



wwPDB EM Validation Summary Report ⓘ

Nov 12, 2022 – 05:13 PM EST

PDB ID : 6PVQ
EMDB ID : EMD-20497
Title : Cryo-EM structure of mouse TRPV3-Y564A in intermediate state at 37 degrees Celsius
Authors : Singh, A.K.; McGoldrick, L.L.; Sobolevsky, A.I.
Deposited on : 2019-07-21
Resolution : 4.75 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev43
MolProbity : 4.02b-467
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

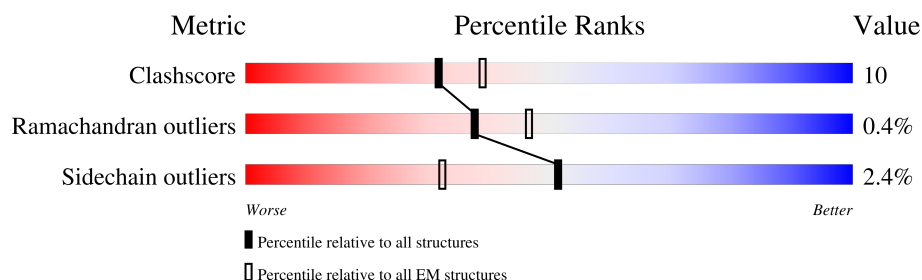
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.75 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	808	<div> <div>16%</div> <div>51%</div> <div>17%</div> <div>30%</div> </div>
1	B	808	<div> <div>15%</div> <div>47%</div> <div>16%</div> <div>36%</div> </div>
1	C	808	<div> <div>16%</div> <div>51%</div> <div>18%</div> <div>30%</div> </div>
1	D	808	<div> <div>15%</div> <div>46%</div> <div>17%</div> <div>36%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 17640 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 Transient receptor potential cation channel subfamily V member 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	562	Total	C	N	O	S	0	0
			4581	2938	781	837	25		
1	B	518	Total	C	N	O	S	0	0
			4239	2736	717	761	25		
1	C	562	Total	C	N	O	S	0	0
			4581	2938	781	837	25		
1	D	518	Total	C	N	O	S	0	0
			4239	2736	717	761	25		

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	564	ALA	TYR	engineered mutation	UNP Q8K424
A	792	LEU	-	expression tag	UNP Q8K424
A	793	VAL	-	expression tag	UNP Q8K424
A	794	PRO	-	expression tag	UNP Q8K424
A	795	ARG	-	expression tag	UNP Q8K424
A	796	GLY	-	expression tag	UNP Q8K424
A	797	SER	-	expression tag	UNP Q8K424
A	798	ALA	-	expression tag	UNP Q8K424
A	799	ALA	-	expression tag	UNP Q8K424
A	800	ALA	-	expression tag	UNP Q8K424
A	801	TRP	-	expression tag	UNP Q8K424
A	802	SER	-	expression tag	UNP Q8K424
A	803	HIS	-	expression tag	UNP Q8K424
A	804	PRO	-	expression tag	UNP Q8K424
A	805	GLN	-	expression tag	UNP Q8K424
A	806	PHE	-	expression tag	UNP Q8K424
A	807	GLU	-	expression tag	UNP Q8K424
A	808	LYS	-	expression tag	UNP Q8K424
B	564	ALA	TYR	engineered mutation	UNP Q8K424
B	792	LEU	-	expression tag	UNP Q8K424
B	793	VAL	-	expression tag	UNP Q8K424

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Chain	Residue	Modelled	Actual	Comment	Reference
B	794	PRO	-	expression tag	UNP Q8K424
B	795	ARG	-	expression tag	UNP Q8K424
B	796	GLY	-	expression tag	UNP Q8K424
B	797	SER	-	expression tag	UNP Q8K424
B	798	ALA	-	expression tag	UNP Q8K424
B	799	ALA	-	expression tag	UNP Q8K424
B	800	ALA	-	expression tag	UNP Q8K424
B	801	TRP	-	expression tag	UNP Q8K424
B	802	SER	-	expression tag	UNP Q8K424
B	803	HIS	-	expression tag	UNP Q8K424
B	804	PRO	-	expression tag	UNP Q8K424
B	805	GLN	-	expression tag	UNP Q8K424
B	806	PHE	-	expression tag	UNP Q8K424
B	807	GLU	-	expression tag	UNP Q8K424
B	808	LYS	-	expression tag	UNP Q8K424
C	564	ALA	TYR	engineered mutation	UNP Q8K424
C	792	LEU	-	expression tag	UNP Q8K424
C	793	VAL	-	expression tag	UNP Q8K424
C	794	PRO	-	expression tag	UNP Q8K424
C	795	ARG	-	expression tag	UNP Q8K424
C	796	GLY	-	expression tag	UNP Q8K424
C	797	SER	-	expression tag	UNP Q8K424
C	798	ALA	-	expression tag	UNP Q8K424
C	799	ALA	-	expression tag	UNP Q8K424
C	800	ALA	-	expression tag	UNP Q8K424
C	801	TRP	-	expression tag	UNP Q8K424
C	802	SER	-	expression tag	UNP Q8K424
C	803	HIS	-	expression tag	UNP Q8K424
C	804	PRO	-	expression tag	UNP Q8K424
C	805	GLN	-	expression tag	UNP Q8K424
C	806	PHE	-	expression tag	UNP Q8K424
C	807	GLU	-	expression tag	UNP Q8K424
C	808	LYS	-	expression tag	UNP Q8K424
D	564	ALA	TYR	engineered mutation	UNP Q8K424
D	792	LEU	-	expression tag	UNP Q8K424
D	793	VAL	-	expression tag	UNP Q8K424
D	794	PRO	-	expression tag	UNP Q8K424
D	795	ARG	-	expression tag	UNP Q8K424
D	796	GLY	-	expression tag	UNP Q8K424
D	797	SER	-	expression tag	UNP Q8K424
D	798	ALA	-	expression tag	UNP Q8K424
D	799	ALA	-	expression tag	UNP Q8K424

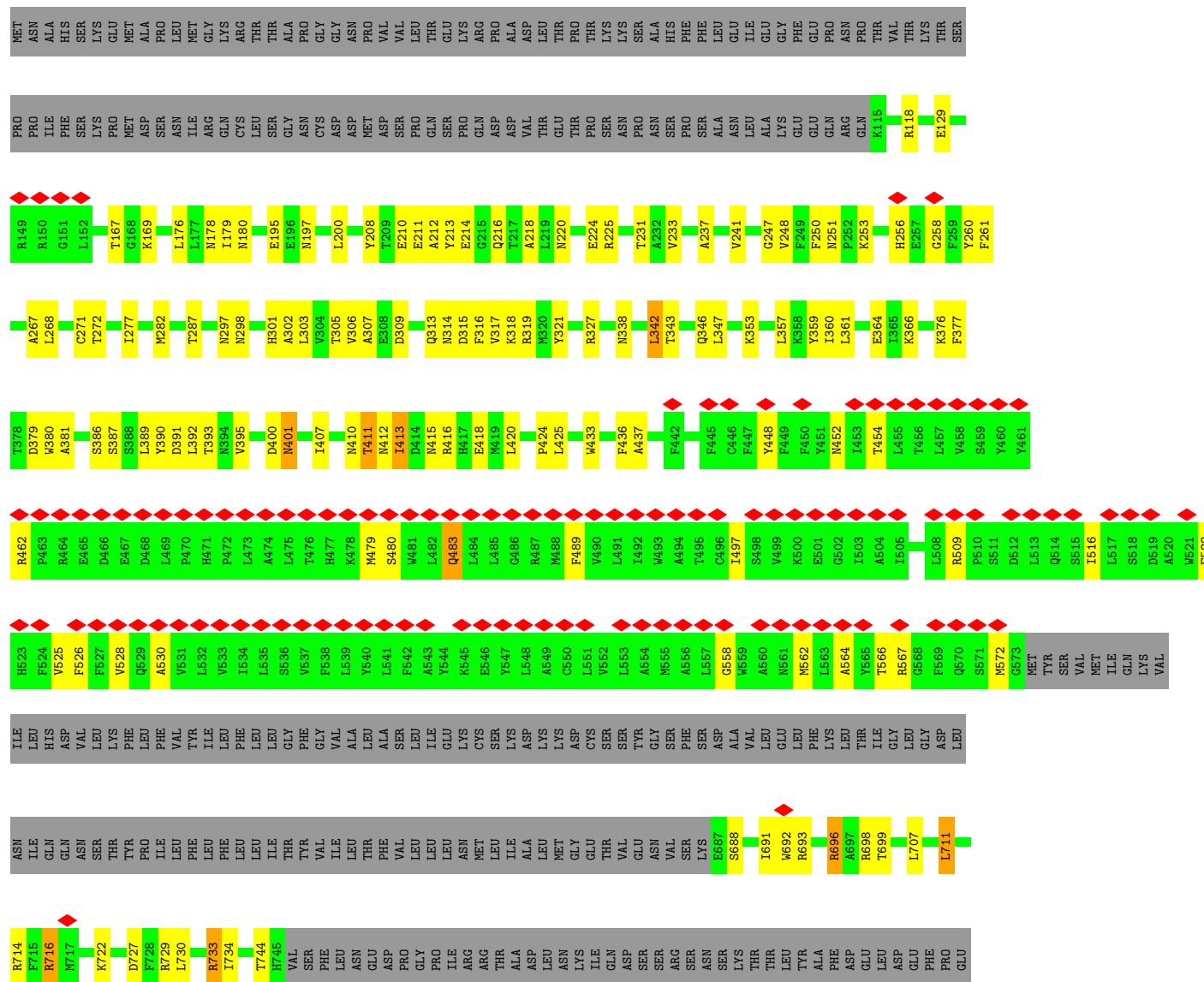
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Chain	Residue	Modelled	Actual	Comment	Reference
D	800	ALA	-	expression tag	UNP Q8K424
D	801	TRP	-	expression tag	UNP Q8K424
D	802	SER	-	expression tag	UNP Q8K424
D	803	HIS	-	expression tag	UNP Q8K424
D	804	PRO	-	expression tag	UNP Q8K424
D	805	GLN	-	expression tag	UNP Q8K424
D	806	PHE	-	expression tag	UNP Q8K424
D	807	GLU	-	expression tag	UNP Q8K424
D	808	LYS	-	expression tag	UNP Q8K424

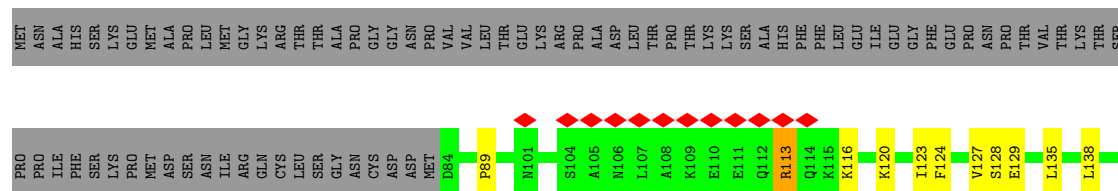
• Molecule 1: Transient receptor potential cation channel subfamily V member 3

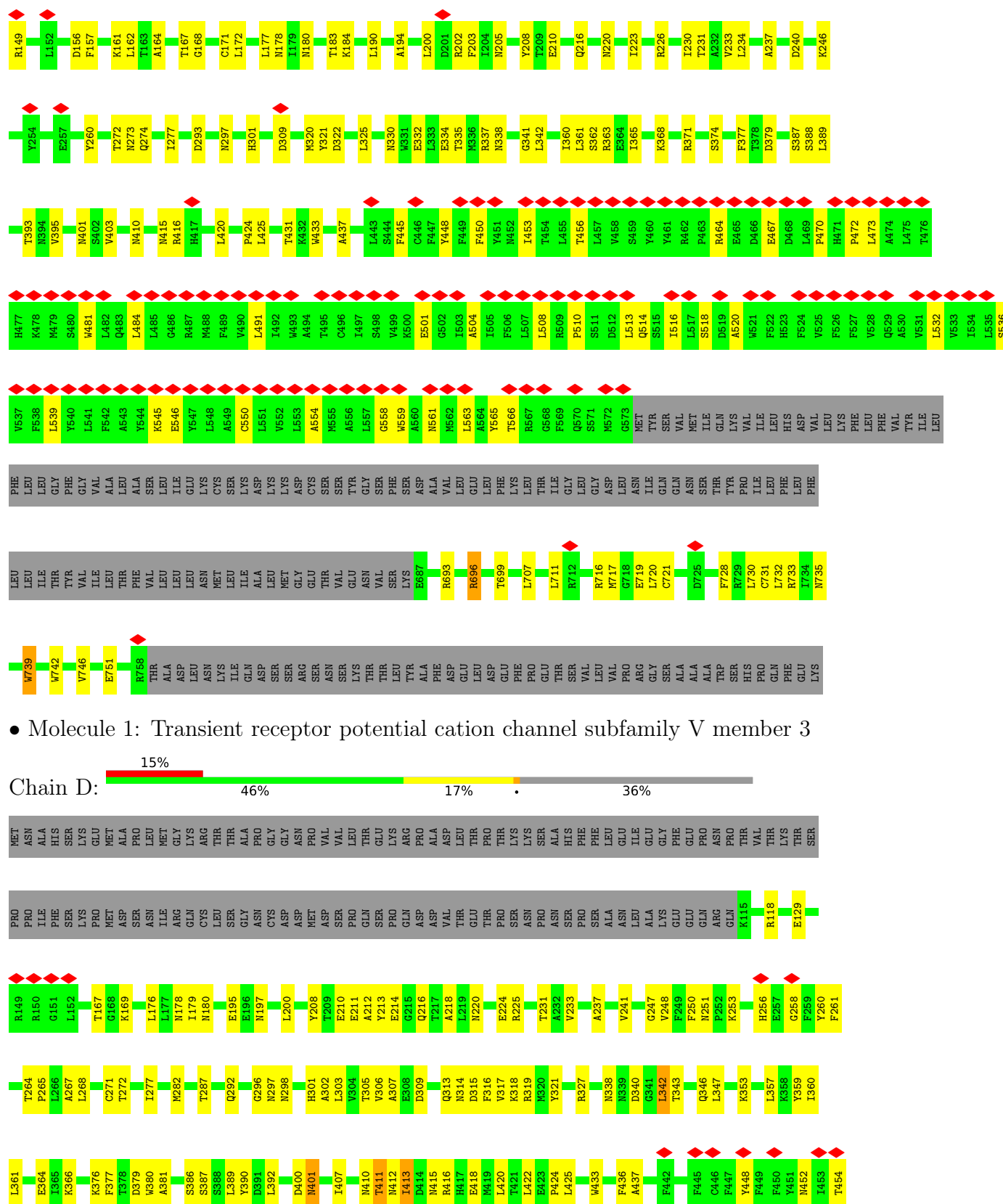
Chain B: 



• Molecule 1: Transient receptor potential cation channel subfamily V member 3

Chain C: 





GLU	LEU	ASP	PHE	PRO	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	IS16	L455
ASP	GLU	PHE	PRO	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	IS17	T456
GLU	GLU	PRO	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	SS18	L457
SER	SER	VAL	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	LYS	DS19	V458
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	AS20	S459
LEU	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	WS21	Y460
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	F522	Y461
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	HS23	P462
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	F524	R463
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	V525	R464
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	F526	E465
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	F527	D466
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	V528	E467
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	Q529	D468
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	AS30	L469
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	V531	P470
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	L532	H471
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	V533	P472
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	IS34	L473
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	L535	A474
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	SS36	L475
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	V537	T476
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	F538	H477
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	L539	K478
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	Y540	M479
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	F542	S480
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	AS43	M481
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	Y544	L482
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	K545	Q483
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	ES46	L484
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	Y547	L485
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	L548	G486
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	AS49	R487
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	C550	M488
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	L551	F489
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	V552	V490
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	L553	L491
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	AS54	I492
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	M555	M493
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	AS56	A494
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	L557	T495
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	C558	C496
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	M559	I497
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	AS60	S498
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	M561	V499
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	M562	X500
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	L563	E501
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	AS64	G502
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	Y565	I503
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	T566	A504
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	R567	
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	G568	I505
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	F569	
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	Q570	L508
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	SS71	R509
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	M572	P510
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	G573	SS11
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	MET	DS12
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE	TYR	L513
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE		Q514
VAL	VAL	LEU	THR	VAL	LEU	GLN	ASN	GLN	ASP	SER	VAL	ILE		SS15

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	101814	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 ($e^-/\text{\AA}^2$)	57.0	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	0.087	Depositor
Minimum map value	-0.052	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.005	Depositor
Recommended contour level	0.0176	Depositor
Map size (Å)	233.19998, 233.19998, 233.19998	wwPDB
Map dimensions	220, 220, 220	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.06, 1.06, 1.06	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.43	0/4682	0.66	2/6339 (0.0%)
1	B	0.42	0/4332	0.67	1/5859 (0.0%)
1	C	0.43	0/4682	0.66	2/6339 (0.0%)
1	D	0.42	0/4332	0.67	2/5859 (0.0%)
All	All	0.42	0/18028	0.67	7/24396 (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	5
1	B	0	3
1	C	0	2
1	D	0	3
All	All	0	13

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	190	LEU	CB-CG-CD2	-6.15	100.55	111.00
1	A	190	LEU	CB-CG-CD2	-6.10	100.63	111.00
1	C	513	LEU	CA-CB-CG	5.46	127.87	115.30
1	A	513	LEU	CA-CB-CG	5.46	127.85	115.30
1	D	711	LEU	CA-CB-CG	5.41	127.75	115.30

There are no chirality outliers.

5 of 13 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	293	ASP	Peptide
1	A	463	PRO	Peptide
1	A	465	GLU	Peptide
1	A	739	TRP	Peptide
1	A	89	PRO	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	4581	0	4575	86	0
1	B	4239	0	4261	84	0
1	C	4581	0	4575	91	0
1	D	4239	0	4261	84	0
All	All	17640	0	17672	337	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 10.

The worst 5 of 337 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:A:368:LYS:HA	1:A:371:ARG:HE	1.57	0.70
1:B:722:LYS:HD2	1:B:727:ASP:HB3	1.74	0.69
1:D:722:LYS:HD2	1:D:727:ASP:HB3	1.74	0.68
1:C:368:LYS:HA	1:C:371:ARG:HE	1.57	0.68
1:A:293:ASP:H	1:A:297:ASN:H	1.42	0.67

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	558/808 (69%)	475 (85%)	82 (15%)	1 (0%)	47	81
1	B	514/808 (64%)	438 (85%)	73 (14%)	3 (1%)	25	65
1	C	558/808 (69%)	476 (85%)	81 (14%)	1 (0%)	47	81
1	D	514/808 (64%)	438 (85%)	73 (14%)	3 (1%)	25	65
All	All	2144/3232 (66%)	1827 (85%)	309 (14%)	8 (0%)	38	72

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	412	ASN
1	D	412	ASN
1	A	113	ARG
1	C	113	ARG
1	B	411	THR

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	499/720 (69%)	491 (98%)	8 (2%)	62	79
1	B	458/720 (64%)	443 (97%)	15 (3%)	38	61
1	C	499/720 (69%)	491 (98%)	8 (2%)	62	79
1	D	458/720 (64%)	443 (97%)	15 (3%)	38	61
All	All	1914/2880 (66%)	1868 (98%)	46 (2%)	51	69

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

Mol	Chain	Res	Type
1	C	545	LYS
1	D	272	THR
1	C	696	ARG

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Mol	Chain	Res	Type
1	D	178	ASN
1	D	314	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 27 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	205	ASN
1	C	401	ASN
1	D	314	ASN
1	C	297	ASN
1	D	178	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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.

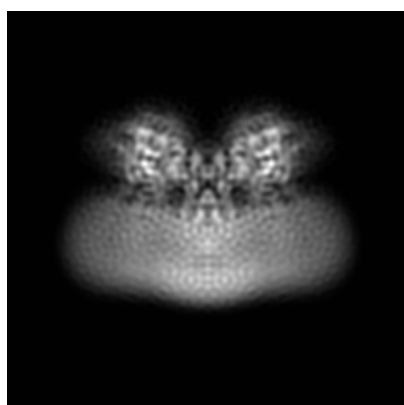
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20497. These allow visual inspection of the internal detail of the map and identification of artifacts.

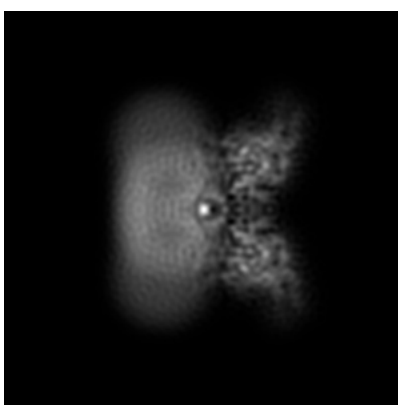
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

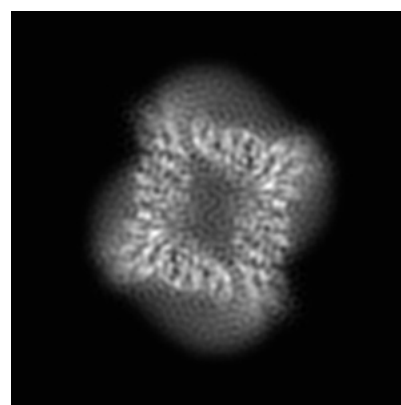
6.1.1 Primary map



X



Y

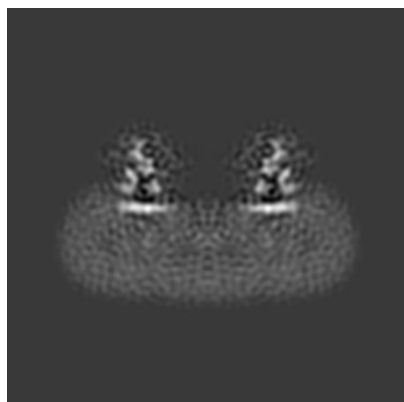


Z

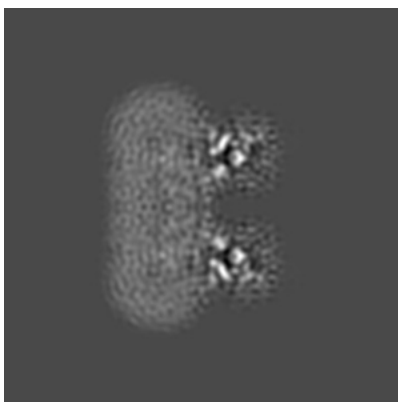
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

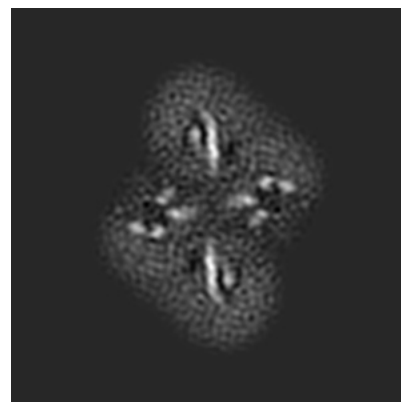
6.2.1 Primary map



X Index: 110



Y Index: 110

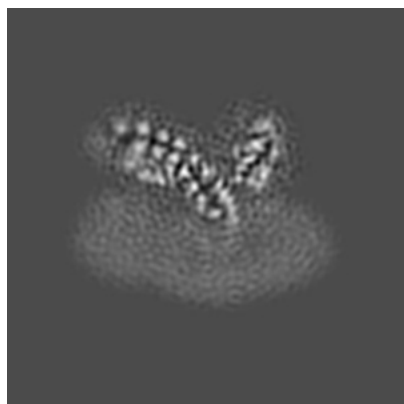


Z Index: 110

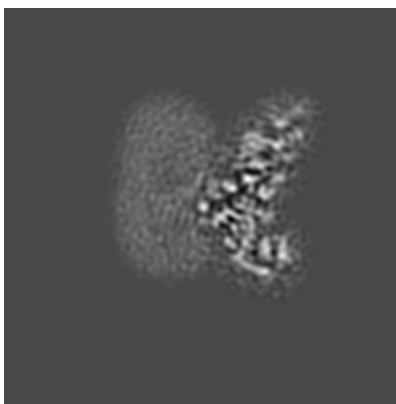
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

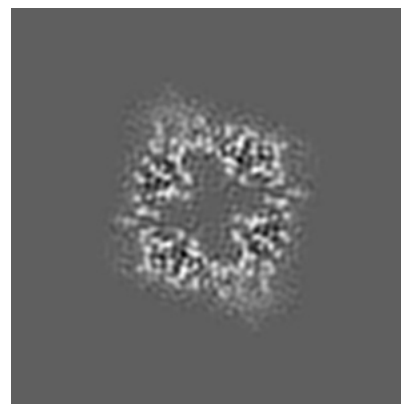
6.3.1 Primary map



X Index: 138



Y Index: 144



Z Index: 135

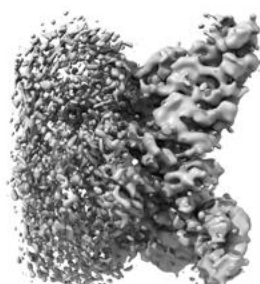
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views [i](#)

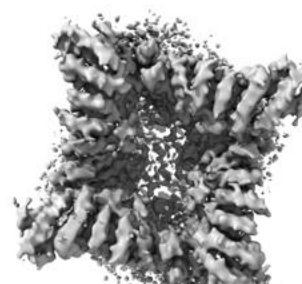
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0176. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

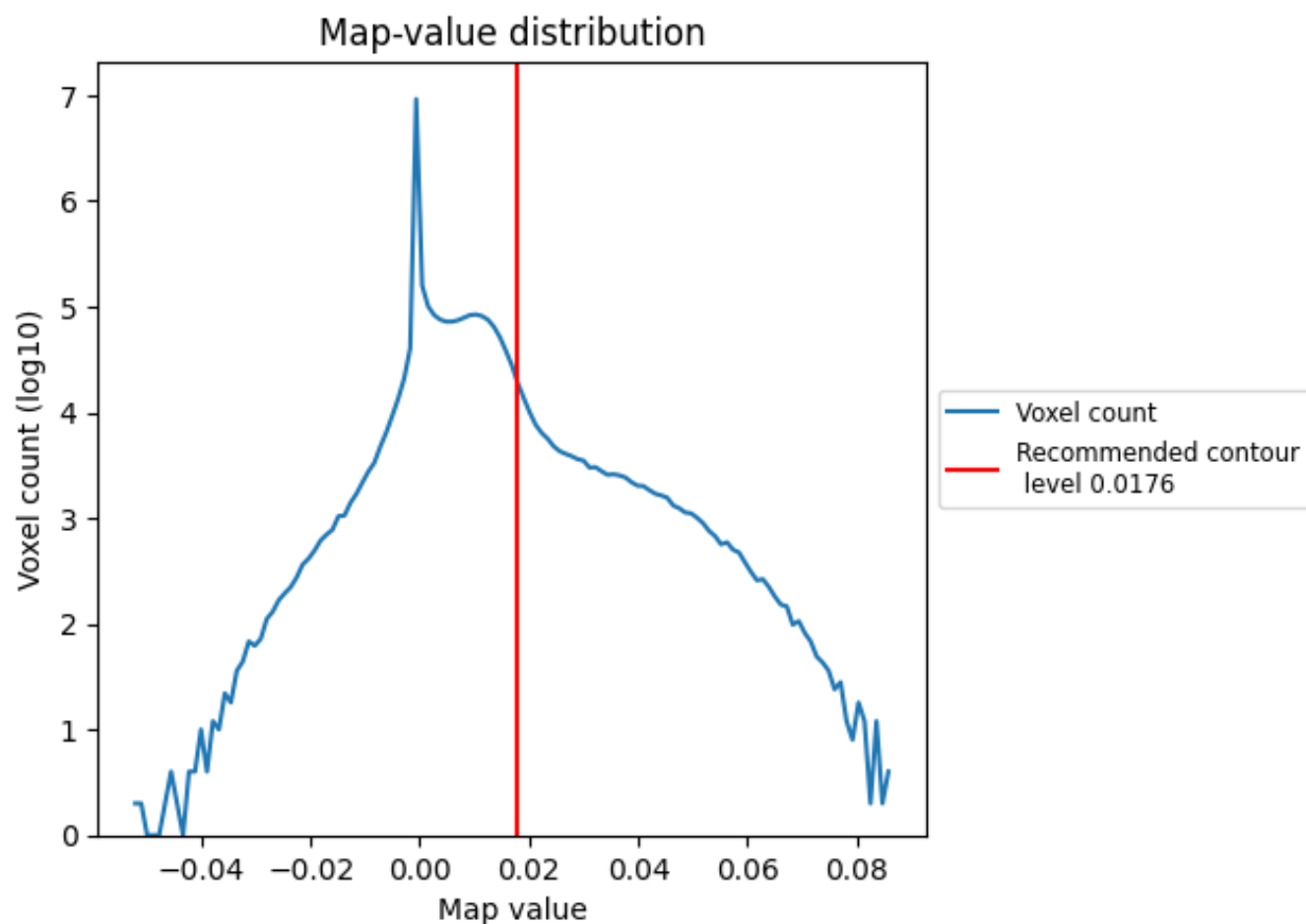
6.5 Mask visualisation

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

7 Map analysis [i](#)

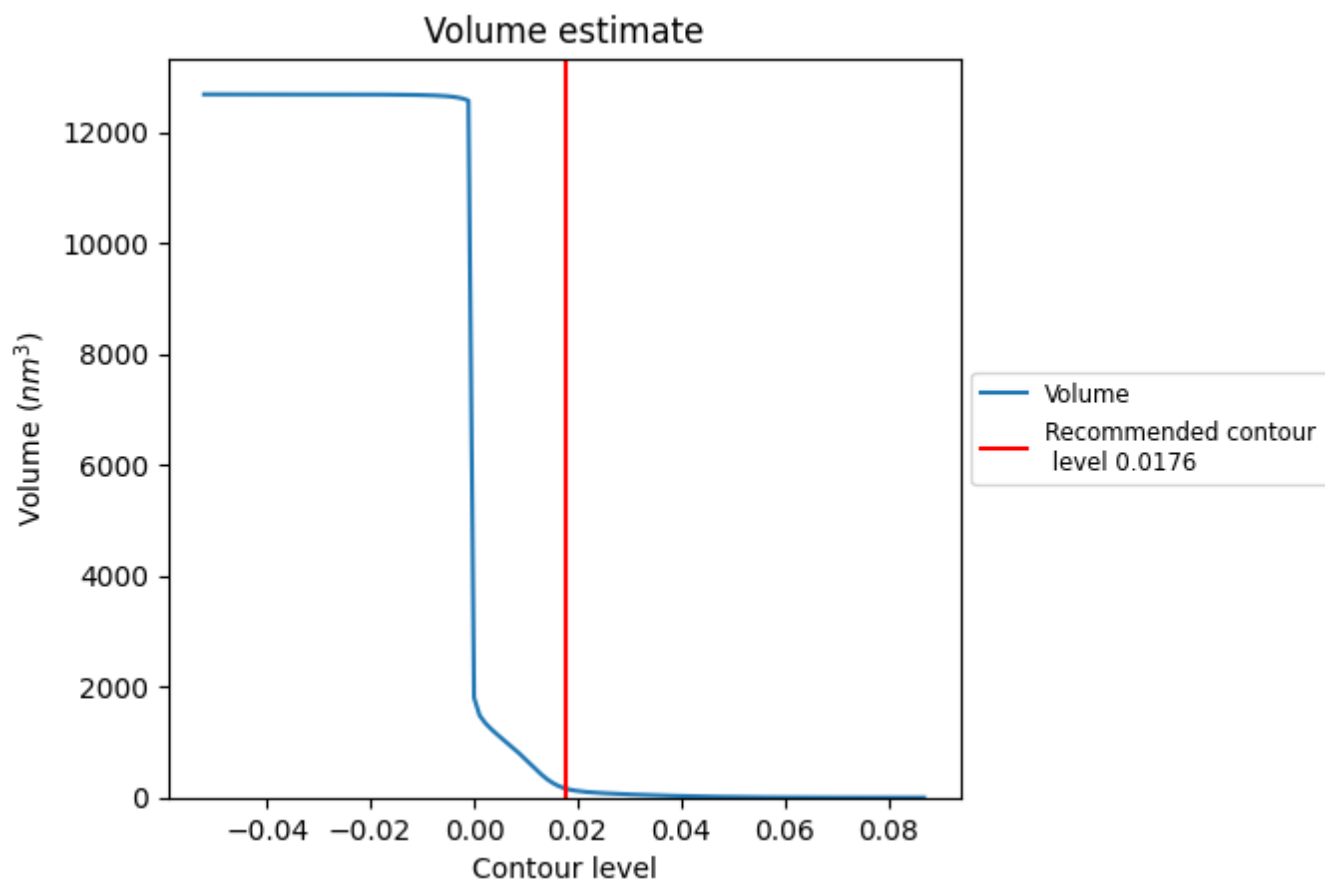
This section contains the results of statistical analysis of the map.

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.

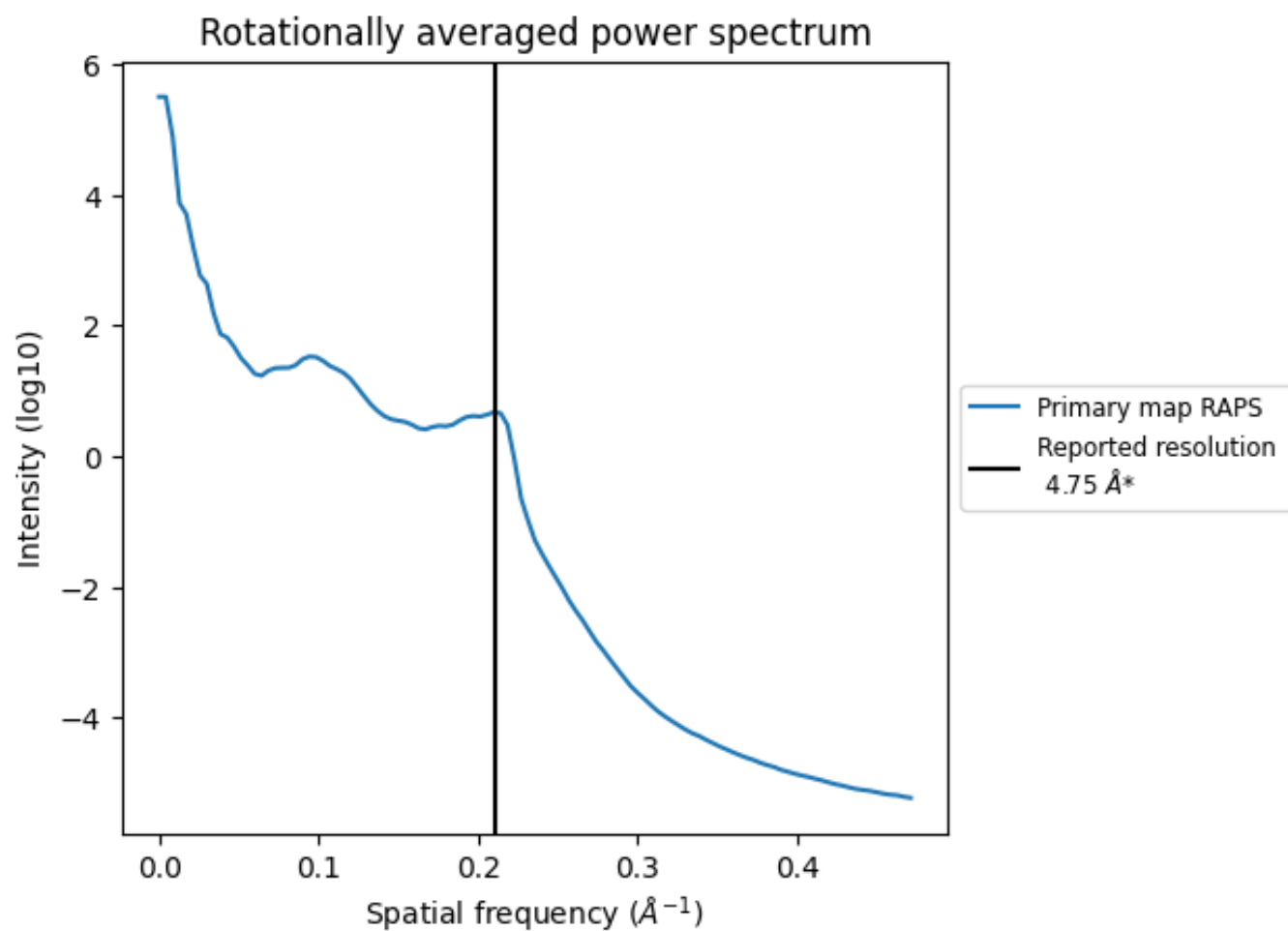
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 166 nm³; this corresponds to an approximate mass of 150 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.211 \AA^{-1}

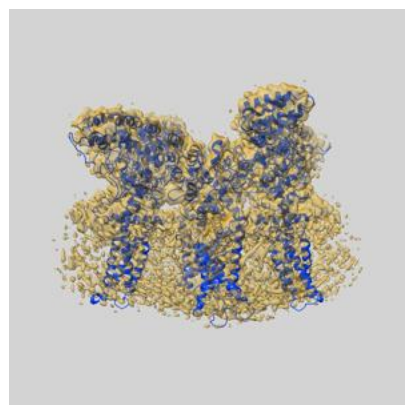
8 Fourier-Shell correlation

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

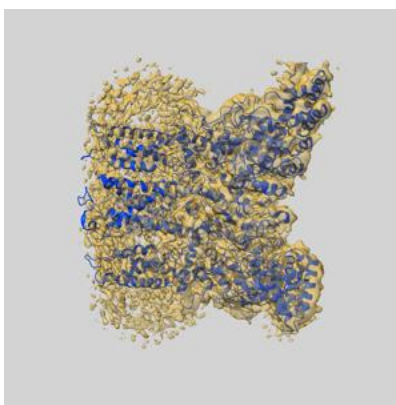
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-20497 and PDB model 6PVQ. Per-residue inclusion information can be found in section [3](#) on page [6](#).

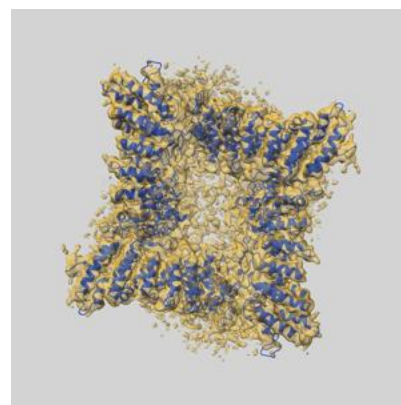
9.1 Map-model overlay [i](#)



X



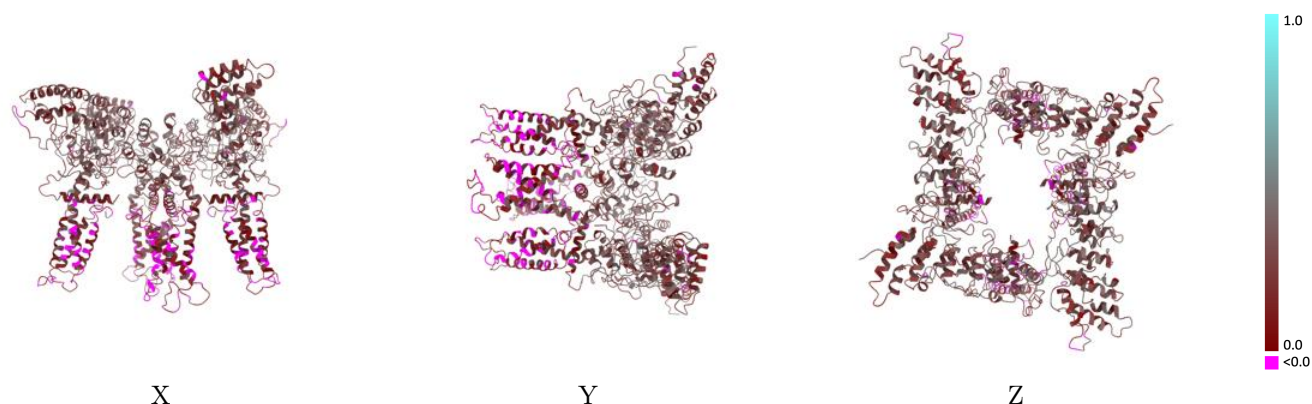
Y



Z

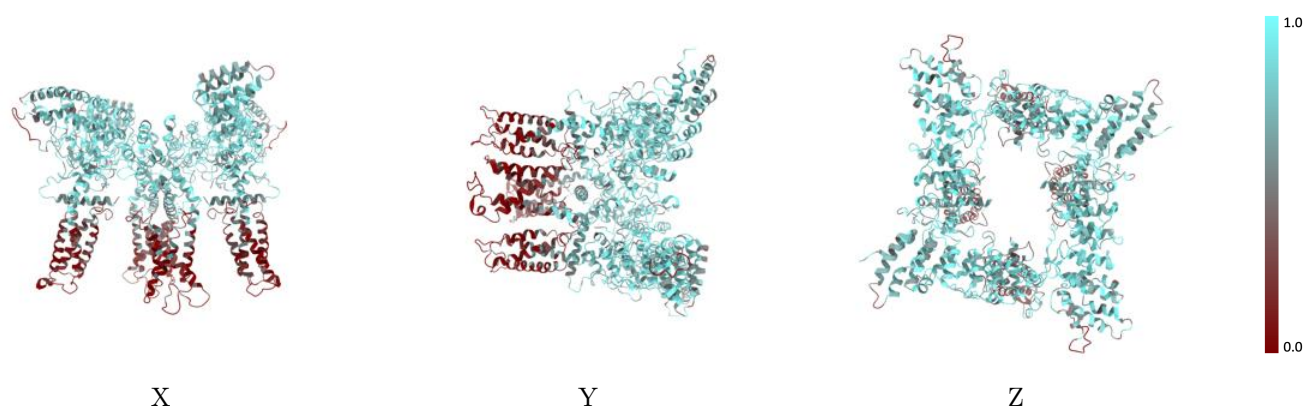
The images above show the 3D surface view of the map at the recommended contour level 0.0176 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



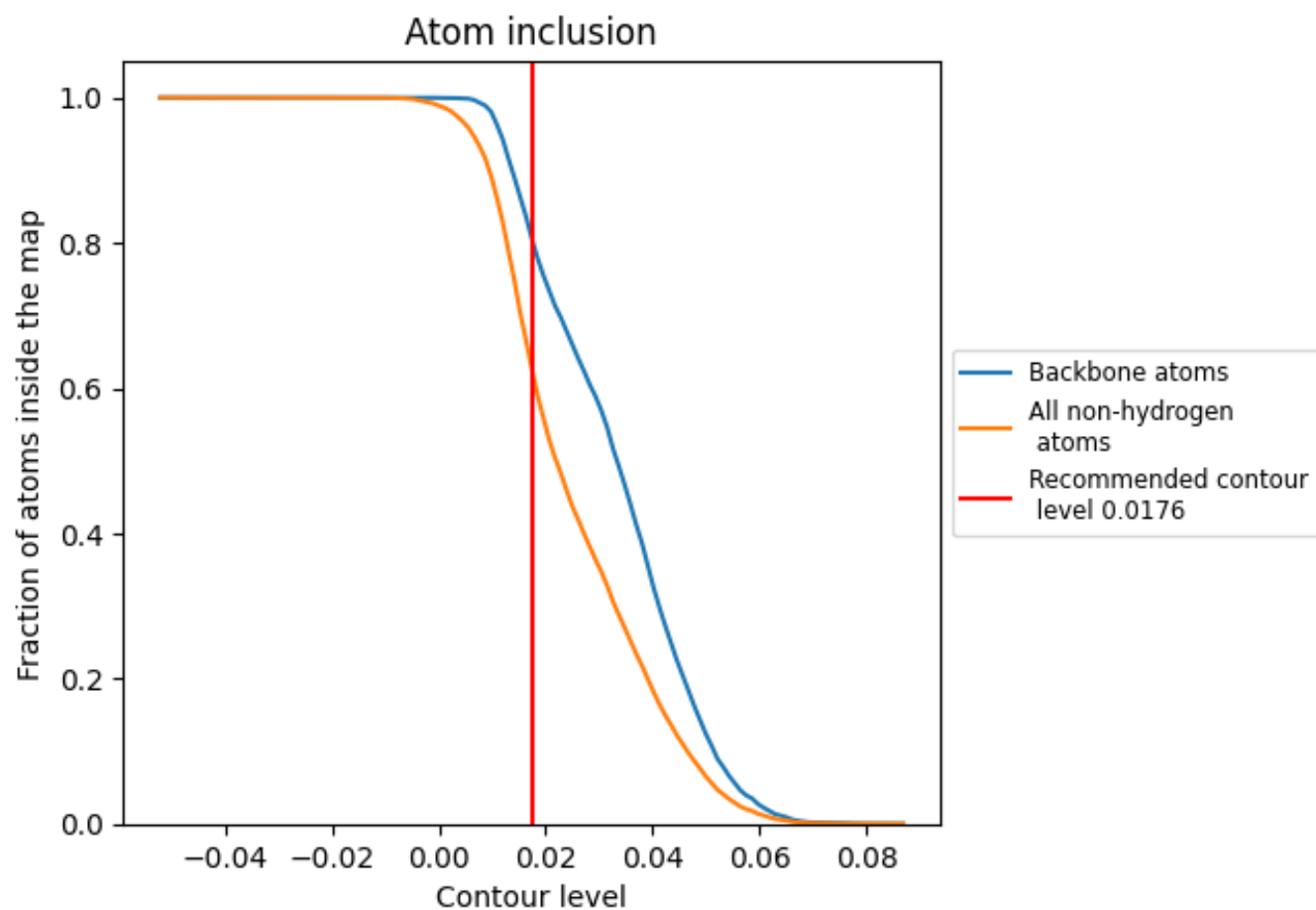
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0176).

9.4 Atom inclusion [i](#)



At the recommended contour level, 80% of all backbone atoms, 62% 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 (0.0176) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.6199	<div></div> 0.2400
A	<div></div> 0.6203	<div></div> 0.2420
B	<div></div> 0.6194	<div></div> 0.2380
C	<div></div> 0.6205	<div></div> 0.2430
D	<div></div> 0.6194	<div></div> 0.2370

