



## Full wwPDB EM Validation Report ⓘ

Dec 10, 2022 – 11:30 pm GMT

PDB ID : 6R7X  
EMDB ID : EMD-4746  
Title : CryoEM structure of calcium-bound human TMEM16K / Anoctamin 10 in detergent (2mM Ca<sup>2+</sup>, closed form)  
Authors : Pike, A.C.W.; Bushell, S.R.; Shintre, C.A.; Tessitore, A.; Baronina, A.; Chu, A.; Mukhopadhyay, S.; Shrestha, L.; Chalk, R.; Burgess-Brown, N.A.; Love, J.; Huiskonen, J.T.; Edwards, A.M.; Arrowsmith, C.H.; Bountra, C.; Carpenter, E.P.; Structural Genomics Consortium (SGC)  
Deposited on : 2019-03-29  
Resolution : 3.47 Å (reported)  
Based on initial model : 5OC9

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 : 1.8.4, 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)

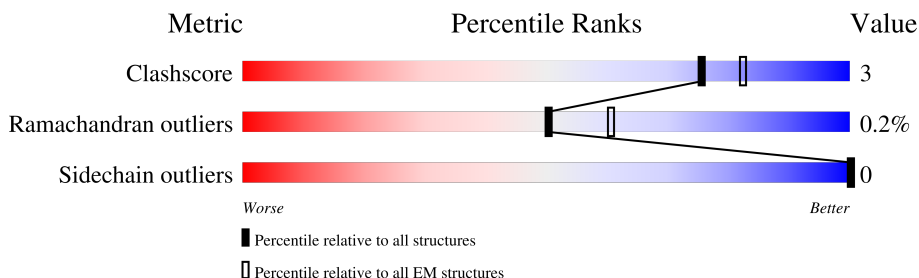
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.47 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	667	 5% 86% 8% 6%
1	B	667	 5% 87% 8% 6%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10630 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 Anoctamin-10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	629	5060	3341	821	871	27	0	0
1	B	629	5060	3341	821	871	27	0	0

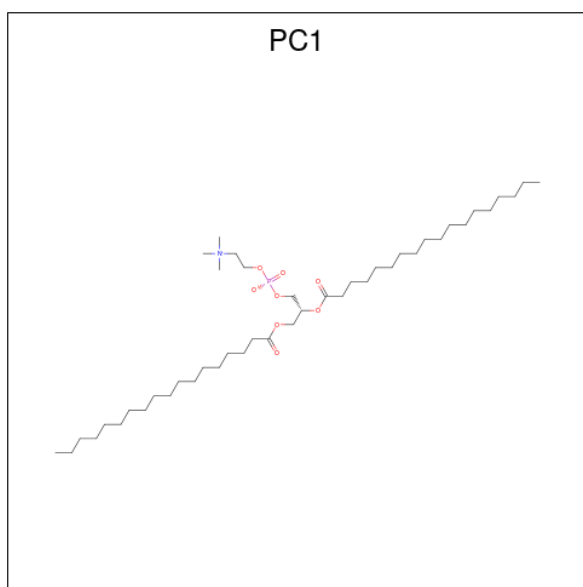
There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	661	ALA	-	expression tag	UNP Q9NW15
A	662	GLU	-	expression tag	UNP Q9NW15
A	663	ASN	-	expression tag	UNP Q9NW15
A	664	LEU	-	expression tag	UNP Q9NW15
A	665	TYR	-	expression tag	UNP Q9NW15
A	666	PHE	-	expression tag	UNP Q9NW15
A	667	GLN	-	expression tag	UNP Q9NW15
B	661	ALA	-	expression tag	UNP Q9NW15
B	662	GLU	-	expression tag	UNP Q9NW15
B	663	ASN	-	expression tag	UNP Q9NW15
B	664	LEU	-	expression tag	UNP Q9NW15
B	665	TYR	-	expression tag	UNP Q9NW15
B	666	PHE	-	expression tag	UNP Q9NW15
B	667	GLN	-	expression tag	UNP Q9NW15

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

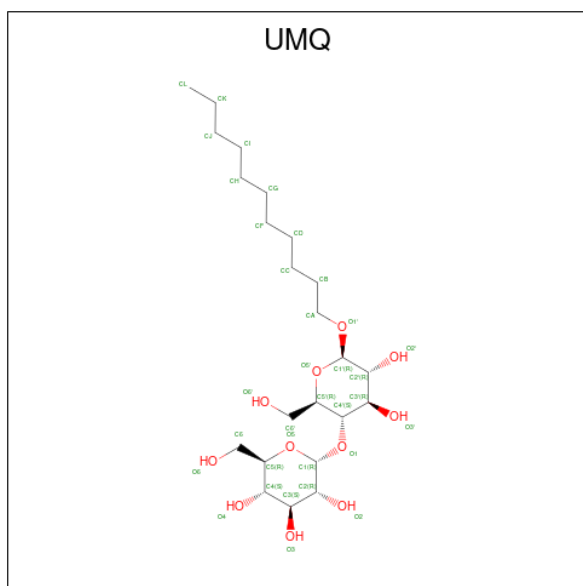
Mol	Chain	Residues	Atoms		AltConf
2	A	3	Total	Ca	0
			3	3	
2	B	3	Total	Ca	0
			3	3	

- Molecule 3 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: PC1) (formula: C<sub>44</sub>H<sub>88</sub>NO<sub>8</sub>P).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
3	A	1	47	37	1	8	1	0
3	B	1	47	37	1	8	1	0

- Molecule 4 is UNDECYL-MALTOSE (three-letter code: UMQ) (formula:  $C_{23}H_{44}O_{11}$ ).



Mol	Chain	Residues	Atoms			AltConf
			Total	C	O	
4	A	1	205	158	47	0
4	A	1	205	158	47	0

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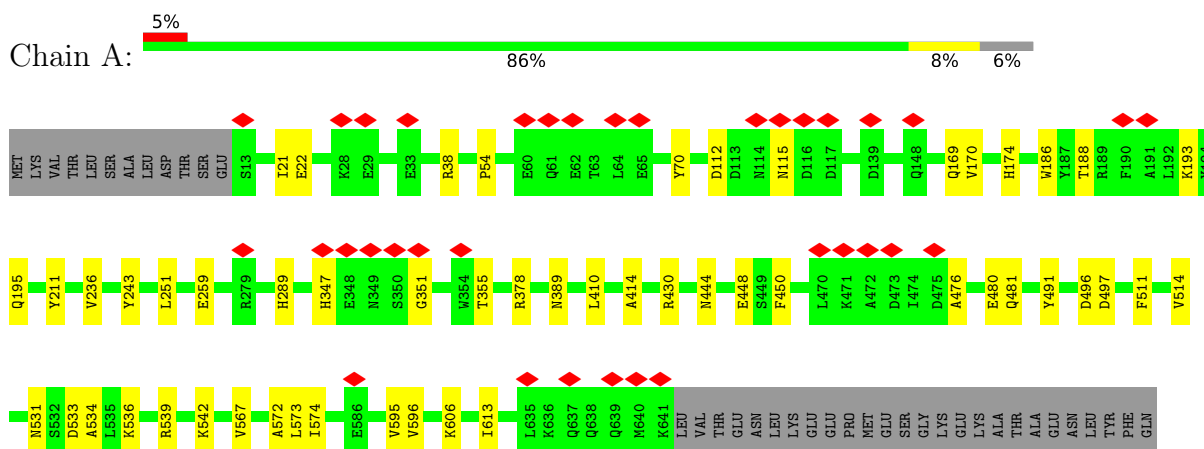
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Mol	Chain	Residues	Atoms			AltConf
			Total	C	O	
4	A	1	205	158	47	0
4	A	1	205	158	47	0
4	A	1	205	158	47	0
4	A	1	205	158	47	0
4	A	1	205	158	47	0
4	A	1	205	158	47	0
4	A	1	205	158	47	0
4	A	1	205	158	47	0
4	A	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0
4	B	1	205	158	47	0

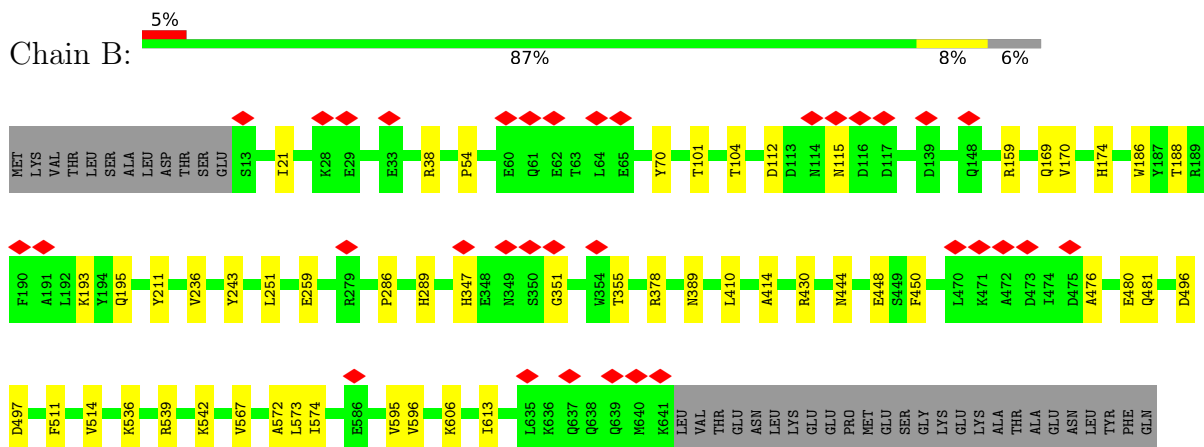
### 3 Residue-property plots

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: Anoctamin-10



- Molecule 1: Anoctamin-10



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	70940	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 ( $e^-/\text{\AA}^2$ )	50.9	Depositor
Minimum defocus (nm)	1250	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.128	Depositor
Minimum map value	-0.091	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.005	Depositor
Recommended contour level	0.017	Depositor
Map size (Å)	187.416, 187.416, 187.416	wwPDB
Map dimensions	228, 228, 228	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.822, 0.822, 0.822	Depositor

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PC1, CA, UMQ

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	0.29	0/5187	0.46	0/7039
1	B	0.29	0/5187	0.46	0/7039
All	All	0.29	0/10374	0.46	0/14078

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

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	5060	0	5091	34	0
1	B	5060	0	5091	32	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	47	0	68	0	0
3	B	47	0	68	1	0
4	A	205	0	305	4	0
4	B	205	0	305	4	0
All	All	10630	0	10928	66	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 3.



All (66) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:476:ALA:O	1:B:480:GLU:HB3	1.83	0.77
1:A:476:ALA:O	1:A:480:GLU:HB3	1.83	0.77
1:A:186:TRP:HE1	1:A:195:GLN:HE21	1.44	0.65
1:B:186:TRP:HE1	1:B:195:GLN:HE21	1.44	0.64
1:A:243:TYR:HB3	1:A:574:ILE:HG12	1.87	0.57
1:B:243:TYR:HB3	1:B:574:ILE:HG12	1.87	0.57
1:A:112:ASP:HB2	1:A:115:ASN:HB2	1.88	0.56
1:A:378:ARG:NH1	1:A:496:ASP:OD2	2.39	0.56
1:B:169:GLN:OE1	1:B:542:LYS:NZ	2.39	0.55
1:A:169:GLN:OE1	1:A:542:LYS:NZ	2.39	0.55
1:B:378:ARG:NH1	1:B:496:ASP:OD2	2.39	0.55
1:B:112:ASP:HB2	1:B:115:ASN:HB2	1.88	0.54
1:B:289:HIS:O	1:B:389:ASN:ND2	2.41	0.54
1:A:289:HIS:O	1:A:389:ASN:ND2	2.41	0.53
1:A:430:ARG:NH1	1:A:514:VAL:O	2.43	0.51
1:B:430:ARG:NH1	1:B:514:VAL:O	2.43	0.51
1:A:497:ASP:OD1	1:A:536:LYS:NZ	2.44	0.51
1:B:251:LEU:HD13	1:B:567:VAL:HG22	1.93	0.51
1:B:497:ASP:OD1	1:B:536:LYS:NZ	2.44	0.50
1:B:174:HIS:O	1:B:481:GLN:NE2	2.46	0.49
1:A:251:LEU:HD13	1:A:567:VAL:HG22	1.93	0.49
1:B:539:ARG:HG3	4:B:908:UMQ:H2'1	1.96	0.48
1:A:21:ILE:HG12	1:A:170:VAL:HG22	1.94	0.48
1:A:174:HIS:O	1:A:481:GLN:NE2	2.46	0.48
1:A:539:ARG:HG3	4:A:707:UMQ:H2'1	1.96	0.48
1:B:21:ILE:HG12	1:B:170:VAL:HG22	1.94	0.48
1:A:595:VAL:HG11	4:A:705:UMQ:HH2	1.96	0.47
1:B:444:ASN:O	1:B:448:GLU:HB3	2.14	0.47
1:B:572:ALA:HA	4:B:906:UMQ:HH1	1.96	0.47
1:A:444:ASN:O	1:A:448:GLU:HB3	2.14	0.47
1:A:572:ALA:HA	4:A:705:UMQ:HH1	1.96	0.46
1:B:448:GLU:OE2	1:B:536:LYS:NZ	2.47	0.46
1:A:347:HIS:ND1	1:A:355:THR:O	2.47	0.46
1:B:347:HIS:ND1	1:B:355:THR:O	2.47	0.46
1:B:450:PHE:HZ	4:B:908:UMQ:HK1	1.81	0.46
1:A:450:PHE:HZ	4:A:707:UMQ:HK1	1.81	0.46
1:B:595:VAL:HG11	4:B:906:UMQ:HH2	1.96	0.46
1:B:211:TYR:CZ	1:B:497:ASP:HB3	2.52	0.45
1:A:211:TYR:CZ	1:A:497:ASP:HB3	2.52	0.45
1:A:596:VAL:HG21	1:B:596:VAL:HG21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:THR:HG22	1:B:193:LYS:HD3	2.00	0.44
1:A:347:HIS:HA	1:A:355:THR:HB	1.99	0.44
1:B:347:HIS:HA	1:B:355:THR:HB	1.99	0.44
1:A:54:PRO:HA	1:A:70:TYR:HD1	1.83	0.43
1:A:448:GLU:OE2	1:A:536:LYS:NZ	2.47	0.43
1:B:351:GLY:H	1:B:355:THR:HG21	1.83	0.43
1:B:511:PHE:HD1	1:B:573:LEU:HD11	1.83	0.43
1:B:54:PRO:HA	1:B:70:TYR:HD1	1.83	0.43
1:A:188:THR:HG22	1:A:193:LYS:HD3	2.00	0.43
1:A:511:PHE:HD1	1:A:573:LEU:HD11	1.83	0.43
1:B:38:ARG:HD2	1:B:38:ARG:HA	1.88	0.43
1:A:410:LEU:O	1:A:414:ALA:HB2	2.19	0.42
1:A:211:TYR:OH	1:A:533:ASP:OD2	2.31	0.42
1:B:101:THR:HG1	1:B:104:THR:HG1	1.68	0.42
1:B:606:LYS:HB2	1:B:606:LYS:HE2	1.89	0.42
1:A:351:GLY:H	1:A:355:THR:HG21	1.83	0.42
1:A:606:LYS:HB2	1:A:606:LYS:HE2	1.89	0.41
1:A:259:GLU:HG2	1:A:613:ILE:HB	2.02	0.41
1:B:410:LEU:O	1:B:414:ALA:HB2	2.20	0.41
1:A:491:TYR:OH	1:A:497:ASP:OD2	2.33	0.41
1:B:159:ARG:HD2	1:B:286:PRO:HB3	2.03	0.41
3:B:905:PC1:H143	3:B:905:PC1:H112	1.88	0.40
1:A:22:GLU:OE1	1:A:169:GLN:NE2	2.43	0.40
1:B:259:GLU:HG2	1:B:613:ILE:HB	2.02	0.40
1:A:531:ASN:O	1:A:534:ALA:HB3	2.21	0.40
1:A:38:ARG:HD2	1:A:38:ARG:HA	1.89	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	627/667 (94%)	611 (97%)	15 (2%)	1 (0%)	47	80
1	B	627/667 (94%)	611 (97%)	15 (2%)	1 (0%)	47	80
All	All	1254/1334 (94%)	1222 (97%)	30 (2%)	2 (0%)	50	80

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	VAL
1	B	236	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	530/591 (90%)	530 (100%)	0	100	100
1	B	530/591 (90%)	530 (100%)	0	100	100
All	All	1060/1182 (90%)	1060 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	195	GLN
1	A	504	GLN
1	B	66	ASN
1	B	195	GLN
1	B	504	GLN

### 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](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 28 ligands modelled in this entry, 6 are monoatomic - leaving 22 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
3	PC1	B	905	-	46,46,53	0.31	0	52,54,61	0.38	0
4	UMQ	B	907	-	23,23,35	0.15	0	28,28,46	0.24	0
4	UMQ	B	908	-	35,35,35	0.14	0	46,46,46	0.25	0
4	UMQ	A	708	-	23,23,35	0.15	0	28,28,46	0.26	0
4	UMQ	A	705	-	35,35,35	0.16	0	46,46,46	0.58	2 (4%)
4	UMQ	B	913	-	9,9,35	0.14	0	8,8,46	0.13	0
4	UMQ	B	906	-	35,35,35	0.16	0	46,46,46	0.57	2 (4%)
4	UMQ	B	911	-	12,12,35	0.20	0	11,11,46	0.12	0
4	UMQ	A	706	-	23,23,35	0.15	0	28,28,46	0.25	0
4	UMQ	B	901	-	35,35,35	0.14	0	46,46,46	0.40	0
4	UMQ	A	710	-	12,12,35	0.21	0	11,11,46	0.12	0
4	UMQ	B	912	-	10,10,35	0.10	0	9,9,46	0.15	0
4	UMQ	B	909	-	23,23,35	0.14	0	28,28,46	0.26	0
4	UMQ	A	712	-	9,9,35	0.15	0	8,8,46	0.13	0
4	UMQ	A	714	-	35,35,35	0.14	0	46,46,46	0.40	0
4	UMQ	A	713	-	11,11,35	0.12	0	10,10,46	0.19	0
4	UMQ	B	914	-	11,11,35	0.12	0	10,10,46	0.19	0
4	UMQ	B	910	-	10,10,35	0.12	0	9,9,46	0.10	0
4	UMQ	A	707	-	35,35,35	0.13	0	46,46,46	0.25	0
4	UMQ	A	709	-	10,10,35	0.13	0	9,9,46	0.11	0
3	PC1	A	704	-	46,46,53	0.31	0	52,54,61	0.38	0
4	UMQ	A	711	-	10,10,35	0.10	0	9,9,46	0.15	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	PC1	B	905	-	-	10/50/50/57	-
4	UMQ	B	907	-	-	4/14/34/60	0/1/1/2
4	UMQ	B	908	-	-	3/20/60/60	0/2/2/2
4	UMQ	A	708	-	-	0/14/34/60	0/1/1/2
4	UMQ	A	705	-	-	10/20/60/60	0/2/2/2
4	UMQ	B	913	-	-	0/7/7/60	-
4	UMQ	B	906	-	-	10/20/60/60	0/2/2/2
4	UMQ	B	911	-	-	0/10/10/60	-
4	UMQ	A	706	-	-	4/14/34/60	0/1/1/2
4	UMQ	B	901	-	-	5/20/60/60	0/2/2/2
4	UMQ	A	710	-	-	0/10/10/60	-
4	UMQ	B	912	-	-	0/8/8/60	-
4	UMQ	B	909	-	-	0/14/34/60	0/1/1/2
4	UMQ	A	712	-	-	0/7/7/60	-
4	UMQ	A	714	-	-	5/20/60/60	0/2/2/2
4	UMQ	A	713	-	-	1/9/9/60	-
4	UMQ	B	914	-	-	1/9/9/60	-
4	UMQ	B	910	-	-	0/8/8/60	-
4	UMQ	A	707	-	-	3/20/60/60	0/2/2/2
4	UMQ	A	709	-	-	0/8/8/60	-
3	PC1	A	704	-	-	10/50/50/57	-
4	UMQ	A	711	-	-	0/8/8/60	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	906	UMQ	C1-O1-C4'	2.34	123.77	117.96
4	A	705	UMQ	C1-O1-C4'	2.33	123.74	117.96
4	A	705	UMQ	O1-C4'-C5'	2.25	115.61	109.45
4	B	906	UMQ	O1-C4'-C5'	2.24	115.57	109.45

There are no chirality outliers.

All (66) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	704	PC1	C11-O13-P-O12
3	A	704	PC1	C11-O13-P-O14
3	A	704	PC1	C11-O13-P-O11
3	A	704	PC1	C1-O11-P-O14
3	B	905	PC1	C11-O13-P-O12
3	B	905	PC1	C11-O13-P-O14
3	B	905	PC1	C11-O13-P-O11
3	B	905	PC1	C1-O11-P-O14
4	A	706	UMQ	C2'-C1'-O1'-CA
4	A	706	UMQ	O5'-C1'-O1'-CA
4	A	706	UMQ	CB-CA-O1'-C1'
4	A	714	UMQ	C2'-C1'-O1'-CA
4	A	714	UMQ	O5'-C1'-O1'-CA
4	A	714	UMQ	CB-CA-O1'-C1'
4	B	901	UMQ	C2'-C1'-O1'-CA
4	B	901	UMQ	O5'-C1'-O1'-CA
4	B	901	UMQ	CB-CA-O1'-C1'
4	B	907	UMQ	C2'-C1'-O1'-CA
4	B	907	UMQ	O5'-C1'-O1'-CA
4	B	907	UMQ	CB-CA-O1'-C1'
4	A	705	UMQ	C5'-C4'-O1-C1
4	B	906	UMQ	C5'-C4'-O1-C1
4	A	713	UMQ	CF-CG-CH-CI
4	B	914	UMQ	CF-CG-CH-CI
4	A	705	UMQ	C2'-C1'-O1'-CA
4	B	906	UMQ	C2'-C1'-O1'-CA
4	A	705	UMQ	CB-CA-O1'-C1'
4	B	906	UMQ	CB-CA-O1'-C1'
4	A	705	UMQ	O5-C5-C6-O6
4	B	906	UMQ	O5-C5-C6-O6
3	A	704	PC1	C1-C2-C3-O31
3	B	905	PC1	C1-C2-C3-O31
4	B	906	UMQ	O1'-CA-CB-CC
4	A	705	UMQ	O1'-CA-CB-CC
4	B	906	UMQ	O5-C1-O1-C4'
4	A	705	UMQ	O5-C1-O1-C4'
4	B	907	UMQ	O1'-CA-CB-CC
4	A	706	UMQ	O1'-CA-CB-CC
3	A	704	PC1	O21-C2-C3-O31
3	B	905	PC1	O21-C2-C3-O31
4	A	705	UMQ	CC-CD-CF-CG
4	B	906	UMQ	CC-CD-CF-CG
3	A	704	PC1	C1-O11-P-O13

*Continued on next page...*

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	B	905	PC1	C1-O11-P-O13
4	A	705	UMQ	C2-C1-O1-C4'
4	B	906	UMQ	C2-C1-O1-C4'
4	B	901	UMQ	CH-CI-CJ-CK
4	A	714	UMQ	CH-CI-CJ-CK
4	B	908	UMQ	CB-CC-CD-CF
4	A	707	UMQ	CB-CC-CD-CF
3	A	704	PC1	C3A-C3B-C3C-C3D
3	B	905	PC1	C3A-C3B-C3C-C3D
4	A	705	UMQ	CF-CG-CH-CI
4	B	906	UMQ	CF-CG-CH-CI
3	A	704	PC1	O31-C31-C32-C33
3	B	905	PC1	O31-C31-C32-C33
3	A	704	PC1	O32-C31-C32-C33
3	B	905	PC1	O32-C31-C32-C33
4	A	707	UMQ	C2'-C1'-O1'-CA
4	B	908	UMQ	C2'-C1'-O1'-CA
4	A	707	UMQ	O5'-C1'-O1'-CA
4	B	908	UMQ	O5'-C1'-O1'-CA
4	A	714	UMQ	CG-CH-CI-CJ
4	B	901	UMQ	CG-CH-CI-CJ
4	A	705	UMQ	CA-CB-CC-CD
4	B	906	UMQ	CA-CB-CC-CD

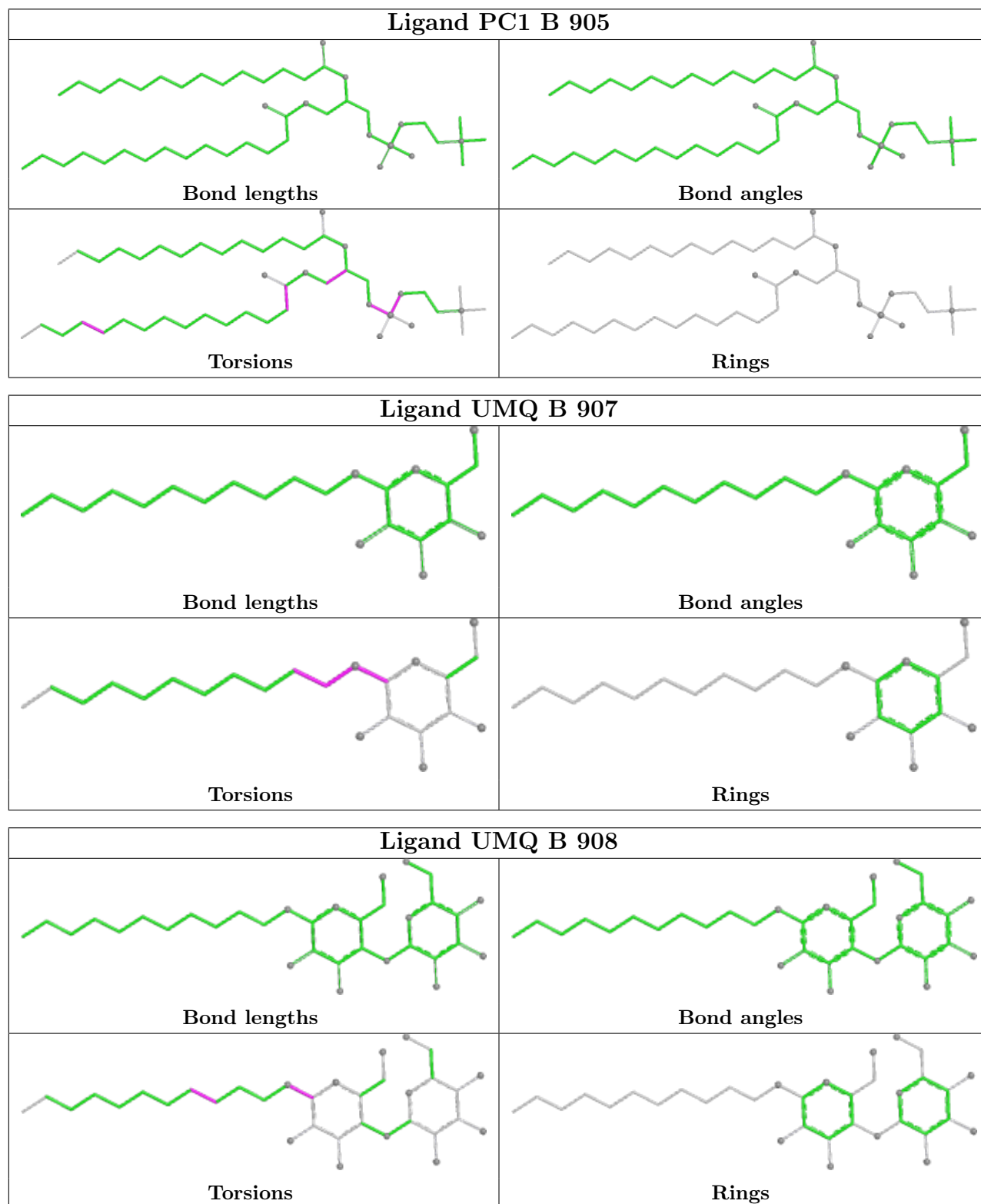
There are no ring outliers.

5 monomers are involved in 9 short contacts:

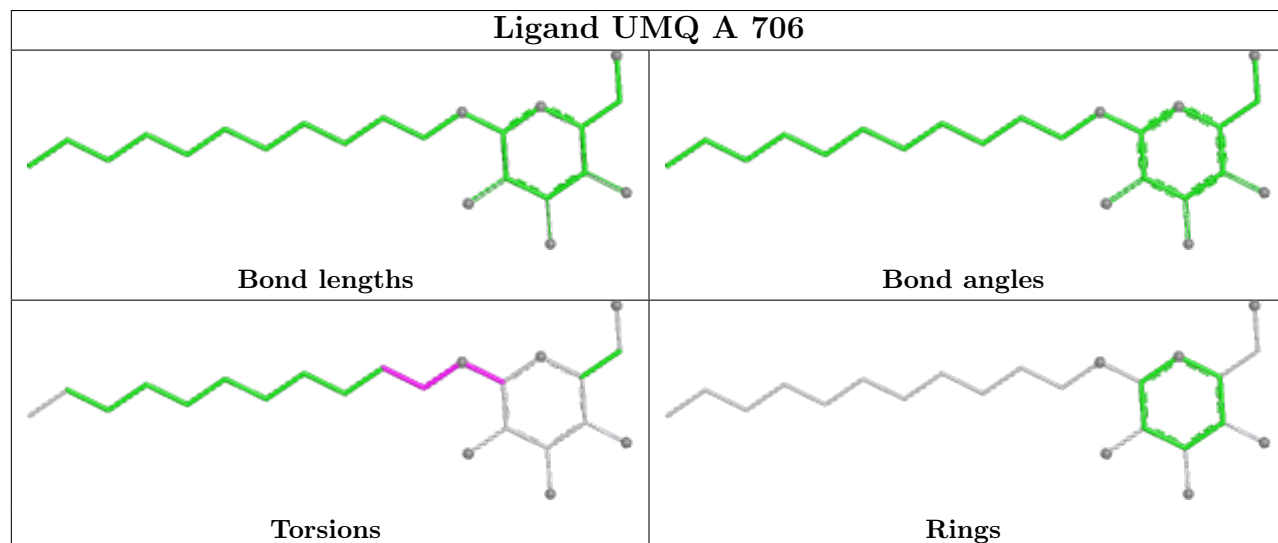
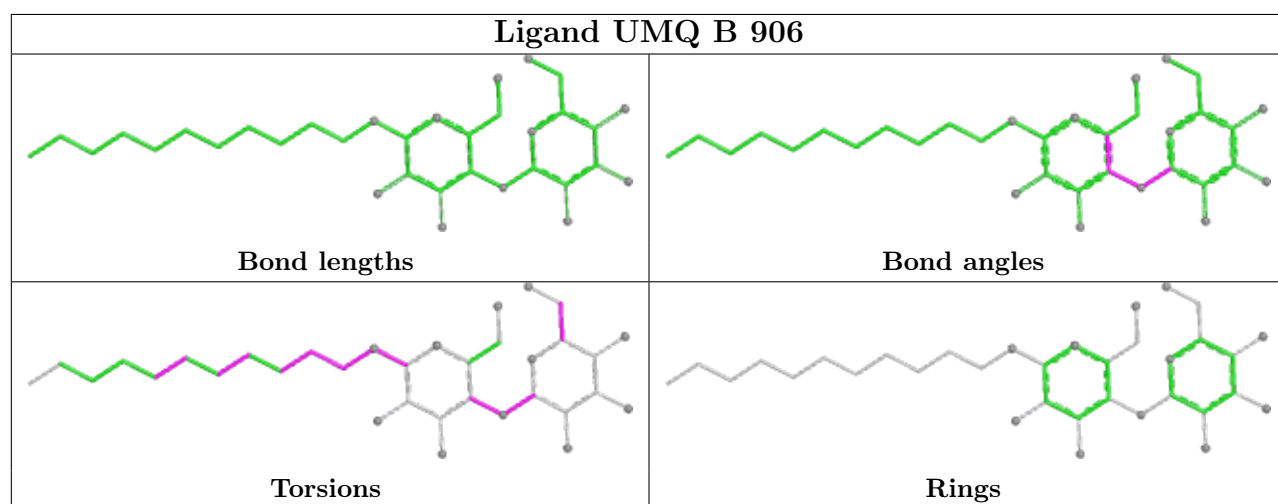
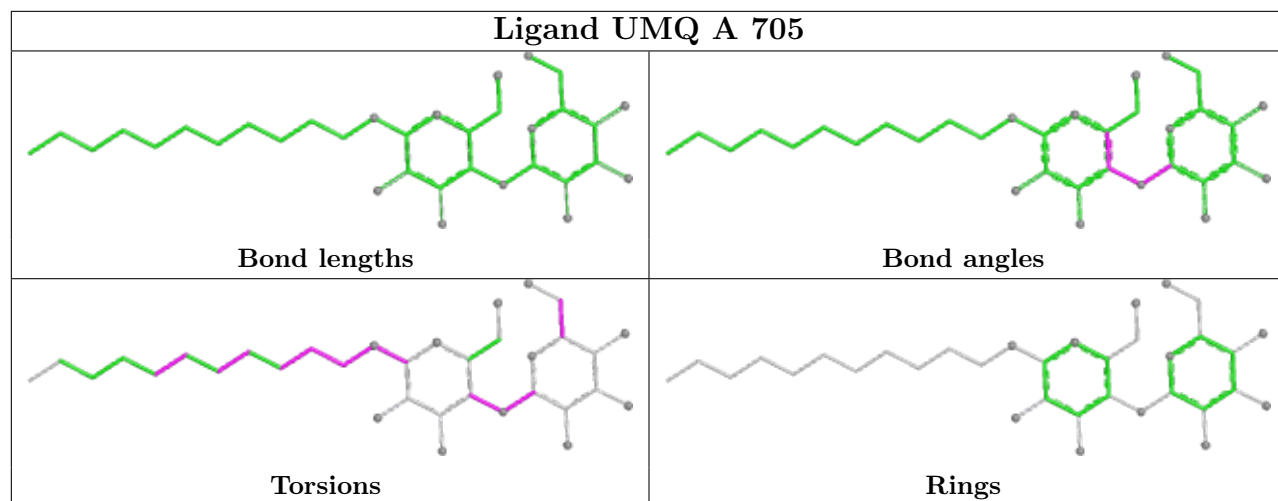
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	905	PC1	1	0
4	B	908	UMQ	2	0
4	A	705	UMQ	2	0
4	B	906	UMQ	2	0
4	A	707	UMQ	2	0

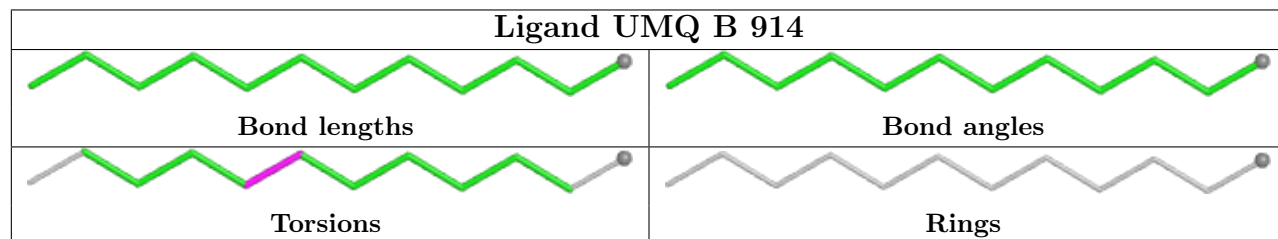
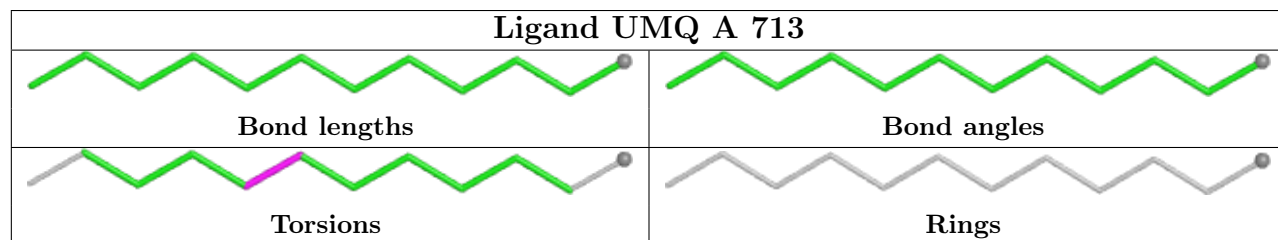
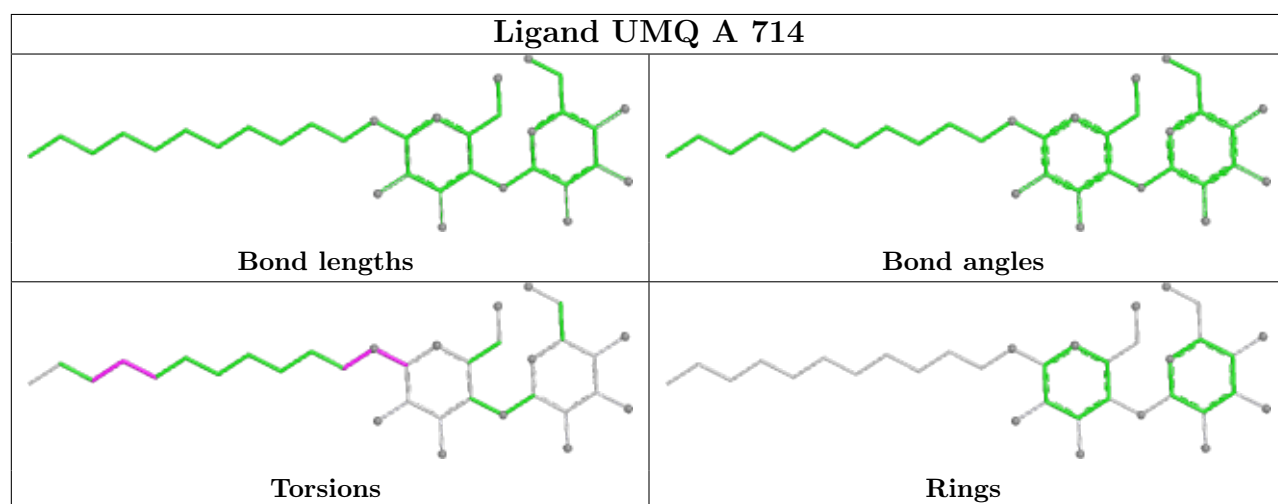
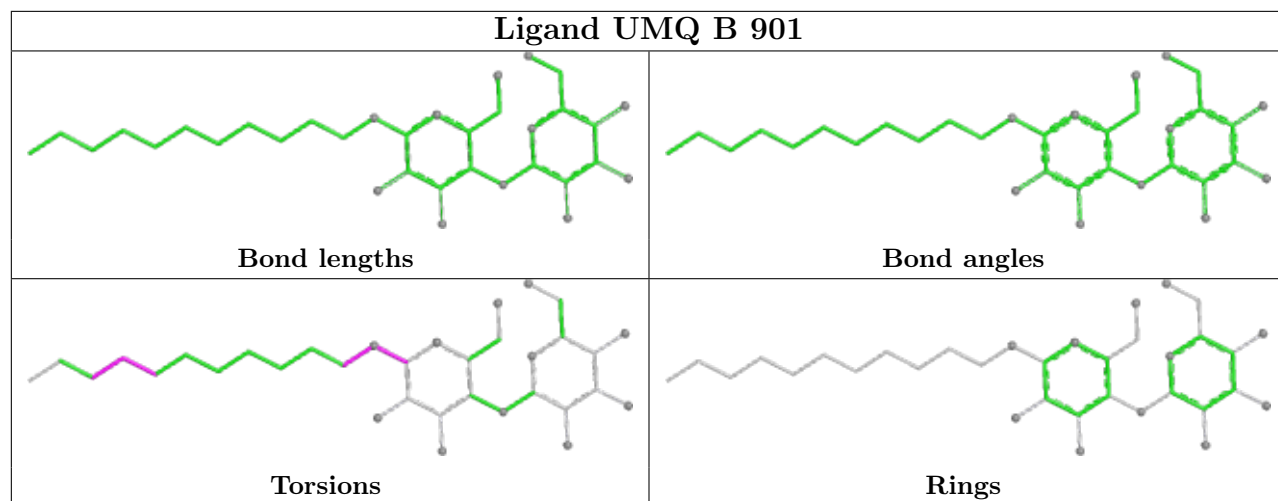
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

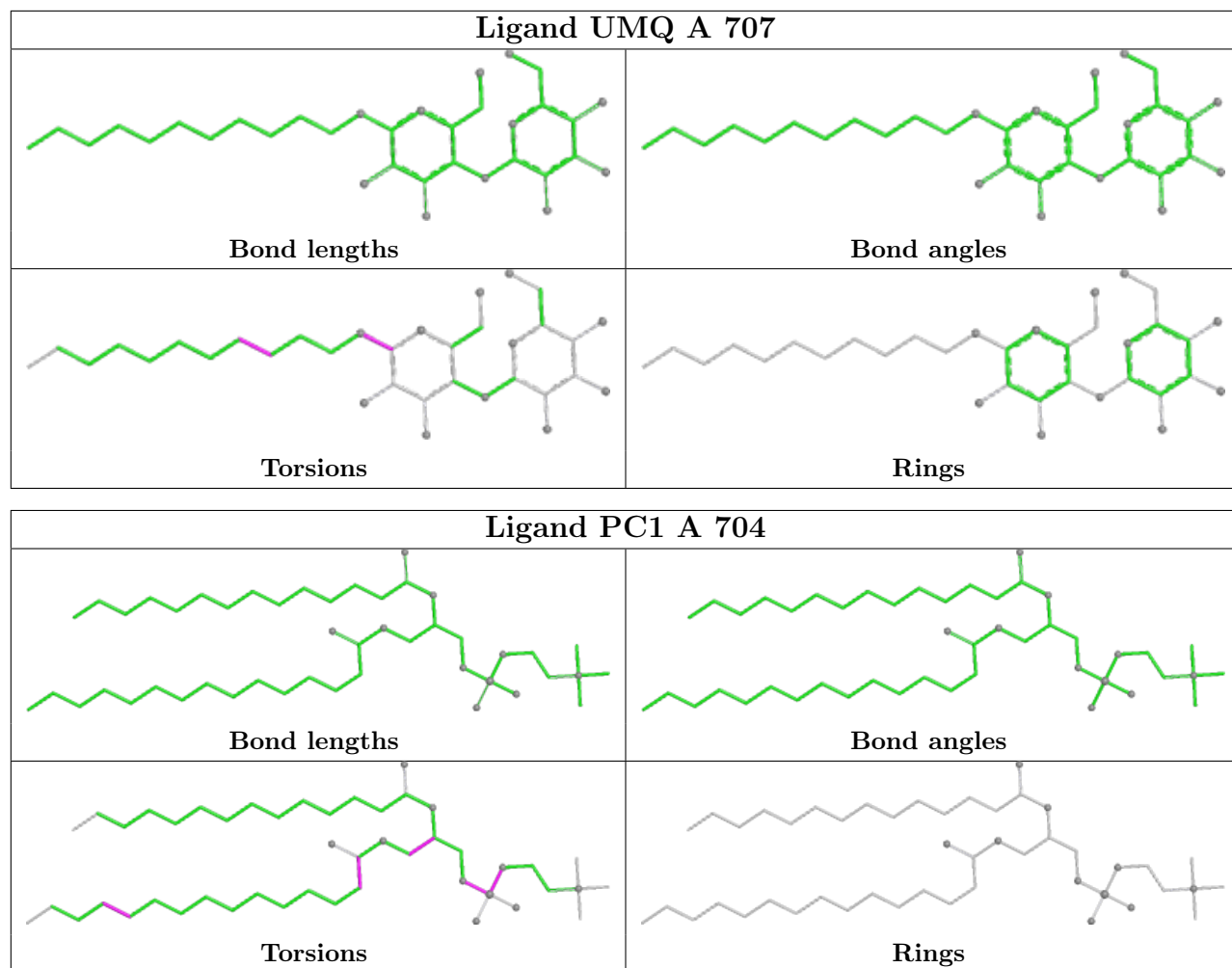
average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents 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.

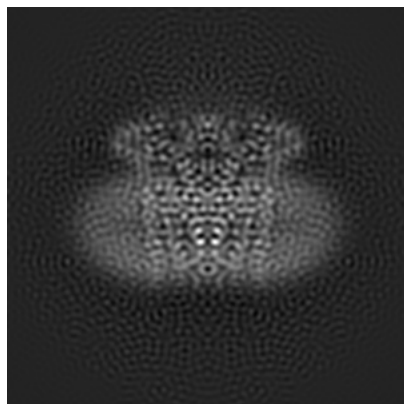
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-4746. These allow visual inspection of the internal detail of the map and identification of artifacts.

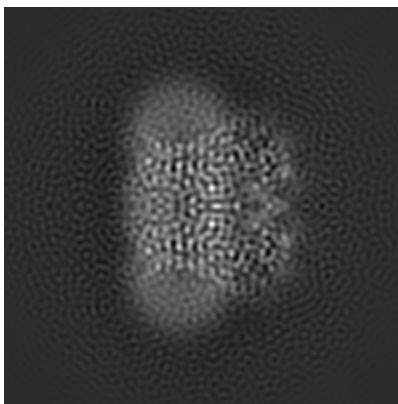
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

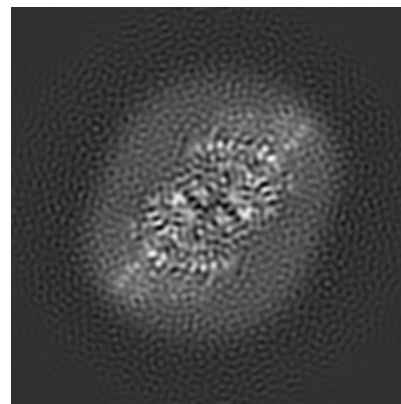
#### 6.1.1 Primary map



X

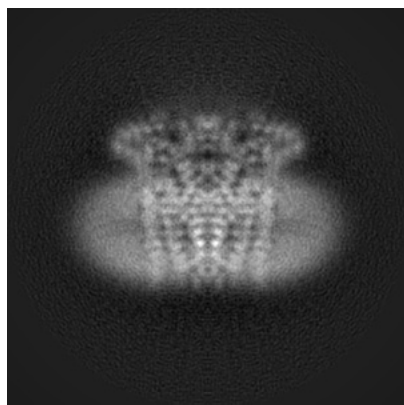


Y

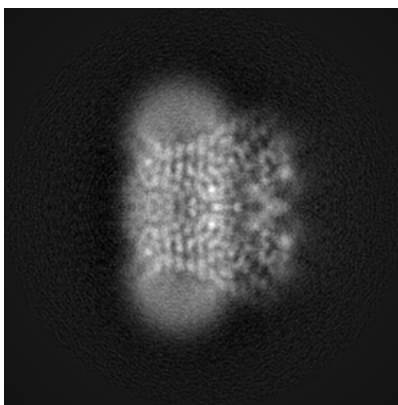


Z

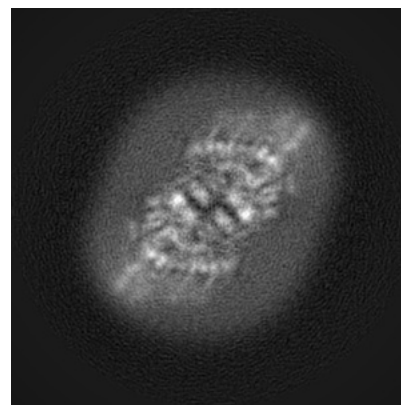
#### 6.1.2 Raw map



X



Y

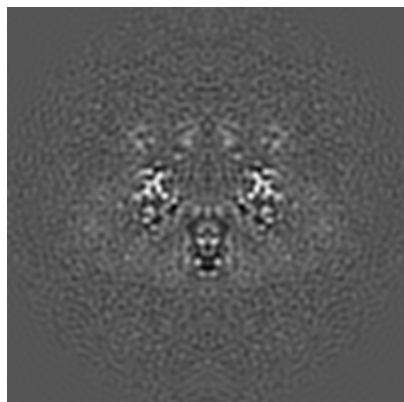


Z

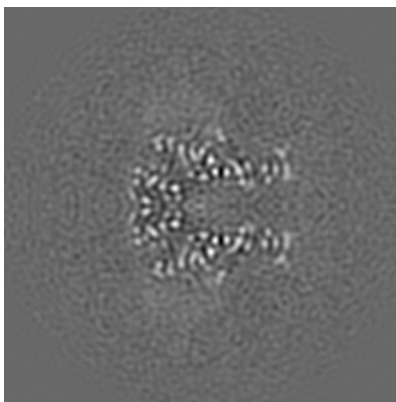
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

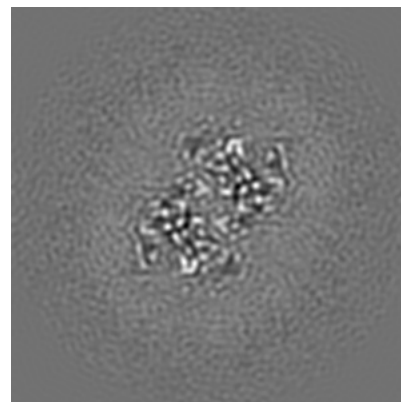
### 6.2.1 Primary map



X Index: 114

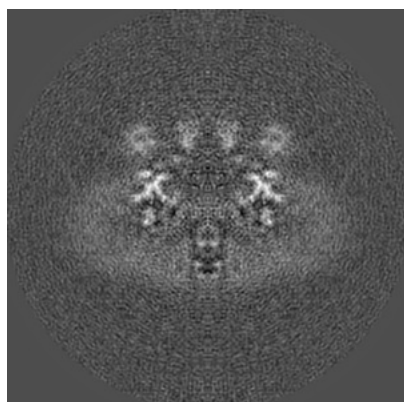


Y Index: 114

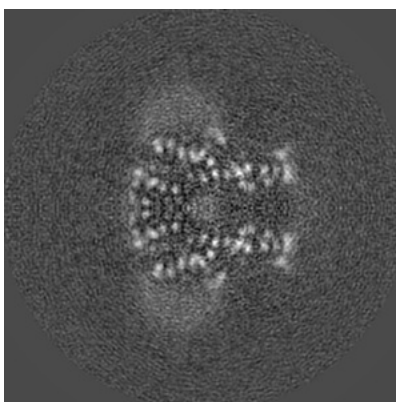


Z Index: 114

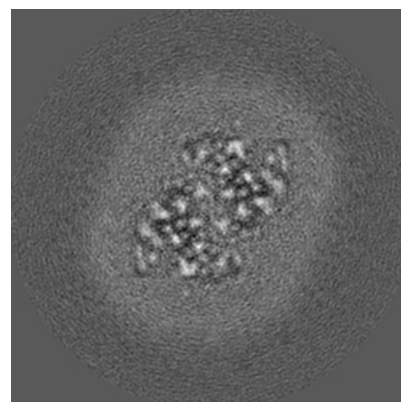
### 6.2.2 Raw map



X Index: 114



Y Index: 114

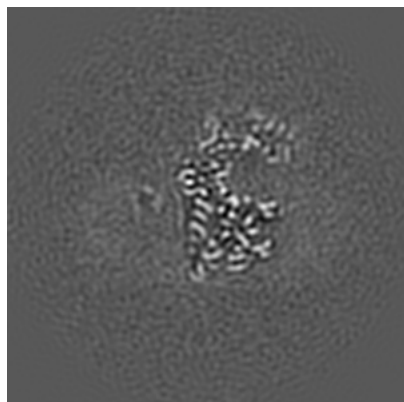


Z Index: 114

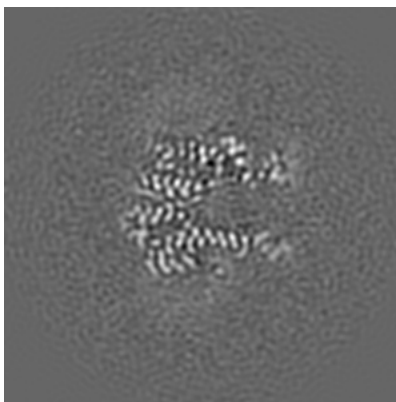
The images above show central slices of the map in three orthogonal directions.

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

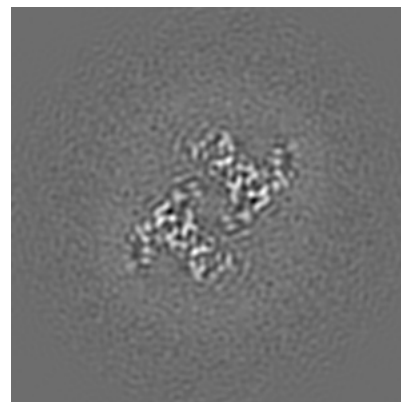
### 6.3.1 Primary map



X Index: 130

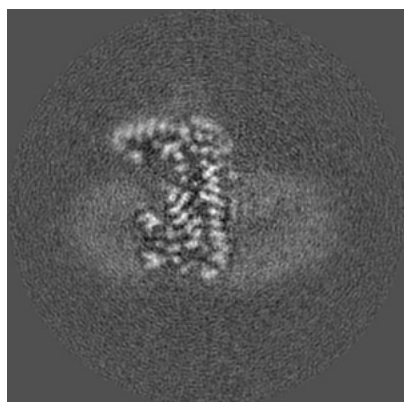


Y Index: 118

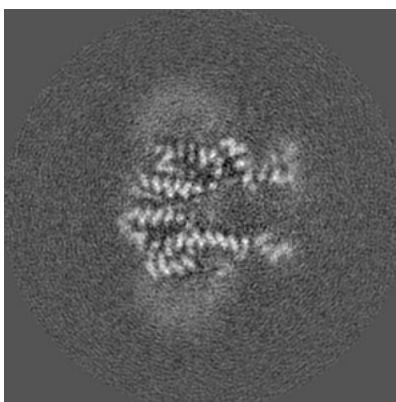


Z Index: 118

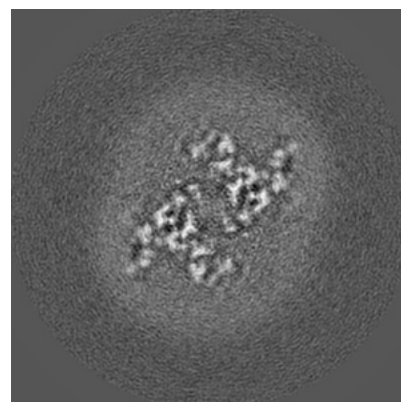
### 6.3.2 Raw map



X Index: 94



Y Index: 118



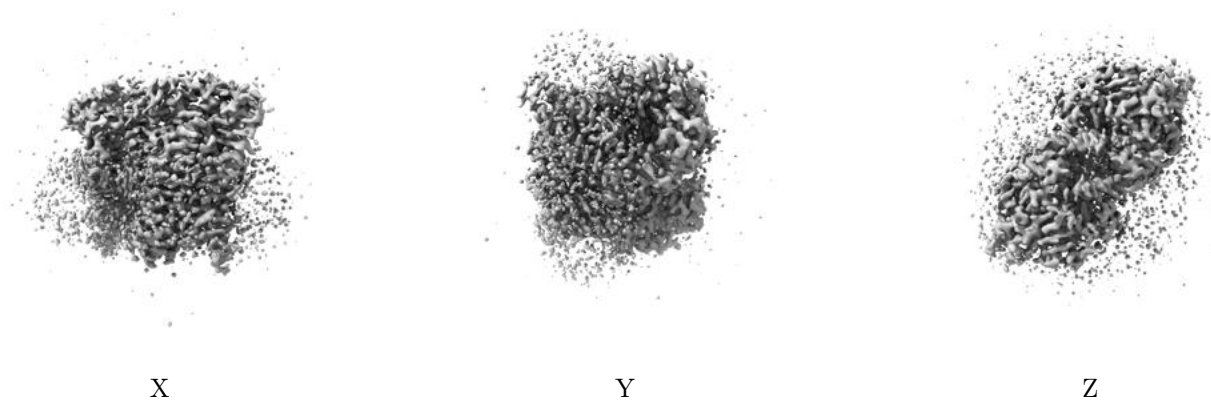
Z Index: 118

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



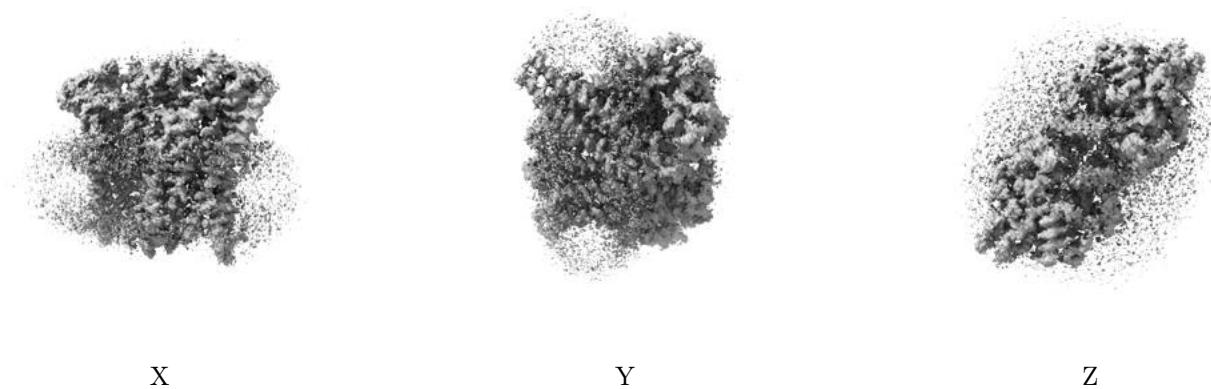
## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.017. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.4.2 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.

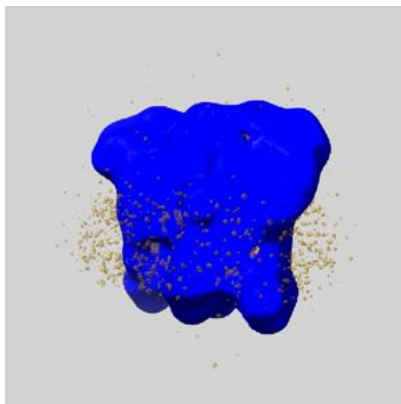
## 6.5 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

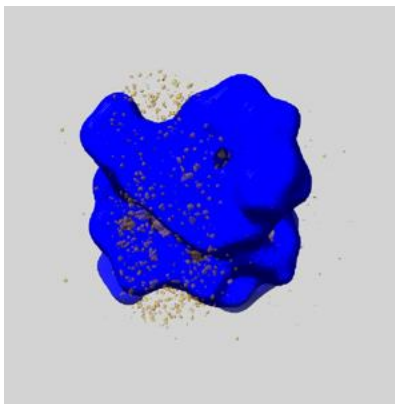
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

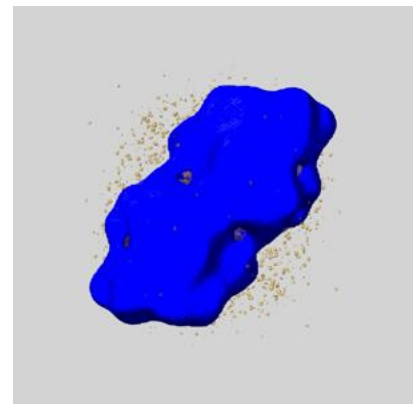
### 6.5.1 emd\_4746\_msk\_1.map [i](#)



X



Y



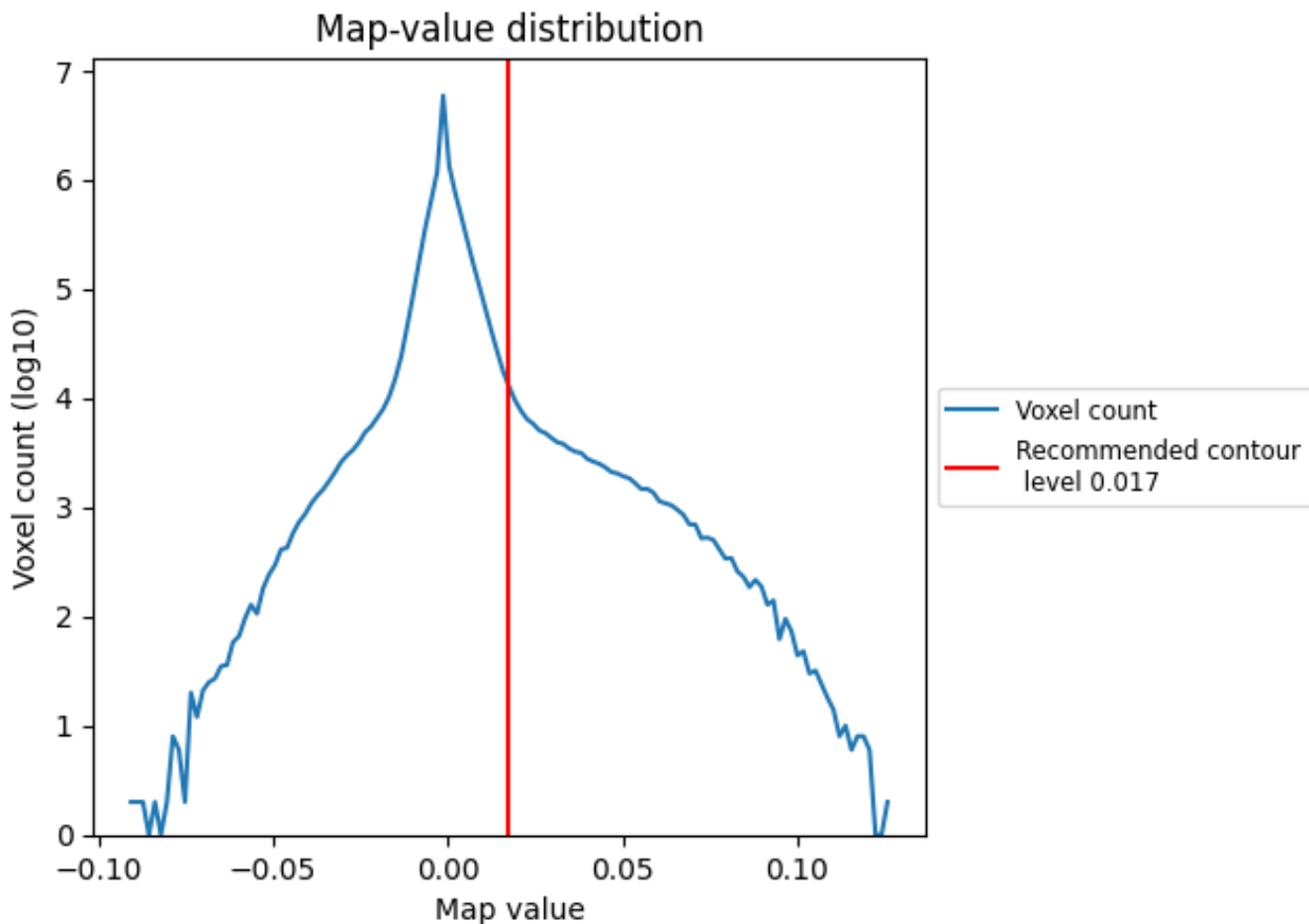
Z



## 7 Map analysis [i](#)

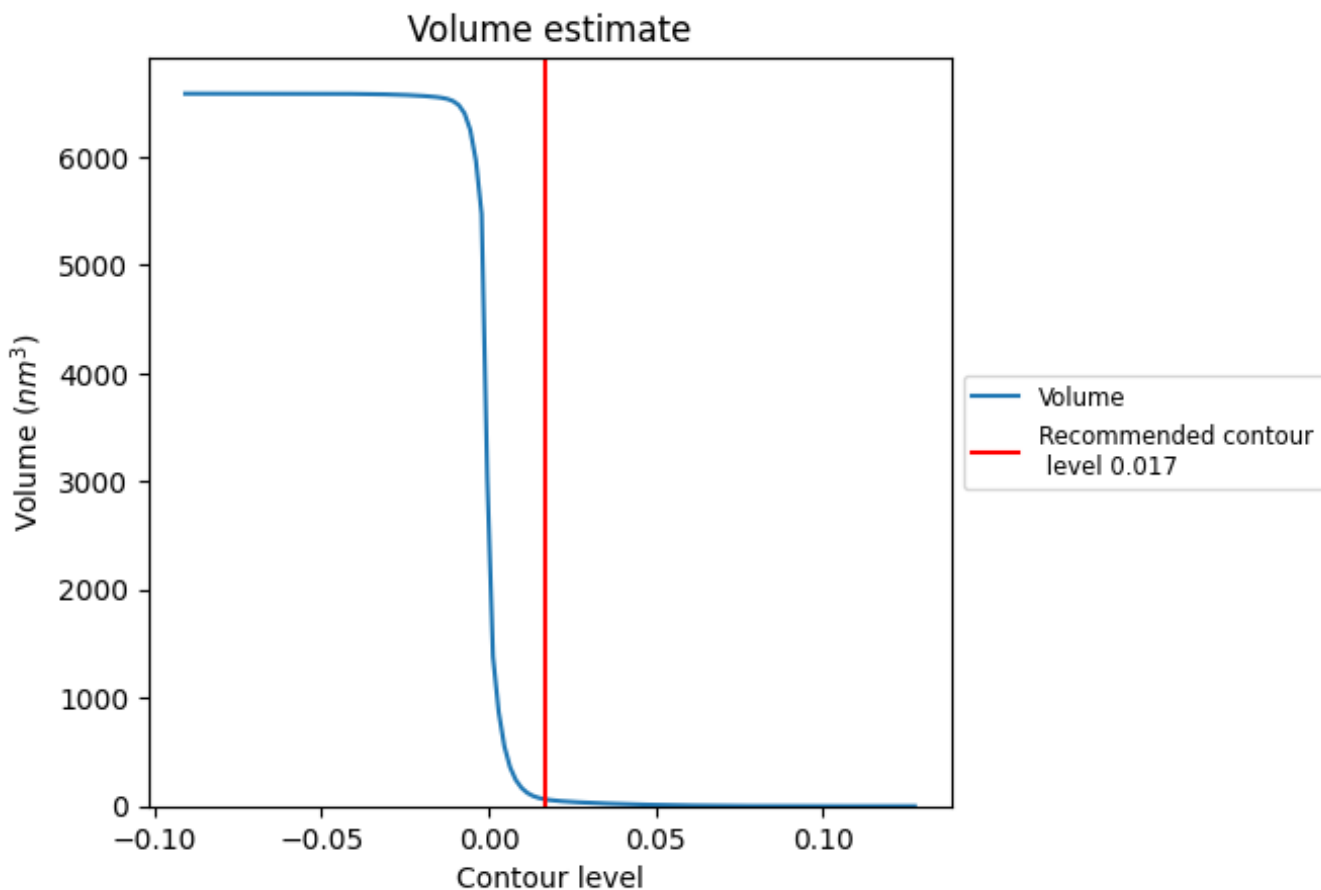
This section contains the results of statistical analysis of the map.

### 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.

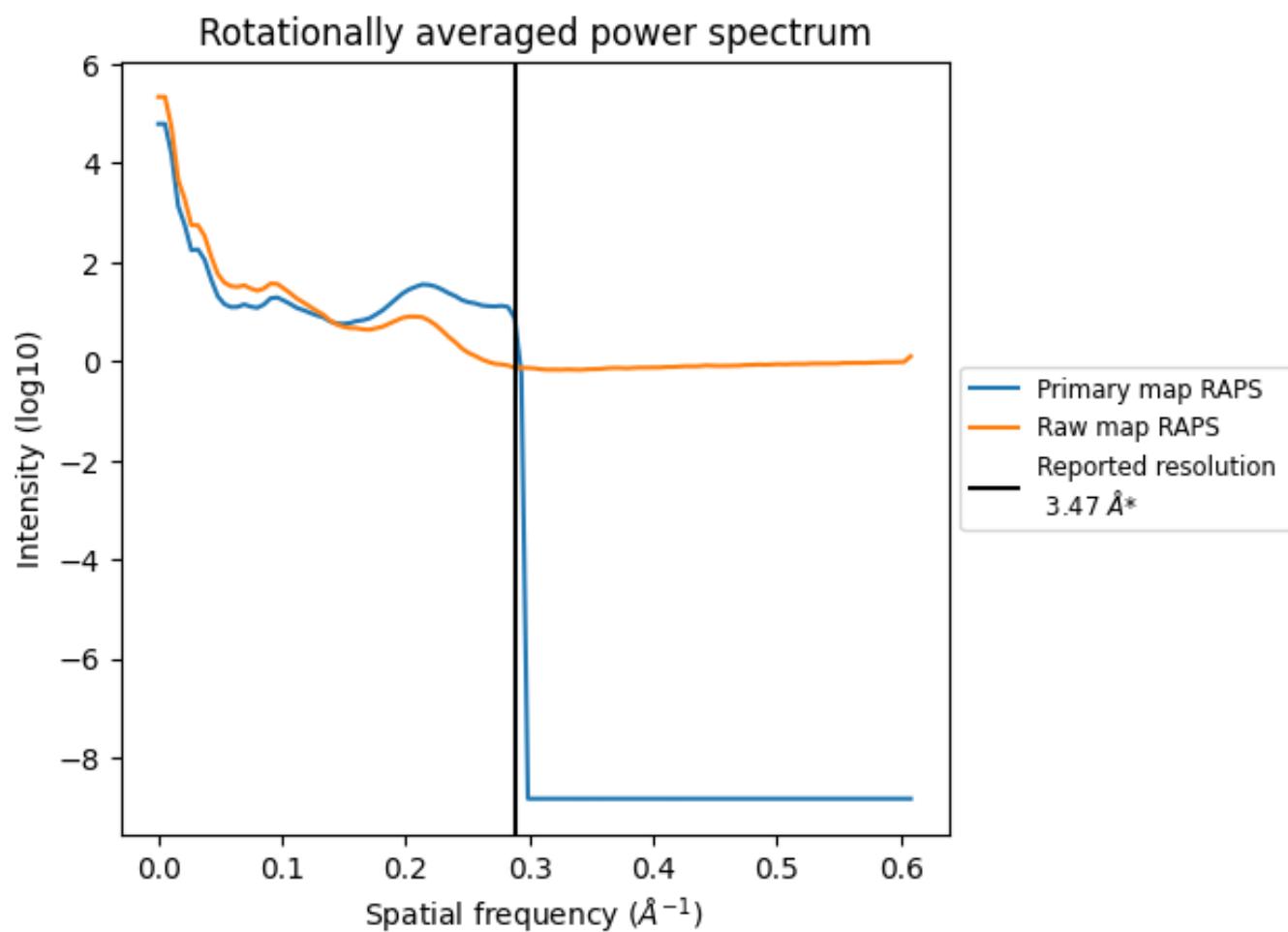
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 63 nm<sup>3</sup>; this corresponds to an approximate mass of 57 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

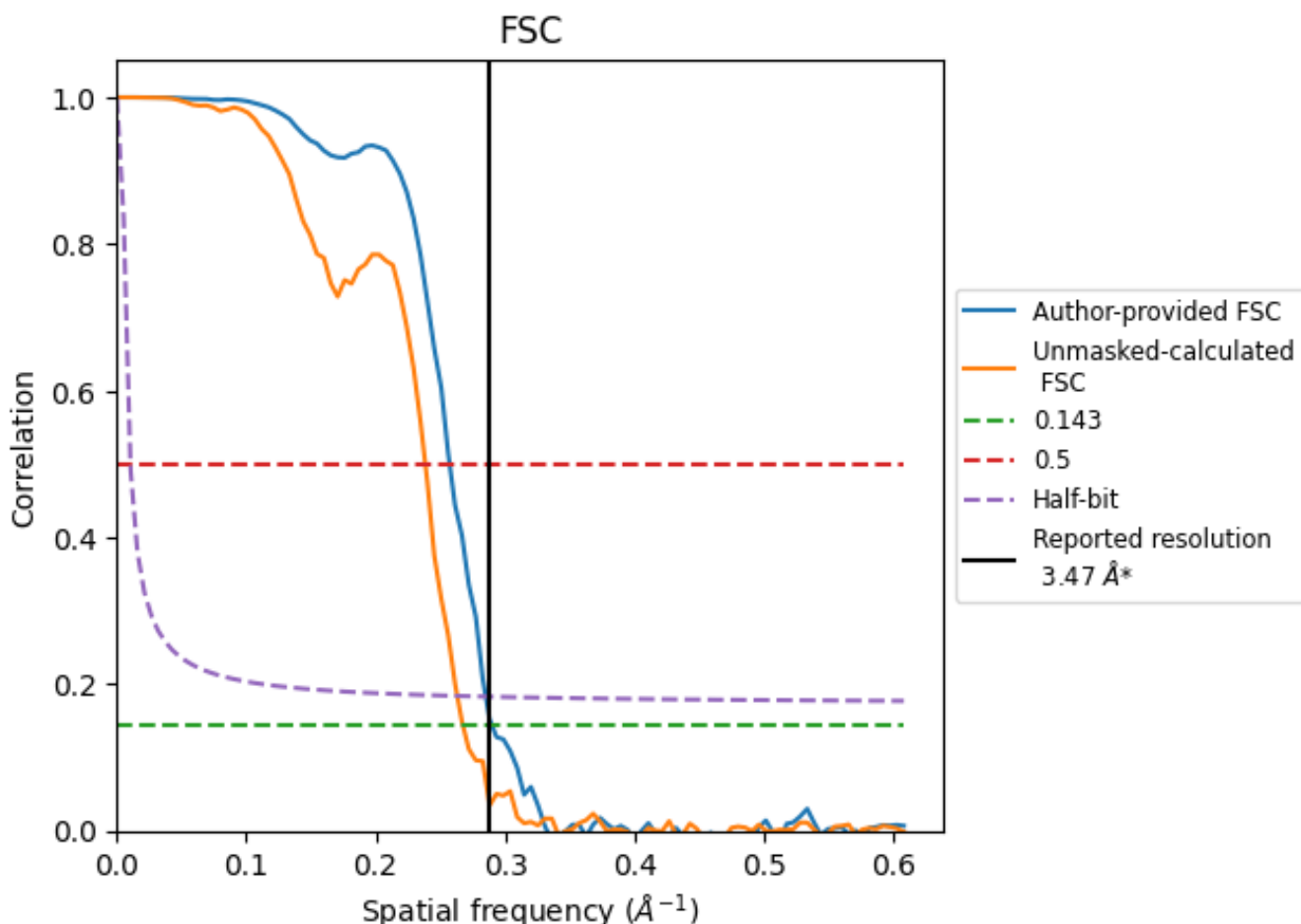


\*Reported resolution corresponds to spatial frequency of 0.288 Å<sup>-1</sup>

## 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 0.288 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

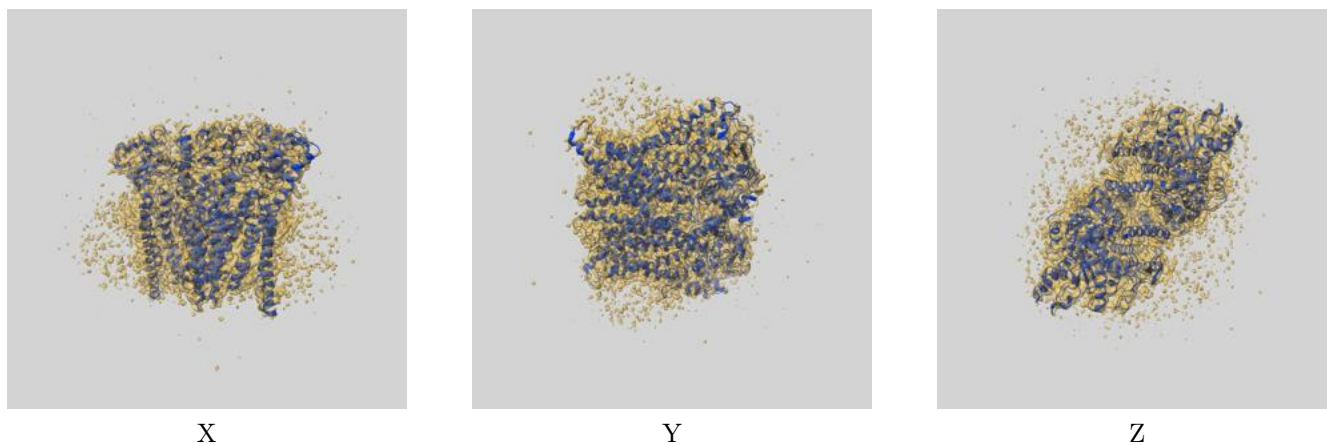
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.47	-	-
Author-provided FSC curve	3.45	3.89	3.51
Unmasked-calculated*	3.74	4.19	3.80

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

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

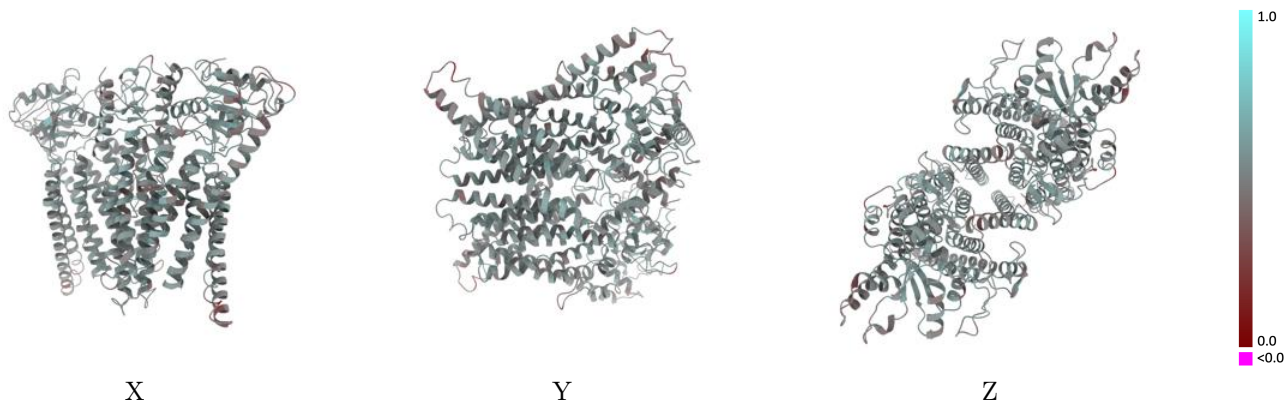
This section contains information regarding the fit between EMDB map EMD-4746 and PDB model 6R7X. Per-residue inclusion information can be found in section 3 on page 6.

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



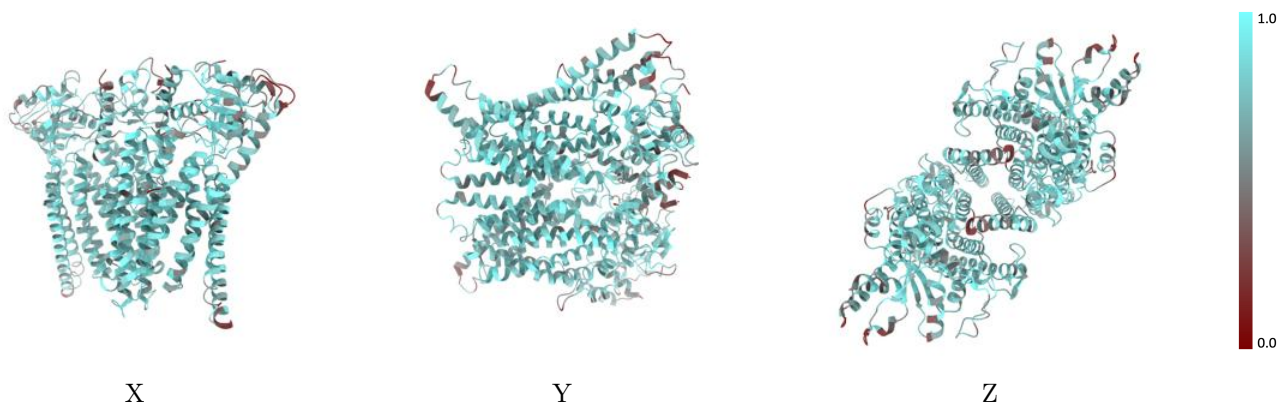
The images above show the 3D surface view of the map at the recommended contour level 0.017 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



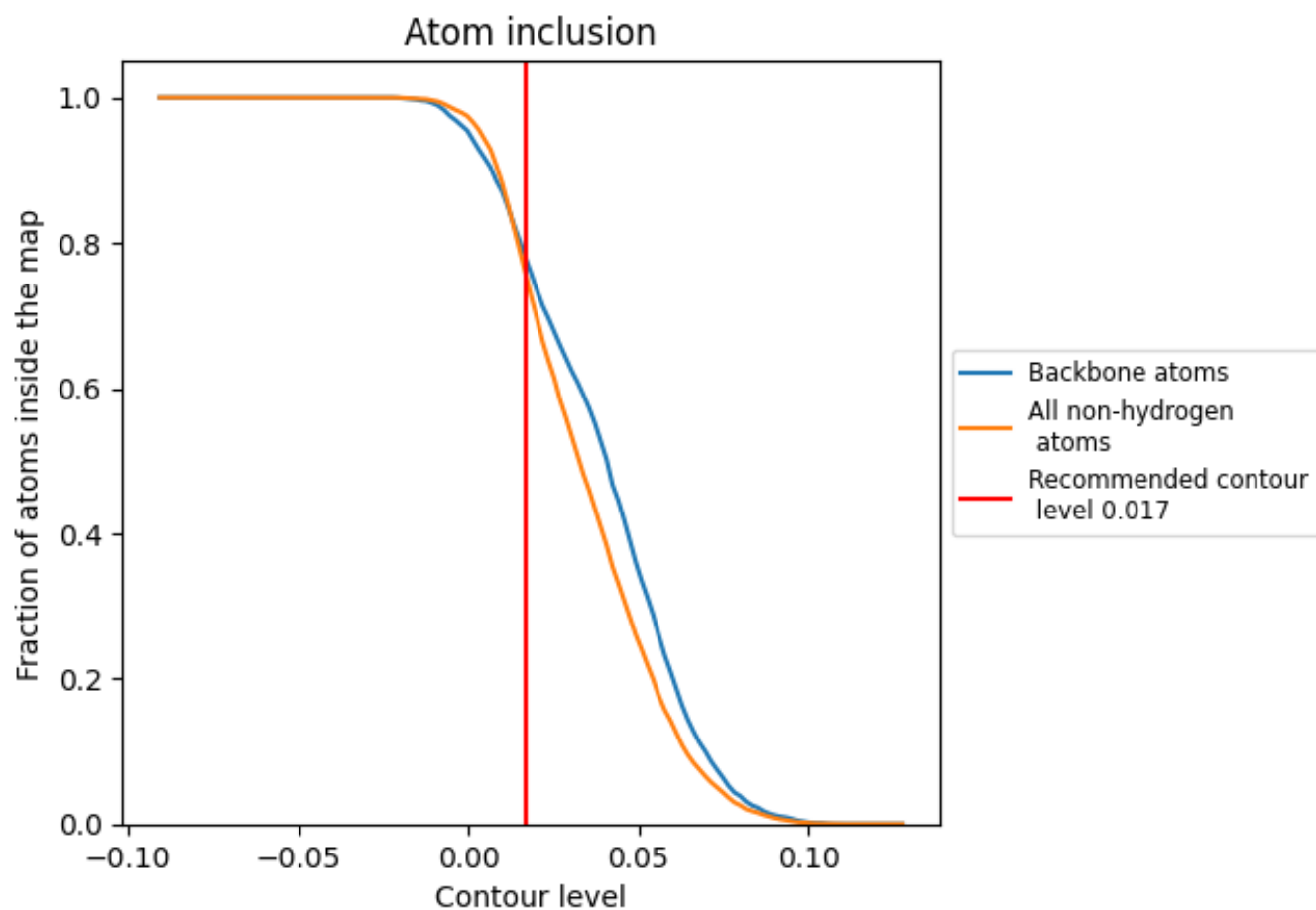
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.017).

## 9.4 Atom inclusion [i](#)







At the recommended contour level, 78% of all backbone atoms, 75% 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 (0.017) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7524	 0.5160
A	 0.7528	 0.5170
B	 0.7521	 0.5160

