



## Full wwPDB EM Validation Report ⓘ

Nov 20, 2022 – 03:14 AM EST

PDB ID : 7M3G  
EMDB ID : EMD-23654  
Title : Asymmetric Activation of the Calcium Sensing Receptor Homodimer  
Authors : Gao, Y.; Robertson, M.J.; Zhang, C.; Meyerowitz, J.G.; Panova, O.; Skiniotis, G.  
Deposited on : 2021-03-18  
Resolution : 2.50 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev43  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
MolProbity : 4.02b-467  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

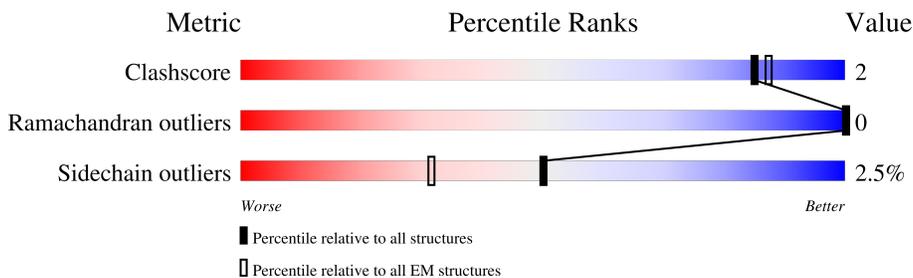
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 2.50 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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  $\geq 3$ , 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	902	
1	B	902	
2	C	9	
2	D	9	
3	E	2	
3	F	2	
3	G	2	
3	H	2	

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Mol	Chain	Length	Quality of chain
3	I	2	

## 2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 13298 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 Extracellular calcium-sensing receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	821	6485	4212	1056	1180	37	0	0
1	B	804	6375	4144	1033	1162	36	0	0

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP P41180
A	-6	LYS	-	expression tag	UNP P41180
A	-5	THR	-	expression tag	UNP P41180
A	-4	ILE	-	expression tag	UNP P41180
A	-3	ILE	-	expression tag	UNP P41180
A	-2	ALA	-	expression tag	UNP P41180
A	-1	LEU	-	expression tag	UNP P41180
A	0	SER	-	expression tag	UNP P41180
A	1	TYR	-	expression tag	UNP P41180
A	2	ILE	-	expression tag	UNP P41180
A	3	PHE	-	expression tag	UNP P41180
A	4	CYS	-	expression tag	UNP P41180
A	5	LEU	-	expression tag	UNP P41180
A	6	VAL	-	expression tag	UNP P41180
A	7	PHE	-	expression tag	UNP P41180
A	8	ALA	-	expression tag	UNP P41180
A	9	ASP	-	expression tag	UNP P41180
A	10	TYR	-	expression tag	UNP P41180
A	11	LYS	-	expression tag	UNP P41180
A	12	ASP	-	expression tag	UNP P41180
A	13	ASP	-	expression tag	UNP P41180
A	14	ASP	-	expression tag	UNP P41180
A	15	ASP	-	expression tag	UNP P41180
A	16	LYS	-	expression tag	UNP P41180
A	17	ALA	-	expression tag	UNP P41180
A	18	ALA	-	expression tag	UNP P41180

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Chain	Residue	Modelled	Actual	Comment	Reference
A	19	ALA	-	expression tag	UNP P41180
B	-7	MET	-	initiating methionine	UNP P41180
B	-6	LYS	-	expression tag	UNP P41180
B	-5	THR	-	expression tag	UNP P41180
B	-4	ILE	-	expression tag	UNP P41180
B	-3	ILE	-	expression tag	UNP P41180
B	-2	ALA	-	expression tag	UNP P41180
B	-1	LEU	-	expression tag	UNP P41180
B	0	SER	-	expression tag	UNP P41180
B	1	TYR	-	expression tag	UNP P41180
B	2	ILE	-	expression tag	UNP P41180
B	3	PHE	-	expression tag	UNP P41180
B	4	CYS	-	expression tag	UNP P41180
B	5	LEU	-	expression tag	UNP P41180
B	6	VAL	-	expression tag	UNP P41180
B	7	PHE	-	expression tag	UNP P41180
B	8	ALA	-	expression tag	UNP P41180
B	9	ASP	-	expression tag	UNP P41180
B	10	TYR	-	expression tag	UNP P41180
B	11	LYS	-	expression tag	UNP P41180
B	12	ASP	-	expression tag	UNP P41180
B	13	ASP	-	expression tag	UNP P41180
B	14	ASP	-	expression tag	UNP P41180
B	15	ASP	-	expression tag	UNP P41180
B	16	LYS	-	expression tag	UNP P41180
B	17	ALA	-	expression tag	UNP P41180
B	18	ALA	-	expression tag	UNP P41180
B	19	ALA	-	expression tag	UNP P41180

- Molecule 2 is a protein (with D amino acids) called etelcalcetide.

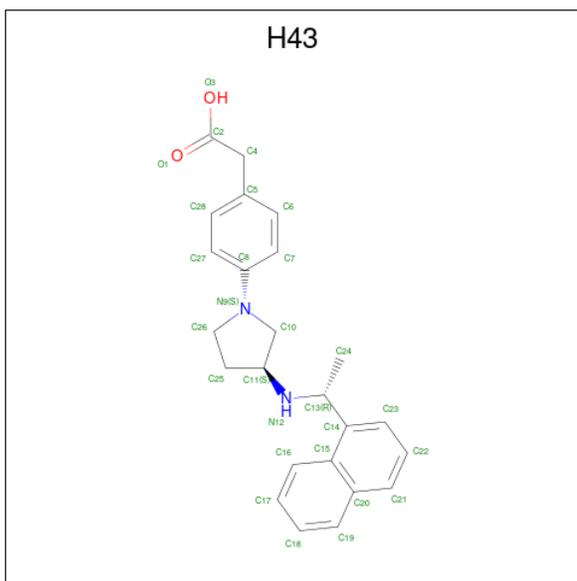
Mol	Chain	Residues	Atoms				AltConf	Trace	
2	C	9	Total	C	N	O	S	0	1
			64	35	20	8	1		
2	D	9	Total	C	N	O	S	0	1
			64	35	20	8	1		

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



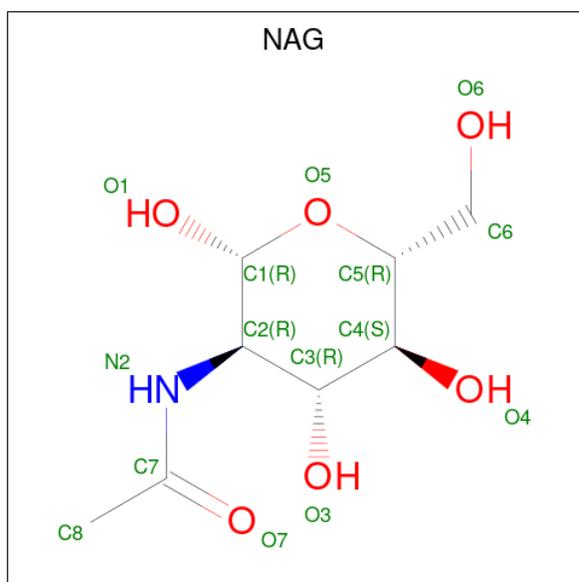
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
3	E	2	28	16	2	10	0	0
3	F	2	28	16	2	10	0	0
3	G	2	28	16	2	10	0	0
3	H	2	28	16	2	10	0	0
3	I	2	28	16	2	10	0	0

- Molecule 4 is 2-[4-[(3S)-3-[(1R)-1-naphthalen-1-ylethyl]amino]pyrrolidin-1-yl]phenyl]ethanoic acid (three-letter code: H43) (formula: C<sub>24</sub>H<sub>26</sub>N<sub>2</sub>O<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



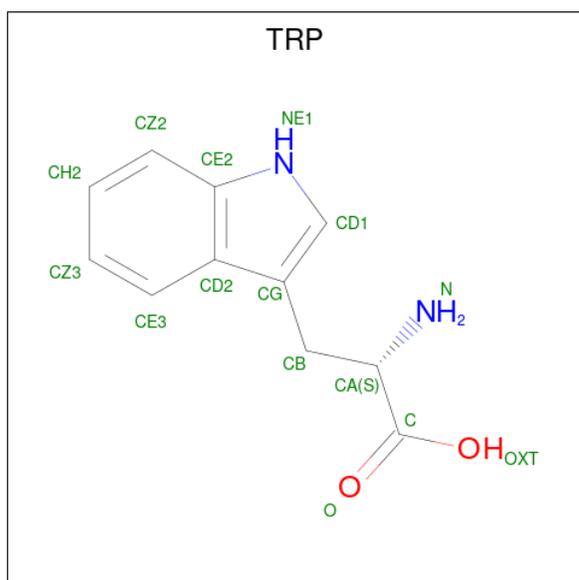
Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
4	A	1	28	24	2	2	0
4	B	1	28	24	2	2	0

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
5	A	1	42	24	3	15	0
5	A	1	42	24	3	15	0
5	A	1	42	24	3	15	0
5	B	1	28	16	2	10	0
5	B	1	28	16	2	10	0

- Molecule 6 is TRYPTOPHAN (three-letter code: TRP) (formula:  $C_{11}H_{12}N_2O_2$ ).

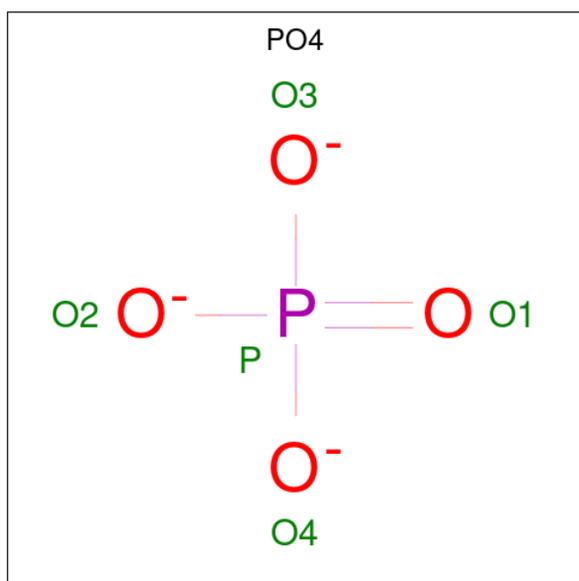


Mol	Chain	Residues	Atoms			AltConf	
6	A	1	Total	C	N	O	0
			15	11	2	2	
6	B	1	Total	C	N	O	0
			15	11	2	2	

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
7	A	3	Total	Ca	0
			3	3	
7	B	1	Total	Ca	0
			1	1	

- Molecule 8 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			AltConf
8	A	1	Total	O	P	0
			5	4	1	
8	B	1	Total	O	P	0
			5	4	1	



There are no outlier residues recorded for this chain.

- Molecule 2: etelcalcetide

Chain D:  78% 22%



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

Chain E:  100%



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

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 H:  50% 50%



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

Chain I:  50% 50%



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	482939	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 ( $e^-/\text{\AA}^2$ )	64.9	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	14.580	Depositor
Minimum map value	-1.291	Depositor
Average map value	0.007	Depositor
Map value standard deviation	0.145	Depositor
Recommended contour level	0.9	Depositor
Map size ( $\text{\AA}$ )	399.6, 399.6, 399.6	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.11, 1.11, 1.11	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: PO4, DCY, NAG, NH2, DAL, ACE, H43, CA, DAR

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	0/6648	0.56	1/9022 (0.0%)
1	B	0.34	0/6537	0.58	0/8871
All	All	0.34	0/13185	0.57	1/17893 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	480	ASP	CB-CG-OD2	5.22	123.00	118.30

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	6485	0	6346	32	0
1	B	6375	0	6241	25	0
2	C	64	0	64	0	0
2	D	64	0	67	2	0
3	E	28	0	25	0	0
3	F	28	0	25	0	0
3	G	28	0	25	0	0
3	H	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	I	28	0	25	0	0
4	A	28	0	0	0	0
4	B	28	0	0	0	0
5	A	42	0	39	1	0
5	B	28	0	26	1	0
6	A	15	0	9	0	0
6	B	15	0	9	0	0
7	A	3	0	0	0	0
7	B	1	0	0	0	0
8	A	5	0	0	0	0
8	B	5	0	0	1	0
All	All	13298	0	12926	57	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 2.

All (57) 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:91:LEU:HD12	1:A:91:LEU:O	1.97	0.64
1:A:251:GLU:HG2	2:D:4:DAR:HD2	1.81	0.63
1:A:480:ASP:OD1	1:A:484:ASP:N	2.29	0.59
1:A:69:ARG:NH1	1:A:407:PRO:O	2.36	0.58
1:A:831:LYS:C	1:A:832:PHE:CD2	2.77	0.57
1:B:69:ARG:NH1	1:B:407:PRO:O	2.37	0.57
1:B:213:ALA:HB2	1:B:223:ILE:HG13	1.85	0.57
1:B:696:LEU:HD13	1:B:787:CYS:CB	2.36	0.56
1:A:673:GLN:OE1	1:A:675:TRP:NE1	2.38	0.56
1:B:69:ARG:NH2	8:B:906:PO4:O4	2.39	0.56
1:A:831:LYS:O	1:A:832:PHE:CD2	2.59	0.56
1:A:831:LYS:O	1:A:832:PHE:CG	2.58	0.55
1:B:227:ARG:NH1	1:B:231:GLU:OE2	2.39	0.55
1:A:661:CYS:SG	1:A:845:SER:OG	2.65	0.55
1:A:396:THR:HG23	1:A:398:ASP:H	1.72	0.54
1:B:696:LEU:HD13	1:B:787:CYS:HA	1.90	0.54
1:B:163:PRO:HG3	1:B:461:LEU:HB2	1.90	0.53
1:B:640:THR:O	1:B:644:LYS:NZ	2.43	0.52
1:B:696:LEU:HD13	1:B:787:CYS:HB2	1.91	0.51
1:A:647:ASN:O	1:A:647:ASN:ND2	2.40	0.51
1:A:27:GLN:HG3	1:A:96:ARG:HG2	1.93	0.50
1:A:548:ALA:O	1:A:574:SER:OG	2.28	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:551:ARG:NH2	1:A:552:LYS:O	2.45	0.49
1:A:91:LEU:HD12	1:A:91:LEU:C	2.33	0.48
1:A:748:PRO:HB3	1:A:767:GLU:HA	1.95	0.48
1:B:117:GLN:HG2	1:B:135:ILE:HB	1.96	0.48
1:A:478:THR:HG21	5:A:902:NAG:H82	1.95	0.48
1:A:51:LEU:HD12	1:A:55:PRO:HG3	1.97	0.47
1:A:737:VAL:HA	1:A:740:VAL:HG12	1.97	0.46
1:A:755:GLU:N	1:A:755:GLU:OE2	2.49	0.46
1:B:349:LYS:HG3	1:B:360:LEU:HD23	1.99	0.45
1:A:831:LYS:C	1:A:832:PHE:CG	2.90	0.45
1:A:570:ASP:O	1:A:591:SER:OG	2.33	0.44
1:B:478:THR:HG21	5:B:902:NAG:H82	2.00	0.44
1:A:87:LEU:HD12	1:A:87:LEU:HA	1.90	0.43
1:A:322:LEU:O	1:A:415:ARG:NH1	2.43	0.43
1:B:212:ILE:HD11	1:B:259:ILE:HD11	2.01	0.42
1:A:592:ASN:OD1	1:A:592:ASN:N	2.53	0.42
1:B:696:LEU:HD13	1:B:787:CYS:CA	2.49	0.42
1:B:589:PHE:HB3	1:B:598:CYS:HB3	2.02	0.42
1:B:422:LEU:O	1:B:426:SER:OG	2.34	0.42
1:A:36:GLY:HA3	1:A:97:ILE:HG23	2.02	0.41
1:B:588:ASP:N	1:B:588:ASP:OD1	2.53	0.41
1:B:587:ASP:OD1	1:B:587:ASP:N	2.53	0.41
1:B:108:LEU:HD23	1:B:108:LEU:HA	1.90	0.41
1:A:211:THR:HB	1:A:223:ILE:HD11	2.02	0.41
1:A:547:LEU:HD23	1:A:547:LEU:HA	1.95	0.41
1:A:32:ILE:HG23	1:A:139:ILE:HG12	2.02	0.41
1:A:293:TRP:HB2	1:A:317:THR:HG23	2.02	0.41
1:A:247:SER:OG	2:D:5:DAR:NH1	2.41	0.41
1:B:757:GLU:HG2	1:B:760:ILE:HG13	2.02	0.41
1:A:212:ILE:HD11	1:A:259:ILE:HD11	2.02	0.40
1:B:54:ARG:HA	1:B:55:PRO:HD3	1.95	0.40
1:B:229:GLU:OE1	1:B:233:ARG:NH1	2.55	0.40
1:B:276:LEU:HD23	1:B:305:ILE:HD13	2.04	0.40
1:B:544:ARG:HD3	1:B:544:ARG:HA	1.91	0.40
1:B:613:GLY:O	1:B:617:THR:OG1	2.33	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	813/902 (90%)	787 (97%)	26 (3%)	0	100	100
1	B	796/902 (88%)	773 (97%)	23 (3%)	0	100	100
All	All	1609/1804 (89%)	1560 (97%)	49 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	701/780 (90%)	680 (97%)	21 (3%)	41	68
1	B	694/780 (89%)	680 (98%)	14 (2%)	55	79
All	All	1395/1560 (89%)	1360 (98%)	35 (2%)	50	73

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	124	ASN
1	A	150	SER
1	A	270	PHE
1	A	463	HIS
1	A	516	LYS
1	A	570	ASP
1	A	576	GLU
1	A	583	ASN

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Mol	Chain	Res	Type
1	A	587	ASP
1	A	592	ASN
1	A	605	PHE
1	A	612	PHE
1	A	637	PHE
1	A	647	ASN
1	A	661	CYS
1	A	687	SER
1	A	691	CYS
1	A	790	PHE
1	A	793	LYS
1	A	795	ARG
1	A	820	SER
1	B	84	SER
1	B	253	GLN
1	B	270	PHE
1	B	441	ARG
1	B	499	GLU
1	B	588	ASP
1	B	594	ASN
1	B	723	LEU
1	B	743	LEU
1	B	752	ARG
1	B	788	PHE
1	B	789	PHE
1	B	790	PHE
1	B	845	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

14 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	DCY	D	2	2,1	4,5,6	0.71	0	1,5,7	0.88	0
2	DAR	D	4	2	9,10,11	0.41	0	5,11,13	0.24	0
2	DAL	C	3	2	3,4,5	0.60	0	2,4,6	0.87	0
2	DAR	D	5	2	9,10,11	0.42	0	5,11,13	0.45	0
2	DAR	D	6	2	9,10,11	0.37	0	5,11,13	0.34	0
2	DAL	C	7	2	3,4,5	0.63	0	2,4,6	0.64	0
2	DAL	D	7	2	3,4,5	0.55	0	2,4,6	0.86	0
2	DAR	C	6	2	9,10,11	0.56	0	5,11,13	0.48	0
2	DCY	C	2	2,1	4,5,6	0.99	0	1,5,7	1.27	0
2	DAR	C	8	2	9,10,11	0.43	0	5,11,13	0.23	0
2	DAR	D	8	2	9,10,11	0.39	0	5,11,13	0.24	0
2	DAR	C	4	2	9,10,11	0.42	0	5,11,13	0.25	0
2	DAR	C	5	2	9,10,11	0.44	0	5,11,13	0.41	0
2	DAL	D	3	2	3,4,5	0.61	0	2,4,6	0.88	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
2	DCY	D	2	2,1	-	0/1/4/6	-
2	DAR	D	4	2	-	4/8/9/11	-
2	DAL	C	3	2	-	0/0/2/4	-
2	DAR	D	5	2	-	1/8/9/11	-
2	DAR	D	6	2	-	0/8/9/11	-
2	DAL	C	7	2	-	0/0/2/4	-
2	DAL	D	7	2	-	0/0/2/4	-
2	DAR	C	6	2	-	0/8/9/11	-
2	DCY	C	2	2,1	-	0/1/4/6	-
2	DAR	C	8	2	-	0/8/9/11	-
2	DAR	D	8	2	-	2/8/9/11	-
2	DAR	C	4	2	-	1/8/9/11	-
2	DAR	C	5	2	-	1/8/9/11	-
2	DAL	D	3	2	-	0/0/2/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	4	DAR	N-CA-CB-CG
2	D	4	DAR	C-CA-CB-CG
2	D	8	DAR	O-C-CA-CB
2	D	4	DAR	NE-CD-CG-CB
2	D	8	DAR	NE-CD-CG-CB
2	C	5	DAR	CA-CB-CG-CD
2	D	5	DAR	CA-CB-CG-CD
2	D	4	DAR	CA-CB-CG-CD
2	C	4	DAR	NE-CD-CG-CB

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	4	DAR	1	0
2	D	5	DAR	1	0

## 5.5 Carbohydrates [i](#)

10 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	3,1	14,14,15	0.30	0	17,19,21	0.51	0
3	NAG	E	2	3	14,14,15	0.38	0	17,19,21	0.40	0
3	NAG	F	1	3,1	14,14,15	0.33	0	17,19,21	0.48	0
3	NAG	F	2	3	14,14,15	0.23	0	17,19,21	0.55	0
3	NAG	G	1	3,1	14,14,15	0.34	0	17,19,21	0.56	0
3	NAG	G	2	3	14,14,15	0.47	0	17,19,21	0.96	1 (5%)
3	NAG	H	1	3,1	14,14,15	0.32	0	17,19,21	0.71	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	H	2	3	14,14,15	0.38	0	17,19,21	0.52	0
3	NAG	I	1	3,1	14,14,15	0.23	0	17,19,21	0.63	1 (5%)
3	NAG	I	2	3	14,14,15	0.30	0	17,19,21	0.50	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	3,1	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	NAG	F	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	NAG	G	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	1/6/23/26	0/1/1/1
3	NAG	H	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1
3	NAG	I	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	2	NAG	C2-N2-C7	3.05	127.25	122.90
3	H	1	NAG	C1-O5-C5	2.51	115.59	112.19
3	I	1	NAG	C1-O5-C5	2.18	115.14	112.19

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	2	NAG	C4-C5-C6-O6
3	E	2	NAG	O5-C5-C6-O6
3	I	2	NAG	O5-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
3	I	2	NAG	C4-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6

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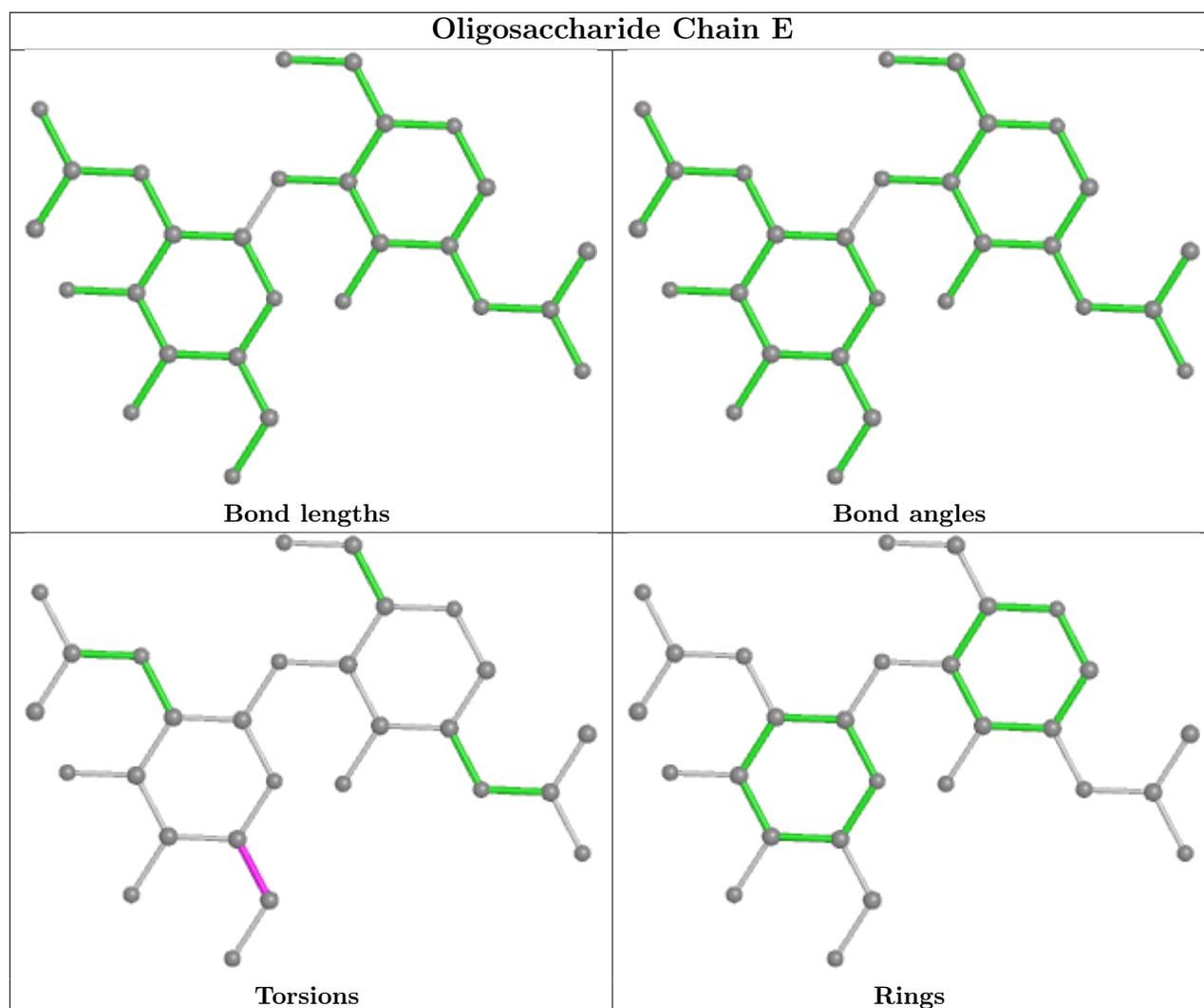
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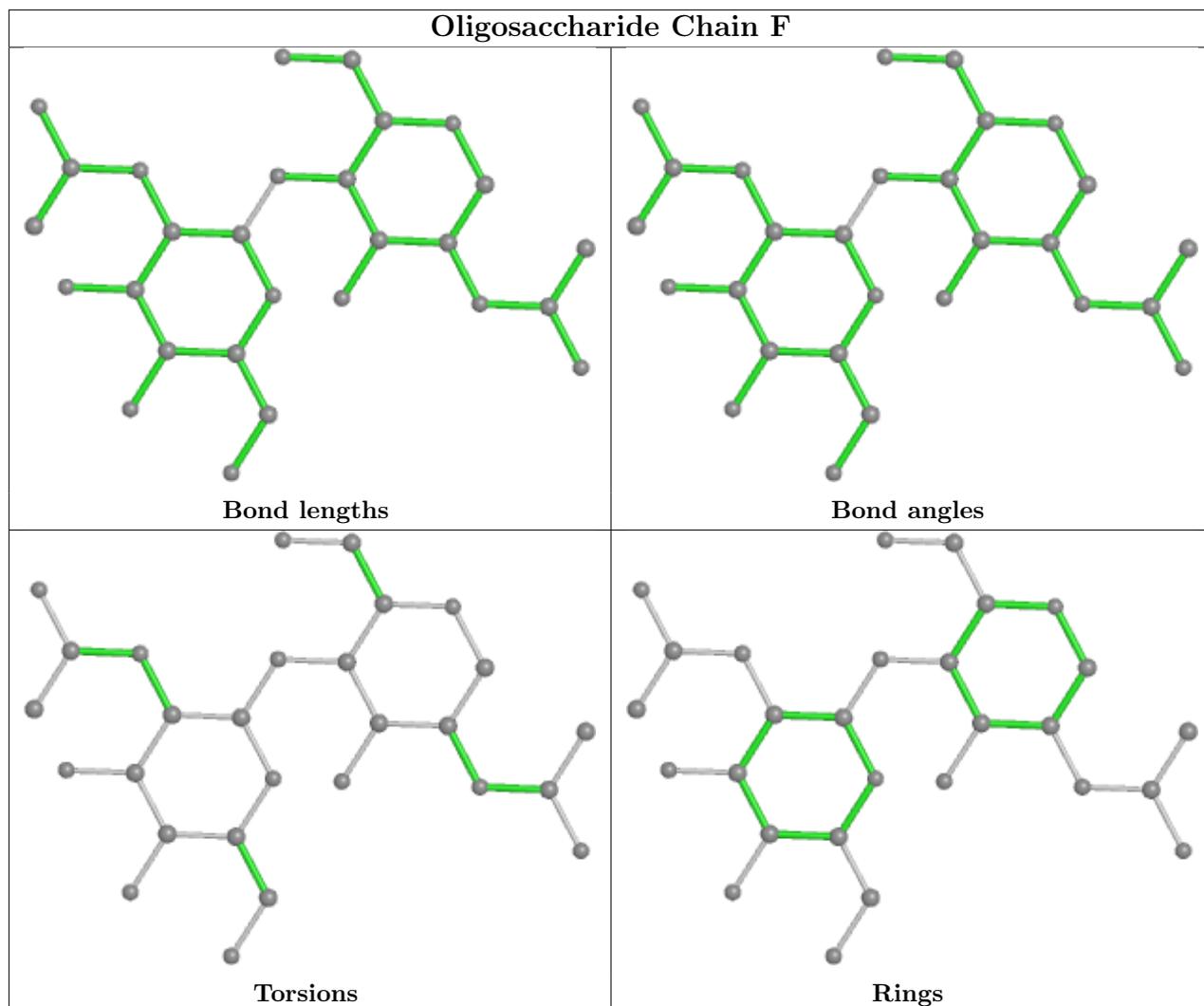
Mol	Chain	Res	Type	Atoms
3	H	1	NAG	C4-C5-C6-O6
3	G	2	NAG	C3-C2-N2-C7

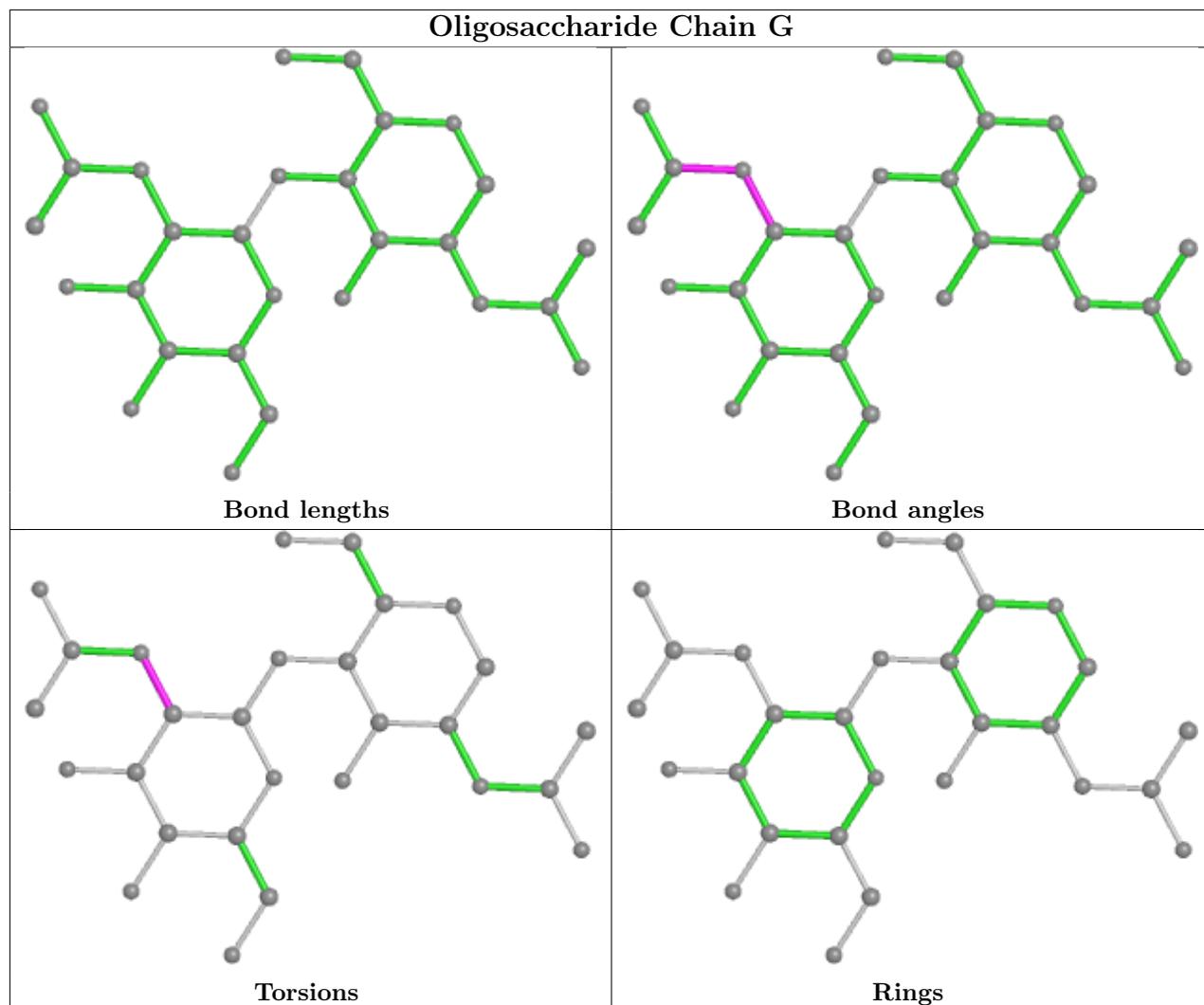
There are no ring outliers.

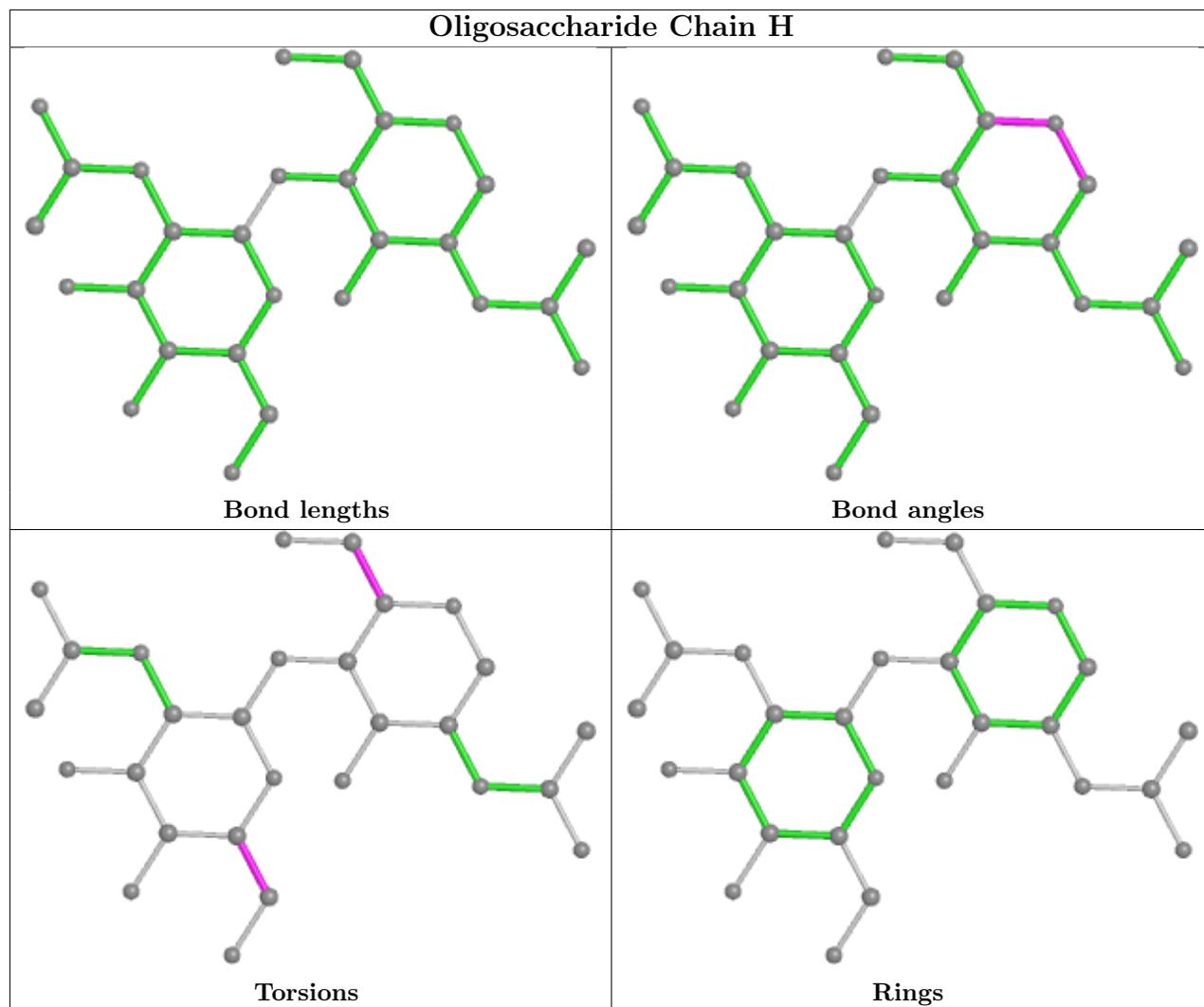
No monomer is involved in short contacts.

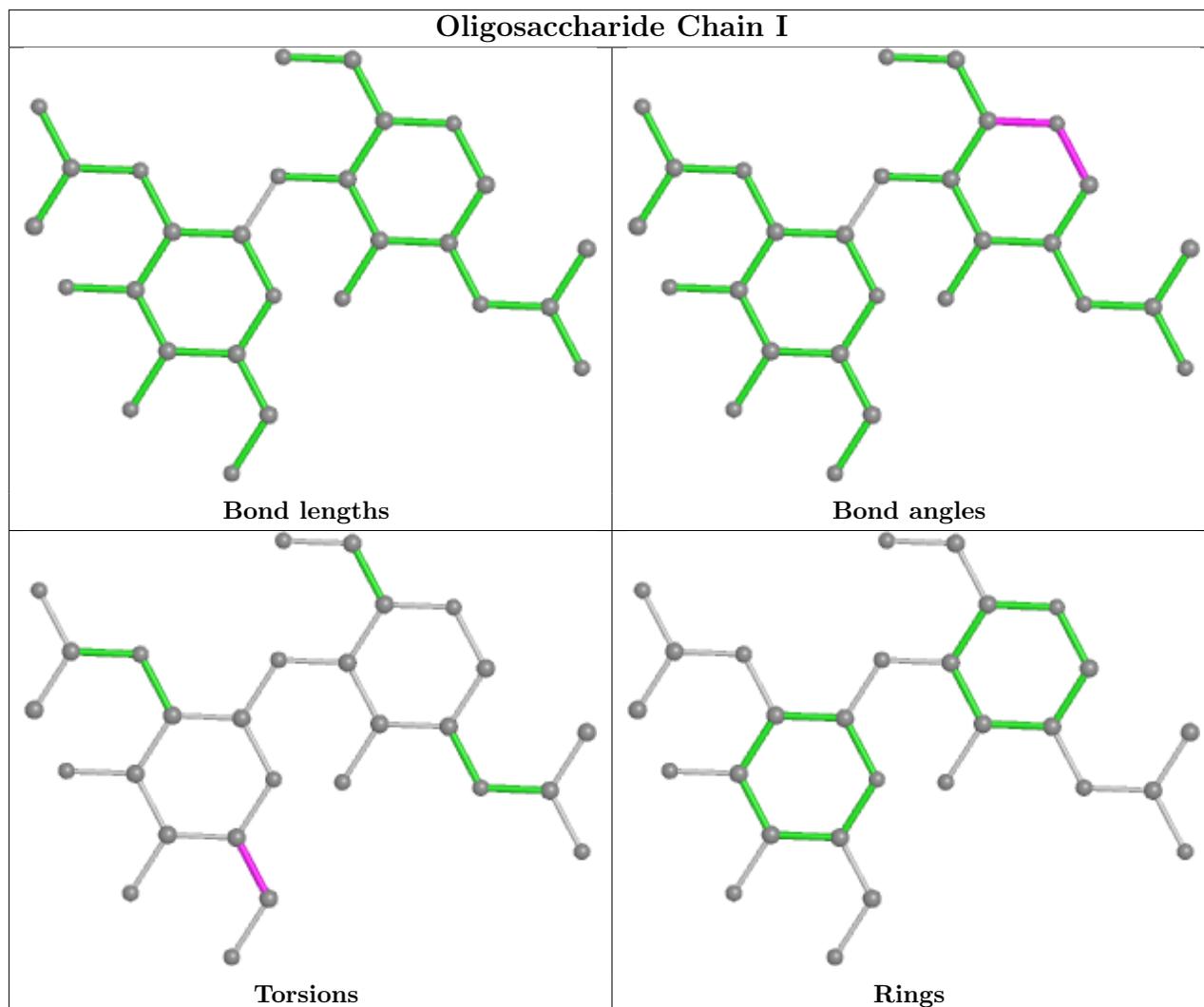
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











## 5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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	$\# Z  > 2$
5	NAG	A	904	1	14,14,15	0.39	0	17,19,21	0.56	0
5	NAG	B	903	1	14,14,15	0.30	0	17,19,21	0.42	0
8	PO4	A	908	-	4,4,4	1.24	0	6,6,6	0.52	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	TRP	B	904	-	14,16,16	0.88	1 (7%)	16,22,22	1.18	2 (12%)
4	H43	A	901	-	31,31,31	1.86	6 (19%)	39,43,43	1.30	5 (12%)
5	NAG	A	903	1	14,14,15	0.45	0	17,19,21	0.77	1 (5%)
6	TRP	A	905	-	14,16,16	0.89	1 (7%)	16,22,22	1.15	2 (12%)
8	PO4	B	906	-	4,4,4	1.23	0	6,6,6	0.51	0
4	H43	B	901	-	31,31,31	1.67	5 (16%)	39,43,43	1.32	4 (10%)
5	NAG	B	902	1	14,14,15	0.39	0	17,19,21	0.51	0
5	NAG	A	902	1	14,14,15	0.56	0	17,19,21	0.66	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	904	1	-	2/6/23/26	0/1/1/1
5	NAG	B	903	1	-	2/6/23/26	0/1/1/1
6	TRP	B	904	-	-	0/7/8/8	0/2/2/2
4	H43	A	901	-	-	8/16/35/35	0/4/4/4
5	NAG	A	903	1	-	0/6/23/26	0/1/1/1
6	TRP	A	905	-	-	0/7/8/8	0/2/2/2
4	H43	B	901	-	-	3/16/35/35	0/4/4/4
5	NAG	B	902	1	-	2/6/23/26	0/1/1/1
5	NAG	A	902	1	-	2/6/23/26	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	H43	C11-N12	6.09	1.57	1.48
4	B	901	H43	C11-N12	4.76	1.55	1.48
4	B	901	H43	C4-C2	4.10	1.59	1.51
4	A	901	H43	C4-C2	3.99	1.59	1.51
4	A	901	H43	C13-N12	3.94	1.56	1.47
4	B	901	H43	C13-N12	3.42	1.55	1.47
4	A	901	H43	C8-N9	2.98	1.47	1.38
4	B	901	H43	C8-N9	2.93	1.46	1.38
4	A	901	H43	C10-C11	2.48	1.57	1.53
6	A	905	TRP	OXT-C	-2.33	1.22	1.30
6	B	904	TRP	OXT-C	-2.30	1.23	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	H43	O3-C2	-2.30	1.23	1.30
4	B	901	H43	O3-C2	-2.29	1.23	1.30

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	901	H43	C25-C26-N9	3.88	107.91	103.35
4	A	901	H43	C25-C26-N9	3.72	107.72	103.35
4	B	901	H43	O1-C2-C4	-3.00	114.48	123.04
4	A	901	H43	O1-C2-C4	-2.95	114.62	123.04
6	A	905	TRP	OXT-C-O	-2.84	117.65	124.09
6	B	904	TRP	OXT-C-O	-2.75	117.84	124.09
5	A	903	NAG	C1-O5-C5	2.75	115.91	112.19
4	B	901	H43	C6-C5-C28	-2.46	114.30	118.17
4	A	901	H43	C6-C5-C28	-2.37	114.43	118.17
6	B	904	TRP	OXT-C-CA	2.36	121.41	113.38
5	A	902	NAG	C1-O5-C5	2.34	115.36	112.19
4	B	901	H43	C19-C20-C21	-2.33	117.75	123.19
4	A	901	H43	C19-C20-C21	-2.27	117.88	123.19
6	A	905	TRP	OXT-C-CA	2.24	121.00	113.38
4	A	901	H43	C27-C8-C7	-2.03	115.04	119.16

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	901	H43	C24-C13-N12-C11
4	A	901	H43	C14-C13-N12-C11
4	A	901	H43	C10-C11-N12-C13
4	A	901	H43	C25-C11-N12-C13
4	A	901	H43	C7-C8-N9-C10
4	A	901	H43	C27-C8-N9-C10
4	B	901	H43	C25-C11-N12-C13
5	A	902	NAG	O5-C5-C6-O6
5	A	904	NAG	O5-C5-C6-O6
5	B	903	NAG	O5-C5-C6-O6
5	A	902	NAG	C4-C5-C6-O6
4	A	901	H43	C27-C8-N9-C26
5	A	904	NAG	C4-C5-C6-O6
5	B	903	NAG	C4-C5-C6-O6
4	A	901	H43	C7-C8-N9-C26
4	B	901	H43	C27-C8-N9-C10

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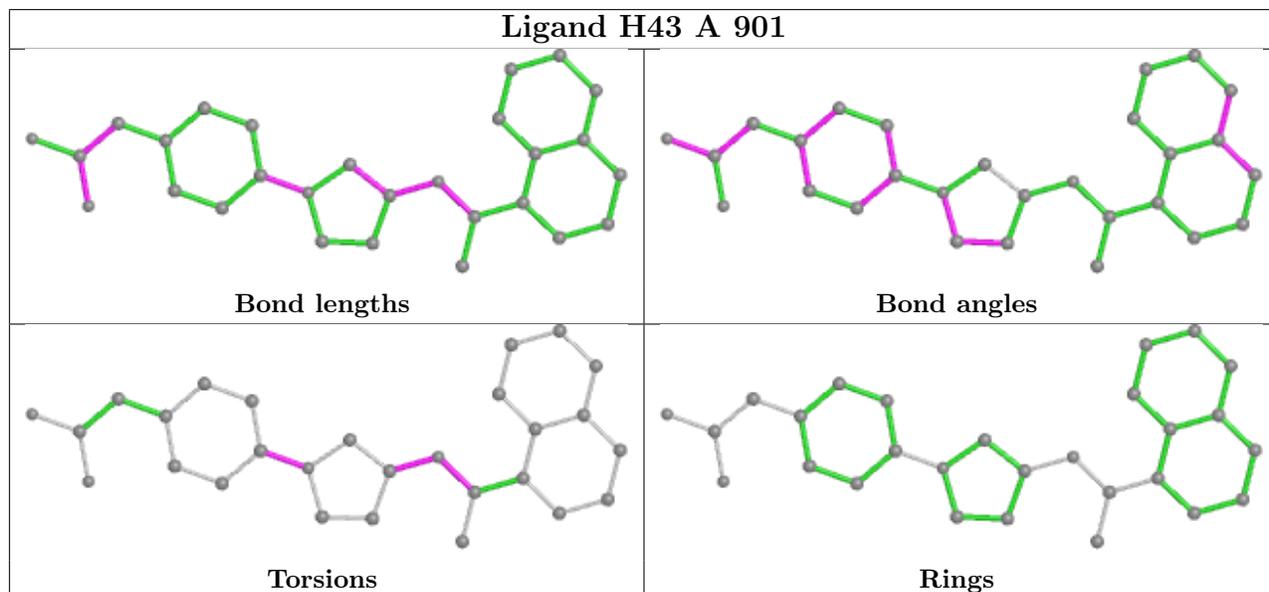
Mol	Chain	Res	Type	Atoms
5	B	902	NAG	C4-C5-C6-O6
4	B	901	H43	C7-C8-N9-C10
5	B	902	NAG	O5-C5-C6-O6

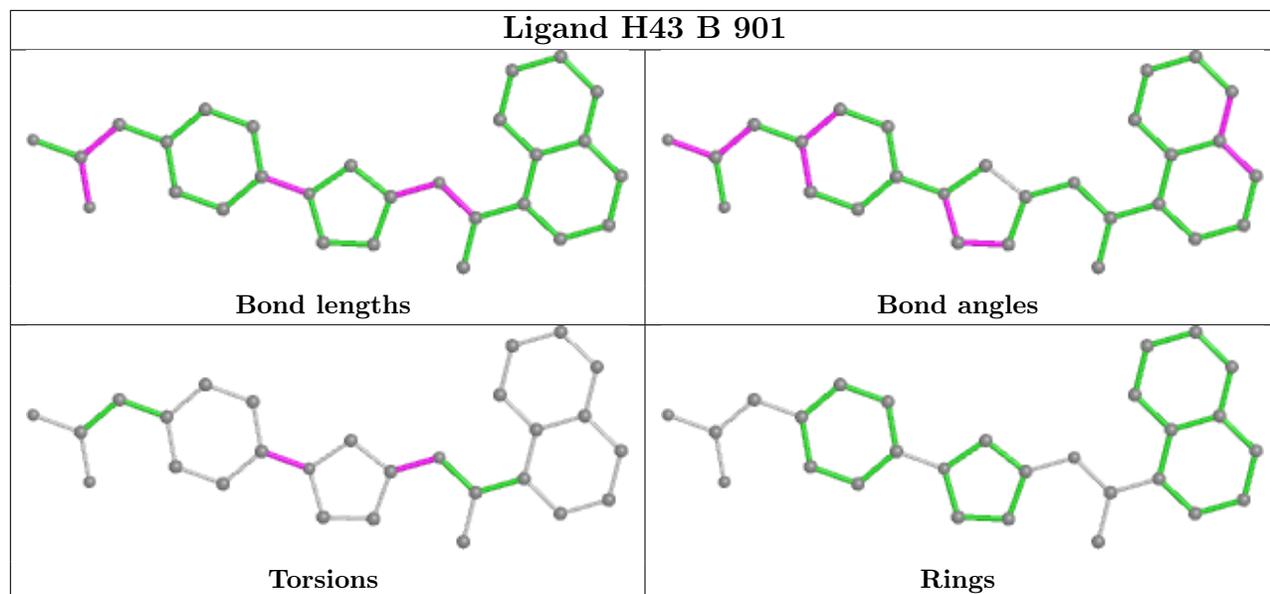
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	906	PO4	1	0
5	B	902	NAG	1	0
5	A	902	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 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.

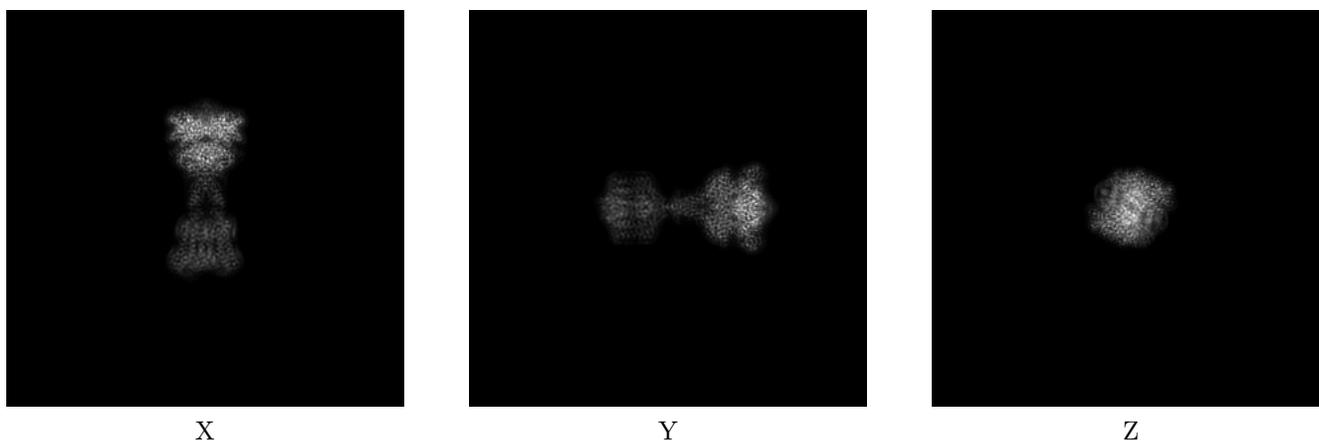
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23654. 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](#)

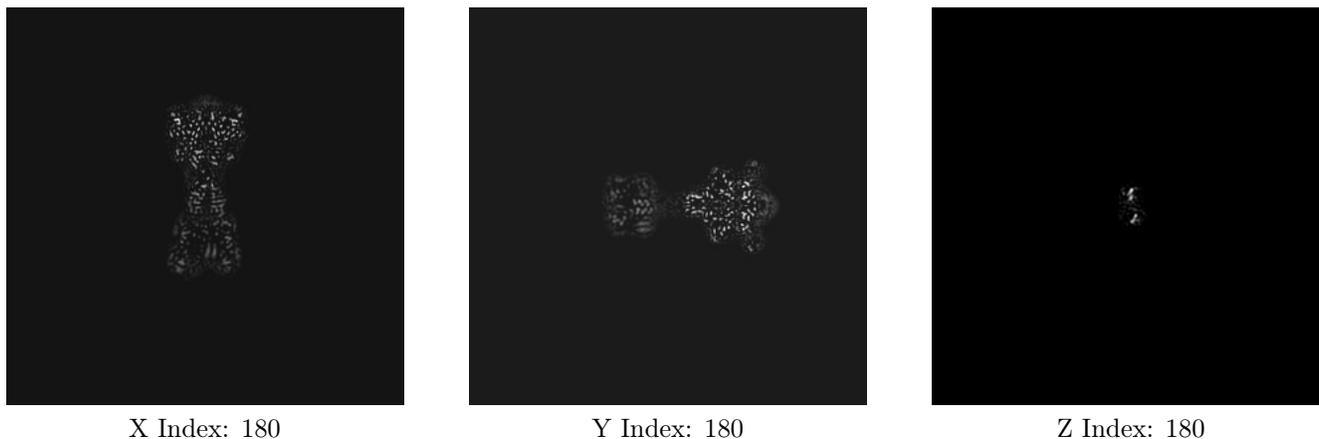
#### 6.1.1 Primary map



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

### 6.2 Central slices [i](#)

#### 6.2.1 Primary map



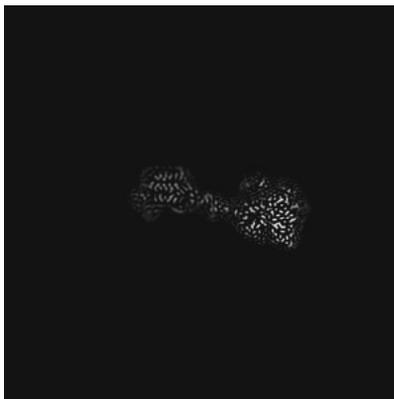
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

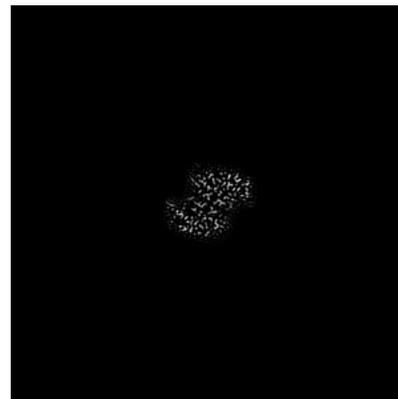
### 6.3.1 Primary map



X Index: 183



Y Index: 168



Z Index: 253

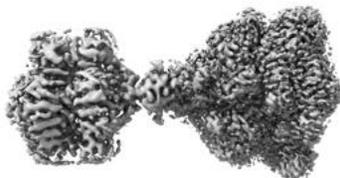
The images above show the largest variance slices of the map in three orthogonal directions.

## 6.4 Orthogonal surface views [i](#)

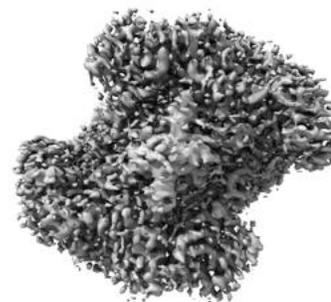
### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.9. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

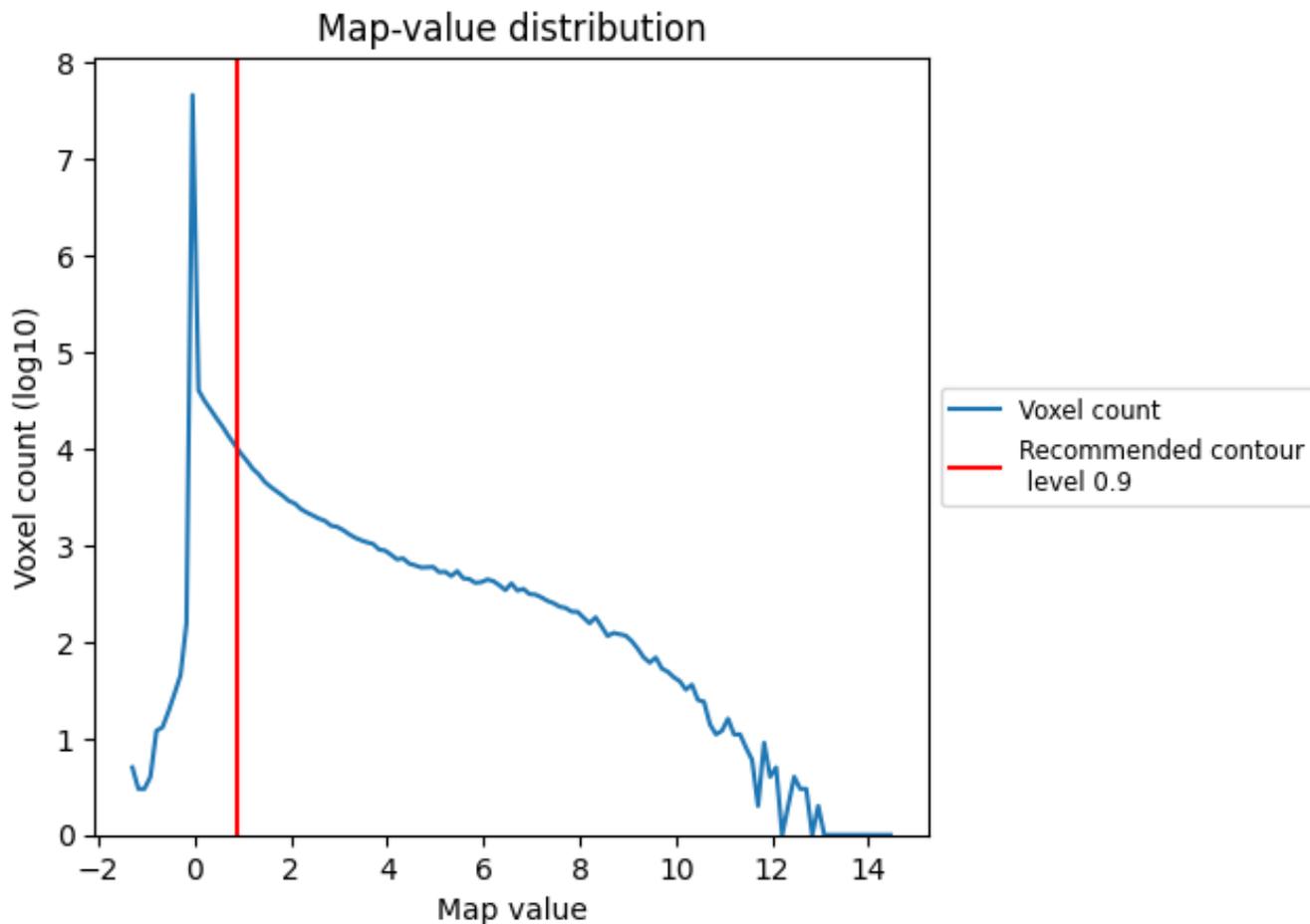
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

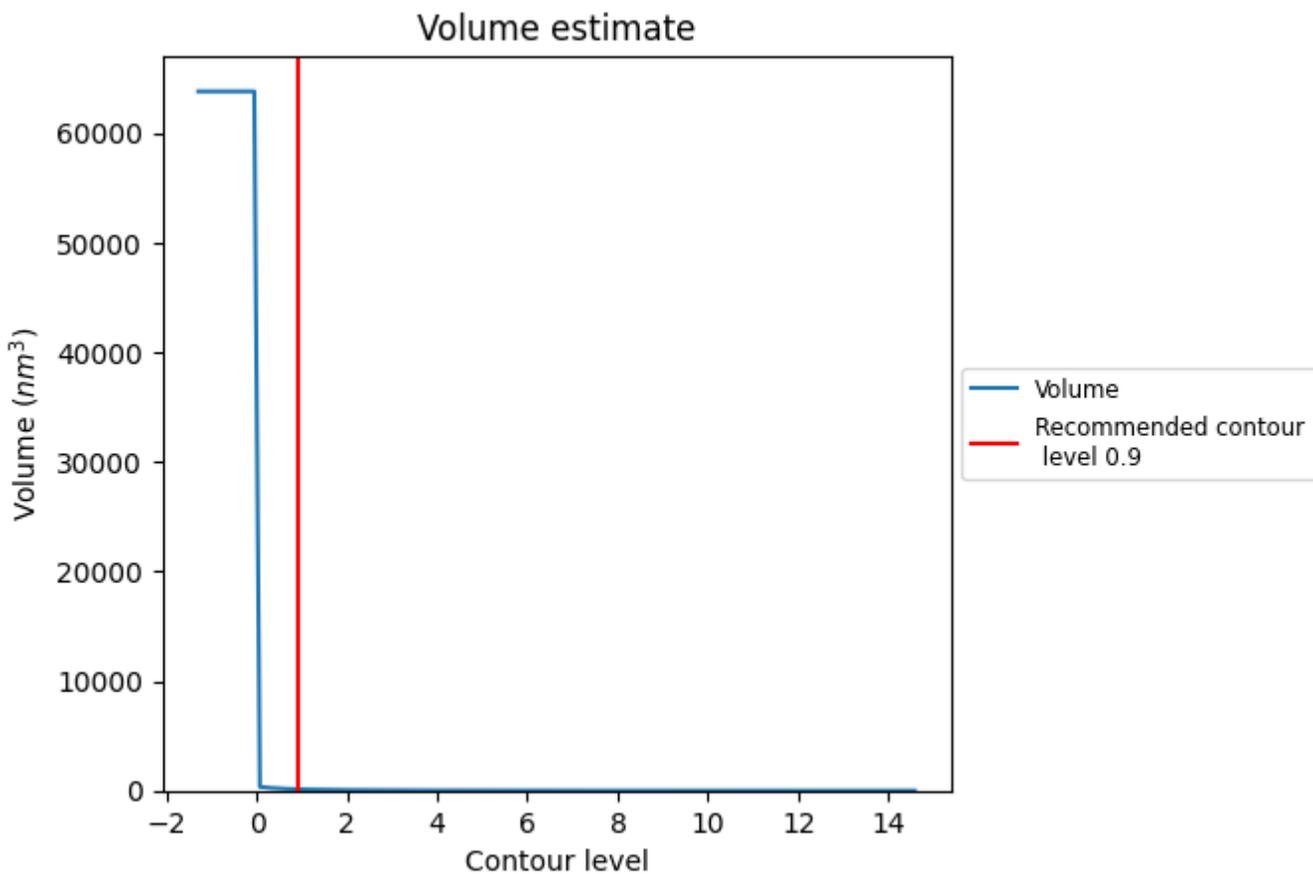
This section contains the results of statistical analysis of the map.

### 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.

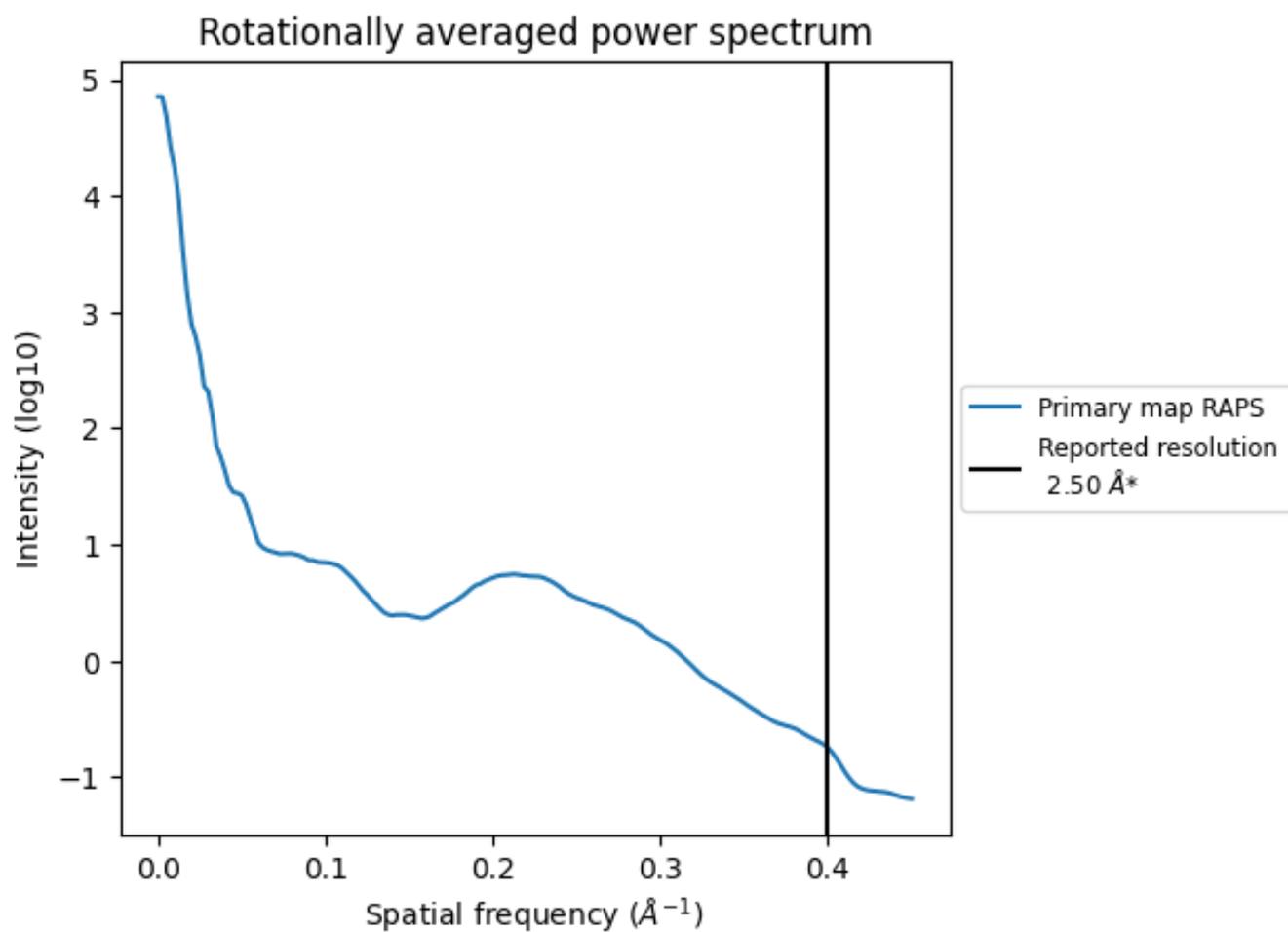
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 127  $\text{nm}^3$ ; this corresponds to an approximate mass of 115 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 0.400 Å<sup>-1</sup>

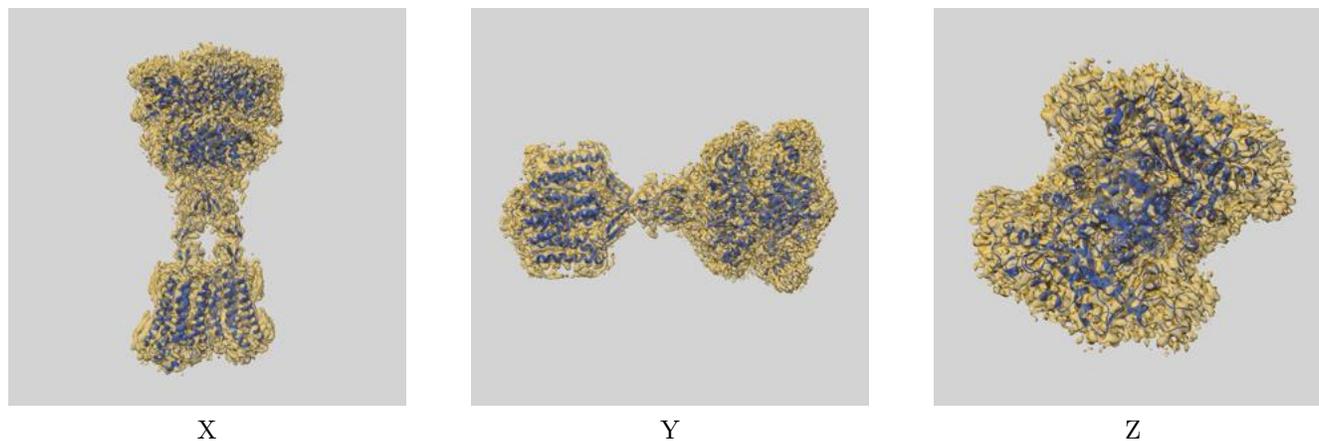
## 8 Fourier-Shell correlation

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

## 9 Map-model fit [i](#)

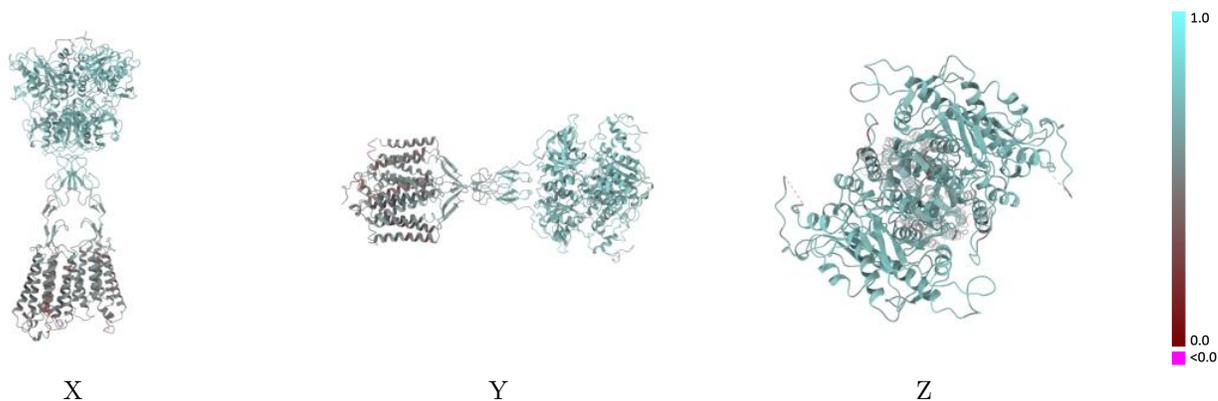
This section contains information regarding the fit between EMDB map EMD-23654 and PDB model 7M3G. Per-residue inclusion information can be found in section 3 on page 9.

### 9.1 Map-model overlay [i](#)



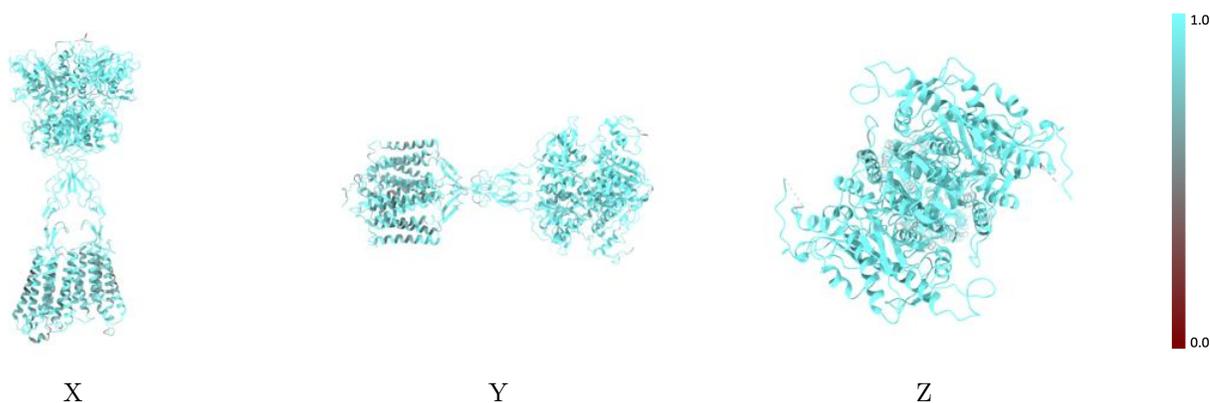
The images above show the 3D surface view of the map at the recommended contour level 0.9 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



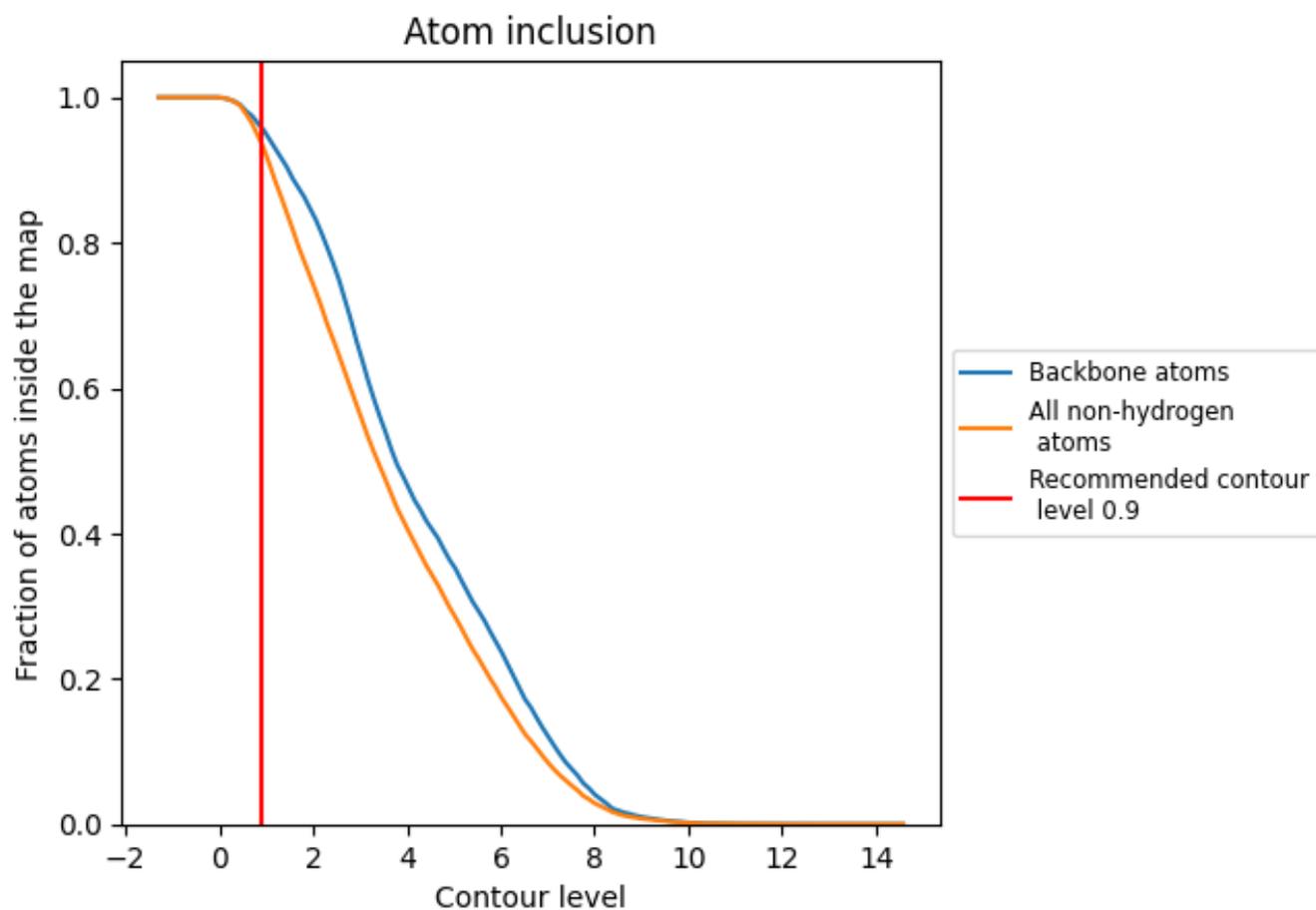
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.9).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 94% 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 (0.9) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9363	 0.6070
A	 0.9312	 0.6040
B	 0.9431	 0.6130
C	 0.8545	 0.5600
D	 0.9636	 0.6210
E	 0.7857	 0.4650
F	 1.0000	 0.5870
G	 0.7143	 0.4060
H	 0.9286	 0.5650
I	 0.9643	 0.5730

