



Full wwPDB EM Validation Report ⓘ

Aug 11, 2022 – 10:11 am BST

PDB ID : 7PQT
EMDB ID : EMD-13604
Title : Apo human Kv3.1 cryo-EM structure
Authors : Botte, M.; Huber, S.; Bucher, D.; Klint, J.K.; Rodriguez, D.; Tagmose, L.;
Chami, M.; Cheng, R.; Hennig, M.; Abdul Rhaman, W.
Deposited on : 2021-09-20
Resolution : 2.65 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at <http://www.wwpdb.org/validation/2017/FAQs#types>.

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

EMDB validation analysis : 0.0.1.dev8
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.29

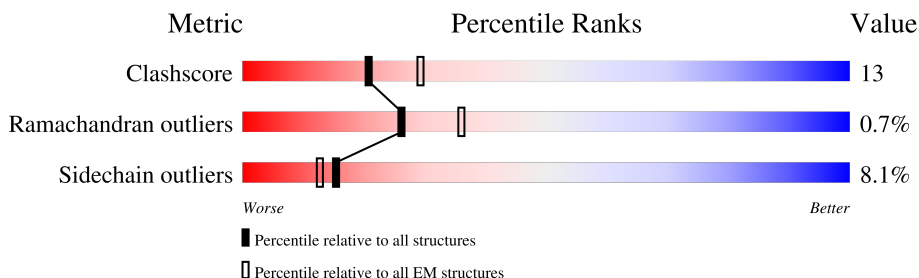
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.65 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	519	
1	B	519	
1	C	519	
1	D	519	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10464 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 Potassium voltage-gated channel subfamily C member 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	318	Total	C	N	O	S	4	0
			2613	1719	433	444	17		
1	B	318	Total	C	N	O	S	4	0
			2613	1719	433	444	17		
1	C	318	Total	C	N	O	S	4	0
			2613	1719	433	444	17		
1	D	318	Total	C	N	O	S	4	0
			2613	1719	433	444	17		

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	512	GLY	-	expression tag	UNP P48547
A	513	SER	-	expression tag	UNP P48547
A	514	LEU	-	expression tag	UNP P48547
A	515	GLU	-	expression tag	UNP P48547
A	516	VAL	-	expression tag	UNP P48547
A	517	LEU	-	expression tag	UNP P48547
A	518	PHE	-	expression tag	UNP P48547
A	519	GLN	-	expression tag	UNP P48547
B	512	GLY	-	expression tag	UNP P48547
B	513	SER	-	expression tag	UNP P48547
B	514	LEU	-	expression tag	UNP P48547
B	515	GLU	-	expression tag	UNP P48547
B	516	VAL	-	expression tag	UNP P48547
B	517	LEU	-	expression tag	UNP P48547
B	518	PHE	-	expression tag	UNP P48547
B	519	GLN	-	expression tag	UNP P48547
C	512	GLY	-	expression tag	UNP P48547
C	513	SER	-	expression tag	UNP P48547
C	514	LEU	-	expression tag	UNP P48547
C	515	GLU	-	expression tag	UNP P48547
C	516	VAL	-	expression tag	UNP P48547
C	517	LEU	-	expression tag	UNP P48547

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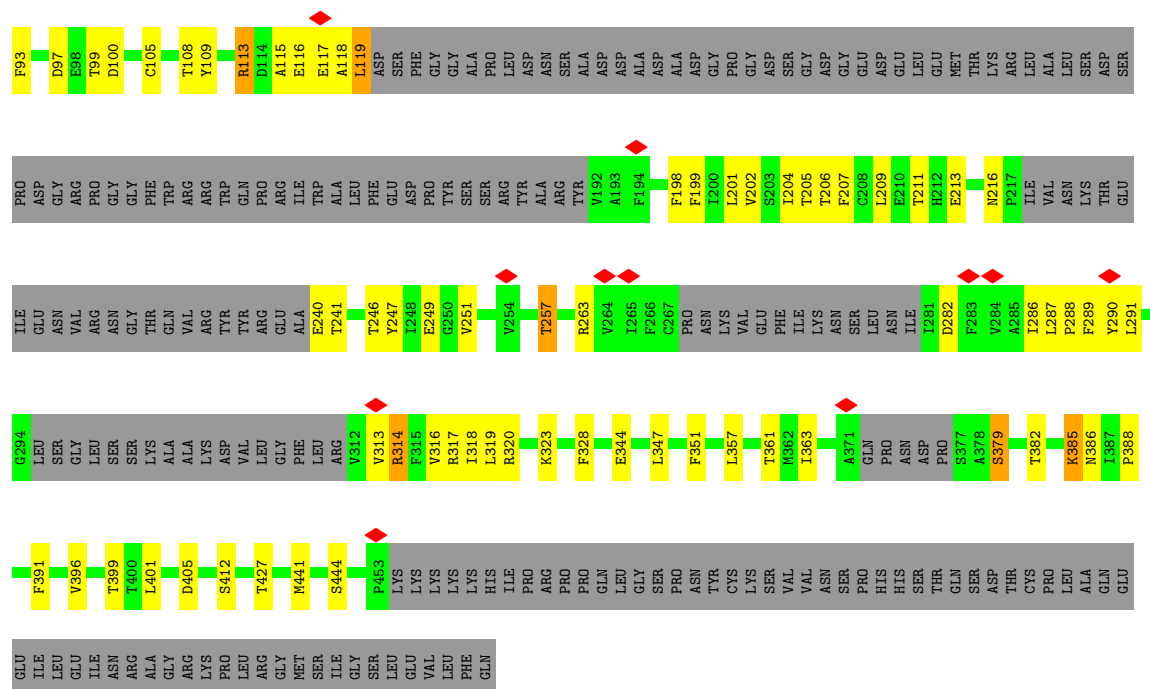
Chain	Residue	Modelled	Actual	Comment	Reference
C	518	PHE	-	expression tag	UNP P48547
C	519	GLN	-	expression tag	UNP P48547
D	512	GLY	-	expression tag	UNP P48547
D	513	SER	-	expression tag	UNP P48547
D	514	LEU	-	expression tag	UNP P48547
D	515	GLU	-	expression tag	UNP P48547
D	516	VAL	-	expression tag	UNP P48547
D	517	LEU	-	expression tag	UNP P48547
D	518	PHE	-	expression tag	UNP P48547
D	519	GLN	-	expression tag	UNP P48547

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K) (labeled as "Ligand of Interest" by depositor).

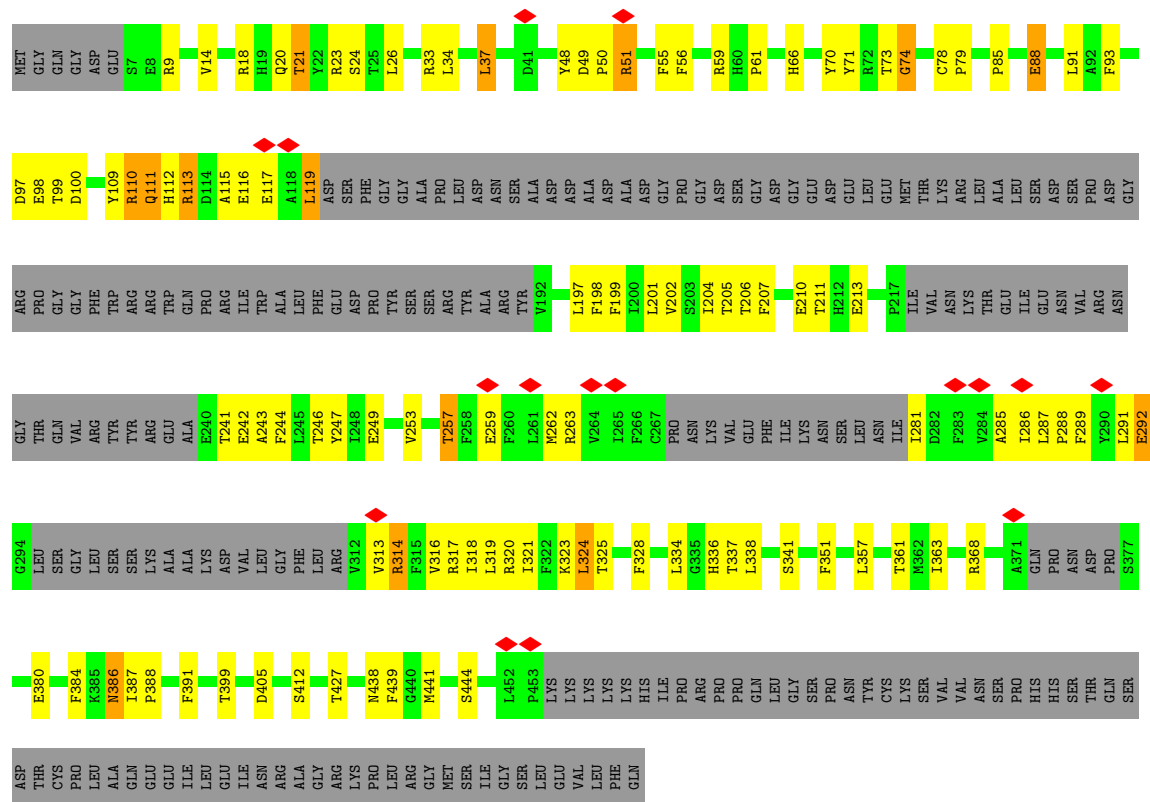
Mol	Chain	Residues	Atoms	AltConf
2	A	3	Total K 3 3	0
2	B	1	Total K 1 1	0

- Molecule 3 is water.

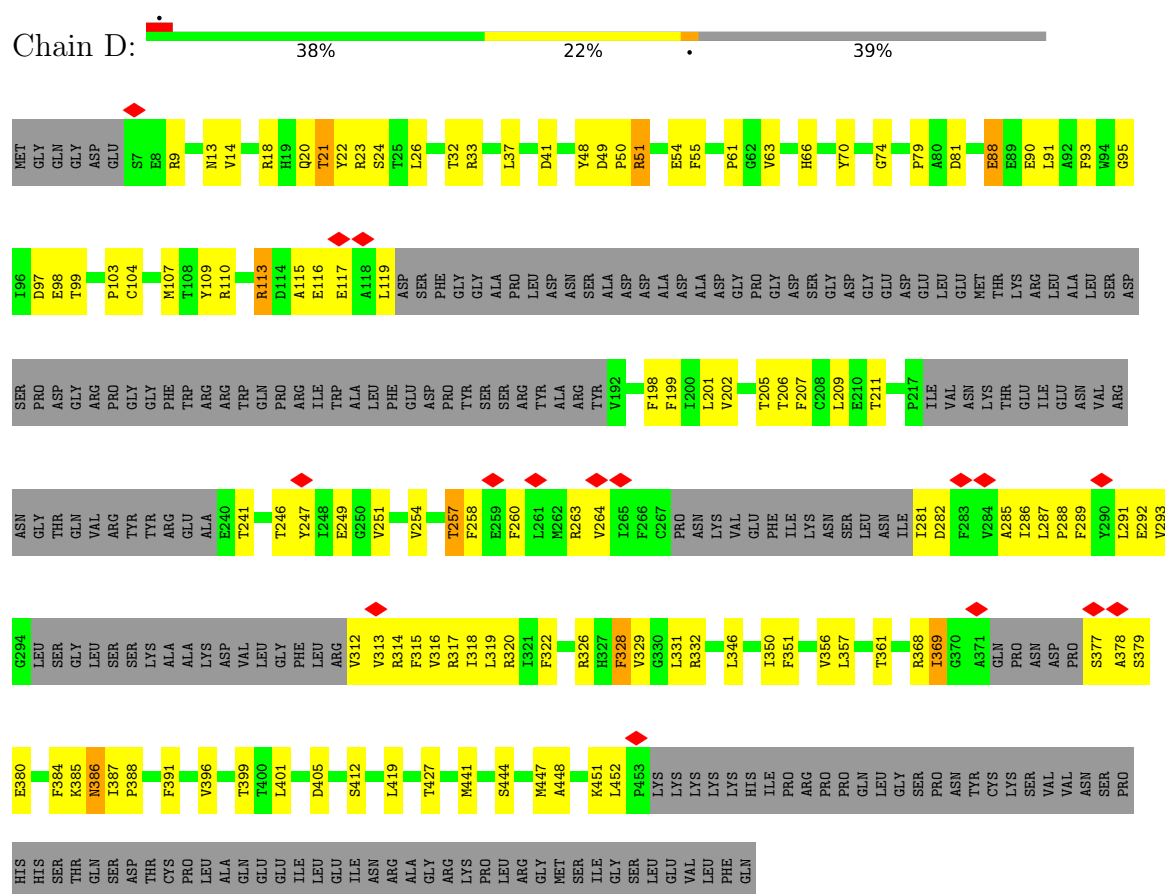
Mol	Chain	Residues	Atoms	AltConf
3	A	2	Total O 2 2	0
3	B	2	Total O 2 2	0
3	C	2	Total O 2 2	0
3	D	2	Total O 2 2	0



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



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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	362349	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 ($e^-/\text{\AA}^2$)	70	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.480	Depositor
Minimum map value	-1.277	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.040	Depositor
Recommended contour level	0.356	Depositor
Map size (\AA)	287.0, 287.0, 287.0	wwPDB
Map dimensions	350, 350, 350	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.82, 0.82, 0.82	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

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	0.56	0/2688	0.53	0/3652
1	B	0.57	0/2688	0.54	0/3652
1	C	0.56	0/2688	0.54	0/3652
1	D	0.57	0/2688	0.53	0/3652
All	All	0.56	0/10752	0.53	0/14608

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	2613	0	2574	64	0
1	B	2613	0	2574	70	0
1	C	2613	0	2574	76	0
1	D	2613	0	2574	76	0
2	A	3	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	2	0	0	2	0
All	All	10464	0	10296	278	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 13.

All (278) 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:D:107:MET:SD	3:D:601:HOH:O	2.36	0.83
1:D:282:ASP:HA	1:D:286:ILE:HD12	1.63	0.80
1:C:317:ARG:HG3	1:C:320:ARG:HH21	1.53	0.73
1:B:291:LEU:HD21	1:B:313:VAL:HA	1.69	0.72
1:D:313:VAL:HG12	1:D:314:ARG:HG3	1.72	0.72
1:B:33:ARG:NH2	1:B:93:PHE:O	2.19	0.72
1:C:33:ARG:NH2	1:C:93:PHE:O	2.21	0.71
1:C:49:ASP:HB3	1:C:51:ARG:NH1	2.06	0.71
1:C:386:ASN:HB2	1:C:388:PRO:HD2	1.71	0.71
1:D:386:ASN:HB2	1:D:388:PRO:HD2	1.72	0.70
1:A:8:GLU:OE1	1:A:23:ARG:NH2	2.22	0.70
1:A:205:THR:O	1:A:209:LEU:HD12	1.92	0.69
1:A:33:ARG:NH2	1:A:93:PHE:O	2.25	0.67
1:D:49:ASP:HB3	1:D:51:ARG:NH1	2.09	0.67
1:C:88:GLU:HA	1:C:91:LEU:HD12	1.75	0.67
1:A:398:MET:HE3	1:A:424:GLY:HA2	1.77	0.66
1:B:109:TYR:O	1:B:113[A]:ARG:NH2	2.28	0.66
1:D:97:ASP:OD2	1:D:99:THR:OG1	2.13	0.66
1:D:368[B]:ARG:NH1	1:D:379:SER:OG	2.29	0.66
1:A:316:VAL:HA	1:A:319:LEU:HD23	1.78	0.65
1:B:257:THR:HG1	1:B:289:PHE:HZ	1.44	0.65
1:C:291:LEU:HD21	1:C:313:VAL:HA	1.77	0.65
1:A:341:SER:OG	1:A:438:ASN:ND2	2.29	0.65
1:D:107:MET:SD	1:D:107:MET:N	2.66	0.65
1:C:34:LEU:HD12	1:C:34:LEU:H	1.63	0.64
1:D:116:GLU:O	1:D:119:LEU:N	2.30	0.64
1:A:212:HIS:O	1:A:216:ASN:ND2	2.31	0.64
1:A:318:ILE:HD13	1:B:361:THR:HG21	1.80	0.64
1:D:33:ARG:NH2	1:D:93:PHE:O	2.26	0.64
1:D:292:GLU:OE1	1:D:293:VAL:HG23	1.98	0.64
1:C:49:ASP:HB3	1:C:51:ARG:HH12	1.63	0.64
1:B:316:VAL:HA	1:B:319:LEU:HD23	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:THR:O	1:B:209:LEU:HD12	1.97	0.63
1:B:34:LEU:H	1:B:34:LEU:HD12	1.63	0.63
1:A:213:GLU:HA	1:A:216:ASN:HD22	1.63	0.63
1:B:116:GLU:O	1:B:119:LEU:N	2.32	0.63
1:B:97:ASP:OD1	1:B:99:THR:OG1	2.17	0.62
1:C:201:LEU:O	1:C:205:THR:HG23	1.98	0.62
1:C:116:GLU:O	1:C:119:LEU:N	2.33	0.62
1:A:37:LEU:HB3	1:A:55:PHE:HE2	1.65	0.61
1:C:288:PRO:HA	1:C:291:LEU:HB3	1.82	0.61
1:B:201:LEU:O	1:B:205:THR:HG23	2.01	0.61
1:D:88:GLU:HA	1:D:91:LEU:HD12	1.81	0.60
1:A:109:TYR:O	1:A:113[A]:ARG:NH2	2.34	0.60
1:B:291:LEU:HD11	1:B:313:VAL:HG22	1.83	0.60
1:C:109:TYR:O	1:C:113[A]:ARG:NH2	2.35	0.60
1:D:207:PHE:HB2	1:D:317:ARG:HH21	1.67	0.59
1:C:97:ASP:OD1	1:C:99:THR:OG1	2.17	0.59
1:C:242:GLU:HG3	1:C:243:ALA:H	1.66	0.59
1:D:49:ASP:HB3	1:D:51:ARG:HH12	1.65	0.59
1:D:81:ASP:OD1	1:D:81:ASP:N	2.36	0.59
1:B:204:ILE:HD12	1:B:318:ILE:HD13	1.83	0.59
1:D:205:THR:O	1:D:209:LEU:HD12	2.03	0.59
1:A:97:ASP:OD2	1:A:99:THR:OG1	2.21	0.58
1:A:116:GLU:O	1:A:119:LEU:N	2.35	0.58
1:A:48:TYR:CD1	1:A:55:PHE:HE1	2.22	0.58
1:A:288:PRO:HA	1:A:291:LEU:HG	1.84	0.58
1:D:357:LEU:O	1:D:361:THR:HG22	2.04	0.57
1:D:314:ARG:HA	1:D:317:ARG:HG3	1.86	0.57
1:C:48:TYR:CD1	1:C:55:PHE:HE1	2.22	0.57
1:C:341:SER:OG	1:C:438:ASN:ND2	2.37	0.57
1:D:246:THR:HA	1:D:249:GLU:OE1	2.05	0.57
1:D:316:VAL:HA	1:D:319:LEU:HD23	1.87	0.56
1:B:12:ILE:HG22	1:B:14:VAL:HG23	1.86	0.56
1:C:316:VAL:HA	1:C:319:LEU:HD23	1.87	0.56
1:D:109:TYR:O	1:D:113[A]:ARG:NH2	2.38	0.56
1:B:37:LEU:HD22	1:B:55:PHE:CD2	2.41	0.56
1:C:318:ILE:HG12	1:C:319:LEU:HD22	1.88	0.55
1:A:287:LEU:HA	1:A:290:TYR:HD2	1.69	0.55
1:B:48:TYR:CD1	1:B:55:PHE:HE1	2.24	0.55
1:B:351:PHE:HB3	1:B:427:THR:HG23	1.87	0.55
1:D:48:TYR:HD1	1:D:55:PHE:CE1	2.25	0.55
1:A:387:ILE:HG13	1:D:211:THR:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:THR:O	1:A:385:LYS:HE2	2.07	0.54
1:B:8:GLU:OE1	1:B:23:ARG:NH2	2.35	0.54
1:D:260:PHE:HA	1:D:263:ARG:HD3	1.88	0.54
1:C:253:VAL:HG21	1:C:292:GLU:HG2	1.88	0.54
1:A:351:PHE:HB3	1:A:427:THR:HG23	1.89	0.54
1:D:351:PHE:HB3	1:D:427:THR:HG23	1.88	0.54
1:C:246:THR:HA	1:C:249:GLU:OE1	2.08	0.54
1:B:344:GLU:OE1	1:B:344:GLU:N	2.40	0.54
1:C:351:PHE:HB3	1:C:427:THR:HG23	1.88	0.54
1:B:39:GLU:OE1	1:B:39:GLU:N	2.39	0.53
1:C:98:GLU:OE2	1:C:110:ARG:NH1	2.40	0.53
1:B:246:THR:HA	1:B:249:GLU:OE1	2.09	0.53
1:A:37:LEU:HD22	1:A:55:PHE:CD2	2.44	0.53
1:C:37:LEU:HD22	1:C:55:PHE:CD2	2.44	0.53
1:A:287:LEU:N	1:A:288:PRO:HD2	2.24	0.52
1:B:382:THR:O	1:B:385:LYS:HE2	2.08	0.52
1:C:388:PRO:HA	1:C:391:PHE:CE2	2.44	0.52
1:D:388:PRO:HA	1:D:391:PHE:CE2	2.44	0.52
1:C:207:PHE:HE1	1:C:314:ARG:HG3	1.74	0.52
1:D:37:LEU:HD22	1:D:55:PHE:CD2	2.44	0.52
1:C:337:THR:HG21	1:C:439:PHE:HA	1.91	0.52
1:D:98:GLU:OE1	1:D:109:TYR:OH	2.20	0.52
1:D:356:VAL:HG12	1:D:391:PHE:HD1	1.75	0.52
1:A:333:VAL:O	1:A:337:THR:HG23	2.10	0.52
1:B:202:VAL:O	1:B:206:THR:HG22	2.10	0.51
1:B:287:LEU:N	1:B:288:PRO:HD2	2.26	0.51
1:C:37:LEU:HB3	1:C:55:PHE:HE2	1.74	0.51
1:C:48:TYR:HD1	1:C:55:PHE:CE1	2.28	0.51
1:D:247:TYR:O	1:D:251:VAL:HG23	2.10	0.51
1:A:216:ASN:HA	1:A:240:GLU:O	2.11	0.51
1:D:21:THR:HG22	1:D:22:TYR:H	1.74	0.51
1:D:377:SER:OG	1:D:380:GLU:OE2	2.28	0.51
1:A:326:ARG:HG2	1:A:326:ARG:HH11	1.76	0.50
1:D:63:VAL:HG21	1:D:90:GLU:HG3	1.93	0.50
1:B:386:ASN:HB2	1:B:388:PRO:HD2	1.93	0.50
1:C:207:PHE:CE1	1:C:314:ARG:HG3	2.46	0.50
1:C:357:LEU:O	1:C:361:THR:HG22	2.12	0.50
1:B:14:VAL:O	1:B:61:PRO:HB3	2.12	0.50
1:C:73:THR:OG1	1:C:74:GLY:N	2.44	0.49
1:D:377:SER:OG	1:D:378:ALA:N	2.45	0.49
1:C:324:LEU:HG	1:C:325:THR:HG23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:ARG:HH11	1:B:263:ARG:HG3	1.77	0.49
1:C:286:ILE:HG22	1:C:286:ILE:O	2.13	0.49
1:D:198:PHE:HD2	1:D:199:PHE:CD1	2.29	0.49
1:A:246:THR:HA	1:A:249:GLU:OE1	2.12	0.49
1:D:202:VAL:O	1:D:206:THR:HG22	2.13	0.49
1:A:323:LYS:O	1:A:326:ARG:NH1	2.44	0.49
1:A:115:ALA:HB2	1:A:441:MET:SD	2.52	0.49
1:A:48:TYR:HD1	1:A:55:PHE:CE1	2.31	0.49
1:B:379:SER:O	1:B:379:SER:OG	2.29	0.49
1:C:287:LEU:N	1:C:288:PRO:HD2	2.28	0.49
1:D:257:THR:OG1	1:D:289:PHE:HZ	1.96	0.48
1:D:448:ALA:O	1:D:452:LEU:HD23	2.13	0.48
1:D:314:ARG:CA	1:D:317:ARG:HG3	2.43	0.48
1:A:263:ARG:HD3	1:A:282:ASP:OD1	2.12	0.48
1:B:286:ILE:HD11	1:B:323:LYS:NZ	2.27	0.48
1:C:100:ASP:N	1:C:100:ASP:OD1	2.47	0.48
1:B:37:LEU:HD22	1:B:55:PHE:HD2	1.79	0.48
1:D:48:TYR:CD1	1:D:55:PHE:HE1	2.32	0.48
1:C:110:ARG:HH21	1:C:113[A]:ARG:HB3	1.77	0.48
1:B:37:LEU:HB3	1:B:55:PHE:HE2	1.79	0.48
1:C:56:PHE:CE2	1:D:9:ARG:HD3	2.48	0.48
1:C:48:TYR:CD1	1:C:55:PHE:CE1	3.02	0.48
1:D:312:VAL:HA	1:D:315:PHE:HB3	1.95	0.47
1:A:318:ILE:O	1:A:321:ILE:HG12	2.14	0.47
1:C:119:LEU:HG	1:C:336:HIS:HB3	1.96	0.47
1:D:48:TYR:CE2	1:D:50:PRO:HB3	2.49	0.47
1:C:14:VAL:O	1:C:61:PRO:HB3	2.15	0.47
1:D:291:LEU:HD11	1:D:313:VAL:HA	1.96	0.47
1:A:337:THR:HG21	1:A:439:PHE:HA	1.97	0.47
1:C:202:VAL:O	1:C:206:THR:HG22	2.13	0.47
1:D:405:ASP:OD1	1:D:405:ASP:N	2.48	0.47
1:A:14:VAL:O	1:A:61:PRO:HB3	2.14	0.47
1:B:9:ARG:HG2	1:B:21:THR:O	2.14	0.47
1:D:113[B]:ARG:HA	1:D:113[B]:ARG:HD3	1.77	0.47
1:C:59:ARG:HE	1:C:59:ARG:HB3	1.49	0.47
1:C:211:THR:HG21	1:D:387:ILE:HG13	1.95	0.47
1:D:287:LEU:N	1:D:288:PRO:HD2	2.29	0.47
1:C:368[B]:ARG:NH2	1:C:384:PHE:O	2.45	0.47
1:B:66:HIS:CG	1:B:79:PRO:HG3	2.50	0.46
1:A:388:PRO:HA	1:A:391:PHE:CE2	2.50	0.46
1:B:73:THR:OG1	1:B:74:GLY:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:THR:HG21	1:C:387:ILE:HG13	1.97	0.46
1:A:37:LEU:HB3	1:A:55:PHE:CE2	2.49	0.46
1:C:66:HIS:CG	1:C:79:PRO:HG3	2.50	0.46
1:D:48:TYR:CD1	1:D:55:PHE:CE1	3.04	0.46
1:A:66:HIS:CG	1:A:79:PRO:HG3	2.50	0.46
1:A:88:GLU:HA	1:A:91:LEU:HD12	1.96	0.46
1:D:346:LEU:O	1:D:350:ILE:HG13	2.16	0.46
1:B:388:PRO:HA	1:B:391:PHE:CE2	2.51	0.46
1:D:14:VAL:O	1:D:61:PRO:HB3	2.15	0.46
1:A:207:PHE:CE1	1:A:314:ARG:HG3	2.50	0.46
1:B:9:ARG:HD2	1:B:20:GLN:OE1	2.15	0.46
1:A:346:LEU:HD23	1:A:346:LEU:HA	1.83	0.46
1:C:9:ARG:HG2	1:C:21:THR:O	2.16	0.46
1:A:328:PHE:HB3	1:A:331:LEU:HG	1.97	0.45
1:C:287:LEU:HG	1:C:288:PRO:HD3	1.97	0.45
1:C:405:ASP:OD1	1:C:405:ASP:N	2.49	0.45
1:A:199:PHE:CD2	1:A:252:CYS:HA	2.51	0.45
1:C:56:PHE:CD2	1:D:9:ARG:HD3	2.51	0.45
1:B:88:GLU:HA	1:B:91:LEU:HD12	1.98	0.45
1:C:198:PHE:HD2	1:C:199:PHE:CD1	2.35	0.45
1:D:198:PHE:HD2	1:D:199:PHE:HD1	1.65	0.45
1:A:81:ASP:OD1	1:A:81:ASP:N	2.49	0.45
1:B:115:ALA:HB2	1:B:441:MET:SD	2.57	0.45
1:B:216:ASN:HA	1:B:240:GLU:O	2.17	0.45
1:A:48:TYR:CD1	1:A:55:PHE:CE1	3.03	0.45
1:A:9:ARG:HG2	1:A:21:THR:O	2.17	0.45
1:A:323:LYS:HB2	1:A:323:LYS:HE2	1.72	0.45
1:C:9:ARG:HD2	1:C:20:GLN:OE1	2.16	0.44
1:D:328:PHE:HB3	1:D:331:LEU:HG	2.00	0.44
1:A:59:ARG:HE	1:A:59:ARG:HB3	1.51	0.44
1:A:253:VAL:HG21	1:A:292:GLU:HG3	1.98	0.44
1:A:344:GLU:OE1	1:A:344:GLU:N	2.51	0.44
1:B:247:TYR:O	1:B:251:VAL:HG13	2.18	0.44
1:C:111:GLN:HG3	1:C:112:HIS:N	2.33	0.44
1:D:26:LEU:HD23	1:D:26:LEU:HA	1.80	0.44
1:D:328:PHE:O	1:D:332:ARG:HG3	2.18	0.44
1:C:51:ARG:HD3	1:C:51:ARG:N	2.33	0.44
1:A:21:THR:HG22	1:A:22:TYR:H	1.83	0.44
1:B:263:ARG:HD3	1:B:282:ASP:OD1	2.17	0.44
1:D:37:LEU:HD22	1:D:55:PHE:HD2	1.83	0.44
1:D:66:HIS:CG	1:D:79:PRO:HG3	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:281:ILE:O	1:C:285:ALA:HB3	2.18	0.43
1:B:207:PHE:CE1	1:B:314:ARG:HB2	2.53	0.43
1:D:13:ASN:ND2	1:D:18:ARG:HE	2.15	0.43
1:A:63:VAL:HG21	1:A:90:GLU:HG3	2.00	0.43
1:A:396:VAL:HG13	1:A:401:LEU:HD23	2.01	0.43
1:B:59:ARG:HE	1:B:59:ARG:HB3	1.52	0.43
1:C:34:LEU:HD13	1:C:71:TYR:OH	2.17	0.43
1:C:198:PHE:HD2	1:C:199:PHE:HD1	1.65	0.43
1:C:213:GLU:OE1	1:D:386:ASN:ND2	2.51	0.43
1:D:207:PHE:CE1	1:D:314:ARG:HB2	2.53	0.43
1:A:253:VAL:HG13	1:A:289:PHE:HE1	1.83	0.43
1:A:284:VAL:HG12	1:A:284:VAL:O	2.18	0.43
1:B:48:TYR:OH	1:B:53:ASP:OD1	2.31	0.43
1:C:257:THR:OG1	1:C:289:PHE:HZ	2.01	0.43
1:D:9:ARG:HD2	1:D:20:GLN:OE1	2.18	0.43
1:B:68:LEU:HD12	1:B:68:LEU:HA	1.83	0.43
1:D:115:ALA:HB2	1:D:441:MET:SD	2.58	0.43
1:D:419:LEU:HD23	1:D:419:LEU:HA	1.87	0.43
1:B:34:LEU:HD13	1:B:71:TYR:OH	2.19	0.43
1:B:405:ASP:OD1	1:B:405:ASP:N	2.47	0.43
1:C:48:TYR:CE2	1:C:50:PRO:HB3	2.53	0.43
1:D:51:ARG:HD3	1:D:51:ARG:N	2.33	0.43
1:C:334:LEU:O	1:C:338:LEU:HG	2.18	0.43
1:B:48:TYR:CD1	1:B:55:PHE:CE1	3.05	0.43
1:B:257:THR:OG1	1:B:289:PHE:HZ	2.00	0.42
1:B:317:ARG:HG3	1:B:320:ARG:HH21	1.84	0.42
1:D:201:LEU:O	1:D:205:THR:HG23	2.18	0.42
1:A:39:GLU:H	1:A:39:GLU:CD	2.22	0.42
1:B:314:ARG:HE	1:B:317:ARG:NH2	2.16	0.42
1:B:396:VAL:HG13	1:B:401:LEU:HD23	2.01	0.42
1:A:100:ASP:N	1:A:100:ASP:OD1	2.52	0.42
1:A:116:GLU:O	1:A:118:ALA:N	2.52	0.42
1:A:9:ARG:HD2	1:A:20:GLN:OE1	2.20	0.42
1:A:106:TRP:CZ2	1:A:110:ARG:HD2	2.55	0.42
1:A:405:ASP:OD1	1:A:405:ASP:N	2.53	0.42
1:B:48:TYR:HD1	1:B:55:PHE:CE1	2.38	0.42
1:A:111:GLN:HG3	1:A:112:HIS:N	2.34	0.42
1:B:357:LEU:O	1:B:361:THR:HG22	2.20	0.42
1:C:115:ALA:HB2	1:C:441:MET:SD	2.60	0.42
1:D:103:PRO:O	3:D:601:HOH:O	2.22	0.42
1:D:322:PHE:O	1:D:326:ARG:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:THR:HG21	1:A:249:GLU:HG3	2.02	0.41
1:B:198:PHE:HD2	1:B:199:PHE:HD1	1.67	0.41
1:D:281:ILE:O	1:D:285:ALA:HB3	2.20	0.41
1:D:18:ARG:NE	1:D:54:GLU:OE2	2.53	0.41
1:B:288:PRO:HA	1:B:291:LEU:HB3	2.02	0.41
1:B:100:ASP:OD1	1:B:100:ASP:N	2.53	0.41
1:C:37:LEU:HB3	1:C:55:PHE:CE2	2.55	0.41
1:C:286:ILE:HD11	1:C:323:LYS:NZ	2.35	0.41
1:B:21:THR:HG22	1:B:22:TYR:H	1.86	0.41
1:B:287:LEU:HA	1:B:290:TYR:HD2	1.84	0.41
1:C:26:LEU:HD23	1:C:26:LEU:HA	1.77	0.41
1:C:48:TYR:HD1	1:C:55:PHE:HE1	1.63	0.41
1:A:337:THR:HG22	1:A:442:TYR:CD2	2.55	0.41
1:B:198:PHE:HD2	1:B:199:PHE:CD1	2.38	0.41
1:C:110:ARG:HH21	1:C:113[B]:ARG:HB3	1.84	0.41
1:C:197:LEU:HD13	1:C:324:LEU:HD13	2.02	0.41
1:A:286:ILE:HD11	1:A:323:LYS:NZ	2.36	0.41
1:B:213:GLU:OE1	1:C:386:ASN:ND2	2.47	0.41
1:D:318:ILE:HG12	1:D:319:LEU:HD22	2.02	0.41
1:B:207:PHE:HD1	1:B:314:ARG:HD3	1.86	0.41
1:C:242:GLU:HG2	1:C:244:PHE:CE1	2.56	0.41
1:C:363:ILE:HD12	1:C:363:ILE:HA	1.90	0.41
1:D:396:VAL:HG13	1:D:401:LEU:HD23	2.02	0.41
1:A:352:LEU:O	1:A:356:VAL:HG23	2.21	0.40
1:A:422:LEU:H	1:A:422:LEU:HG	1.76	0.40
1:B:11:VAL:HG22	1:B:20:GLN:HG2	2.01	0.40
1:B:323:LYS:HE2	1:B:323:LYS:HB2	1.61	0.40
1:C:204:ILE:HG13	1:C:321:ILE:HD11	2.03	0.40
1:D:32:THR:HG22	1:D:95:GLY:O	2.21	0.40
1:D:368[B]:ARG:NH2	1:D:384:PHE:O	2.49	0.40
1:B:78:CYS:HB2	1:B:105:CYS:HB3	2.03	0.40
1:C:85:PRO:O	1:C:88:GLU:HG3	2.21	0.40
1:C:243:ALA:O	1:C:246:THR:OG1	2.32	0.40
1:C:259:GLU:HG3	1:C:259:GLU:O	2.21	0.40
1:B:26:LEU:HD22	1:B:34:LEU:HB3	2.03	0.40
1:B:116:GLU:O	1:B:118:ALA:N	2.55	0.40
1:B:363:ILE:HD12	1:B:363:ILE:HA	1.95	0.40
1:C:78:CYS:SG	1:C:109:TYR:HB2	2.62	0.40
1:B:347:LEU:HD12	1:B:347:LEU:HA	1.86	0.40

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	310/519 (60%)	287 (93%)	21 (7%)	2 (1%)	25	37
1	B	310/519 (60%)	285 (92%)	23 (7%)	2 (1%)	25	37
1	C	310/519 (60%)	288 (93%)	20 (6%)	2 (1%)	25	37
1	D	310/519 (60%)	283 (91%)	25 (8%)	2 (1%)	25	37
All	All	1240/2076 (60%)	1143 (92%)	89 (7%)	8 (1%)	26	37

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	117	GLU
1	B	117	GLU
1	C	117	GLU
1	A	74	GLY
1	B	74	GLY
1	C	74	GLY
1	D	74	GLY
1	D	117	GLU

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	276/443 (62%)	256 (93%)	20 (7%)	14	22
1	B	276/443 (62%)	257 (93%)	19 (7%)	15	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	276/443 (62%)	247 (90%)	29 (10%)	7	10
1	D	276/443 (62%)	248 (90%)	28 (10%)	7	10
All	All	1104/1772 (62%)	1008 (91%)	96 (9%)	15	15

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ARG
1	A	51	ARG
1	A	70	TYR
1	A	110	ARG
1	A	113[A]	ARG
1	A	113[B]	ARG
1	A	119	LEU
1	A	203	SER
1	A	210	GLU
1	A	241[A]	THR
1	A	241[B]	THR
1	A	257	THR
1	A	314	ARG
1	A	328	PHE
1	A	380	GLU
1	A	385	LYS
1	A	399	THR
1	A	412	SER
1	A	444	SER
1	A	450	GLN
1	B	21	THR
1	B	23	ARG
1	B	37	LEU
1	B	51	ARG
1	B	70	TYR
1	B	108	THR
1	B	113[A]	ARG
1	B	113[B]	ARG
1	B	119	LEU
1	B	241[A]	THR
1	B	241[B]	THR
1	B	257	THR
1	B	314	ARG
1	B	328	PHE

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Mol	Chain	Res	Type
1	B	379	SER
1	B	385	LYS
1	B	399	THR
1	B	412	SER
1	B	444	SER
1	C	18	ARG
1	C	21	THR
1	C	23	ARG
1	C	24	SER
1	C	37	LEU
1	C	51	ARG
1	C	70	TYR
1	C	88	GLU
1	C	110	ARG
1	C	111	GLN
1	C	113[A]	ARG
1	C	113[B]	ARG
1	C	119	LEU
1	C	210	GLU
1	C	241[A]	THR
1	C	241[B]	THR
1	C	247	TYR
1	C	257	THR
1	C	262	MET
1	C	263	ARG
1	C	292	GLU
1	C	314	ARG
1	C	324	LEU
1	C	328	PHE
1	C	380	GLU
1	C	386	ASN
1	C	399	THR
1	C	412	SER
1	C	444	SER
1	D	21	THR
1	D	23	ARG
1	D	24	SER
1	D	41	ASP
1	D	51	ARG
1	D	70	TYR
1	D	88	GLU
1	D	104	CYS

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Mol	Chain	Res	Type
1	D	110	ARG
1	D	113[A]	ARG
1	D	113[B]	ARG
1	D	241[A]	THR
1	D	241[B]	THR
1	D	254	VAL
1	D	257	THR
1	D	258	PHE
1	D	264	VAL
1	D	320	ARG
1	D	328	PHE
1	D	329	VAL
1	D	369	ILE
1	D	385	LYS
1	D	386	ASN
1	D	399	THR
1	D	412	SER
1	D	444	SER
1	D	447	MET
1	D	451	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	216	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13604. These allow visual inspection of the internal detail of the map and identification of artifacts.

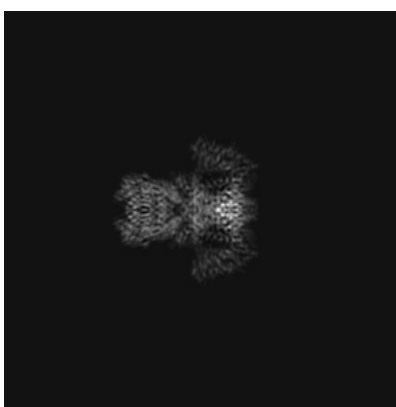
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

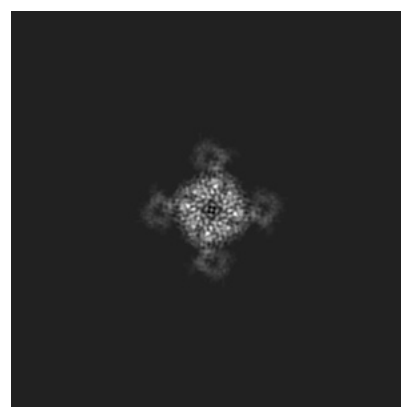
6.1.1 Primary map



X



Y



Z

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

6.2 Central slices [i](#)

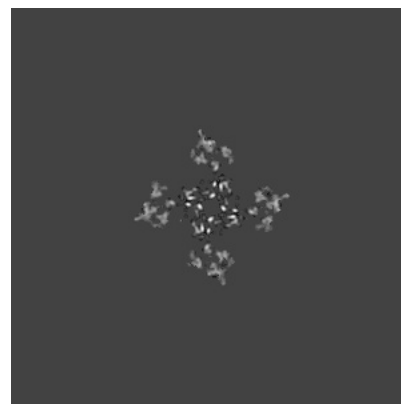
6.2.1 Primary map



X Index: 175



Y Index: 175



Z Index: 175

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: 180



Y Index: 170



Z Index: 200

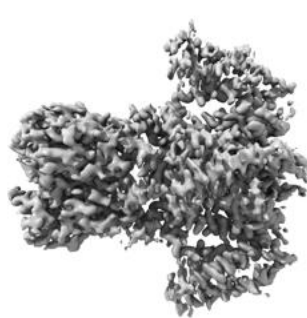
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views [i](#)

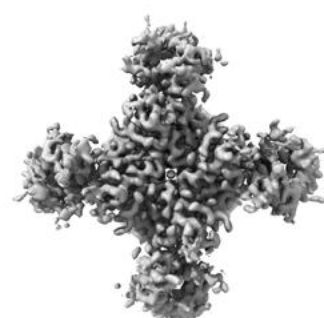
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.356. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

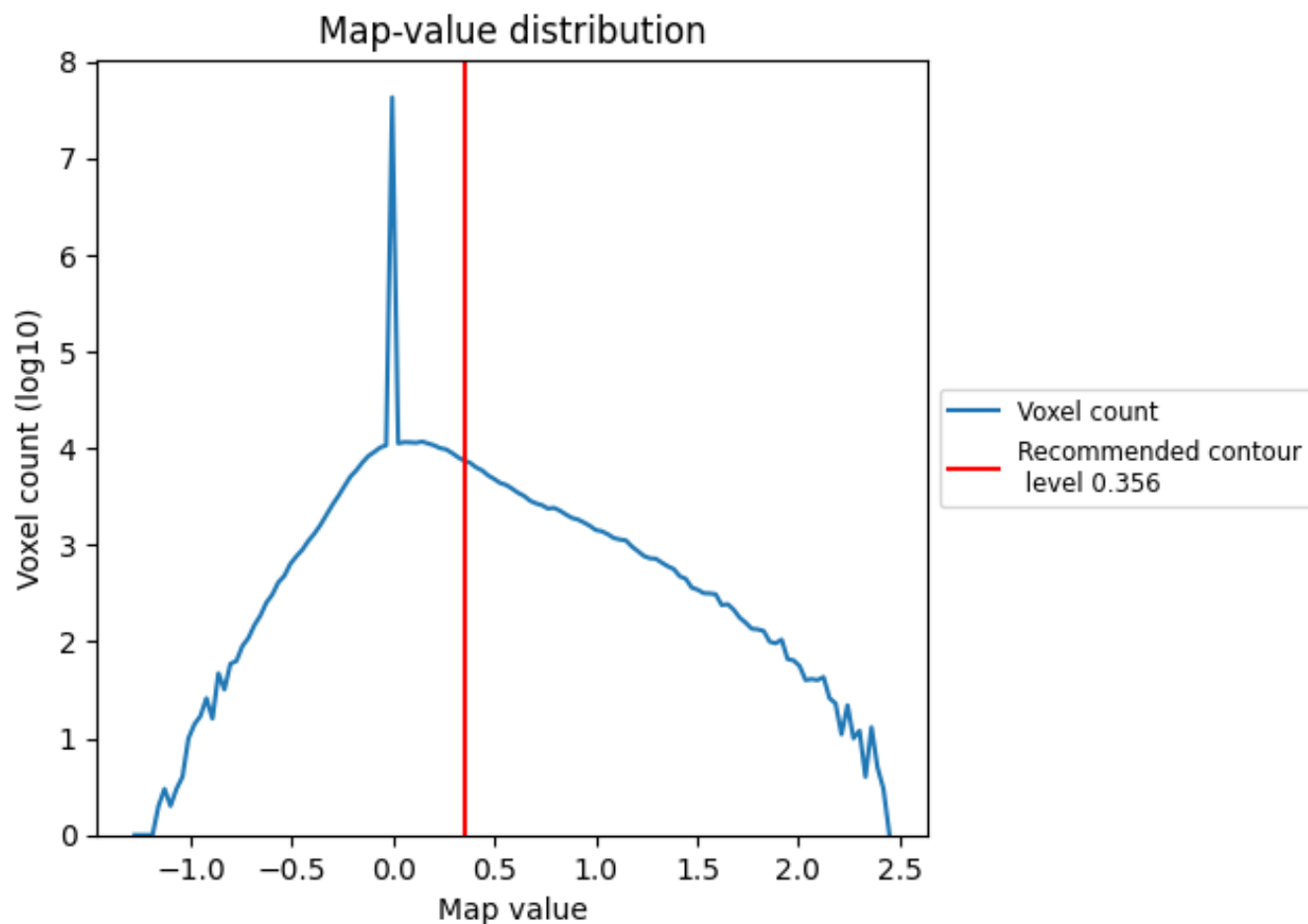
6.5 Mask visualisation

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

7 Map analysis [i](#)

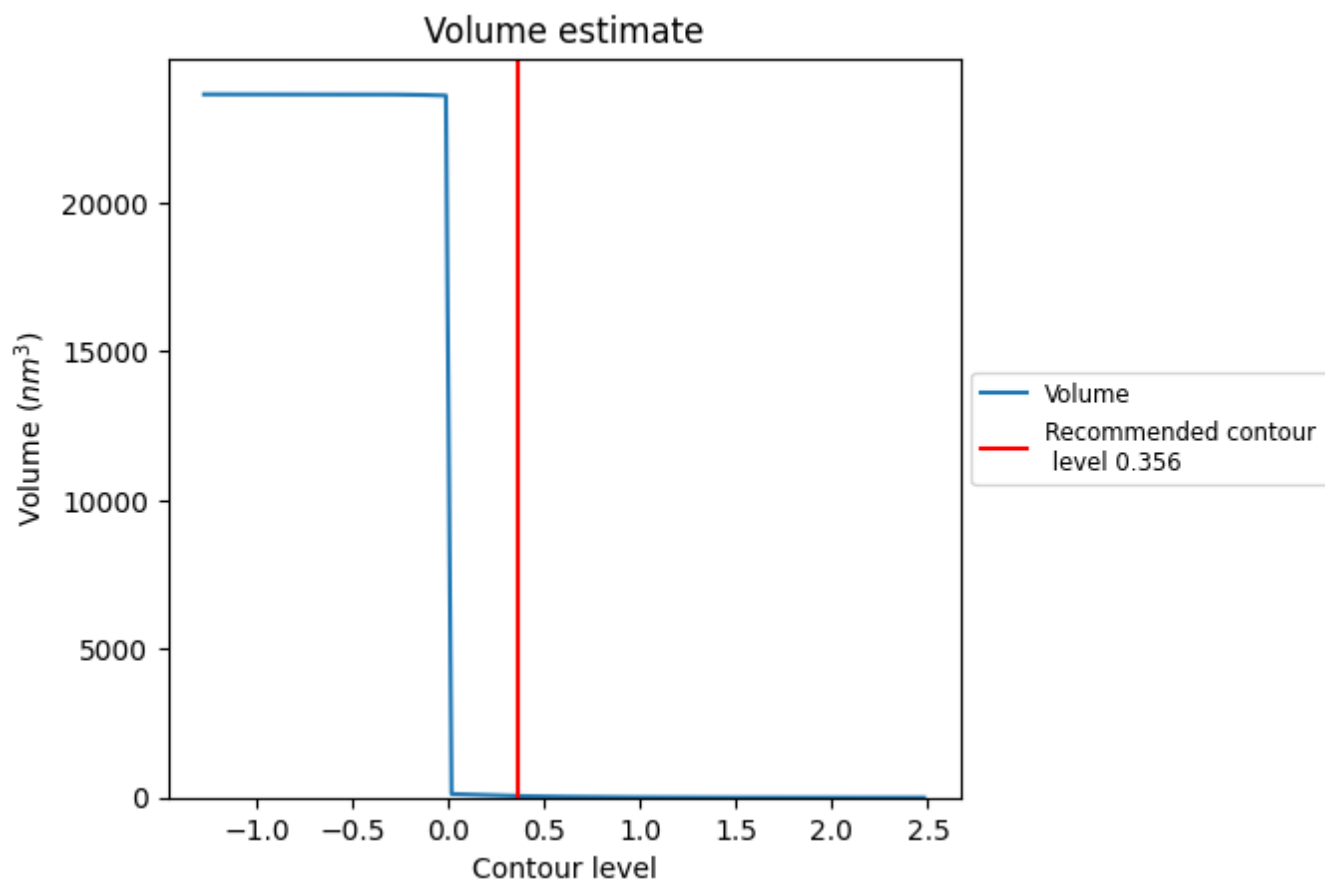
This section contains the results of statistical analysis of the map.

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.

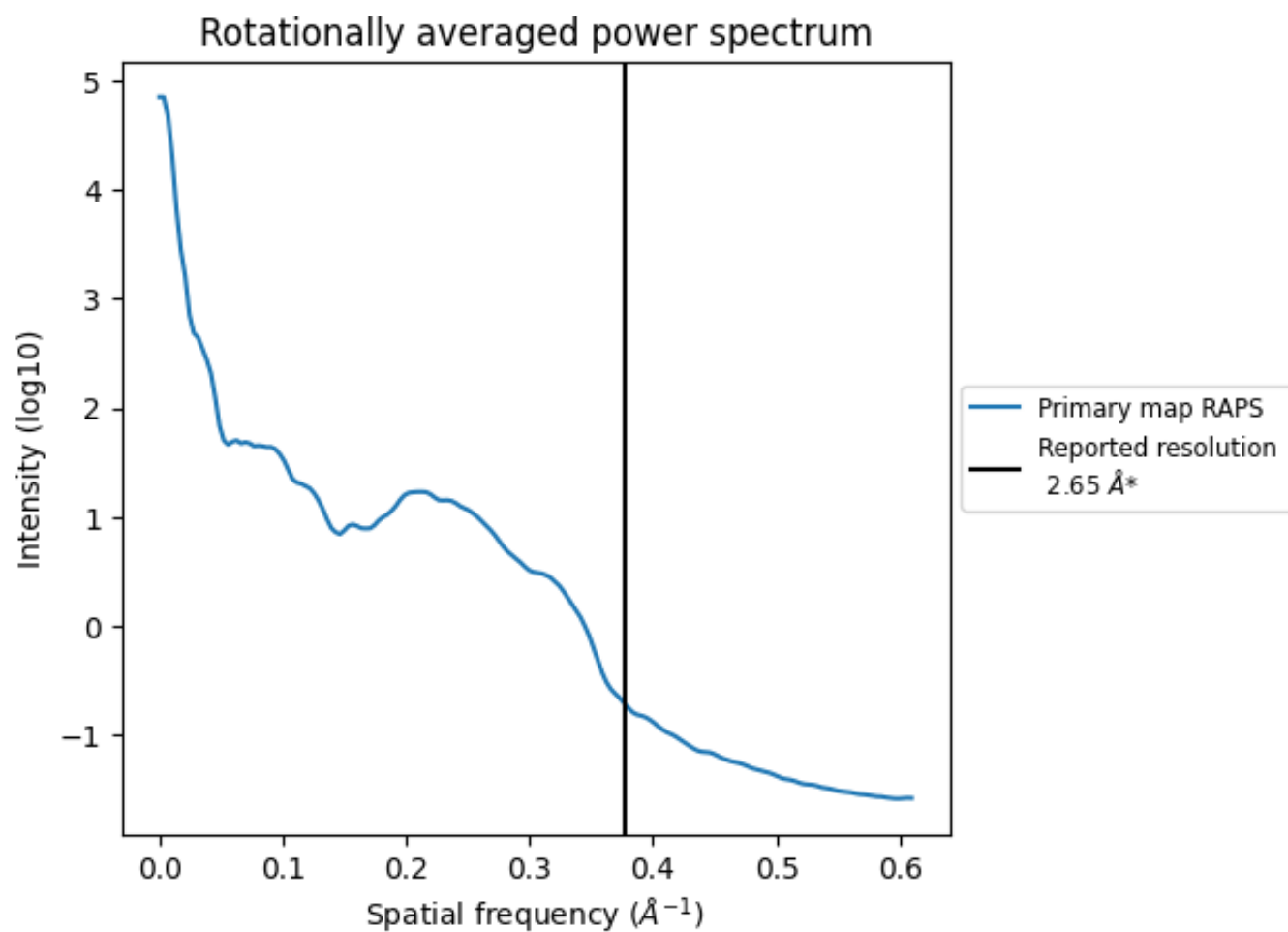
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 53 nm^3 ; this corresponds to an approximate mass of 48 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 0.377 Å⁻¹

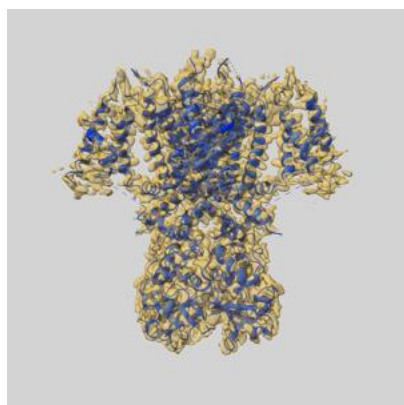
8 Fourier-Shell correlation

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

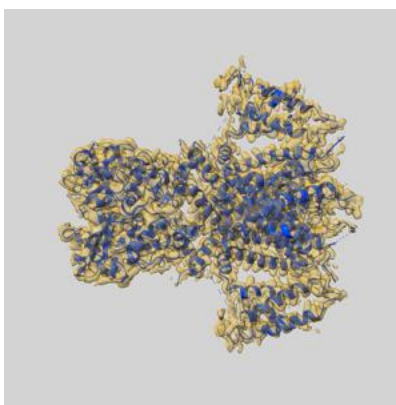
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-13604 and PDB model 7PQT. Per-residue inclusion information can be found in section 3 on page 5.

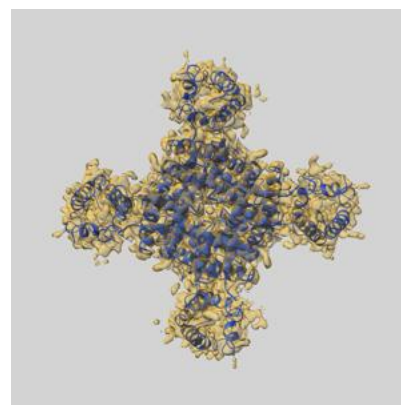
9.1 Map-model overlay [i](#)



X



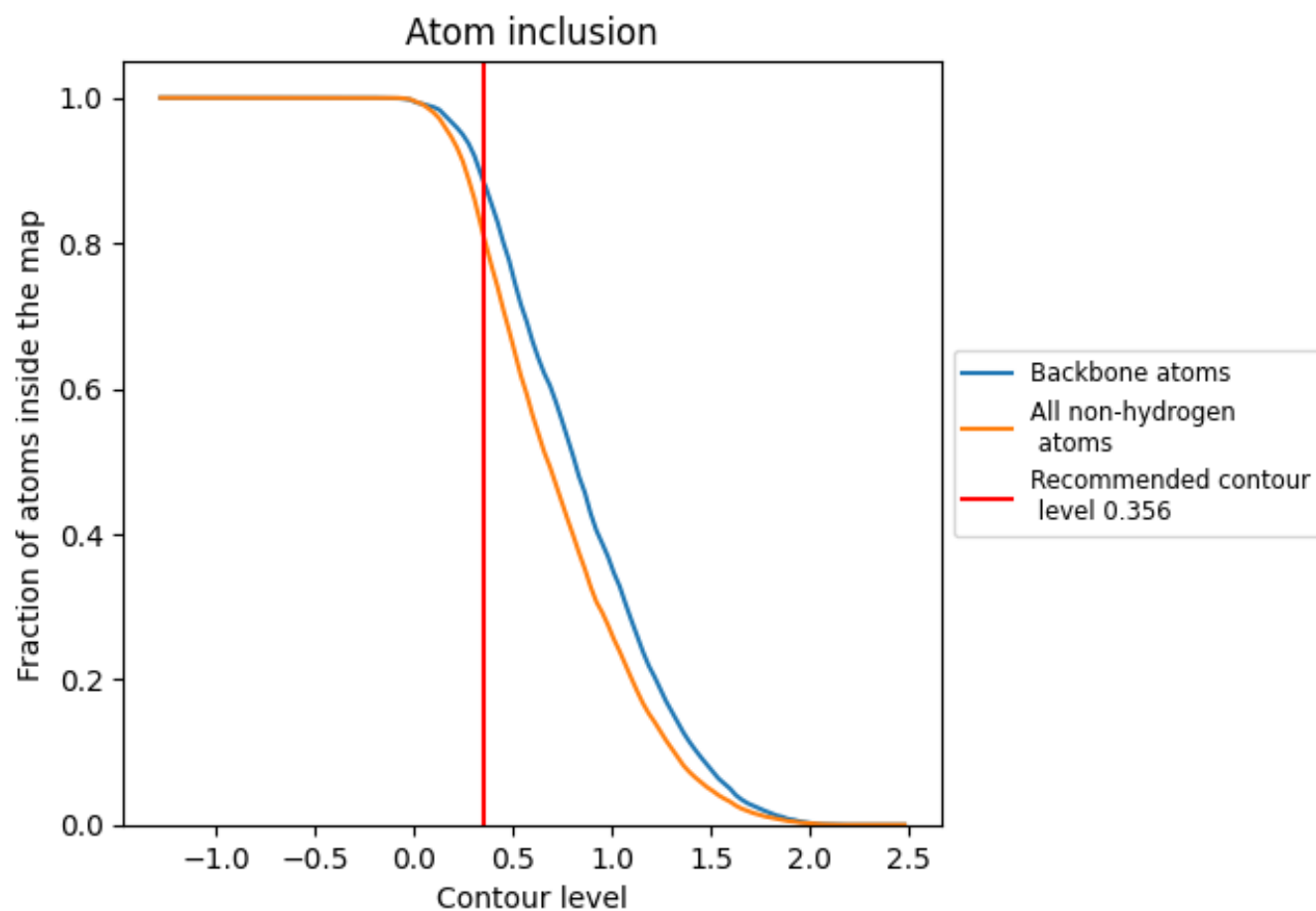
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.356 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 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 81% of all non-hydrogen atoms, are inside the map.