



Full wwPDB EM Validation Report ⓘ

Mar 13, 2023 – 08:12 PM JST

PDB ID : 7YRQ
EMDB ID : EMD-34064
Title : Cryo-EM structure of human Peroxisomal ABC Transporter ABCD1
Authors : Chao, X.; Li-Na, J.; Lin, T.
Deposited on : 2022-08-10
Resolution : 3.35 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

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<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.dev43
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.32.1

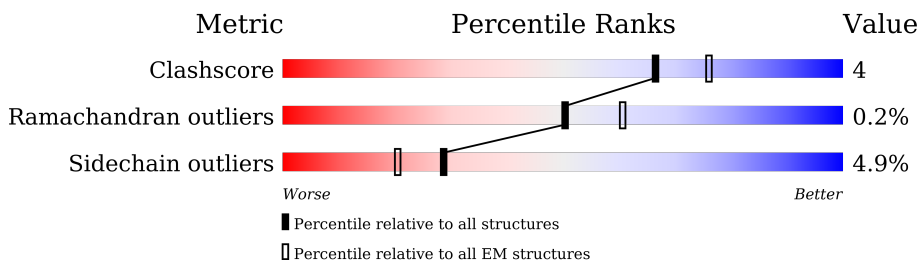
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.35 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 ≥ 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	745	<div style="display: flex; align-items: center;"> <div style="width: 56%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green);"></div> <div style="margin-left: 10px;">56%</div> </div> <div style="display: flex; align-items: center; margin-top: 5px;"> <div style="width: 72%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green);"></div> <div style="margin-left: 10px;">72%</div> <div style="width: 11%; height: 10px; background-color: grey; margin-left: 10px;"></div> <div style="margin-left: 10px;">11%</div> <div style="width: 16%; height: 10px; background-color: grey; margin-left: 10px;"></div> <div style="margin-left: 10px;">16%</div> </div>
1	B	745	<div style="display: flex; align-items: center;"> <div style="width: 60%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green);"></div> <div style="margin-left: 10px;">60%</div> </div> <div style="display: flex; align-items: center; margin-top: 5px;"> <div style="width: 72%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green);"></div> <div style="margin-left: 10px;">72%</div> <div style="width: 11%; height: 10px; background-color: grey; margin-left: 10px;"></div> <div style="margin-left: 10px;">11%</div> <div style="width: 16%; height: 10px; background-color: grey; margin-left: 10px;"></div> <div style="margin-left: 10px;">16%</div> </div>

2 Entry composition i

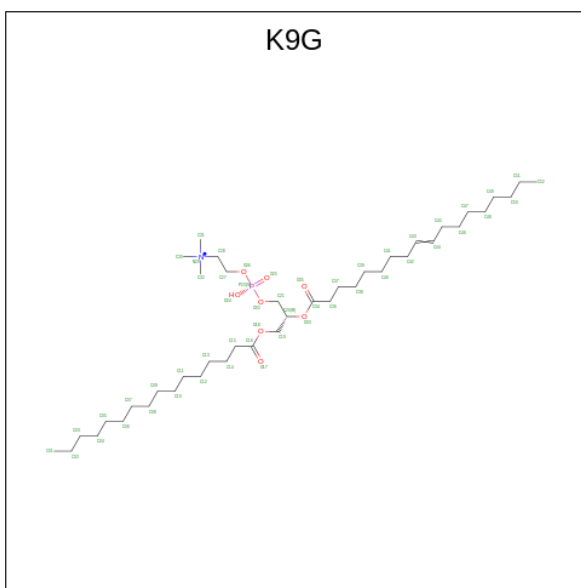
There are 2 unique types of molecules in this entry. The entry contains 9812 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ATP-binding cassette sub-family D member 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	624	Total	C	N	O	S	0	0
			4887	3130	862	873	22		
1	B	623	Total	C	N	O	S	0	0
			4883	3128	861	872	22		

- Molecule 2 is [(2 {R})-1-hexadecanoyloxy-3-[oxidanyl-[2-(trimethyl- l^4 -azanylethoxy]phosphoryl]oxy-propan-2-yl] octadec-9-enoate (three-letter code: K9G) (formula: $C_{42}H_{83}NO_8P$).

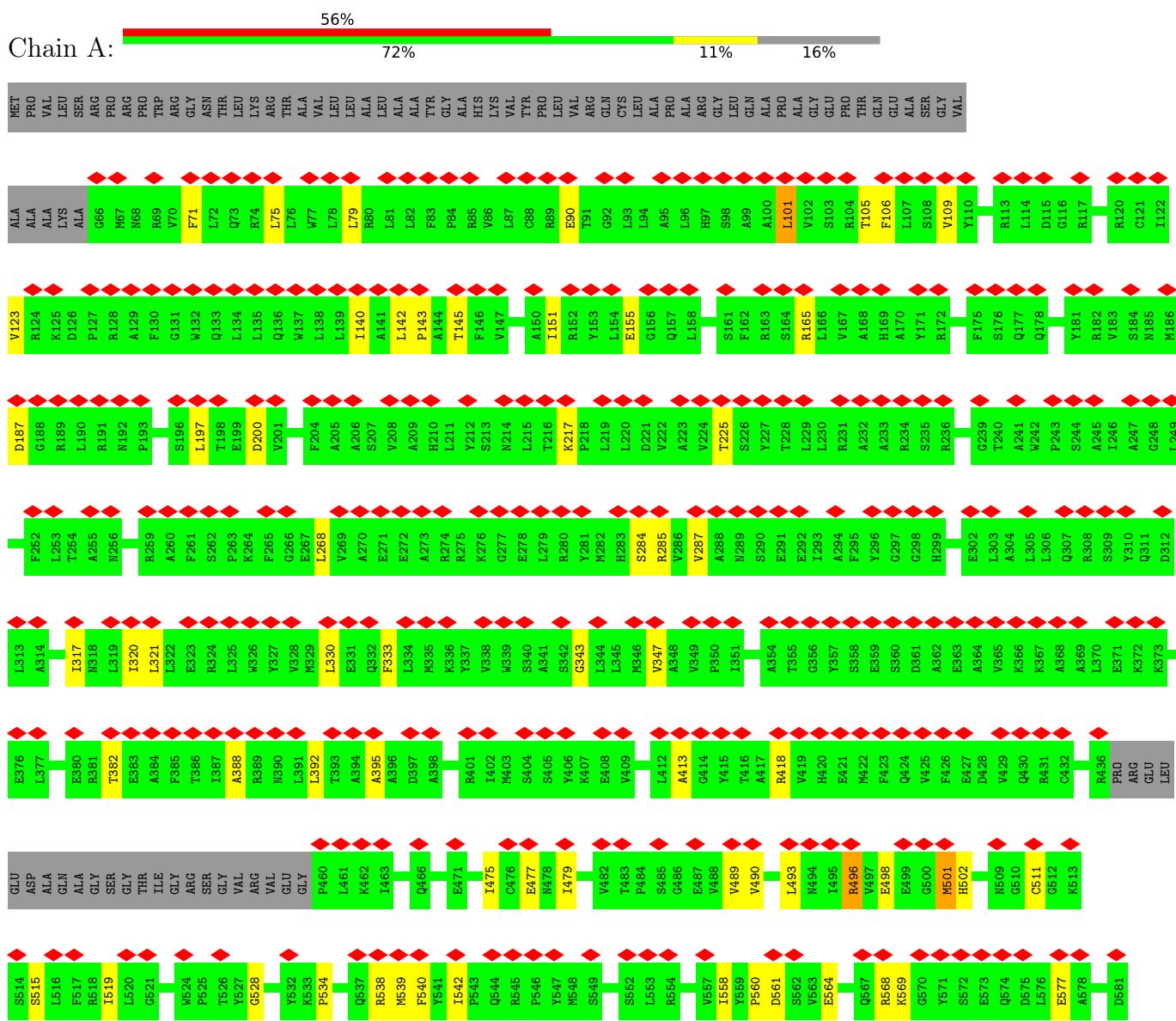


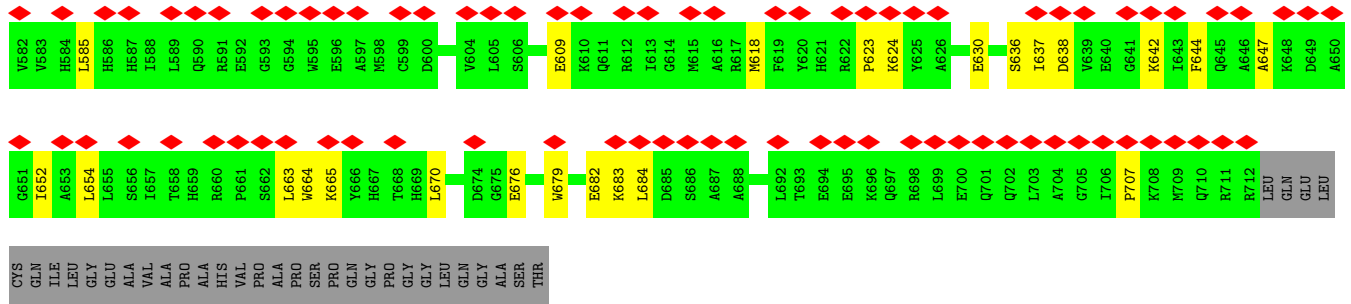
Mol	Chain	Residues	Atoms				AltConf
			Total	C	O	P	
2	A	1	Total	C	O	P	0
			21	13	7	1	
2	B	1	Total	C	O	P	0
			21	13	7	1	

3 Residue-property plots

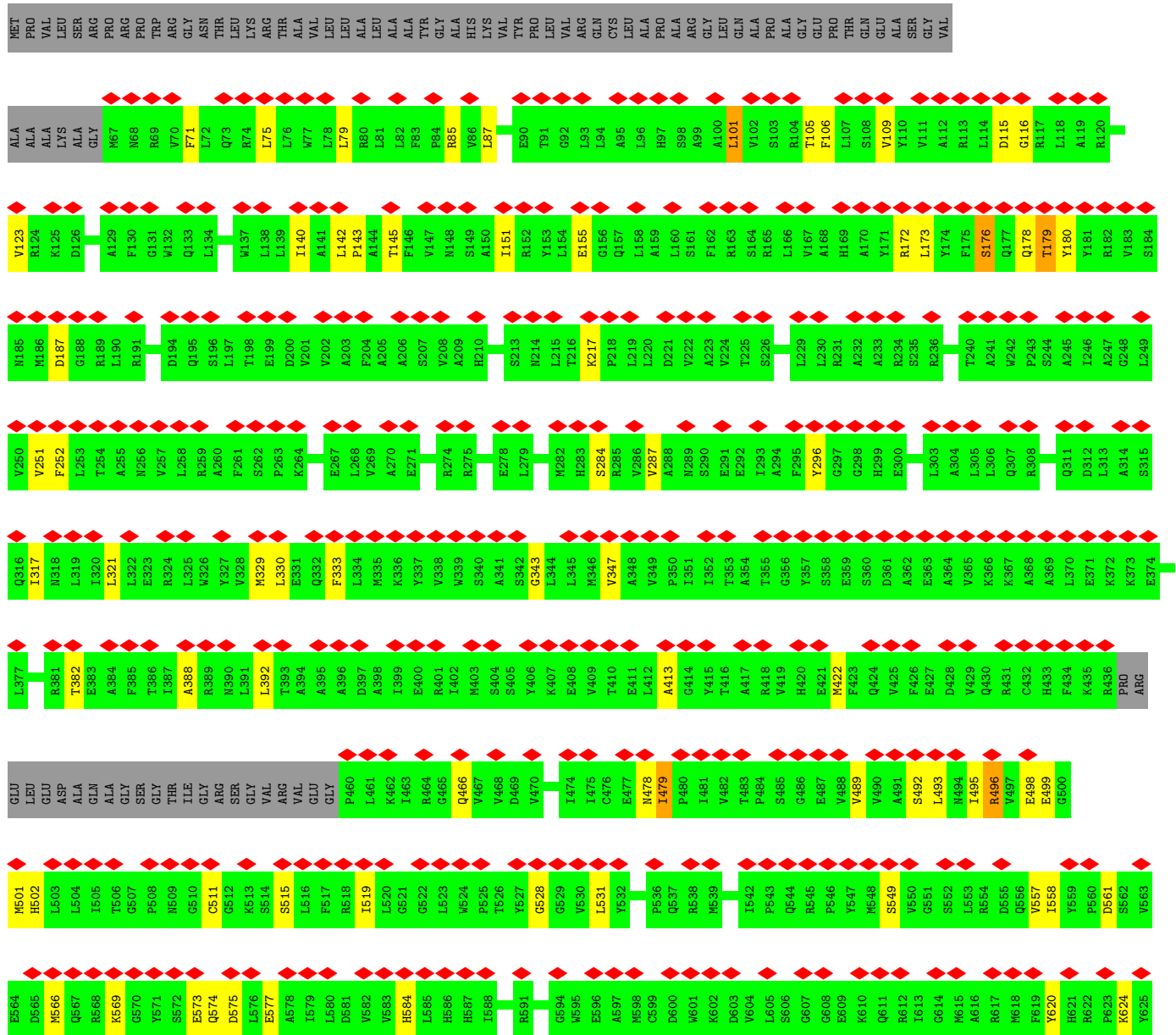
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: ATP-binding cassette sub-family D member 1





• Molecule 1: ATP-binding cassette sub-family D member 1



A626	L627	L628	D629	E630	C631	T632	S633	A634	V635	S636	I637	D638	V639	E640	G641	F644	Q645	A646	A647	K648	I652	A653	L654	L655	S656	I657	T658	H659	R660	P661	S662	L663	W664	K665	Y666	H667	T668	H669	L670	L671	Q672	F673	D674	G675	E676	G677	G678	W679	K680	F681	E682	K683	L684	D685	S686	A687	A688
R689	L690	S691	L692	T693	E694	E695	K696	Q697	R698	L699	E700	Q701	Q702	L703	A704	G705	I706	P707	K708	M709	Q710	R711	R712	LEU	GLN	GLU	LEU	CYS	GLN	ILE	LEU	GLY	GLU	ALA	VAL	ALA	PRO	ALA	HIS	VAL	PRO	ALA	SER	PRO	GLN	GLY	PRO	GLY	GLY	LEU	GLN	GLY	ALA	SER	THR		

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	291703	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 ($e^-/\text{\AA}^2$)	66.5	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	FEI EAGLE (4k x 4k)	Depositor
Maximum map value	1.263	Depositor
Minimum map value	-0.821	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.09	Depositor
Map size (Å)	298.8, 298.8, 298.8	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	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: K9G

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	0.31	0/4991	0.58	2/6763 (0.0%)
1	B	0.31	0/4987	0.58	2/6758 (0.0%)
All	All	0.31	0/9978	0.58	4/13521 (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	A	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	707	PRO	N-CA-CB	6.54	111.15	103.30
1	A	155	GLU	CB-CA-C	-5.90	98.61	110.40
1	B	707	PRO	N-CA-CB	5.75	110.19	103.30
1	B	155	GLU	CB-CA-C	-5.55	99.31	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	684	LEU	Peptide

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	4887	0	4865	40	0
1	B	4883	0	4862	45	0
2	A	21	0	0	0	0
2	B	21	0	0	0	0
All	All	9812	0	9727	79	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 (79) 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:640:GLU:HG2	1:B:663:LEU:HD21	1.70	0.72
1:A:542:ILE:HD11	1:A:618:MET:HB2	1.76	0.67
1:B:498:GLU:H	1:B:501:MET:HG3	1.60	0.66
1:B:101:LEU:HD13	1:B:151:ILE:HD12	1.81	0.63
1:A:101:LEU:HD13	1:A:151:ILE:HD12	1.82	0.61
1:B:251:VAL:O	1:B:251:VAL:HG12	2.01	0.61
1:B:636:SER:HB2	1:B:660:ARG:HH21	1.66	0.60
1:B:176:SER:O	1:B:179:THR:HG23	2.01	0.60
1:B:479:ILE:HD13	1:B:531:LEU:HD13	1.86	0.57
1:A:142:LEU:HB3	1:A:143:PRO:CD	2.36	0.55
1:A:284:SER:HA	1:A:287:VAL:HG12	1.87	0.54
1:A:560:PRO:HG3	1:B:296:TYR:HB3	1.90	0.54
1:B:284:SER:HA	1:B:287:VAL:HG12	1.90	0.52
1:A:623:PRO:HD2	1:A:652:ILE:HG13	1.90	0.52
1:A:145:THR:HG21	1:B:333:PHE:CA	2.41	0.51
1:A:515:SER:O	1:A:519:ILE:HG12	2.11	0.51
1:A:333:PHE:CA	1:B:145:THR:HG21	2.40	0.51
1:A:498:GLU:O	1:A:501:MET:HB2	2.12	0.50
1:B:493:LEU:HD23	1:B:495:ILE:HD12	1.93	0.50
1:A:489:VAL:HG13	1:A:511:CYS:O	2.13	0.49
1:B:115:ASP:OD1	1:B:116:GLY:N	2.46	0.49
1:A:490:VAL:HG21	1:A:493:LEU:HD11	1.96	0.48
1:A:333:PHE:HA	1:B:145:THR:HG21	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:479:ILE:O	1:A:479:ILE:HG13	2.13	0.48
1:B:478:ASN:HA	1:B:492:SER:HB2	1.96	0.48
1:B:382:THR:HG22	1:B:382:THR:O	2.14	0.48
1:B:142:LEU:HB2	1:B:143:PRO:CD	2.44	0.47
1:A:268:LEU:HB3	1:A:320:ILE:HG22	1.96	0.47
1:B:637:ILE:H	1:B:637:ILE:HG13	1.64	0.47
1:B:109:VAL:HG12	1:B:109:VAL:O	2.15	0.47
1:B:557:VAL:O	1:B:558:ILE:HD13	2.14	0.47
1:A:109:VAL:O	1:A:109:VAL:HG12	2.15	0.47
1:B:515:SER:O	1:B:519:ILE:HG12	2.16	0.46
1:B:343:GLY:HA3	1:B:392:LEU:HD21	1.98	0.46
1:A:496:ARG:HE	1:A:496:ARG:HB3	1.54	0.46
1:A:343:GLY:HA3	1:A:392:LEU:HD21	1.98	0.45
1:B:123:VAL:HG21	1:B:382:THR:HG21	1.97	0.45
1:B:178:GLN:HA	1:B:180:TYR:CE1	2.51	0.45
1:B:489:VAL:HG13	1:B:511:CYS:O	2.16	0.45
1:B:140:ILE:HG22	1:B:140:ILE:O	2.16	0.45
1:A:90:GLU:OE1	1:A:165:ARG:NH2	2.49	0.45
1:A:647:ALA:HB1	1:A:652:ILE:HG21	1.98	0.45
1:A:145:THR:HG21	1:B:333:PHE:HA	1.97	0.45
1:A:140:ILE:HG22	1:A:140:ILE:O	2.16	0.45
1:A:123:VAL:HG21	1:A:382:THR:HG21	1.99	0.44
1:B:105:THR:HG23	1:B:106:PHE:CD2	2.52	0.44
1:A:151:ILE:HG21	1:A:217:LYS:HE2	1.99	0.44
1:B:142:LEU:HB2	1:B:143:PRO:HD3	1.99	0.44
1:A:285:ARG:NH1	1:B:549:SER:HA	2.33	0.44
1:B:151:ILE:HG21	1:B:217:LYS:HE2	2.00	0.44
1:B:558:ILE:HG23	1:B:620:TYR:CE2	2.52	0.44
1:B:566:MET:O	1:B:566:MET:SD	2.76	0.44
1:A:75:LEU:O	1:A:79:LEU:HD23	2.18	0.43
1:B:75:LEU:O	1:B:79:LEU:HD23	2.18	0.43
1:B:644:PHE:HE2	1:B:663:LEU:HD13	1.83	0.43
1:A:347:VAL:HG13	1:A:388:ALA:CB	2.48	0.43
1:A:538:ARG:HH11	1:A:538:ARG:HG3	1.83	0.43
1:A:540:PHE:HB2	1:A:623:PRO:HB3	2.01	0.43
1:B:347:VAL:HG13	1:B:388:ALA:CB	2.49	0.43
1:B:496:ARG:HE	1:B:496:ARG:HB3	1.54	0.43
1:A:105:THR:HG23	1:A:106:PHE:CD2	2.53	0.42
1:A:317:ILE:O	1:A:321:LEU:HD13	2.19	0.42
1:A:534:PRO:HD2	1:A:539:MET:HE2	2.02	0.42
1:B:558:ILE:HG22	1:B:561:ASP:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:647:ALA:HB1	1:B:652:ILE:HG21	2.01	0.42
1:A:71:PHE:CE1	1:A:413:ALA:HA	2.55	0.41
1:A:558:ILE:HG22	1:A:561:ASP:HB2	2.02	0.41
1:A:475:ILE:HG22	1:A:477:GLU:HG3	2.02	0.41
1:A:585:LEU:HD21	1:A:609:GLU:HG3	2.02	0.41
1:B:317:ILE:O	1:B:321:LEU:HD13	2.20	0.41
1:B:329:MET:O	1:B:329:MET:HG2	2.20	0.41
1:A:225:THR:OG1	1:A:395:ALA:HB2	2.21	0.41
1:B:71:PHE:CE1	1:B:413:ALA:HA	2.56	0.41
1:B:624:LYS:HD2	1:B:624:LYS:HA	1.92	0.41
1:A:644:PHE:HE2	1:A:663:LEU:HD13	1.85	0.41
1:B:172:ARG:O	1:B:176:SER:HB3	2.21	0.41
1:B:502:HIS:HA	1:B:654:LEU:O	2.21	0.41
1:A:502:HIS:HA	1:A:654:LEU:O	2.21	0.40
1:A:200:ASP:OD2	1:A:418:ARG:HB3	2.21	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	620/745 (83%)	574 (93%)	45 (7%)	1 (0%)	47 78
1	B	619/745 (83%)	575 (93%)	42 (7%)	2 (0%)	41 73
All	All	1239/1490 (83%)	1149 (93%)	87 (7%)	3 (0%)	50 78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	528	GLY
1	B	528	GLY
1	B	573	GLU

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	500/610 (82%)	477 (95%)	23 (5%)	27	59
1	B	500/610 (82%)	474 (95%)	26 (5%)	23	55
All	All	1000/1220 (82%)	951 (95%)	49 (5%)	29	57

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	LEU
1	A	187	ASP
1	A	197	LEU
1	A	330	LEU
1	A	496	ARG
1	A	501	MET
1	A	564	GLU
1	A	568	ARG
1	A	569	LYS
1	A	577	GLU
1	A	624	LYS
1	A	630	GLU
1	A	636	SER
1	A	637	ILE
1	A	638	ASP
1	A	642	LYS
1	A	664	TRP
1	A	665	LYS
1	A	670	LEU
1	A	676	GLU
1	A	679	TRP
1	A	682	GLU
1	A	683	LYS
1	B	85	ARG
1	B	87	LEU
1	B	101	LEU
1	B	173	LEU

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Mol	Chain	Res	Type
1	B	176	SER
1	B	179	THR
1	B	187	ASP
1	B	252	PHE
1	B	330	LEU
1	B	422	MET
1	B	466	GLN
1	B	479	ILE
1	B	496	ARG
1	B	499	GLU
1	B	569	LYS
1	B	574	GLN
1	B	575	ASP
1	B	577	GLU
1	B	584	HIS
1	B	630	GLU
1	B	635	VAL
1	B	637	ILE
1	B	638	ASP
1	B	640	GLU
1	B	662	SER
1	B	664	TRP

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	K9G	A	801	-	20,20,51	1.48	4 (20%)	22,24,59	0.76	1 (4%)
2	K9G	B	801	-	20,20,51	1.54	4 (20%)	22,24,59	0.82	1 (4%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	K9G	A	801	-	-	10/22/22/55	-
2	K9G	B	801	-	-	13/22/22/55	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	K9G	O18-C16	4.21	1.45	1.33
2	A	801	K9G	O18-C16	3.98	1.45	1.33
2	B	801	K9G	P23-O26	2.72	1.69	1.59
2	B	801	K9G	P23-O22	2.66	1.70	1.59
2	A	801	K9G	P23-O22	2.65	1.70	1.59
2	A	801	K9G	P23-O26	2.64	1.68	1.59
2	B	801	K9G	C15-C16	2.18	1.57	1.50
2	A	801	K9G	C15-C16	2.02	1.56	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	K9G	O18-C16-C15	2.44	119.55	111.91
2	A	801	K9G	O18-C16-C15	2.16	118.70	111.91

There are no chirality outliers.

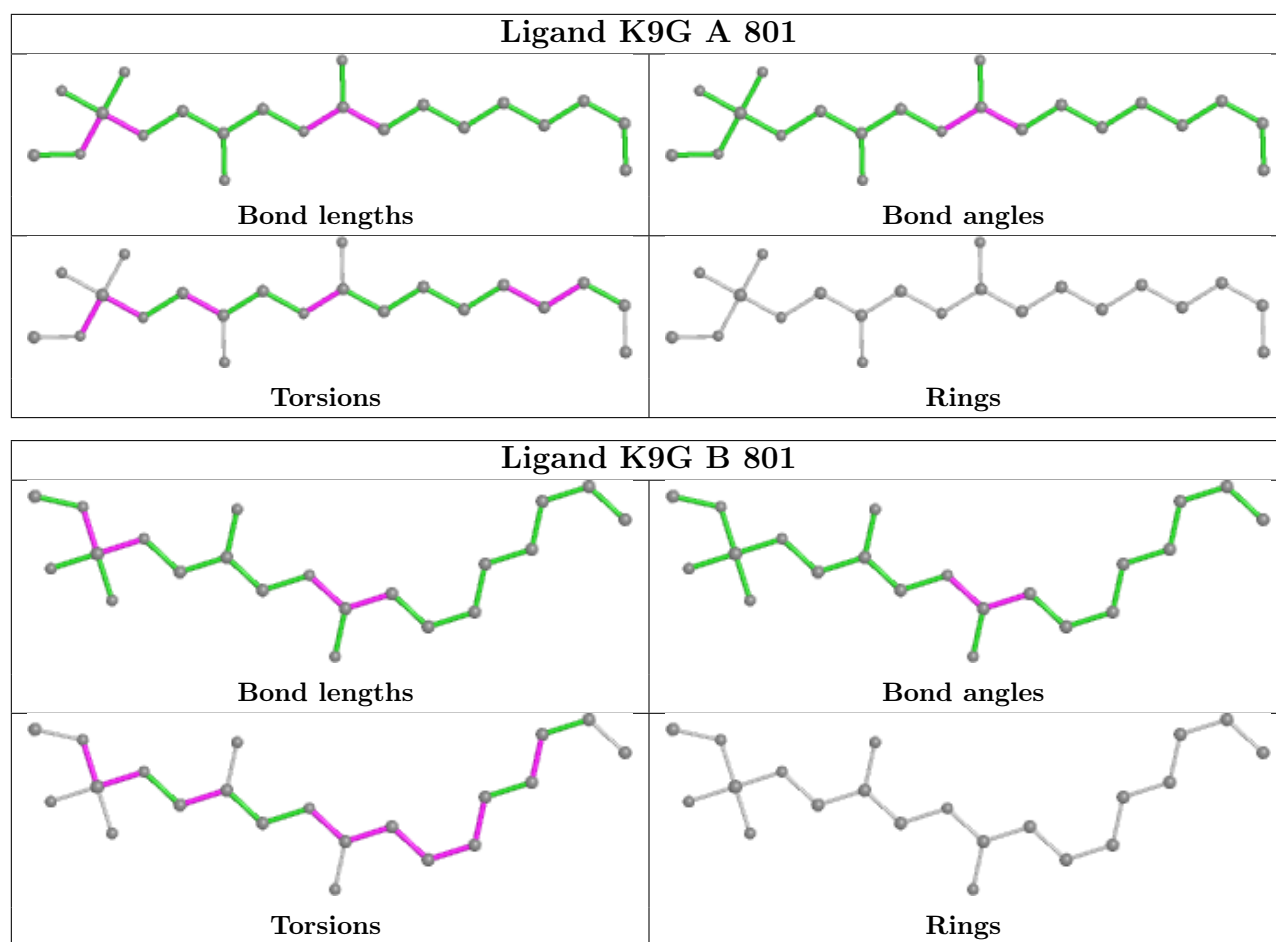
All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	K9G	C21-O22-P23-O24
2	A	801	K9G	C21-O22-P23-O25
2	A	801	K9G	C21-O22-P23-O26
2	B	801	K9G	C21-O22-P23-O24
2	B	801	K9G	C21-O22-P23-O25
2	B	801	K9G	C21-O22-P23-O26
2	B	801	K9G	O17-C16-O18-C19
2	B	801	K9G	C12-C13-C14-C15
2	A	801	K9G	O17-C16-O18-C19
2	A	801	K9G	C15-C16-O18-C19
2	B	801	K9G	C15-C16-O18-C19
2	B	801	K9G	C13-C14-C15-C16
2	A	801	K9G	O33-C20-C21-O22
2	B	801	K9G	O33-C20-C21-O22
2	A	801	K9G	C10-C11-C12-C13
2	A	801	K9G	C19-C20-C21-O22
2	B	801	K9G	C19-C20-C21-O22
2	A	801	K9G	C27-O26-P23-O25
2	B	801	K9G	C27-O26-P23-O25
2	B	801	K9G	C11-C12-C13-C14
2	A	801	K9G	C09-C10-C11-C12
2	B	801	K9G	C09-C10-C11-C12
2	B	801	K9G	C14-C15-C16-O18

There are no ring outliers.

No monomer is involved in short contacts.

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

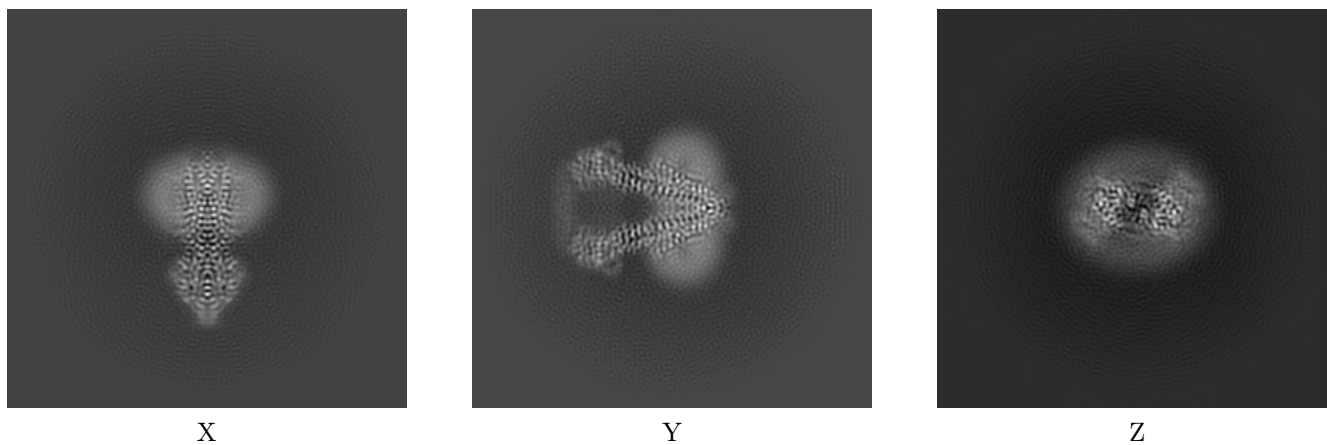
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34064. These allow visual inspection of the internal detail of the map and identification of artifacts.

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

6.1 Orthogonal projections [i](#)

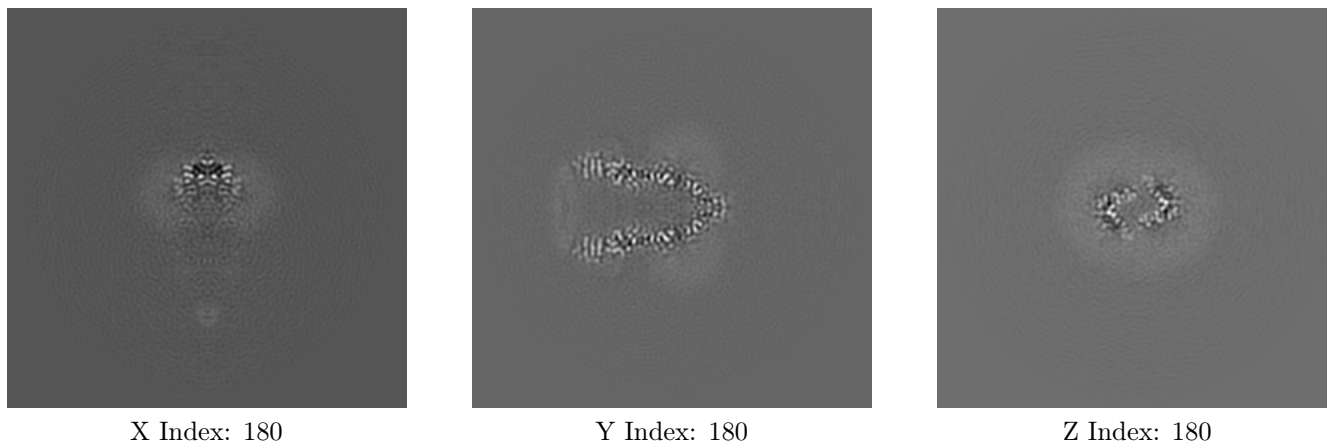
6.1.1 Primary map



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

6.2 Central slices [i](#)

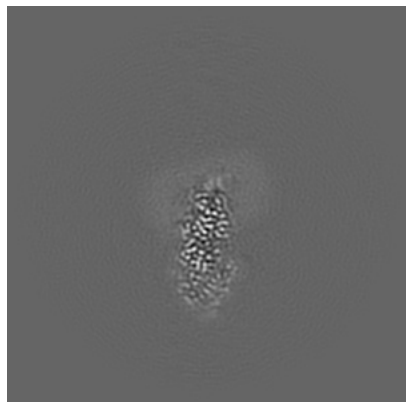
6.2.1 Primary map



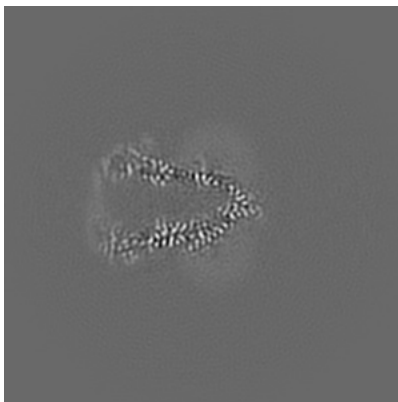
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

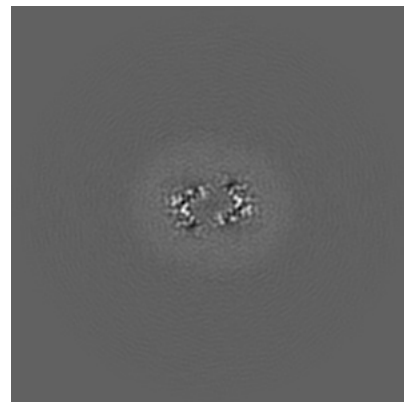
6.3.1 Primary map



X Index: 208



Y Index: 185



Z Index: 179

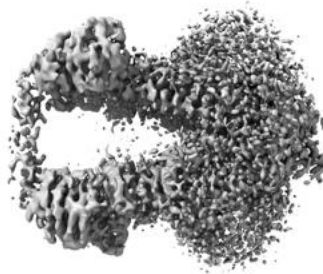
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views [i](#)

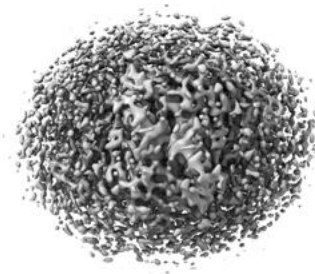
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.09. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

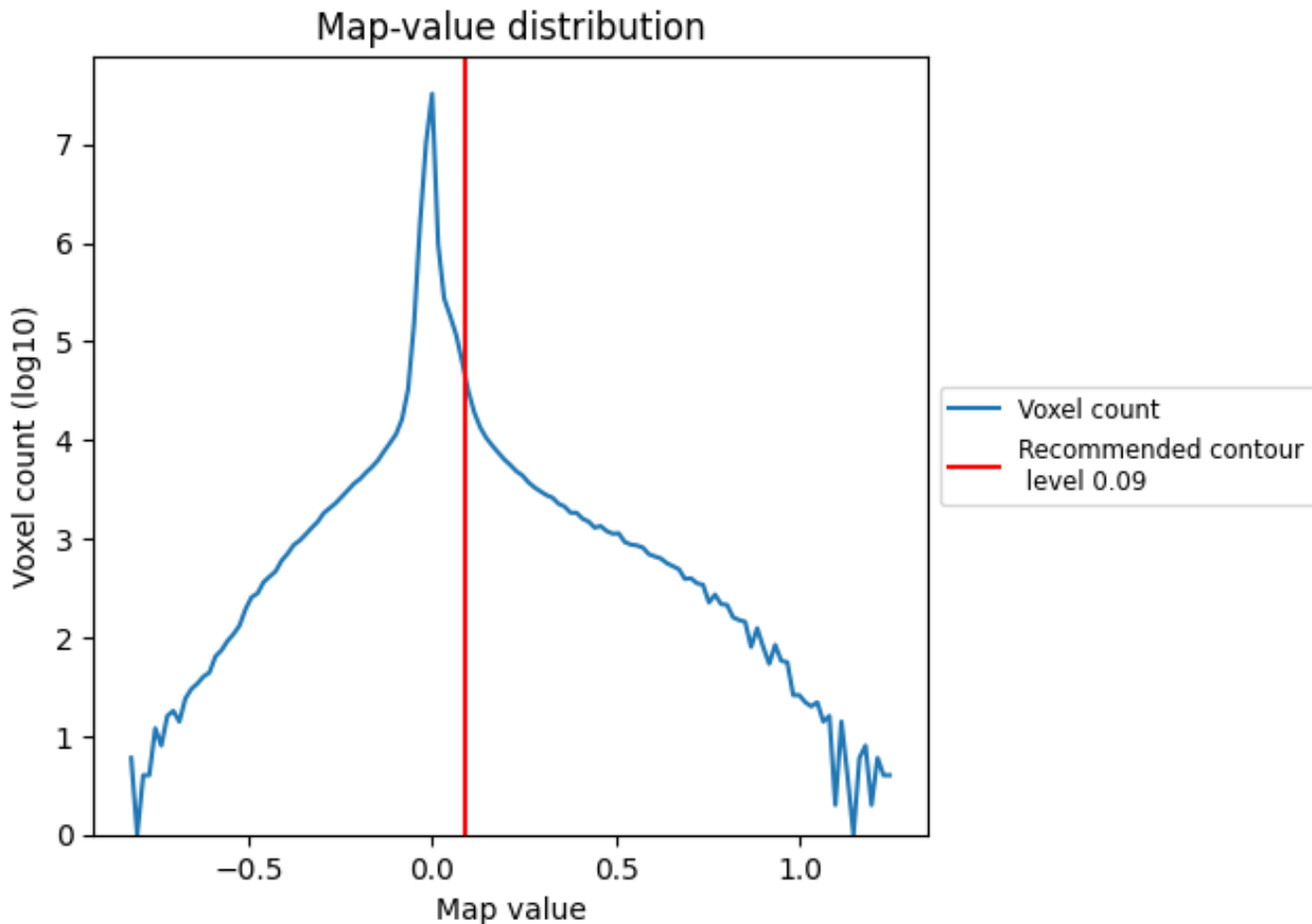
6.5 Mask visualisation

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

7 Map analysis [i](#)

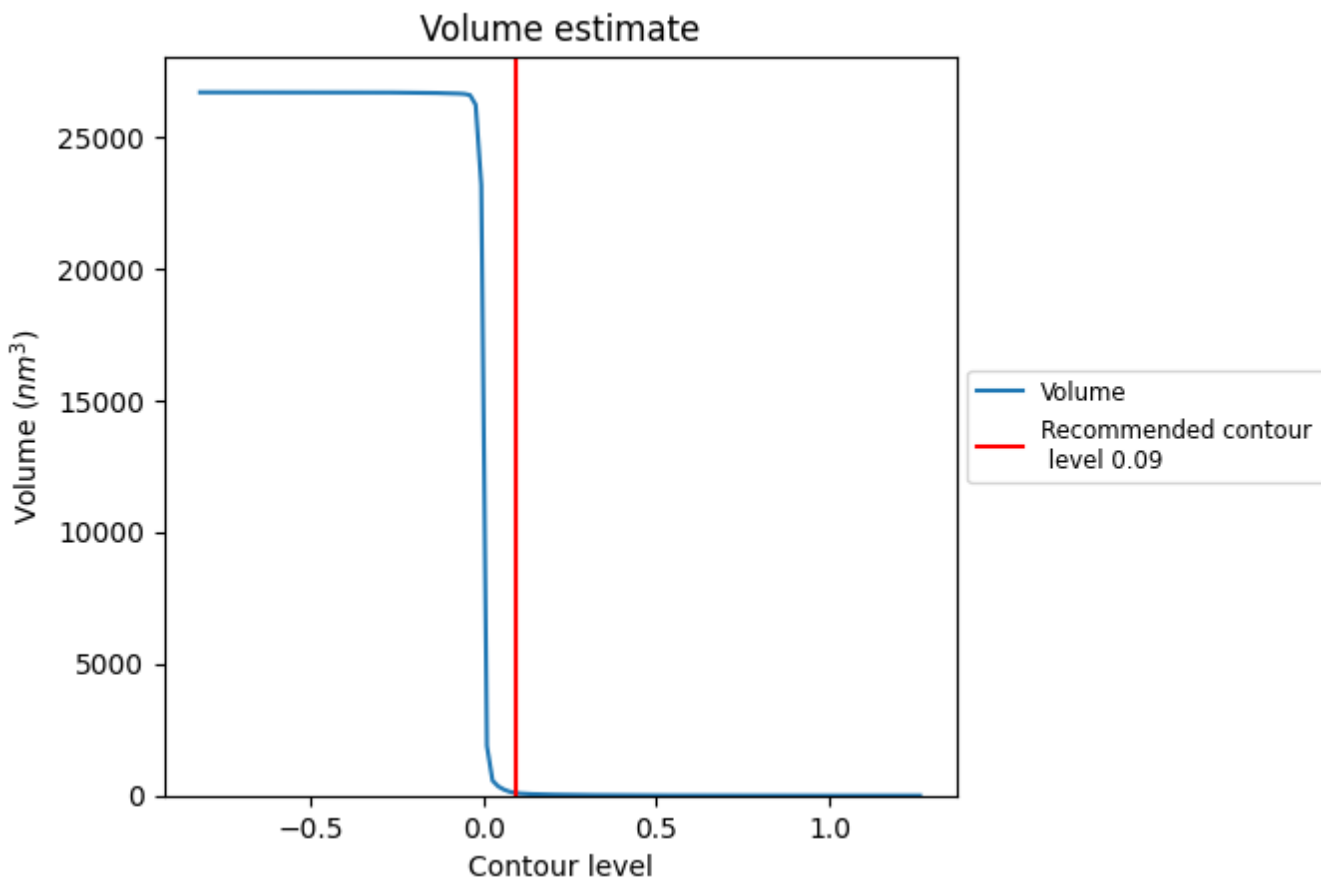
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

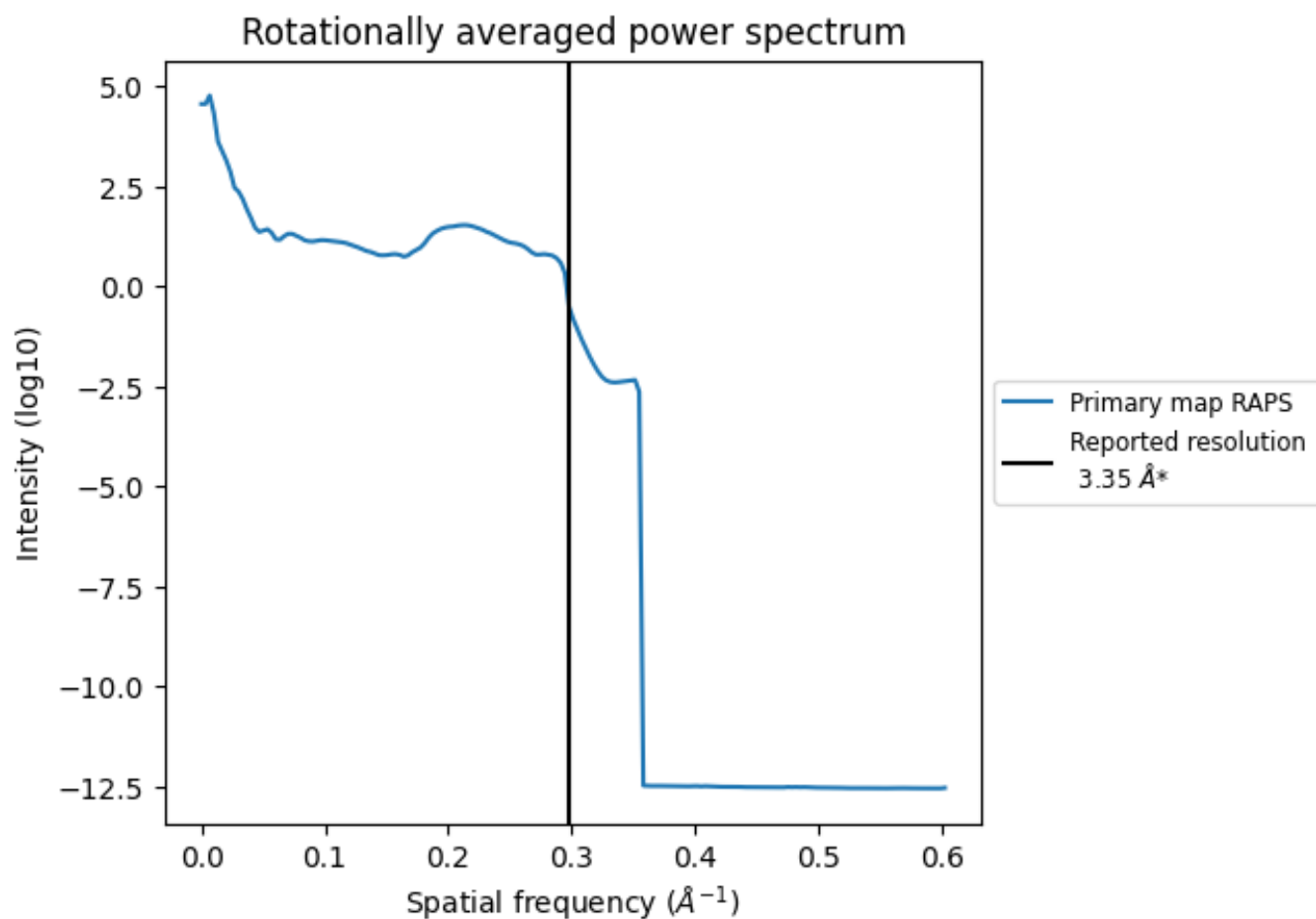
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 103 nm³; this corresponds to an approximate mass of 93 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 0.299 Å⁻¹

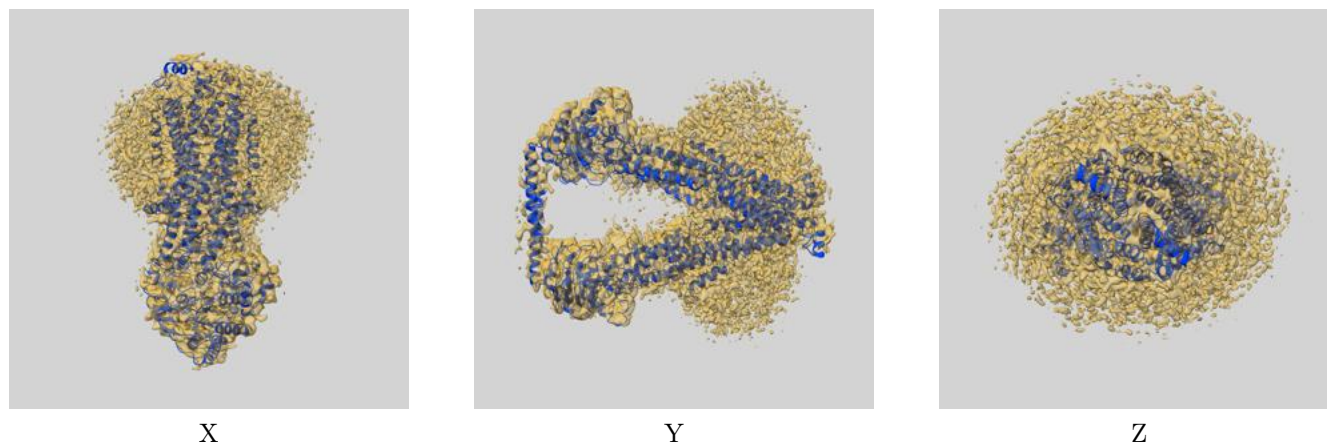
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

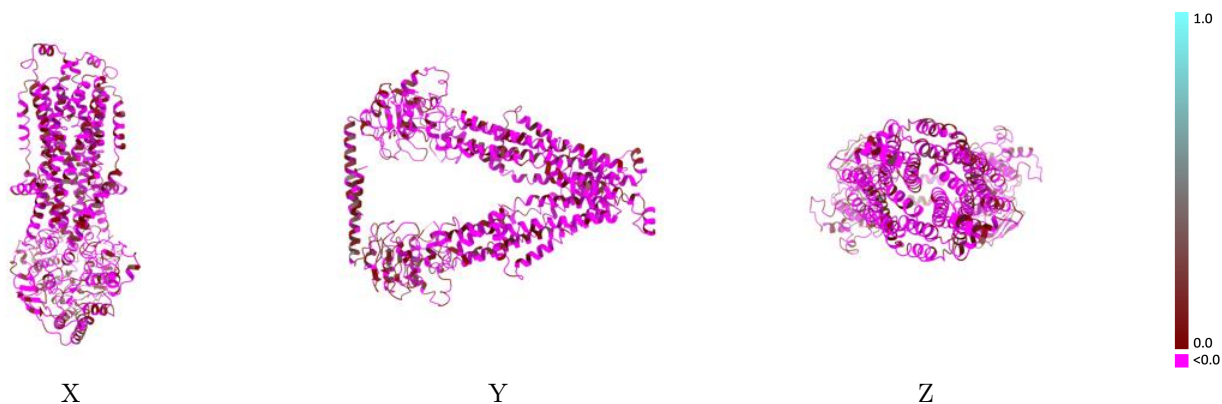
This section contains information regarding the fit between EMDB map EMD-34064 and PDB model 7YRQ. Per-residue inclusion information can be found in section [3](#) on page [4](#).

9.1 Map-model overlay [i](#)



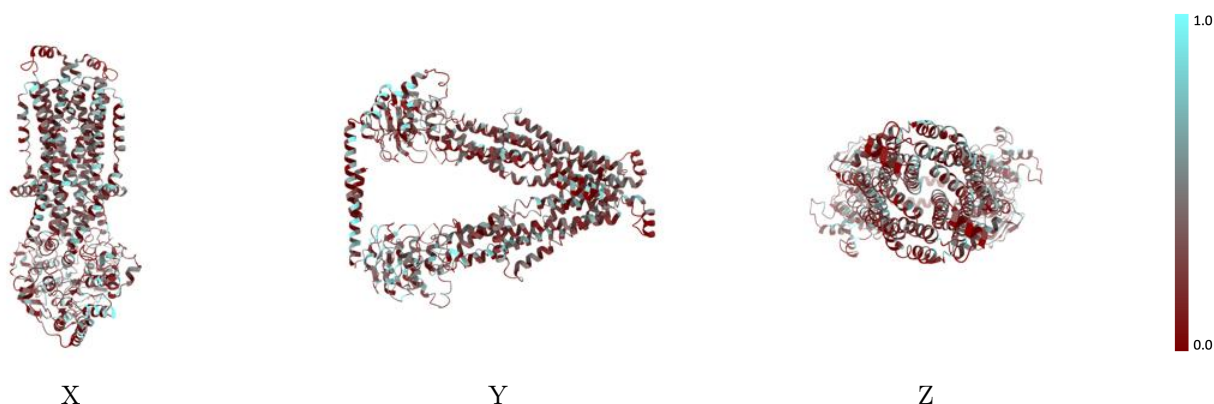
The images above show the 3D surface view of the map at the recommended contour level 0.09 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



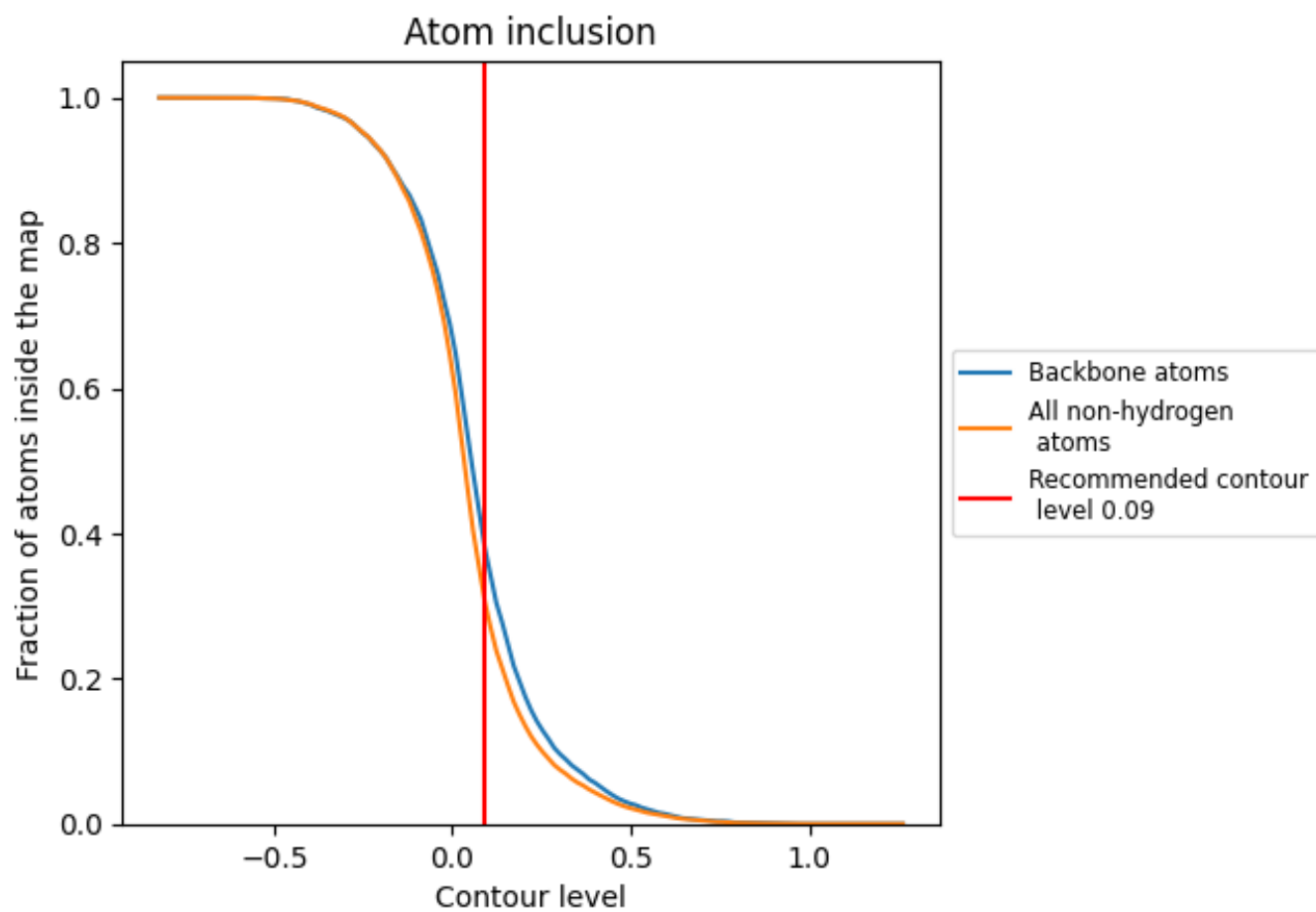
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.09).







9.4 Atom inclusion [i](#)



At the recommended contour level, 38% of all backbone atoms, 31% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary [i](#)

The table lists the average atom inclusion at the recommended contour level (0.09) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.3081	 -0.0660
A	 0.3230	 -0.0480
B	 0.2931	 -0.0840

