



wwPDB EM Validation Summary Report ⓘ

Nov 19, 2022 – 04:17 pm GMT

PDB ID : 5NG5
EMDB ID : EMD-3636
Title : multi-drug efflux; membrane transport; RND superfamily; Drug resistance
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Chiu, W.; Luisi, B.F.; Du, D.
Deposited on : 2017-03-16
Resolution : 6.50 Å(reported)

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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 : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

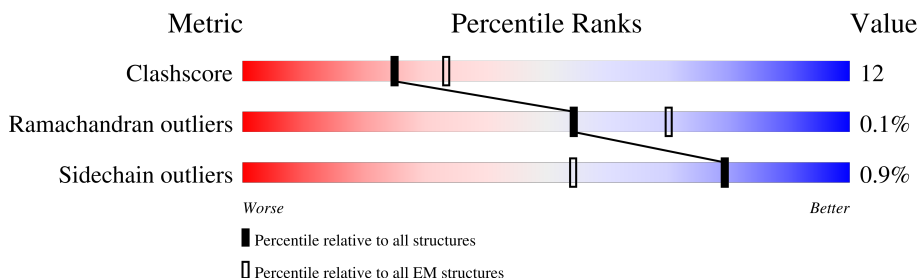
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 6.50 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	373	 69% 21% 9%
1	B	373	 69% 21% 9%
1	D	373	 5% 60% 28% 9%
1	E	373	 6% 63% 25% 9%
1	G	373	 5% 61% 28% 9%
1	H	373	 67% 23% 9%
2	C	493	 66% 21% 13%
2	F	493	 66% 21% 13%

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Mol	Chain	Length	Quality of chain
2	I	493	
3	J	1049	
3	K	1049	
3	L	1049	
4	M	54	
4	N	54	
4	O	54	

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	5QF	J	1101	-	-	X	-
5	5QF	K	1101	-	-	X	-
5	5QF	L	1101	-	-	X	-

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 49920 atoms, of which 114 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 Multidrug efflux pump subunit AcrA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	E	340	2553	1591	451	506	5	0	0
1	D	340	2553	1591	451	506	5	0	0
1	G	340	2553	1591	451	506	5	0	0
1	H	340	2553	1591	451	506	5	0	0
1	A	340	2553	1591	451	506	5	0	0
1	B	340	2553	1591	451	506	5	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	223	MET	PHE	conflict	UNP P0AE06
E	224	MET	LEU	conflict	UNP P0AE06
E	287	MET	LEU	conflict	UNP P0AE06
E	288	MET	LEU	conflict	UNP P0AE06
D	223	MET	PHE	conflict	UNP P0AE06
D	224	MET	LEU	conflict	UNP P0AE06
D	287	MET	LEU	conflict	UNP P0AE06
D	288	MET	LEU	conflict	UNP P0AE06
G	223	MET	PHE	conflict	UNP P0AE06
G	224	MET	LEU	conflict	UNP P0AE06
G	287	MET	LEU	conflict	UNP P0AE06
G	288	MET	LEU	conflict	UNP P0AE06
H	223	MET	PHE	conflict	UNP P0AE06
H	224	MET	LEU	conflict	UNP P0AE06
H	287	MET	LEU	conflict	UNP P0AE06
H	288	MET	LEU	conflict	UNP P0AE06
A	223	MET	PHE	conflict	UNP P0AE06
A	224	MET	LEU	conflict	UNP P0AE06

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Chain	Residue	Modelled	Actual	Comment	Reference
A	287	MET	LEU	conflict	UNP P0AE06
A	288	MET	LEU	conflict	UNP P0AE06
B	223	MET	PHE	conflict	UNP P0AE06
B	224	MET	LEU	conflict	UNP P0AE06
B	287	MET	LEU	conflict	UNP P0AE06
B	288	MET	LEU	conflict	UNP P0AE06

- Molecule 2 is a protein called Outer membrane protein TolC.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	F	428	Total	C	N	O	S	0	0
			3304	2037	586	676	5		
2	I	428	Total	C	N	O	S	0	0
			3304	2037	586	676	5		
2	C	428	Total	C	N	O	S	0	0
			3304	2037	586	676	5		

- Molecule 3 is a protein called Multidrug efflux pump subunit AcrB.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	K	1037	Total	C	N	O	S	0	0
			7819	5032	1290	1453	44		
3	L	1037	Total	C	N	O	S	0	0
			7819	5032	1290	1453	44		
3	J	1033	Total	C	N	O	S	0	0
			7849	5052	1295	1458	44		

- Molecule 4 is a protein called Multidrug efflux pump accessory protein AcrZ.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	M	46	Total	C	N	O	S	0	0
			326	219	48	56	3		
4	N	46	Total	C	N	O	S	0	0
			326	219	48	56	3		
4	O	46	Total	C	N	O	S	0	0
			326	219	48	56	3		

There are 15 discrepancies between the modelled and reference sequences:

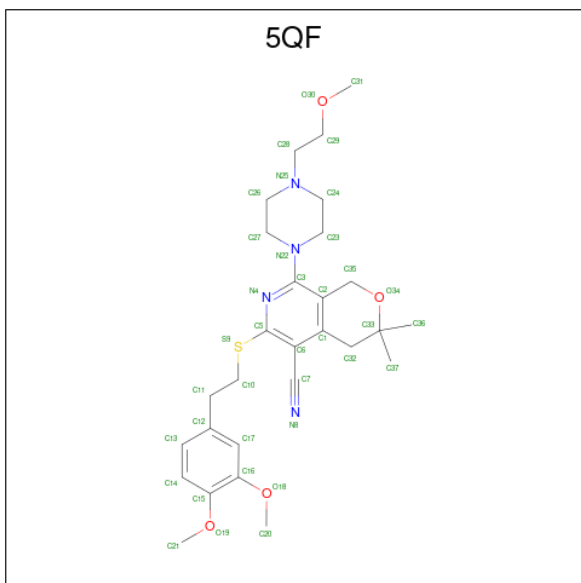
Chain	Residue	Modelled	Actual	Comment	Reference
M	50	HIS	-	expression tag	UNP P0AAW9
M	51	HIS	-	expression tag	UNP P0AAW9

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Chain	Residue	Modelled	Actual	Comment	Reference
M	52	HIS	-	expression tag	UNP P0AAW9
M	53	HIS	-	expression tag	UNP P0AAW9
M	54	HIS	-	expression tag	UNP P0AAW9
N	50	HIS	-	expression tag	UNP P0AAW9
N	51	HIS	-	expression tag	UNP P0AAW9
N	52	HIS	-	expression tag	UNP P0AAW9
N	53	HIS	-	expression tag	UNP P0AAW9
N	54	HIS	-	expression tag	UNP P0AAW9
O	50	HIS	-	expression tag	UNP P0AAW9
O	51	HIS	-	expression tag	UNP P0AAW9
O	52	HIS	-	expression tag	UNP P0AAW9
O	53	HIS	-	expression tag	UNP P0AAW9
O	54	HIS	-	expression tag	UNP P0AAW9

- Molecule 5 is 6-[2-(3,4-dimethoxyphenyl)ethylsulfanyl]-8-[4-(2-methoxyethyl)piperazin-1-yl]-3,3-dimethyl-1,4-dihydropyranopyridine-5-carbonitrile (three-letter code: 5QF) (formula: C₂₈H₃₈N₄O₄S).

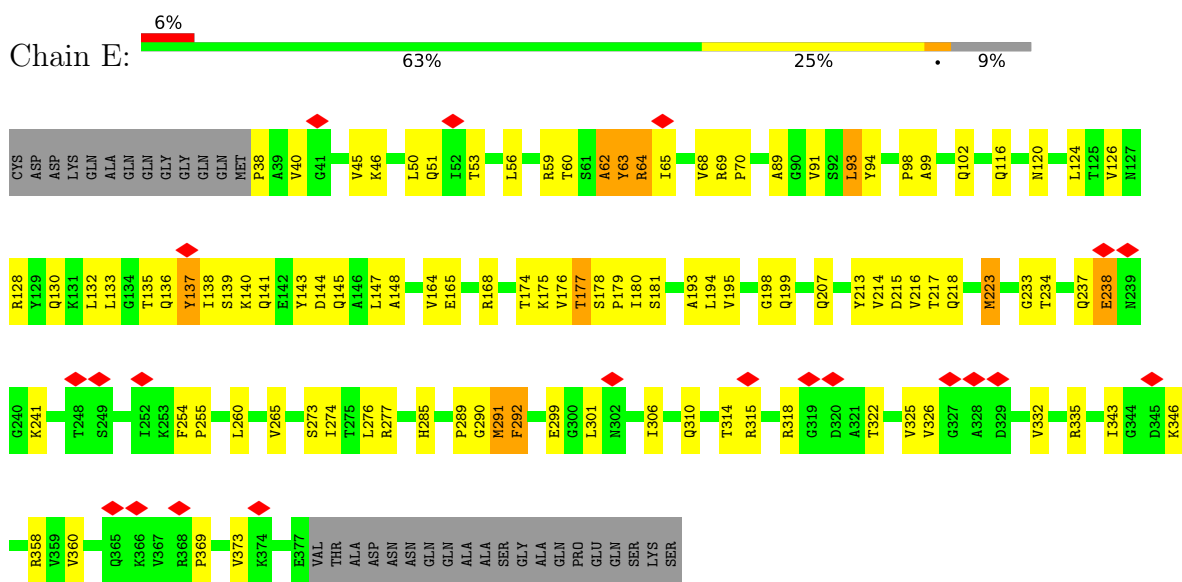


Mol	Chain	Residues	Atoms					AltConf	
			Total	C	H	N	O		S
5	K	1	Total	C	H	N	O	S	0
			75	28	38	4	4	1	
5	L	1	Total	C	H	N	O	S	0
			75	28	38	4	4	1	
5	J	1	Total	C	H	N	O	S	0
			75	28	38	4	4	1	

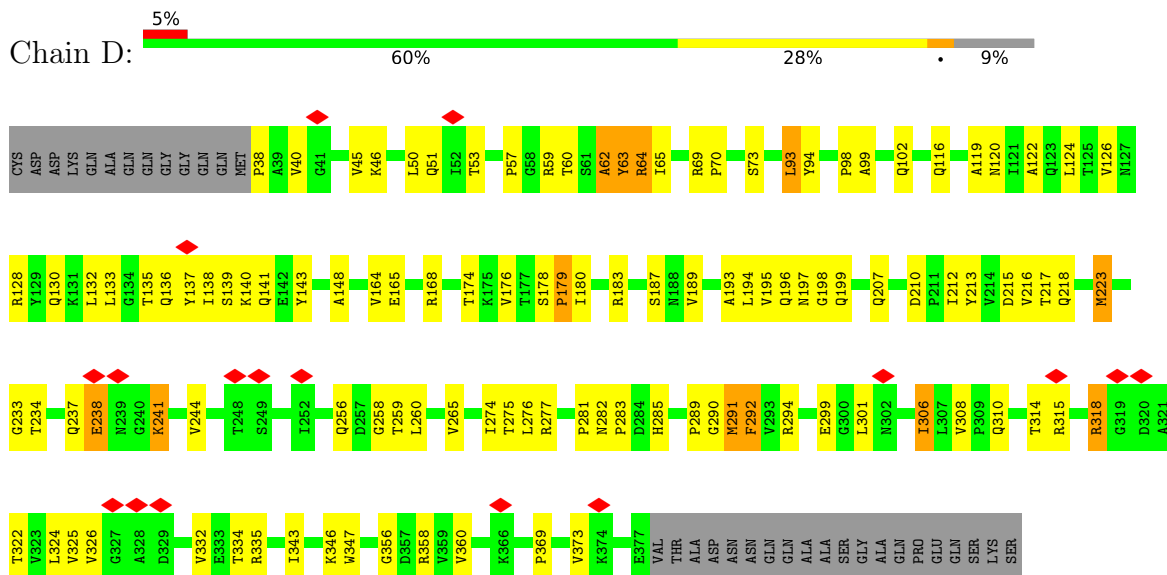
3 Residue-property plots

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.

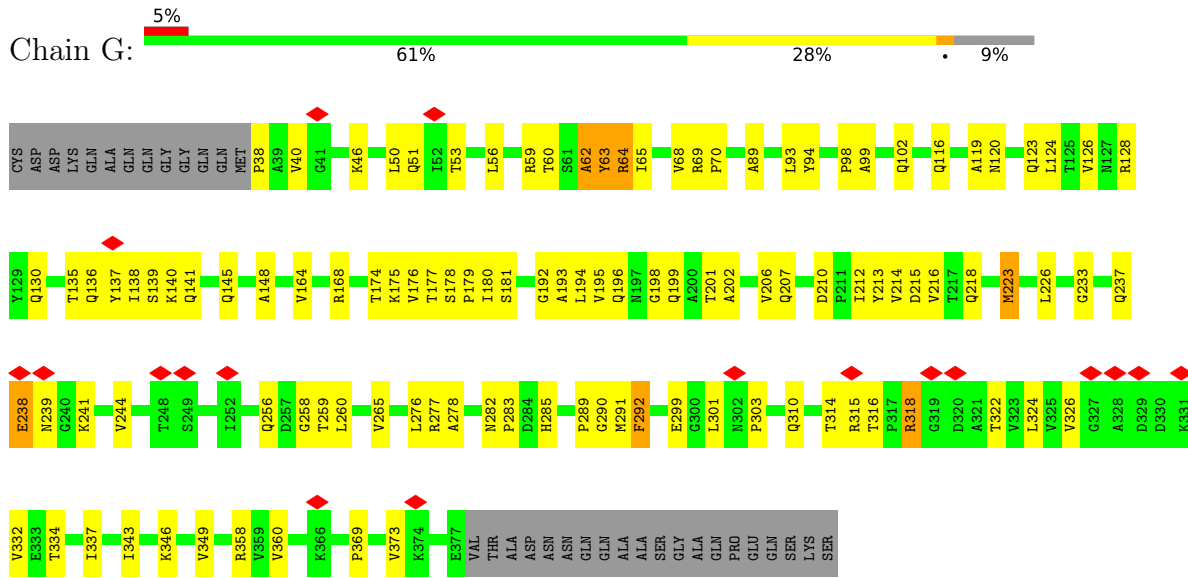
- Molecule 1: Multidrug efflux pump subunit AcrA



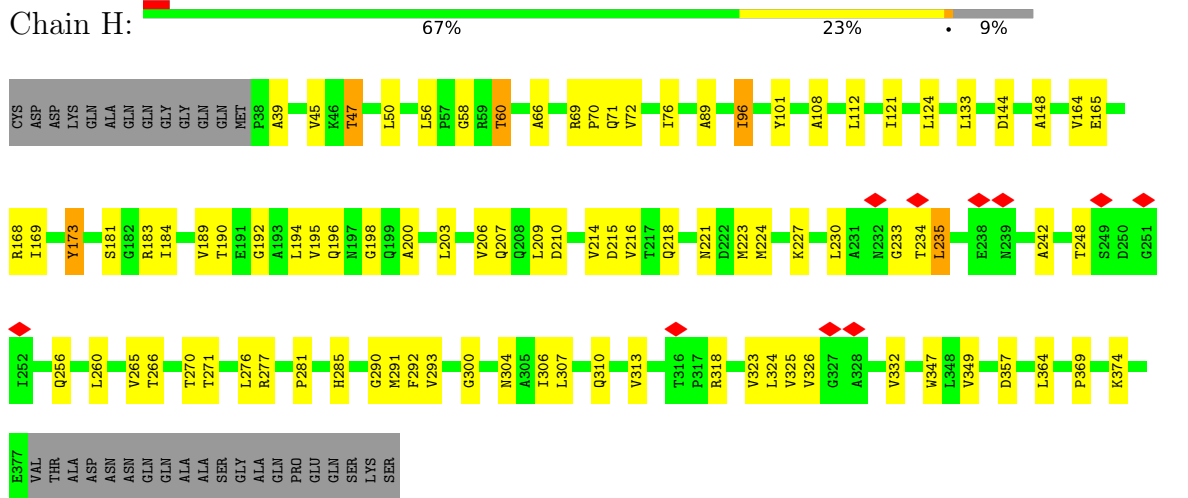
- Molecule 1: Multidrug efflux pump subunit AcrA



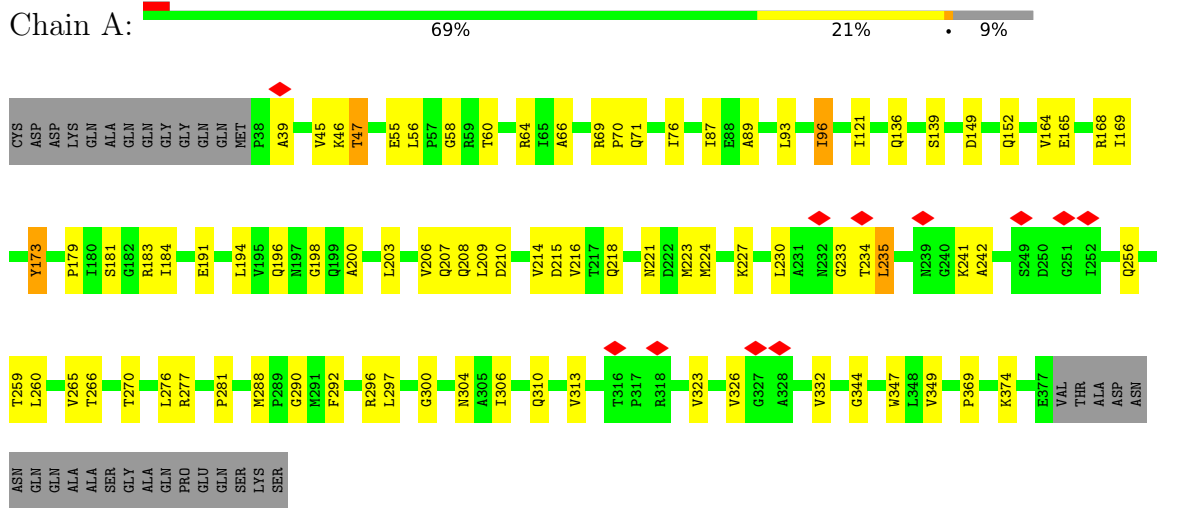
- Molecule 1: Multidrug efflux pump subunit AcrA



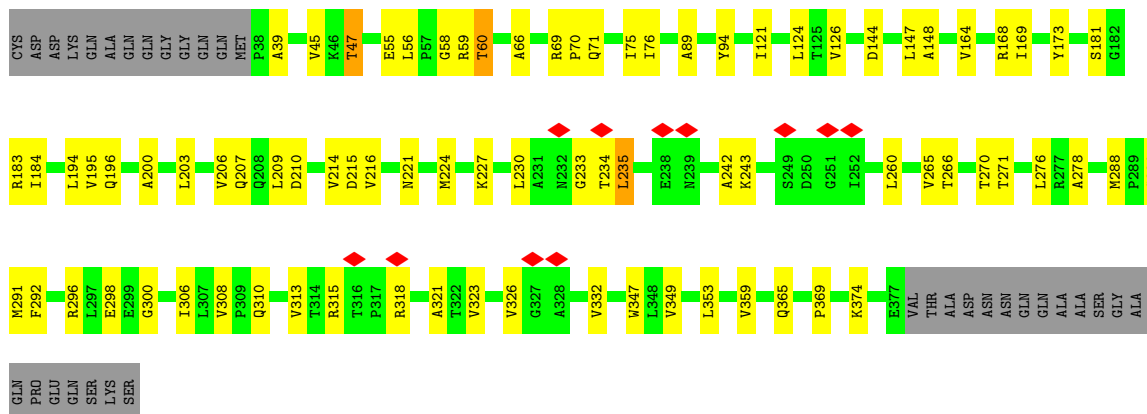
• Molecule 1: Multidrug efflux pump subunit AcrA



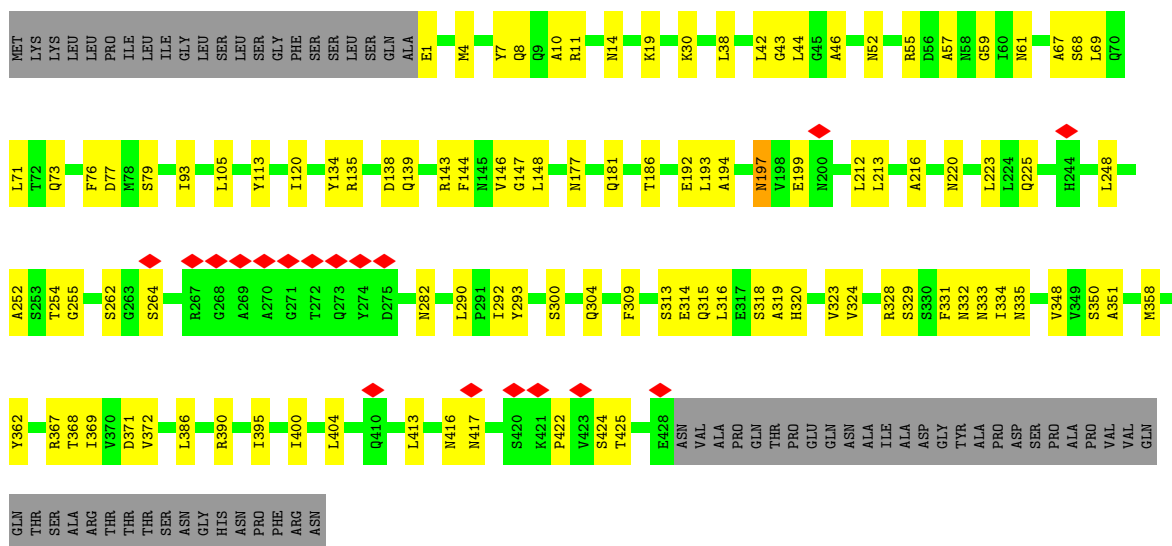
• Molecule 1: Multidrug efflux pump subunit AcrA



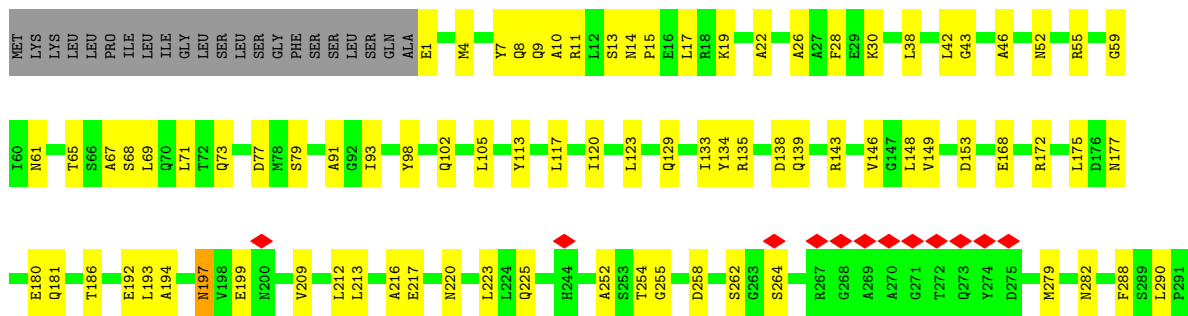
• Molecule 1: Multidrug efflux pump subunit AcrA

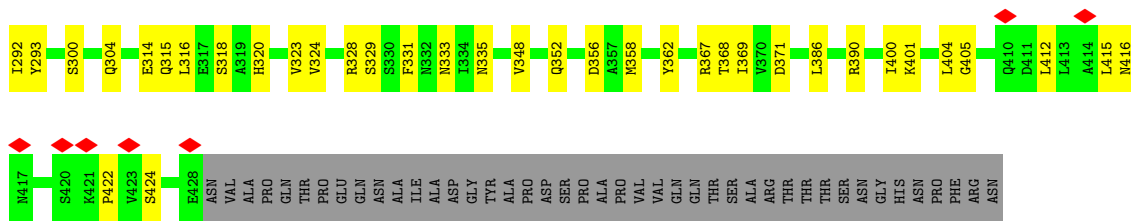


• Molecule 2: Outer membrane protein TolC

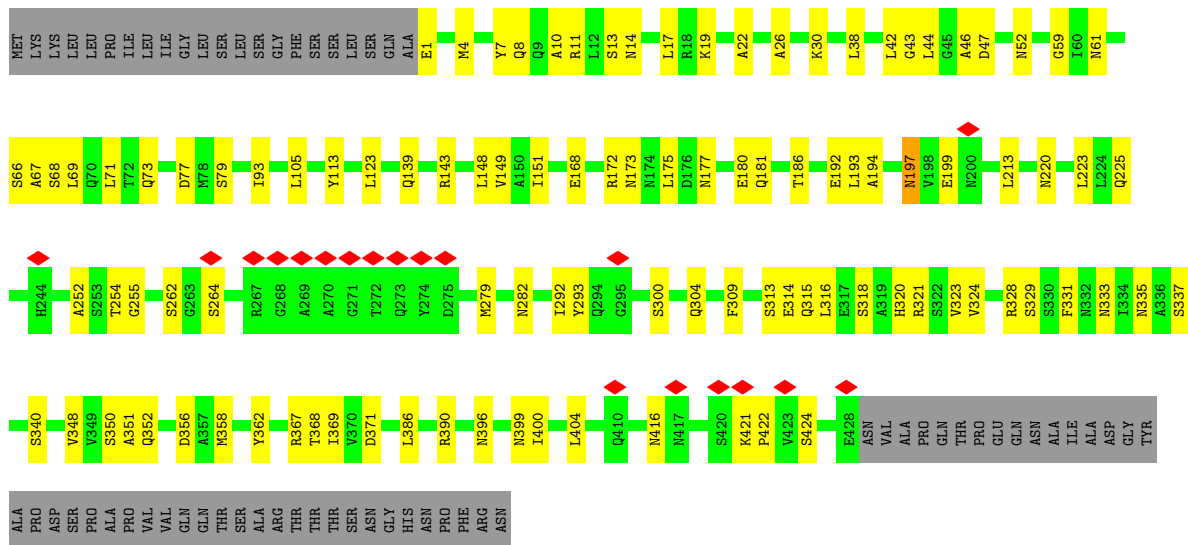


• Molecule 2: Outer membrane protein TolC

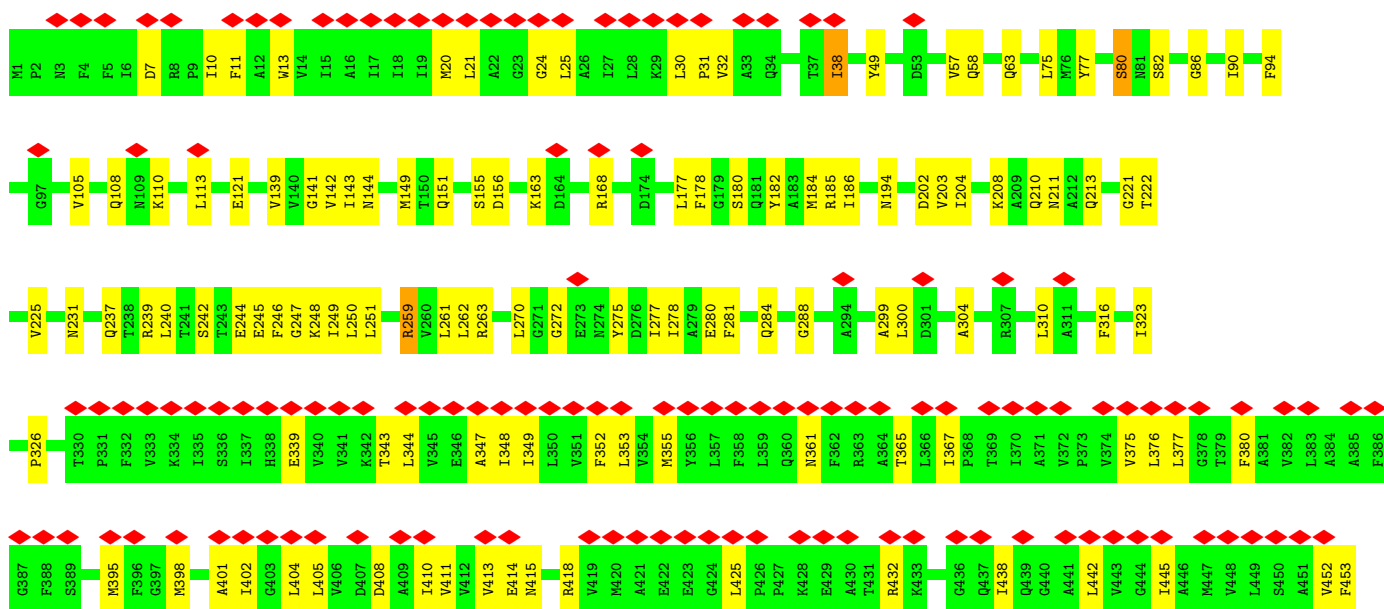


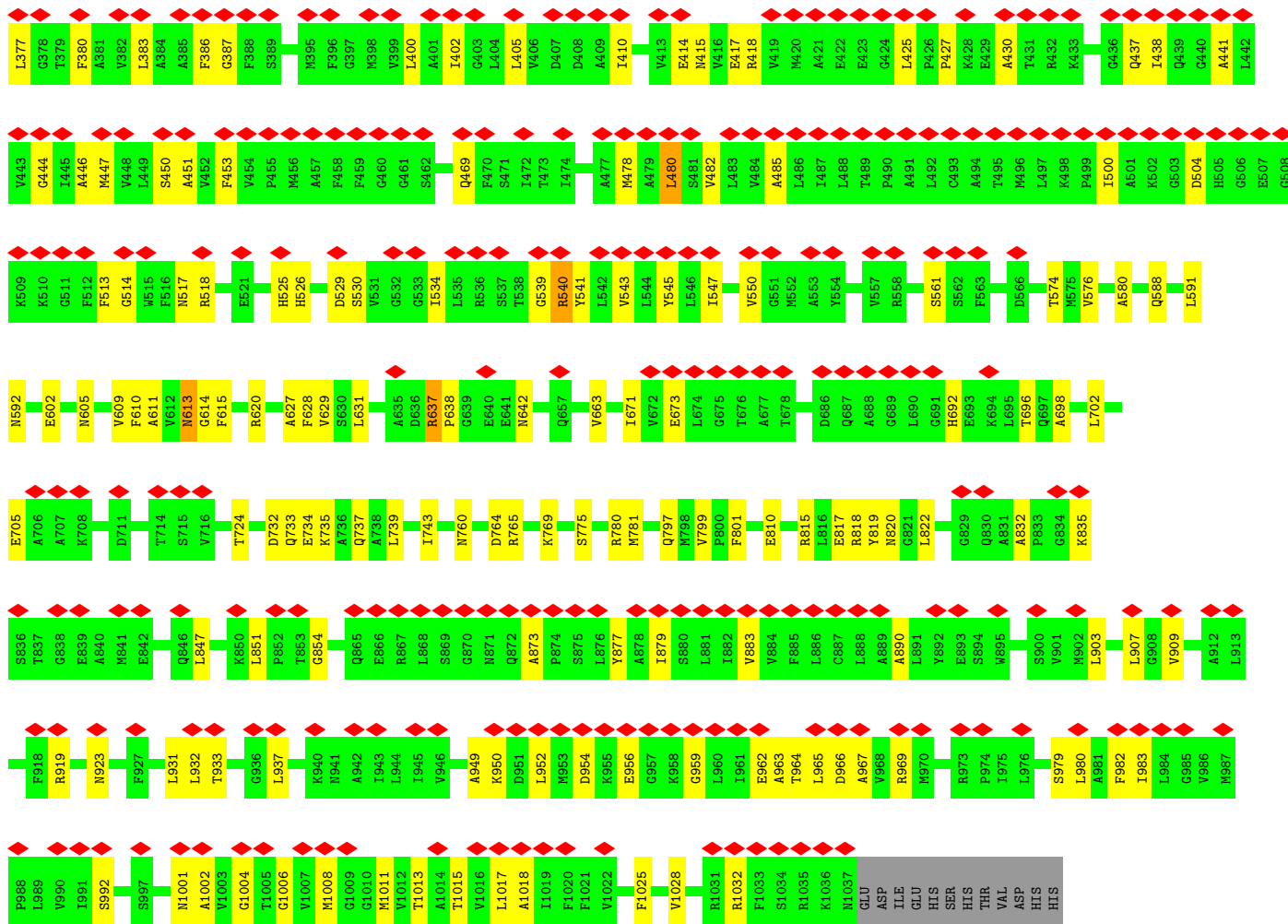


• Molecule 2: Outer membrane protein TolC

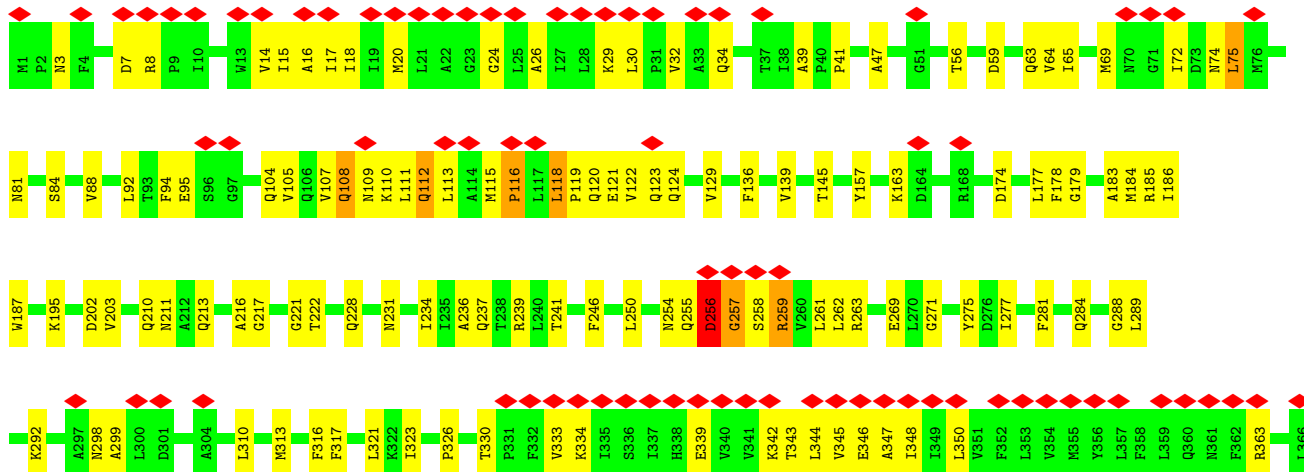


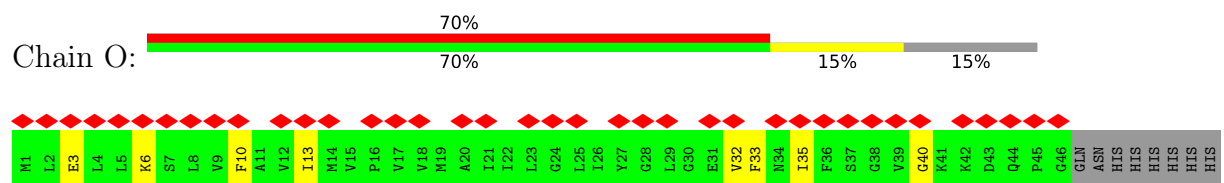
• Molecule 3: Multidrug efflux pump subunit AcrB





• Molecule 3: Multidrug efflux pump subunit AcrB





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	13544	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.143	Depositor
Minimum map value	-0.070	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.026	Depositor
Map size (\AA)	399.9, 399.9, 399.9	wwPDB
Map dimensions	310, 310, 310	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.29, 1.29, 1.29	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: 5QF

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	0.57	0/2585	0.74	1/3512 (0.0%)
1	B	0.58	0/2585	0.75	1/3512 (0.0%)
1	D	0.56	0/2585	0.79	3/3512 (0.1%)
1	E	0.59	1/2585 (0.0%)	0.78	3/3512 (0.1%)
1	G	0.57	1/2585 (0.0%)	0.80	3/3512 (0.1%)
1	H	0.57	0/2585	0.76	2/3512 (0.1%)
2	C	0.43	0/3345	0.60	0/4544
2	F	0.44	1/3345 (0.0%)	0.60	0/4544
2	I	0.43	0/3345	0.60	0/4544
3	J	0.49	0/7999	0.71	10/10863 (0.1%)
3	K	0.51	2/7968 (0.0%)	0.69	7/10826 (0.1%)
3	L	0.49	0/7968	0.68	2/10826 (0.0%)
4	M	0.29	0/330	0.61	0/448
4	N	0.30	0/330	0.62	0/448
4	O	0.31	0/330	0.61	0/448
All	All	0.51	5/50470 (0.0%)	0.70	32/68563 (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	6
1	B	0	4
1	D	0	10
1	E	0	10
1	G	0	6
1	H	0	5
3	J	0	9
3	K	0	4

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Mol	Chain	#Chirality outliers	#Planarity outliers
3	L	0	3
All	All	0	57

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	68	VAL	C-N	-7.85	1.16	1.34
3	K	805	SER	CA-CB	-7.13	1.42	1.52
1	G	68	VAL	C-N	-5.13	1.22	1.34
3	K	789	TRP	CB-CG	-5.05	1.41	1.50
2	F	146	VAL	CB-CG1	-5.04	1.42	1.52

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	256	ASP	CB-CG-OD1	12.60	129.64	118.30
1	H	56	LEU	CB-CG-CD1	-8.22	97.02	111.00
1	A	56	LEU	CB-CG-CD1	-7.29	98.62	111.00
3	J	113	LEU	CA-CB-CG	7.00	131.40	115.30
1	B	56	LEU	CB-CG-CD1	-6.76	99.51	111.00

There are no chirality outliers.

5 of 57 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	135	THR	Peptide
1	E	174	THR	Peptide
1	E	177	THR	Peptide
1	E	62	ALA	Peptide
1	E	93	LEU	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2553	0	2610	59	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2553	0	2610	55	0
1	D	2553	0	2610	80	0
1	E	2553	0	2609	68	0
1	G	2553	0	2610	76	0
1	H	2553	0	2610	61	0
2	C	3304	0	3254	78	0
2	F	3304	0	3254	78	0
2	I	3304	0	3254	88	0
3	J	7849	0	8001	226	0
3	K	7819	0	7919	207	0
3	L	7819	0	7919	204	0
4	M	326	0	346	7	0
4	N	326	0	346	4	0
4	O	326	0	346	6	0
5	J	37	38	0	49	0
5	K	37	38	0	43	0
5	L	37	38	0	42	0
All	All	49806	114	50298	1161	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 12.

The worst 5 of 1161 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)
3:J:628:PHE:HE2	5:J:1101:5QF:C5	1.07	1.64
3:J:628:PHE:CE2	5:J:1101:5QF:C5	1.84	1.59
3:K:277:ILE:HG21	5:K:1101:5QF:C16	1.31	1.57
3:L:178:PHE:HB3	5:L:1101:5QF:C15	1.39	1.49
3:J:628:PHE:HE2	5:J:1101:5QF:N4	1.18	1.39

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	338/373 (91%)	299 (88%)	39 (12%)	0	100	100
1	B	338/373 (91%)	300 (89%)	38 (11%)	0	100	100
1	D	338/373 (91%)	288 (85%)	48 (14%)	2 (1%)	25	66
1	E	338/373 (91%)	286 (85%)	50 (15%)	2 (1%)	25	66
1	G	338/373 (91%)	283 (84%)	53 (16%)	2 (1%)	25	66
1	H	338/373 (91%)	297 (88%)	41 (12%)	0	100	100
2	C	426/493 (86%)	407 (96%)	19 (4%)	0	100	100
2	F	426/493 (86%)	407 (96%)	19 (4%)	0	100	100
2	I	426/493 (86%)	408 (96%)	18 (4%)	0	100	100
3	J	1031/1049 (98%)	937 (91%)	93 (9%)	1 (0%)	51	85
3	K	1035/1049 (99%)	925 (89%)	110 (11%)	0	100	100
3	L	1035/1049 (99%)	935 (90%)	100 (10%)	0	100	100
4	M	44/54 (82%)	39 (89%)	5 (11%)	0	100	100
4	N	44/54 (82%)	40 (91%)	4 (9%)	0	100	100
4	O	44/54 (82%)	41 (93%)	3 (7%)	0	100	100
All	All	6539/7026 (93%)	5892 (90%)	640 (10%)	7 (0%)	54	85

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	140	LYS
1	D	140	LYS
1	G	140	LYS
1	D	179	PRO
3	J	257	GLY

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	274/299 (92%)	273 (100%)	1 (0%)	91	94
1	B	274/299 (92%)	273 (100%)	1 (0%)	91	94
1	D	274/299 (92%)	270 (98%)	4 (2%)	65	80
1	E	274/299 (92%)	269 (98%)	5 (2%)	59	77
1	G	274/299 (92%)	270 (98%)	4 (2%)	65	80
1	H	274/299 (92%)	272 (99%)	2 (1%)	84	90
2	C	358/412 (87%)	356 (99%)	2 (1%)	86	92
2	F	358/412 (87%)	356 (99%)	2 (1%)	86	92
2	I	358/412 (87%)	356 (99%)	2 (1%)	86	92
3	J	839/855 (98%)	832 (99%)	7 (1%)	81	89
3	K	826/855 (97%)	821 (99%)	5 (1%)	86	92
3	L	826/855 (97%)	815 (99%)	11 (1%)	69	82
4	M	34/46 (74%)	34 (100%)	0	100	100
4	N	34/46 (74%)	34 (100%)	0	100	100
4	O	34/46 (74%)	34 (100%)	0	100	100
All	All	5311/5733 (93%)	5265 (99%)	46 (1%)	79	88

5 of 46 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	L	231	ASN
3	L	637	ARG
3	L	255	GLN
3	L	418	ARG
3	L	801	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 30 such sidechains are listed below:

Mol	Chain	Res	Type
3	K	108	GLN
3	J	109	ASN
3	K	213	GLN
3	J	733	GLN
3	L	176	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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
5	5QF	L	1101	3	40,40,40	1.94	12 (30%)	49,56,56	2.43	15 (30%)
5	5QF	K	1101	3	40,40,40	1.96	10 (25%)	49,56,56	2.09	12 (24%)
5	5QF	J	1101	-	40,40,40	2.03	10 (25%)	49,56,56	2.15	12 (24%)

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
5	5QF	L	1101	3	-	11/19/41/41	0/4/4/4
5	5QF	K	1101	3	-	12/19/41/41	0/4/4/4
5	5QF	J	1101	-	-	10/19/41/41	0/4/4/4

The worst 5 of 32 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	K	1101	5QF	C5-S9	4.77	1.83	1.76
5	K	1101	5QF	C3-N22	4.74	1.49	1.37
5	J	1101	5QF	C3-N22	4.42	1.48	1.37
5	L	1101	5QF	C2-C1	-4.37	1.33	1.40
5	K	1101	5QF	C6-C7	4.25	1.51	1.44

The worst 5 of 39 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	1101	5QF	C10-C11-C12	-7.65	100.22	113.97
5	K	1101	5QF	C10-S9-C5	7.01	112.87	101.82
5	L	1101	5QF	C10-S9-C5	6.81	112.55	101.82
5	J	1101	5QF	O18-C16-C15	6.59	124.60	115.41
5	K	1101	5QF	C6-C5-N4	-4.94	119.37	123.12

There are no chirality outliers.

5 of 33 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	L	1101	5QF	C11-C10-S9-C5
5	J	1101	5QF	S9-C10-C11-C12
5	L	1101	5QF	C16-C15-O19-C21
5	K	1101	5QF	C15-C16-O18-C20
5	J	1101	5QF	C14-C15-O19-C21

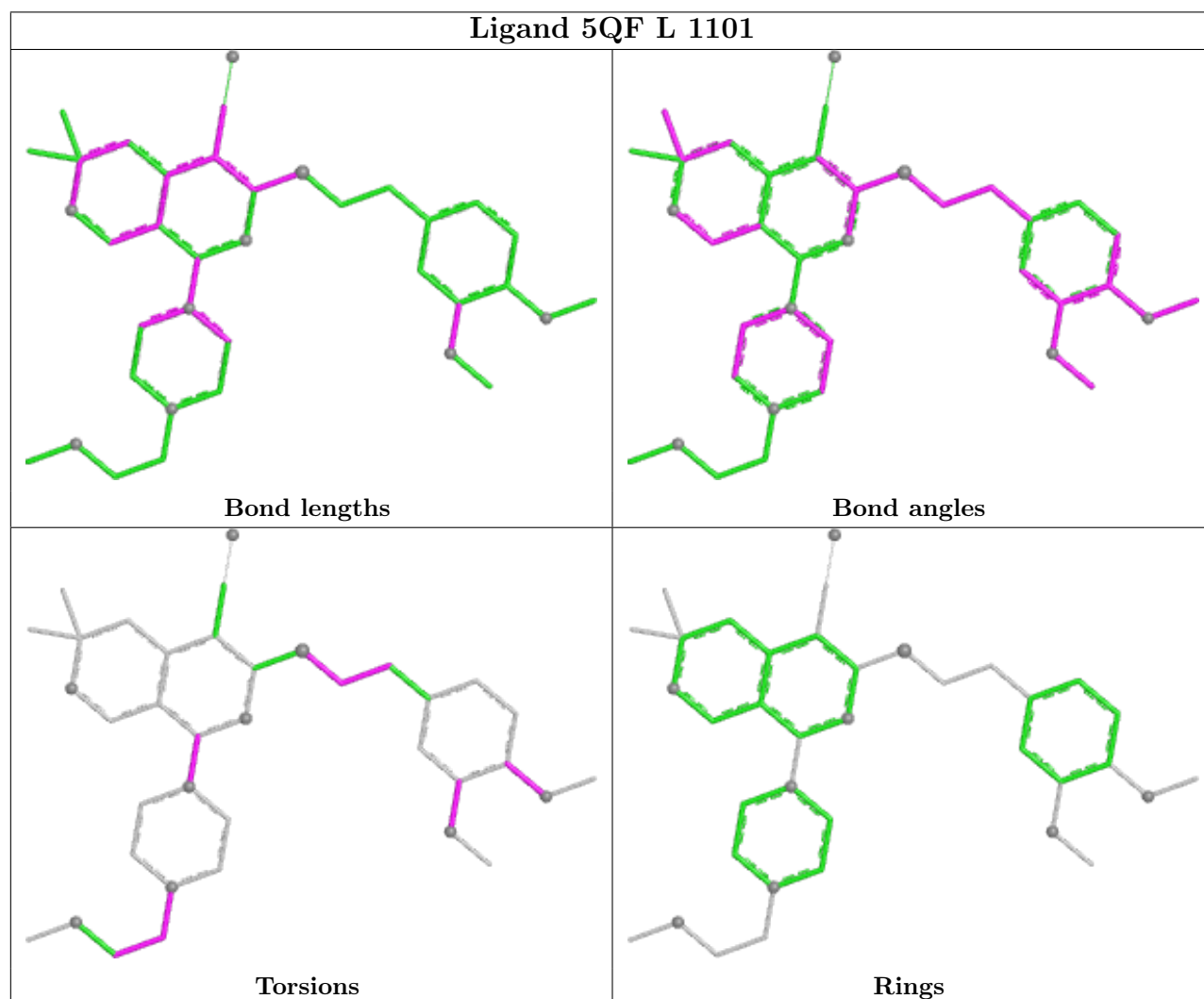
There are no ring outliers.

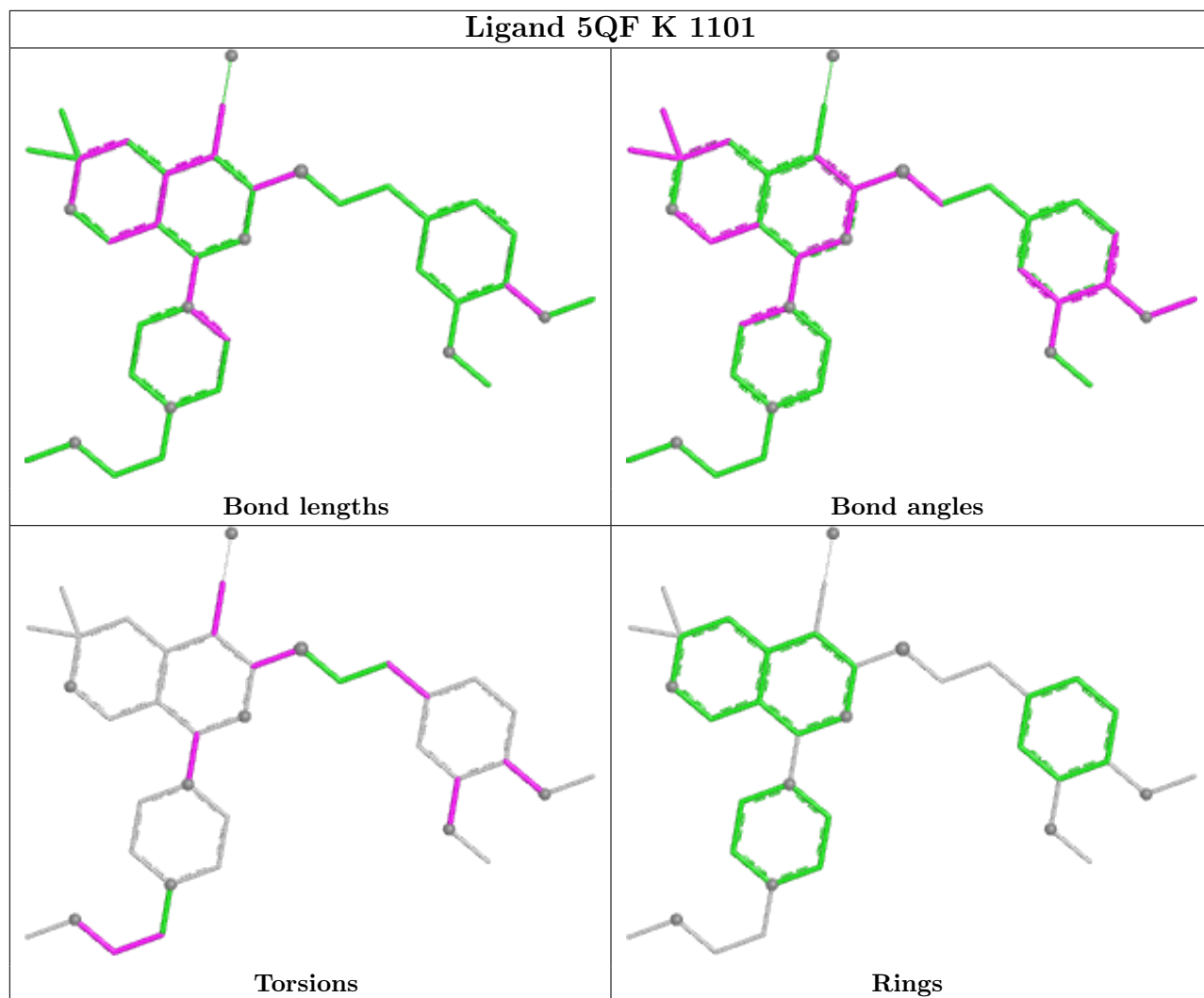
3 monomers are involved in 134 short contacts:

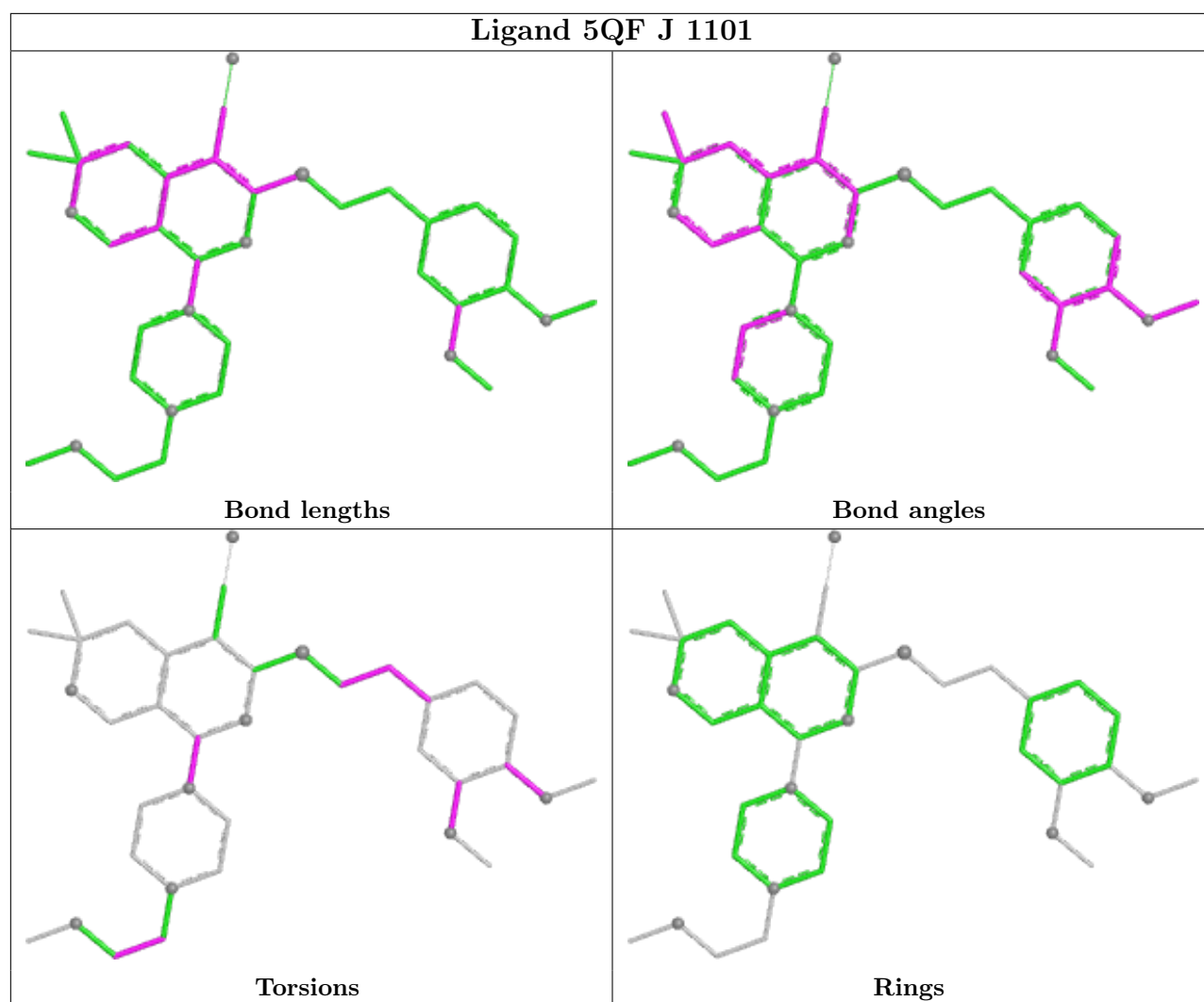
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	L	1101	5QF	42	0
5	K	1101	5QF	43	0
5	J	1101	5QF	49	0

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	E	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	68:VAL	C	69:ARG	N	1.16

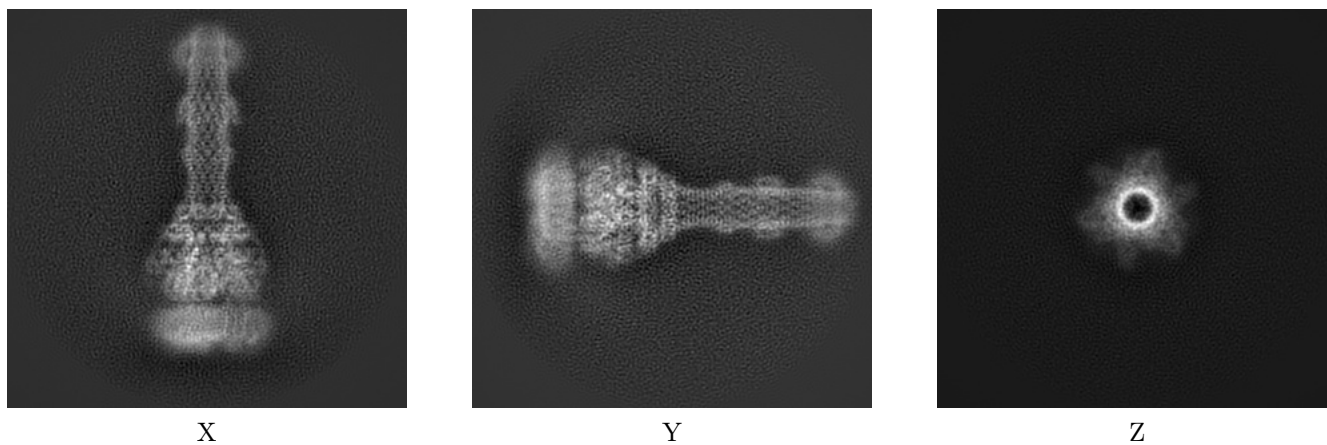
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-3636. 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](#)

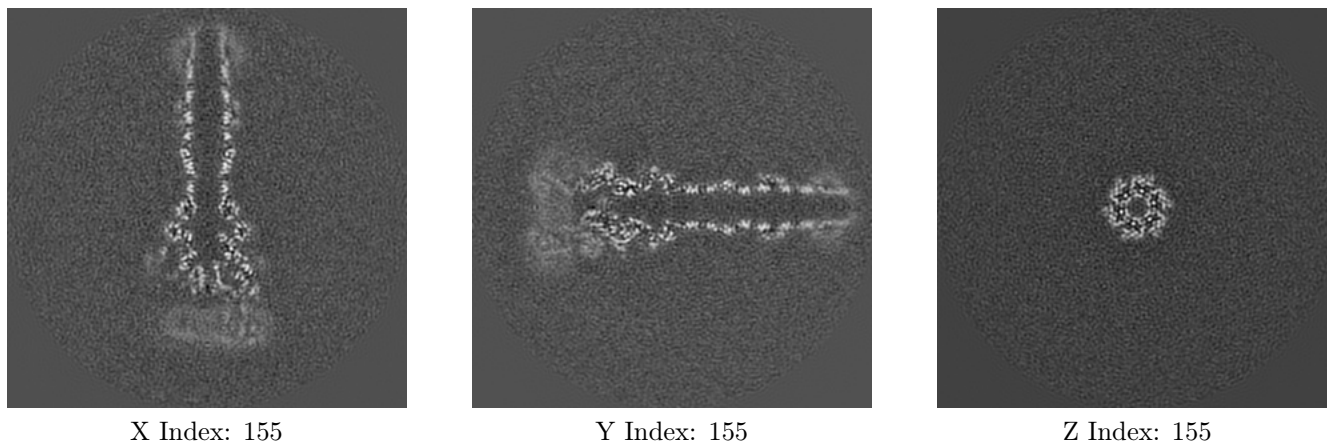
6.1.1 Primary map



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

6.2 Central slices [i](#)

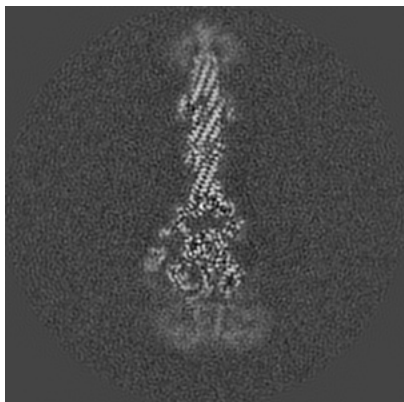
6.2.1 Primary map



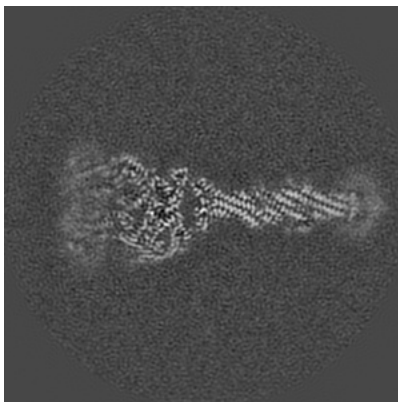
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

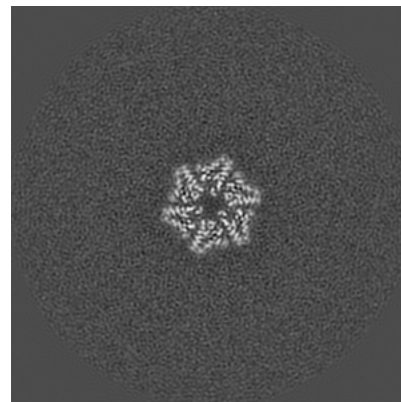
6.3.1 Primary map



X Index: 142



Y Index: 142



Z Index: 134

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 0.026. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

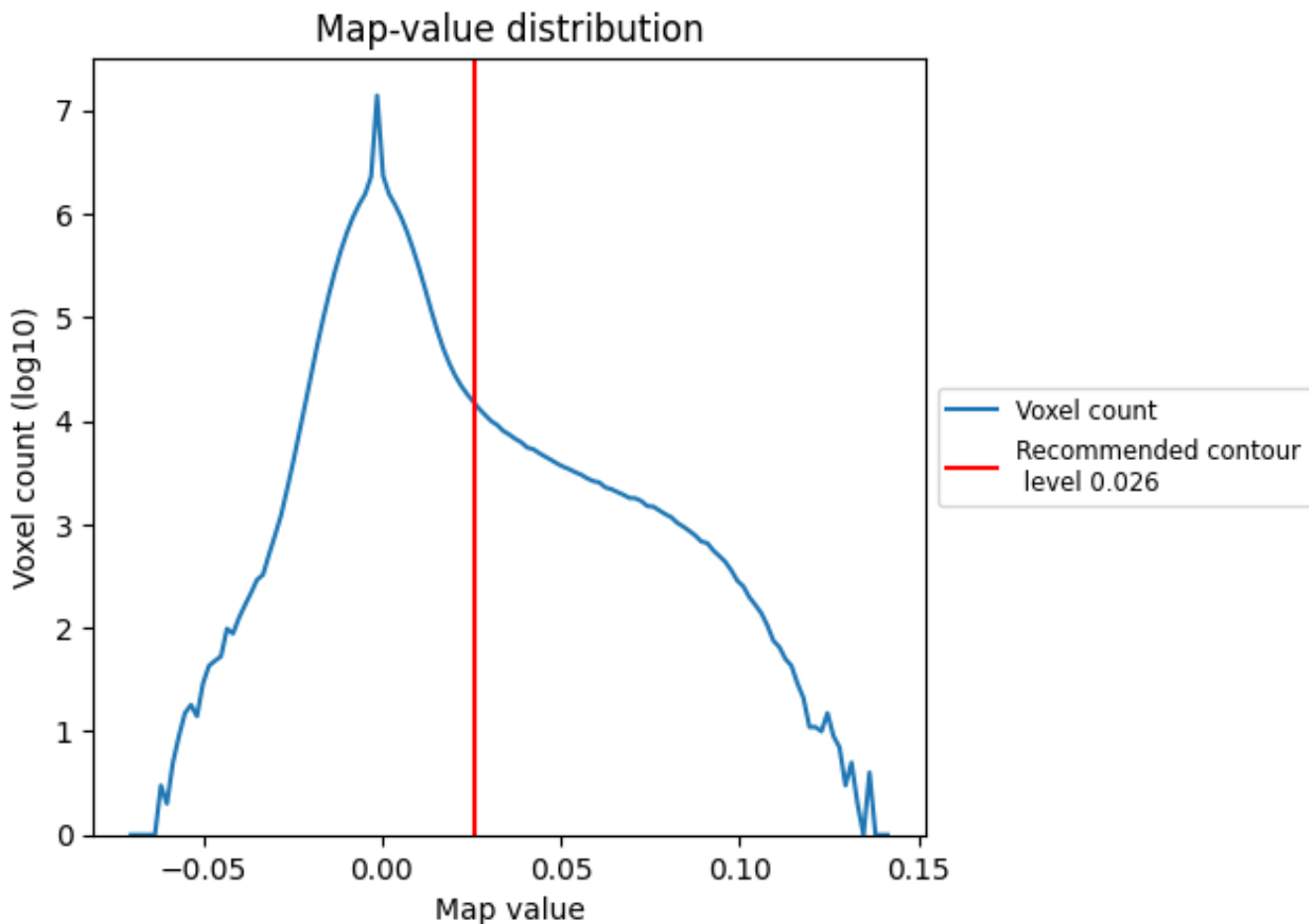
6.5 Mask visualisation

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

7 Map analysis [i](#)

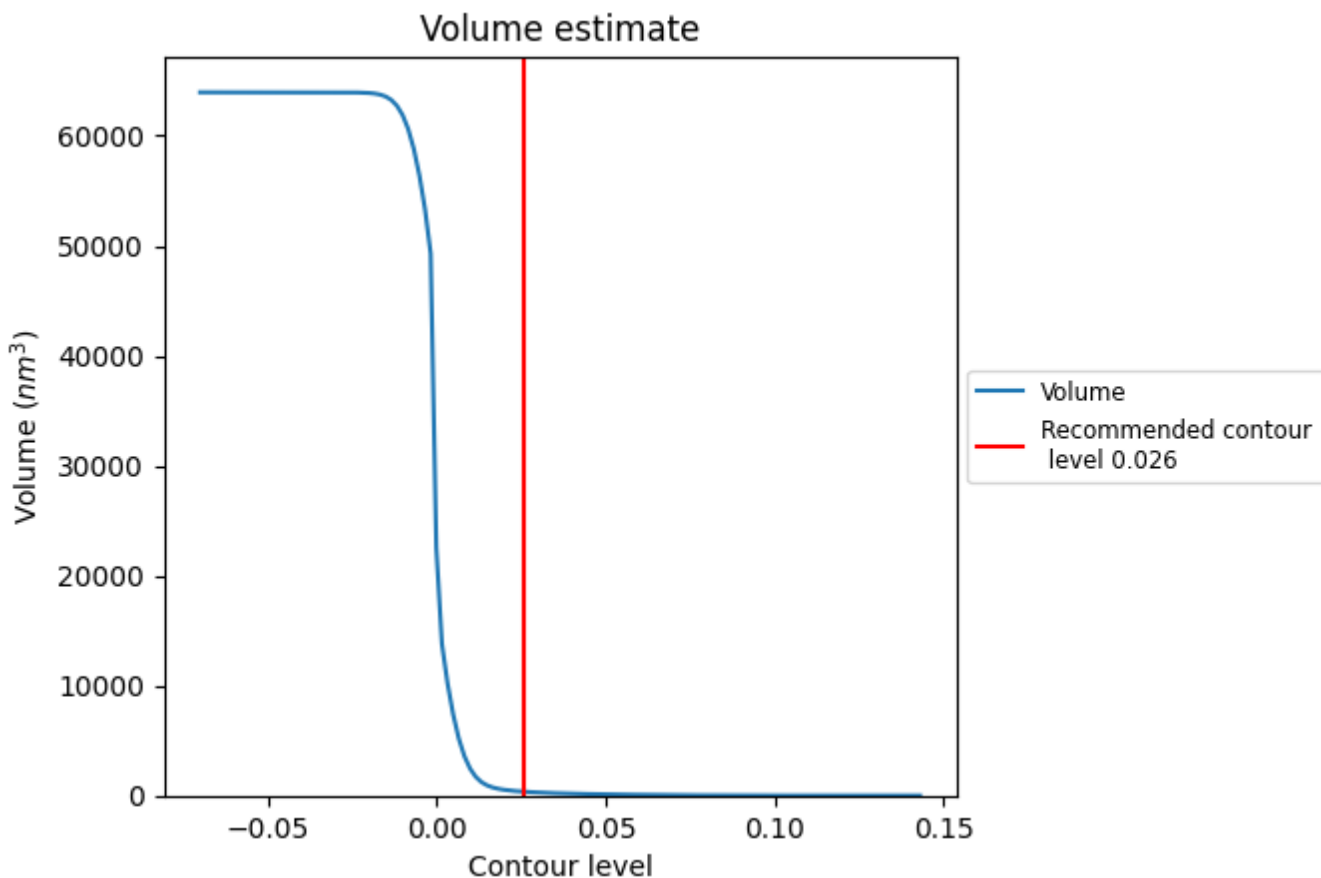
This section contains the results of statistical analysis of the map.

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.

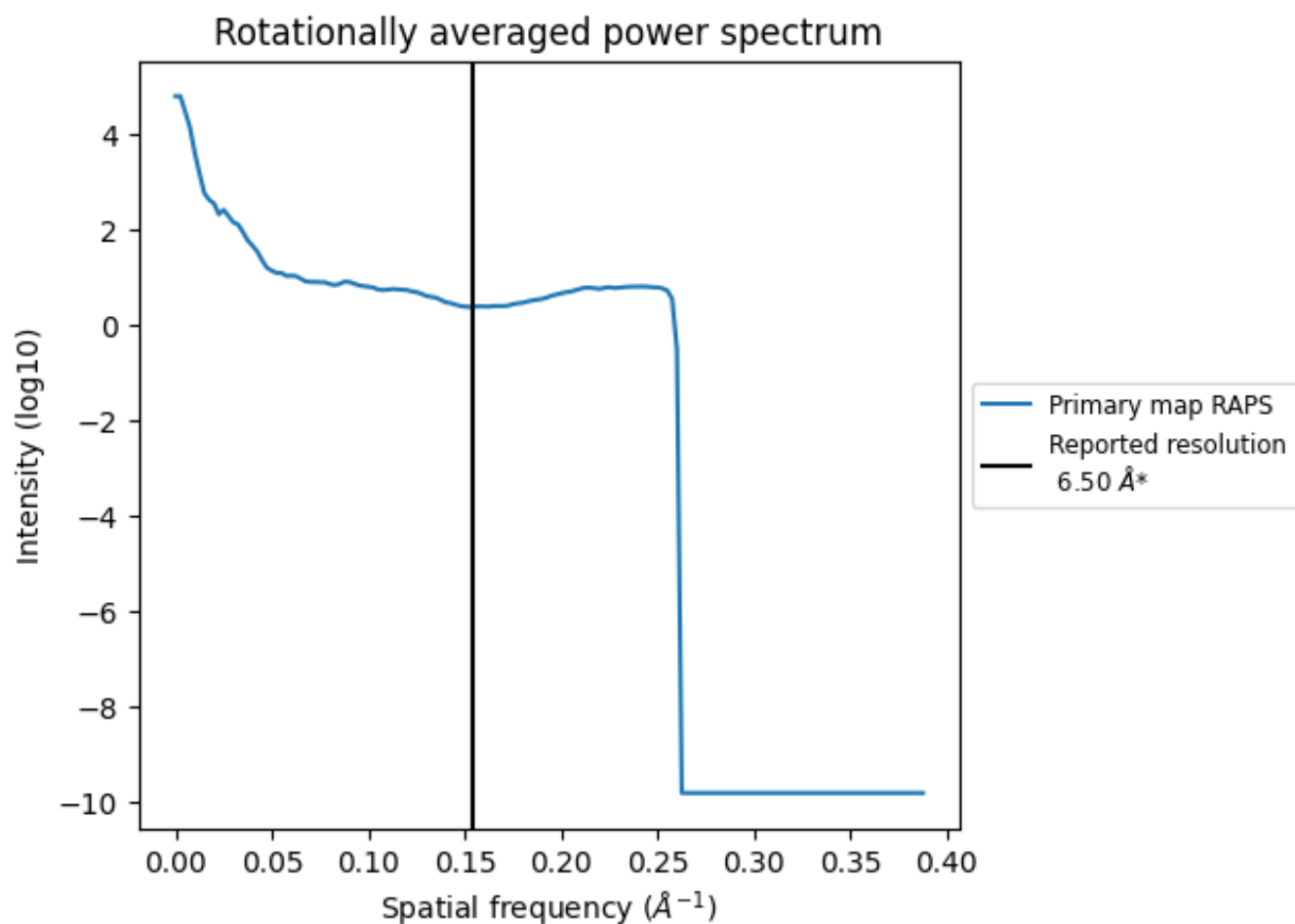
7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 346 nm³; this corresponds to an approximate mass of 312 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 [\(i\)](#)

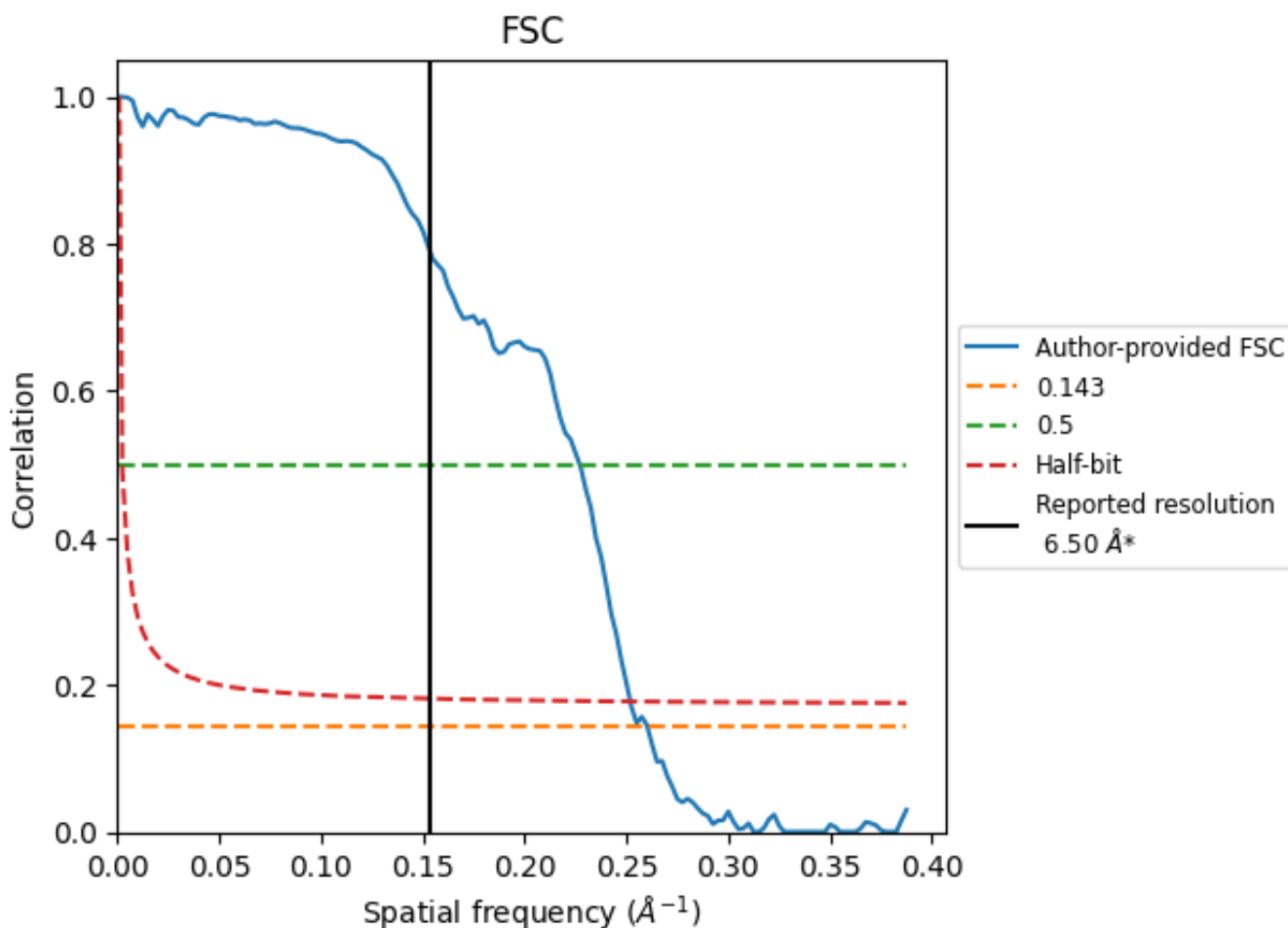


*Reported resolution corresponds to spatial frequency of 0.154 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.154 Å⁻¹

8.2 Resolution estimates [i](#)

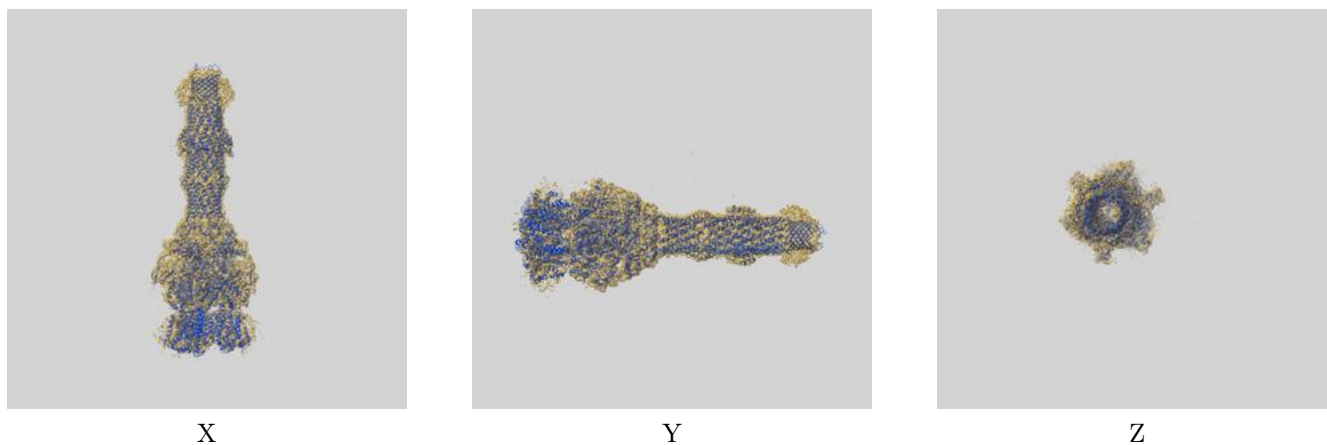
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	6.50	-	-
Author-provided FSC curve	3.84	4.41	3.97
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from author-provided FSC intersecting FSC 0.143 CUT-OFF 3.84 differs from the reported value 6.5 by more than 10 %

9 Map-model fit [i](#)

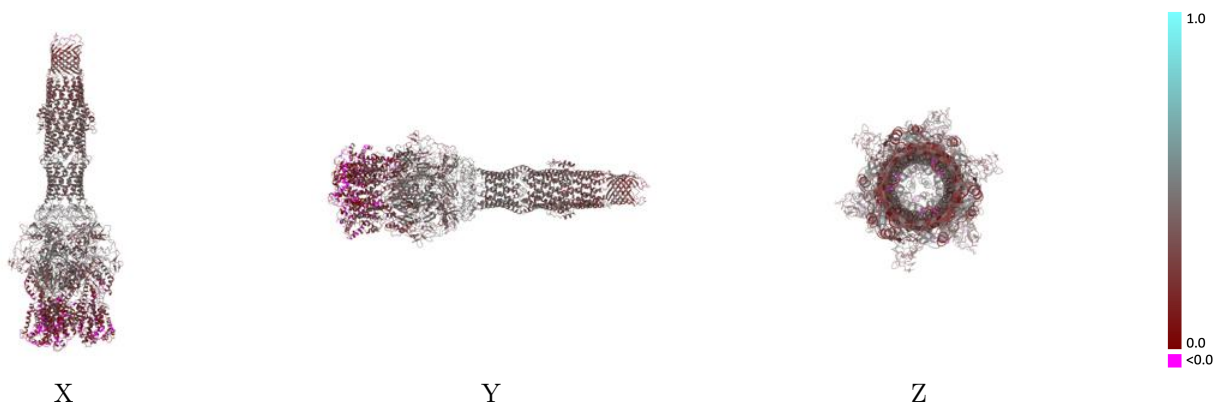
This section contains information regarding the fit between EMDB map EMD-3636 and PDB model 5NG5. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)



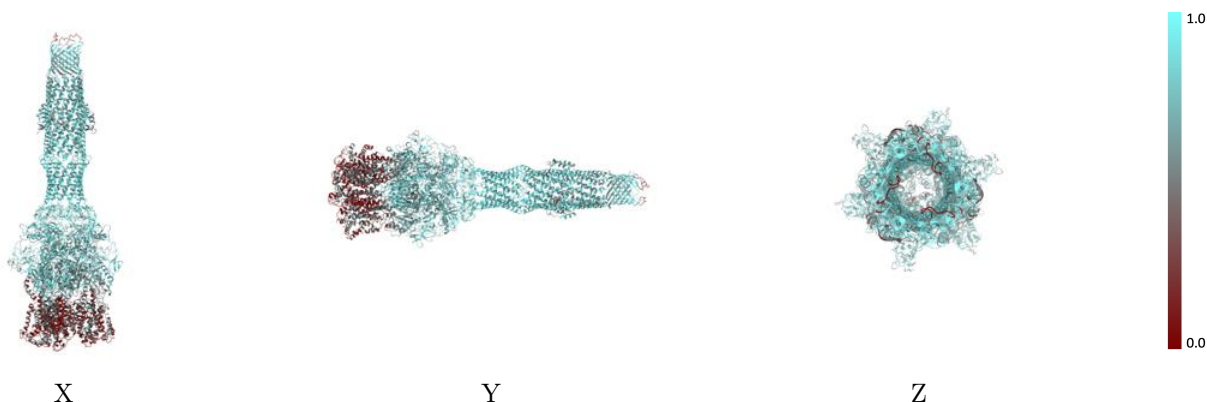
The images above show the 3D surface view of the map at the recommended contour level 0.026 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



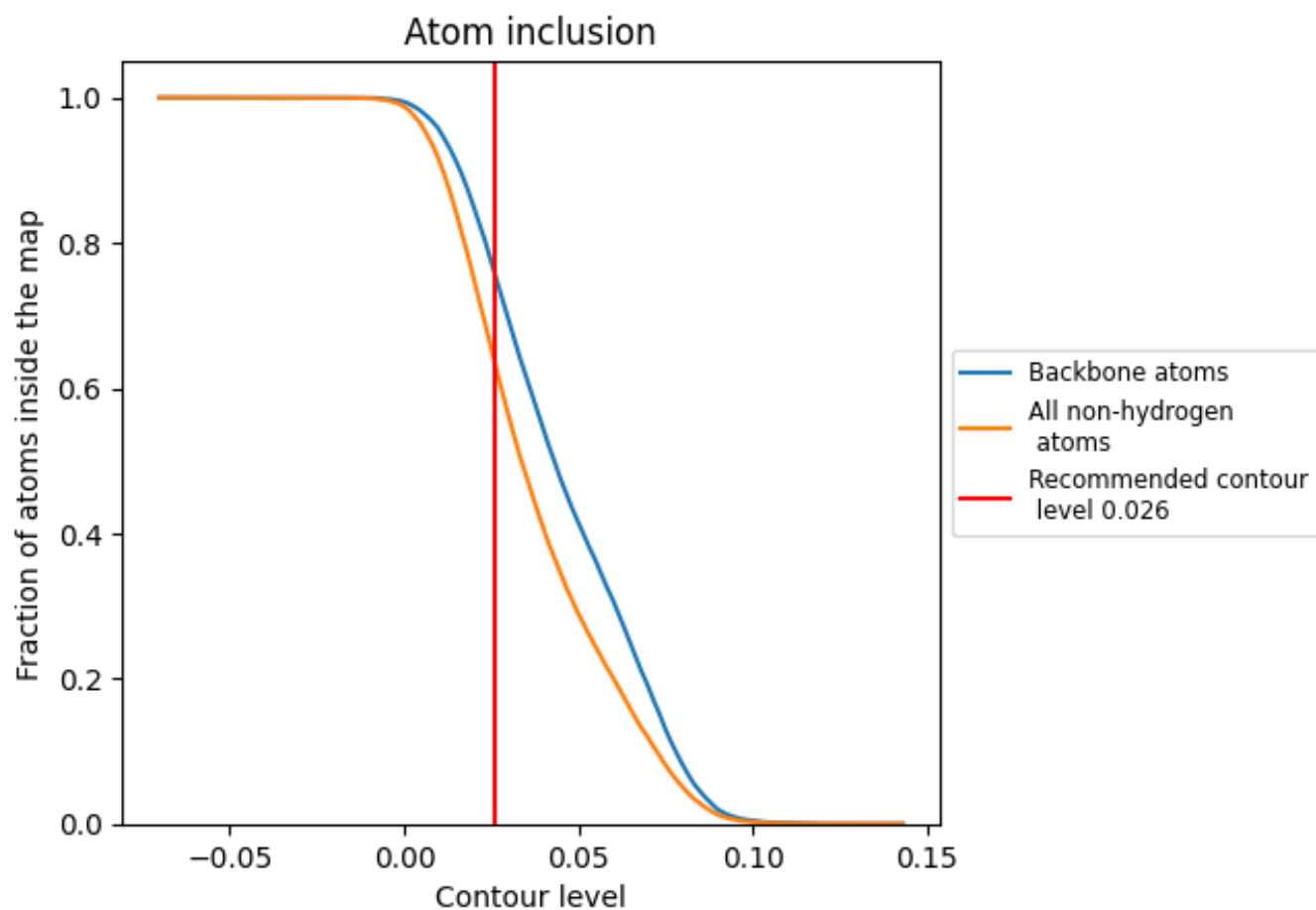
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.026).

































9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 63% 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 (0.026) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6350	 0.3260
A	 0.7670	 0.4010
B	 0.7698	 0.4040
C	 0.7286	 0.3410
D	 0.7519	 0.3740
E	 0.7547	 0.3730
F	 0.7274	 0.3400
G	 0.7595	 0.3750
H	 0.7714	 0.4020
I	 0.7283	 0.3390
J	 0.5296	 0.2900
K	 0.5334	 0.2900
L	 0.5408	 0.2920
M	 0.1938	 0.0740
N	 0.1877	 0.0960
O	 0.1969	 0.0670

