



wwPDB EM Validation Summary Report ⓘ

Mar 9, 2026 – 08:13 PM UTC

PDB ID : 8E92 / pdb_00008e92
EMDB ID : EMD-27953
Title : D-cycloserine and glutamate bound Human GluN1a-GluN2C NMDA receptor
in intact conformation
Authors : Chou, T.-H.; Furukawa, H.
Deposited on : 2022-08-26
Resolution : 3.96 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:

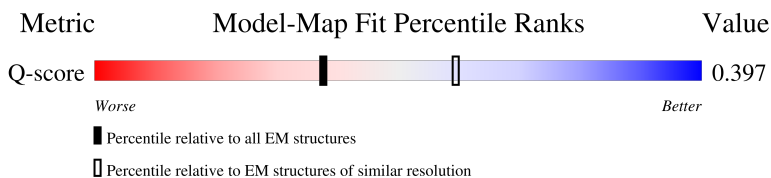
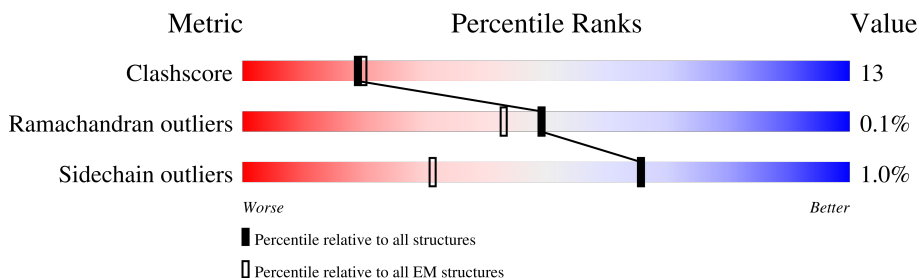
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 3.96 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	7646 (3.46 - 4.46)

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	847	<p>62% 14% 23%</p>
1	C	847	<p>62% 15% 23%</p>
2	B	880	<p>55% 16% 28%</p>
2	D	880	<p>54% 16% 28%</p>

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Mol	Chain	Length	Quality of chain
3	E	2	<p>50% 50%</p>
3	F	2	<p>100%</p>
3	G	2	<p>50% 50%</p>
3	I	2	<p>50% 50%</p>
3	L	2	<p>50% 50%</p>
3	N	2	<p>100%</p>
4	H	3	<p>33% 67% 33%</p>
4	J	3	<p>67% 33%</p>
4	K	3	<p>67% 33% 33%</p>
4	M	3	<p>33% 33% 67%</p>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18011 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 Glutamate receptor ionotropic, NMDA 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	654	4526	2890	786	829	21	0	0
1	C	655	4468	2868	764	818	18	0	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	HIS	ARG	conflict	UNP Q05586
A	9	PHE	LEU	conflict	UNP Q05586
A	17	PHE	VAL	conflict	UNP Q05586
A	22	SER	CYS	conflict	UNP Q05586
A	844	ASN	ARG	conflict	UNP Q05586
A	845	GLY	ARG	conflict	UNP Q05586
A	846	ALA	LYS	conflict	UNP Q05586
C	5	HIS	ARG	conflict	UNP Q05586
C	9	PHE	LEU	conflict	UNP Q05586
C	17	PHE	VAL	conflict	UNP Q05586
C	22	SER	CYS	conflict	UNP Q05586
C	844	ASN	ARG	conflict	UNP Q05586
C	845	GLY	ARG	conflict	UNP Q05586
C	846	ALA	LYS	conflict	UNP Q05586

- Molecule 2 is a protein called Glutamate receptor ionotropic, NMDA 2C.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	633	4229	2731	709	772	17	0	0
2	D	638	4366	2820	739	791	16	0	0

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-30	MET	-	expression tag	UNP Q14957
B	-29	GLY	-	expression tag	UNP Q14957
B	-28	THR	-	expression tag	UNP Q14957
B	-27	MET	-	expression tag	UNP Q14957
B	-26	ARG	-	expression tag	UNP Q14957
B	-25	LEU	-	expression tag	UNP Q14957
B	-24	PHE	-	expression tag	UNP Q14957
B	-23	LEU	-	expression tag	UNP Q14957
B	-22	LEU	-	expression tag	UNP Q14957
B	-21	ALA	-	expression tag	UNP Q14957
B	-20	VAL	-	expression tag	UNP Q14957
B	-19	LEU	-	expression tag	UNP Q14957
B	-18	PHE	-	expression tag	UNP Q14957
B	-17	LEU	-	expression tag	UNP Q14957
B	-16	PHE	-	expression tag	UNP Q14957
B	-15	SER	-	expression tag	UNP Q14957
B	-14	PHE	-	expression tag	UNP Q14957
B	-13	ALA	-	expression tag	UNP Q14957
B	-12	ARG	-	expression tag	UNP Q14957
B	-11	ALA	-	expression tag	UNP Q14957
B	-10	THR	-	expression tag	UNP Q14957
B	-9	GLY	-	expression tag	UNP Q14957
B	-8	TRP	-	expression tag	UNP Q14957
B	-7	SER	-	expression tag	UNP Q14957
B	-6	HIS	-	expression tag	UNP Q14957
B	-5	PRO	-	expression tag	UNP Q14957
B	-4	GLN	-	expression tag	UNP Q14957
B	-3	PHE	-	expression tag	UNP Q14957
B	-2	GLU	-	expression tag	UNP Q14957
B	-1	LYS	-	expression tag	UNP Q14957
B	0	GLY	-	expression tag	UNP Q14957
B	1	GLY	-	expression tag	UNP Q14957
B	2	GLY	-	expression tag	UNP Q14957
B	3	SER	-	expression tag	UNP Q14957
B	4	GLY	-	expression tag	UNP Q14957
B	5	GLY	-	expression tag	UNP Q14957
B	6	GLY	-	expression tag	UNP Q14957
B	7	SER	-	expression tag	UNP Q14957
B	8	GLY	-	expression tag	UNP Q14957
B	9	GLY	-	expression tag	UNP Q14957
B	10	SER	-	expression tag	UNP Q14957
B	11	ALA	-	expression tag	UNP Q14957
B	12	TRP	-	expression tag	UNP Q14957

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Chain	Residue	Modelled	Actual	Comment	Reference
B	13	SER	-	expression tag	UNP Q14957
B	14	HIS	-	expression tag	UNP Q14957
B	15	PRO	-	expression tag	UNP Q14957
B	16	GLN	-	expression tag	UNP Q14957
B	17	PHE	-	expression tag	UNP Q14957
B	18	GLU	-	expression tag	UNP Q14957
B	19	LYS	-	expression tag	UNP Q14957
B	20	GLY	-	expression tag	UNP Q14957
B	21	ALA	-	expression tag	UNP Q14957
B	22	LEU	-	expression tag	UNP Q14957
B	23	VAL	-	expression tag	UNP Q14957
B	24	PRO	-	expression tag	UNP Q14957
B	25	ARG	-	expression tag	UNP Q14957
D	-30	MET	-	expression tag	UNP Q14957
D	-29	GLY	-	expression tag	UNP Q14957
D	-28	THR	-	expression tag	UNP Q14957
D	-27	MET	-	expression tag	UNP Q14957
D	-26	ARG	-	expression tag	UNP Q14957
D	-25	LEU	-	expression tag	UNP Q14957
D	-24	PHE	-	expression tag	UNP Q14957
D	-23	LEU	-	expression tag	UNP Q14957
D	-22	LEU	-	expression tag	UNP Q14957
D	-21	ALA	-	expression tag	UNP Q14957
D	-20	VAL	-	expression tag	UNP Q14957
D	-19	LEU	-	expression tag	UNP Q14957
D	-18	PHE	-	expression tag	UNP Q14957
D	-17	LEU	-	expression tag	UNP Q14957
D	-16	PHE	-	expression tag	UNP Q14957
D	-15	SER	-	expression tag	UNP Q14957
D	-14	PHE	-	expression tag	UNP Q14957
D	-13	ALA	-	expression tag	UNP Q14957
D	-12	ARG	-	expression tag	UNP Q14957
D	-11	ALA	-	expression tag	UNP Q14957
D	-10	THR	-	expression tag	UNP Q14957
D	-9	GLY	-	expression tag	UNP Q14957
D	-8	TRP	-	expression tag	UNP Q14957
D	-7	SER	-	expression tag	UNP Q14957
D	-6	HIS	-	expression tag	UNP Q14957
D	-5	PRO	-	expression tag	UNP Q14957
D	-4	GLN	-	expression tag	UNP Q14957
D	-3	PHE	-	expression tag	UNP Q14957
D	-2	GLU	-	expression tag	UNP Q14957

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	LYS	-	expression tag	UNP Q14957
D	0	GLY	-	expression tag	UNP Q14957
D	1	GLY	-	expression tag	UNP Q14957
D	2	GLY	-	expression tag	UNP Q14957
D	3	SER	-	expression tag	UNP Q14957
D	4	GLY	-	expression tag	UNP Q14957
D	5	GLY	-	expression tag	UNP Q14957
D	6	GLY	-	expression tag	UNP Q14957
D	7	SER	-	expression tag	UNP Q14957
D	8	GLY	-	expression tag	UNP Q14957
D	9	GLY	-	expression tag	UNP Q14957
D	10	SER	-	expression tag	UNP Q14957
D	11	ALA	-	expression tag	UNP Q14957
D	12	TRP	-	expression tag	UNP Q14957
D	13	SER	-	expression tag	UNP Q14957
D	14	HIS	-	expression tag	UNP Q14957
D	15	PRO	-	expression tag	UNP Q14957
D	16	GLN	-	expression tag	UNP Q14957
D	17	PHE	-	expression tag	UNP Q14957
D	18	GLU	-	expression tag	UNP Q14957
D	19	LYS	-	expression tag	UNP Q14957
D	20	GLY	-	expression tag	UNP Q14957
D	21	ALA	-	expression tag	UNP Q14957
D	22	LEU	-	expression tag	UNP Q14957
D	23	VAL	-	expression tag	UNP Q14957
D	24	PRO	-	expression tag	UNP Q14957
D	25	ARG	-	expression tag	UNP Q14957

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



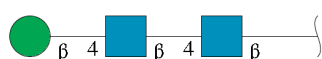
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
3	E	2	28	16	2	10	0	0
3	F	2	28	16	2	10	0	0
3	G	2	28	16	2	10	0	0

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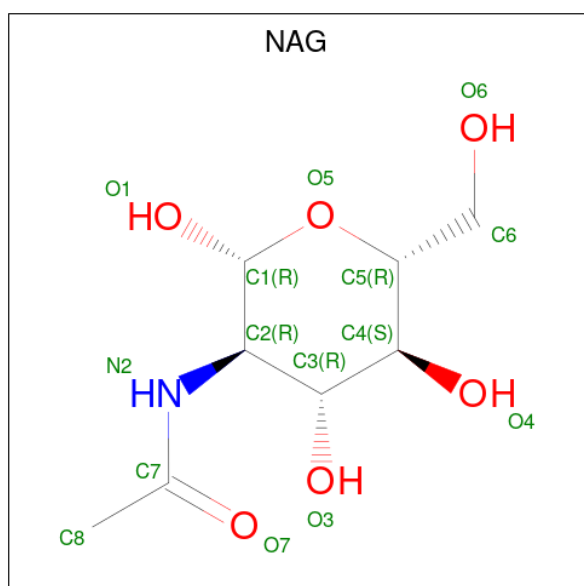
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
3	I	2	28	16	2	10	0	0
3	L	2	28	16	2	10	0	0
3	N	2	28	16	2	10	0	0

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

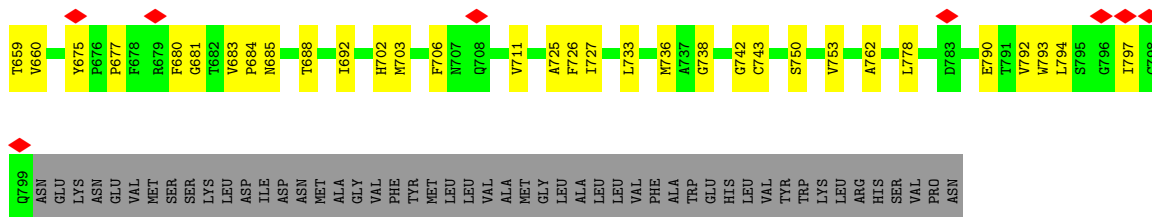


Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
4	H	3	39	22	2	15	0	0
4	J	3	39	22	2	15	0	0
4	K	3	39	22	2	15	0	0
4	M	3	39	22	2	15	0	0

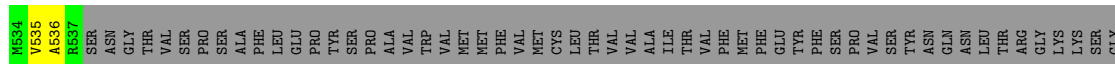
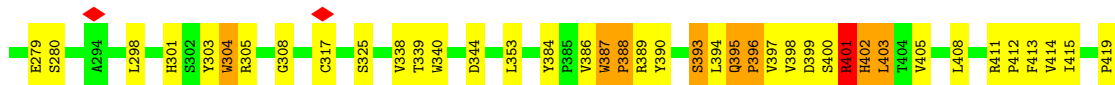
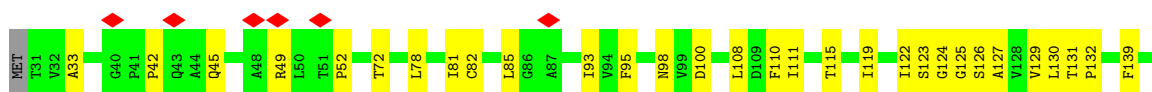
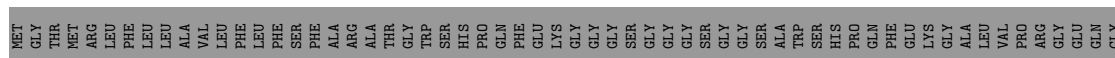
- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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



• Molecule 2: Glutamate receptor ionotropic, NMDA 2C



Chain F:  100%



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

Chain G:  50% 50%



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

Chain I:  50% 50%



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

Chain L:  50% 50%



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

Chain N:  100%



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

Chain H:  33% 67% 33%



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

Chain J:  67% 33%



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



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



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	43478	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	57.6	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	2800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	11.080	Depositor
Minimum map value	-6.534	Depositor
Average map value	0.008	Depositor
Map value standard deviation	0.472	Depositor
Recommended contour level	1.85	Depositor
Map size (\AA)	342.4, 342.4, 342.4	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.856, 0.856, 0.856	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: 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	0.34	6/4628 (0.1%)	0.51	3/6364 (0.0%)
1	C	0.23	1/4570 (0.0%)	0.44	0/6302
2	B	0.30	2/4330 (0.0%)	0.61	4/5985 (0.1%)
2	D	0.53	4/4475 (0.1%)	0.91	18/6175 (0.3%)
All	All	0.37	13/18003 (0.1%)	0.64	25/24826 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	C	0	1
2	B	0	1
2	D	0	2
All	All	0	6

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	677	PRO	CB-CG	21.21	2.55	1.49
2	D	677	PRO	CG-CD	-14.66	1.00	1.50
2	B	388	PRO	CG-CD	-14.10	1.02	1.50
1	A	454	CYS	CA-CB	9.54	1.67	1.53
1	C	519	ILE	CG1-CD1	-9.18	1.16	1.51

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	677	PRO	CB-CG-CD	-33.15	0.03	106.10
2	B	388	PRO	N-CD-CG	-21.55	70.88	103.20
2	D	677	PRO	CA-N-CD	-20.38	83.47	112.00
2	B	388	PRO	CA-CB-CG	-16.20	73.72	104.50
2	D	677	PRO	N-CA-CB	-15.67	86.80	103.25

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	438	GLY	Peptide
1	A	454	CYS	Peptide
2	B	204	GLY	Peptide
1	C	299	GLU	Peptide
2	D	325	SER	Peptide

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	4526	0	4023	100	0
1	C	4468	0	3913	100	0
2	B	4229	0	3776	103	0
2	D	4366	0	3958	135	0
3	E	28	0	25	1	0
3	F	28	0	25	0	0
3	G	28	0	25	1	0
3	I	28	0	25	1	0
3	L	28	0	25	1	0
3	N	28	0	25	0	0
4	H	39	0	34	2	0
4	J	39	0	32	2	0
4	K	39	0	34	1	0
4	M	39	0	34	2	0
5	A	56	0	52	1	0
5	C	28	0	26	0	0
5	D	14	0	13	0	0
All	All	18011	0	16045	436	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 13.

The worst 5 of 436 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:TYR:HE2	1:A:263:PRO:HD2	1.12	1.12
1:A:232:TYR:HE2	1:A:263:PRO:CD	1.64	1.10
2:D:248:PRO:HB3	2:D:390:TYR:O	1.57	1.02
2:D:461:ILE:O	2:D:465:LEU:HB2	1.58	1.01
2:B:81:ILE:HD11	2:B:111:ILE:HD13	1.48	0.94

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	650/847 (77%)	573 (88%)	77 (12%)	0	100	100
1	C	651/847 (77%)	603 (93%)	48 (7%)	0	100	100
2	B	625/880 (71%)	560 (90%)	63 (10%)	2 (0%)	36	69
2	D	632/880 (72%)	568 (90%)	63 (10%)	1 (0%)	43	75
All	All	2558/3454 (74%)	2304 (90%)	251 (10%)	3 (0%)	49	81

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	470	LYS
2	B	205	PRO
2	B	99	VAL

5.3.2 Protein sidechains [i](#)

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	405/728 (56%)	402 (99%)	3 (1%)	76	79
1	C	388/728 (53%)	388 (100%)	0	100	100
2	B	371/735 (50%)	370 (100%)	1 (0%)	86	85
2	D	394/735 (54%)	382 (97%)	12 (3%)	36	57
All	All	1558/2926 (53%)	1542 (99%)	16 (1%)	65	76

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	471	PHE
2	D	469	VAL
2	D	398	VAL
2	D	468	VAL
2	D	395	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	402	HIS
2	D	218	GLN
1	C	28	ASN
2	B	116	HIS
1	C	171	HIS

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

24 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
3	NAG	E	1	1,3	14,14,15	0.25	0	17,19,21	0.54	0
3	NAG	E	2	3	14,14,15	0.27	0	17,19,21	0.41	0
3	NAG	F	1	1,3	14,14,15	0.23	0	17,19,21	0.45	0
3	NAG	F	2	3	14,14,15	0.22	0	17,19,21	0.40	0
3	NAG	G	1	1,3	14,14,15	0.49	0	17,19,21	1.43	2 (11%)
3	NAG	G	2	3	14,14,15	0.43	0	17,19,21	0.49	0
4	NAG	H	1	4,2	14,14,15	0.79	1 (7%)	17,19,21	0.84	1 (5%)
4	NAG	H	2	4	14,14,15	0.57	0	17,19,21	0.48	0
4	BMA	H	3	4	11,11,12	0.64	0	15,15,17	0.80	0
3	NAG	I	1	3,2	14,14,15	0.44	0	17,19,21	0.70	1 (5%)
3	NAG	I	2	3	14,14,15	0.24	0	17,19,21	0.46	0
4	NAG	J	1	4,1	14,14,15	0.65	0	17,19,21	0.85	1 (5%)
4	NAG	J	2	4	14,14,15	0.18	0	17,19,21	0.46	0
4	BMA	J	3	4	11,11,12	0.59	0	15,15,17	0.68	0
4	NAG	K	1	4,1	14,14,15	0.49	0	17,19,21	0.50	0
4	NAG	K	2	4	14,14,15	0.42	0	17,19,21	1.39	2 (11%)
4	BMA	K	3	4	11,11,12	0.61	0	15,15,17	0.99	1 (6%)
3	NAG	L	1	1,3	14,14,15	0.52	0	17,19,21	0.75	0
3	NAG	L	2	3	14,14,15	0.33	0	17,19,21	0.40	0
4	NAG	M	1	4,1	14,14,15	0.69	1 (7%)	17,19,21	1.00	1 (5%)
4	NAG	M	2	4	14,14,15	0.24	0	17,19,21	0.62	1 (5%)
4	BMA	M	3	4	11,11,12	0.59	0	15,15,17	0.67	0
3	NAG	N	1	3,2	14,14,15	0.37	0	17,19,21	0.35	0
3	NAG	N	2	3	14,14,15	0.40	0	17,19,21	0.37	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
3	NAG	E	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	NAG	G	1	1,3	-	6/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
4	NAG	H	1	4,2	-	2/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1
4	BMA	H	3	4	-	1/2/19/22	0/1/1/1
3	NAG	I	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	1/6/23/26	0/1/1/1
4	NAG	J	1	4,1	-	3/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1
4	BMA	J	3	4	-	0/2/19/22	0/1/1/1
4	NAG	K	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	K	2	4	-	6/6/23/26	0/1/1/1
4	BMA	K	3	4	-	2/2/19/22	0/1/1/1
3	NAG	L	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	L	2	3	-	0/6/23/26	0/1/1/1
4	NAG	M	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	M	2	4	-	2/6/23/26	0/1/1/1
4	BMA	M	3	4	-	0/2/19/22	0/1/1/1
3	NAG	N	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	N	2	3	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	1	NAG	O5-C1	-2.49	1.39	1.43
4	M	1	NAG	O5-C1	2.26	1.47	1.43

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	1	NAG	C2-N2-C7	4.69	129.19	122.90
4	K	2	NAG	C2-N2-C7	4.56	129.01	122.90
4	M	1	NAG	C1-O5-C5	3.82	117.31	112.19
3	G	1	NAG	C1-C2-N2	2.45	114.29	110.43
4	K	2	NAG	C1-C2-N2	2.31	114.07	110.43

There are no chirality outliers.

5 of 42 torsion outliers are listed below:

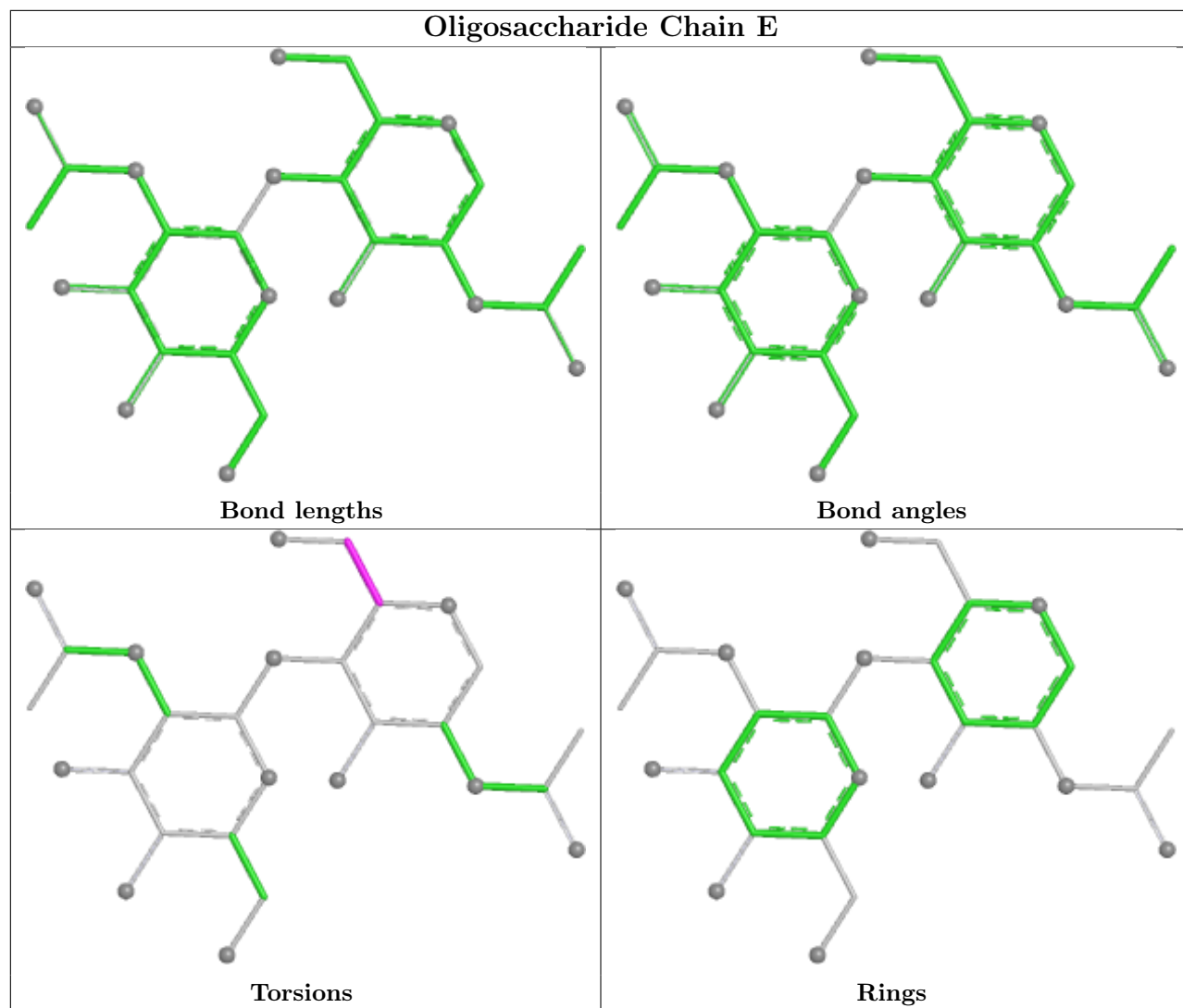
Mol	Chain	Res	Type	Atoms
4	J	1	NAG	C1-C2-N2-C7
4	K	2	NAG	C4-C5-C6-O6
4	H	1	NAG	C4-C5-C6-O6
3	N	2	NAG	C4-C5-C6-O6
4	K	2	NAG	O5-C5-C6-O6

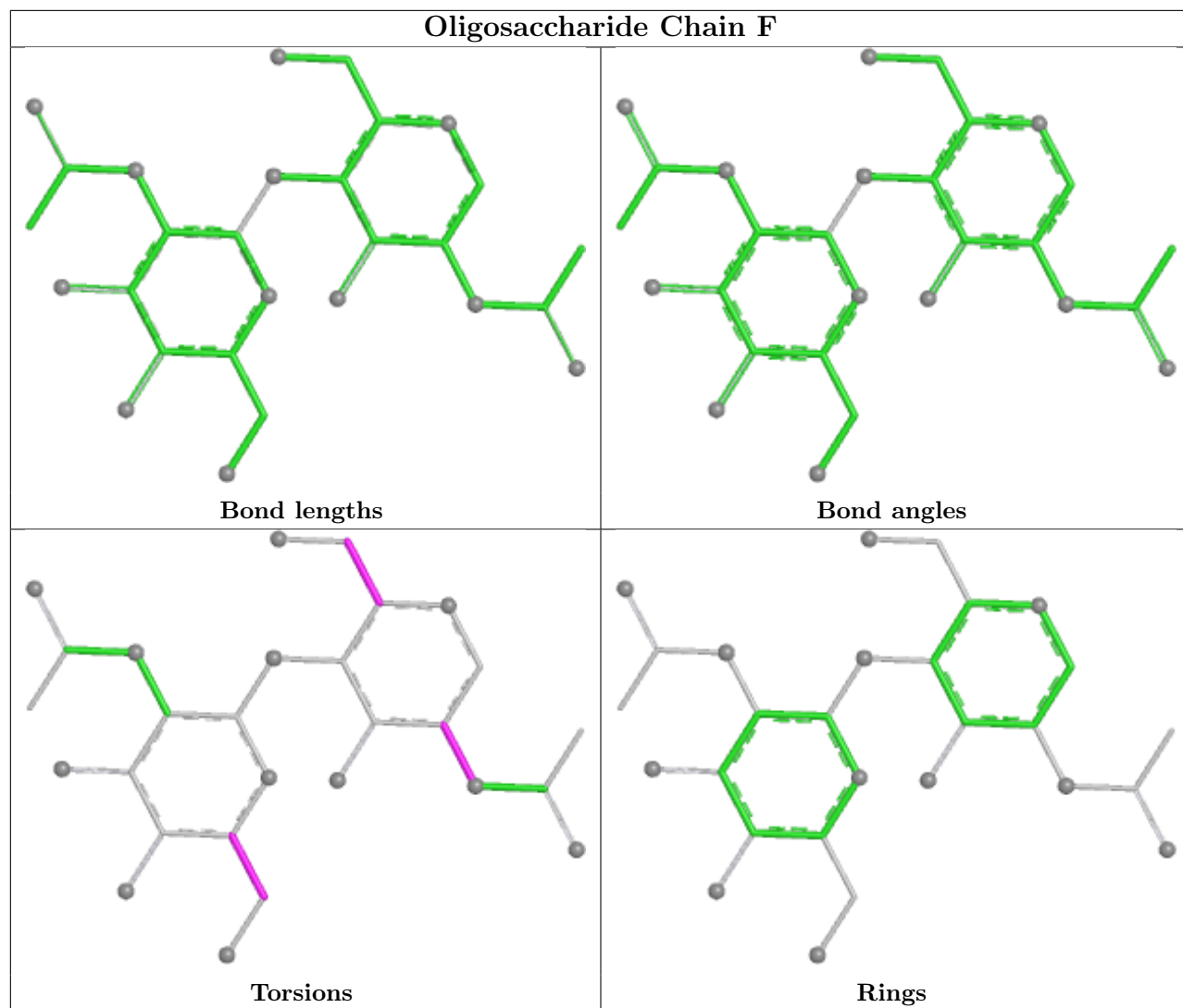
There are no ring outliers.

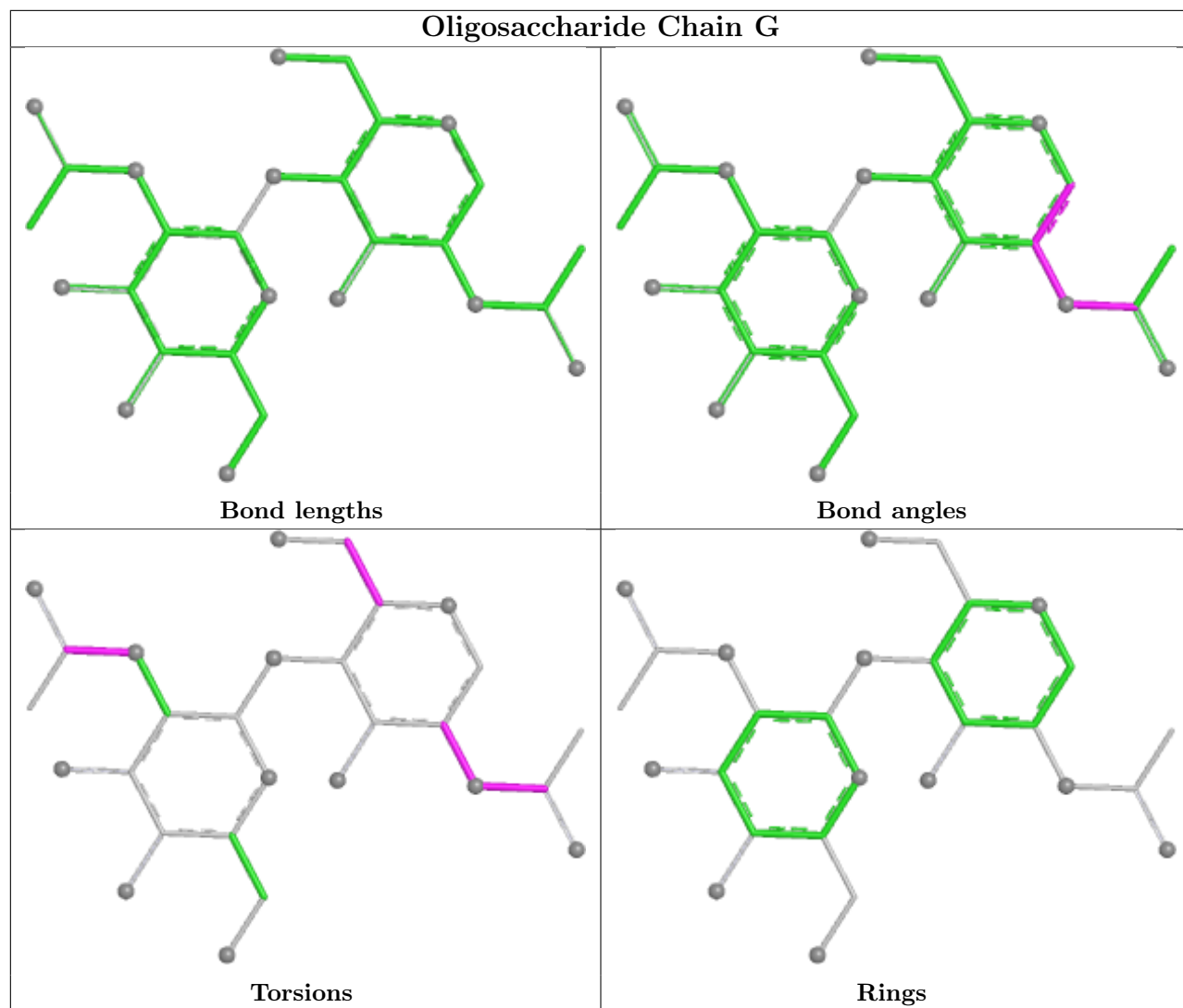
11 monomers are involved in 11 short contacts:

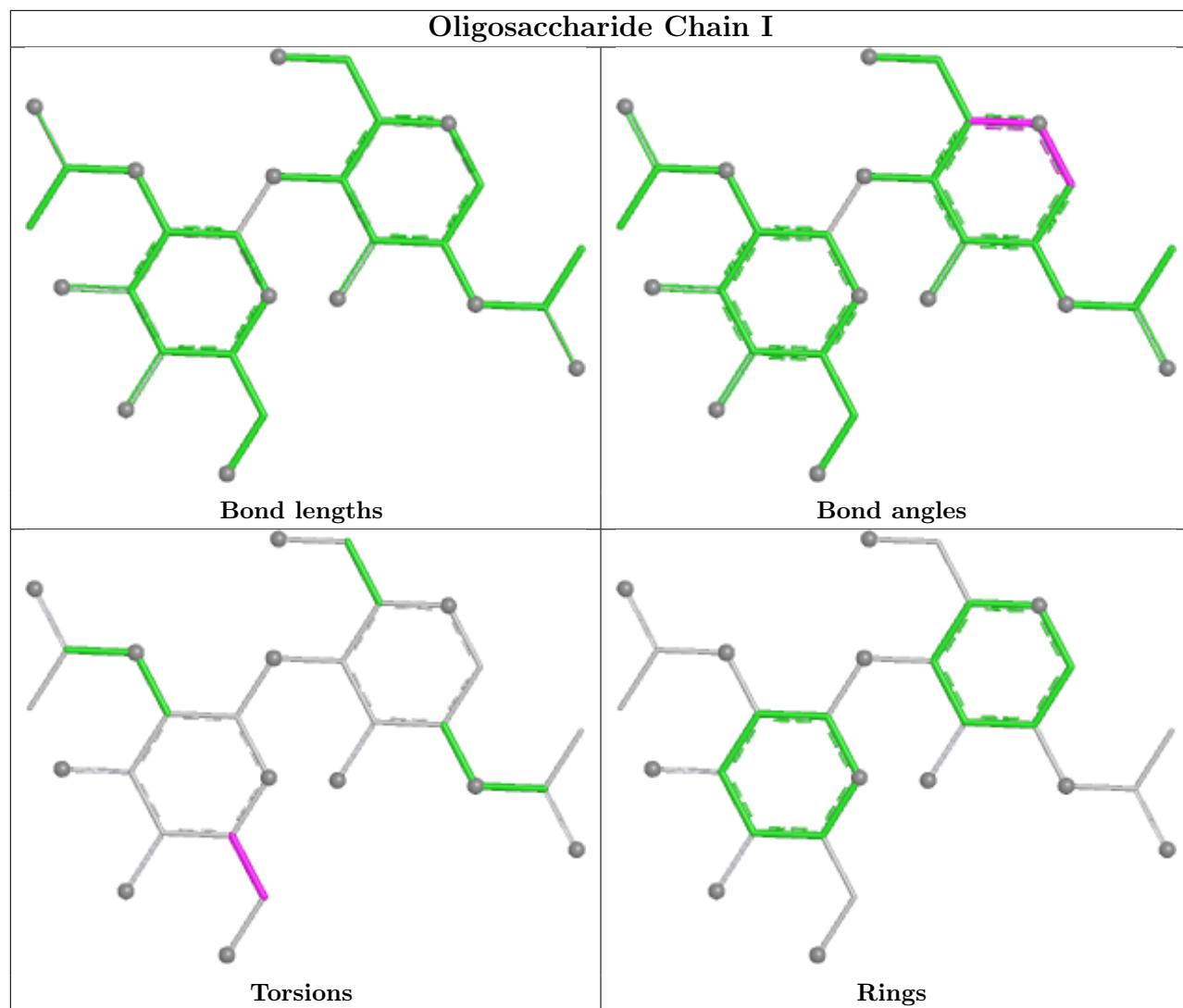
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	1	NAG	1	0
4	H	3	BMA	1	0
4	M	1	NAG	2	0
3	I	1	NAG	1	0
3	G	1	NAG	1	0
4	H	1	NAG	1	0
4	J	1	NAG	2	0
4	K	2	NAG	1	0
4	H	2	NAG	1	0
3	E	1	NAG	1	0
4	M	2	NAG	2	0

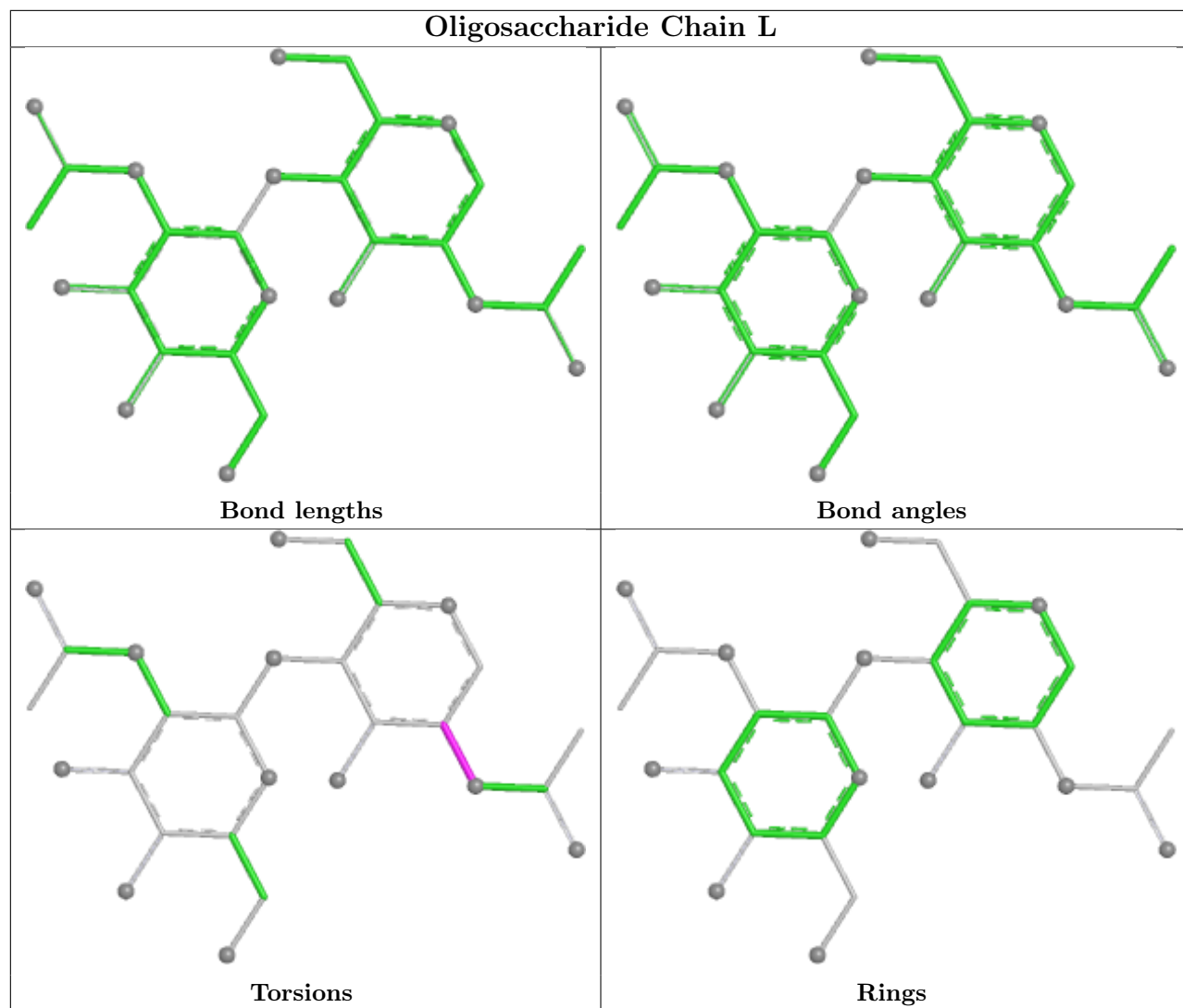
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

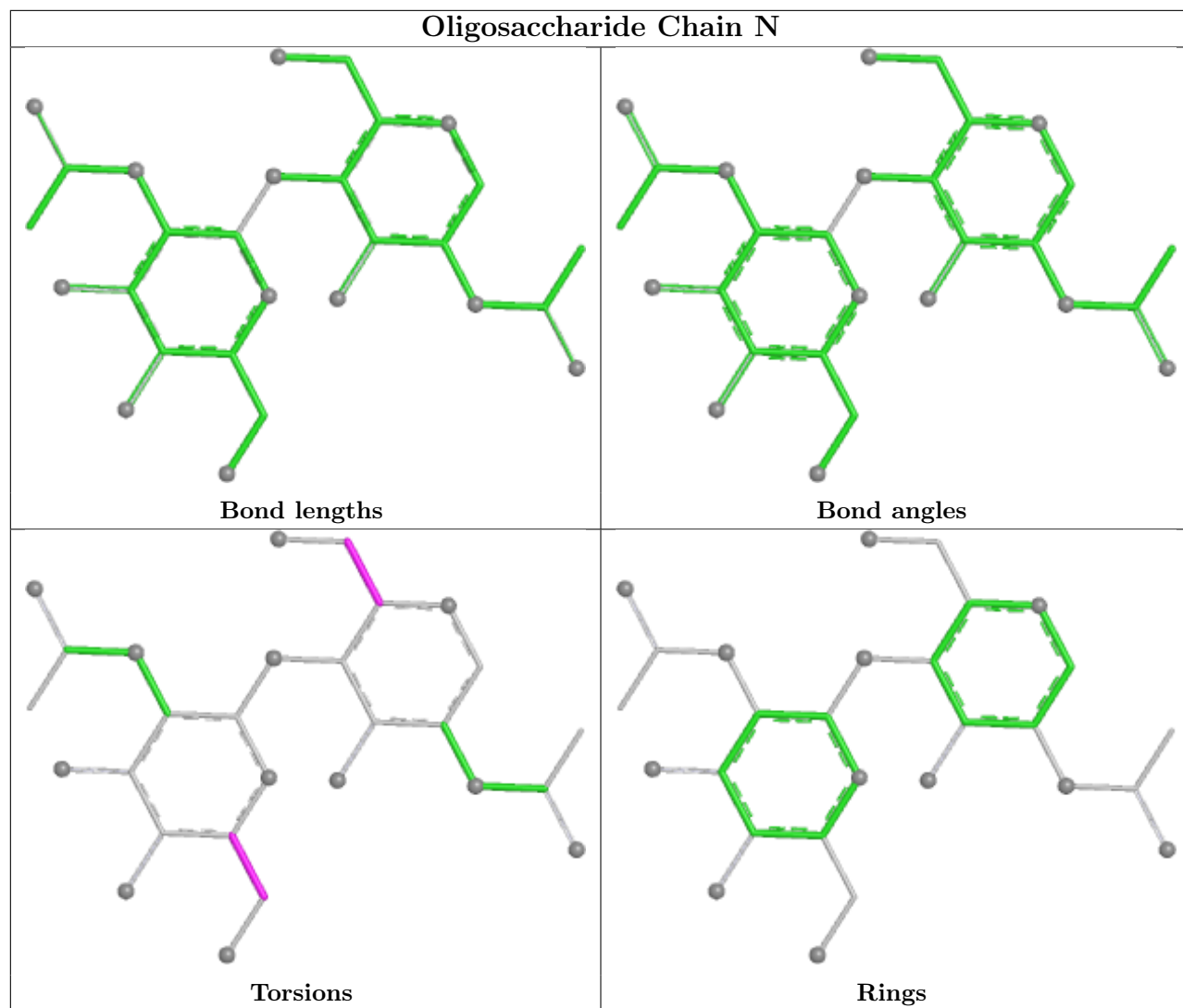


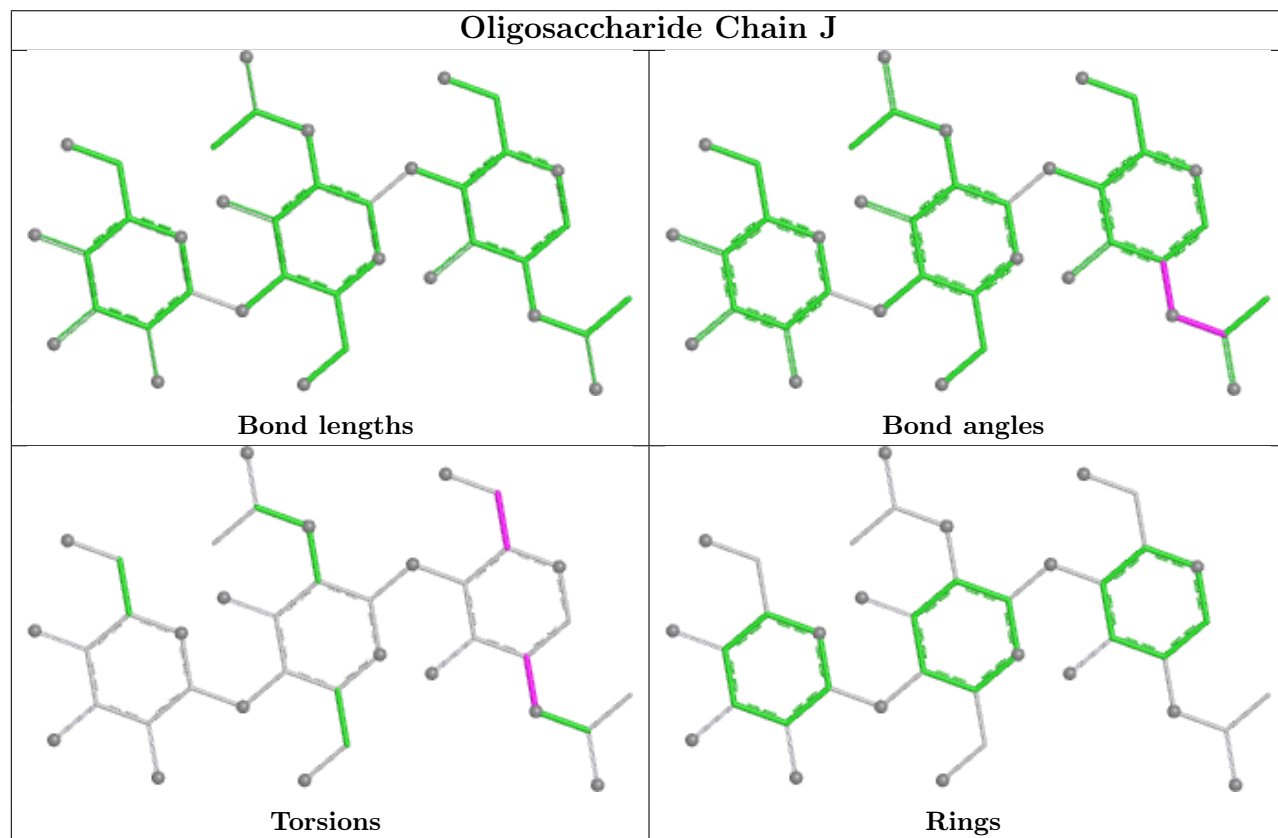
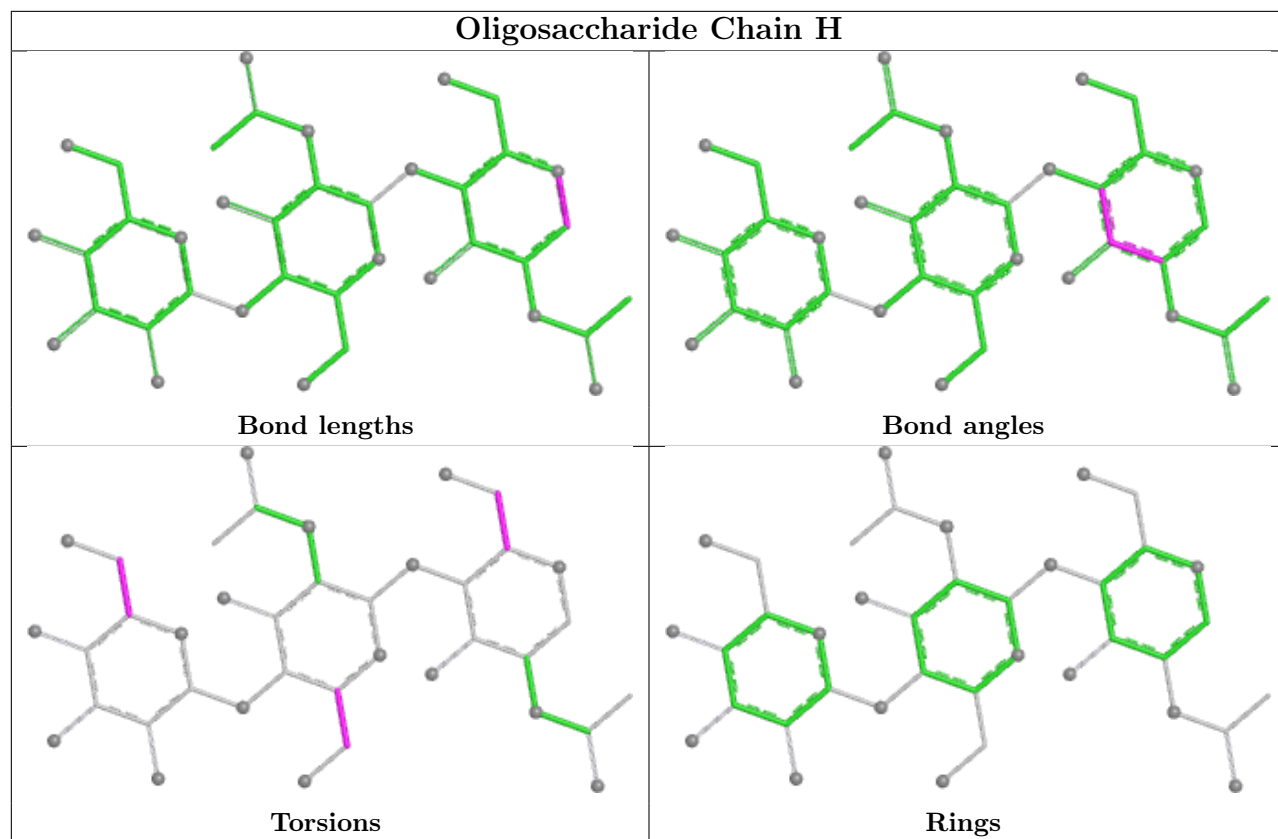


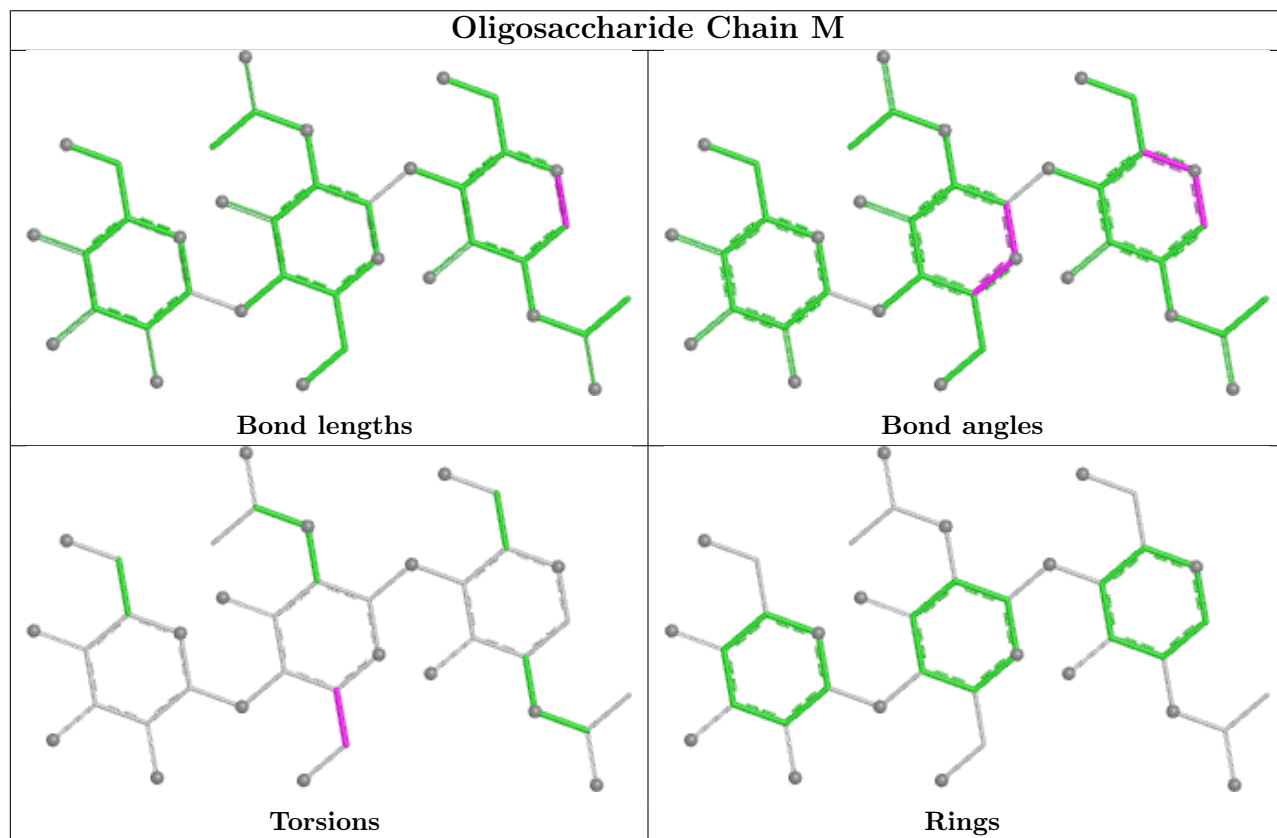
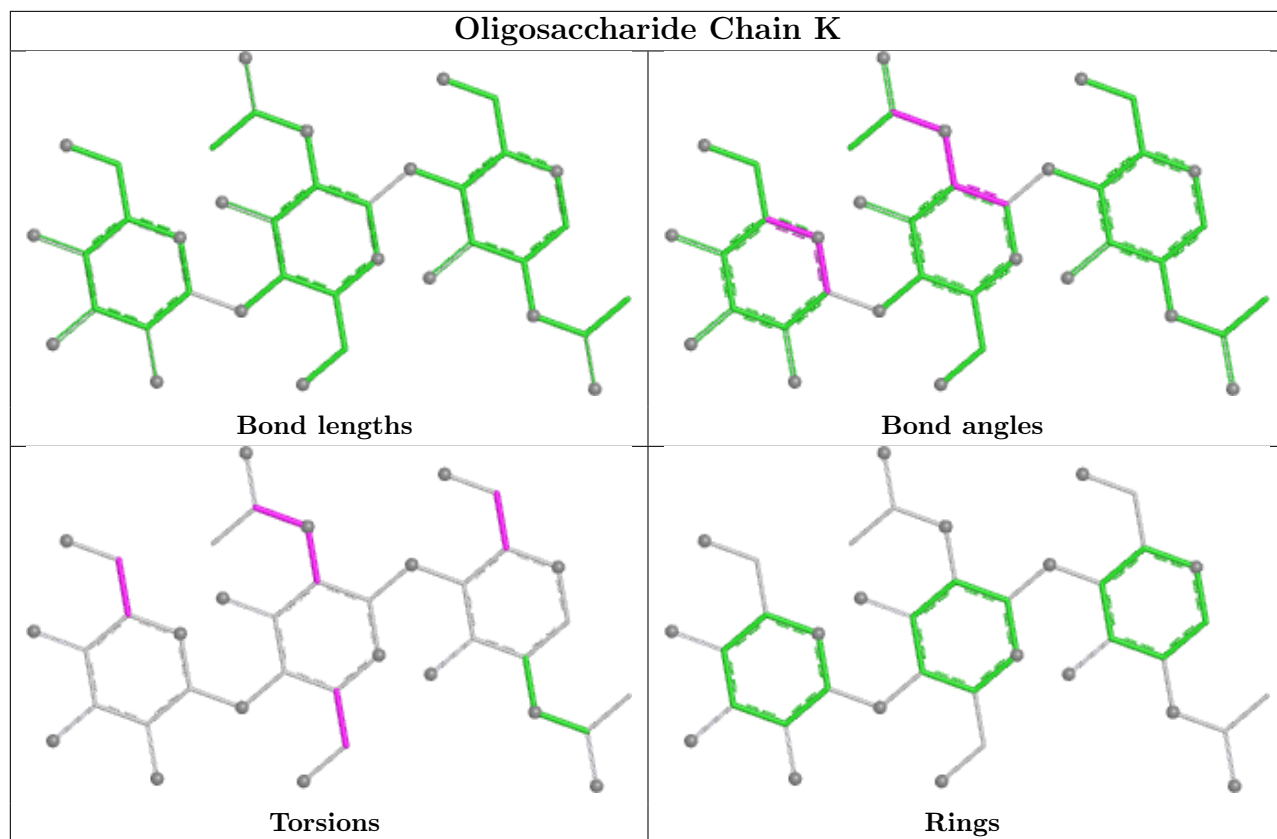












5.6 Ligand geometry

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	C	902	1	14,14,15	0.22	0	17,19,21	0.54	0
5	NAG	D	901	2	14,14,15	0.25	0	17,19,21	0.54	0
5	NAG	A	901	1	14,14,15	0.22	0	17,19,21	0.42	0
5	NAG	A	902	-	14,14,15	0.62	0	17,19,21	0.78	1 (5%)
5	NAG	A	903	1	14,14,15	0.23	0	17,19,21	0.48	0
5	NAG	C	901	1	14,14,15	0.23	0	17,19,21	0.41	0
5	NAG	A	904	-	14,14,15	0.21	0	17,19,21	0.43	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
5	NAG	C	902	1	-	4/6/23/26	0/1/1/1
5	NAG	D	901	2	-	4/6/23/26	0/1/1/1
5	NAG	A	901	1	-	2/6/23/26	0/1/1/1
5	NAG	A	902	-	-	3/6/23/26	0/1/1/1
5	NAG	A	903	1	-	2/6/23/26	0/1/1/1
5	NAG	C	901	1	-	2/6/23/26	0/1/1/1
5	NAG	A	904	-	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	902	NAG	C1-O5-C5	2.91	116.08	112.19

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	901	NAG	C4-C5-C6-O6
5	A	903	NAG	O5-C5-C6-O6
5	D	901	NAG	O5-C5-C6-O6
5	C	901	NAG	O5-C5-C6-O6
5	A	901	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	901	NAG	1	0
5	A	902	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

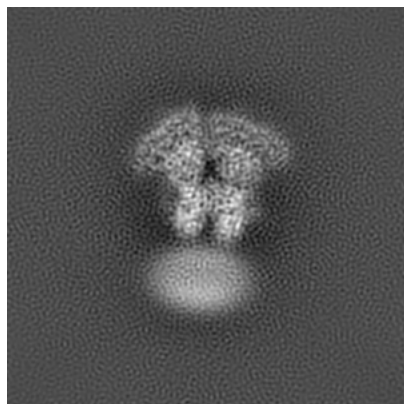
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-27953. These allow visual inspection of the internal detail of the map and identification of artifacts.

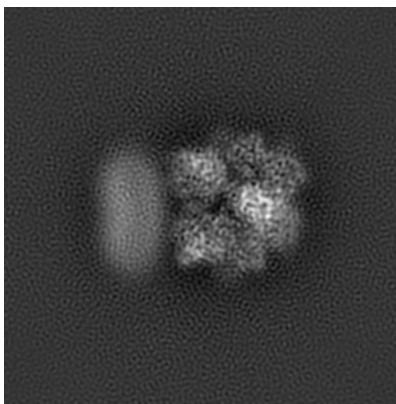
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

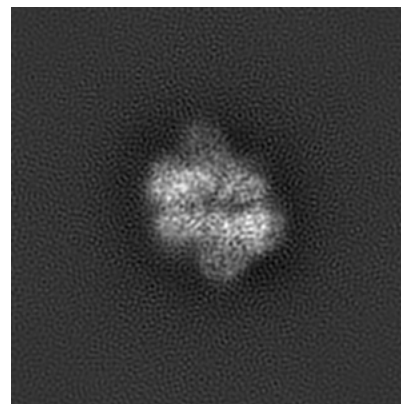
6.1.1 Primary map



X

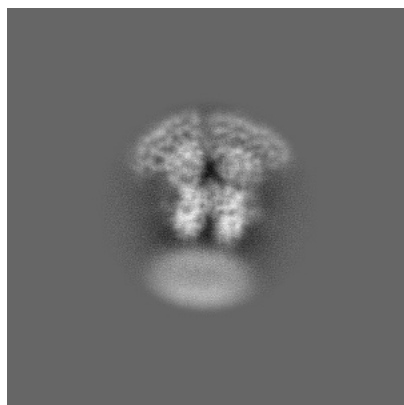


Y

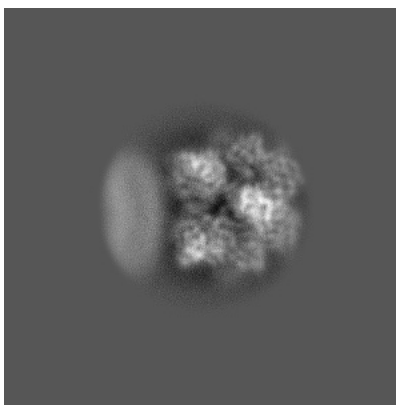


Z

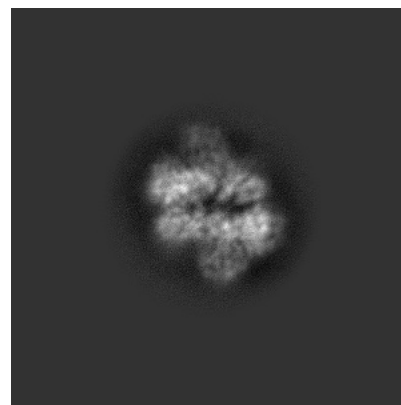
6.1.2 Raw map



X



Y

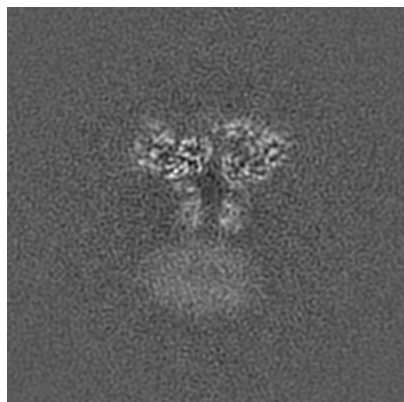


Z

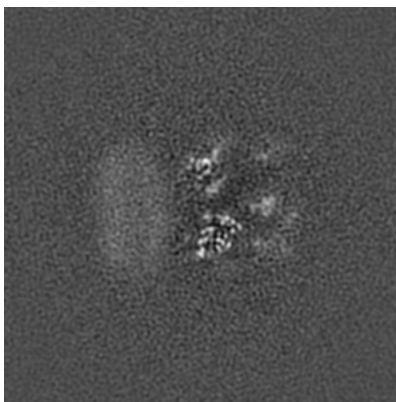
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

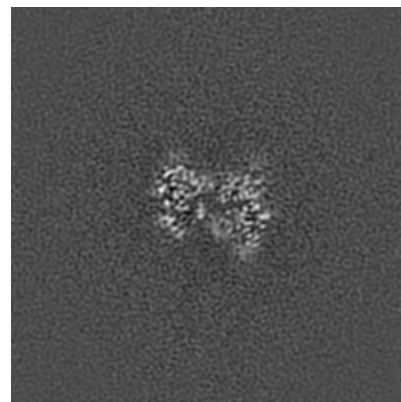
6.2.1 Primary map



X Index: 200

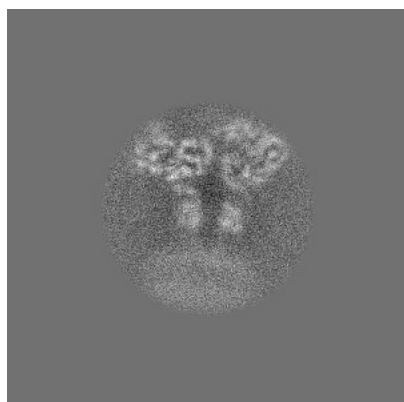


Y Index: 200

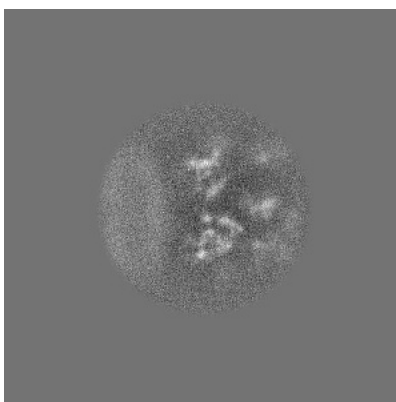


Z Index: 200

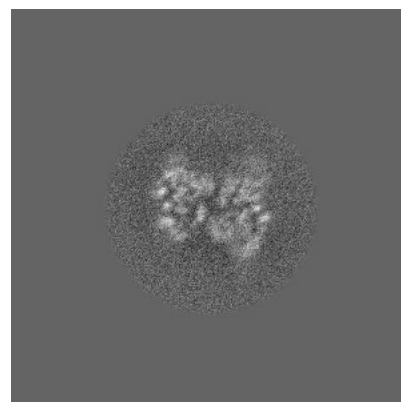
6.2.2 Raw map



X Index: 200



Y Index: 200

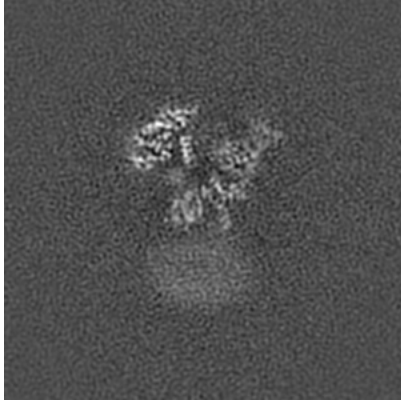


Z Index: 200

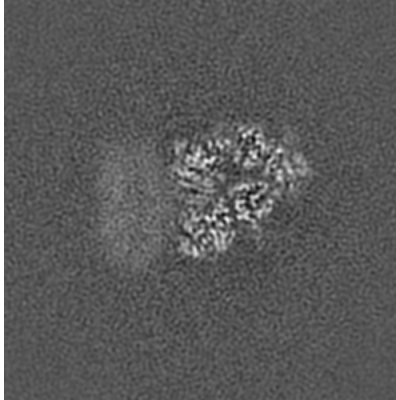
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

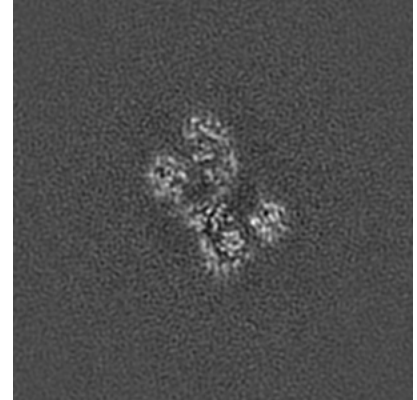
6.3.1 Primary map



X Index: 215

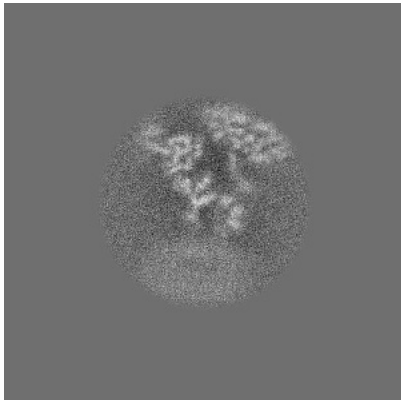


Y Index: 184

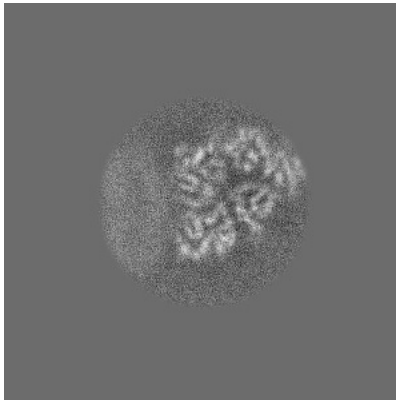


Z Index: 253

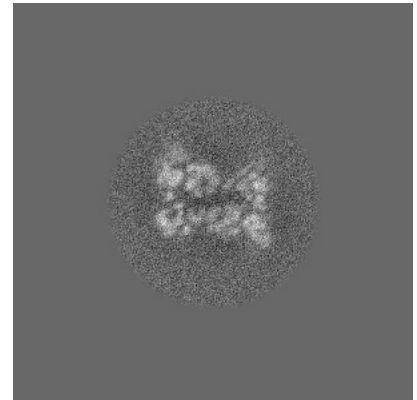
6.3.2 Raw map



X Index: 189



Y Index: 186

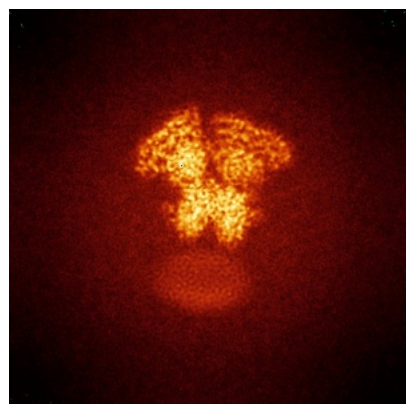


Z Index: 190

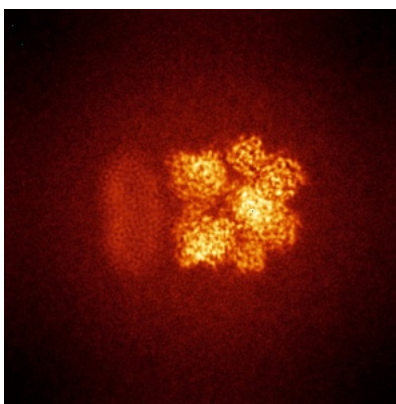
The images above show the largest variance slices of the map in three orthogonal directions.

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

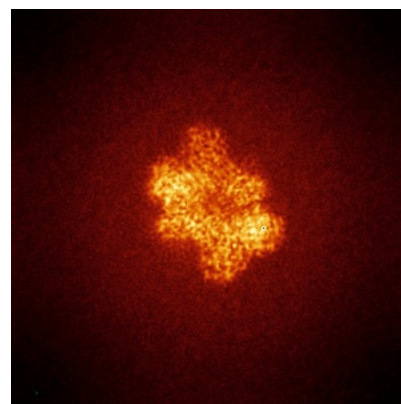
6.4.1 Primary map



X

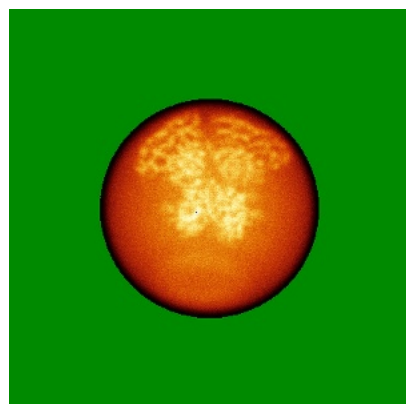


Y

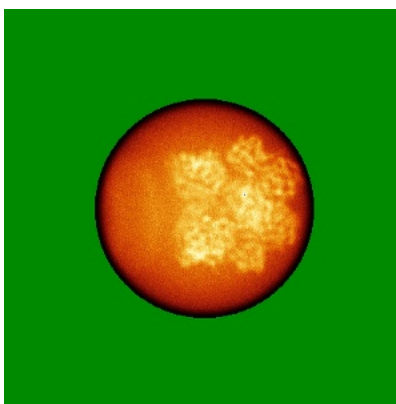


Z

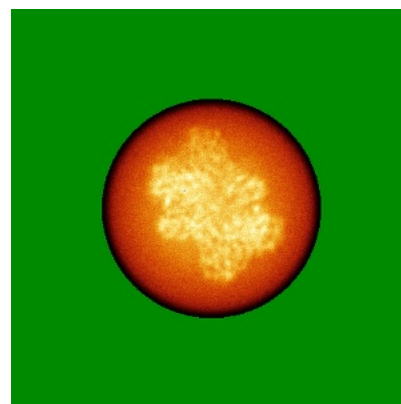
6.4.2 Raw map



X



Y

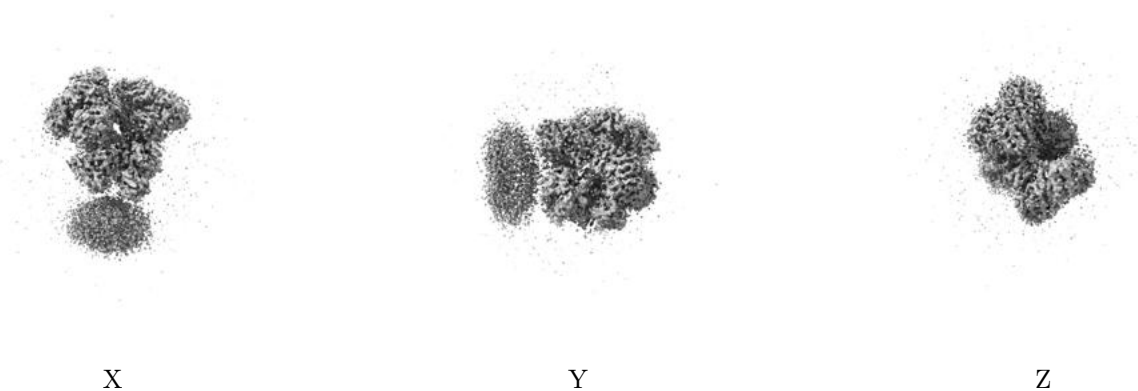


Z

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.

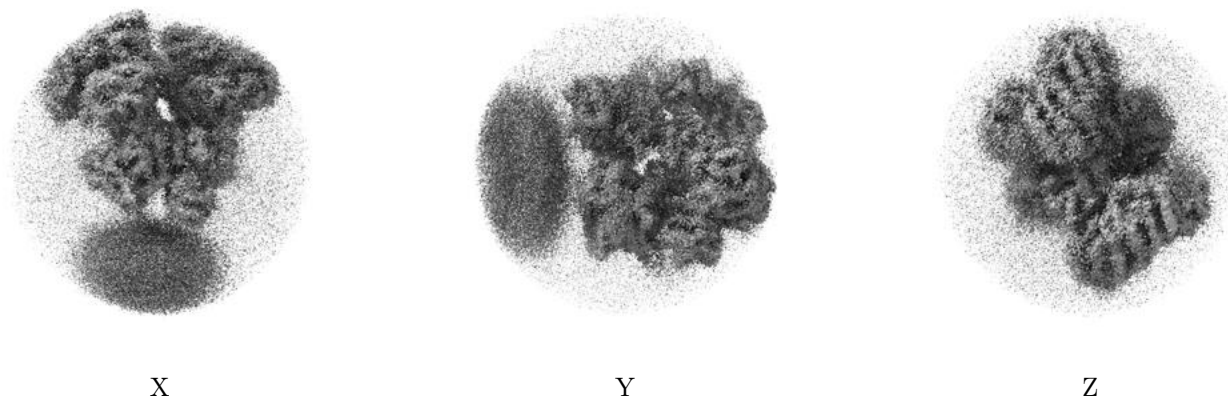
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 1.85. 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.

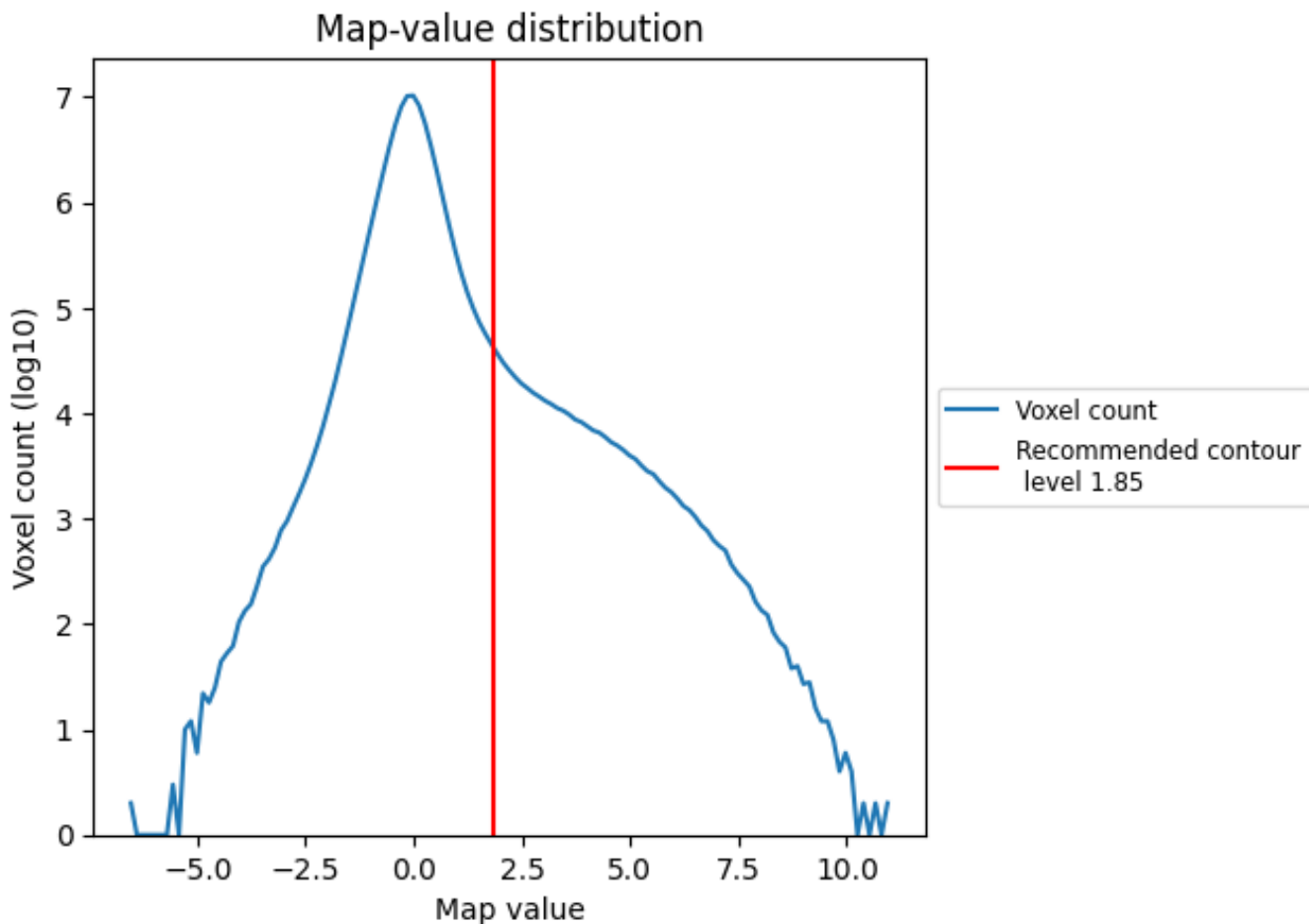
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

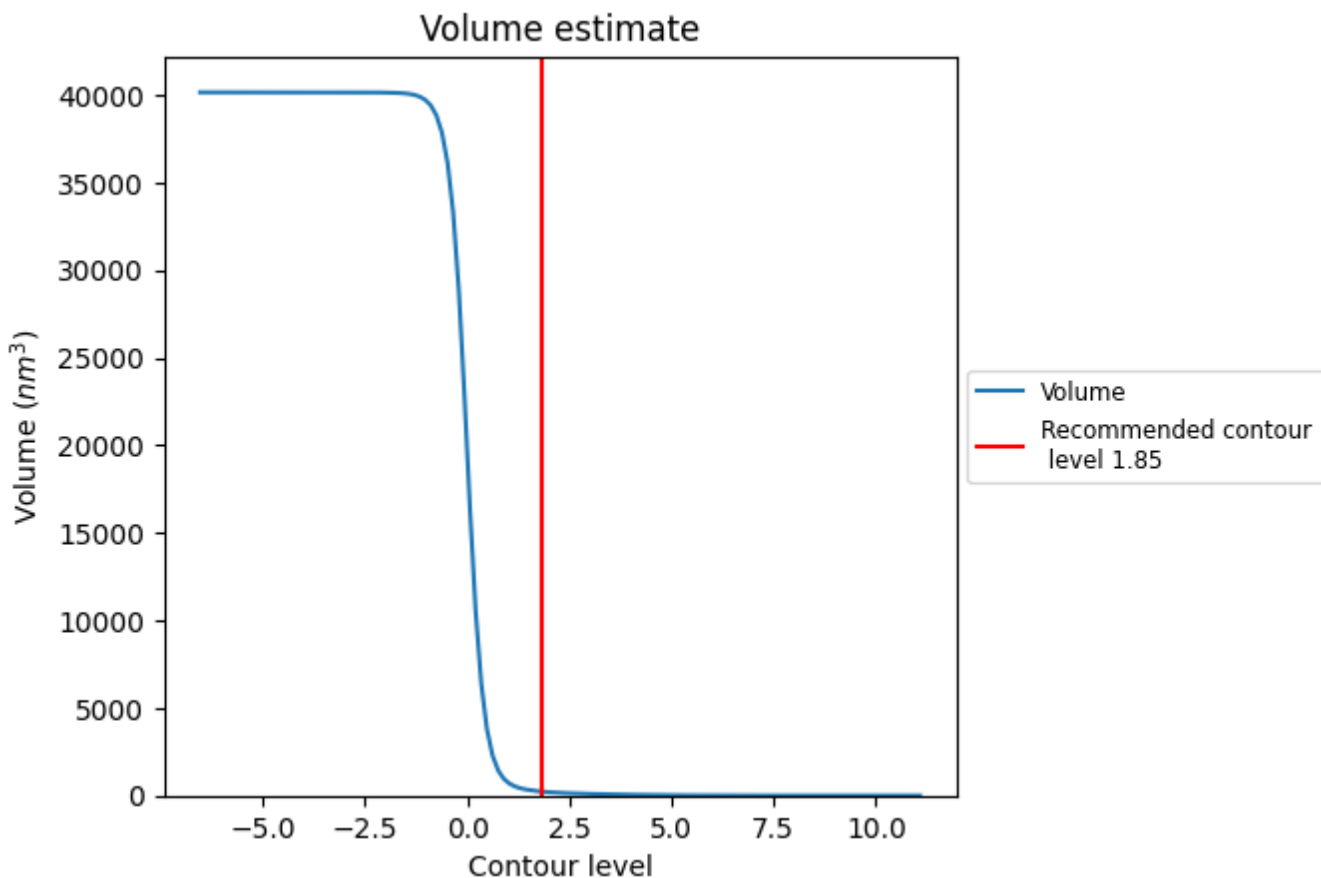
This section contains the results of statistical analysis of the map.

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.

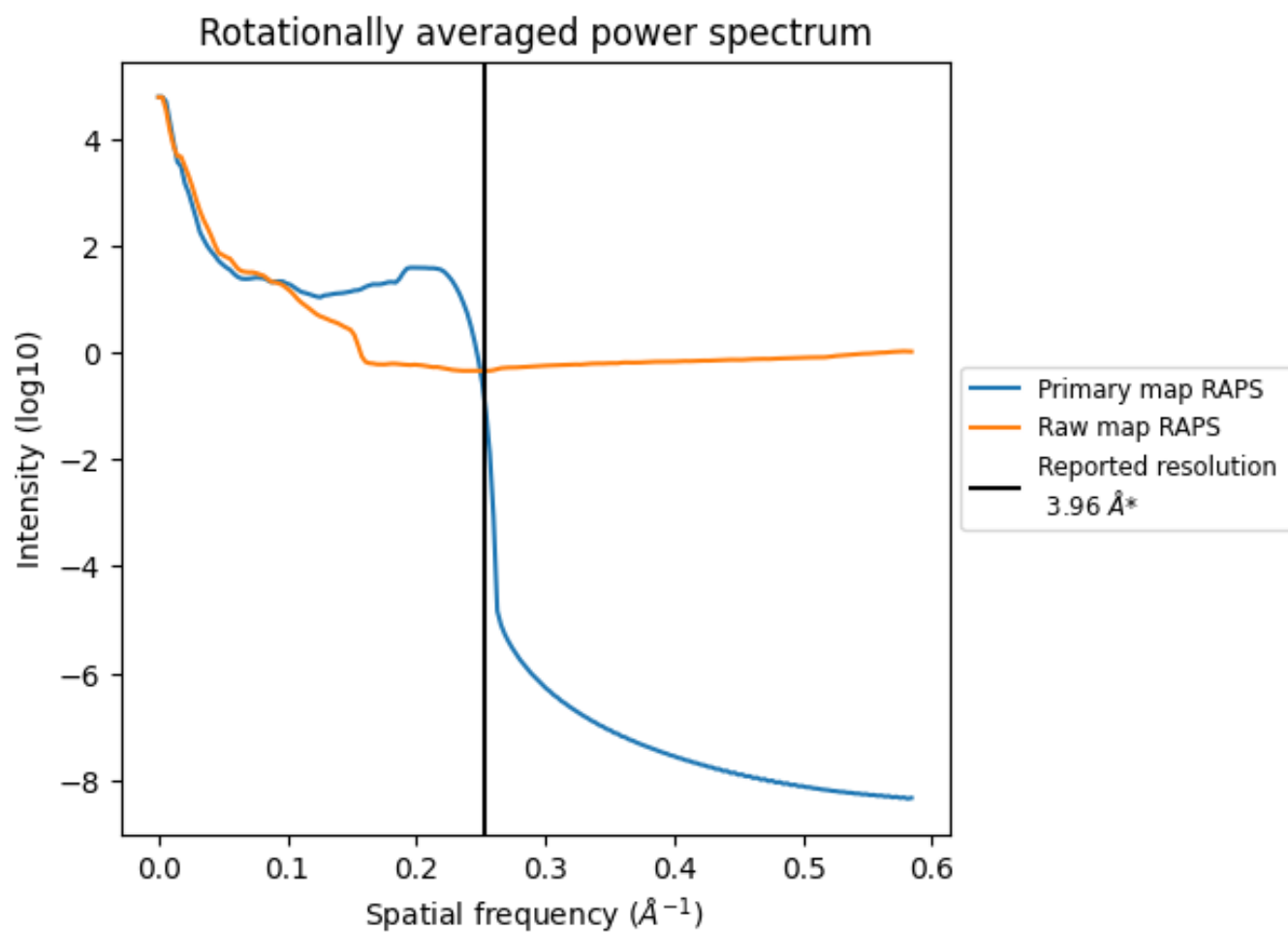
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 222 nm³; this corresponds to an approximate mass of 200 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 [i](#)

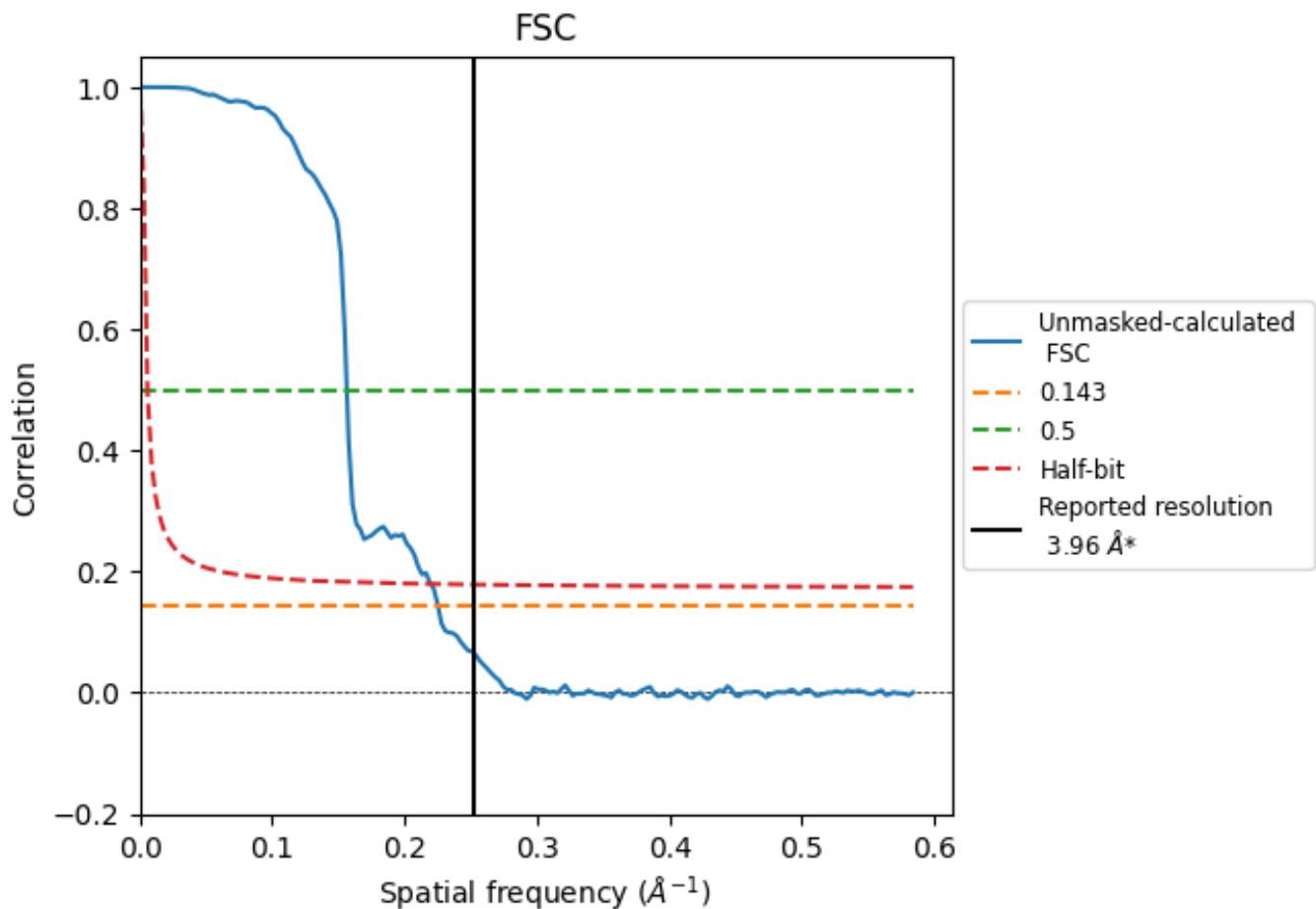


*Reported resolution corresponds to spatial frequency of 0.253 Å⁻¹

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 0.253 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.96	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.44	6.40	4.55

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.44 differs from the reported value 3.96 by more than 10 %

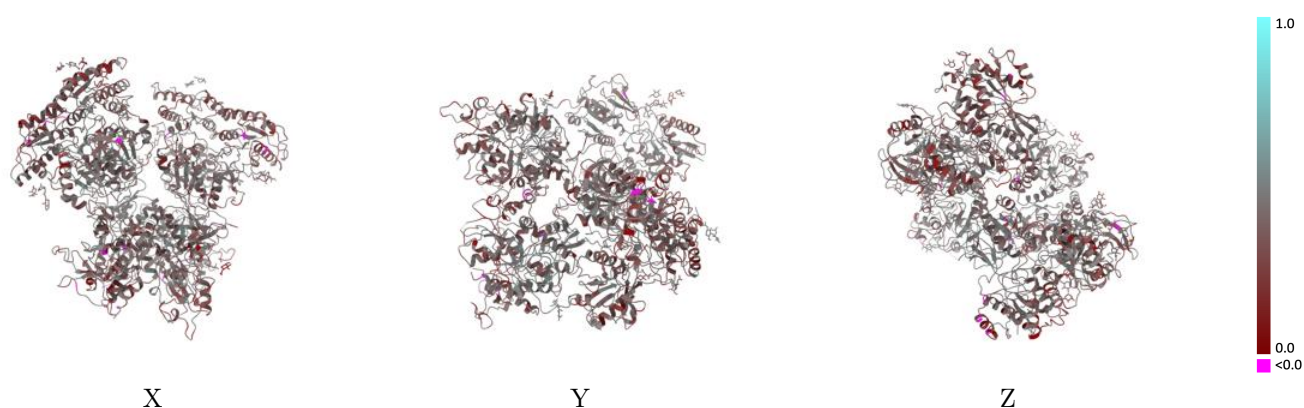
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-27953 and PDB model 8E92. Per-residue inclusion information can be found in section 3 on page 10.

9.1 Map-model overlay [i](#)

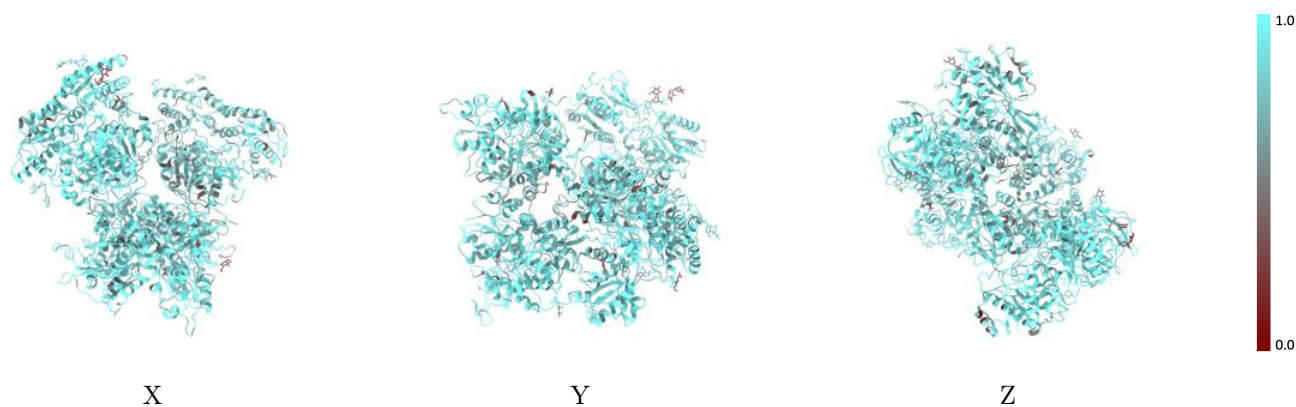
This section was not generated.

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



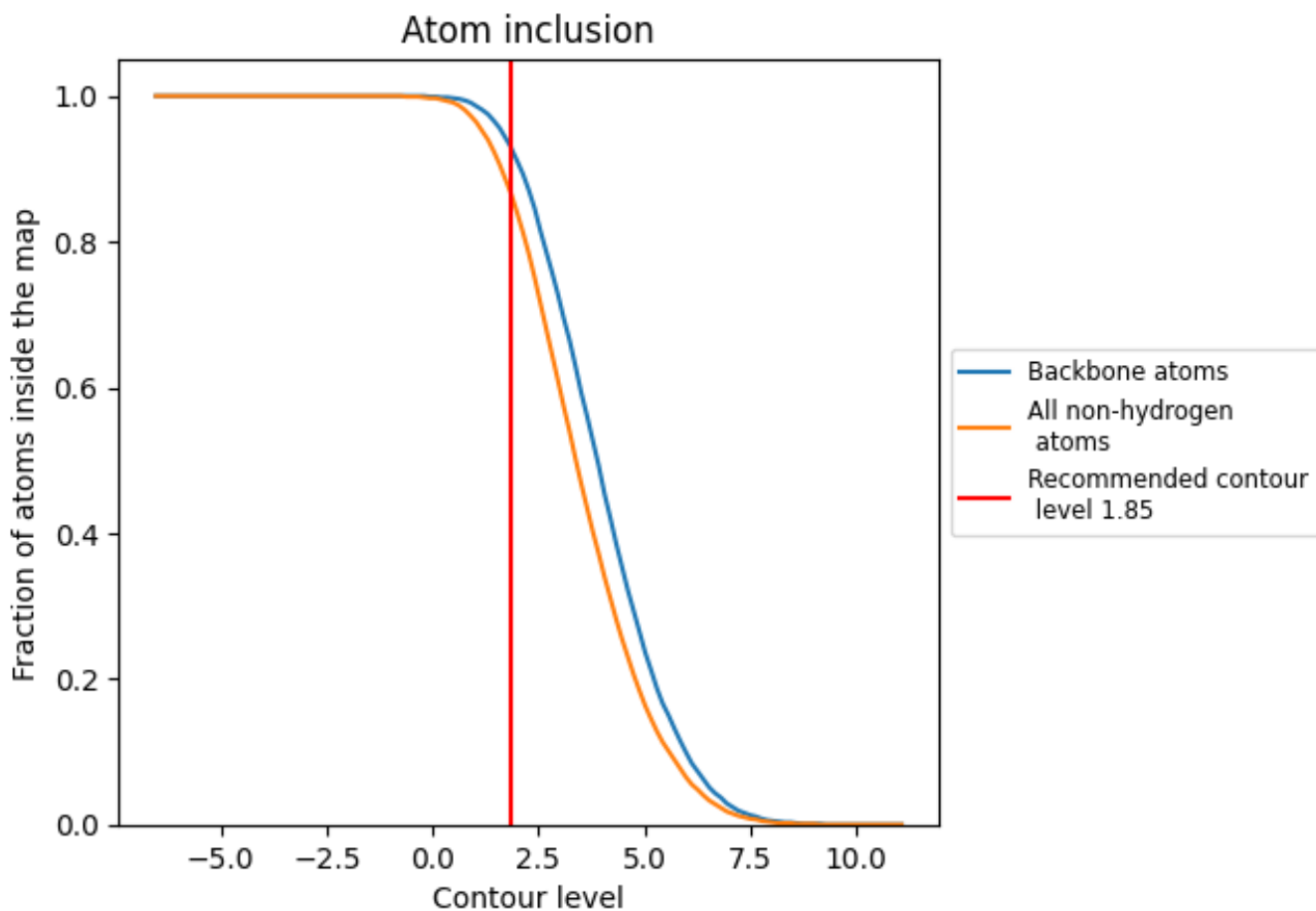
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (1.85).































9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 87% 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 (1.85) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8660	 0.3970
A	 0.8780	 0.3930
B	 0.8220	 0.3850
C	 0.8870	 0.3960
D	 0.8920	 0.4200
E	 0.6070	 0.3820
F	 0.7860	 0.4530
G	 0.3930	 0.2860
H	 0.6150	 0.3440
I	 0.6430	 0.2500
J	 0.8720	 0.3170
K	 0.4870	 0.3050
L	 0.6790	 0.3120
M	 0.4870	 0.3810
N	 0.8570	 0.4020

