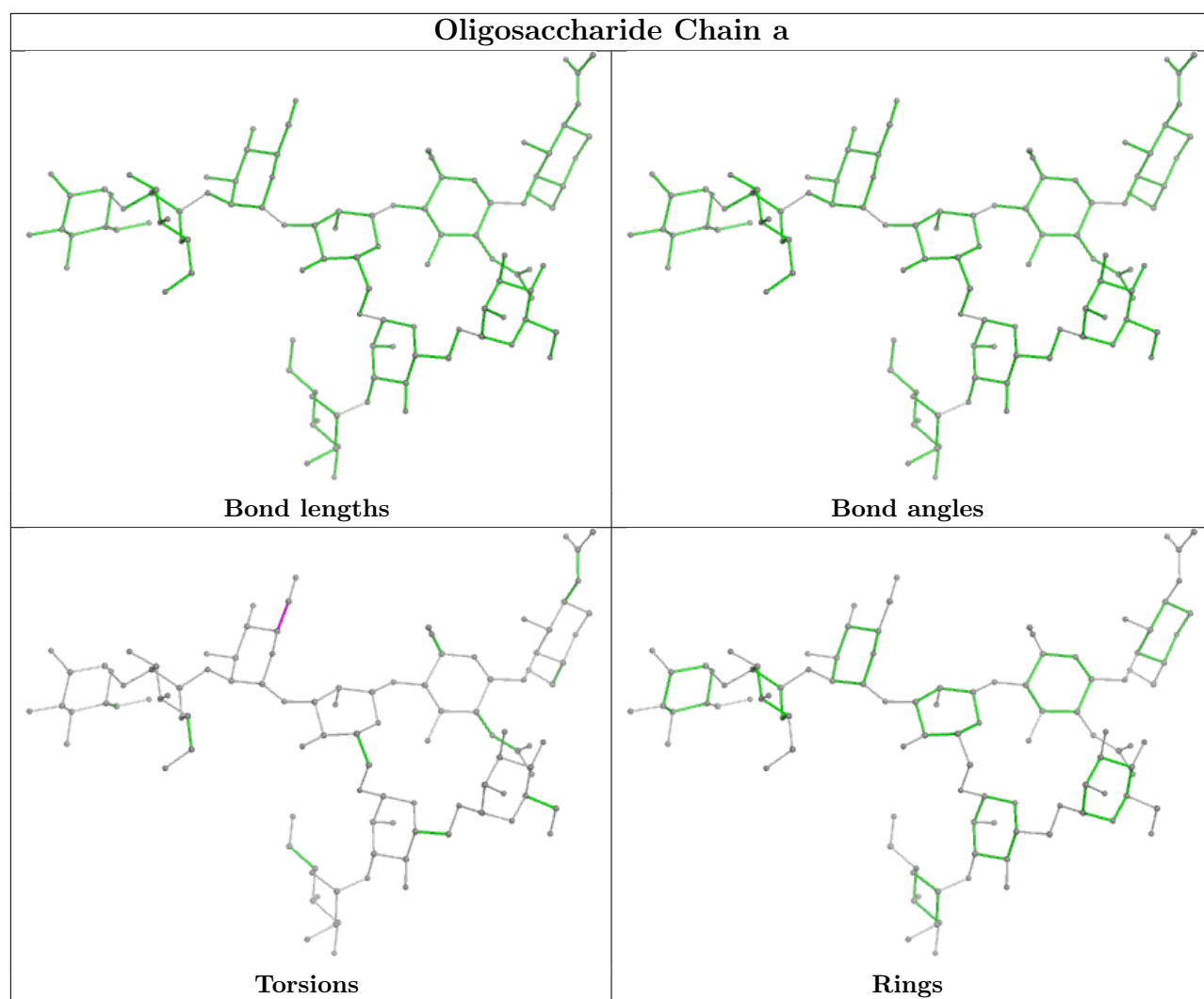
5 of 6 torsion outliers are listed below:

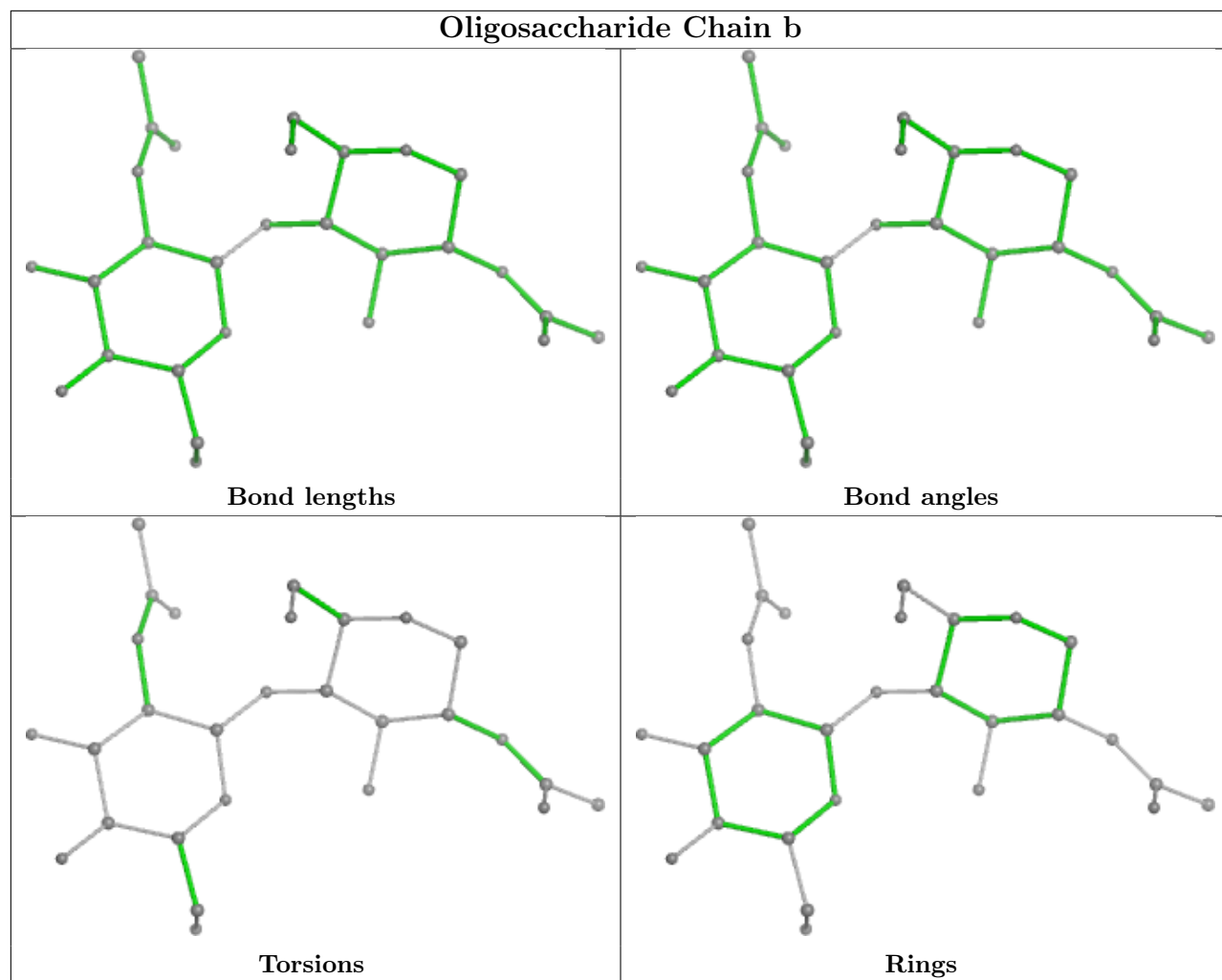
Mol	Chain	Res	Type	Atoms
7	F	3	BMA	C4-C5-C6-O6
7	F	3	BMA	O5-C5-C6-O6
7	H	5	MAN	O5-C5-C6-O6
5	a	4	MAN	O5-C5-C6-O6
7	F	4	MAN	O5-C5-C6-O6

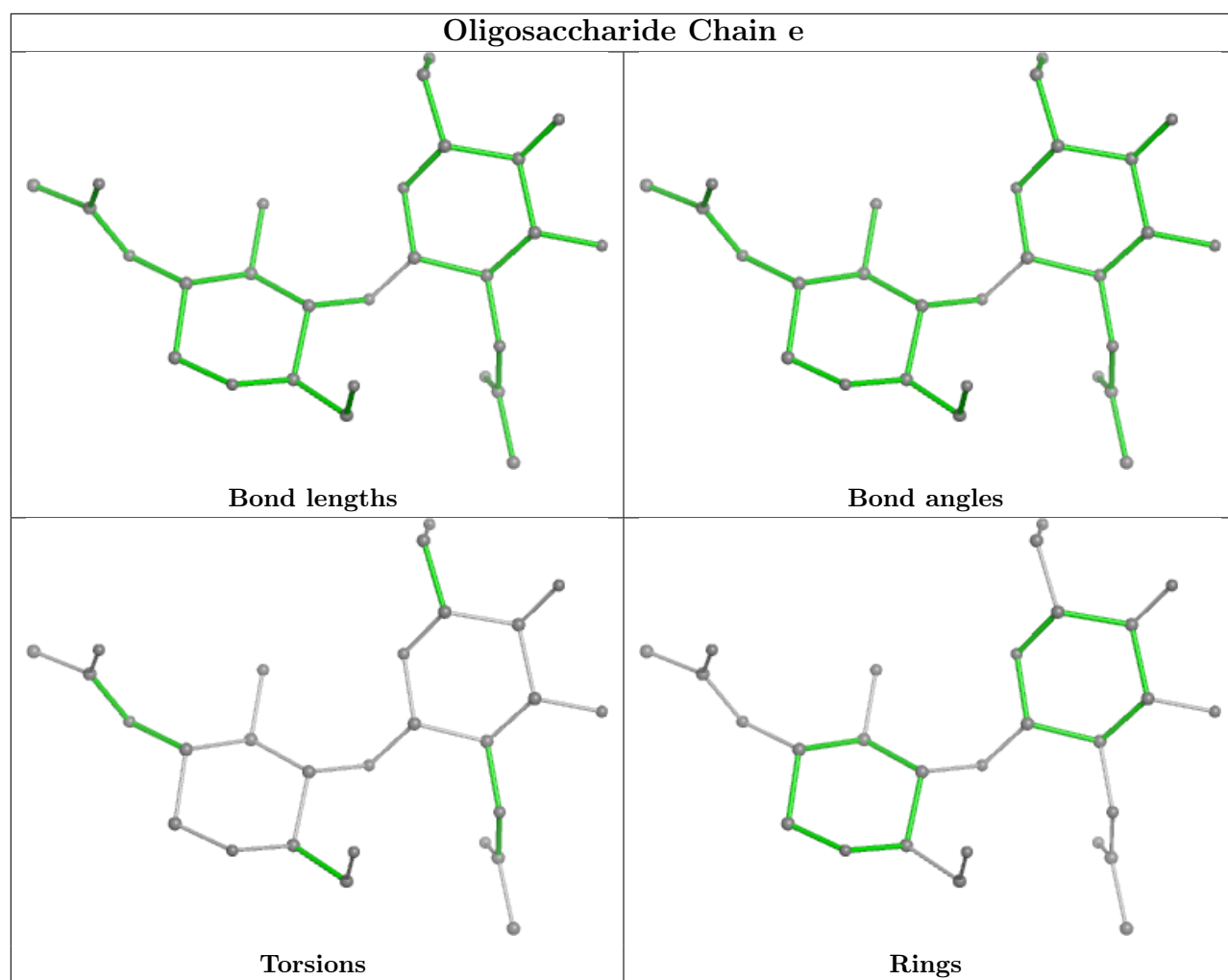
There are no ring outliers.

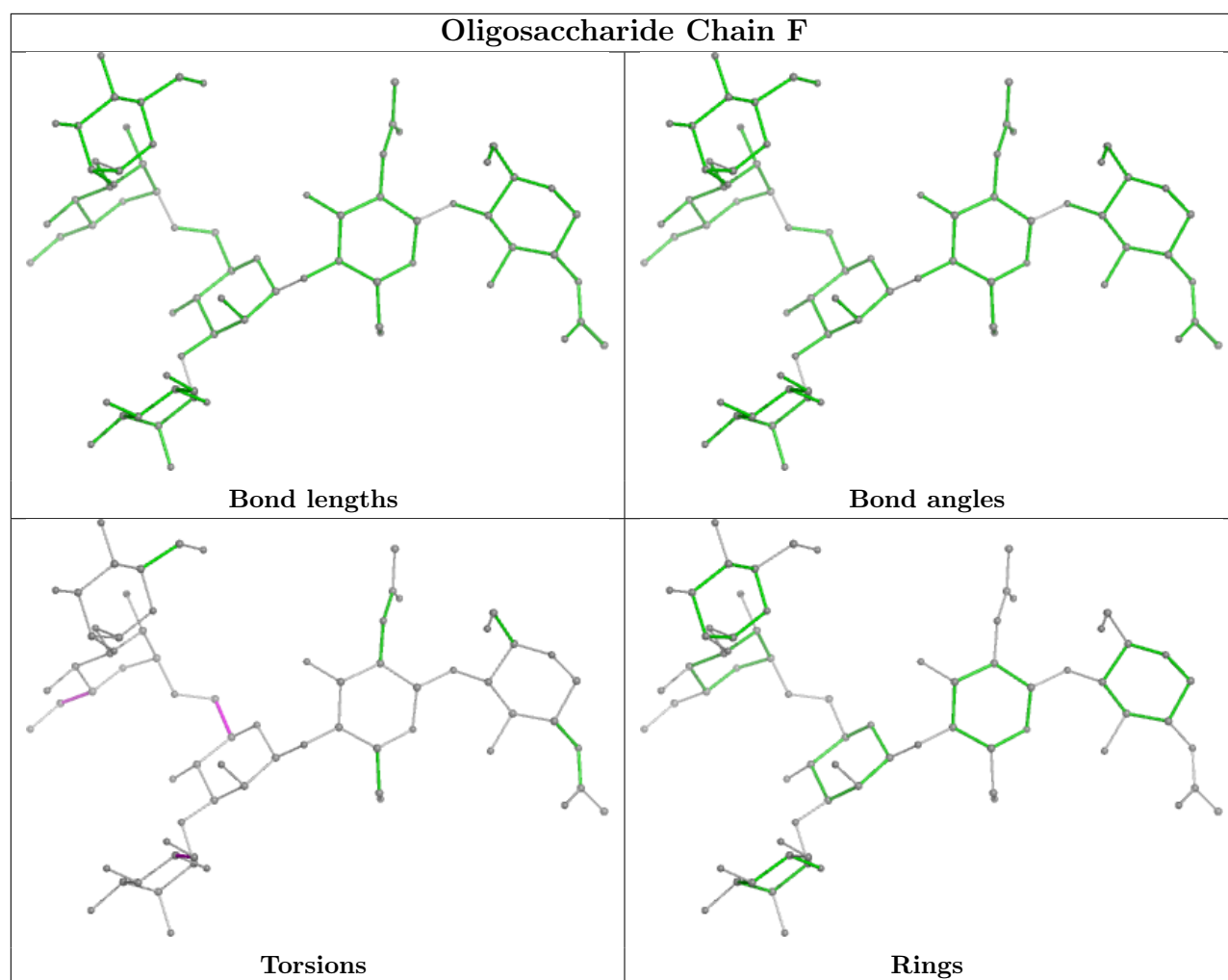
No monomer is involved in short contacts.

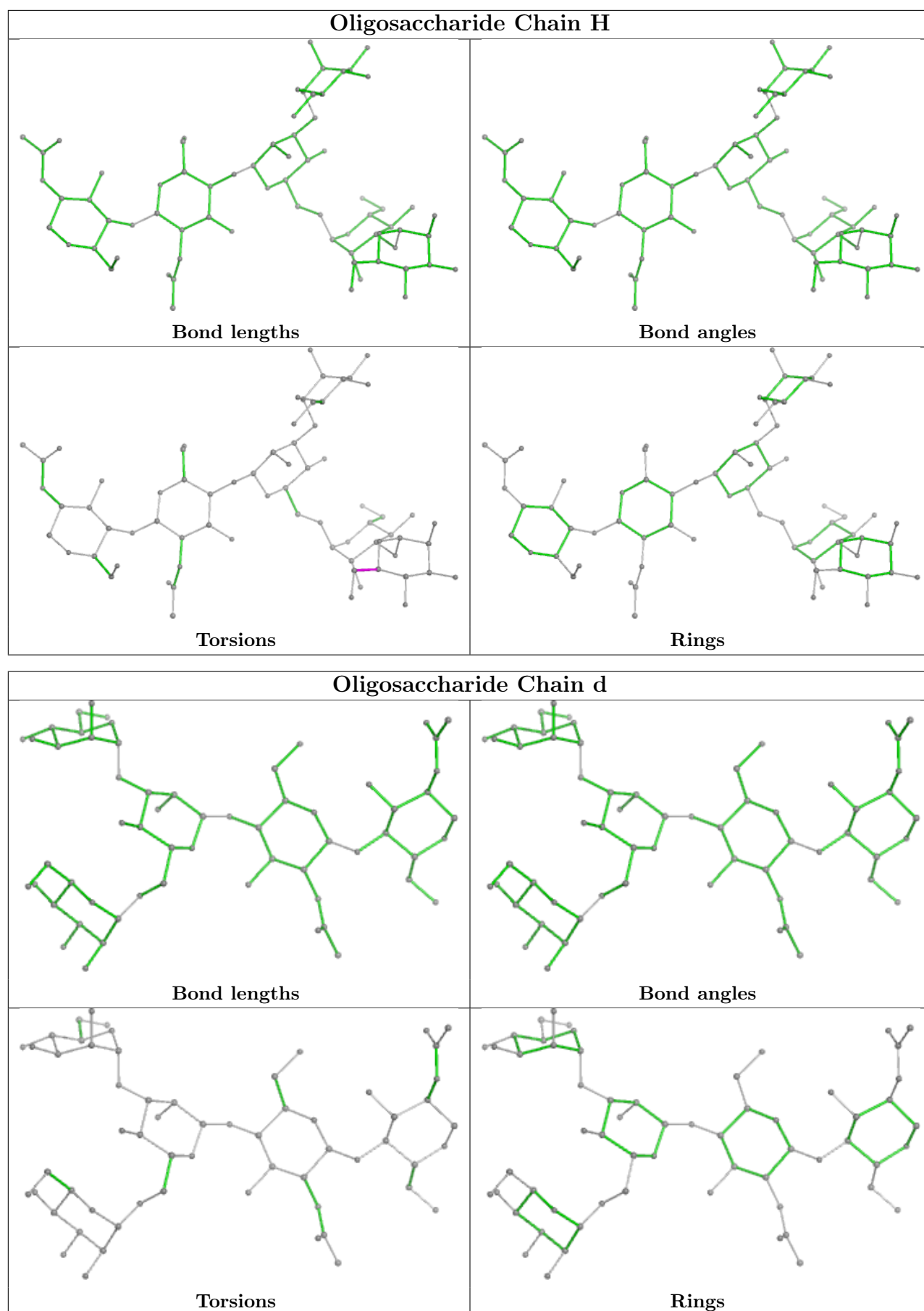
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry

Of 12 ligands modelled in this entry, 3 are monoatomic - leaving 9 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	D10	E	3201	-	9,9,9	0.14	0	8,8,8	0.11	0
14	EIE	D	502	-	23,26,26	0.88	2 (8%)	27,37,37	1.16	3 (11%)
11	D10	C	502	-	9,9,9	0.13	0	8,8,8	0.12	0
11	D10	A	503	-	9,9,9	0.14	0	8,8,8	0.11	0
9	PIO	A	501	-	47,47,47	0.32	0	61,65,65	0.24	0
13	NAG	C	503	3	14,14,15	0.24	0	17,19,21	0.48	0
10	OCT	A	502	-	7,7,7	0.12	0	6,6,6	0.11	0
11	D10	B	3101	-	9,9,9	0.13	0	8,8,8	0.16	0
9	PIO	D	503	-	47,47,47	0.32	0	61,65,65	0.26	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	D10	E	3201	-	-	0/7/7/7	-
14	EIE	D	502	-	-	0/10/26/26	0/2/3/3
11	D10	C	502	-	-	0/7/7/7	-
11	D10	A	503	-	-	0/7/7/7	-
9	PIO	A	501	-	-	2/44/68/68	0/1/1/1
13	NAG	C	503	3	-	2/6/23/26	0/1/1/1
10	OCT	A	502	-	-	0/5/5/5	-
11	D10	B	3101	-	-	0/7/7/7	-
9	PIO	D	503	-	-	2/44/68/68	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	D	502	EIE	C5-C4	-2.52	1.48	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	D	502	EIE	C14-N5	-2.15	1.31	1.35

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	D	502	EIE	C4-C5-N	-3.22	107.71	111.85
14	D	502	EIE	C12-C13-C8	-3.00	120.21	122.95
14	D	502	EIE	C4-C3-C	-2.20	127.23	131.50

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

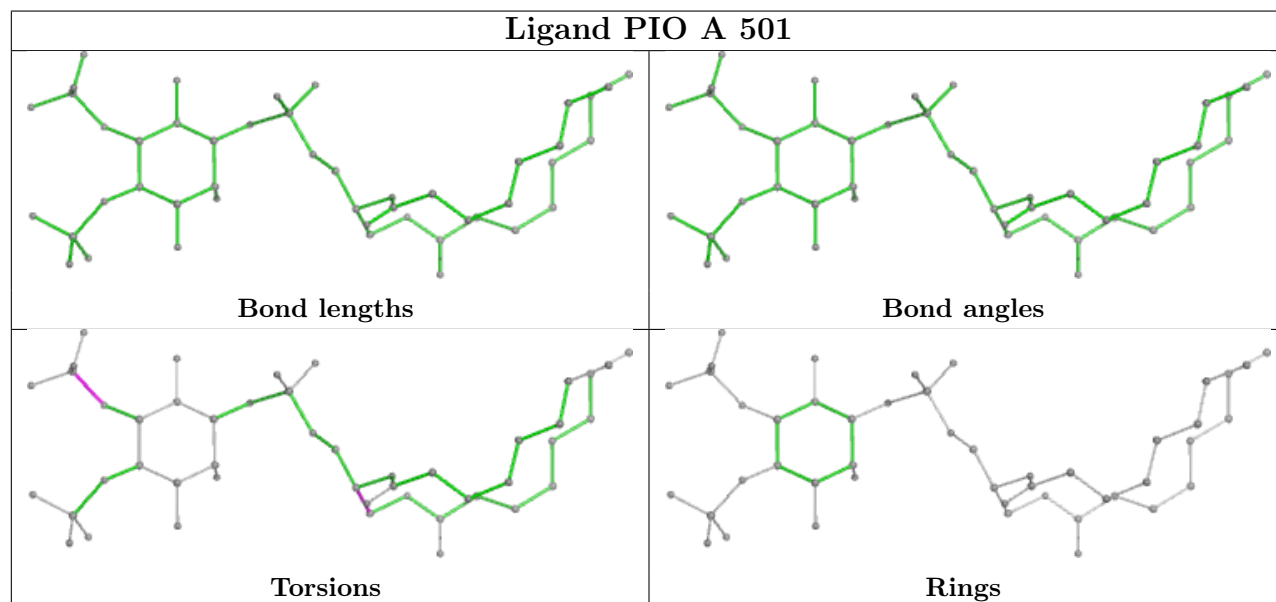
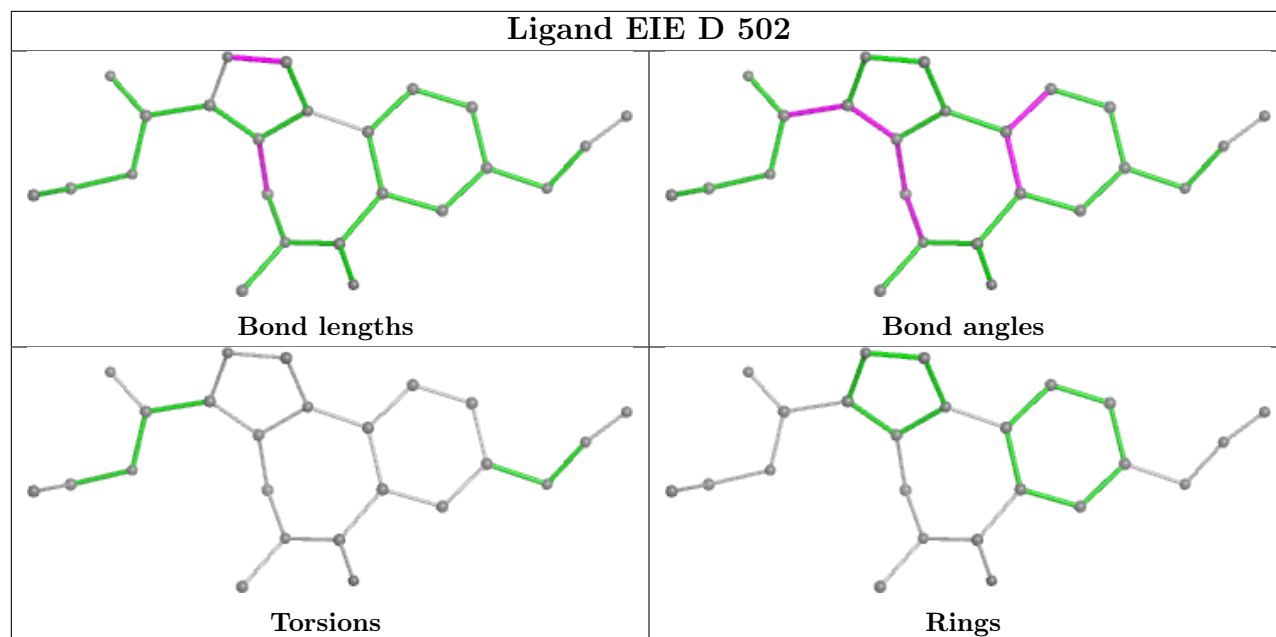
Mol	Chain	Res	Type	Atoms
9	D	503	PIO	C4-O4-P4-O42
9	A	501	PIO	O2C-C2C-C3C-O3C
9	A	501	PIO	C5-O5-P5-O53
9	D	503	PIO	C4-O4-P4-O43
13	C	503	NAG	C3-C2-N2-C7

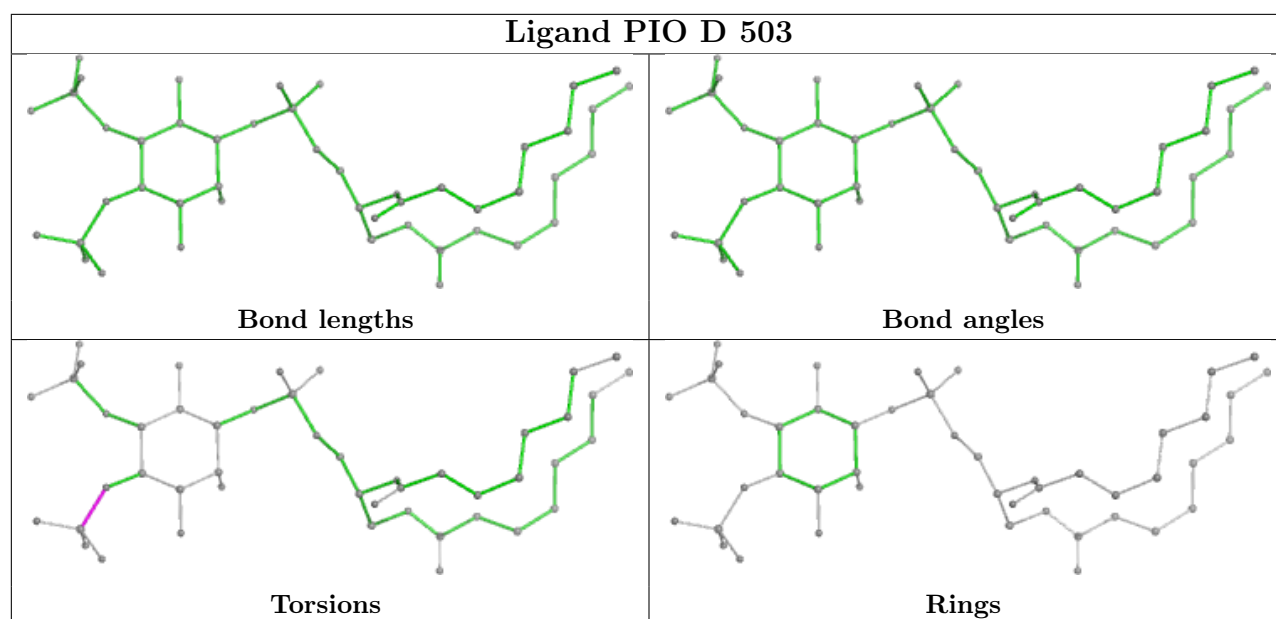
There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	D	502	EIE	1	0
11	A	503	D10	1	0
9	A	501	PIO	4	0
11	B	3101	D10	1	0
9	D	503	PIO	2	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

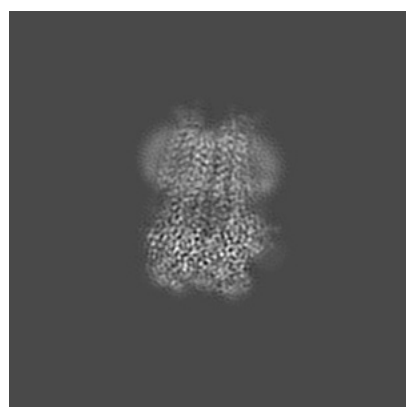
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-14076. These allow visual inspection of the internal detail of the map and identification of artifacts.

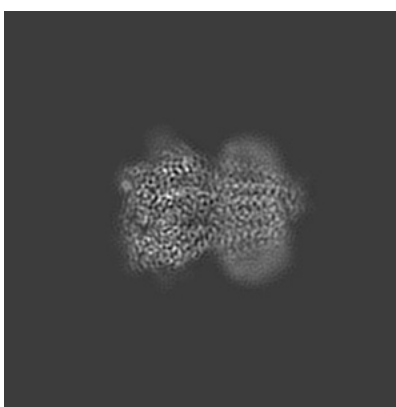
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

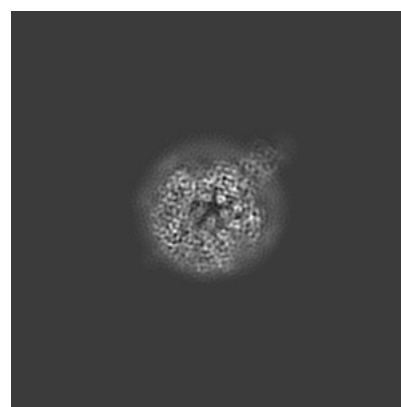
6.1.1 Primary map



X



Y



Z

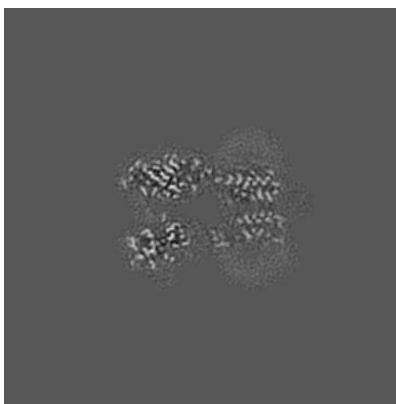
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

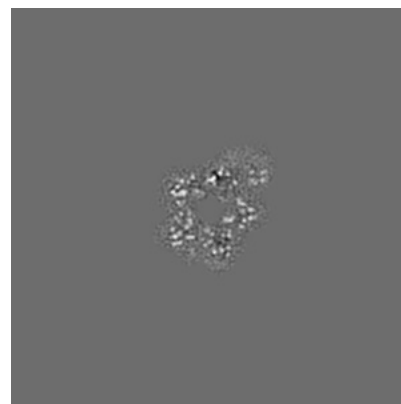
6.2.1 Primary map



X Index: 145



Y Index: 145



Z Index: 145

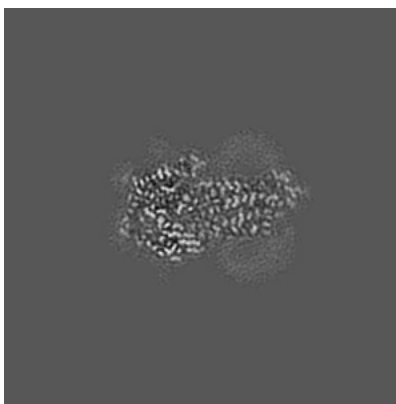
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

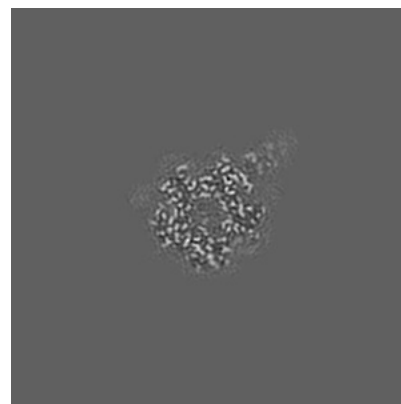
6.3.1 Primary map



X Index: 155



Y Index: 162

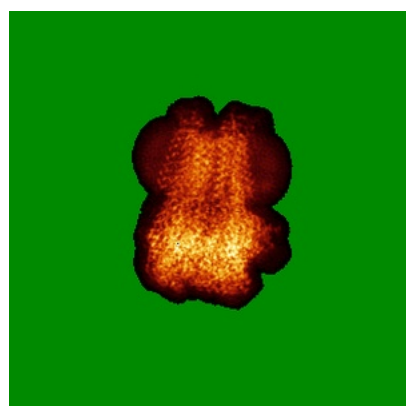


Z Index: 118

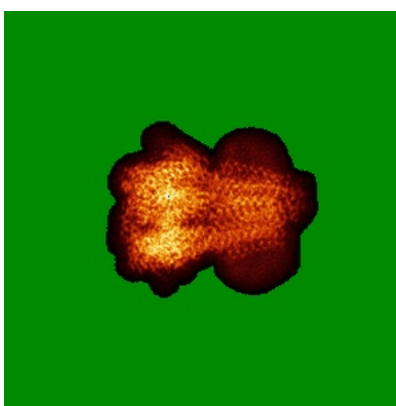
The images above show the largest variance slices of the map in three orthogonal directions.

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

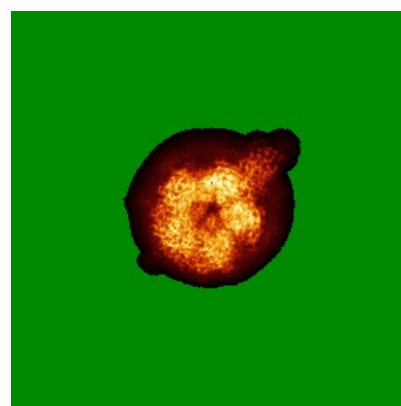
6.4.1 Primary map



X



Y

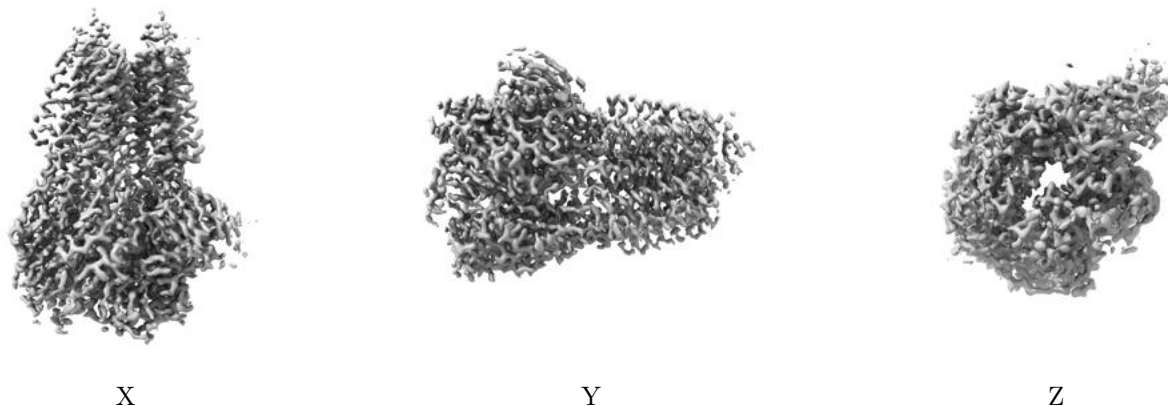


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



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

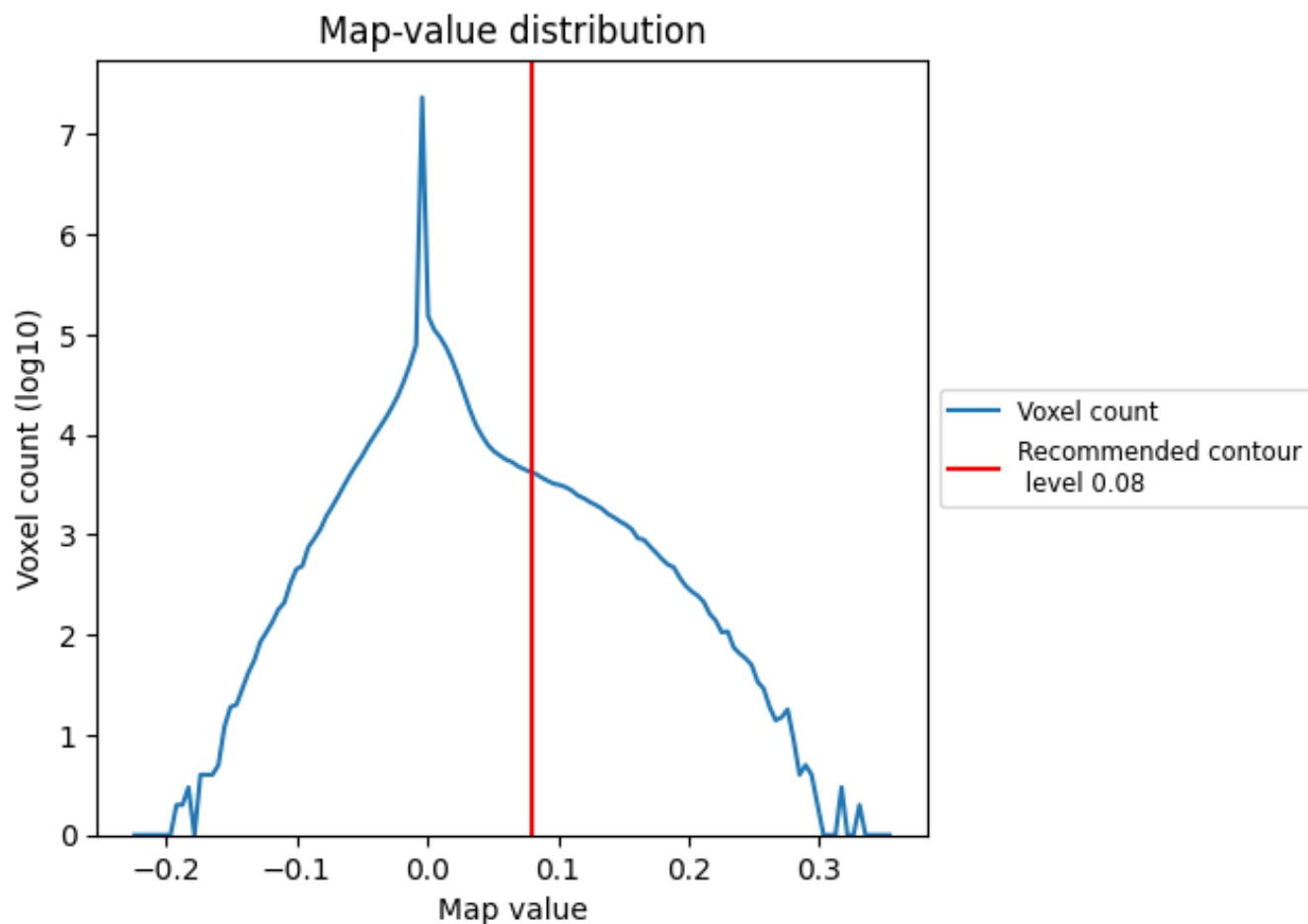
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

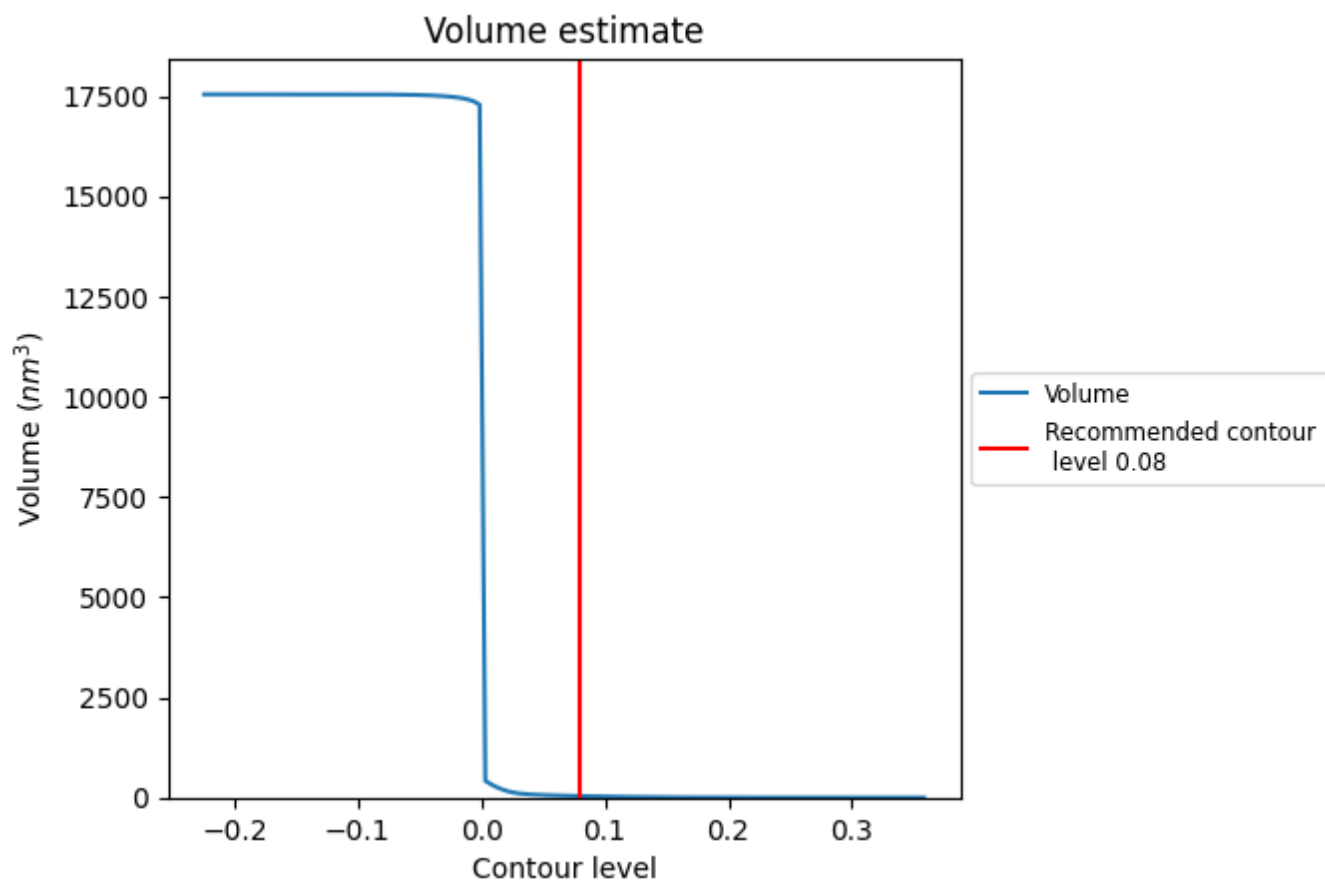
This section contains the results of statistical analysis of the map.

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.

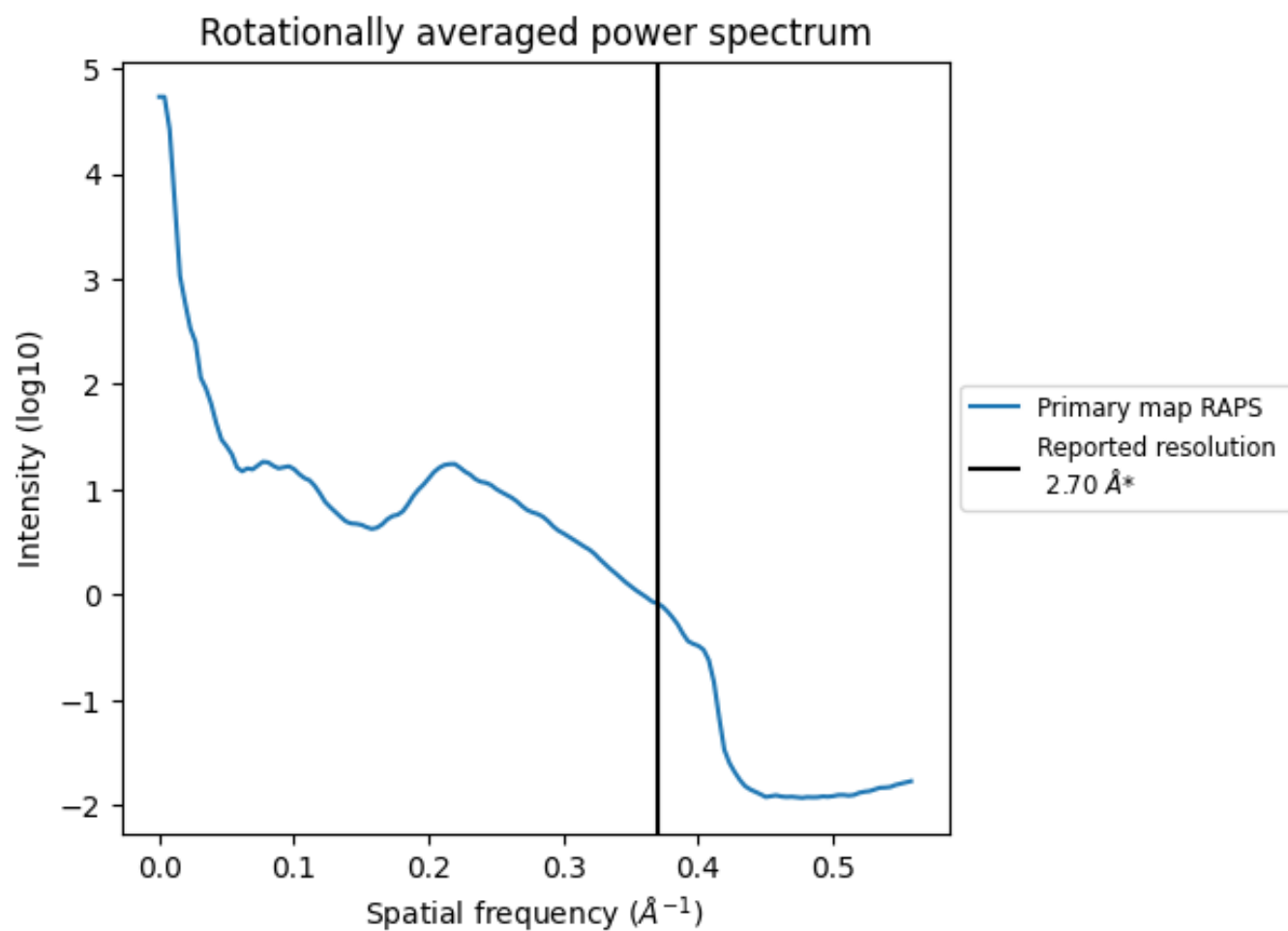
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 36 nm³; this corresponds to an approximate mass of 33 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 0.370 Å⁻¹

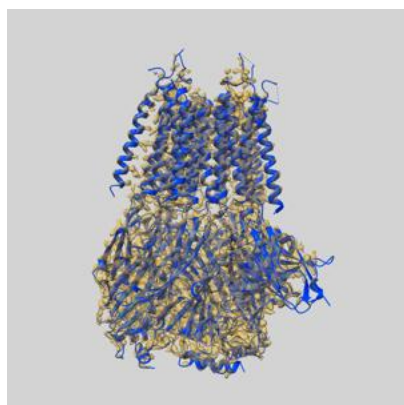
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

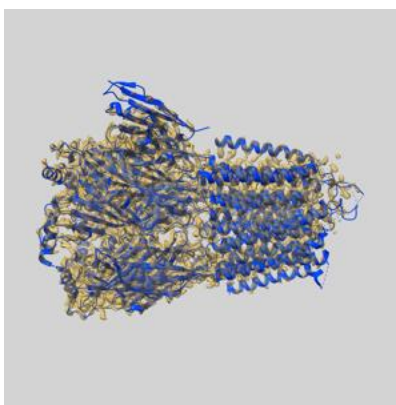
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-14076 and PDB model 7QNE. Per-residue inclusion information can be found in section [3](#) on page [11](#).

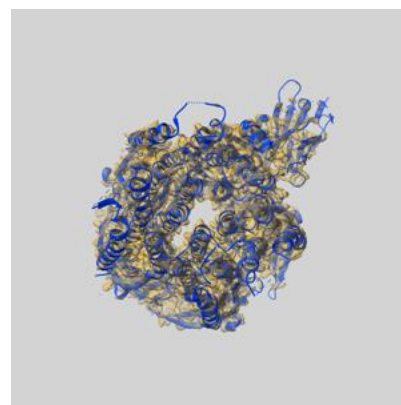
9.1 Map-model overlay [i](#)



X



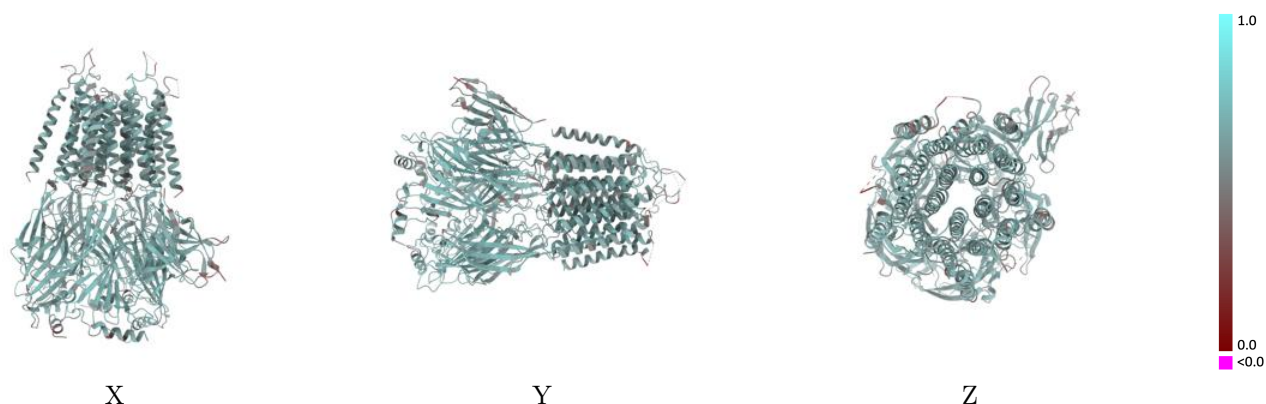
Y



Z

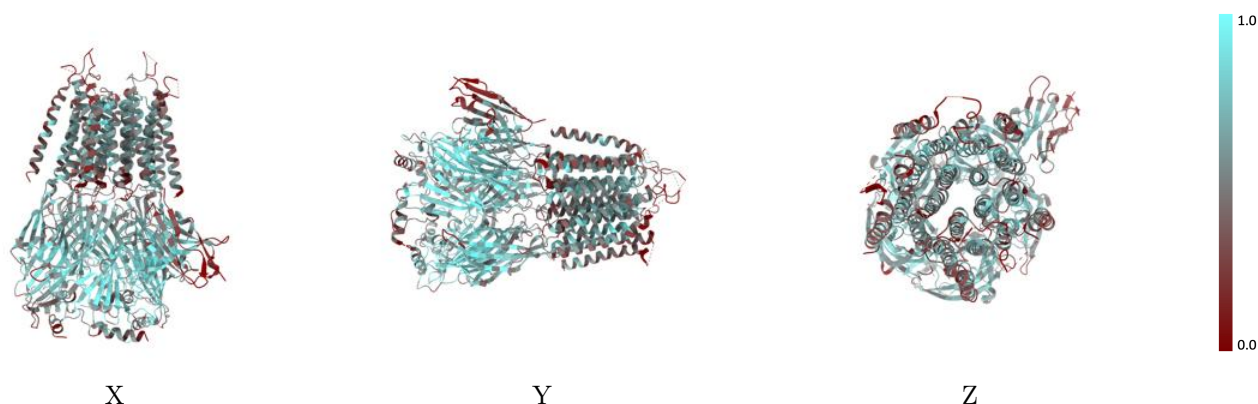
The images above show the 3D surface view of the map at the recommended contour level 0.08 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



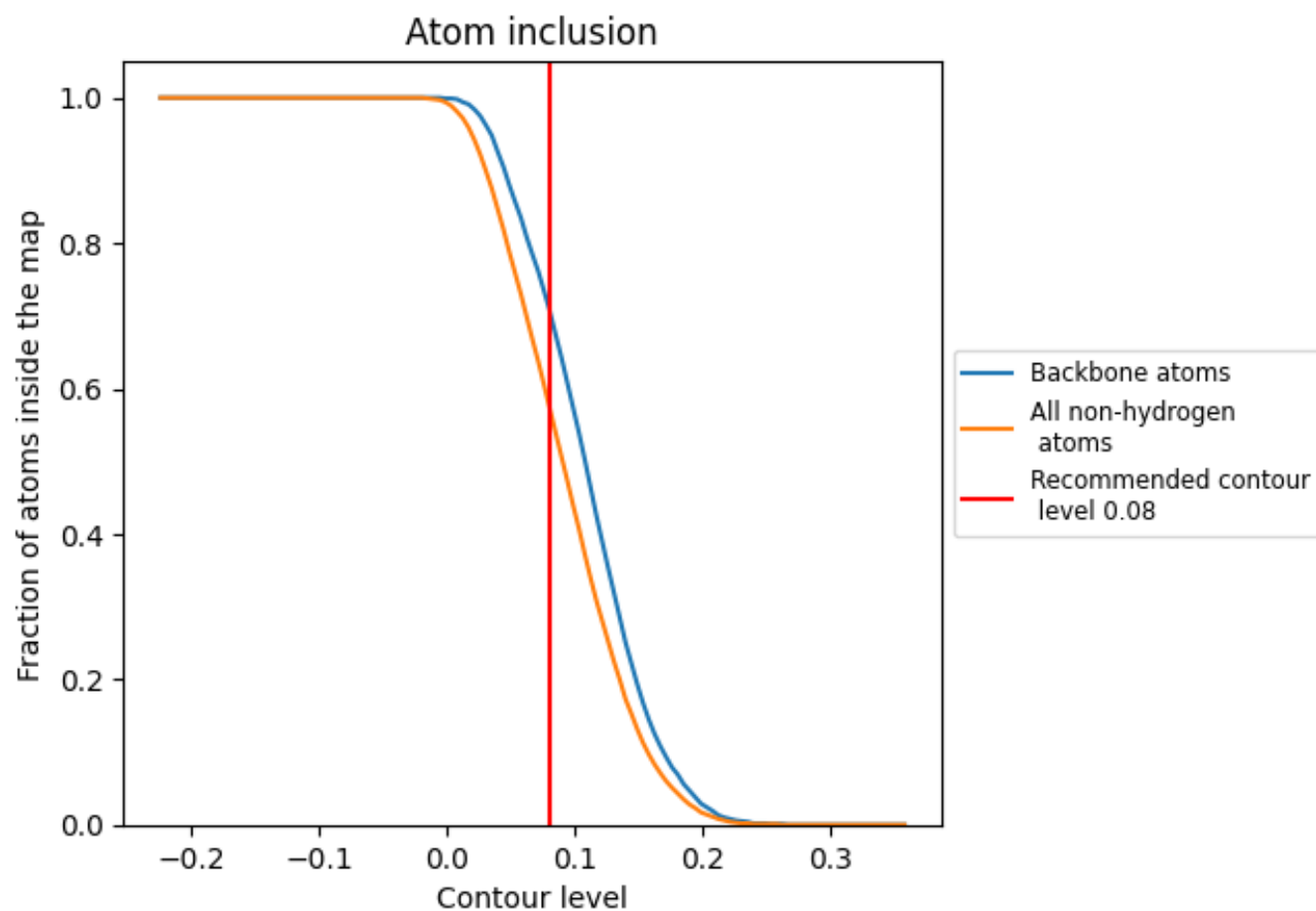
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.08).

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 58% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.08) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.5760	<div></div> 0.6060
A	<div></div> 0.6140	<div></div> 0.6100
B	<div></div> 0.6230	<div></div> 0.6220
C	<div></div> 0.6050	<div></div> 0.6180
D	<div></div> 0.5470	<div></div> 0.6030
E	<div></div> 0.5970	<div></div> 0.6130
F	<div></div> 0.1810	<div></div> 0.4930
G	<div></div> 0.4520	<div></div> 0.5710
H	<div></div> 0.1810	<div></div> 0.3770
a	<div></div> 0.0950	<div></div> 0.4430
b	<div></div> 0.1790	<div></div> 0.5350
d	<div></div> 0.0000	<div></div> 0.3230
e	<div></div> 0.0360	<div></div> 0.3730

