



Full wwPDB EM Validation Report ⓘ

Mar 28, 2026 – 03:09 AM UTC

PDB ID : 7Q1U / pdb_00007q1u
EMDB ID : EMD-13764
Title : Structure of Hedgehog acyltransferase (HHAT) in complex with megabody 177 bound to non-hydrolysable palmitoyl-CoA (Composite Map)
Authors : Coupland, C.; Carrique, L.; Siebold, C.
Deposited on : 2021-10-21
Resolution : 2.70 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

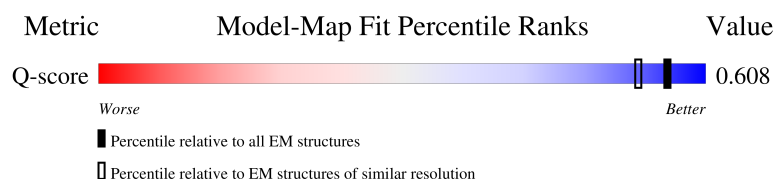
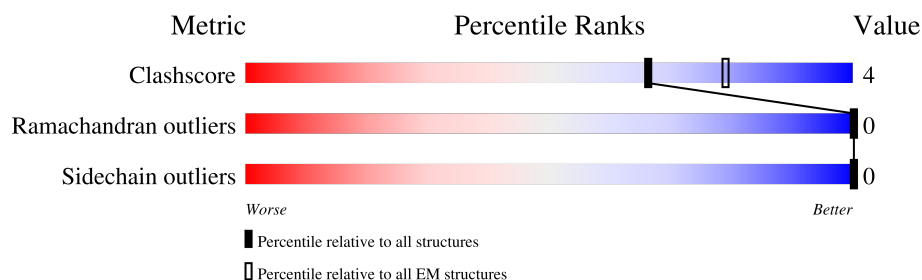
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.70 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10327 (2.20 - 3.20)

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

Mol	Chain	Length	Quality of chain
1	A	503	
2	B	919	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	HD6	A	603	X	-	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 21944 atoms, of which 10829 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 Protein-cysteine N-palmitoyltransferase HHAT.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
1	A	485	7979	2670	3972	638	668	31	0	0

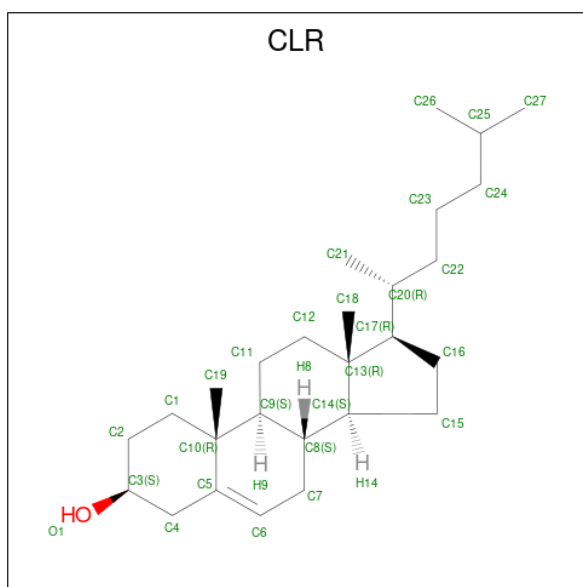
There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	182	ASN	SER	variant	UNP Q5VTY9
A	494	GLY	-	expression tag	UNP Q5VTY9
A	495	THR	-	expression tag	UNP Q5VTY9
A	496	GLU	-	expression tag	UNP Q5VTY9
A	497	THR	-	expression tag	UNP Q5VTY9
A	498	SER	-	expression tag	UNP Q5VTY9
A	499	GLN	-	expression tag	UNP Q5VTY9
A	500	VAL	-	expression tag	UNP Q5VTY9
A	501	ALA	-	expression tag	UNP Q5VTY9
A	502	PRO	-	expression tag	UNP Q5VTY9
A	503	ALA	-	expression tag	UNP Q5VTY9

- Molecule 2 is a protein called Megabody 177.

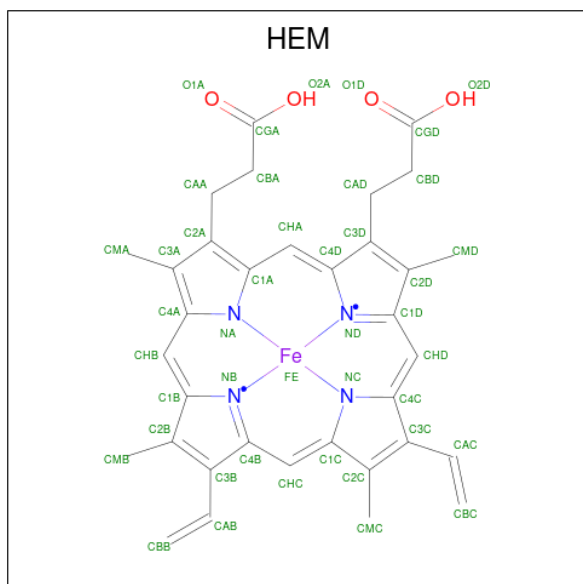
Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
2	B	872	13621	4398	6685	1200	1317	21	0	0

- Molecule 3 is CHOLESTEROL (CCD ID: CLR) (formula: C₂₇H₄₆O).



Mol	Chain	Residues	Atoms				AltConf
3	A	1	Total	C	H	O	0
			74	27	46	1	

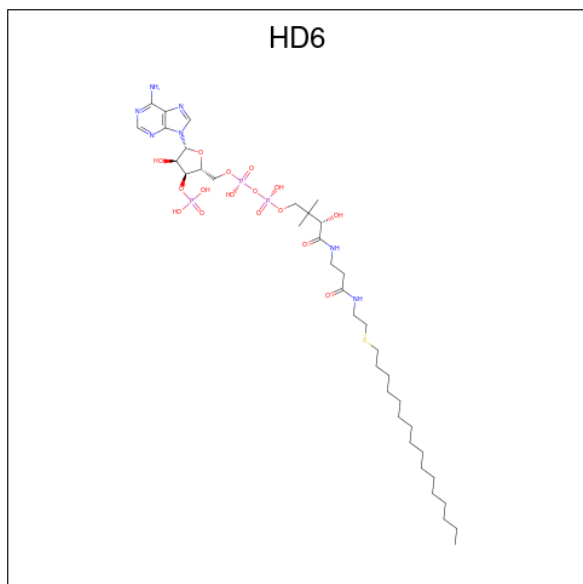
- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
4	A	1	Total	C	Fe	N	O	0
			43	34	1	4	4	

- Molecule 5 is $[[[(2 \{R\}, 3 \{S\}, 4 \{R\}, 5 \{R\})-5-(6\text{-aminopurin-9-yl})-4\text{-oxidanyl-3-phosphonooxyl-oxolan-2-yl}]methoxy-oxidanyl-phosphoryl] [(3 \{S\})-4-[[3-(2\text{-hexadecylsulfanylethylamino}$

)-3-oxidanylidene-propyl]amino]-2,2-dimethyl-3-oxidanyl-4-oxidanylidene-butyl] hydrogen phosphate (CCD ID: HD6) (formula: $C_{37}H_{68}N_7O_{16}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							AltConf
			Total	C	H	N	O	P	S	
5	A	1	128	37	64	7	16	3	1	0

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
6	A	1	Total	Mg	0
			1	1	

- Molecule 7 is PALMITIC ACID (CCD ID: PLM) (formula: $C_{16}H_{32}O_2$).



Mol	Chain	Residues	Atoms				AltConf
7	A	1	Total 49	C 16	H 31	O 2	0
7	A	1	Total 49	C 16	H 31	O 2	0

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	238000	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 ($e^-/\text{\AA}^2$)	54.8	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	61.287	Depositor
Minimum map value	-31.064	Depositor
Average map value	0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	10.8	Depositor
Map size ($\text{\AA}$)	331.59, 331.59, 331.59	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.1053, 1.1053, 1.1053	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CLR, PLM, HEM, HD6, MG, P1L

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	0.15	0/4119	0.28	0/5615
2	B	0.14	0/7119	0.26	0/9674
All	All	0.15	0/11238	0.26	0/15289

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	4007	3972	3959	31	0
2	B	6936	6685	6670	53	0
3	A	28	46	46	1	0
4	A	43	0	30	2	0
5	A	64	64	0	0	0
6	A	1	0	0	0	0
7	A	36	62	62	0	0
All	All	11115	10829	10767	86	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 4.

All (86) 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:B:409:THR:HG21	2:B:536:LEU:HD12	1.64	0.79
1:A:176:ARG:NH2	1:A:213:VAL:O	2.24	0.71
2:B:580:LEU:HD13	2:B:586:VAL:HA	1.76	0.65
2:B:355:LYS:O	2:B:630:ARG:NH2	2.30	0.64
1:A:317:THR:O	1:A:317:THR:HG23	1.99	0.63
2:B:643:TRP:CZ2	2:B:688:ILE:HD12	2.33	0.62
1:A:207:TYR:OH	1:A:220:LEU:O	2.14	0.61
1:A:213:VAL:CG1	1:A:311:MET:HE2	2.31	0.61
1:A:142:LEU:O	1:A:146:THR:HG23	2.00	0.61
1:A:34:GLU:OE1	1:A:34:GLU:N	2.34	0.61
2:B:409:THR:HG21	2:B:536:LEU:CD1	2.30	0.60
2:B:682:ALA:O	2:B:754:VAL:HG11	2.01	0.59
2:B:512:TYR:HB2	2:B:530:VAL:HG21	1.84	0.58
2:B:609:GLN:O	2:B:613:GLU:OE1	2.22	0.57
2:B:643:TRP:NE1	2:B:688:ILE:HG23	2.20	0.57
1:A:248:LEU:O	1:A:252:LEU:HD23	2.05	0.57
4:A:602:HEM:HBC2	4:A:602:HEM:HHD	1.87	0.56
2:B:390:HIS:CD2	2:B:460:LEU:HD21	2.41	0.55
2:B:752:TYR:O	2:B:756:GLN:N	2.33	0.55
1:A:280:VAL:O	1:A:459:ARG:NH1	2.38	0.55
2:B:448:LYS:NZ	2:B:457:GLU:OE2	2.39	0.55
2:B:261:PHE:CE2	2:B:265:LEU:HD11	2.42	0.54
2:B:390:HIS:CG	2:B:460:LEU:HD21	2.41	0.54
1:A:11:LEU:HA	1:A:482:VAL:HG21	1.89	0.54
2:B:890:HIS:NE2	2:B:902:ASN:OD1	2.35	0.54
2:B:670:TRP:CZ2	2:B:718:LEU:HD12	2.43	0.53
1:A:21:SER:O	1:A:25:VAL:HG23	2.10	0.52
1:A:75:LEU:HD22	1:A:310:LEU:CD1	2.40	0.52
2:B:25:VAL:HG22	2:B:821:ALA:HB2	1.91	0.52
2:B:550:LEU:HD12	2:B:573:LEU:HD22	1.92	0.51
1:A:460:ILE:O	1:A:469:THR:OG1	2.25	0.51
4:A:602:HEM:HHD	4:A:602:HEM:CBC	2.42	0.50
2:B:823:PHE:O	2:B:865:ARG:NH2	2.45	0.50
2:B:367:MET:HB2	2:B:505:ILE:HD11	1.93	0.50
2:B:212:ALA:O	2:B:271:TYR:OH	2.24	0.50
1:A:7:LEU:HD11	1:A:485:ALA:HB1	1.94	0.48
1:A:271:TYR:HB3	2:B:897:MET:HE1	1.95	0.48
1:A:79:MET:HE2	1:A:313:LEU:CD1	2.44	0.48
2:B:791:ALA:HB1	2:B:797:HIS:CE1	2.48	0.48
1:A:157:ARG:NH2	1:A:163:GLU:OE2	2.42	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:249:ASP:OD1	2:B:302:ASN:ND2	2.45	0.47
2:B:315:SER:OG	2:B:642:ASN:ND2	2.35	0.47
1:A:7:LEU:HD11	1:A:485:ALA:CB	2.45	0.47
2:B:683:HIS:NE2	2:B:750:GLU:OE2	2.48	0.46
2:B:259:ASP:HB3	2:B:316:TRP:HE1	1.81	0.46
1:A:144:LEU:HD21	1:A:378:TRP:HA	1.97	0.46
1:A:14:SER:HA	1:A:17:PHE:CE1	2.51	0.46
1:A:307:PRO:O	1:A:311:MET:HG3	2.14	0.46
1:A:335:TRP:NE1	1:A:392:ASN:OD1	2.49	0.46
2:B:519:THR:HG22	2:B:520:ARG:N	2.31	0.46
2:B:464:THR:CG2	2:B:596:ILE:HD11	2.47	0.45
2:B:172:THR:HG21	2:B:176:TYR:OH	2.17	0.45
1:A:219:ILE:HG23	1:A:219:ILE:O	2.17	0.45
2:B:674:THR:HG21	2:B:696:VAL:HG21	1.99	0.44
1:A:75:LEU:HD13	1:A:306:VAL:HG11	2.00	0.44
2:B:59:GLU:OE2	2:B:64:VAL:CG2	2.66	0.44
2:B:331:LYS:HB3	2:B:625:THR:HG23	2.00	0.43
1:A:177:CYS:HA	1:A:180:TYR:CZ	2.54	0.43
2:B:139:MET:HG2	2:B:761:VAL:HG13	2.00	0.43
1:A:98:MET:HE3	1:A:102:MET:HE2	2.01	0.43
2:B:656:THR:HG21	2:B:666:GLY:O	2.18	0.43
2:B:408:LEU:HD22	2:B:665:SER:OG	2.19	0.43
2:B:325:TYR:CG	2:B:632:ALA:HB2	2.54	0.42
2:B:845:ILE:HG23	2:B:845:ILE:O	2.20	0.42
2:B:238:THR:HG23	2:B:255:ARG:O	2.19	0.42
2:B:718:LEU:HD22	2:B:733:ASN:O	2.20	0.42
1:A:317:THR:O	1:A:317:THR:CG2	2.66	0.42
3:A:601:CLR:H121	3:A:601:CLR:H212	2.02	0.41
2:B:149:ALA:HB1	2:B:153:ARG:NH1	2.34	0.41
1:A:162:THR:HG23	1:A:165:GLU:H	1.85	0.41
1:A:323:ARG:HB2	1:A:328:MET:HE3	2.01	0.41
2:B:38:GLU:OE1	2:B:38:GLU:N	2.47	0.41
2:B:409:THR:O	2:B:411:GLU:N	2.53	0.41
1:A:108:VAL:HG23	1:A:109:LEU:HG	2.02	0.41
2:B:409:THR:O	2:B:410:GLU:C	2.63	0.41
2:B:149:ALA:HB1	2:B:153:ARG:HH12	1.86	0.41
2:B:523:LEU:HD13	2:B:554:THR:HG21	2.03	0.41
2:B:643:TRP:HE1	2:B:653:ASN:HB3	1.86	0.41
2:B:740:LYS:HB3	2:B:741:PRO:CD	2.52	0.41
2:B:877:ASN:O	2:B:878:SER:OG	2.39	0.40
2:B:701:ILE:HB	2:B:711:ASP:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:PHE:CE2	1:A:290:LEU:HD11	2.57	0.40
2:B:169:ASP:O	2:B:173:GLN:N	2.47	0.40
1:A:98:MET:HE3	1:A:102:MET:CE	2.51	0.40
2:B:65:ALA:HB3	2:B:791:ALA:HB2	2.02	0.40
2:B:772:HIS:O	2:B:776:LEU:HD13	2.22	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	480/503 (95%)	466 (97%)	14 (3%)	0	100	100
2	B	868/919 (94%)	846 (98%)	22 (2%)	0	100	100
All	All	1348/1422 (95%)	1312 (97%)	36 (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	422/435 (97%)	422 (100%)	0	100	100
2	B	726/752 (96%)	726 (100%)	0	100	100
All	All	1148/1187 (97%)	1148 (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 (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	GLN
1	A	428	HIS
1	A	443	ASN
2	B	44	ASN
2	B	92	ASN
2	B	125	GLN
2	B	157	GLN
2	B	158	GLN
2	B	173	GLN
2	B	313	ASN
2	B	320	HIS
2	B	379	ASN
2	B	477	ASN
2	B	677	GLN
2	B	692	ASN
2	B	781	HIS
2	B	875	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	P1L	A	387	1	21,22,23	0.73	0	19,23,25	1.41	3 (15%)

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
1	P1L	A	387	1	-	3/20/22/24	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	387	P1L	C8-C7-SG	-4.00	108.64	113.40
1	A	387	P1L	CB-SG-C7	3.08	104.99	100.76
1	A	387	P1L	O7-C7-SG	2.63	126.03	122.68

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	387	P1L	CA-CB-SG-C7
1	A	387	P1L	C9-C10-C11-C12
1	A	387	P1L	C13-C14-C15-C16

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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
7	PLM	A	606	-	17,17,17	0.96	1 (5%)	17,17,17	0.73	2 (11%)
4	HEM	A	602	1	50,50,50	1.44	6 (12%)	67,82,82	1.64	11 (16%)
5	HD6	A	603	6	64,66,66	2.69	19 (29%)	86,91,91	2.34	19 (22%)
3	CLR	A	601	-	31,31,31	0.11	0	48,48,48	0.21	0
7	PLM	A	605	-	17,17,17	0.92	1 (5%)	17,17,17	0.73	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
7	PLM	A	606	-	-	3/15/15/15	-
4	HEM	A	602	1	-	4/14/54/54	-
5	HD6	A	603	6	1/1/13/13	19/64/80/80	0/3/3/3
3	CLR	A	601	-	-	2/10/68/68	0/4/4/4
7	PLM	A	605	-	-	6/15/15/15	-

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	603	HD6	C5-N7	7.66	1.53	1.39
5	A	603	HD6	CBV-NBU	6.99	1.50	1.33
5	A	603	HD6	PCG-OCF	6.93	1.67	1.59
5	A	603	HD6	CBZ-NBY	6.71	1.49	1.33
5	A	603	HD6	C8-N9	-6.22	1.27	1.37
5	A	603	HD6	PCE-OCF	5.61	1.65	1.59
5	A	603	HD6	C4-N9	-5.55	1.26	1.37
5	A	603	HD6	C6-N6	4.52	1.45	1.34
4	A	602	HEM	FE-NB	4.45	2.08	1.94
5	A	603	HD6	OCK-CCL	3.63	1.50	1.42
5	A	603	HD6	CAP-CAQ	-3.63	1.45	1.53
5	A	603	HD6	PAM-OAN	3.46	1.65	1.59
4	A	602	HEM	FE-NC	3.41	2.06	1.95
4	A	602	HEM	C4D-ND	-3.38	1.34	1.40
7	A	605	PLM	C2-C1	3.08	1.57	1.50
7	A	606	PLM	C2-C1	3.05	1.57	1.50
5	A	603	HD6	C8-N7	2.91	1.37	1.31
5	A	603	HD6	OCK-CCJ	2.91	1.51	1.45
4	A	602	HEM	FE-NA	2.74	2.04	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	602	HEM	C4C-NC	-2.56	1.34	1.39
5	A	603	HD6	PAM-OAK	-2.37	1.46	1.54
5	A	603	HD6	OAZ-CBV	-2.30	1.18	1.23
5	A	603	HD6	PCE-OAU	-2.18	1.45	1.55
5	A	603	HD6	PAM-OAJ	-2.15	1.46	1.54
5	A	603	HD6	C4-N3	-2.14	1.30	1.34
4	A	602	HEM	C1C-NC	-2.10	1.35	1.39
5	A	603	HD6	OAX-CCA	-2.01	1.38	1.42

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	603	HD6	C4-N9-C8	14.19	120.64	105.74
5	A	603	HD6	C4-C5-N7	-5.64	104.14	110.58
5	A	603	HD6	N3-C4-N9	4.90	135.50	127.17
5	A	603	HD6	C5-C4-N3	-4.82	120.08	126.72
4	A	602	HEM	CHD-C4C-NC	4.69	129.56	124.45
5	A	603	HD6	N9-C8-N7	-4.62	107.39	113.94
4	A	602	HEM	CHB-C1B-NB	4.01	129.33	124.37
4	A	602	HEM	CHC-C4B-NB	3.66	128.36	124.42
5	A	603	HD6	N1-C2-N3	-3.64	123.08	128.58
4	A	602	HEM	C1B-NB-C4B	3.51	109.37	105.21
4	A	602	HEM	C4D-ND-C1D	3.40	109.23	105.21
5	A	603	HD6	CCL-N9-C8	-3.23	119.93	127.09
5	A	603	HD6	CCJ-OCK-CCL	-3.11	102.60	109.47
5	A	603	HD6	C4-N9-CCL	-3.08	119.44	126.63
4	A	602	HEM	CHA-C4D-ND	3.05	128.14	124.37
4	A	602	HEM	CHA-C1A-NA	2.76	128.86	123.86
5	A	603	HD6	C2-N3-C4	2.73	118.50	111.83
5	A	603	HD6	CAQ-CAP-CCL	2.49	105.36	99.89
5	A	603	HD6	OAS-PCG-OAR	-2.37	101.40	112.44
5	A	603	HD6	OAU-PCE-OAT	-2.37	101.40	112.44
4	A	602	HEM	CHC-C1C-NC	2.34	127.00	124.45
5	A	603	HD6	CAP-CAQ-CCJ	2.31	107.28	103.24
4	A	602	HEM	C2A-C1A-NA	-2.25	107.65	110.15
5	A	603	HD6	CBW-CBV-NBU	2.19	120.34	116.34
5	A	603	HD6	OAK-PAM-OAN	2.12	114.13	105.85
4	A	602	HEM	C4C-NC-C1C	2.12	109.28	105.82
7	A	606	PLM	O1-C1-O2	2.10	128.72	123.33
7	A	605	PLM	O1-C1-O2	2.10	128.72	123.33
5	A	603	HD6	OAJ-PAM-OAN	2.05	113.84	105.85
7	A	606	PLM	O2-C1-C2	-2.04	116.63	123.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	603	HD6	C6-C5-C4	2.03	119.95	117.18
4	A	602	HEM	CHD-C1D-ND	2.01	126.58	124.42
5	A	603	HD6	CBW-CBX-NBY	-2.00	107.74	112.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	603	HD6	CCA

All (34) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	603	HD6	CCI-OCH-PCG-OCF
5	A	603	HD6	CCI-OCH-PCG-OAR
5	A	603	HD6	OCH-CCI-CCJ-OCK
3	A	601	CLR	C17-C20-C22-C23
3	A	601	CLR	C21-C20-C22-C23
5	A	603	HD6	CBN-CBO-CBP-CBQ
5	A	603	HD6	CBK-CBL-CBM-CBN
7	A	606	PLM	CC-CD-CE-CF
7	A	605	PLM	C3-C4-C5-C6
5	A	603	HD6	OCH-CCI-CCJ-CAQ
5	A	603	HD6	CBM-CBN-CBO-CBP
5	A	603	HD6	CBF-CBG-CBH-CBI
7	A	605	PLM	C7-C8-C9-CA
5	A	603	HD6	SBR-CBS-CBT-NBU
5	A	603	HD6	PCE-OCF-PCG-OCH
7	A	605	PLM	C1-C2-C3-C4
5	A	603	HD6	CBC-CBD-CBE-CBF
5	A	603	HD6	CBP-CBQ-SBR-CBS
5	A	603	HD6	CCC-OCD-PCE-OAT
5	A	603	HD6	CCI-OCH-PCG-OAS
5	A	603	HD6	CBL-CBM-CBN-CBO
7	A	605	PLM	C8-C9-CA-CB
5	A	603	HD6	CBI-CBJ-CBK-CBL
4	A	602	HEM	CAA-CBA-CGA-O2A
4	A	602	HEM	CAA-CBA-CGA-O1A
4	A	602	HEM	CAD-CBD-CGD-O1D
4	A	602	HEM	CAD-CBD-CGD-O2D
7	A	605	PLM	CC-CD-CE-CF
5	A	603	HD6	PCG-OCF-PCE-OAU
7	A	606	PLM	O2-C1-C2-C3

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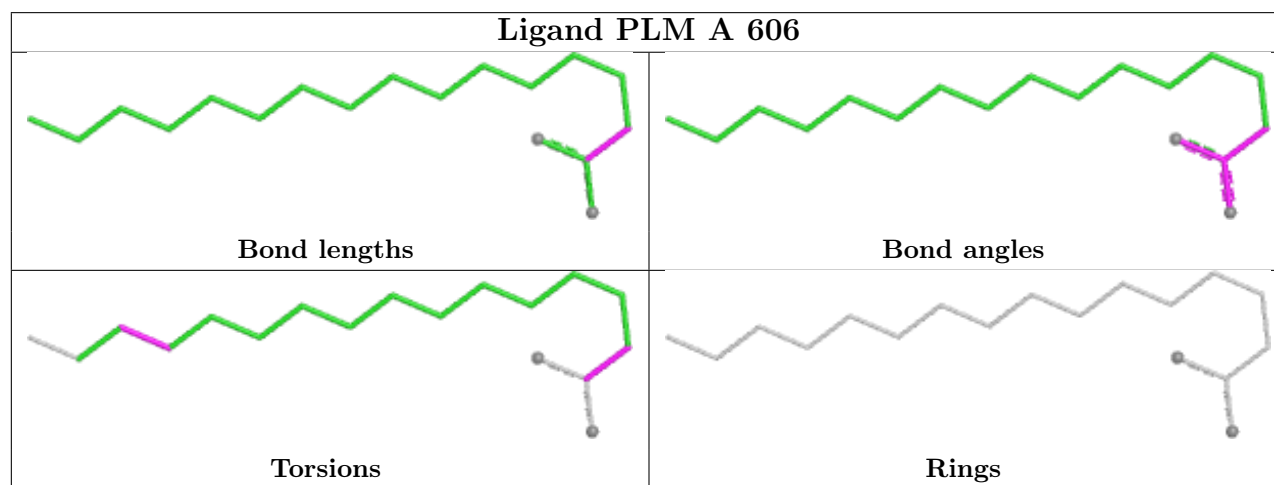
Mol	Chain	Res	Type	Atoms
7	A	606	PLM	O1-C1-C2-C3
7	A	605	PLM	C9-CA-CB-CC
5	A	603	HD6	NBY-CBZ-CCA-OAX
5	A	603	HD6	PCG-OCF-PCE-OAT

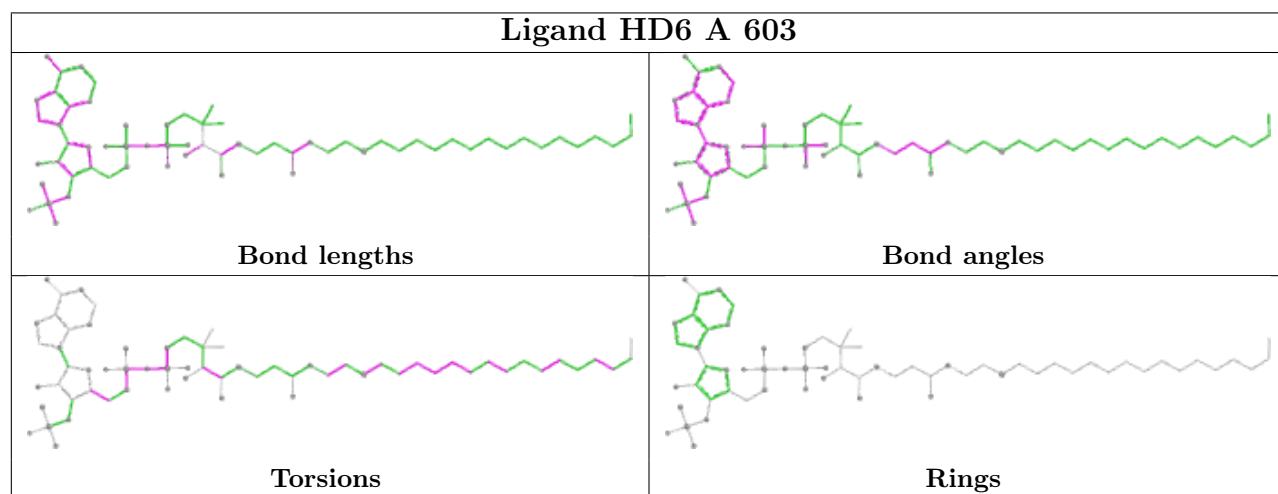
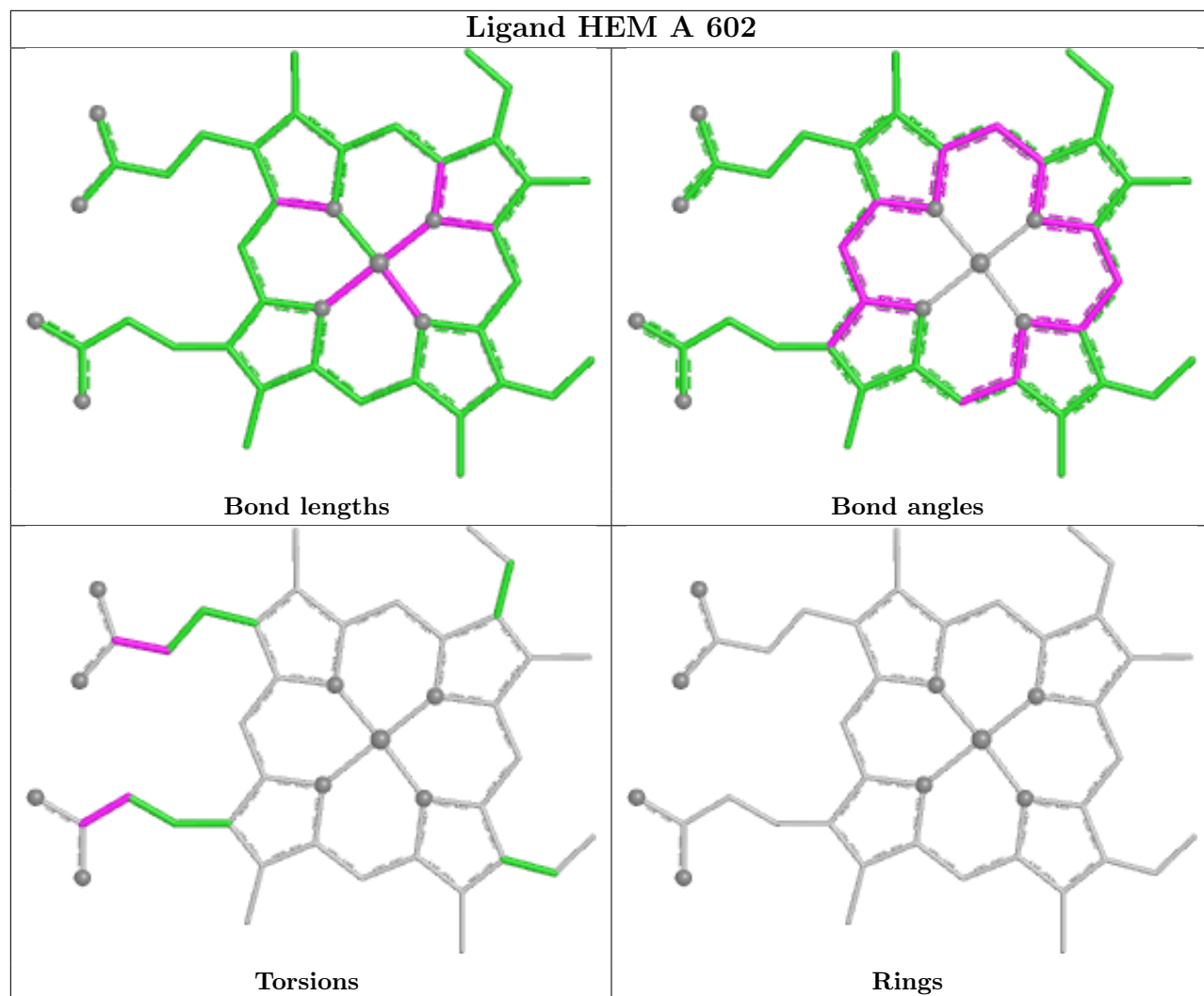
There are no ring outliers.

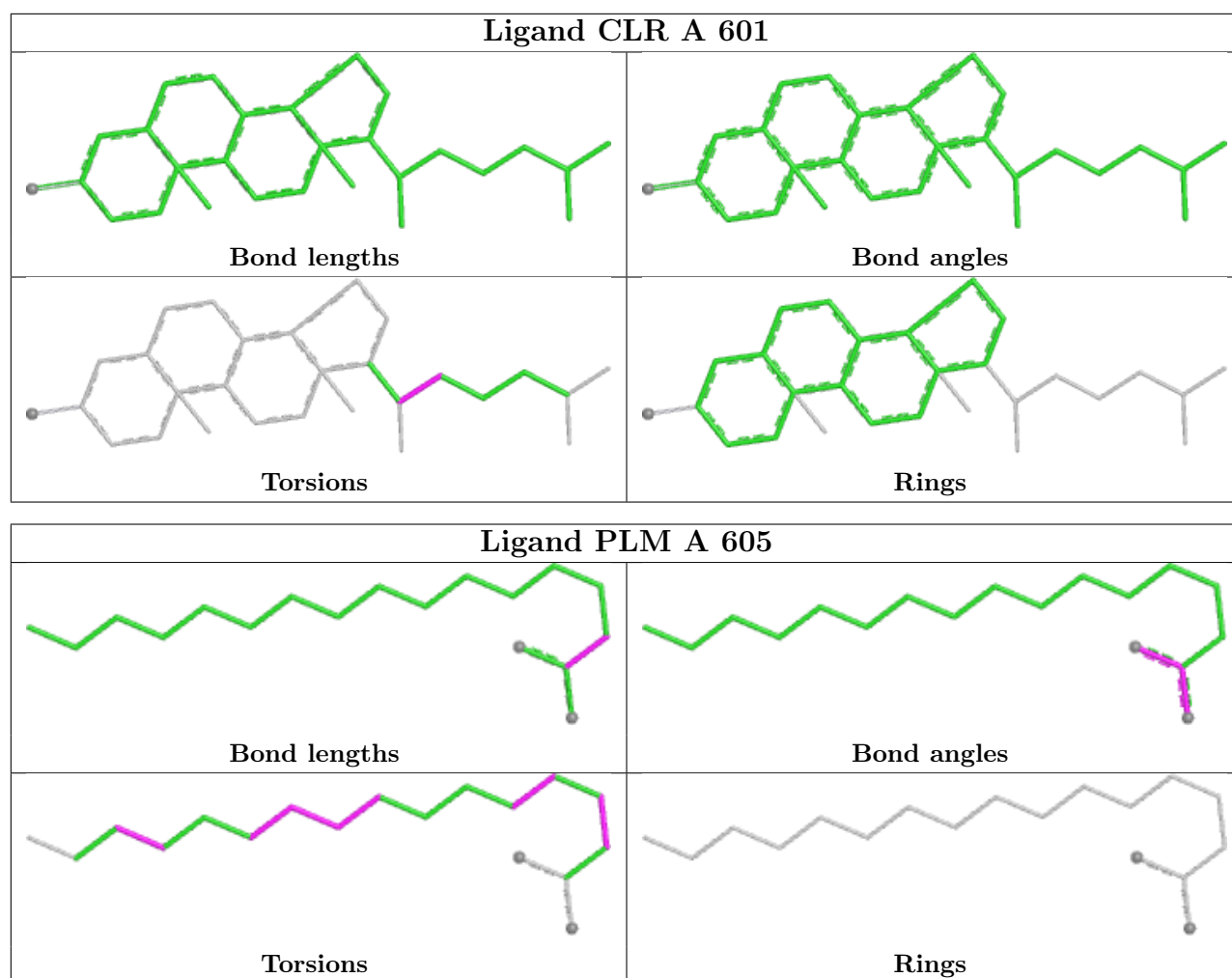
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	602	HEM	2	0
3	A	601	CLR	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.

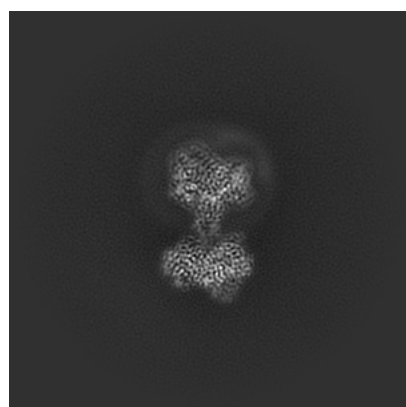
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13764. These allow visual inspection of the internal detail of the map and identification of artifacts.

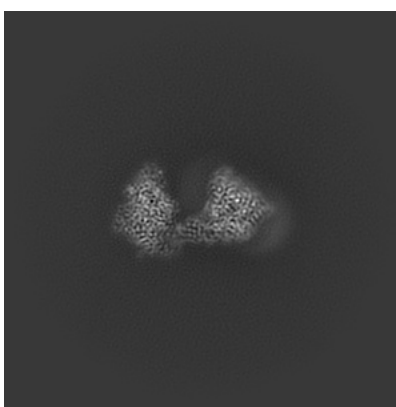
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

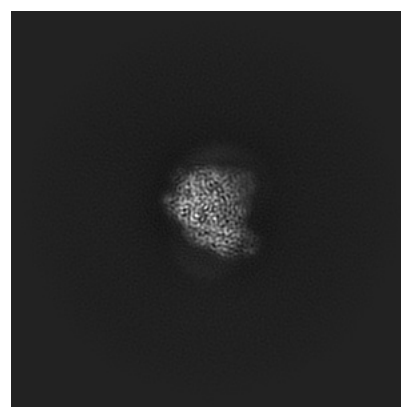
6.1.1 Primary map



X



Y

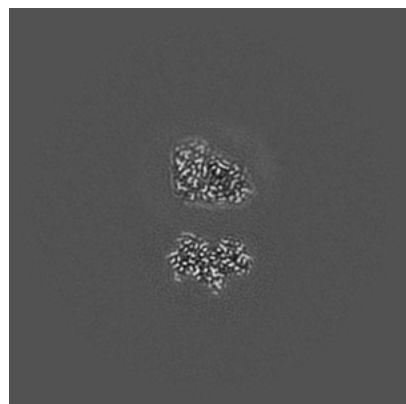


Z

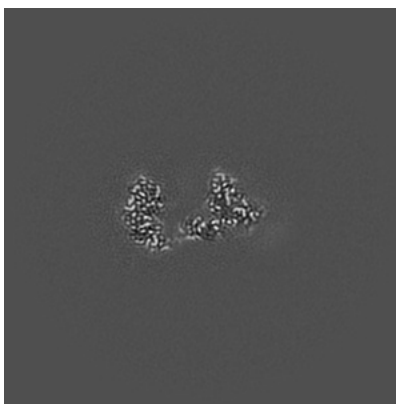
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

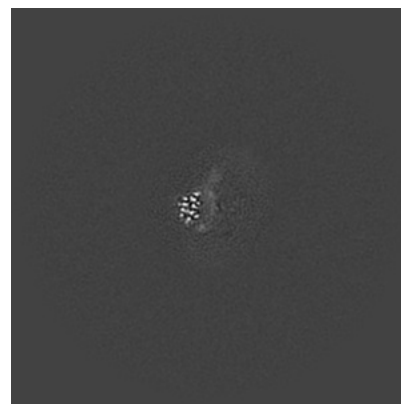
6.2.1 Primary map



X Index: 150



Y Index: 150

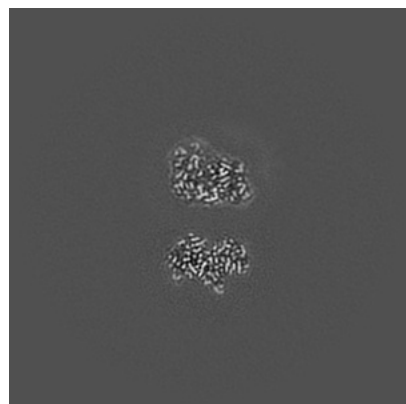


Z Index: 150

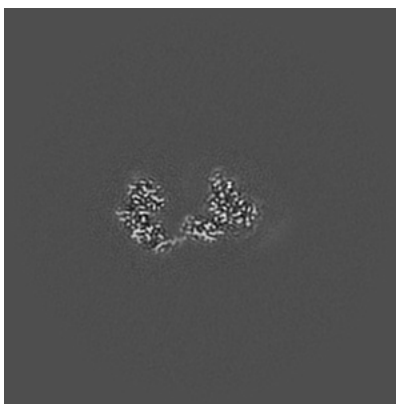
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

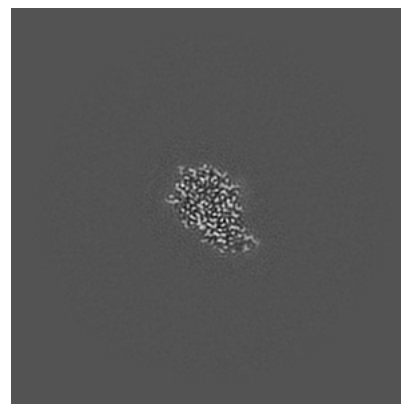
6.3.1 Primary map



X Index: 153



Y Index: 152

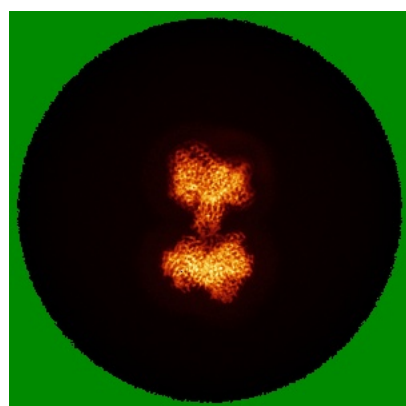


Z Index: 105

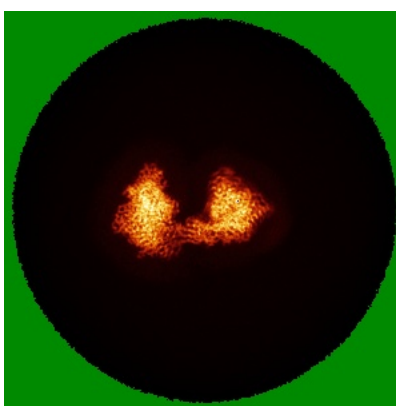
The images above show the largest variance slices of the map in three orthogonal directions.

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

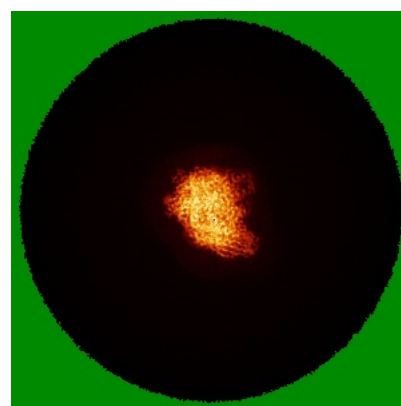
6.4.1 Primary map



X



Y

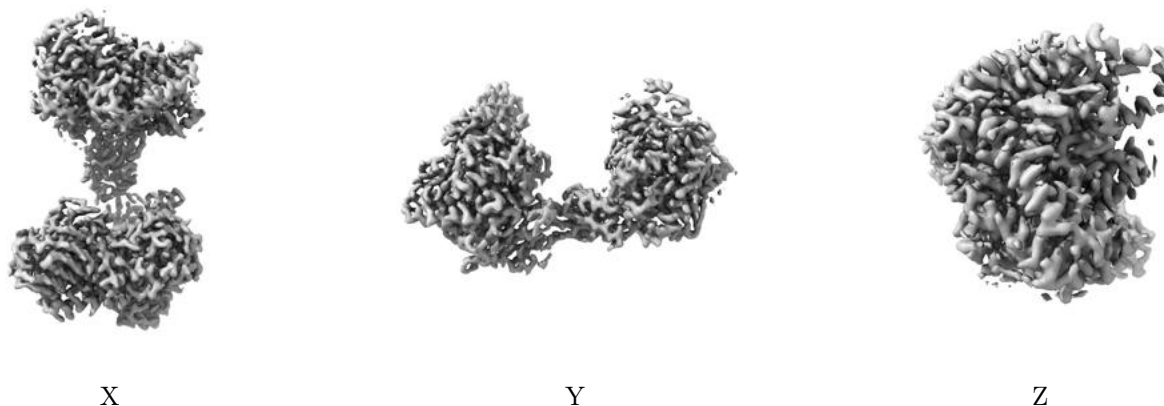


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 10.8. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

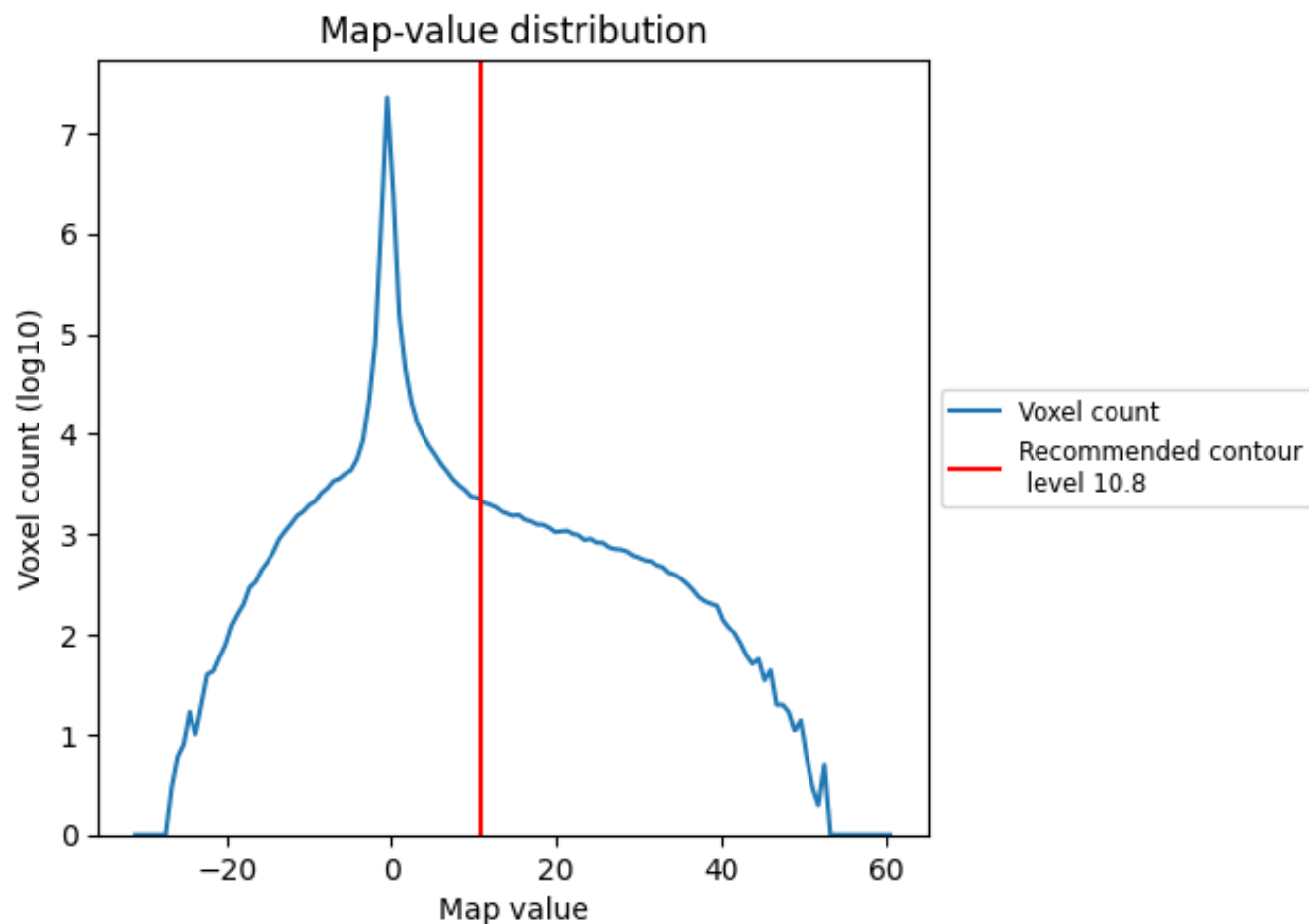
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

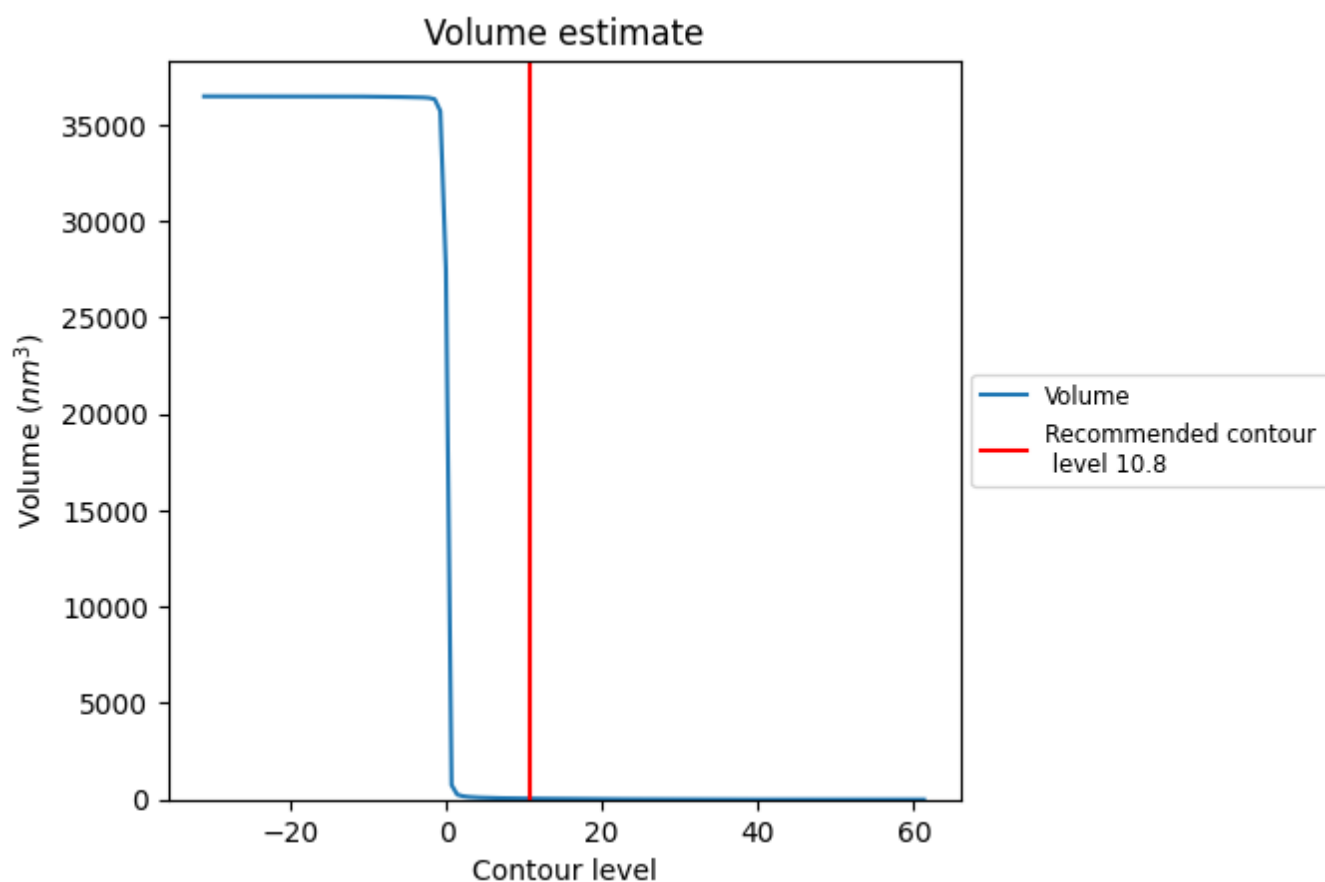
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

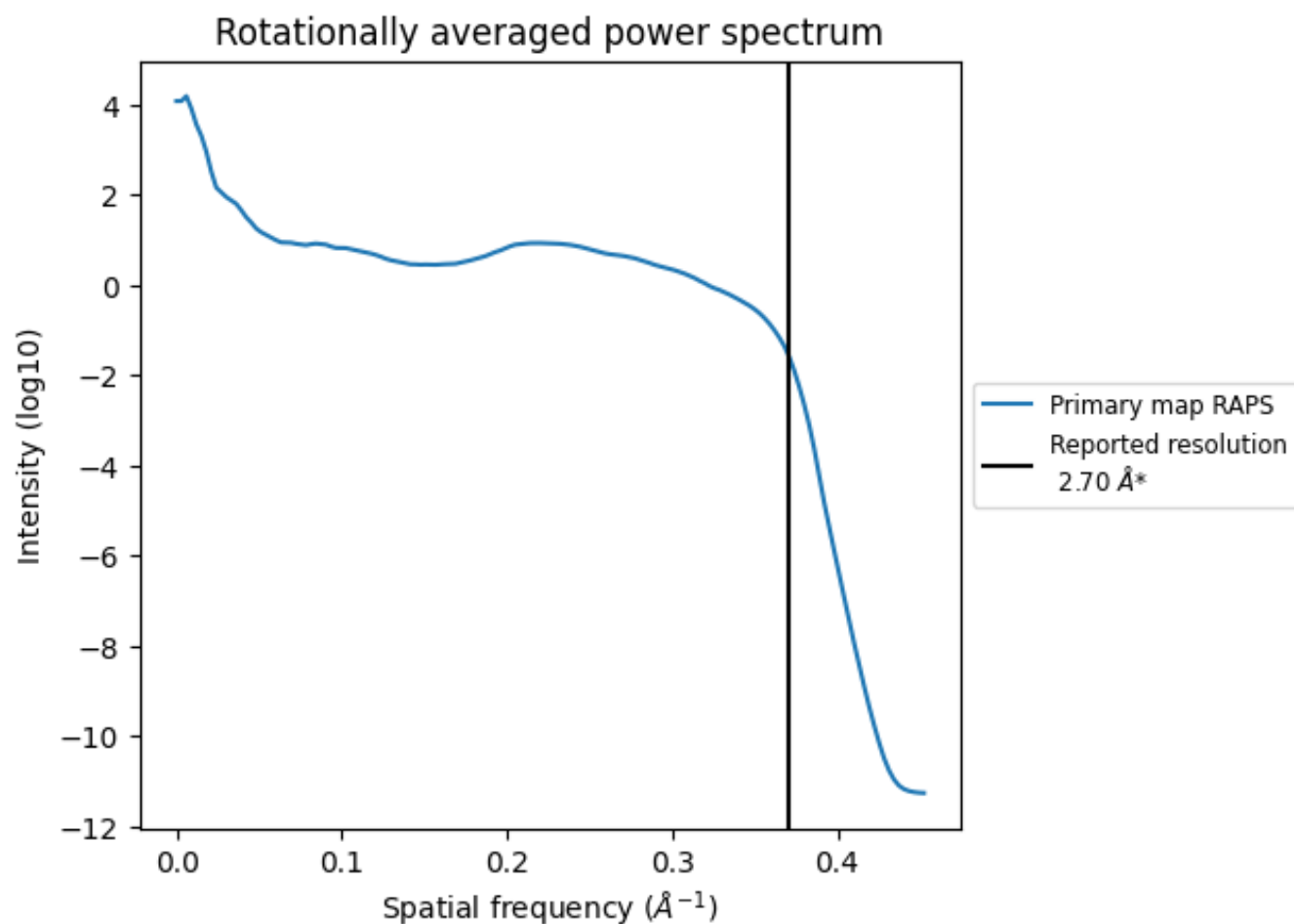
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 51 nm^3 ; this corresponds to an approximate mass of 46 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.370 Å⁻¹

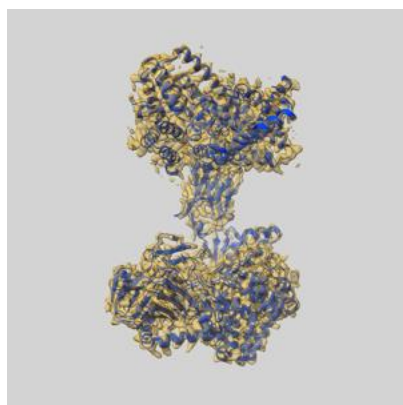
8 Fourier-Shell correlation

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

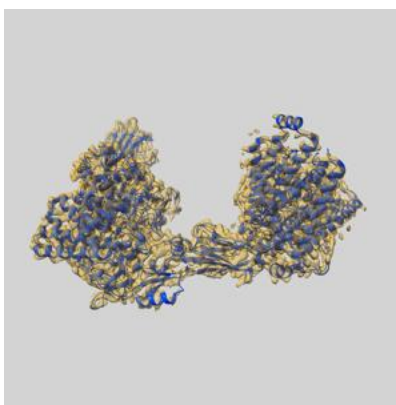
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-13764 and PDB model 7Q1U. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

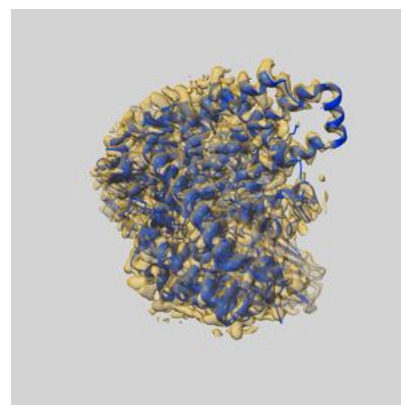
9.1 Map-model overlay [i](#)



X



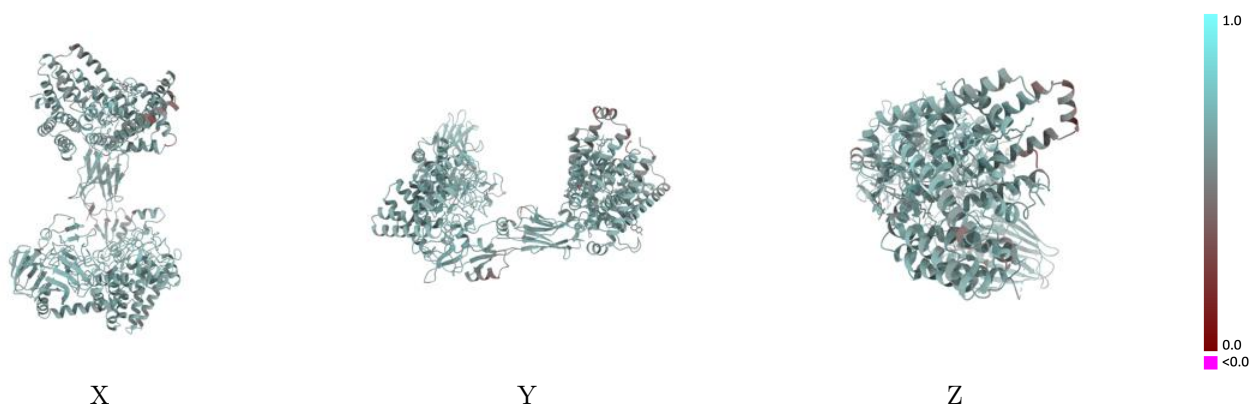
Y



Z

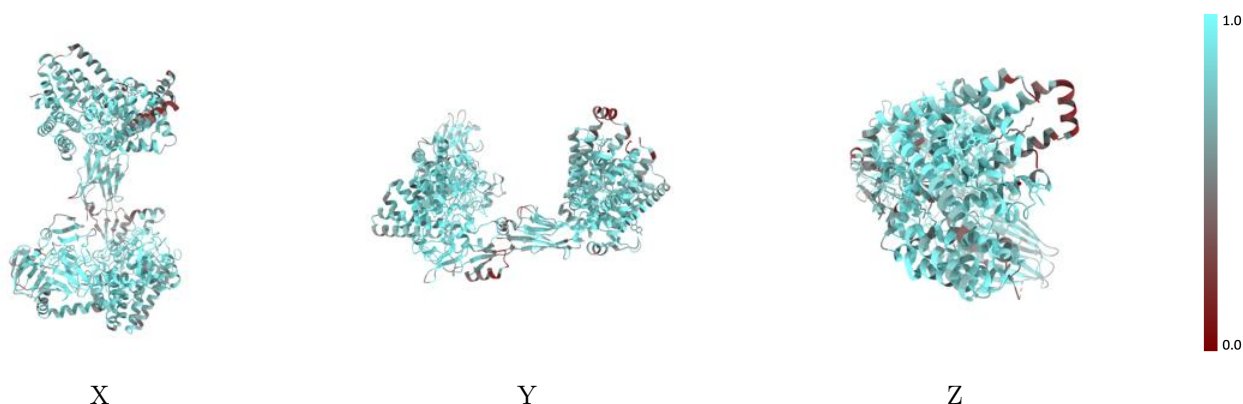
The images above show the 3D surface view of the map at the recommended contour level 10.8 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



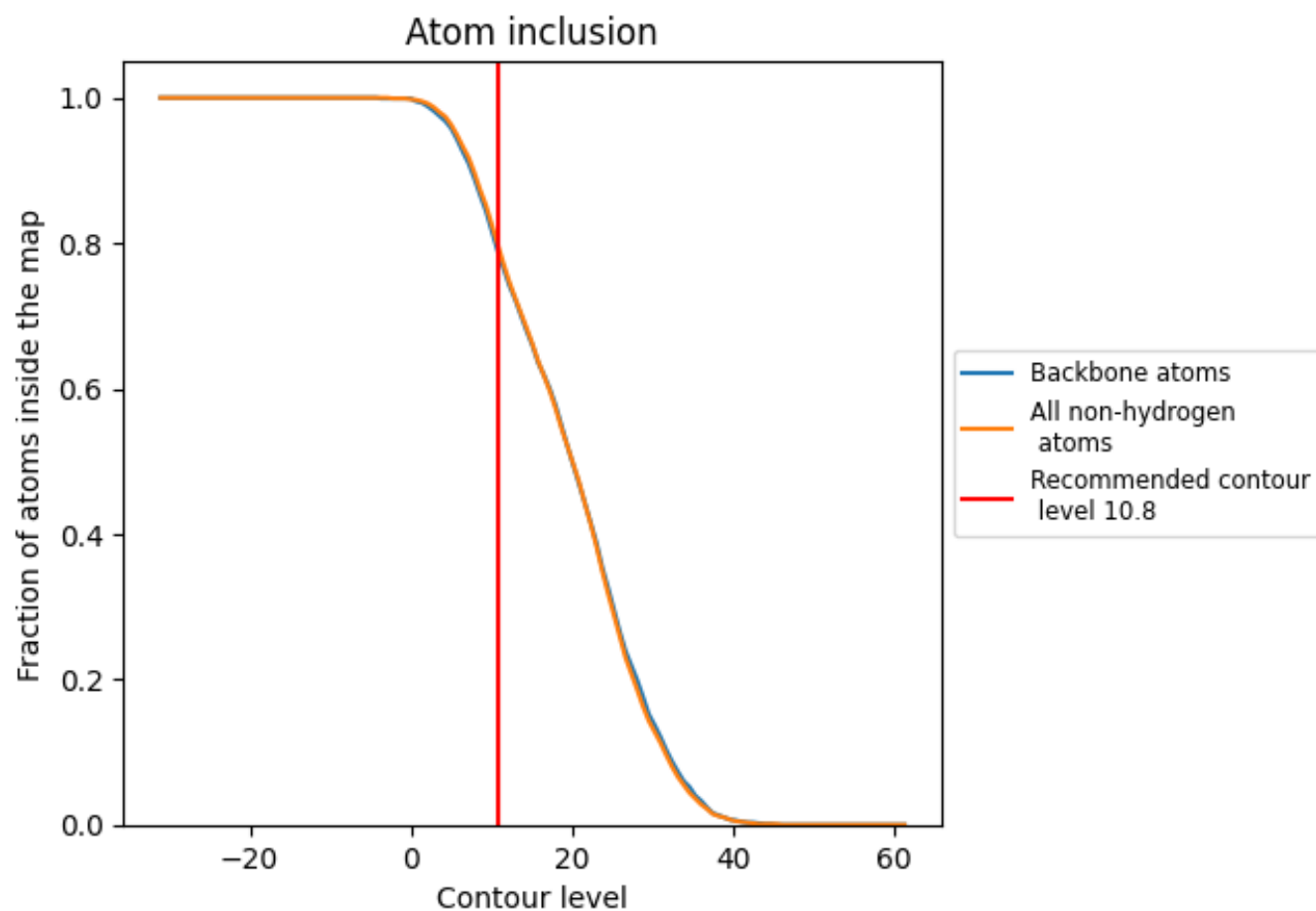
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (10.8).

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 80% 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 (10.8) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7970	0.6080
A	0.7880	0.5940
B	0.8100	0.6170

