



## Full wwPDB EM Validation Report ⓘ

Feb 13, 2023 – 04:16 PM JST

PDB ID : 7XNL  
EMDB ID : EMD-33318  
Title : human KCNQ1-CaM-ML277-PIP2 complex in state A  
Authors : Ma, D.; Guo, J.  
Deposited on : 2022-04-29  
Resolution : 3.10 Å (reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

---

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

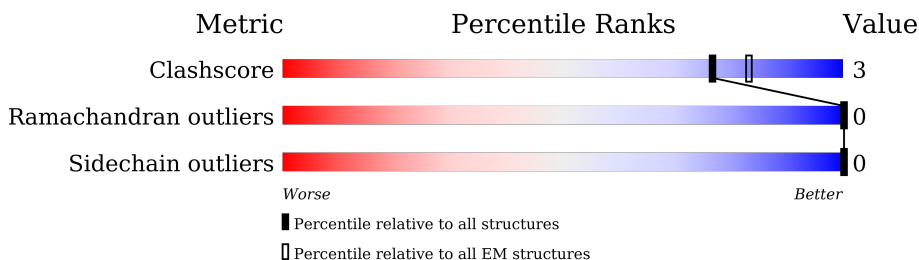
EMDB validation analysis : 0.0.1.dev43  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
MolProbity : 4.02b-467  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.32.1

# 1 Overall quality at a glance

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

The reported resolution of this entry is 3.10 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	692	45% 5% 50%
1	C	692	46% 5% 50%
1	E	692	46% 5% 50%
1	G	692	46% 5% 50%
2	B	177	24% 79% 19%
2	D	177	24% 77% 19%
2	F	177	25% 79% 19%
2	H	177	24% 79% 19%

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 31964 atoms, of which 15928 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 Potassium voltage-gated channel subfamily KQT member 1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
1	A	348	5712	1851	2891	492	467	11	0	0
1	C	348	5712	1851	2891	492	467	11	0	0
1	E	348	5712	1851	2891	492	467	11	0	0
1	G	348	5712	1851	2891	492	467	11	0	0

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	677	LEU	-	expression tag	UNP P51787
A	678	GLU	-	expression tag	UNP P51787
A	679	GLY	-	expression tag	UNP P51787
A	680	GLY	-	expression tag	UNP P51787
A	681	SER	-	expression tag	UNP P51787
A	682	SER	-	expression tag	UNP P51787
A	683	GLY	-	expression tag	UNP P51787
A	684	GLY	-	expression tag	UNP P51787
A	685	TRP	-	expression tag	UNP P51787
A	686	SER	-	expression tag	UNP P51787
A	687	HIS	-	expression tag	UNP P51787
A	688	PRO	-	expression tag	UNP P51787
A	689	GLN	-	expression tag	UNP P51787
A	690	PHE	-	expression tag	UNP P51787
A	691	GLU	-	expression tag	UNP P51787
A	692	LYS	-	expression tag	UNP P51787
C	677	LEU	-	expression tag	UNP P51787
C	678	GLU	-	expression tag	UNP P51787
C	679	GLY	-	expression tag	UNP P51787
C	680	GLY	-	expression tag	UNP P51787
C	681	SER	-	expression tag	UNP P51787
C	682	SER	-	expression tag	UNP P51787

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
C	683	GLY	-	expression tag	UNP P51787
C	684	GLY	-	expression tag	UNP P51787
C	685	TRP	-	expression tag	UNP P51787
C	686	SER	-	expression tag	UNP P51787
C	687	HIS	-	expression tag	UNP P51787
C	688	PRO	-	expression tag	UNP P51787
C	689	GLN	-	expression tag	UNP P51787
C	690	PHE	-	expression tag	UNP P51787
C	691	GLU	-	expression tag	UNP P51787
C	692	LYS	-	expression tag	UNP P51787
E	677	LEU	-	expression tag	UNP P51787
E	678	GLU	-	expression tag	UNP P51787
E	679	GLY	-	expression tag	UNP P51787
E	680	GLY	-	expression tag	UNP P51787
E	681	SER	-	expression tag	UNP P51787
E	682	SER	-	expression tag	UNP P51787
E	683	GLY	-	expression tag	UNP P51787
E	684	GLY	-	expression tag	UNP P51787
E	685	TRP	-	expression tag	UNP P51787
E	686	SER	-	expression tag	UNP P51787
E	687	HIS	-	expression tag	UNP P51787
E	688	PRO	-	expression tag	UNP P51787
E	689	GLN	-	expression tag	UNP P51787
E	690	PHE	-	expression tag	UNP P51787
E	691	GLU	-	expression tag	UNP P51787
E	692	LYS	-	expression tag	UNP P51787
G	677	LEU	-	expression tag	UNP P51787
G	678	GLU	-	expression tag	UNP P51787
G	679	GLY	-	expression tag	UNP P51787
G	680	GLY	-	expression tag	UNP P51787
G	681	SER	-	expression tag	UNP P51787
G	682	SER	-	expression tag	UNP P51787
G	683	GLY	-	expression tag	UNP P51787
G	684	GLY	-	expression tag	UNP P51787
G	685	TRP	-	expression tag	UNP P51787
G	686	SER	-	expression tag	UNP P51787
G	687	HIS	-	expression tag	UNP P51787
G	688	PRO	-	expression tag	UNP P51787
G	689	GLN	-	expression tag	UNP P51787
G	690	PHE	-	expression tag	UNP P51787
G	691	GLU	-	expression tag	UNP P51787
G	692	LYS	-	expression tag	UNP P51787

- Molecule 2 is a protein called Calmodulin-3.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
2	B	144	2130	682	1022	180	237	9	0	0
2	D	144	2130	682	1022	180	237	9	0	0
2	F	144	2130	682	1022	180	237	9	0	0
2	H	144	2130	682	1022	180	237	9	0	0

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	150	LEU	-	expression tag	UNP P0DP25
B	151	GLU	-	expression tag	UNP P0DP25
B	152	GLY	-	expression tag	UNP P0DP25
B	153	GLY	-	expression tag	UNP P0DP25
B	154	SER	-	expression tag	UNP P0DP25
B	155	SER	-	expression tag	UNP P0DP25
B	156	GLY	-	expression tag	UNP P0DP25
B	157	GLY	-	expression tag	UNP P0DP25
B	158	LEU	-	expression tag	UNP P0DP25
B	159	VAL	-	expression tag	UNP P0DP25
B	160	PRO	-	expression tag	UNP P0DP25
B	161	ARG	-	expression tag	UNP P0DP25
B	162	GLY	-	expression tag	UNP P0DP25
B	163	SER	-	expression tag	UNP P0DP25
B	164	GLY	-	expression tag	UNP P0DP25
B	165	GLY	-	expression tag	UNP P0DP25
B	166	SER	-	expression tag	UNP P0DP25
B	167	SER	-	expression tag	UNP P0DP25
B	168	GLY	-	expression tag	UNP P0DP25
B	169	GLY	-	expression tag	UNP P0DP25
B	170	HIS	-	expression tag	UNP P0DP25
B	171	HIS	-	expression tag	UNP P0DP25
B	172	HIS	-	expression tag	UNP P0DP25
B	173	HIS	-	expression tag	UNP P0DP25
B	174	HIS	-	expression tag	UNP P0DP25
B	175	HIS	-	expression tag	UNP P0DP25
B	176	HIS	-	expression tag	UNP P0DP25
B	177	HIS	-	expression tag	UNP P0DP25
D	150	LEU	-	expression tag	UNP P0DP25
D	151	GLU	-	expression tag	UNP P0DP25

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
D	152	GLY	-	expression tag	UNP P0DP25
D	153	GLY	-	expression tag	UNP P0DP25
D	154	SER	-	expression tag	UNP P0DP25
D	155	SER	-	expression tag	UNP P0DP25
D	156	GLY	-	expression tag	UNP P0DP25
D	157	GLY	-	expression tag	UNP P0DP25
D	158	LEU	-	expression tag	UNP P0DP25
D	159	VAL	-	expression tag	UNP P0DP25
D	160	PRO	-	expression tag	UNP P0DP25
D	161	ARG	-	expression tag	UNP P0DP25
D	162	GLY	-	expression tag	UNP P0DP25
D	163	SER	-	expression tag	UNP P0DP25
D	164	GLY	-	expression tag	UNP P0DP25
D	165	GLY	-	expression tag	UNP P0DP25
D	166	SER	-	expression tag	UNP P0DP25
D	167	SER	-	expression tag	UNP P0DP25
D	168	GLY	-	expression tag	UNP P0DP25
D	169	GLY	-	expression tag	UNP P0DP25
D	170	HIS	-	expression tag	UNP P0DP25
D	171	HIS	-	expression tag	UNP P0DP25
D	172	HIS	-	expression tag	UNP P0DP25
D	173	HIS	-	expression tag	UNP P0DP25
D	174	HIS	-	expression tag	UNP P0DP25
D	175	HIS	-	expression tag	UNP P0DP25
D	176	HIS	-	expression tag	UNP P0DP25
D	177	HIS	-	expression tag	UNP P0DP25
F	150	LEU	-	expression tag	UNP P0DP25
F	151	GLU	-	expression tag	UNP P0DP25
F	152	GLY	-	expression tag	UNP P0DP25
F	153	GLY	-	expression tag	UNP P0DP25
F	154	SER	-	expression tag	UNP P0DP25
F	155	SER	-	expression tag	UNP P0DP25
F	156	GLY	-	expression tag	UNP P0DP25
F	157	GLY	-	expression tag	UNP P0DP25
F	158	LEU	-	expression tag	UNP P0DP25
F	159	VAL	-	expression tag	UNP P0DP25
F	160	PRO	-	expression tag	UNP P0DP25
F	161	ARG	-	expression tag	UNP P0DP25
F	162	GLY	-	expression tag	UNP P0DP25
F	163	SER	-	expression tag	UNP P0DP25
F	164	GLY	-	expression tag	UNP P0DP25
F	165	GLY	-	expression tag	UNP P0DP25

*Continued on next page...*

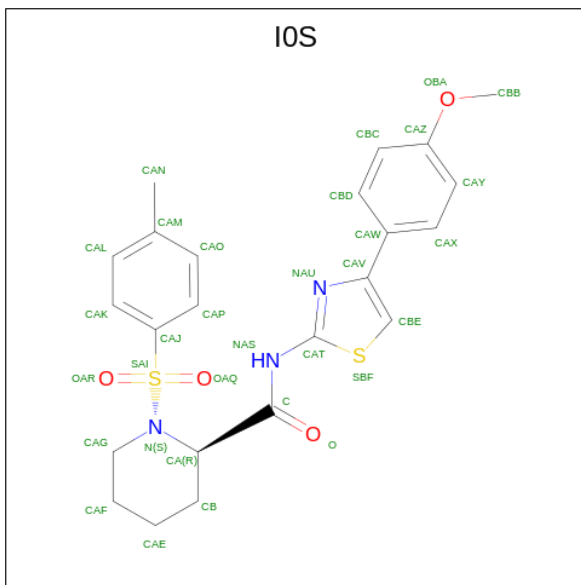
*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
F	166	SER	-	expression tag	UNP P0DP25
F	167	SER	-	expression tag	UNP P0DP25
F	168	GLY	-	expression tag	UNP P0DP25
F	169	GLY	-	expression tag	UNP P0DP25
F	170	HIS	-	expression tag	UNP P0DP25
F	171	HIS	-	expression tag	UNP P0DP25
F	172	HIS	-	expression tag	UNP P0DP25
F	173	HIS	-	expression tag	UNP P0DP25
F	174	HIS	-	expression tag	UNP P0DP25
F	175	HIS	-	expression tag	UNP P0DP25
F	176	HIS	-	expression tag	UNP P0DP25
F	177	HIS	-	expression tag	UNP P0DP25
H	150	LEU	-	expression tag	UNP P0DP25
H	151	GLU	-	expression tag	UNP P0DP25
H	152	GLY	-	expression tag	UNP P0DP25
H	153	GLY	-	expression tag	UNP P0DP25
H	154	SER	-	expression tag	UNP P0DP25
H	155	SER	-	expression tag	UNP P0DP25
H	156	GLY	-	expression tag	UNP P0DP25
H	157	GLY	-	expression tag	UNP P0DP25
H	158	LEU	-	expression tag	UNP P0DP25
H	159	VAL	-	expression tag	UNP P0DP25
H	160	PRO	-	expression tag	UNP P0DP25
H	161	ARG	-	expression tag	UNP P0DP25
H	162	GLY	-	expression tag	UNP P0DP25
H	163	SER	-	expression tag	UNP P0DP25
H	164	GLY	-	expression tag	UNP P0DP25
H	165	GLY	-	expression tag	UNP P0DP25
H	166	SER	-	expression tag	UNP P0DP25
H	167	SER	-	expression tag	UNP P0DP25
H	168	GLY	-	expression tag	UNP P0DP25
H	169	GLY	-	expression tag	UNP P0DP25
H	170	HIS	-	expression tag	UNP P0DP25
H	171	HIS	-	expression tag	UNP P0DP25
H	172	HIS	-	expression tag	UNP P0DP25
H	173	HIS	-	expression tag	UNP P0DP25
H	174	HIS	-	expression tag	UNP P0DP25
H	175	HIS	-	expression tag	UNP P0DP25
H	176	HIS	-	expression tag	UNP P0DP25
H	177	HIS	-	expression tag	UNP P0DP25

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	AltConf
3	A	4	Total K 4 4	0

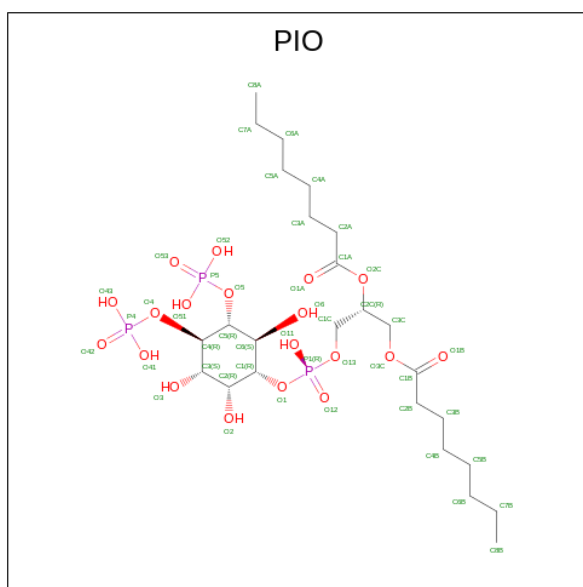
- Molecule 4 is (2R)-N-[4-(4-methoxyphenyl)-1,3-thiazol-2-yl]-1-(4-methylbenzene-1-sulfonyl)piperidine-2-carboxamide (three-letter code: IOS) (formula: C<sub>23</sub>H<sub>25</sub>N<sub>3</sub>O<sub>4</sub>S<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf	
			Total	C	H	N	O		S
4	A	1	Total 57	C 23	H 25	N 3	O 4	S 2	0
4	C	1	Total 57	C 23	H 25	N 3	O 4	S 2	0
4	E	1	Total 57	C 23	H 25	N 3	O 4	S 2	0
4	G	1	Total 57	C 23	H 25	N 3	O 4	S 2	0

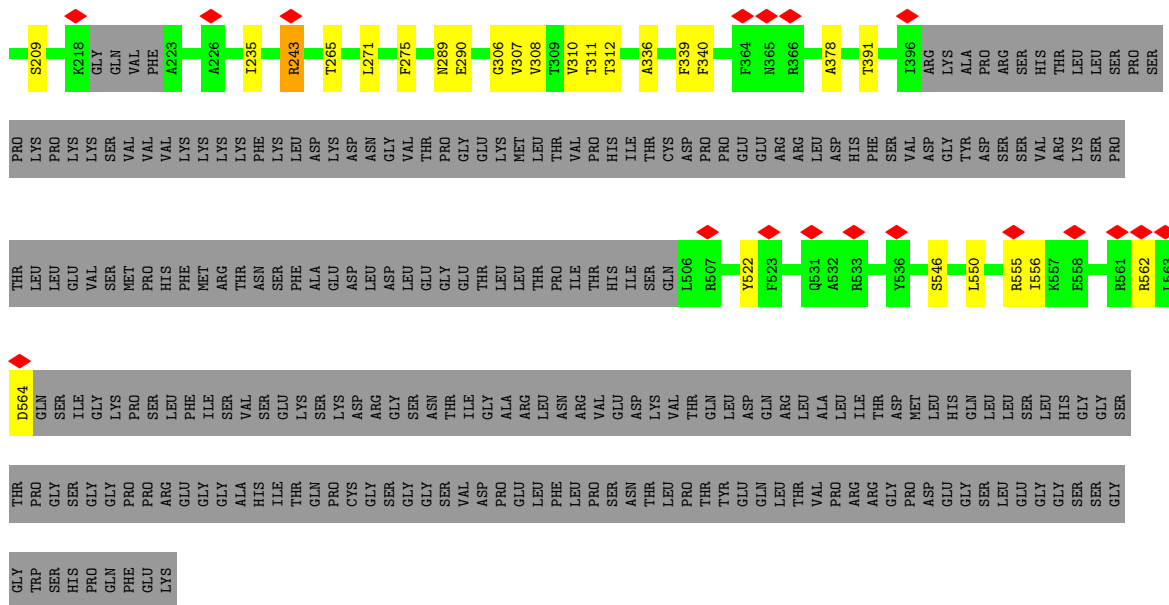
- Molecule 5 is [(2R)-2-octanoyloxy-3-[oxidanyl-[(1R,2R,3S,4R,5R,6S)-2,3,6-tris(oxidanyl)-4,5-diphosphonoxy-cyclohexyl]oxy-phosphoryl]oxy-propyl] octanoate (three-letter code: PIO) (formula: C<sub>25</sub>H<sub>49</sub>O<sub>19</sub>P<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



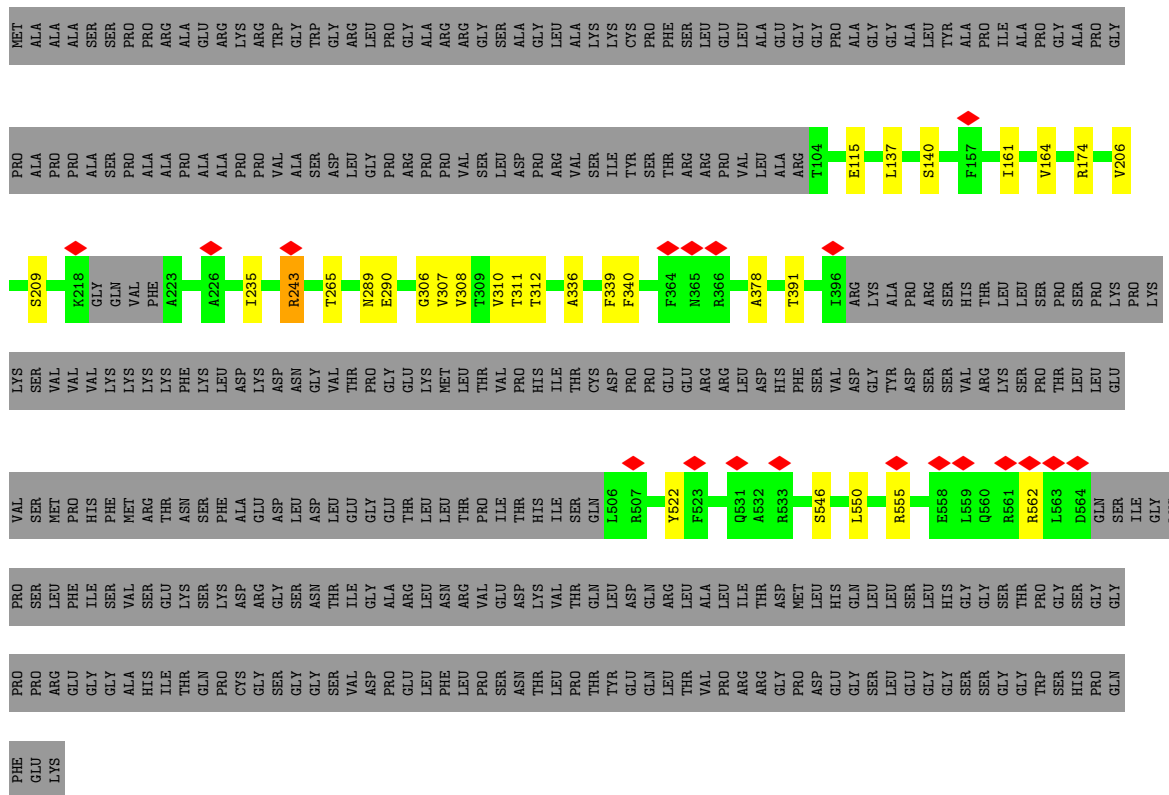


Mol	Chain	Residues	Atoms				AltConf		
			Total	C	H	O		P	
5	B	1	Total	91	25	44	19	3	0
5	D	1	Total	91	25	44	19	3	0
5	F	1	Total	91	25	44	19	3	0
5	H	1	Total	91	25	44	19	3	0



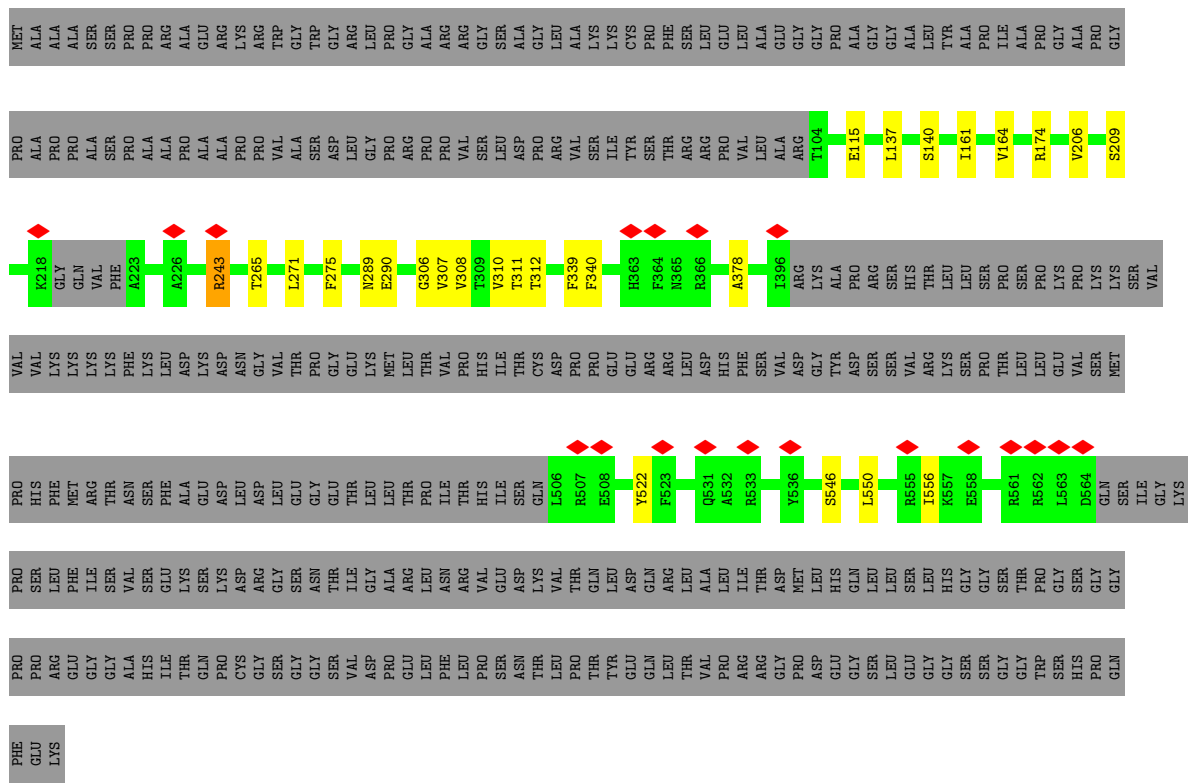


- Molecule 1: Potassium voltage-gated channel subfamily KQT member 1

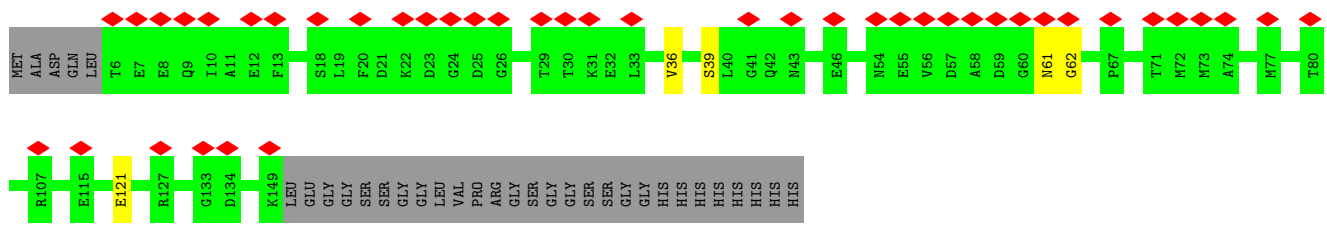
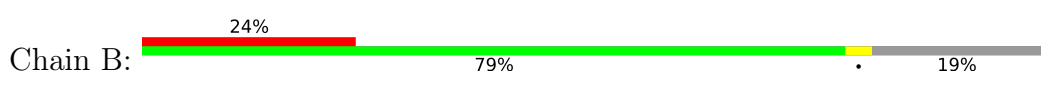


- Molecule 1: Potassium voltage-gated channel subfamily KQT member 1

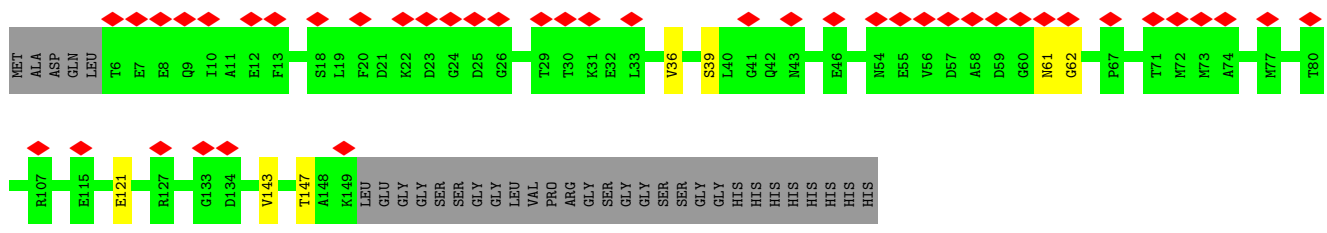
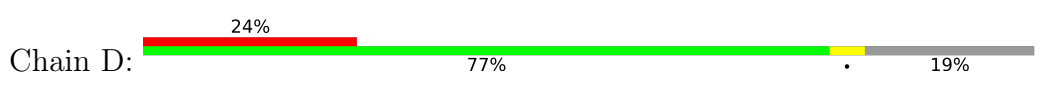




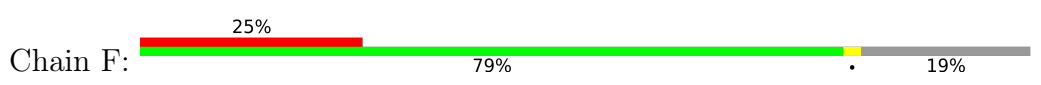
• Molecule 2: Calmodulin-3

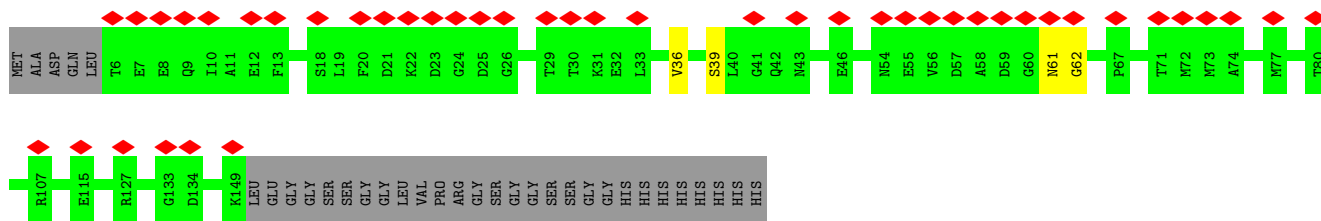


• Molecule 2: Calmodulin-3

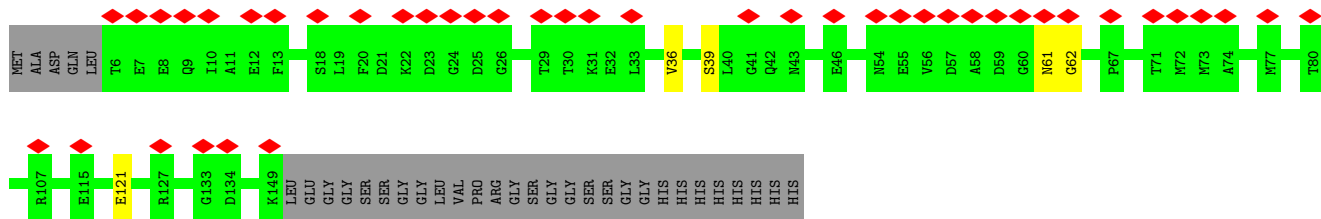
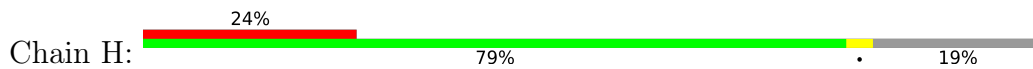


• Molecule 2: Calmodulin-3





• Molecule 2: Calmodulin-3



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	103745	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 ( $e^-/\text{\AA}^2$ )	64	Depositor
Minimum defocus (nm)	-1100	Depositor
Maximum defocus (nm)	-1300	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.076	Depositor
Minimum map value	-0.041	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.0098	Depositor
Map size ( $\text{\AA}$ )	243.36002, 243.36002, 243.36002	wwPDB
Map dimensions	240, 240, 240	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	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: K, I0S, PIO

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	0.69	1/2891 (0.0%)	0.63	1/3913 (0.0%)
1	C	0.69	1/2891 (0.0%)	0.63	1/3913 (0.0%)
1	E	0.69	1/2891 (0.0%)	0.63	1/3913 (0.0%)
1	G	0.69	1/2891 (0.0%)	0.63	1/3913 (0.0%)
2	B	0.38	0/1120	0.55	0/1507
2	D	0.38	0/1120	0.55	0/1507
2	F	0.38	0/1120	0.55	0/1507
2	H	0.38	0/1120	0.55	0/1507
All	All	0.62	4/16044 (0.0%)	0.61	4/21680 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	308	VAL	CB-CG2	-5.35	1.41	1.52
1	A	308	VAL	CB-CG2	-5.32	1.41	1.52
1	G	308	VAL	CB-CG2	-5.30	1.41	1.52
1	C	308	VAL	CB-CG2	-5.27	1.41	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	243	ARG	NE-CZ-NH1	-7.91	116.34	120.30
1	A	243	ARG	NE-CZ-NH1	-7.86	116.37	120.30
1	G	243	ARG	NE-CZ-NH1	-7.86	116.37	120.30
1	C	243	ARG	NE-CZ-NH1	-7.78	116.41	120.30

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	2821	2891	2891	25	0
1	C	2821	2891	2891	26	0
1	E	2821	2891	2891	22	0
1	G	2821	2891	2891	19	0
2	B	1108	1022	1022	3	0
2	D	1108	1022	1022	4	0
2	F	1108	1022	1022	2	0
2	H	1108	1022	1022	3	0
3	A	4	0	0	0	0
4	A	32	25	0	1	0
4	C	32	25	0	1	0
4	E	32	25	0	1	0
4	G	32	25	0	0	0
5	B	47	44	44	1	0
5	D	47	44	44	1	0
5	F	47	44	44	1	0
5	H	47	44	44	1	0
All	All	16036	15928	15828	96	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 (96) 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:E:137:LEU:O	1:E:140:SER:OG	2.08	0.67
1:C:137:LEU:O	1:C:140:SER:OG	2.08	0.66
1:C:275:PHE:CE1	1:E:235:ILE:HG22	2.32	0.65
1:A:137:LEU:O	1:A:140:SER:OG	2.08	0.64
1:C:546:SER:O	1:C:550:LEU:HD23	2.01	0.61
2:D:36:VAL:O	2:D:39:SER:OG	2.15	0.61
2:H:36:VAL:O	2:H:39:SER:OG	2.15	0.61
1:A:546:SER:O	1:A:550:LEU:HD23	2.01	0.61
1:A:235:ILE:HG22	1:G:275:PHE:CE1	2.36	0.60
1:G:137:LEU:O	1:G:140:SER:OG	2.08	0.60

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:546:SER:O	1:E:550:LEU:HD23	2.01	0.59
1:C:206:VAL:O	1:C:209:SER:OG	2.20	0.58
1:G:546:SER:O	1:G:550:LEU:HD23	2.01	0.58
1:E:391:THR:OG1	2:H:121:GLU:OE2	2.17	0.58
1:C:310:VAL:HG21	1:C:340:PHE:CE2	2.39	0.57
2:B:36:VAL:O	2:B:39:SER:OG	2.15	0.57
1:A:206:VAL:O	1:A:209:SER:OG	2.20	0.57
1:G:206:VAL:O	1:G:209:SER:OG	2.20	0.57
2:F:36:VAL:O	2:F:39:SER:OG	2.15	0.57
1:C:564:ASP:OD1	1:E:562:ARG:NH2	2.38	0.56
1:C:115:GLU:OE2	1:C:243:ARG:NH1	2.38	0.56
5:F:601:PIO:O1	5:F:601:PIO:O3	2.21	0.56
1:E:115:GLU:OE2	1:E:243:ARG:NH1	2.38	0.56
1:A:115:GLU:OE2	1:A:243:ARG:NH1	2.39	0.56
1:G:115:GLU:OE2	1:G:243:ARG:NH1	2.39	0.55
1:G:310:VAL:HG21	1:G:340:PHE:CE2	2.41	0.55
1:A:275:PHE:CE1	1:C:235:ILE:HG22	2.41	0.55
1:E:310:VAL:HG21	1:E:340:PHE:CE2	2.42	0.55
1:E:206:VAL:O	1:E:209:SER:OG	2.20	0.55
5:B:601:PIO:O1	5:B:601:PIO:O3	2.21	0.55
1:A:310:VAL:HG21	1:A:340:PHE:CE2	2.42	0.55
1:C:556:ILE:HG21	1:E:555:ARG:HG2	1.89	0.55
5:H:601:PIO:O1	5:H:601:PIO:O3	2.21	0.52
1:C:310:VAL:HG22	1:C:340:PHE:CD2	2.45	0.51
5:D:601:PIO:O1	5:D:601:PIO:O3	2.21	0.51
1:A:310:VAL:HG22	1:A:340:PHE:CD2	2.47	0.50
1:A:555:ARG:HG2	1:G:556:ILE:HG21	1.94	0.50
1:G:310:VAL:HG22	1:G:340:PHE:CD2	2.47	0.49
1:G:311:THR:O	1:G:312:THR:OG1	2.28	0.49
1:E:310:VAL:HG22	1:E:340:PHE:CD2	2.47	0.49
1:C:310:VAL:CG2	1:C:340:PHE:CE2	2.96	0.49
1:A:311:THR:O	1:A:312:THR:OG1	2.28	0.48
1:A:289:ASN:OD1	1:A:290:GLU:N	2.46	0.48
1:C:307:VAL:O	1:C:311:THR:HG22	2.14	0.48
1:E:307:VAL:O	1:E:311:THR:HG22	2.13	0.48
1:A:307:VAL:O	1:A:311:THR:HG22	2.14	0.48
1:G:307:VAL:O	1:G:311:THR:HG22	2.14	0.48
1:A:556:ILE:HG21	1:C:555:ARG:HG2	1.96	0.48
1:E:289:ASN:OD1	1:E:290:GLU:N	2.46	0.48
1:C:391:THR:OG1	2:D:121:GLU:OE2	2.20	0.48
1:C:289:ASN:OD1	1:C:290:GLU:N	2.46	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:THR:OG1	1:A:339:PHE:O	2.32	0.47
1:E:378:ALA:HB2	1:E:522:TYR:HE1	1.80	0.47
1:E:265:THR:OG1	1:E:339:PHE:O	2.32	0.47
1:C:311:THR:O	1:C:312:THR:OG1	2.28	0.47
1:E:336:ALA:HB2	4:E:701:I0S:OAR	2.15	0.47
1:G:378:ALA:HB2	1:G:522:TYR:HE1	1.80	0.47
1:C:336:ALA:HB2	4:C:701:I0S:OAR	2.15	0.47
1:G:265:THR:OG1	1:G:339:PHE:O	2.32	0.47
1:C:378:ALA:HB2	1:C:522:TYR:HE1	1.80	0.46
1:G:310:VAL:CG2	1:G:340:PHE:CE2	2.98	0.46
1:G:289:ASN:OD1	1:G:290:GLU:N	2.46	0.46
1:A:378:ALA:HB2	1:A:522:TYR:HE1	1.80	0.46
1:E:310:VAL:CG2	1:E:340:PHE:CE2	2.99	0.45
2:D:61:ASN:OD1	2:D:62:GLY:N	2.50	0.45
1:A:310:VAL:CG2	1:A:340:PHE:CE2	2.99	0.45
2:B:61:ASN:OD1	2:B:62:GLY:N	2.50	0.45
2:H:61:ASN:OD1	2:H:62:GLY:N	2.50	0.45
1:C:265:THR:OG1	1:C:339:PHE:O	2.32	0.44
1:G:271:LEU:O	1:G:275:PHE:HD2	2.00	0.44
2:F:61:ASN:OD1	2:F:62:GLY:N	2.50	0.43
1:A:271:LEU:O	1:A:275:PHE:HD2	2.02	0.43
1:A:306:GLY:O	1:A:310:VAL:HG12	2.18	0.43
1:A:391:THR:OG1	2:B:121:GLU:OE2	2.23	0.43
1:E:306:GLY:O	1:E:310:VAL:HG12	2.18	0.43
1:A:336:ALA:HB2	4:A:705:I0S:OAR	2.18	0.43
1:C:306:GLY:O	1:C:310:VAL:HG12	2.18	0.43
1:E:311:THR:O	1:E:312:THR:OG1	2.28	0.42
1:G:306:GLY:O	1:G:310:VAL:HG12	2.18	0.42
1:C:378:ALA:HB2	1:C:522:TYR:CE1	2.55	0.42
1:G:161:ILE:HA	1:G:164:VAL:HG12	2.02	0.42
1:A:113:PHE:HD1	1:A:122:CYS:HG	1.63	0.42
1:A:161:ILE:HA	1:A:164:VAL:HG12	2.02	0.42
1:C:174:ARG:HE	1:C:243:ARG:HH12	1.67	0.42
1:A:174:ARG:HE	1:A:243:ARG:HH12	1.68	0.42
1:A:378:ALA:HB2	1:A:522:TYR:CE1	2.55	0.42
1:E:161:ILE:HA	1:E:164:VAL:HG12	2.01	0.42
1:G:378:ALA:HB2	1:G:522:TYR:CE1	2.55	0.42
1:G:174:ARG:HE	1:G:243:ARG:HH12	1.67	0.41
2:D:143:VAL:O	2:D:147:THR:OG1	2.30	0.41
1:E:378:ALA:HB2	1:E:522:TYR:CE1	2.55	0.41
1:C:161:ILE:HA	1:C:164:VAL:HG12	2.02	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:271:LEU:O	1:C:275:PHE:HD2	2.03	0.41
1:E:174:ARG:HE	1:E:243:ARG:HH12	1.68	0.41
1:C:556:ILE:HG21	1:E:555:ARG:CG	2.50	0.40
1:A:564:ASP:OD1	1:C:562:ARG:NH2	2.51	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	342/692 (49%)	330 (96%)	12 (4%)	0	100	100
1	C	342/692 (49%)	331 (97%)	11 (3%)	0	100	100
1	E	342/692 (49%)	330 (96%)	12 (4%)	0	100	100
1	G	342/692 (49%)	330 (96%)	12 (4%)	0	100	100
2	B	142/177 (80%)	135 (95%)	7 (5%)	0	100	100
2	D	142/177 (80%)	135 (95%)	7 (5%)	0	100	100
2	F	142/177 (80%)	135 (95%)	7 (5%)	0	100	100
2	H	142/177 (80%)	135 (95%)	7 (5%)	0	100	100
All	All	1936/3476 (56%)	1861 (96%)	75 (4%)	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	297/579 (51%)	297 (100%)	0	100	100
1	C	297/579 (51%)	297 (100%)	0	100	100
1	E	297/579 (51%)	297 (100%)	0	100	100
1	G	297/579 (51%)	297 (100%)	0	100	100
2	B	115/146 (79%)	115 (100%)	0	100	100
2	D	115/146 (79%)	115 (100%)	0	100	100
2	F	115/146 (79%)	115 (100%)	0	100	100
2	H	115/146 (79%)	115 (100%)	0	100	100
All	All	1648/2900 (57%)	1648 (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. 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 [i](#)

Of 12 ligands modelled in this entry, 4 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
4	I0S	C	701	-	32,35,35	2.26	8 (25%)	43,50,50	2.01	7 (16%)
5	PIO	B	601	-	47,47,47	1.17	6 (12%)	61,65,65	0.83	2 (3%)
5	PIO	F	601	-	47,47,47	1.17	6 (12%)	61,65,65	0.83	2 (3%)
5	PIO	H	601	-	47,47,47	1.17	6 (12%)	61,65,65	0.83	2 (3%)
4	I0S	E	701	-	32,35,35	2.26	7 (21%)	43,50,50	2.01	7 (16%)
5	PIO	D	601	-	47,47,47	1.16	6 (12%)	61,65,65	0.83	2 (3%)
4	I0S	G	701	-	32,35,35	2.26	7 (21%)	43,50,50	2.01	7 (16%)
4	I0S	A	705	-	32,35,35	2.26	7 (21%)	43,50,50	2.01	7 (16%)

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
4	I0S	C	701	-	-	1/24/37/37	0/4/4/4
5	PIO	B	601	-	-	15/44/68/68	0/1/1/1
5	PIO	F	601	-	-	15/44/68/68	0/1/1/1
5	PIO	H	601	-	-	15/44/68/68	0/1/1/1
4	I0S	E	701	-	-	1/24/37/37	0/4/4/4
5	PIO	D	601	-	-	15/44/68/68	0/1/1/1
4	I0S	G	701	-	-	1/24/37/37	0/4/4/4
4	I0S	A	705	-	-	1/24/37/37	0/4/4/4

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	701	I0S	CBE-SBF	-7.45	1.58	1.70
4	G	701	I0S	CBE-SBF	-7.40	1.58	1.70
4	C	701	I0S	CBE-SBF	-7.40	1.58	1.70
4	A	705	I0S	CBE-SBF	-7.37	1.59	1.70
4	A	705	I0S	SAI-N	4.02	1.69	1.63
4	G	701	I0S	CAE-CB	-4.00	1.42	1.53
4	A	705	I0S	CAE-CB	-3.99	1.42	1.53
4	C	701	I0S	CAE-CB	-3.98	1.42	1.53
4	E	701	I0S	CAE-CB	-3.98	1.42	1.53
4	C	701	I0S	SAI-N	3.98	1.69	1.63

Continued on next page...

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	701	I0S	SAI-N	3.97	1.69	1.63
4	E	701	I0S	SAI-N	3.96	1.69	1.63
4	A	705	I0S	CAT-NAS	3.34	1.42	1.36
4	C	701	I0S	CAT-NAS	3.31	1.42	1.36
4	E	701	I0S	CAT-NAS	3.29	1.42	1.36
4	G	701	I0S	CAT-NAS	3.28	1.42	1.36
4	A	705	I0S	O-C	-3.28	1.16	1.23
4	G	701	I0S	O-C	-3.26	1.16	1.23
4	E	701	I0S	O-C	-3.25	1.16	1.23
4	C	701	I0S	O-C	-3.24	1.16	1.23
5	B	601	PIO	P5-O5	3.15	1.65	1.59
5	F	601	PIO	P5-O5	3.15	1.65	1.59
5	D	601	PIO	P5-O5	3.14	1.65	1.59
5	H	601	PIO	P5-O5	3.13	1.65	1.59
4	C	701	I0S	C-NAS	3.07	1.42	1.35
4	A	705	I0S	C-NAS	3.06	1.42	1.35
4	G	701	I0S	C-NAS	3.05	1.42	1.35
4	E	701	I0S	C-NAS	3.04	1.42	1.35
5	F	601	PIO	P4-O4	2.93	1.64	1.59
5	B	601	PIO	P4-O4	2.89	1.64	1.59
5	H	601	PIO	P4-O4	2.88	1.64	1.59
5	D	601	PIO	P4-O4	2.85	1.64	1.59
5	D	601	PIO	O2C-C2C	-2.61	1.40	1.46
5	B	601	PIO	O2C-C2C	-2.60	1.40	1.46
5	H	601	PIO	O2C-C2C	-2.56	1.40	1.46
5	F	601	PIO	O2C-C2C	-2.55	1.40	1.46
5	F	601	PIO	O3C-C3C	-2.49	1.39	1.45
5	H	601	PIO	O3C-C3C	-2.46	1.39	1.45
5	B	601	PIO	O3C-C3C	-2.45	1.39	1.45
5	D	601	PIO	O3C-C3C	-2.44	1.39	1.45
4	C	701	I0S	CAF-CAE	-2.29	1.42	1.51
4	A	705	I0S	CAF-CAE	-2.29	1.42	1.51
4	G	701	I0S	CAF-CAE	-2.27	1.42	1.51
4	E	701	I0S	CAF-CAE	-2.27	1.42	1.51
5	B	601	PIO	O3C-C1B	2.25	1.39	1.33
5	F	601	PIO	O3C-C1B	2.25	1.39	1.33
5	H	601	PIO	O3C-C1B	2.24	1.39	1.33
5	D	601	PIO	O3C-C1B	2.21	1.39	1.33
4	C	701	I0S	CAJ-SAI	2.04	1.79	1.76
5	B	601	PIO	O2C-C1A	2.03	1.40	1.34
5	F	601	PIO	O2C-C1A	2.03	1.40	1.34
5	D	601	PIO	O2C-C1A	2.02	1.40	1.34

*Continued on next page...*

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	601	PIO	O2C-C1A	2.02	1.40	1.34

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	701	I0S	OAR-SAI-OAQ	-9.82	103.60	119.52
4	C	701	I0S	OAR-SAI-OAQ	-9.82	103.61	119.52
4	A	705	I0S	OAR-SAI-OAQ	-9.81	103.62	119.52
4	G	701	I0S	OAR-SAI-OAQ	-9.81	103.62	119.52
5	B	601	PIO	O2C-C1A-C2A	3.21	118.42	111.50
5	D	601	PIO	O2C-C1A-C2A	3.21	118.41	111.50
5	F	601	PIO	O2C-C1A-C2A	3.20	118.40	111.50
5	H	601	PIO	O2C-C1A-C2A	3.20	118.39	111.50
4	G	701	I0S	OAQ-SAI-CAJ	2.99	111.83	108.05
4	A	705	I0S	OAQ-SAI-CAJ	2.99	111.83	108.05
4	C	701	I0S	OAQ-SAI-CAJ	2.97	111.81	108.05
4	E	701	I0S	OAQ-SAI-CAJ	2.96	111.79	108.05
4	C	701	I0S	OAQ-SAI-N	2.84	112.15	106.97
4	E	701	I0S	OAQ-SAI-N	2.84	112.14	106.97
4	G	701	I0S	OAQ-SAI-N	2.83	112.14	106.97
4	A	705	I0S	OAQ-SAI-N	2.81	112.10	106.97
5	D	601	PIO	O3C-C1B-C2B	2.79	120.66	111.91
5	B	601	PIO	O3C-C1B-C2B	2.78	120.62	111.91
5	F	601	PIO	O3C-C1B-C2B	2.77	120.61	111.91
5	H	601	PIO	O3C-C1B-C2B	2.77	120.61	111.91
4	G	701	I0S	CAG-N-CA	-2.69	107.76	114.64
4	A	705	I0S	CAG-N-CA	-2.69	107.77	114.64
4	E	701	I0S	CAG-N-CA	-2.68	107.80	114.64
4	C	701	I0S	CAG-N-CA	-2.67	107.81	114.64
4	E	701	I0S	CB-CA-C	-2.53	107.52	112.12
4	C	701	I0S	CB-CA-C	-2.53	107.53	112.12
4	A	705	I0S	CAF-CAG-N	2.53	114.01	110.31
4	G	701	I0S	CAF-CAG-N	2.52	114.01	110.31
4	G	701	I0S	CB-CA-C	-2.52	107.54	112.12
4	A	705	I0S	CB-CA-C	-2.52	107.55	112.12
4	E	701	I0S	CAF-CAG-N	2.51	113.99	110.31
4	C	701	I0S	CAF-CAG-N	2.50	113.98	110.31
4	A	705	I0S	CAT-NAS-C	-2.19	124.15	129.02
4	C	701	I0S	CAT-NAS-C	-2.15	124.23	129.02
4	G	701	I0S	CAT-NAS-C	-2.15	124.23	129.02
4	E	701	I0S	CAT-NAS-C	-2.14	124.26	129.02

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	601	PIO	C2A-C1A-O2C-C2C
5	D	601	PIO	C2A-C1A-O2C-C2C
5	F	601	PIO	C2A-C1A-O2C-C2C
5	H	601	PIO	C2A-C1A-O2C-C2C
5	B	601	PIO	O1A-C1A-O2C-C2C
5	D	601	PIO	O1A-C1A-O2C-C2C
5	F	601	PIO	O1A-C1A-O2C-C2C
5	H	601	PIO	O1A-C1A-O2C-C2C
5	B	601	PIO	C2B-C1B-O3C-C3C
5	D	601	PIO	C2B-C1B-O3C-C3C
5	F	601	PIO	C2B-C1B-O3C-C3C
5	H	601	PIO	C2B-C1B-O3C-C3C
5	B	601	PIO	O1B-C1B-O3C-C3C
5	D	601	PIO	O1B-C1B-O3C-C3C
5	F	601	PIO	O1B-C1B-O3C-C3C
5	H	601	PIO	O1B-C1B-O3C-C3C
5	B	601	PIO	C1C-O13-P1-O1
5	D	601	PIO	C1C-O13-P1-O1
5	F	601	PIO	C1C-O13-P1-O1
5	H	601	PIO	C1C-O13-P1-O1
5	B	601	PIO	O13-C1C-C2C-C3C
5	D	601	PIO	O13-C1C-C2C-C3C
5	F	601	PIO	O13-C1C-C2C-C3C
5	H	601	PIO	O13-C1C-C2C-C3C
5	B	601	PIO	C5-O5-P5-O53
5	D	601	PIO	C5-O5-P5-O53
5	F	601	PIO	C5-O5-P5-O53
5	H	601	PIO	C5-O5-P5-O53
5	B	601	PIO	O13-C1C-C2C-O2C
5	D	601	PIO	O13-C1C-C2C-O2C
5	F	601	PIO	O13-C1C-C2C-O2C
5	H	601	PIO	O13-C1C-C2C-O2C
5	B	601	PIO	C4-O4-P4-O43
5	D	601	PIO	C4-O4-P4-O43
5	F	601	PIO	C4-O4-P4-O43
5	H	601	PIO	C4-O4-P4-O43
5	B	601	PIO	C1C-O13-P1-O12
5	D	601	PIO	C1C-O13-P1-O12
5	F	601	PIO	C1C-O13-P1-O12
5	H	601	PIO	C1C-O13-P1-O12

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
5	B	601	PIO	C2C-C1C-O13-P1
5	D	601	PIO	C2C-C1C-O13-P1
5	F	601	PIO	C2C-C1C-O13-P1
5	H	601	PIO	C2C-C1C-O13-P1
5	B	601	PIO	C5B-C6B-C7B-C8B
5	D	601	PIO	C5B-C6B-C7B-C8B
5	H	601	PIO	C5B-C6B-C7B-C8B
5	F	601	PIO	C5B-C6B-C7B-C8B
5	B	601	PIO	C5-C4-O4-P4
5	D	601	PIO	C5-C4-O4-P4
5	F	601	PIO	C5-C4-O4-P4
5	H	601	PIO	C5-C4-O4-P4
4	A	705	I0S	CA-N-SAI-OAQ
4	C	701	I0S	CA-N-SAI-OAQ
4	E	701	I0S	CA-N-SAI-OAQ
4	G	701	I0S	CA-N-SAI-OAQ
5	B	601	PIO	C4-O4-P4-O41
5	B	601	PIO	C5-O5-P5-O51
5	D	601	PIO	C4-O4-P4-O41
5	D	601	PIO	C5-O5-P5-O51
5	F	601	PIO	C4-O4-P4-O41
5	F	601	PIO	C5-O5-P5-O51
5	H	601	PIO	C4-O4-P4-O41
5	H	601	PIO	C5-O5-P5-O51

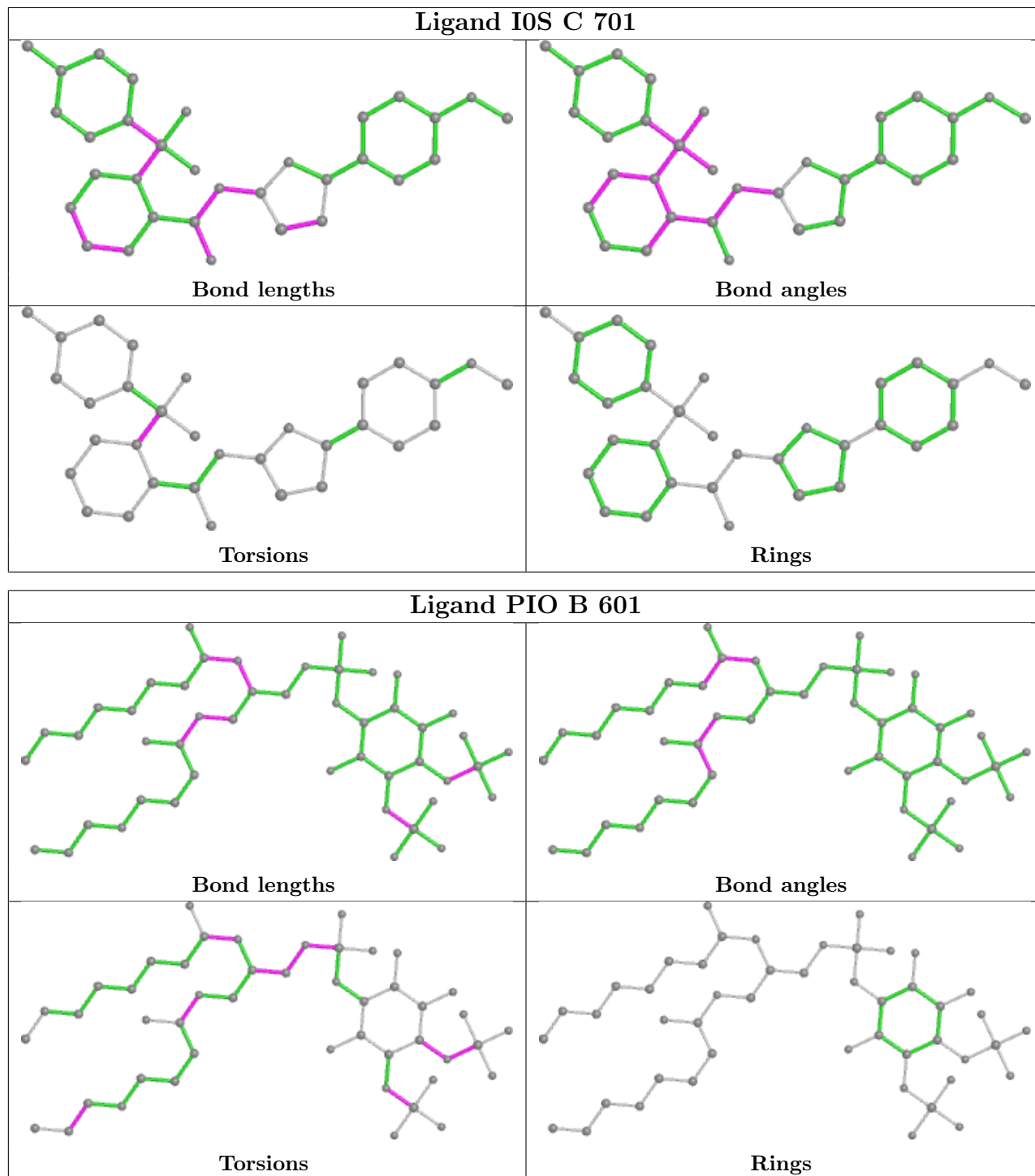
There are no ring outliers.

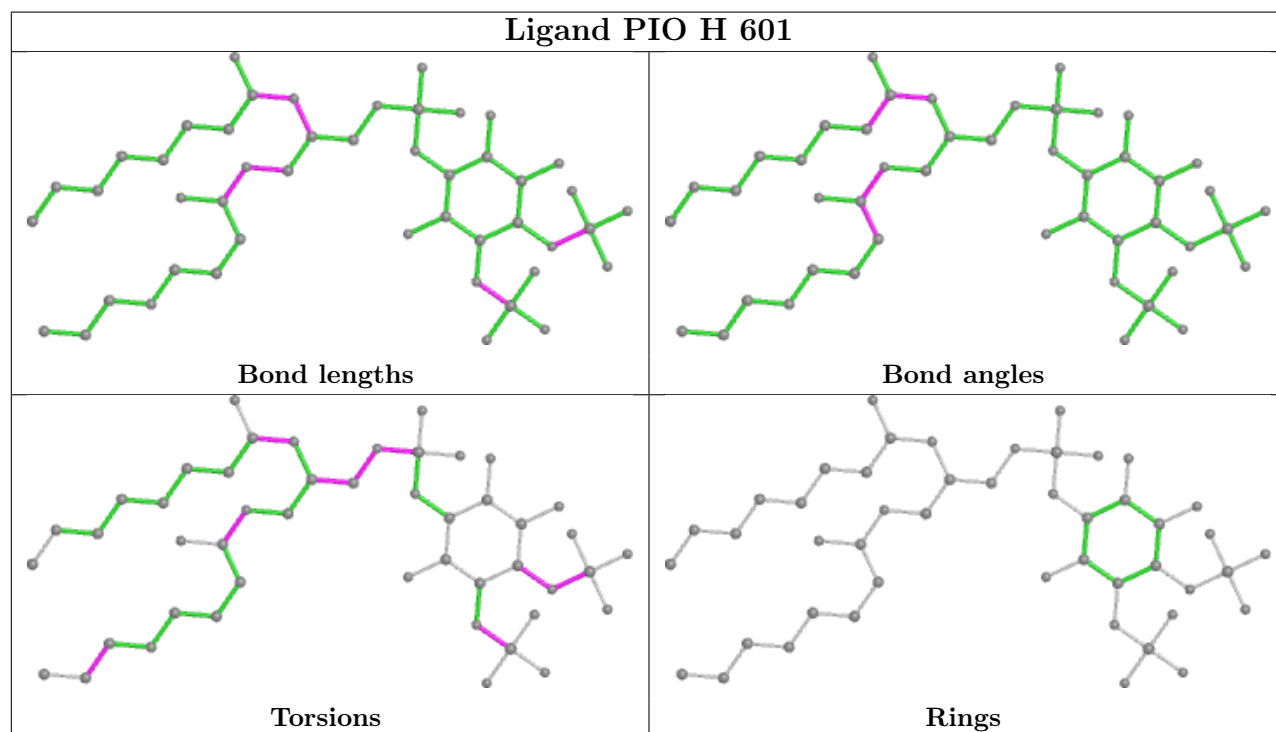
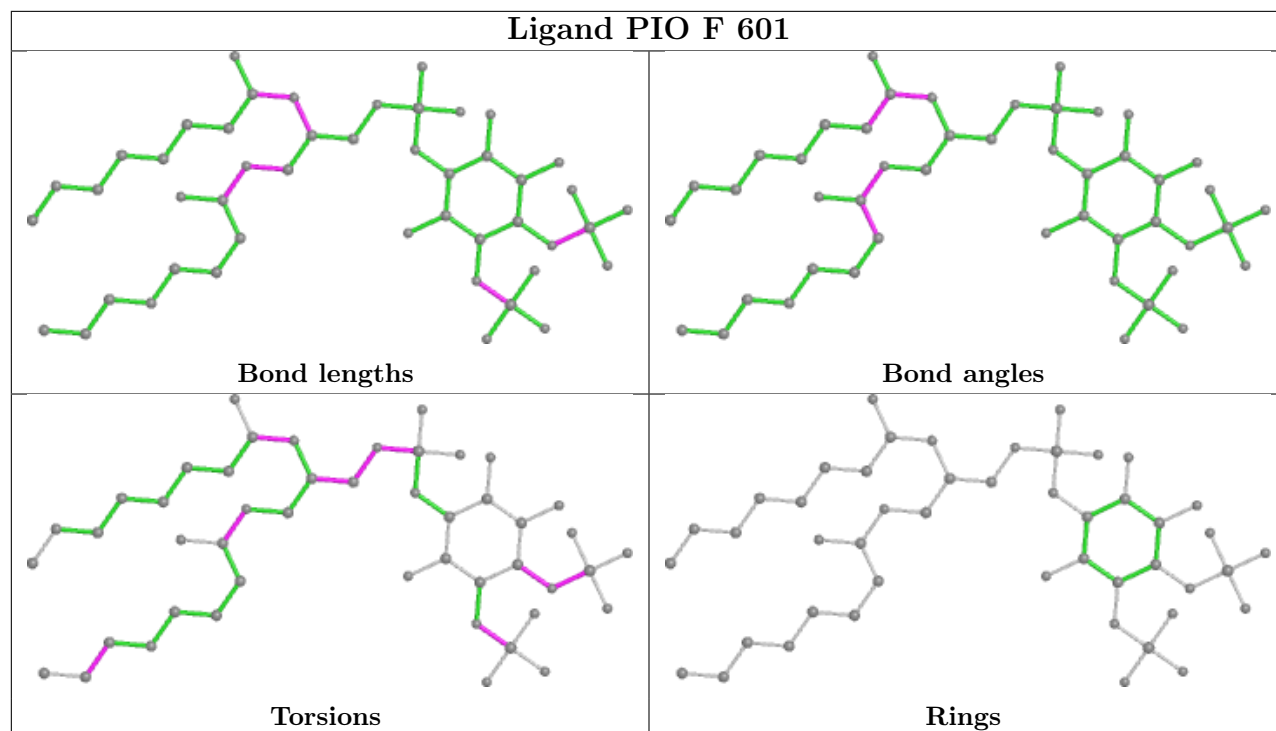
7 monomers are involved in 7 short contacts:

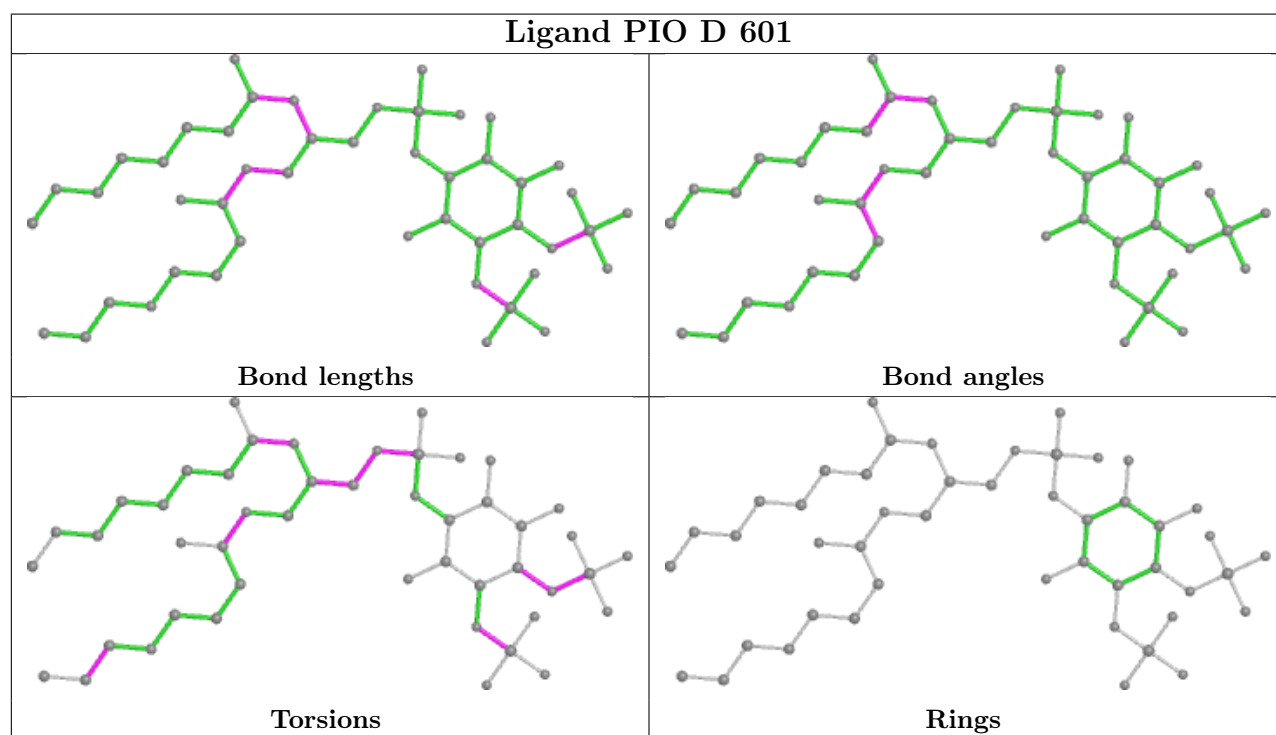
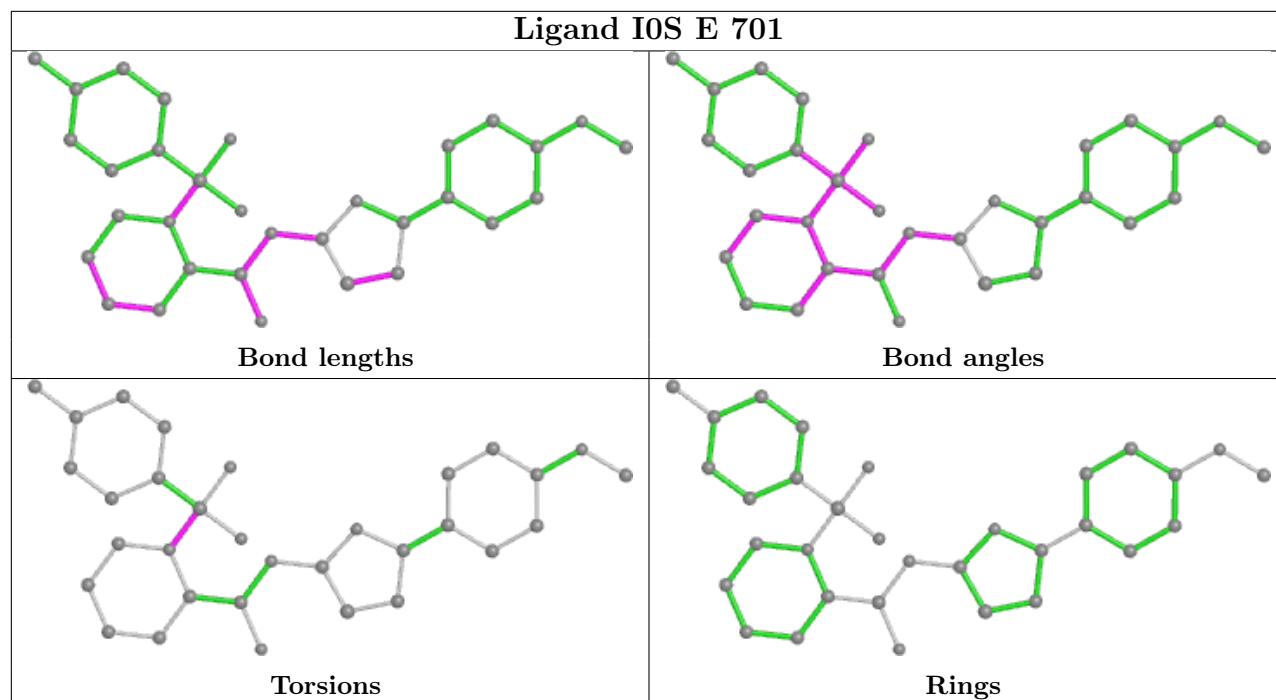
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	701	I0S	1	0
5	B	601	PIO	1	0
5	F	601	PIO	1	0
5	H	601	PIO	1	0
4	E	701	I0S	1	0
5	D	601	PIO	1	0
4	A	705	I0S	1	0

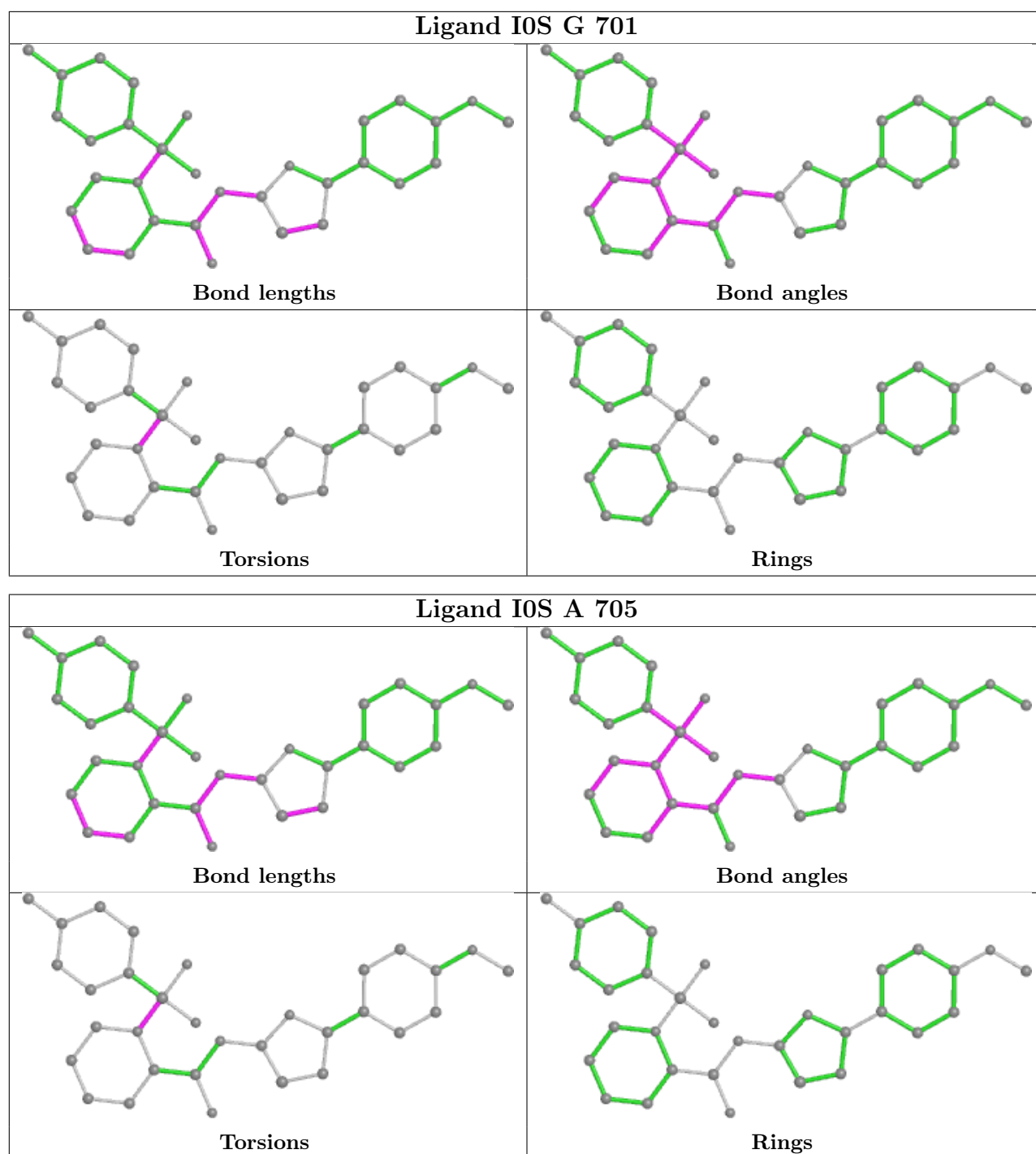
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

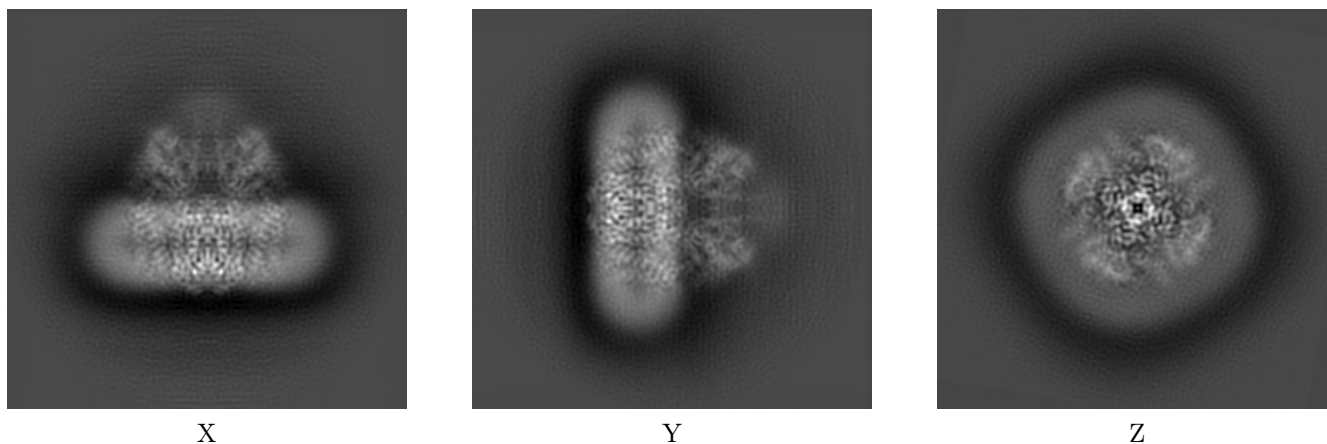
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-33318. These allow visual inspection of the internal detail of the map and identification of artifacts.

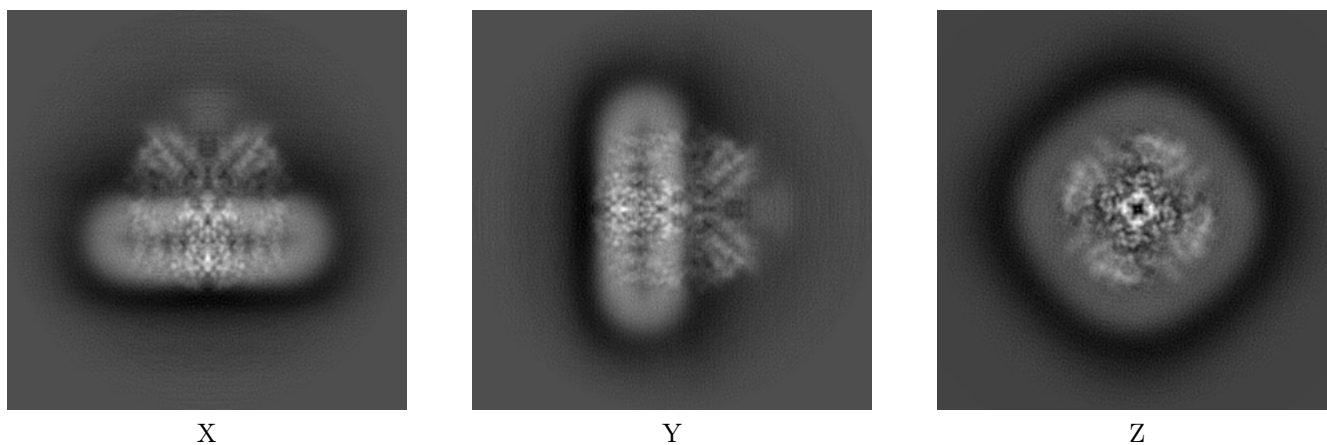
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

#### 6.1.1 Primary map



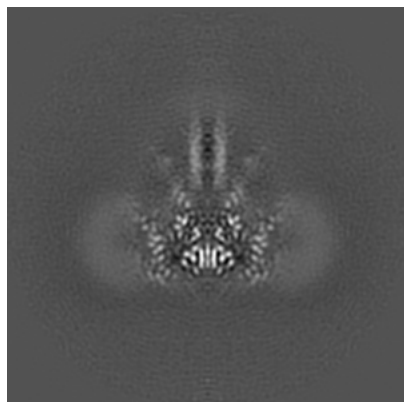
#### 6.1.2 Raw map



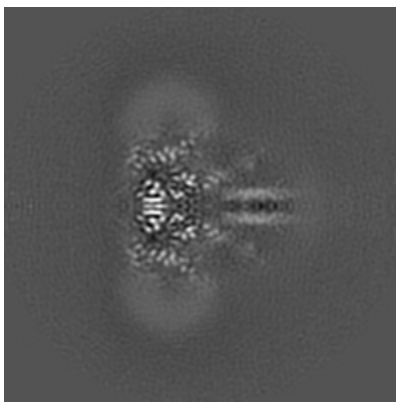
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

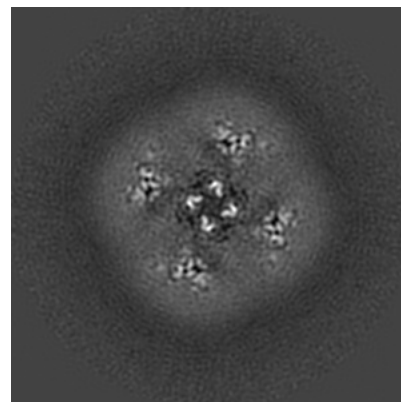
### 6.2.1 Primary map



X Index: 120

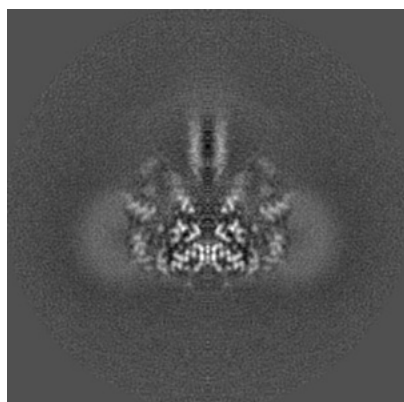


Y Index: 120

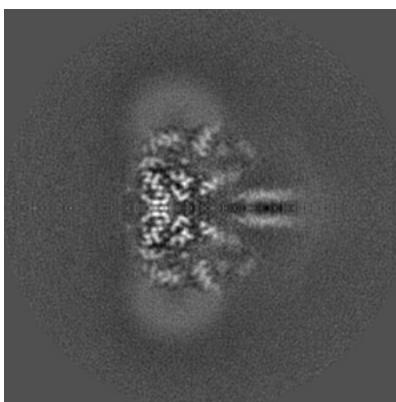


Z Index: 120

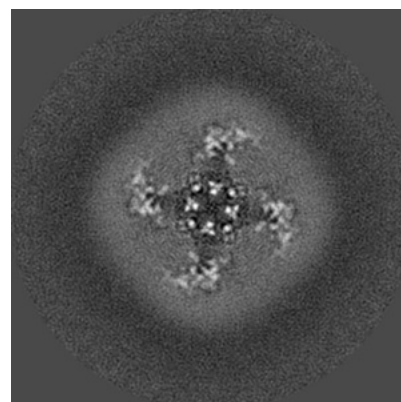
### 6.2.2 Raw map



X Index: 120



Y Index: 120

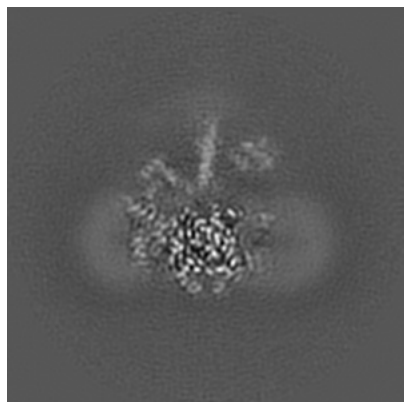


Z Index: 120

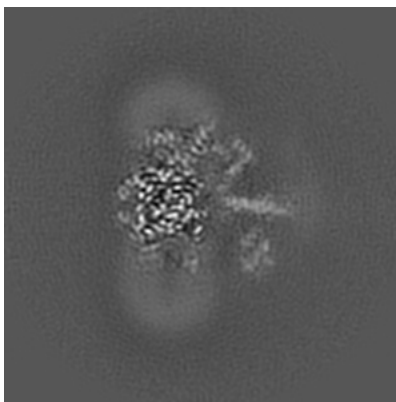
The images above show central slices of the map in three orthogonal directions.

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

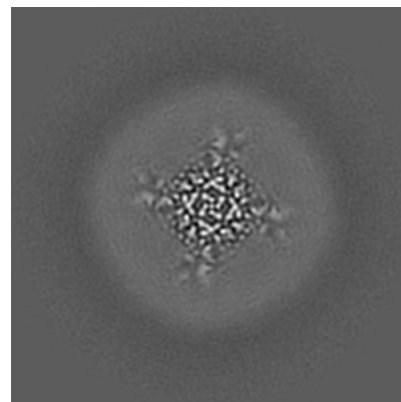
### 6.3.1 Primary map



X Index: 113

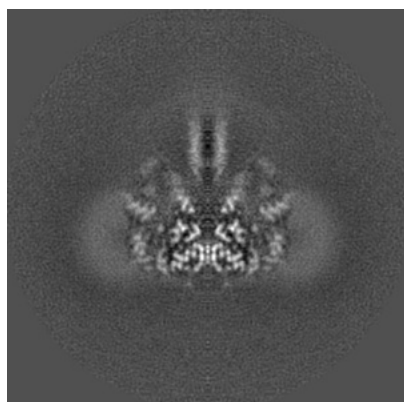


Y Index: 113

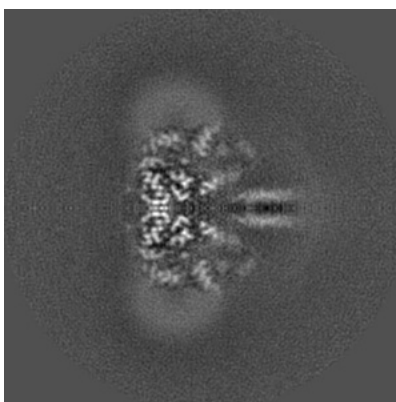


Z Index: 86

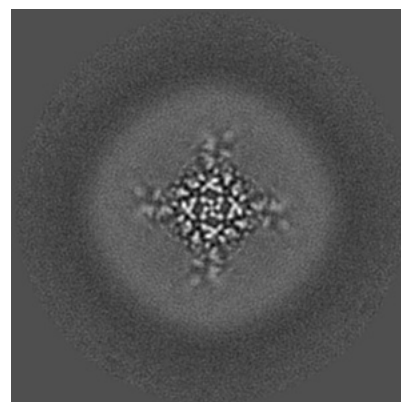
### 6.3.2 Raw map



X Index: 120



Y Index: 120



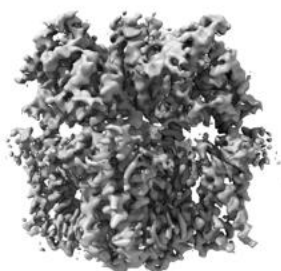
Z Index: 89

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

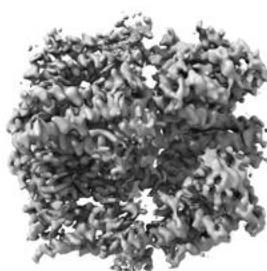


## 6.4 Orthogonal surface views [i](#)

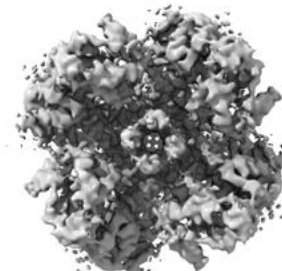
### 6.4.1 Primary map



X



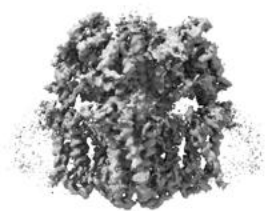
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0098. 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



X



Y



Z

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.

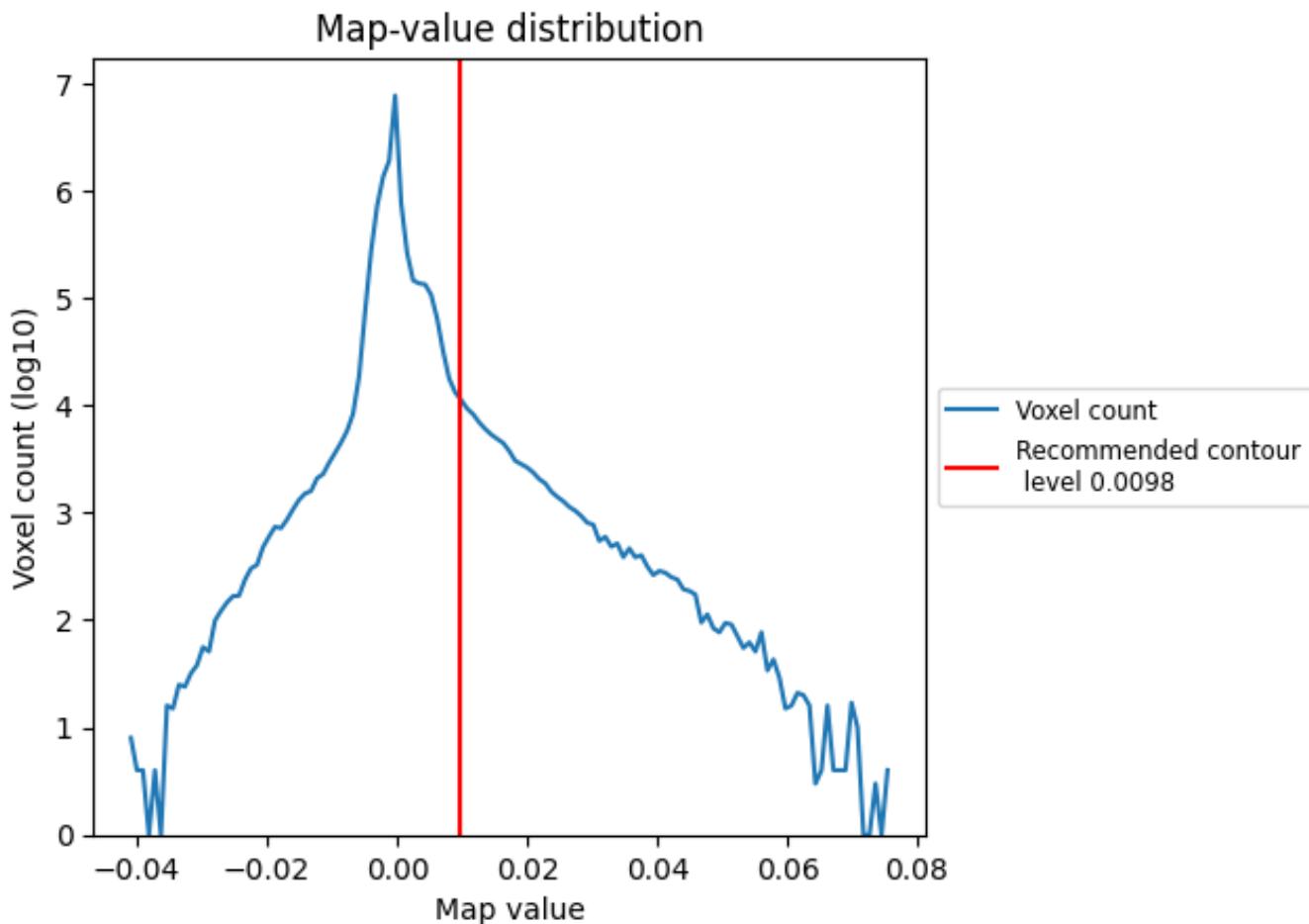
## 6.5 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

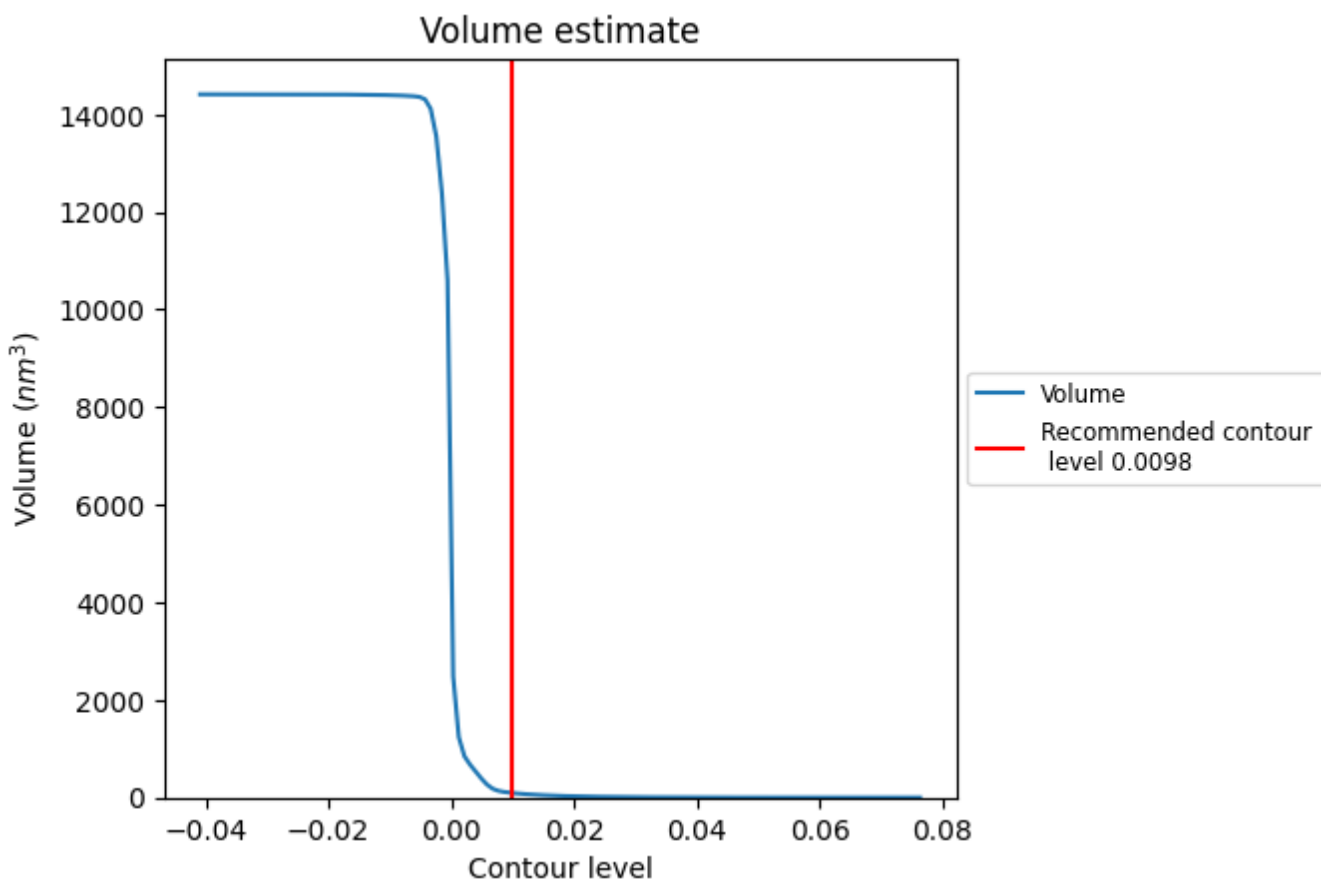
This section contains the results of statistical analysis of the map.

### 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.

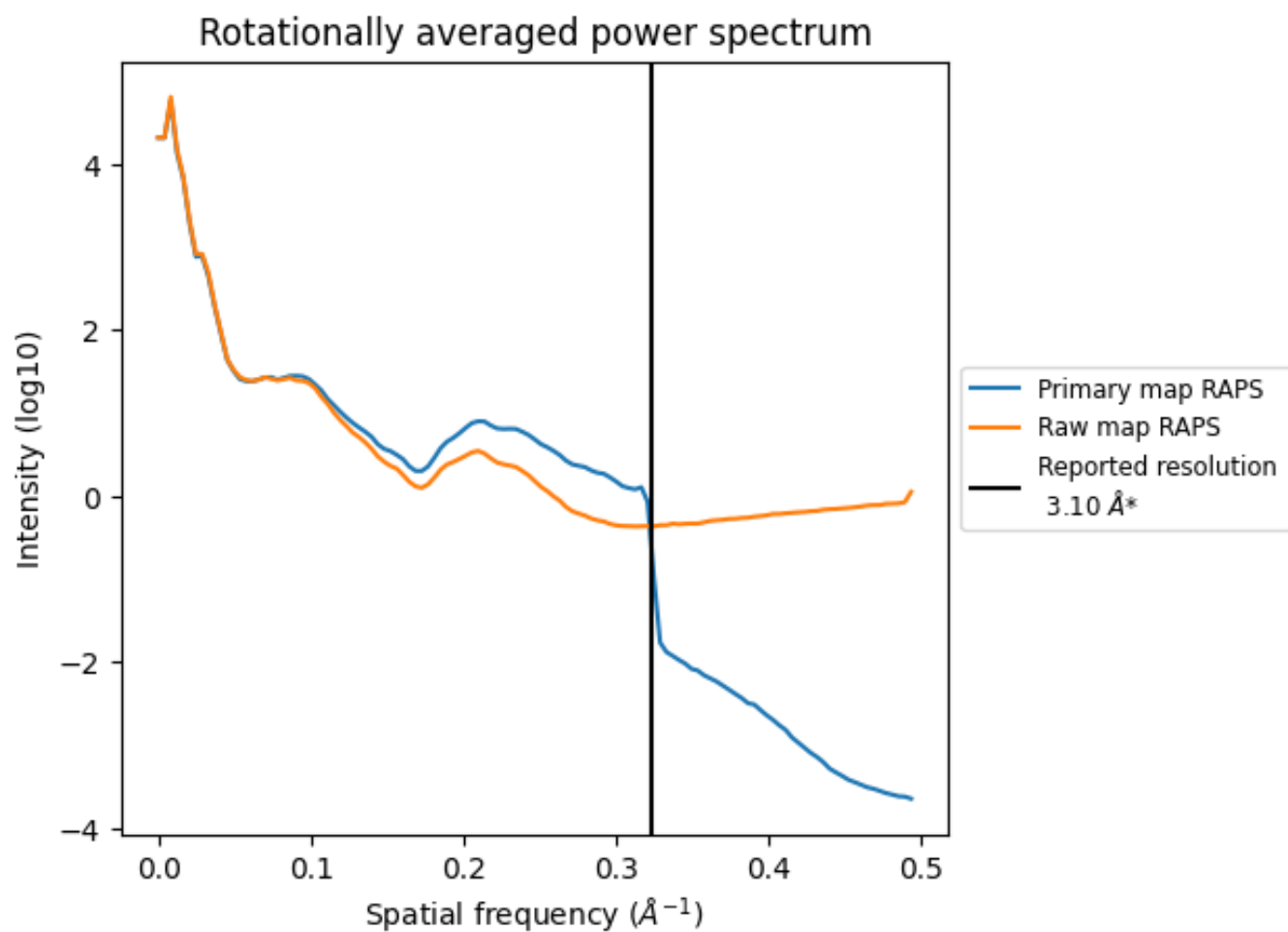
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 96 nm<sup>3</sup>; this corresponds to an approximate mass of 87 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

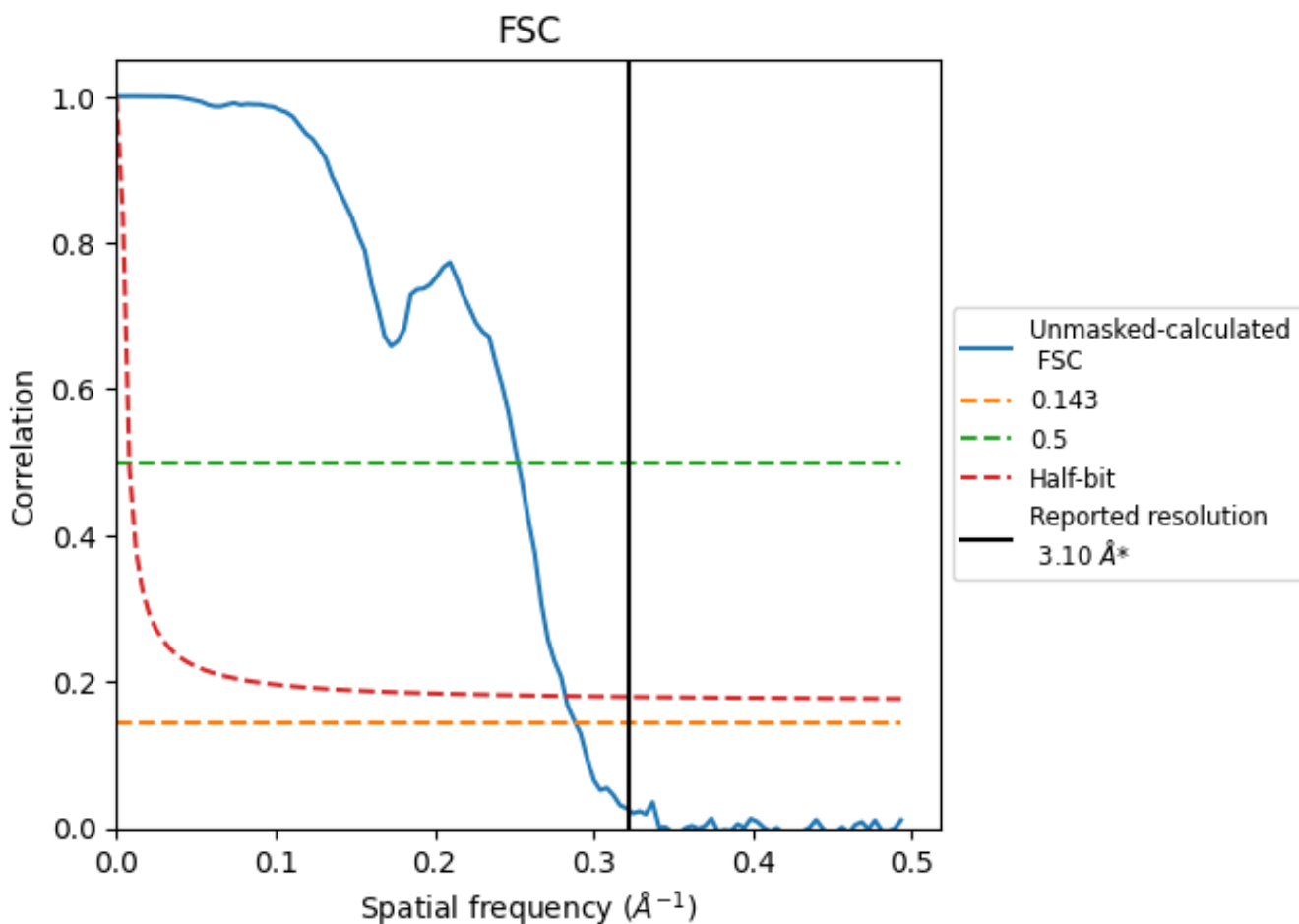


\*Reported resolution corresponds to spatial frequency of  $0.323 \text{ \AA}^{-1}$

## 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 0.323 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

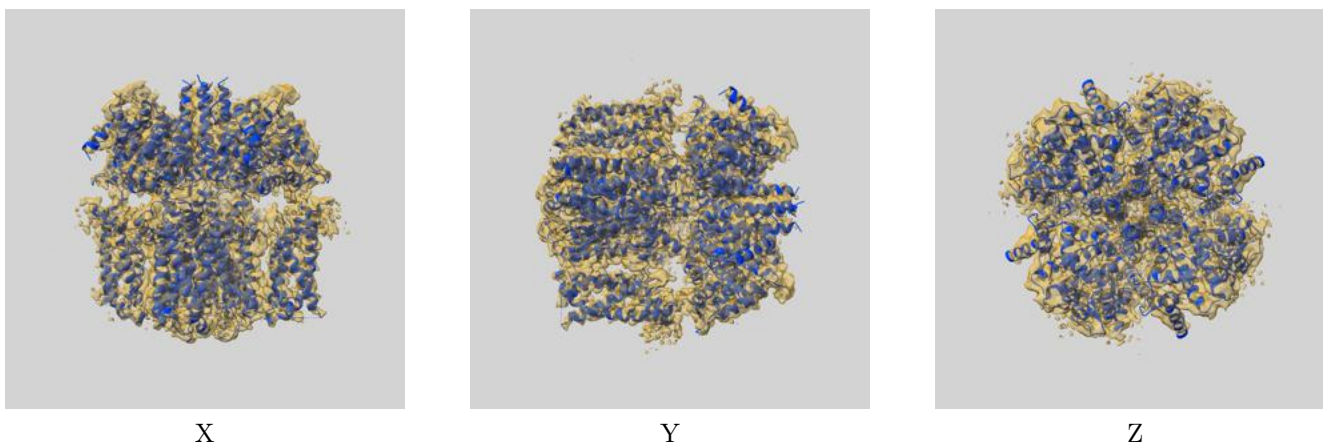
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.47	3.96	3.54

\*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 3.47 differs from the reported value 3.1 by more than 10 %

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

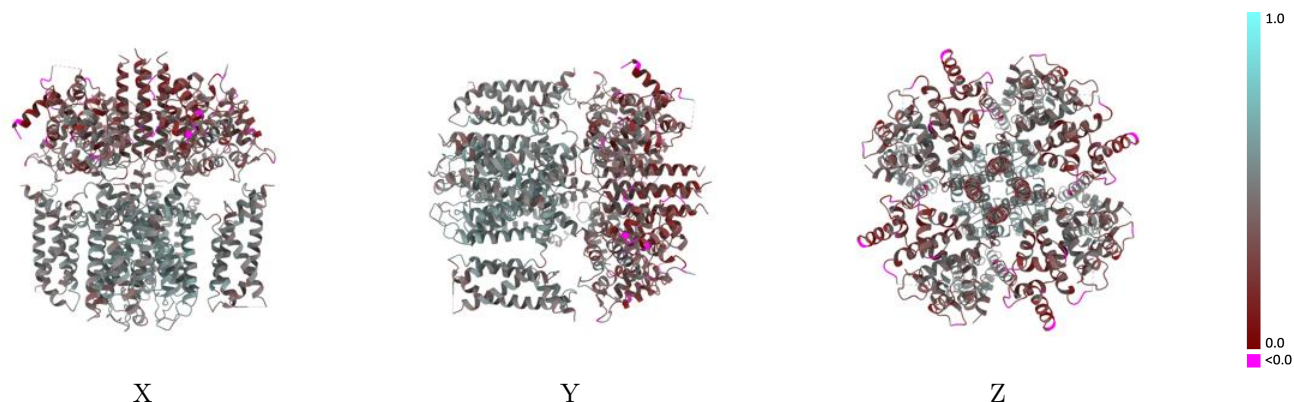
This section contains information regarding the fit between EMDB map EMD-33318 and PDB model 7XNL. Per-residue inclusion information can be found in section 3 on page 10.

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



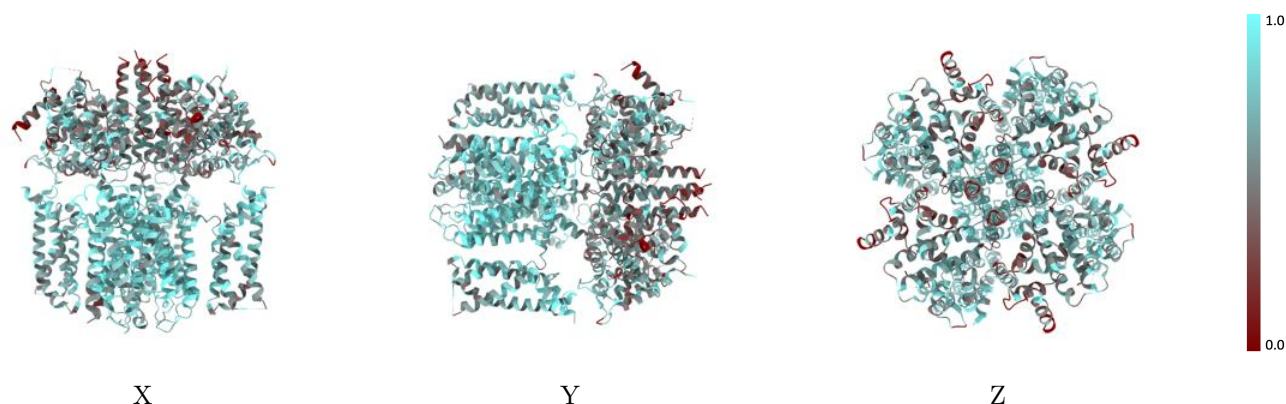
The images above show the 3D surface view of the map at the recommended contour level 0.0098 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

## 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.

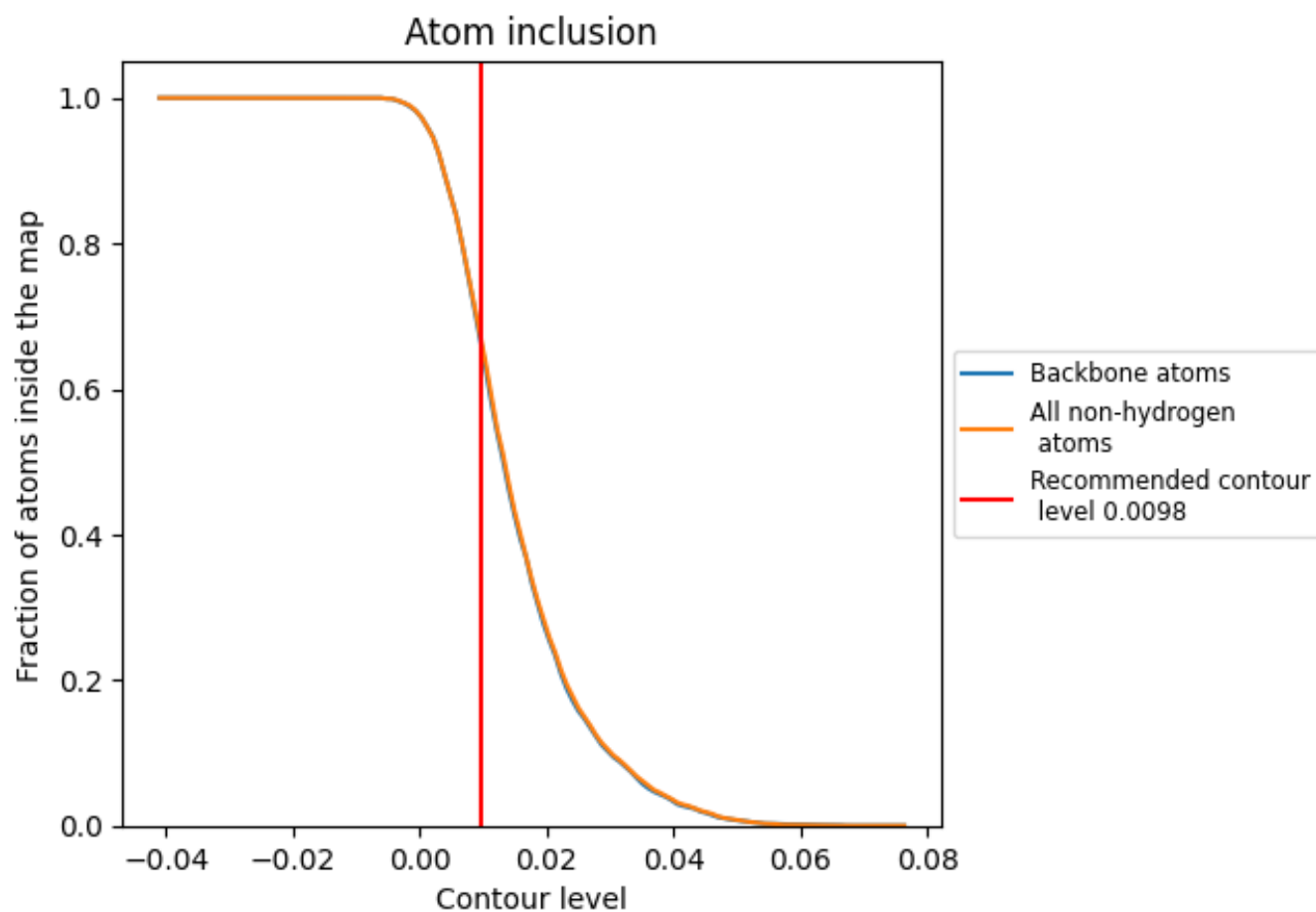
## 9.3 Atom inclusion mapped to coordinate model [i](#)



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 (0.0098).





















## 9.4 Atom inclusion [i](#)



At the recommended contour level, 66% of all backbone atoms, 66% 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 (0.0098) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6633	 0.4120
A	 0.7399	 0.4560
B	 0.5223	 0.3060
C	 0.7396	 0.4570
D	 0.5223	 0.3080
E	 0.7385	 0.4550
F	 0.5223	 0.3030
G	 0.7381	 0.4540
H	 0.5215	 0.3050

