



wwPDB EM Validation Summary Report ⓘ

Nov 10, 2024 – 06:36 am GMT

PDB ID : 8A7D
EMDB ID : EMD-15220
Title : Partial dimer complex of PAPP-A and its inhibitor STC2
Authors : Kobbero, S.D.; Gajhede, M.; Mirza, O.A.; Boesen, T.; Oxvig, C.
Deposited on : 2022-06-20
Resolution : 3.06 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev113
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

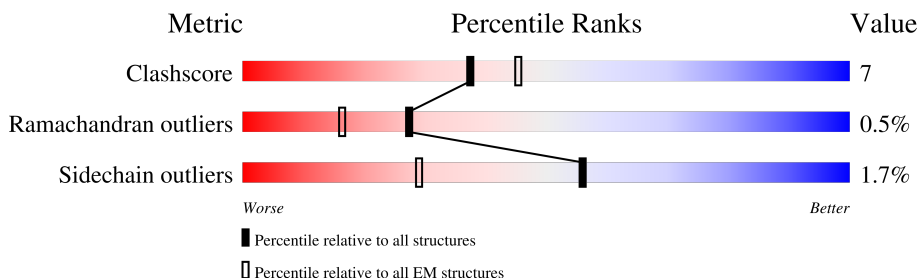
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.06 Å.

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



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	C	1536	
1	Q	1536	
2	P	167	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10771 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 Pappalysin-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	921	Total	C	N	O	S	0	0
			7256	4557	1247	1399	53		
1	Q	273	Total	C	N	O	S	0	0
			2085	1288	370	396	31		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	563	GLN	GLU	engineered mutation	UNP Q13219
Q	563	GLN	GLU	engineered mutation	UNP Q13219

- Molecule 2 is a protein called Stanniocalcin-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	P	167	Total	C	N	O	S	0	0
			1309	819	236	238	16		

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
4	C	1	Total	Zn	0
			1	1	

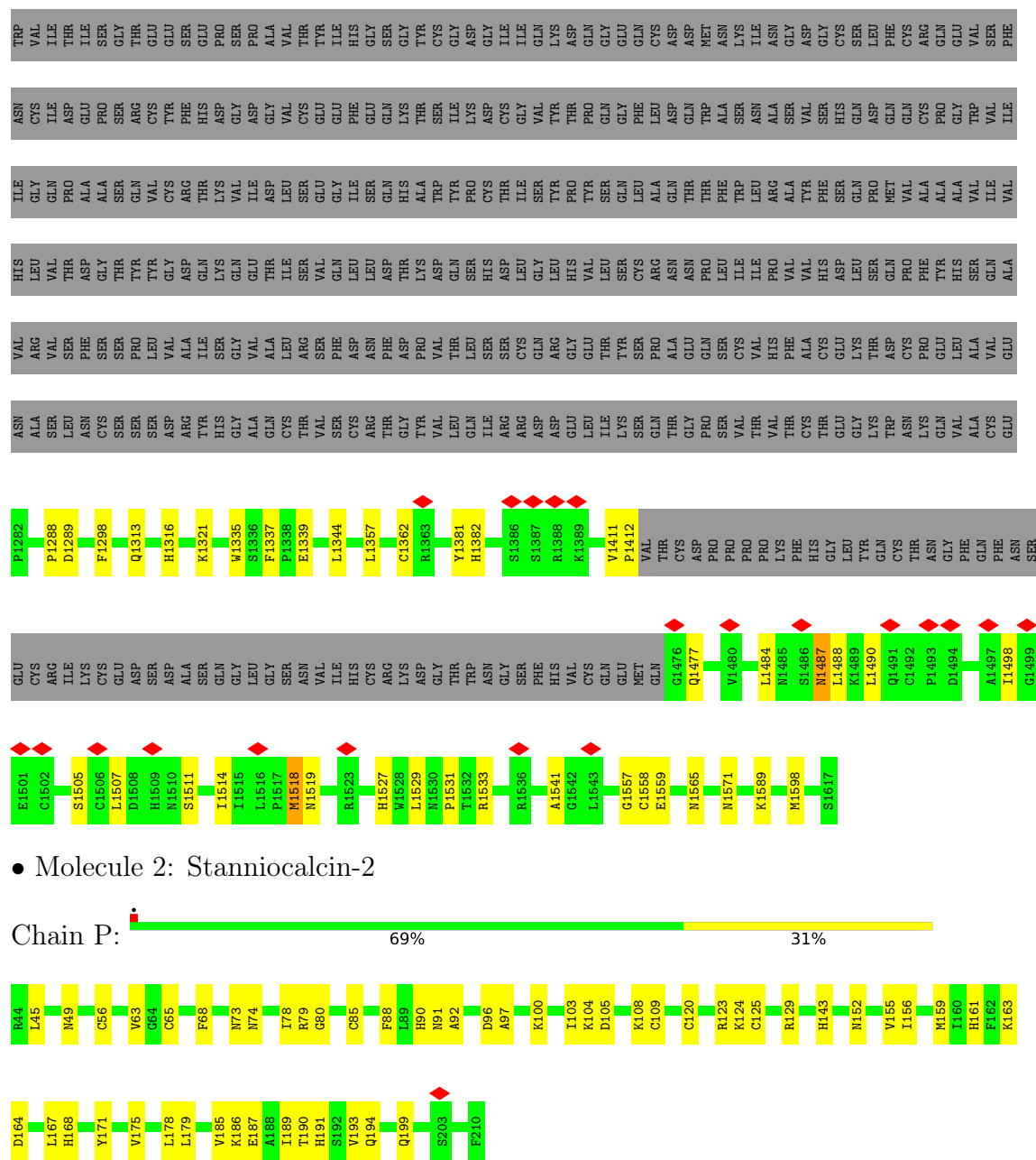
- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
5	C	7	Total 7	Ca 7	0
5	Q	1	Total 1	Ca 1	0

PHE	ARG	GLY	ASP	ASP
PRO	VAL	GLN	PRO	PRO
MET	VAL	CYS	PRO	GLN
SER	CYS	SER	THR	THR
CYS	THR	VAL	PRO	ALA
ASP	ALA	PRO	LYS	ARG
LEU	GLY	ASN	PHE	CYS
GLN	LEU	GLU	HIS	CYS
LEU	LYS	LEU	GLY	GLU
GLY	LYS	LEU	ASN	ASN
ASP	TRP	ASN	LEU	THR
CYS	TYR	SER	TYR	LYS
ALA	PRO	ASN	GLN	HIS
ALA	HIS	LEU	CYS	LYS
CYS	PRO	LYS	CYS	VAL
ARG	ALA	LEU	GLY	GLY
ASP	ALA	LEU	ASN	THR
PRO	LEU	GLN	PHE	SER
GLN	ILE	CYS	GLY	CYS
ALA	HIS	PRO	GLN	CYS
ALA	CYS	ASP	PHE	LYS
GLN	TYR	GLY	ASN	TYR
GLU	VAL	TYR	SER	CYS
HIS	GLY	ALA	GLU	LYS
SER	CYS	ILE	CYS	LYS
	GLU	GLY	ARG	PRO
	PRO	SER	ILE	GLY
	PHE	GLU	LYS	TYR
	MET	CYS	CYS	HIS
	GLY	THR	ASP	VAL
	ASP	THR	GLU	VAL
	ASN	SER	SER	GLY
	ILE	ASN	GLY	SER
	ALA	ASP	GLN	LYS
	ALA	HIS	GLN	LYS
	ILE	ASN	GLY	SER
	ASN	SER	LEU	LYS
	ASN	GLU	GLY	LYS
	ASN	CYS	SER	ARG
	ARG	ALA	ASN	ALA
	ALA	ILE	VAL	PHE
	PHE	ILE	ILE	LYS
	CYS	LEU	ILE	THR
	ASN	PRO	HIS	THR
	TYR	MET	CYS	GLN
	ASP	ASN	ARG	CYS
	GLY	VAL	LYS	THR
	GLY	THR	ASP	GLN
	ASP	VAL	GLY	ASP
	CYS	ARG	THR	GLY
	CYS	ASP	SER	SER
	THR	ILE	TRP	GLN
	SER	PRO	GLY	THR
	THR	HIS	SER	GLU
	VAL	TRP	PHE	GLY
	LYS	LEU	HIS	ALA
	THR	ASN	VAL	CYS
	PRO	PRO	CYS	VAL
	LYS	THR	GLN	PRO
	VAL	ARG	GLU	VAL
	THR	VAL	MET	THR
	PRO	GLN	GLN	CYS

- Molecule 1: Pappalysin-1

[illegible]



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	278982	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 ($e^-/\text{\AA}^2$)	59, 58	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k), GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.150	Depositor
Minimum map value	-0.794	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.027	Depositor
Recommended contour level	0.05	Depositor
Map size (\AA)	223.08002, 223.08002, 223.08002	wwPDB
Map dimensions	220, 220, 220	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.014, 1.014, 1.014	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: ZN, CA, 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	C	0.32	1/7454 (0.0%)	0.58	2/10147 (0.0%)
1	Q	0.27	0/2140	0.47	0/2905
2	P	0.30	0/1331	0.51	0/1791
All	All	0.31	1/10925 (0.0%)	0.55	2/14843 (0.0%)

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	C	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	458	ASP	N-CA	7.95	1.62	1.46

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	457	CYS	O-C-N	-21.00	89.11	122.70
1	C	457	CYS	CA-C-N	14.16	148.36	117.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	457	CYS	Mainchain

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Mol	Chain	Res	Type	Group
1	C	777	GLU	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	C	7256	0	6834	97	0
1	Q	2085	0	1961	26	0
2	P	1309	0	1286	36	0
3	C	98	0	91	1	0
3	Q	14	0	13	0	0
4	C	1	0	0	0	0
5	C	7	0	0	0	0
5	Q	1	0	0	0	0
All	All	10771	0	10185	152	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

The worst 5 of 152 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:74:ASN:HD21	2:P:78:ILE:HB	1.49	0.78
1:C:562:HIS:CE1	1:C:572:HIS:HE2	1.97	0.76
2:P:73:ASN:H	2:P:79:ARG:HH12	1.37	0.73
1:Q:1381:TYR:HA	1:Q:1412:PRO:HA	1.71	0.72
2:P:91:ASN:HD22	2:P:159:MET:HG2	1.53	0.71

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	C	919/1536 (60%)	853 (93%)	59 (6%)	7 (1%)	16	43
1	Q	269/1536 (18%)	261 (97%)	8 (3%)	0	100	100
2	P	165/167 (99%)	160 (97%)	5 (3%)	0	100	100
All	All	1353/3239 (42%)	1274 (94%)	72 (5%)	7 (0%)	27	54

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	430	HIS
1	C	808	ASP
1	C	459	MET
1	C	726	ALA
1	C	737	HIS

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	809/1347 (60%)	798 (99%)	11 (1%)	62	79
1	Q	238/1347 (18%)	232 (98%)	6 (2%)	42	66
2	P	145/145 (100%)	142 (98%)	3 (2%)	48	69
All	All	1192/2839 (42%)	1172 (98%)	20 (2%)	56	75

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	Q	1316	HIS
1	Q	1518	MET
1	Q	1558	CYS
1	Q	1527	HIS
1	C	443	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	905	ASN
1	Q	1553	HIS
2	P	91	ASN
1	Q	1398	GLN
2	P	73	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 9 are monoatomic - leaving 8 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	NAG	C	1704	1	14,14,15	0.28	0	17,19,21	0.50	0
3	NAG	C	1701	1	14,14,15	0.21	0	17,19,21	0.38	0
3	NAG	C	1705	1	14,14,15	0.28	0	17,19,21	0.52	0
3	NAG	C	1702	1	14,14,15	0.17	0	17,19,21	0.36	0
3	NAG	C	1703	1	14,14,15	1.01	1 (7%)	17,19,21	1.12	2 (11%)
3	NAG	Q	1701	1	14,14,15	0.32	0	17,19,21	0.37	0
3	NAG	C	1706	1	14,14,15	0.18	0	17,19,21	0.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	C	1707	1	14,14,15	0.42	0	17,19,21	0.55	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	C	1704	1	-	3/6/23/26	0/1/1/1
3	NAG	C	1701	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1705	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1702	1	-	1/6/23/26	0/1/1/1
3	NAG	C	1703	1	-	4/6/23/26	0/1/1/1
3	NAG	Q	1701	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1706	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1707	1	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1703	NAG	O5-C1	-3.49	1.38	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1703	NAG	C1-C2-N2	-2.50	106.21	110.49
3	C	1703	NAG	C2-N2-C7	2.18	126.01	122.90

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

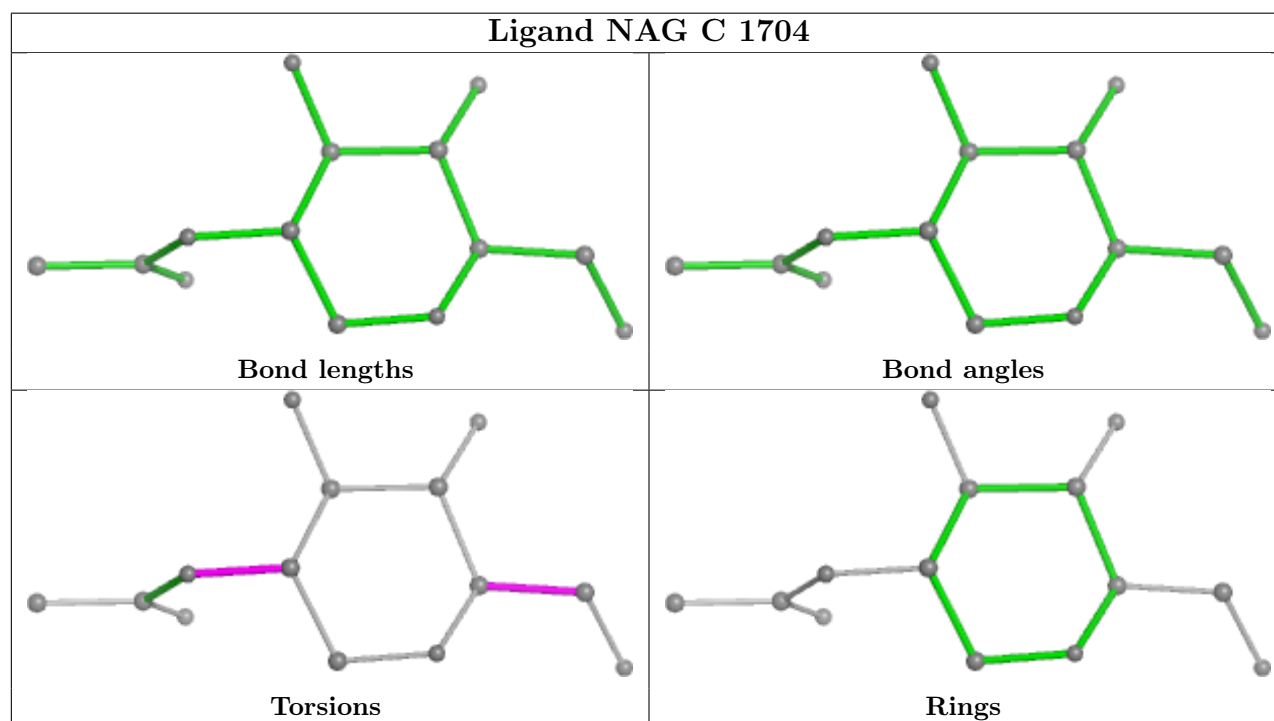
Mol	Chain	Res	Type	Atoms
3	C	1703	NAG	C8-C7-N2-C2
3	C	1703	NAG	O7-C7-N2-C2
3	C	1702	NAG	O5-C5-C6-O6
3	C	1703	NAG	C4-C5-C6-O6
3	C	1703	NAG	O5-C5-C6-O6

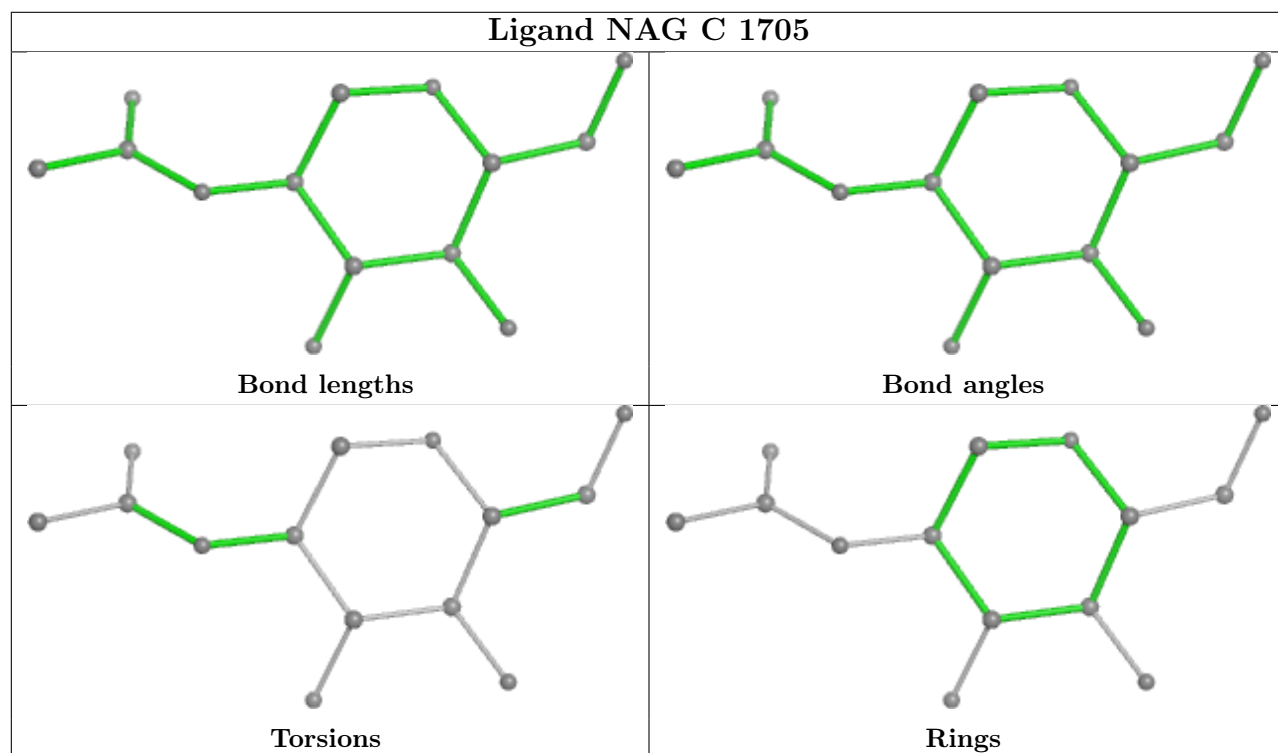
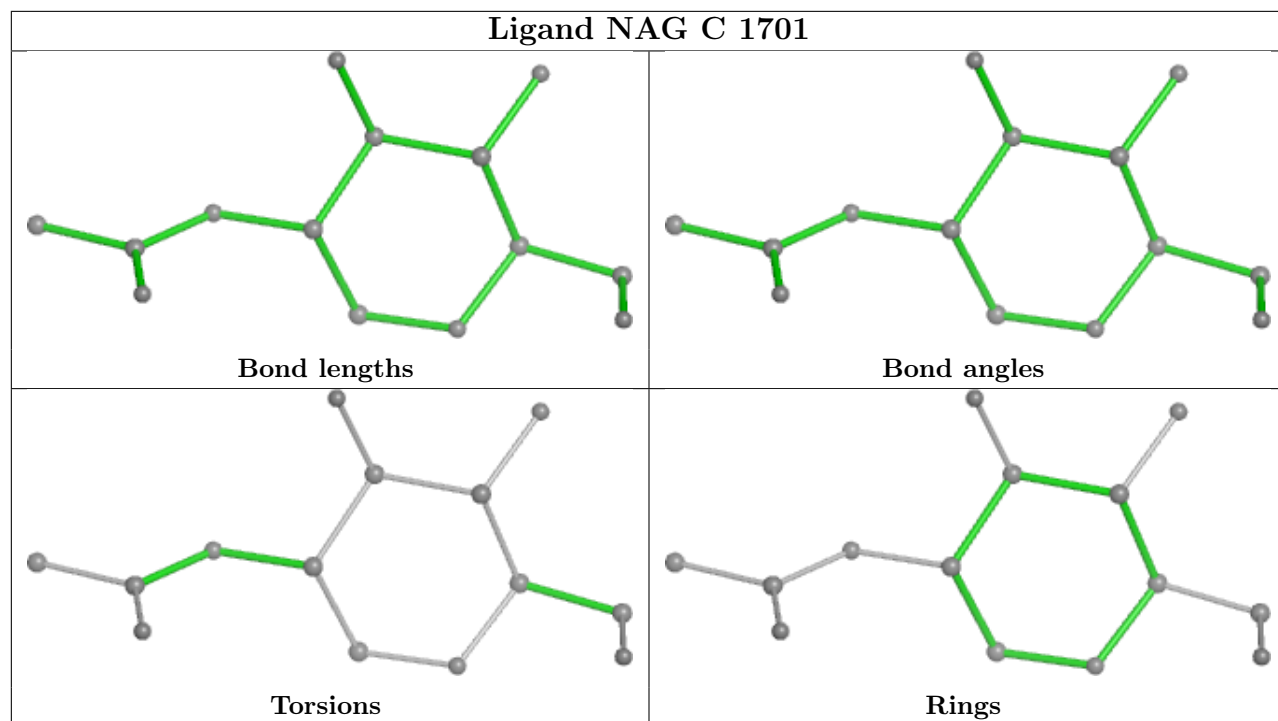
There are no ring outliers.

1 monomer is involved in 1 short contact:

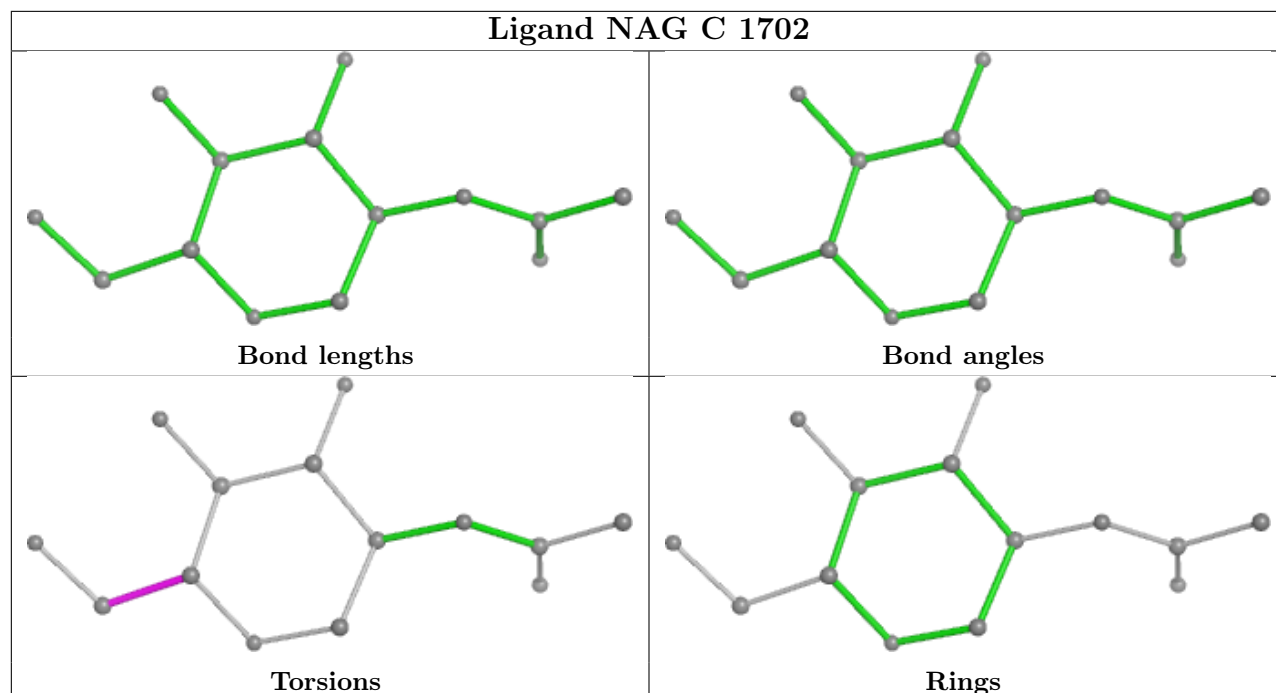
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1703	NAG	1	0

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.

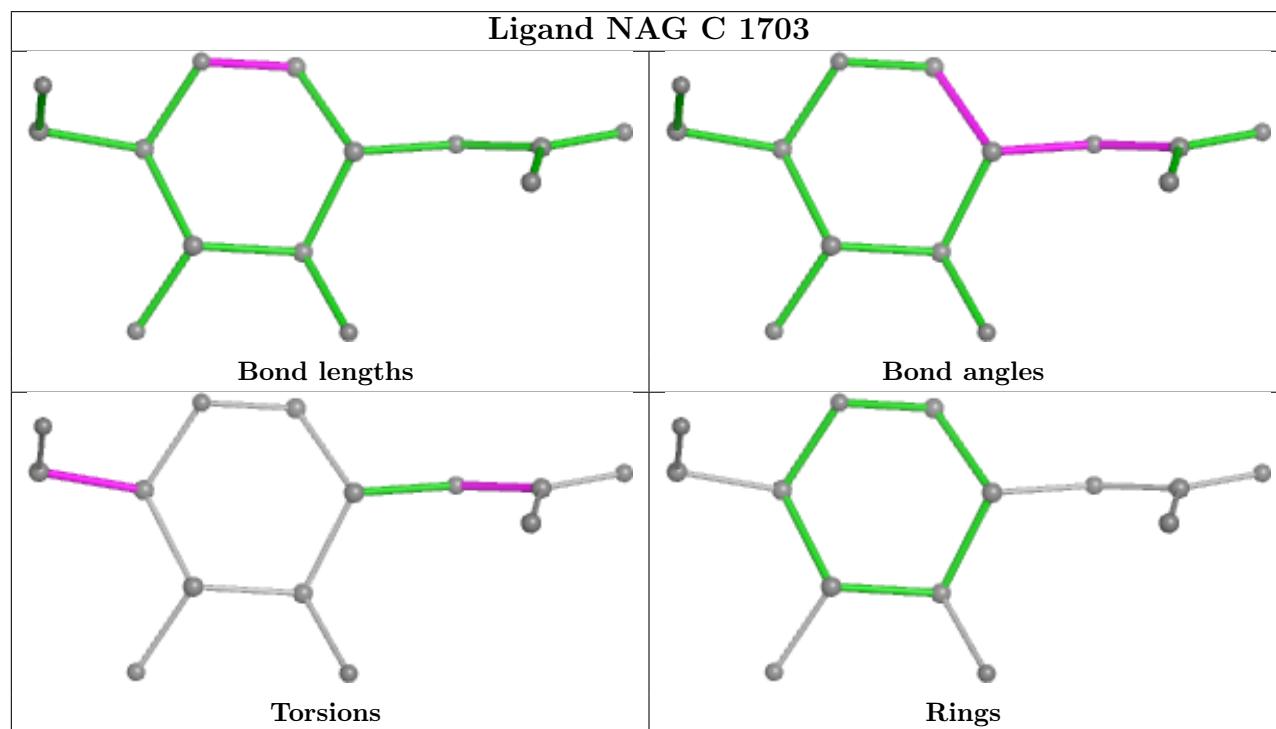


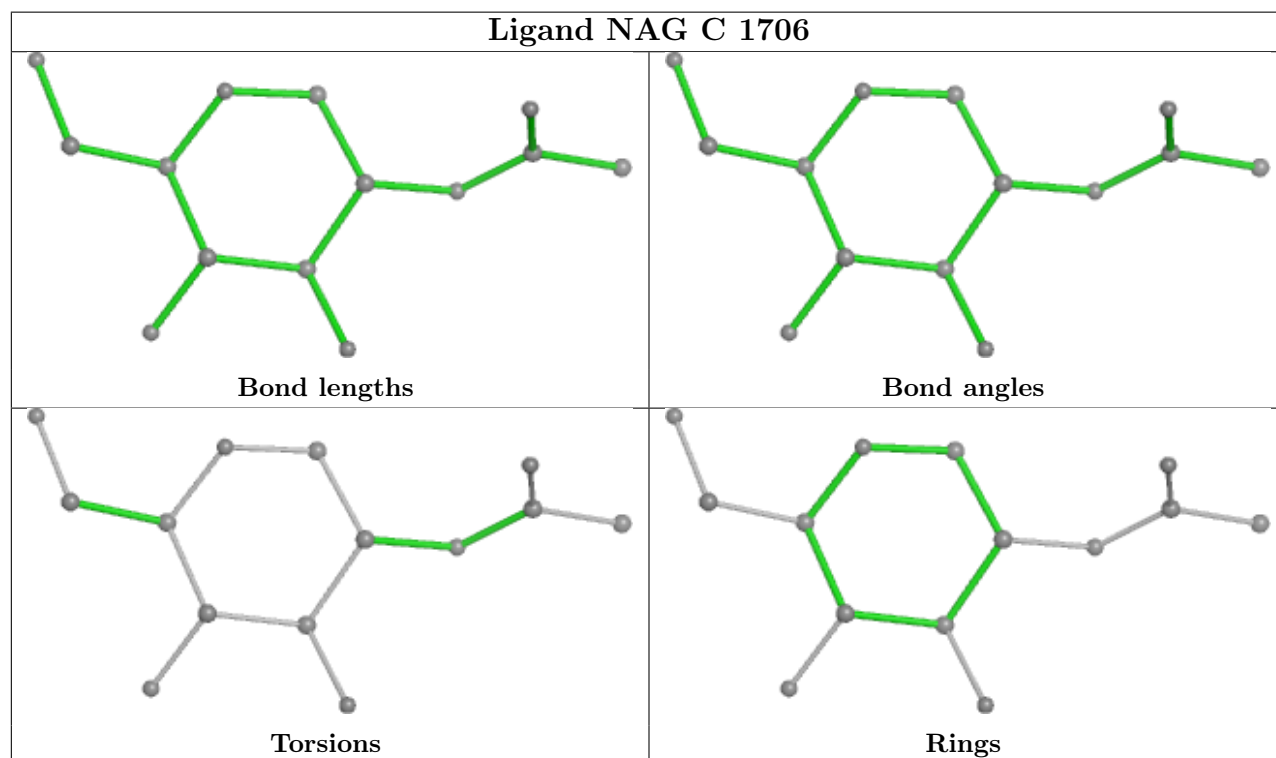
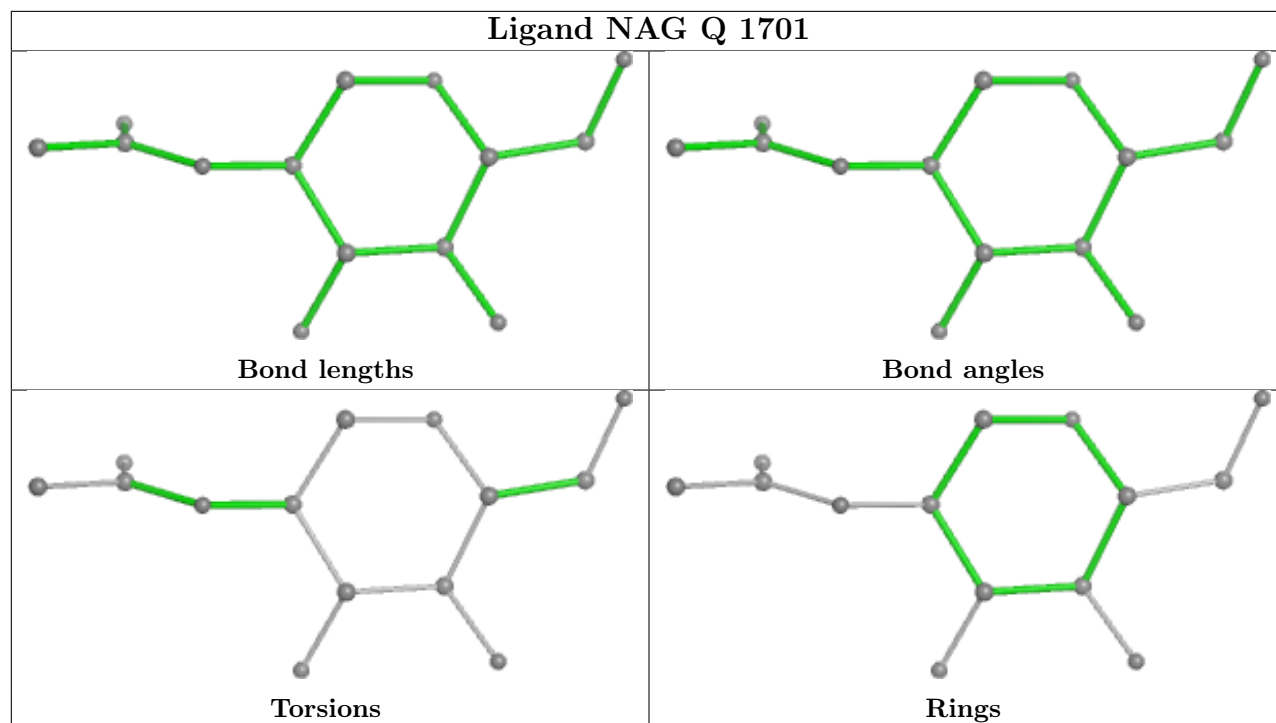


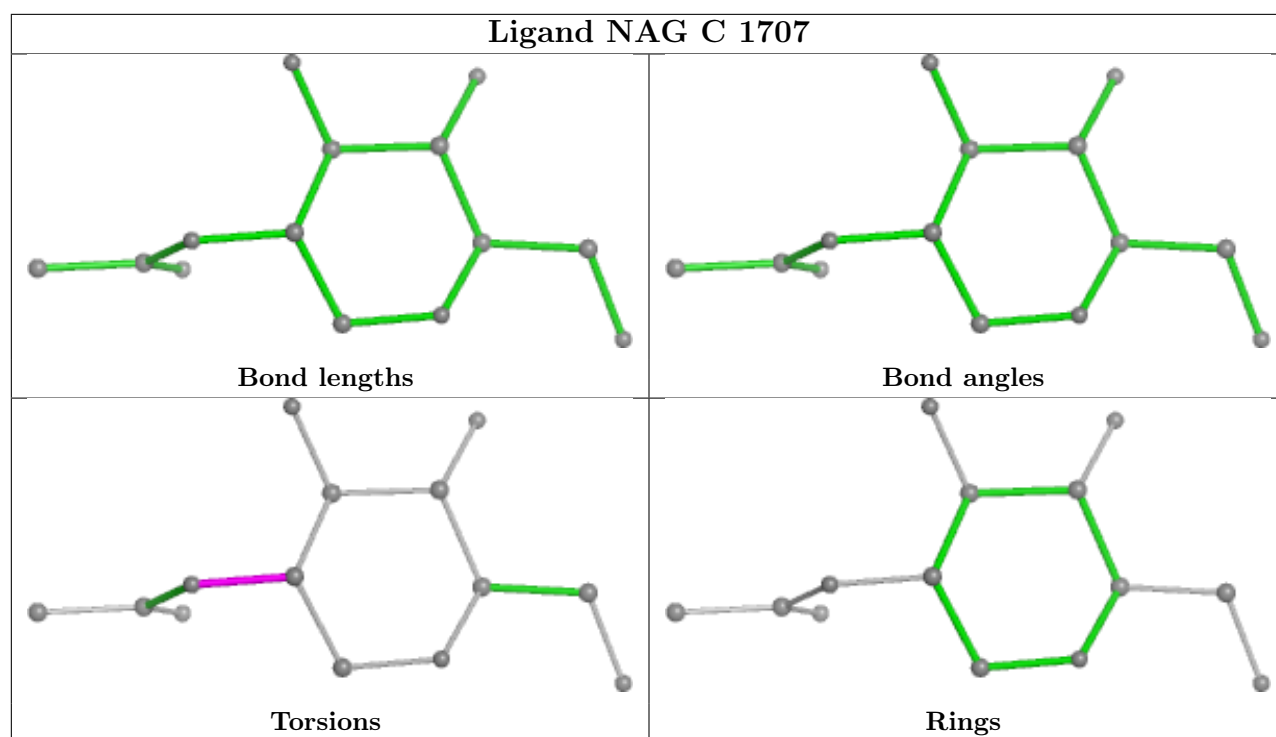
Ligand NAG C 1702



Ligand NAG C 1703







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.

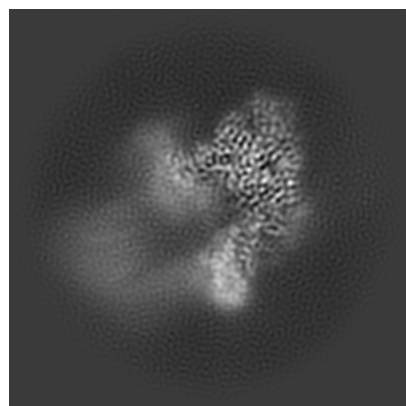
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-15220. These allow visual inspection of the internal detail of the map and identification of artifacts.

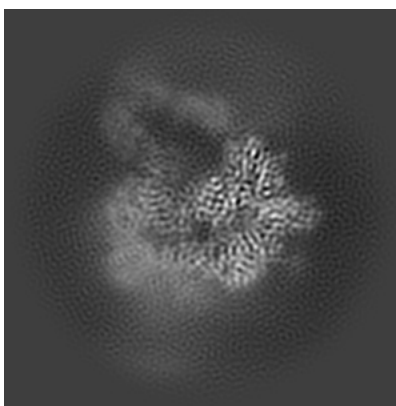
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

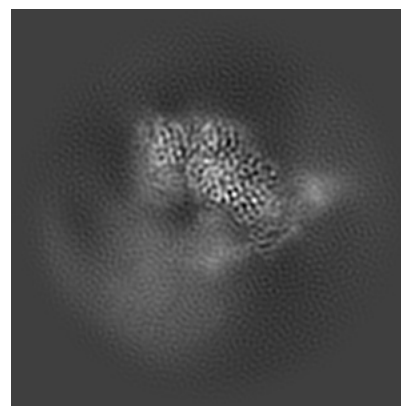
6.1.1 Primary map



X

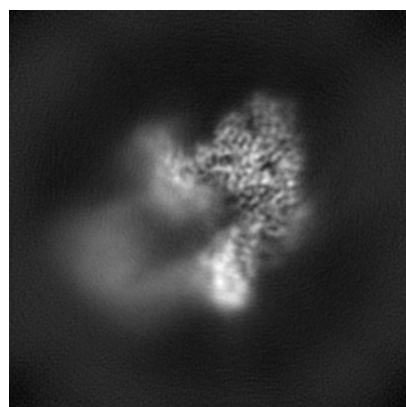


Y

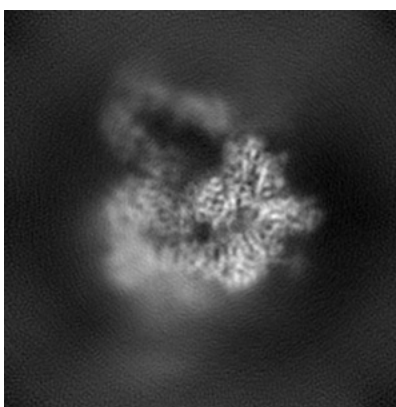


Z

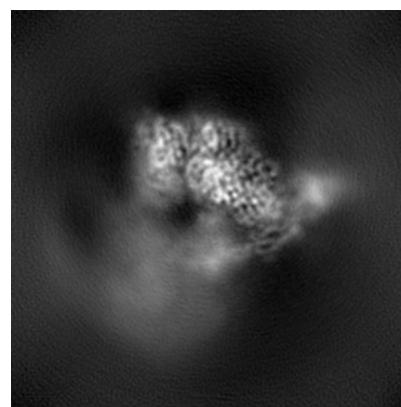
6.1.2 Raw map



X



Y

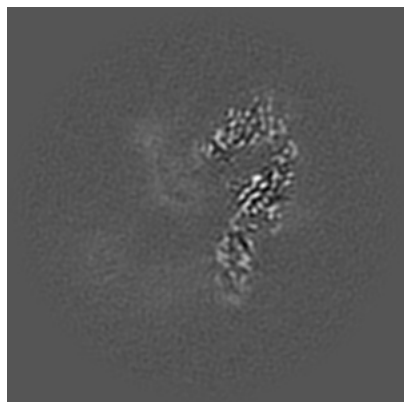


Z

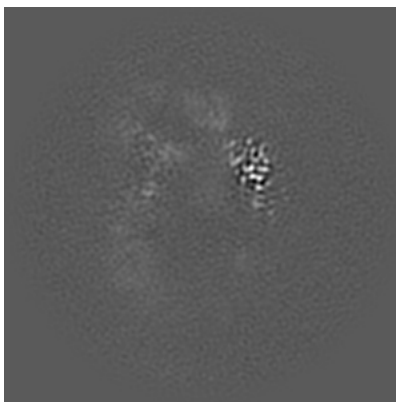
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

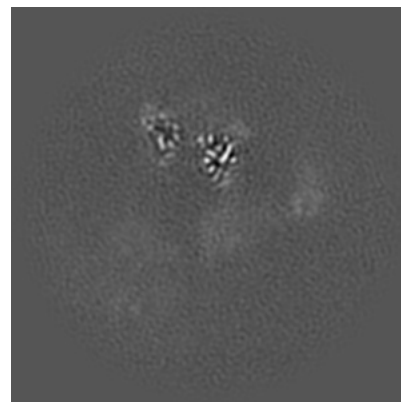
6.2.1 Primary map



X Index: 110

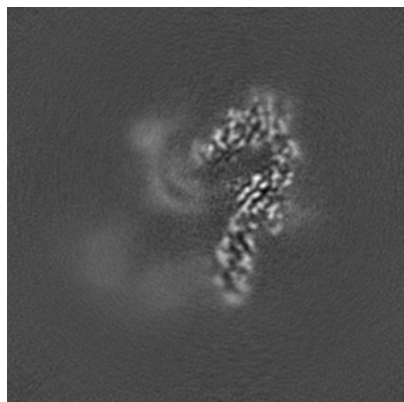


Y Index: 110

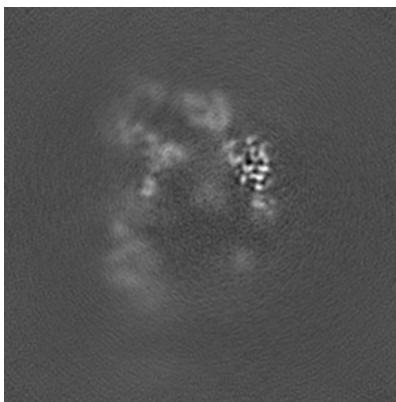


Z Index: 110

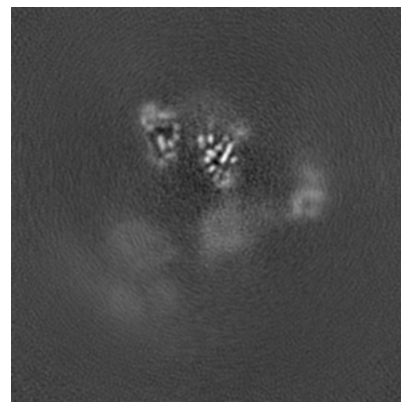
6.2.2 Raw map



X Index: 110



Y Index: 110

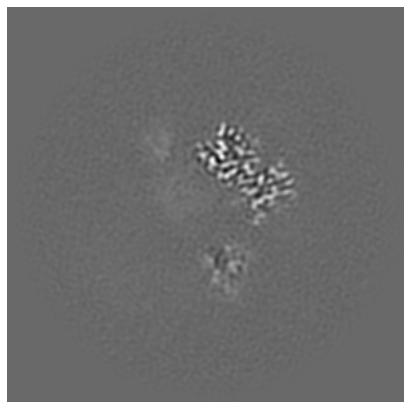


Z Index: 110

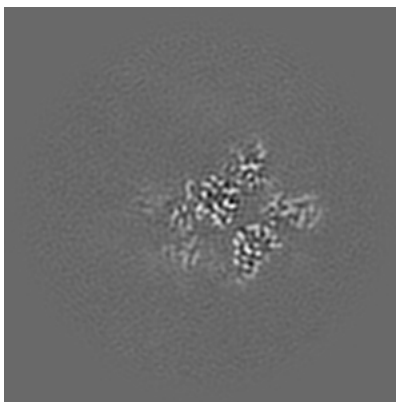
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

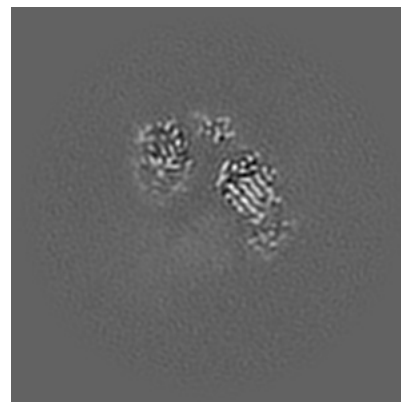
6.3.1 Primary map



X Index: 122

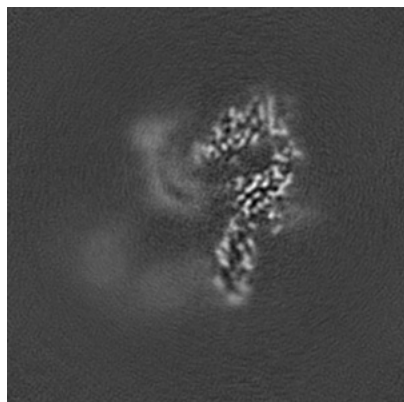


Y Index: 136

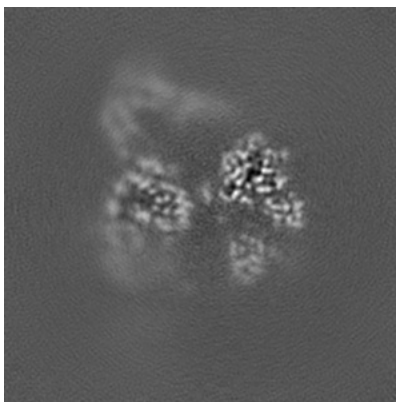


Z Index: 132

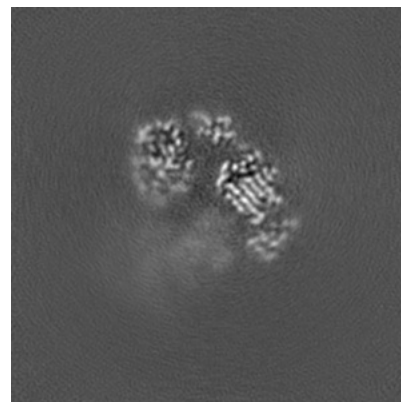
6.3.2 Raw map



X Index: 111



Y Index: 123

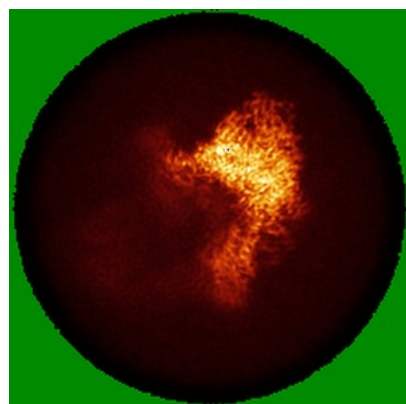


Z Index: 132

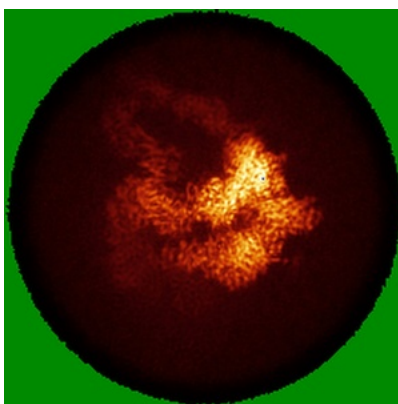
The images above show the largest variance slices of the map in three orthogonal directions.

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

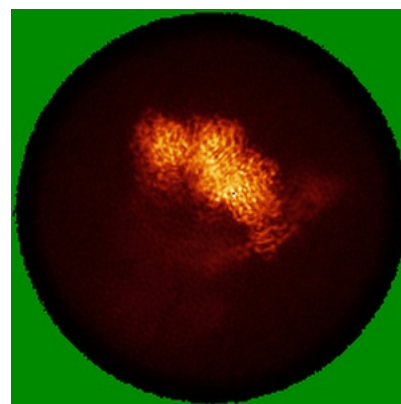
6.4.1 Primary map



X

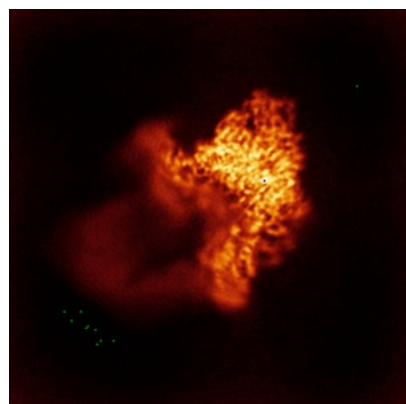


Y

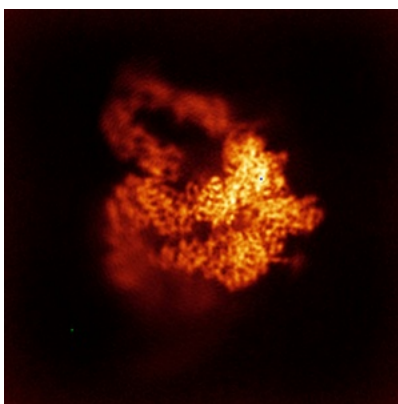


Z

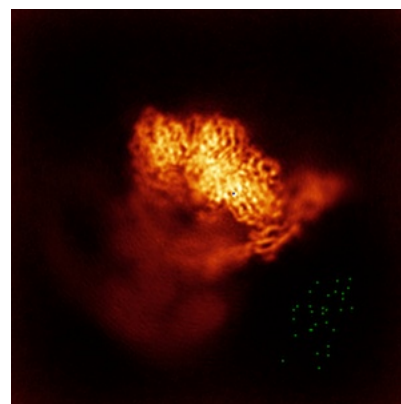
6.4.2 Raw map



X



Y

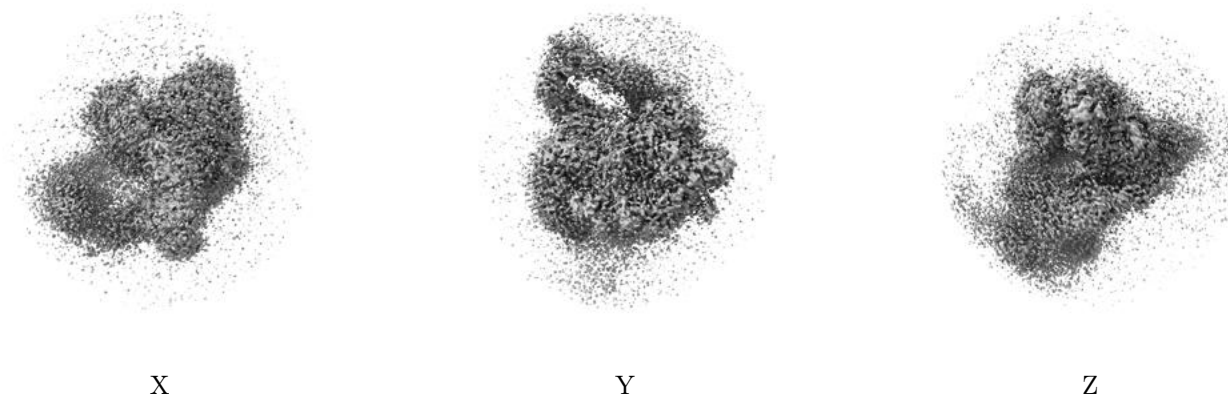


Z

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.

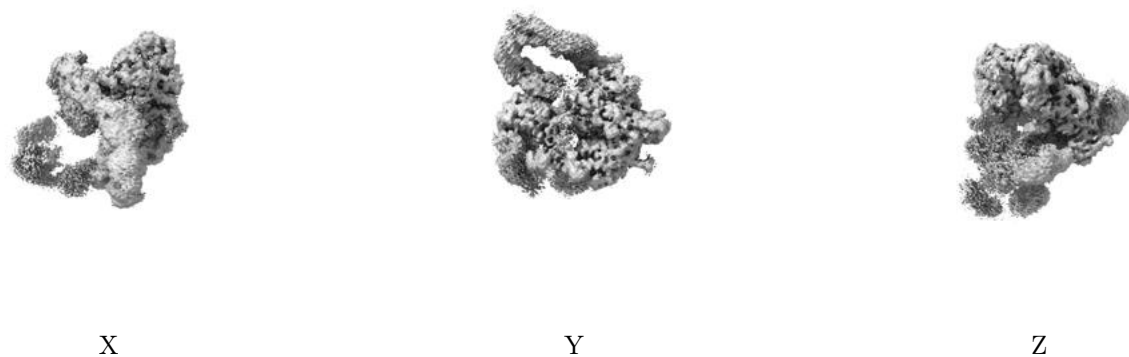
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.05. 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.

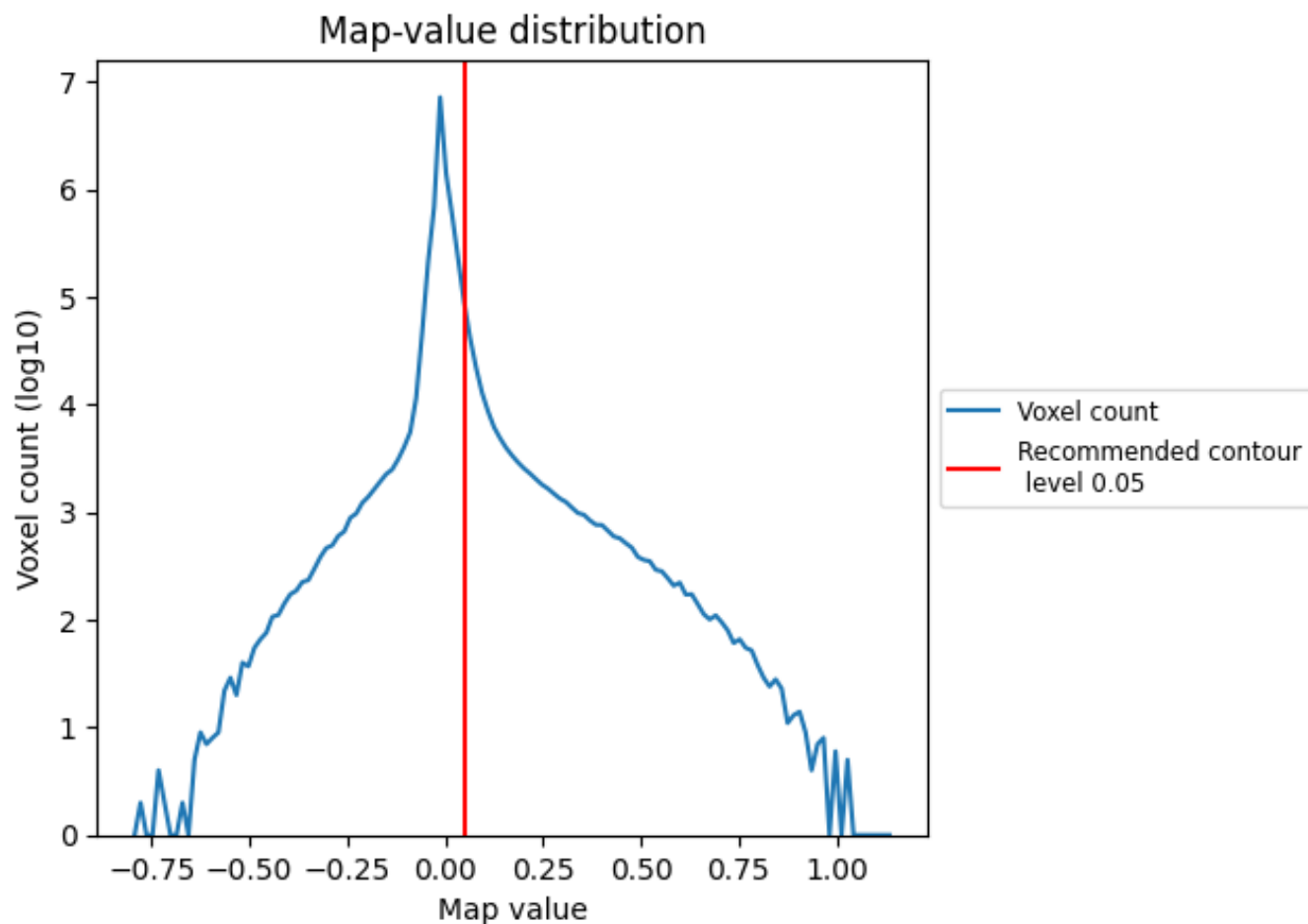
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

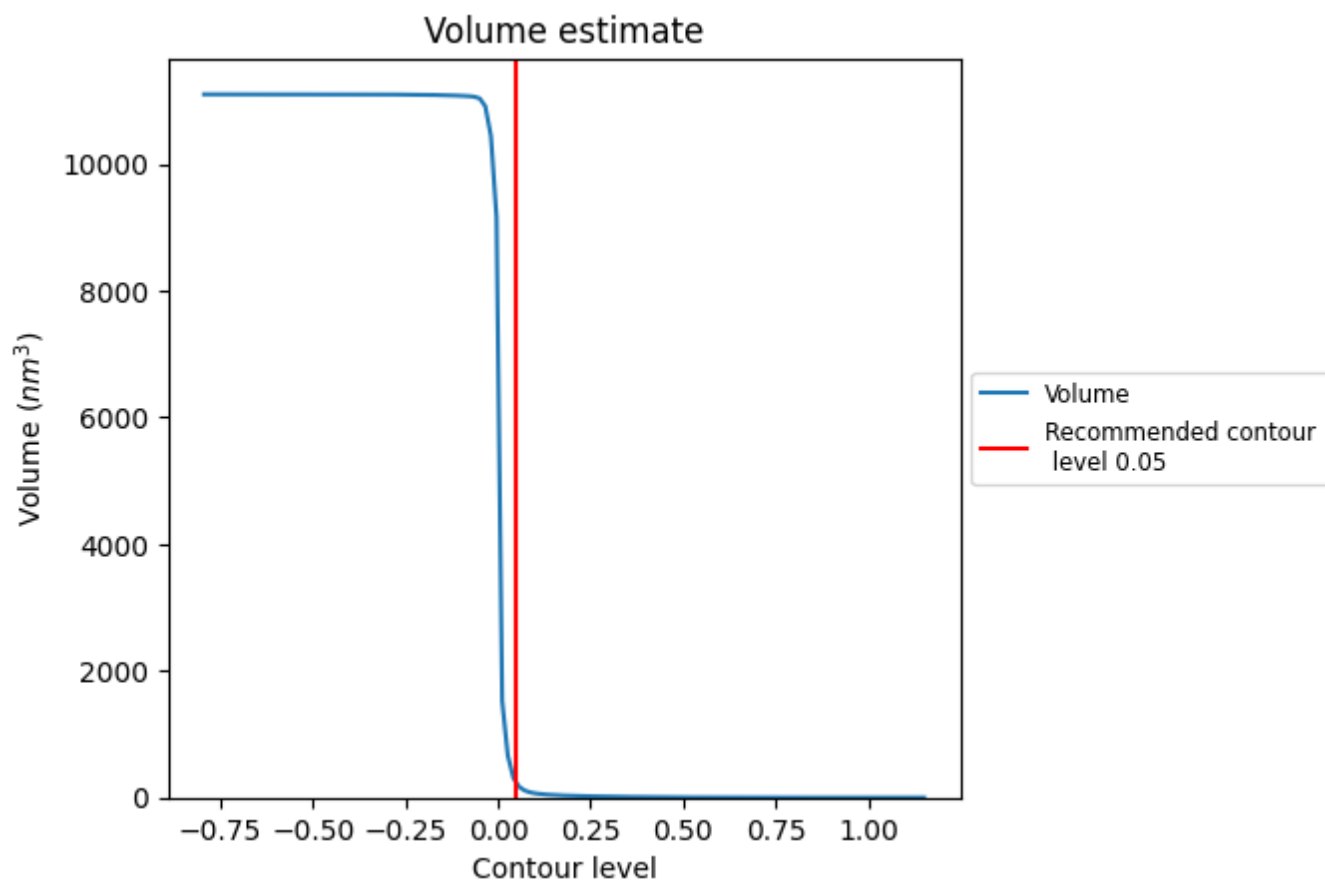
This section contains the results of statistical analysis of the map.

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.

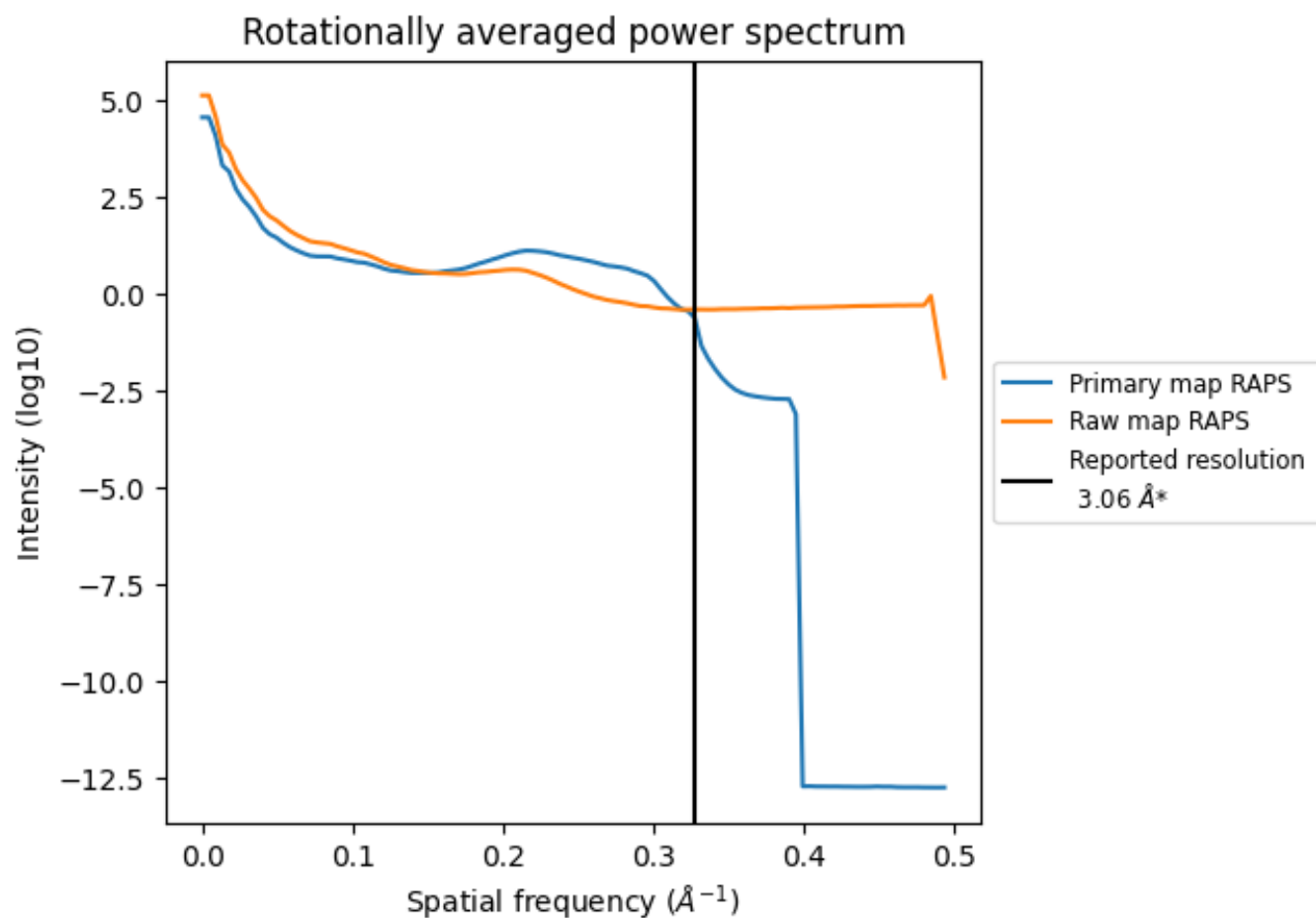
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 231 nm^3 ; this corresponds to an approximate mass of 209 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

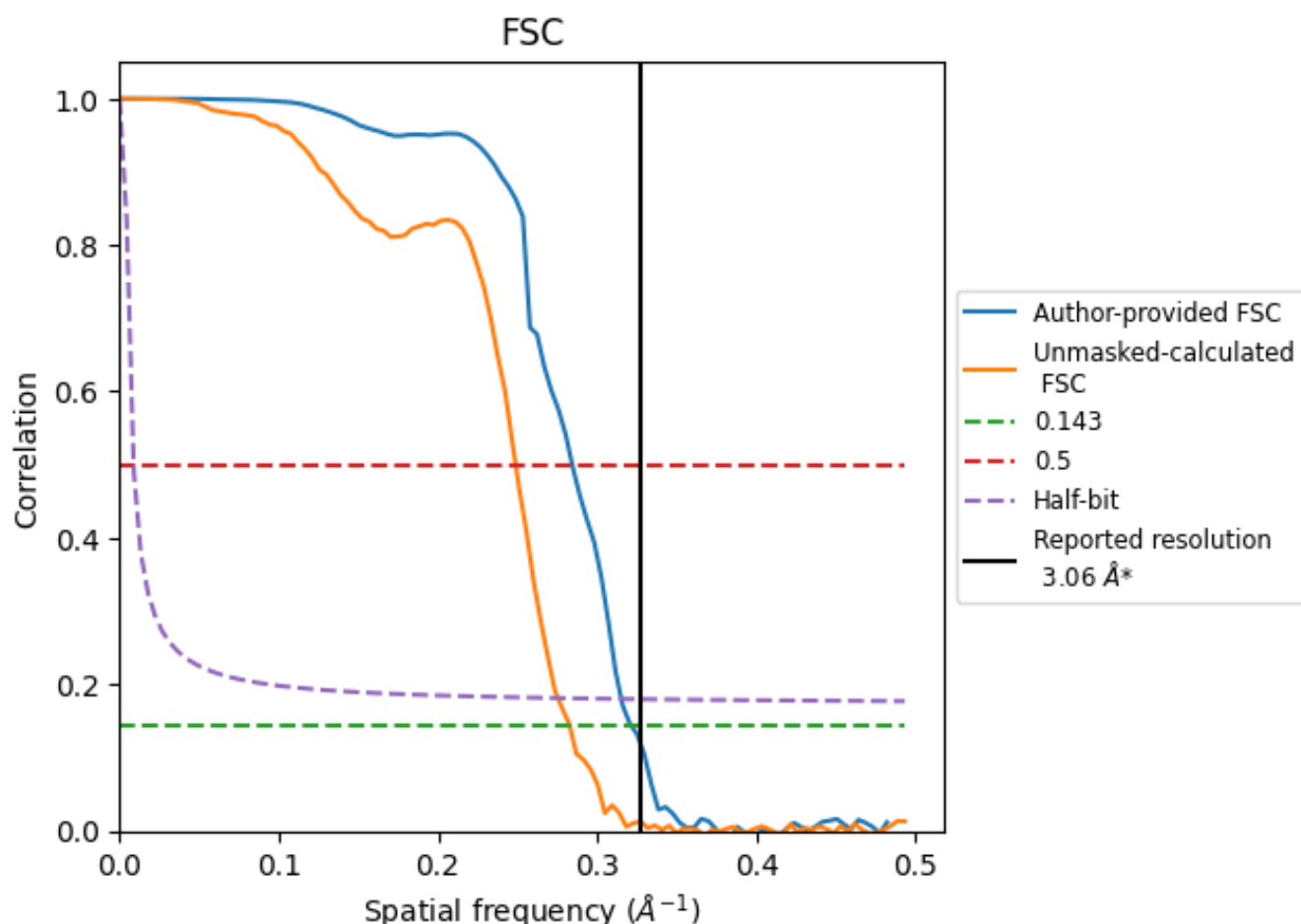


*Reported resolution corresponds to spatial frequency of 0.327 Å⁻¹

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 0.327 Å⁻¹

8.2 Resolution estimates [i](#)

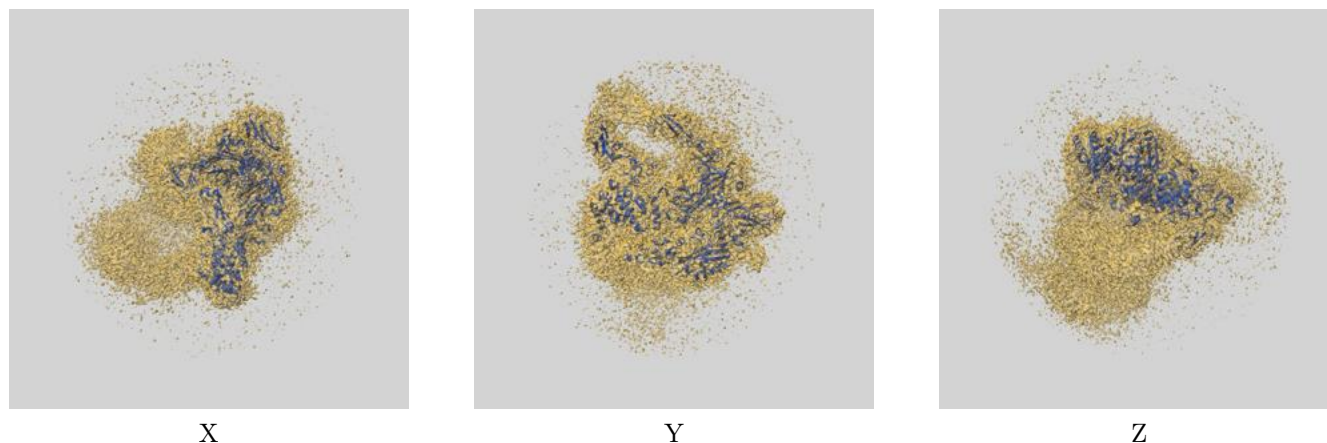
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.06	-	-
Author-provided FSC curve	3.11	3.52	3.17
Unmasked-calculated*	3.54	4.02	3.63

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.54 differs from the reported value 3.06 by more than 10 %

9 Map-model fit [i](#)

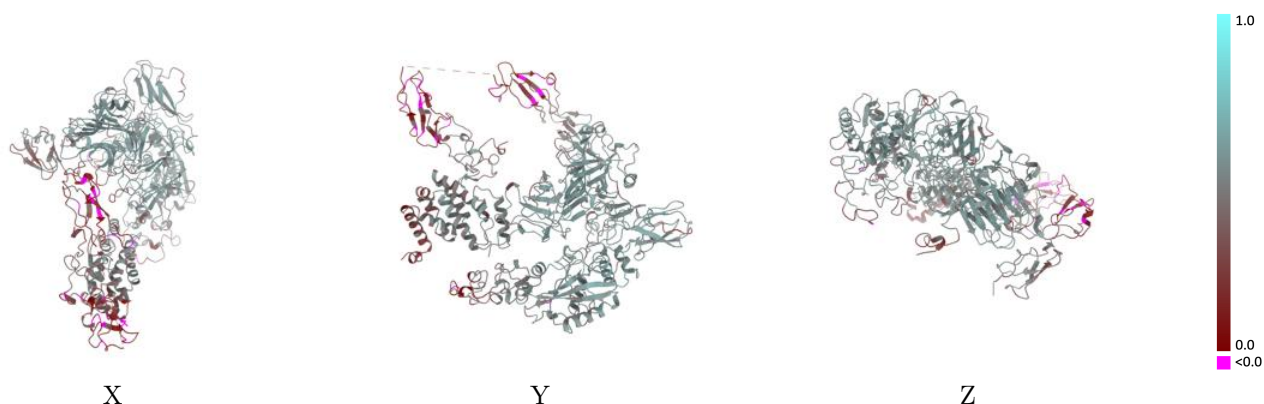
This section contains information regarding the fit between EMDB map EMD-15220 and PDB model 8A7D. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



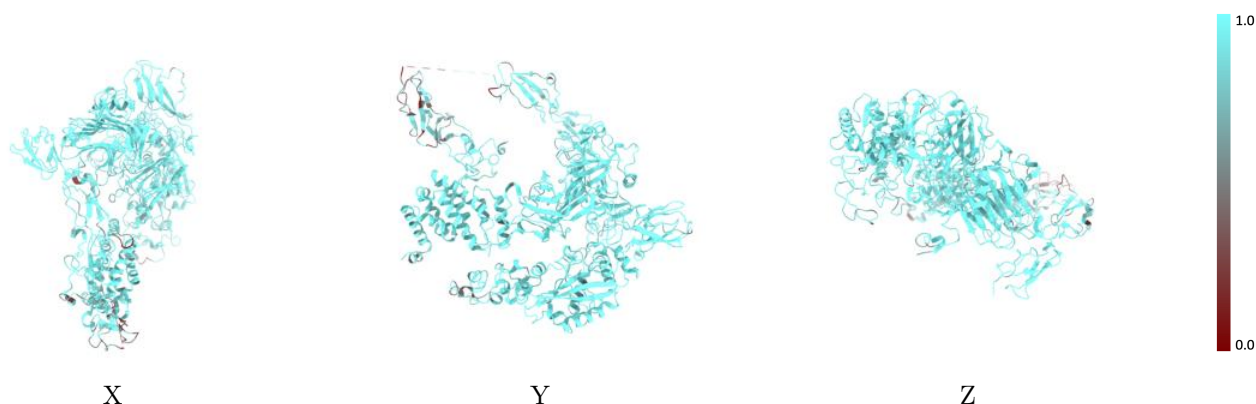
The images above show the 3D surface view of the map at the recommended contour level 0.05 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



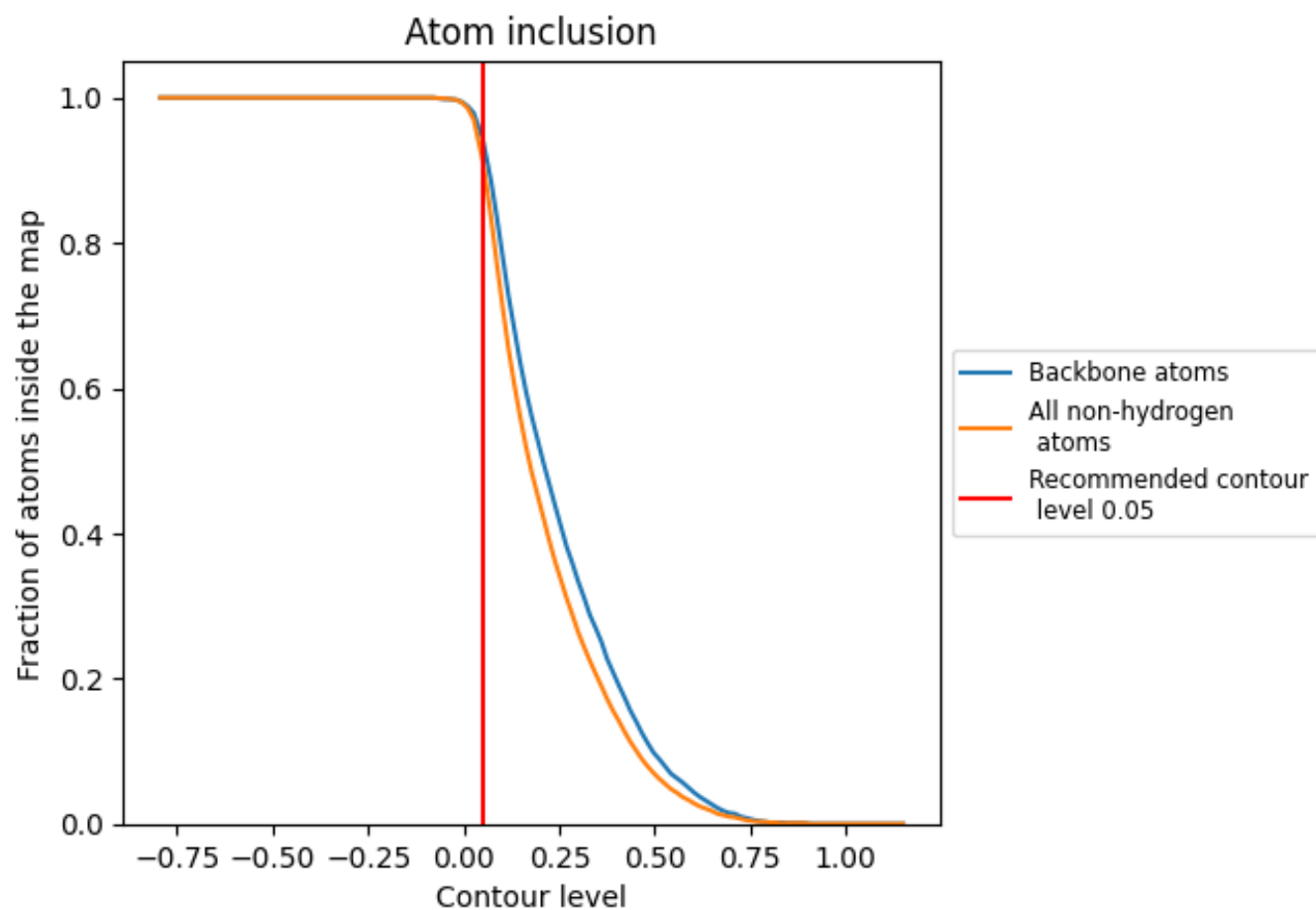
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.05).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 91% 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 (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.9130	<div></div> 0.4570
C	<div></div> 0.9350	<div></div> 0.5070
P	<div></div> 0.9290	<div></div> 0.4320
Q	<div></div> 0.8260	<div></div> 0.2980

