



Full wwPDB EM Validation Report ⓘ

Oct 7, 2024 – 10:07 AM EDT

PDB ID : 7RW6
EMDB ID : EMD-24709
Title : BORF2-APOBEC3Bctd Complex
Authors : Shaban, N.M.; Yan, R.; Shi, K.; McLellan, J.S.; Yu, Z.; Harris, R.S.
Deposited on : 2021-08-19
Resolution : 2.55 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.dev113
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

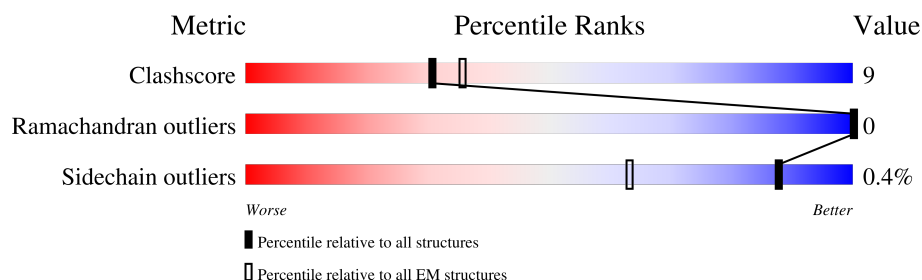
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.55 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	1197	
1	C	1197	
2	B	215	
2	D	215	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13130 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 Maltose/maltodextrin-binding periplasmic protein,Ribonucleoside-diphosphate reductase large subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	642	Total	C	N	O	S	0	0
			5100	3254	871	943	32		
1	C	642	Total	C	N	O	S	0	0
			5100	3254	871	943	32		

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-370	MET	-	initiating methionine	UNP P0AEX9
A	-288	ALA	ASP	engineered mutation	UNP P0AEX9
A	-287	ALA	LYS	engineered mutation	UNP P0AEX9
A	-198	ALA	GLU	engineered mutation	UNP P0AEX9
A	-197	ALA	ASN	engineered mutation	UNP P0AEX9
A	-131	ALA	LYS	engineered mutation	UNP P0AEX9
A	-11	ALA	-	linker	UNP P0AEX9
A	-10	ALA	-	linker	UNP P0AEX9
A	-9	LEU	-	linker	UNP P0AEX9
A	-8	ALA	-	linker	UNP P0AEX9
A	-7	ALA	-	linker	UNP P0AEX9
A	-6	ALA	-	linker	UNP P0AEX9
A	-5	GLN	-	linker	UNP P0AEX9
A	-4	THR	-	linker	UNP P0AEX9
A	-3	ASN	-	linker	UNP P0AEX9
A	-2	ALA	-	linker	UNP P0AEX9
A	-1	ALA	-	linker	UNP P0AEX9
A	0	ALA	-	linker	UNP P0AEX9
C	-370	MET	-	initiating methionine	UNP P0AEX9
C	-288	ALA	ASP	engineered mutation	UNP P0AEX9
C	-287	ALA	LYS	engineered mutation	UNP P0AEX9
C	-198	ALA	GLU	engineered mutation	UNP P0AEX9
C	-197	ALA	ASN	engineered mutation	UNP P0AEX9
C	-131	ALA	LYS	engineered mutation	UNP P0AEX9
C	-11	ALA	-	linker	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	ALA	-	linker	UNP P0AEX9
C	-9	LEU	-	linker	UNP P0AEX9
C	-8	ALA	-	linker	UNP P0AEX9
C	-7	ALA	-	linker	UNP P0AEX9
C	-6	ALA	-	linker	UNP P0AEX9
C	-5	GLN	-	linker	UNP P0AEX9
C	-4	THR	-	linker	UNP P0AEX9
C	-3	ASN	-	linker	UNP P0AEX9
C	-2	ALA	-	linker	UNP P0AEX9
C	-1	ALA	-	linker	UNP P0AEX9
C	0	ALA	-	linker	UNP P0AEX9

- Molecule 2 is a protein called DNA dC->dU-editing enzyme APOBEC-3B.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	175	Total	C	N	O	S	0	0
			1464	940	250	263	11		
2	D	175	Total	C	N	O	S	0	0
			1464	940	250	263	11		

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	383	LYS	-	expression tag	UNP Q9UH17
B	384	LEU	-	expression tag	UNP Q9UH17
B	385	GLY	-	expression tag	UNP Q9UH17
B	386	PRO	-	expression tag	UNP Q9UH17
B	387	GLU	-	expression tag	UNP Q9UH17
B	388	GLN	-	expression tag	UNP Q9UH17
B	389	LYS	-	expression tag	UNP Q9UH17
B	390	LEU	-	expression tag	UNP Q9UH17
B	391	ILE	-	expression tag	UNP Q9UH17
B	392	SER	-	expression tag	UNP Q9UH17
B	393	GLU	-	expression tag	UNP Q9UH17
B	394	GLU	-	expression tag	UNP Q9UH17
B	395	ASP	-	expression tag	UNP Q9UH17
B	396	LEU	-	expression tag	UNP Q9UH17
B	397	ASN	-	expression tag	UNP Q9UH17
B	398	SER	-	expression tag	UNP Q9UH17
B	399	ALA	-	expression tag	UNP Q9UH17
B	400	VAL	-	expression tag	UNP Q9UH17
B	401	ASP	-	expression tag	UNP Q9UH17

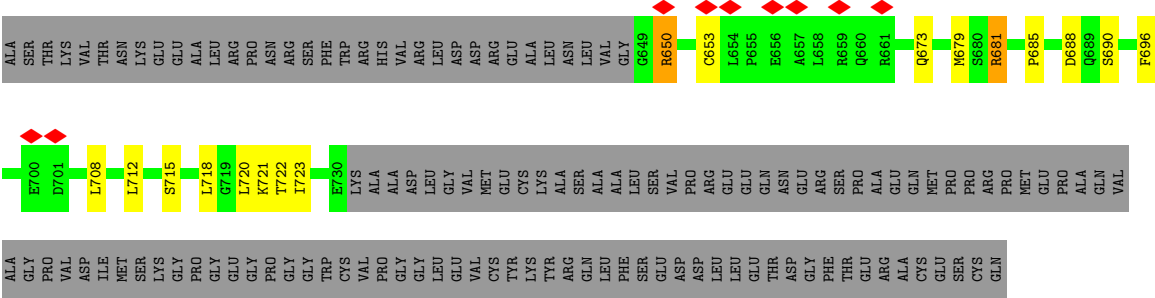
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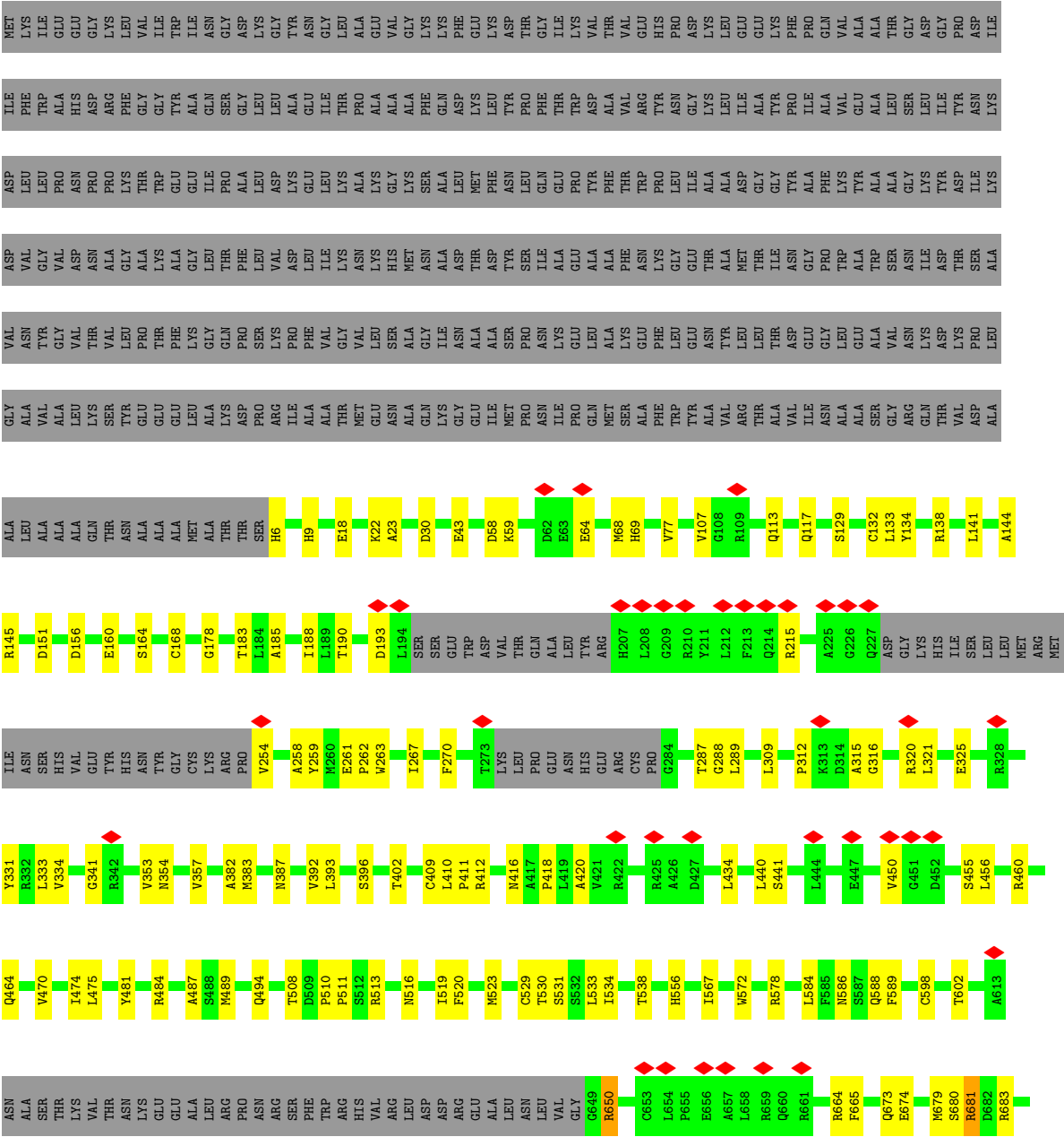
Chain	Residue	Modelled	Actual	Comment	Reference
B	402	HIS	-	expression tag	UNP Q9UH17
B	403	HIS	-	expression tag	UNP Q9UH17
B	404	HIS	-	expression tag	UNP Q9UH17
B	405	HIS	-	expression tag	UNP Q9UH17
B	406	HIS	-	expression tag	UNP Q9UH17
B	407	HIS	-	expression tag	UNP Q9UH17
D	383	LYS	-	expression tag	UNP Q9UH17
D	384	LEU	-	expression tag	UNP Q9UH17
D	385	GLY	-	expression tag	UNP Q9UH17
D	386	PRO	-	expression tag	UNP Q9UH17
D	387	GLU	-	expression tag	UNP Q9UH17
D	388	GLN	-	expression tag	UNP Q9UH17
D	389	LYS	-	expression tag	UNP Q9UH17
D	390	LEU	-	expression tag	UNP Q9UH17
D	391	ILE	-	expression tag	UNP Q9UH17
D	392	SER	-	expression tag	UNP Q9UH17
D	393	GLU	-	expression tag	UNP Q9UH17
D	394	GLU	-	expression tag	UNP Q9UH17
D	395	ASP	-	expression tag	UNP Q9UH17
D	396	LEU	-	expression tag	UNP Q9UH17
D	397	ASN	-	expression tag	UNP Q9UH17
D	398	SER	-	expression tag	UNP Q9UH17
D	399	ALA	-	expression tag	UNP Q9UH17
D	400	VAL	-	expression tag	UNP Q9UH17
D	401	ASP	-	expression tag	UNP Q9UH17
D	402	HIS	-	expression tag	UNP Q9UH17
D	403	HIS	-	expression tag	UNP Q9UH17
D	404	HIS	-	expression tag	UNP Q9UH17
D	405	HIS	-	expression tag	UNP Q9UH17
D	406	HIS	-	expression tag	UNP Q9UH17
D	407	HIS	-	expression tag	UNP Q9UH17

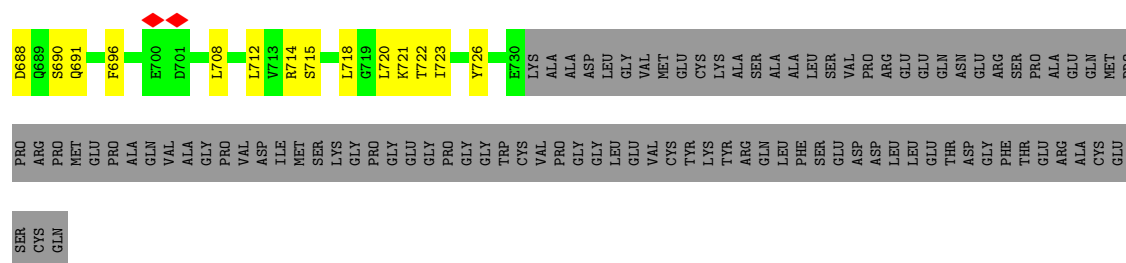
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	AltConf
3	B	1	Total Zn 1 1	0
3	D	1	Total Zn 1 1	0

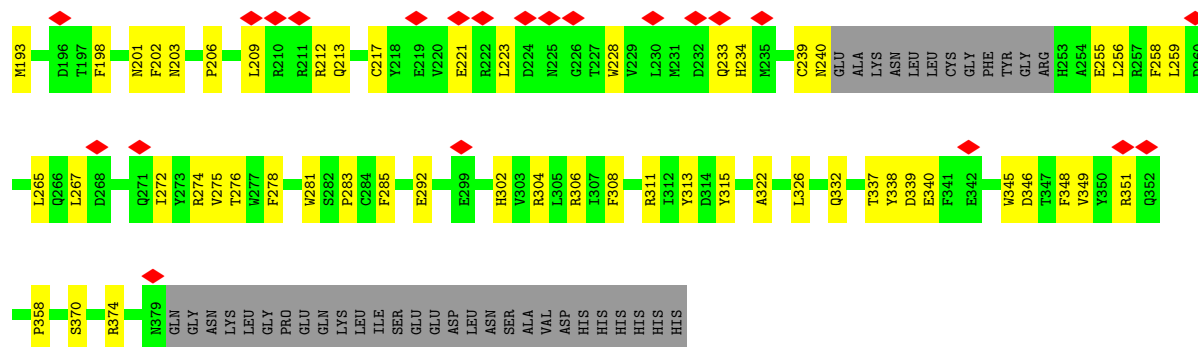


● Molecule 1: Maltose/maltodextrin-binding periplasmic protein,Ribonucleoside-diphosphate reductase large subunit

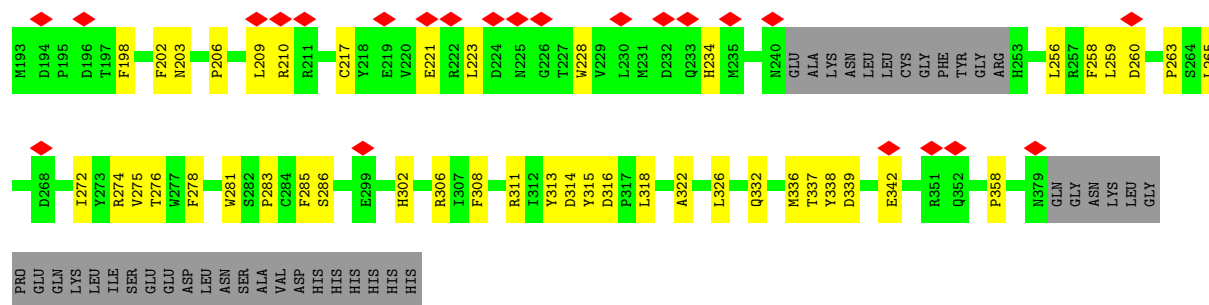




- Molecule 2: DNA dC->dU-editing enzyme APOBEC-3B



- Molecule 2: DNA dC->dU-editing enzyme APOBEC-3B



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	243192	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 ($e^-/\text{\AA}^2$)	1.2	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	72.193	Depositor
Minimum map value	-3.075	Depositor
Average map value	0.013	Depositor
Map value standard deviation	1.289	Depositor
Recommended contour level	10	Depositor
Map size (Å)	283.58398, 283.58398, 283.58398	wwPDB
Map dimensions	336, 336, 336	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.844, 0.844, 0.844	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: ZN

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	0.26	0/5222	0.49	0/7069
1	C	0.26	0/5222	0.49	0/7069
2	B	0.25	0/1507	0.51	0/2048
2	D	0.26	0/1507	0.50	0/2048
All	All	0.26	0/13458	0.49	0/18234

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	5100	0	4954	79	0
1	C	5100	0	4954	86	0
2	B	1464	0	1378	35	0
2	D	1464	0	1378	30	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
All	All	13130	0	12664	225	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 9.

All (225) 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)
2:B:285:PHE:HB2	2:B:322:ALA:HB2	1.58	0.85
2:D:285:PHE:HB2	2:D:322:ALA:HB2	1.62	0.80
1:C:168:CYS:SG	1:C:494:GLN:NE2	2.56	0.79
2:B:306:ARG:HG2	2:B:332:GLN:HE21	1.48	0.78
1:A:681:ARG:HG3	1:A:718:LEU:HB3	1.67	0.76
2:D:306:ARG:HG2	2:D:332:GLN:HE21	1.50	0.75
1:C:168:CYS:HB3	1:C:598:CYS:SG	2.28	0.73
1:C:321:LEU:HD13	1:C:325:GLU:HG3	1.69	0.73
2:D:217:CYS:HB2	2:D:278:PHE:HB2	1.71	0.73
1:C:681:ARG:HG3	1:C:718:LEU:HB3	1.69	0.72
1:C:516:ASN:HD21	1:C:683:ARG:HB3	1.54	0.72
1:A:9:HIS:NE2	1:A:43:GLU:OE2	2.24	0.69
1:C:183:THR:HG22	1:C:185:ALA:H	1.58	0.68
2:B:274:ARG:HG3	2:B:304:ARG:HB2	1.75	0.67
2:D:234:HIS:HD2	2:D:265:LEU:HG	1.60	0.67
1:C:393:LEU:O	1:C:588:GLN:NE2	2.28	0.67
1:A:293:GLU:HB2	1:A:368:GLU:HG3	1.77	0.66
1:A:331:TYR:HA	1:A:334:VAL:HG22	1.78	0.65
1:A:168:CYS:HB3	1:A:598:CYS:SG	2.37	0.65
2:D:198:PHE:O	2:D:202:PHE:HB2	1.97	0.65
1:C:680:SER:OG	1:C:691:GLN:NE2	2.30	0.65
2:D:337:THR:HG22	2:D:338:TYR:H	1.62	0.65
1:C:529:CYS:SG	1:C:578:ARG:NH1	2.70	0.64
1:A:520:PHE:HE1	1:A:592:LEU:HD21	1.63	0.64
1:A:529:CYS:SG	1:A:578:ARG:NH1	2.71	0.64
2:B:223:LEU:HD23	2:B:272:ILE:HD11	1.79	0.63
2:B:346:ASP:HA	2:B:351:ARG:HD3	1.81	0.63
1:C:688:ASP:O	1:C:721:LYS:NZ	2.32	0.63
1:A:183:THR:HG22	1:A:185:ALA:H	1.63	0.63
2:B:221:GLU:OE1	2:B:274:ARG:NH2	2.32	0.62
1:A:429:GLN:HG3	1:A:431:ASP:H	1.64	0.62
2:D:338:TYR:HE1	2:D:358:PRO:HG3	1.64	0.61
2:B:203:ASN:O	2:B:311:ARG:NH1	2.31	0.61
1:C:267:ILE:HD11	1:C:289:LEU:HD11	1.83	0.61
2:B:338:TYR:HE1	2:B:358:PRO:HG3	1.65	0.61
1:A:688:ASP:O	1:A:721:LYS:NZ	2.34	0.61
2:D:209:LEU:O	2:D:315:TYR:OH	2.18	0.61
1:A:138:ARG:HH22	1:A:151:ASP:HB3	1.67	0.60
1:A:376:ARG:NH2	1:A:381:GLU:OE2	2.34	0.59
1:A:393:LEU:O	1:A:588:GLN:NE2	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:674:GLU:HG2	1:C:714:ARG:HH12	1.67	0.59
1:C:331:TYR:HA	1:C:334:VAL:HG22	1.83	0.58
2:B:276:THR:HG22	2:B:306:ARG:HB2	1.85	0.58
2:D:221:GLU:HB3	2:D:228:TRP:HB3	1.84	0.58
2:B:217:CYS:HB2	2:B:278:PHE:HB2	1.86	0.57
2:B:370:SER:OG	2:B:374:ARG:NH2	2.37	0.57
1:C:156:ASP:O	1:C:160:GLU:HG3	2.04	0.57
1:C:690:SER:OG	1:C:722:THR:OG1	2.18	0.57
2:B:193:MET:N	2:B:348:PHE:O	2.37	0.57
1:C:434:LEU:HD22	1:C:460:ARG:HD2	1.85	0.57
2:B:198:PHE:O	2:B:202:PHE:HB2	2.05	0.57
1:C:416:ASN:ND2	1:C:450:VAL:O	2.37	0.57
2:D:308:PHE:HB3	2:D:336:MET:HG2	1.87	0.57
1:A:267:ILE:HD11	1:A:289:LEU:HD11	1.86	0.56
1:C:353:VAL:HG13	1:C:708:LEU:HD23	1.87	0.56
1:A:434:LEU:HD22	1:A:460:ARG:HD2	1.87	0.56
1:C:316:GLY:HA3	1:C:333:LEU:HD11	1.88	0.56
1:A:440:LEU:N	1:A:455:SER:O	2.36	0.56
1:C:440:LEU:N	1:C:455:SER:O	2.37	0.56
1:A:134:TYR:CZ	1:A:475:LEU:HB3	2.41	0.56
2:D:223:LEU:HD23	2:D:272:ILE:HD11	1.87	0.56
1:A:412:ARG:NH2	1:A:418:PRO:O	2.39	0.56
1:C:412:ARG:NH2	1:C:418:PRO:O	2.39	0.56
2:B:339:ASP:OD1	2:B:340:GLU:N	2.40	0.55
2:D:283:PRO:HB2	2:D:326:LEU:HD11	1.89	0.55
1:A:593:MET:HG3	1:A:595:THR:H	1.72	0.55
1:C:261:GLU:OE1	1:C:263:TRP:NE1	2.33	0.55
1:A:88:ARG:NH1	1:A:165:GLN:OE1	2.40	0.55
1:C:132:CYS:SG	1:C:138:ARG:NH2	2.63	0.54
1:A:690:SER:OG	1:A:722:THR:OG1	2.21	0.54
2:B:209:LEU:O	2:B:315:TYR:OH	2.25	0.54
2:D:286:SER:OG	2:D:316:ASP:OD1	2.17	0.54
1:A:215:ARG:HA	1:A:254:VAL:HG11	1.90	0.54
1:A:143:ARG:NH1	1:A:430:GLY:O	2.41	0.54
1:A:353:VAL:HG13	1:A:708:LEU:HD23	1.90	0.54
1:A:321:LEU:HD13	1:A:325:GLU:HG3	1.89	0.53
1:C:393:LEU:HD11	1:C:690:SER:HB2	1.91	0.53
1:A:261:GLU:OE1	1:A:263:TRP:NE1	2.39	0.53
1:A:439:ARG:HG3	1:A:440:LEU:HD23	1.91	0.52
2:B:256:LEU:O	2:B:259:LEU:N	2.42	0.52
1:C:262:PRO:HG3	1:C:289:LEU:HD12	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:GLU:O	1:C:22:LYS:NZ	2.43	0.52
1:A:715:SER:HB2	1:A:723:ILE:HG23	1.92	0.52
2:B:234:HIS:HD2	2:B:265:LEU:HG	1.75	0.52
1:C:409:CYS:HB2	1:C:494:GLN:HE21	1.75	0.52
1:A:316:GLY:HA3	1:A:333:LEU:HD11	1.91	0.51
1:A:584:LEU:HD13	1:A:589:PHE:HZ	1.76	0.51
1:C:9:HIS:NE2	1:C:43:GLU:OE2	2.43	0.51
1:C:584:LEU:HD13	1:C:589:PHE:HZ	1.75	0.51
1:C:178:GLY:HA2	2:D:210:ARG:HH22	1.75	0.51
1:C:134:TYR:CZ	1:C:475:LEU:HB3	2.46	0.51
1:A:18:GLU:O	1:A:22:LYS:NZ	2.43	0.50
1:C:673:GLN:NE2	1:C:696:PHE:O	2.44	0.50
1:A:409:CYS:SG	1:A:494:GLN:NE2	2.83	0.50
2:D:272:ILE:HG22	2:D:302:HIS:HB2	1.93	0.50
1:A:190:THR:HA	1:A:396:SER:HB2	1.94	0.50
1:C:712:LEU:HD22	1:C:723:ILE:HG22	1.94	0.49
1:A:173:PHE:HA	1:A:184:LEU:HD22	1.95	0.49
1:C:164:SER:HB3	1:C:420:ALA:H	1.78	0.49
1:C:320:ARG:O	1:C:382:ALA:N	2.46	0.49
2:D:203:ASN:O	2:D:311:ARG:NH1	2.38	0.48
1:A:712:LEU:HD22	1:A:723:ILE:HG22	1.95	0.48
1:C:144:ALA:HB2	1:C:533:LEU:HD23	1.95	0.48
1:A:437:LEU:HD22	1:A:567:ILE:HD12	1.95	0.48
2:B:201:ASN:ND2	2:B:213:GLN:O	2.38	0.48
1:C:188:ILE:HG12	1:C:402:THR:HG22	1.96	0.48
1:A:46:THR:OG1	1:A:118:HIS:ND1	2.40	0.48
1:A:132:CYS:SG	1:A:138:ARG:NH2	2.79	0.48
1:C:190:THR:HA	1:C:396:SER:HB2	1.96	0.48
2:B:283:PRO:HB2	2:B:326:LEU:HD11	1.95	0.48
1:C:441:SER:HB2	1:C:456:LEU:HB2	1.96	0.47
1:A:104:TYR:HE2	1:A:601:VAL:HG11	1.80	0.47
1:A:320:ARG:O	1:A:382:ALA:N	2.47	0.47
1:C:520:PHE:HA	1:C:523:MET:HG2	1.97	0.47
2:D:258:PHE:CZ	2:D:275:VAL:HG21	2.50	0.47
1:A:520:PHE:HA	1:A:523:MET:HG2	1.97	0.47
2:B:239:CYS:SG	2:B:240:ASN:N	2.86	0.47
2:D:256:LEU:O	2:D:259:LEU:N	2.49	0.46
2:D:260:ASP:O	2:D:263:PRO:HD2	2.15	0.46
1:A:262:PRO:HG3	1:A:289:LEU:HD12	1.98	0.46
2:B:256:LEU:HD11	2:B:292:GLU:HG3	1.97	0.46
1:A:160:GLU:OE1	1:A:425:ARG:NH1	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:534:ILE:O	1:A:538:THR:OG1	2.18	0.46
1:C:6:HIS:O	1:C:6:HIS:ND1	2.49	0.46
2:D:234:HIS:CD2	2:D:265:LEU:HG	2.47	0.46
1:C:141:LEU:O	1:C:145:ARG:N	2.49	0.46
1:A:325:GLU:HB3	1:A:328:ARG:NH2	2.30	0.46
2:B:221:GLU:HB3	2:B:228:TRP:HB3	1.98	0.46
1:C:534:ILE:O	1:C:538:THR:OG1	2.26	0.46
1:C:674:GLU:CG	1:C:714:ARG:HH12	2.29	0.46
2:D:223:LEU:HB2	2:D:228:TRP:CZ3	2.50	0.46
2:D:276:THR:HG22	2:D:306:ARG:HB2	1.97	0.46
1:A:151:ASP:OD1	1:A:151:ASP:N	2.49	0.46
1:A:673:GLN:NE2	1:A:696:PHE:O	2.49	0.46
2:B:256:LEU:HD22	2:B:259:LEU:HD12	1.98	0.45
2:B:281:TRP:HB3	2:B:313:TYR:HB2	1.98	0.45
1:A:64:GLU:O	1:A:68:MET:HG2	2.17	0.45
1:C:312:PRO:HA	1:C:315:ALA:HB2	1.98	0.45
1:A:441:SER:HB2	1:A:456:LEU:HB2	1.99	0.45
1:A:12:LEU:O	1:A:16:ILE:HG12	2.17	0.45
2:B:345:TRP:CZ3	2:B:349:VAL:HG11	2.52	0.45
1:C:410:LEU:HD13	1:C:519:ILE:HD13	1.98	0.45
1:A:258:ALA:O	1:A:288:GLY:N	2.50	0.45
1:A:688:ASP:OD1	1:A:688:ASP:N	2.50	0.45
1:C:470:VAL:HG11	1:C:489:MET:HG3	1.99	0.45
1:A:312:PRO:HA	1:A:315:ALA:HB2	1.98	0.45
1:C:508:THR:HG23	1:C:513:ARG:HH21	1.82	0.44
1:C:64:GLU:O	1:C:68:MET:HG2	2.17	0.44
1:C:151:ASP:OD1	1:C:151:ASP:N	2.49	0.44
1:C:270:PHE:CE2	1:C:287:THR:HG23	2.53	0.44
1:A:23:ALA:HB1	1:A:30:ASP:HB3	1.99	0.44
1:C:481:TYR:CZ	2:D:206:PRO:HG3	2.53	0.44
1:C:513:ARG:HG2	1:C:679:MET:HG2	2.00	0.44
2:D:339:ASP:HA	2:D:342:GLU:HG2	1.99	0.44
1:C:464:GLN:HG2	1:C:530:THR:OG1	2.17	0.44
2:D:206:PRO:HA	2:D:209:LEU:HG	2.00	0.44
2:D:221:GLU:OE1	2:D:274:ARG:NH2	2.29	0.44
1:A:513:ARG:HG2	1:A:679:MET:HG2	1.99	0.44
1:C:681:ARG:HB2	1:C:720:LEU:HD11	1.99	0.44
2:B:258:PHE:CZ	2:B:275:VAL:HG21	2.52	0.43
1:C:23:ALA:HB1	1:C:30:ASP:HB3	2.00	0.43
1:C:258:ALA:O	1:C:288:GLY:N	2.50	0.43
1:A:309:LEU:HB2	1:A:341:GLY:HA3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:518:GLU:HG2	1:A:566:SER:OG	2.18	0.43
2:B:223:LEU:HB2	2:B:228:TRP:CZ3	2.53	0.43
1:C:484:ARG:NH1	2:D:314:ASP:OD2	2.43	0.43
2:B:267:LEU:HD11	2:B:302:HIS:HE1	1.84	0.43
1:A:144:ALA:HB2	1:A:533:LEU:HD23	2.01	0.43
1:A:507:TYR:OH	1:A:608:PHE:O	2.30	0.43
1:C:387:ASN:HB3	1:C:392:VAL:HB	2.01	0.43
1:C:129:SER:O	1:C:133:LEU:HB2	2.18	0.43
1:A:578:ARG:O	1:A:582:ASP:HB2	2.19	0.43
1:A:650:ARG:HB2	1:A:653:CYS:HB3	2.00	0.43
1:C:688:ASP:OD1	1:C:688:ASP:N	2.48	0.42
1:C:715:SER:HB2	1:C:723:ILE:HG23	2.01	0.42
1:C:168:CYS:SG	1:C:602:THR:HG21	2.59	0.42
1:A:141:LEU:O	1:A:145:ARG:N	2.51	0.42
1:A:429:GLN:HB3	1:A:435:LEU:HG	2.01	0.42
1:C:58:ASP:OD1	1:C:59:LYS:HG2	2.20	0.42
1:C:383:MET:CE	1:C:393:LEU:HB3	2.49	0.42
1:A:45:VAL:HG12	1:A:113:GLN:HB3	2.00	0.42
1:A:293:GLU:HB2	1:A:368:GLU:CG	2.48	0.42
1:A:393:LEU:HD11	1:A:690:SER:HB2	2.00	0.42
2:B:337:THR:OG1	2:B:339:ASP:OD1	2.38	0.42
1:C:77:VAL:HG13	1:C:117:GLN:HB2	2.00	0.42
1:C:178:GLY:HA2	2:D:210:ARG:HH12	1.84	0.42
1:C:193:ASP:OD1	1:C:193:ASP:N	2.52	0.42
1:C:567:ILE:HB	1:C:572:TRP:CD1	2.55	0.42
1:A:270:PHE:CE2	1:A:287:THR:HG23	2.54	0.42
1:A:481:TYR:CZ	2:B:206:PRO:HG3	2.55	0.42
1:C:556:HIS:CD2	1:C:688:ASP:HA	2.55	0.42
1:C:533:LEU:HD13	1:C:578:ARG:HH22	1.85	0.41
2:D:285:PHE:CD1	2:D:318:LEU:HD13	2.54	0.41
1:A:375:TRP:CZ3	1:A:376:ARG:HG3	2.55	0.41
1:A:533:LEU:HD13	1:A:578:ARG:HH22	1.85	0.41
2:B:278:PHE:HD1	2:B:308:PHE:HB2	1.85	0.41
1:C:474:ILE:HD13	1:C:487:ALA:HA	2.03	0.41
1:C:664:ARG:HG3	1:C:665:PHE:HD1	1.85	0.41
1:A:681:ARG:O	1:A:685:PRO:HD3	2.21	0.41
2:B:233:GLN:HG2	2:B:234:HIS:ND1	2.36	0.41
1:C:510:PRO:N	1:C:511:PRO:HD2	2.36	0.41
1:A:268:PHE:CZ	1:A:343:VAL:HG21	2.56	0.41
1:C:309:LEU:HB2	1:C:341:GLY:HA3	2.02	0.41
1:A:259:TYR:CD2	1:A:288:GLY:HA3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:510:PRO:N	1:A:511:PRO:HD2	2.36	0.41
1:A:681:ARG:HB2	1:A:720:LEU:HD11	2.03	0.41
1:C:107:VAL:HG22	1:C:113:GLN:HG3	2.02	0.41
1:C:354:ASN:HA	1:C:357:VAL:HG22	2.02	0.41
1:C:409:CYS:SG	1:C:411:PRO:HD2	2.61	0.41
2:B:206:PRO:HA	2:B:209:LEU:HG	2.03	0.41
1:C:259:TYR:CD2	1:C:288:GLY:HA3	2.56	0.41
1:C:531:SER:HG	1:C:586:ASN:H	1.69	0.41
1:A:77:VAL:HG13	1:A:117:GLN:HB2	2.03	0.40
1:C:215:ARG:HA	1:C:254:VAL:HG11	2.02	0.40
2:B:255:GLU:O	2:B:259:LEU:HG	2.21	0.40
1:C:68:MET:HG3	1:C:69:HIS:CD2	2.56	0.40
1:C:138:ARG:NH2	1:C:151:ASP:HB3	2.37	0.40
1:C:650:ARG:O	1:C:650:ARG:HG2	2.21	0.40
1:A:107:VAL:HG22	1:A:113:GLN:HG3	2.03	0.40
2:B:255:GLU:OE1	2:B:255:GLU:N	2.41	0.40
1:C:516:ASN:ND2	1:C:683:ARG:HB3	2.29	0.40
1:A:68:MET:HG3	1:A:69:HIS:CD2	2.56	0.40
1:A:374:PHE:HE1	1:A:588:GLN:HE21	1.68	0.40
1:C:138:ARG:HH22	1:C:151:ASP:HB3	1.85	0.40
2:D:281:TRP:HB3	2:D:313:TYR:HB2	2.04	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	632/1197 (53%)	590 (93%)	42 (7%)	0	100	100
1	C	632/1197 (53%)	592 (94%)	40 (6%)	0	100	100
2	B	171/215 (80%)	166 (97%)	5 (3%)	0	100	100
2	D	171/215 (80%)	163 (95%)	8 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1606/2824 (57%)	1511 (94%)	95 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	541/985 (55%)	539 (100%)	2 (0%)	89	94
1	C	541/985 (55%)	538 (99%)	3 (1%)	84	92
2	B	156/190 (82%)	155 (99%)	1 (1%)	84	92
2	D	156/190 (82%)	156 (100%)	0	100	100
All	All	1394/2350 (59%)	1388 (100%)	6 (0%)	88	94

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	650	ARG
1	A	681	ARG
2	B	212	ARG
1	C	650	ARG
1	C	681	ARG
1	C	726	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	689	GLN
2	B	332	GLN
1	C	494	GLN
1	C	673	GLN
1	C	689	GLN
1	C	691	GLN

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Mol	Chain	Res	Type
2	D	332	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

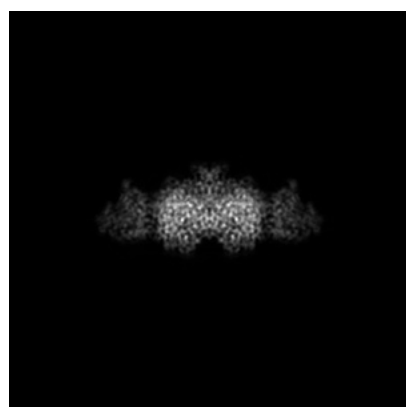
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24709. These allow visual inspection of the internal detail of the map and identification of artifacts.

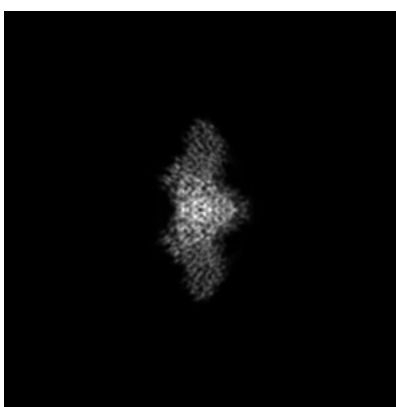
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

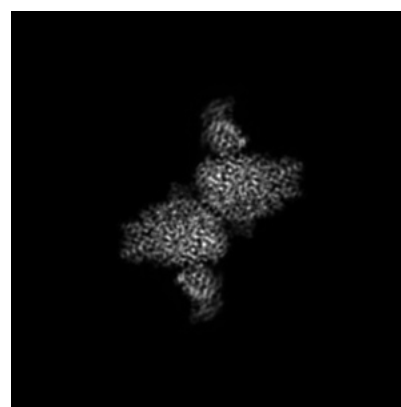
6.1.1 Primary map



X



Y

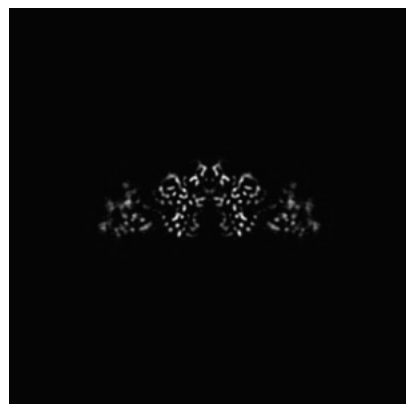


Z

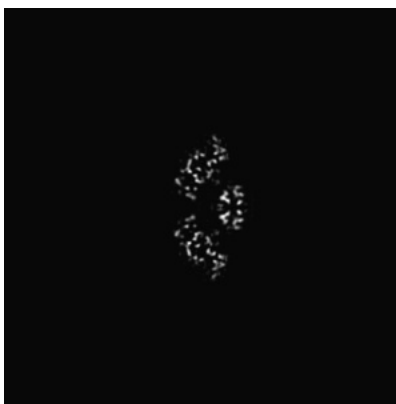
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 168



Y Index: 168



Z Index: 168

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: 168



Y Index: 139

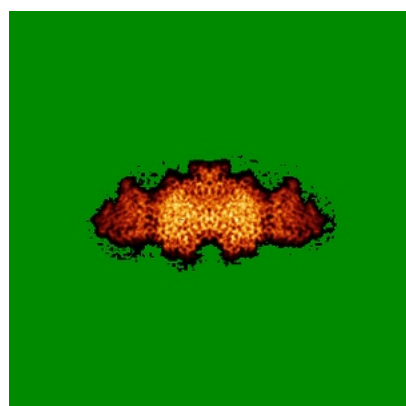


Z Index: 160

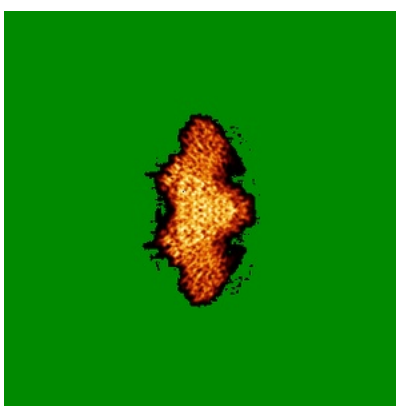
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

6.4.1 Primary map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 10.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

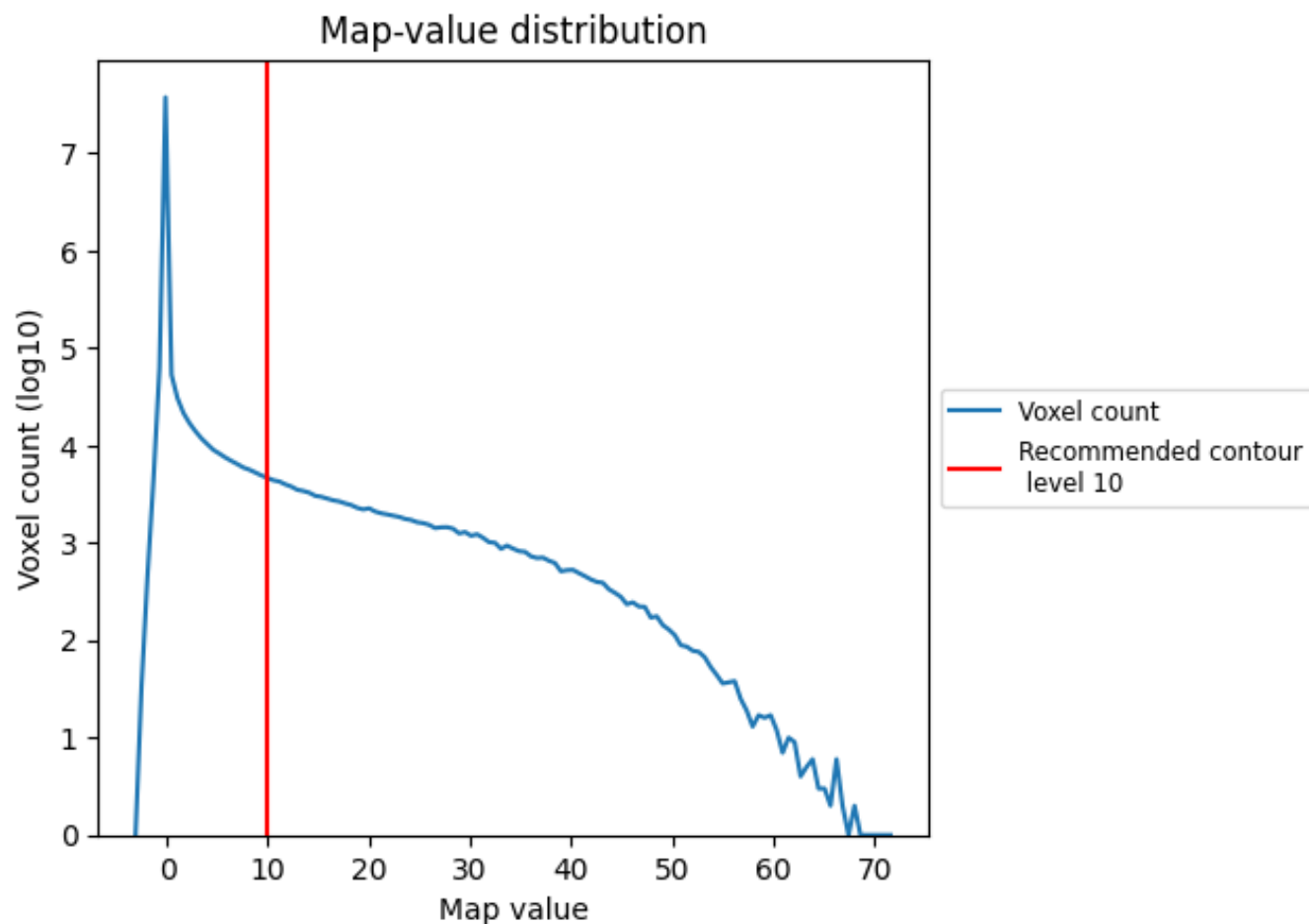
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

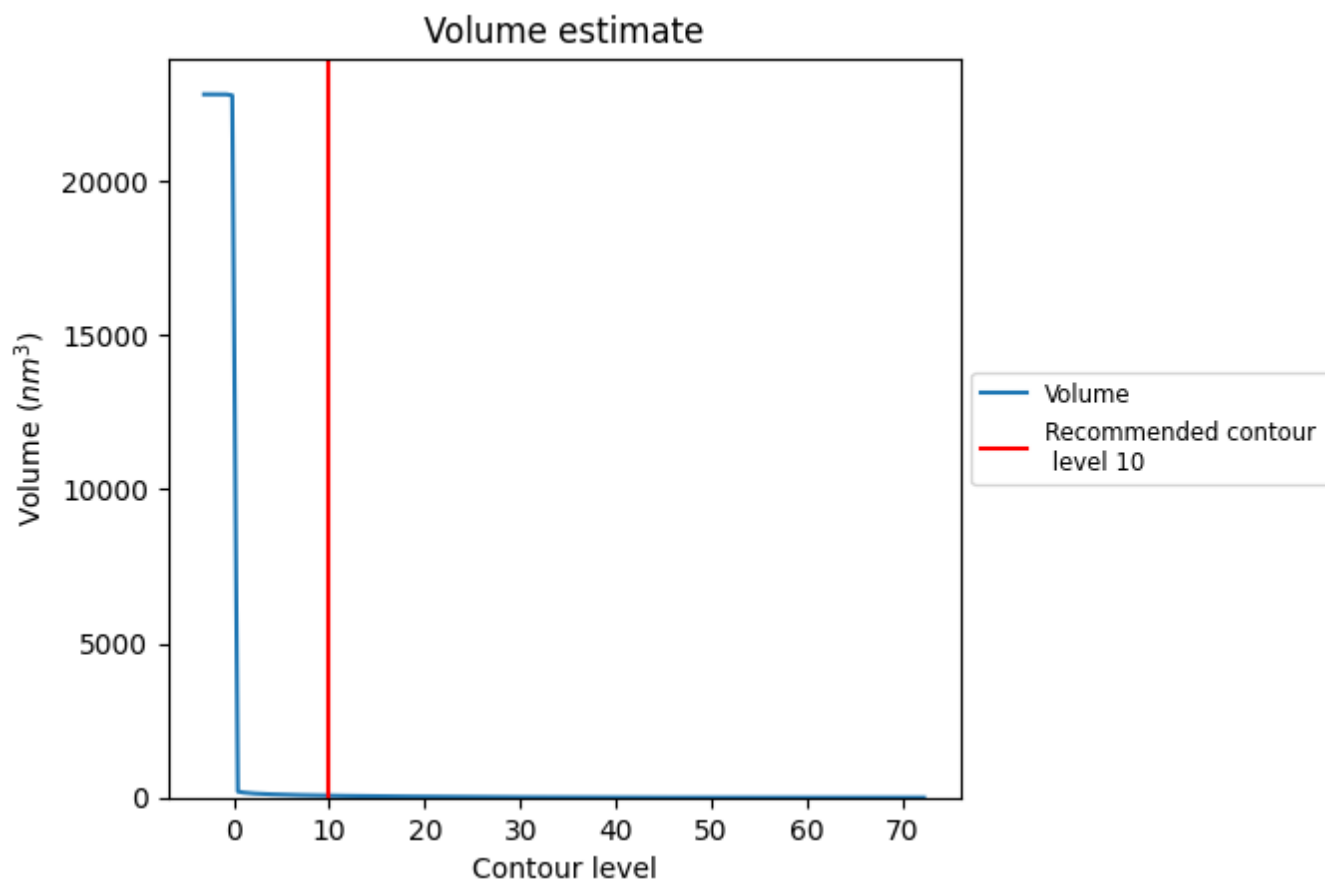
This section contains the results of statistical analysis of the map.

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.

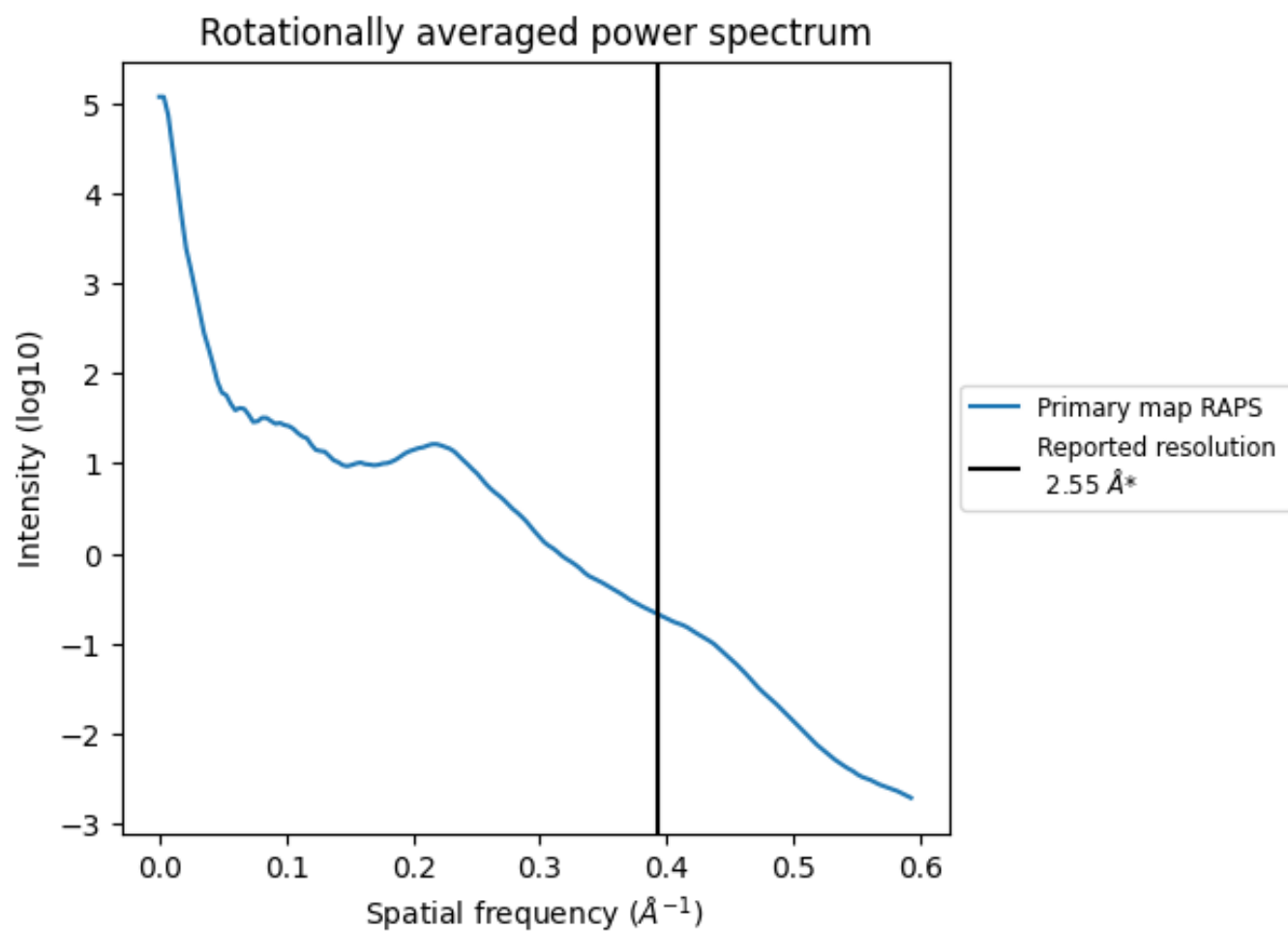
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 62 nm^3 ; this corresponds to an approximate mass of 56 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 0.392 Å⁻¹

8 Fourier-Shell correlation

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

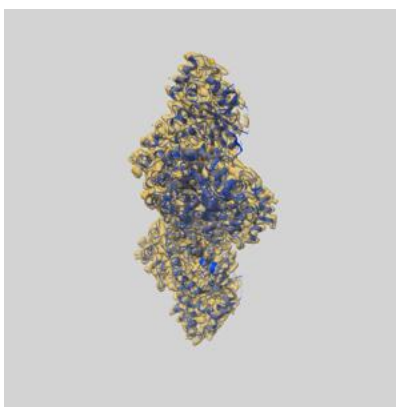
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-24709 and PDB model 7RW6. Per-residue inclusion information can be found in section [3](#) on page [6](#).

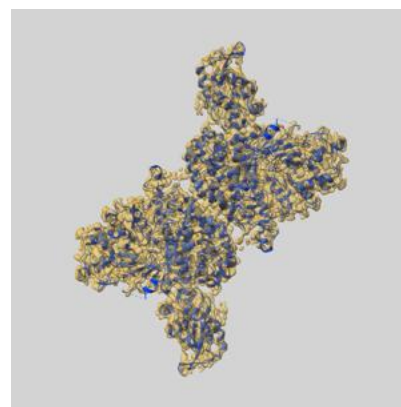
9.1 Map-model overlay [i](#)



X



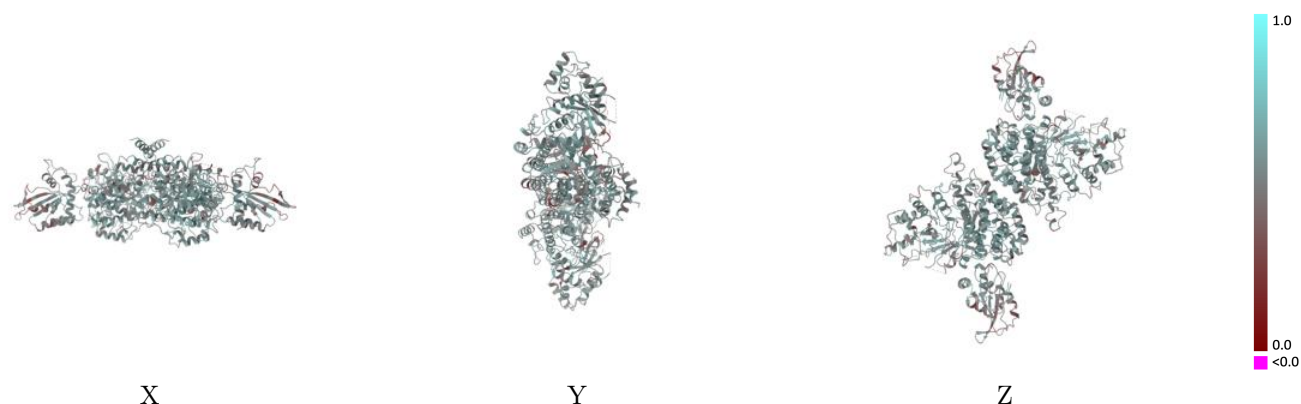
Y



Z

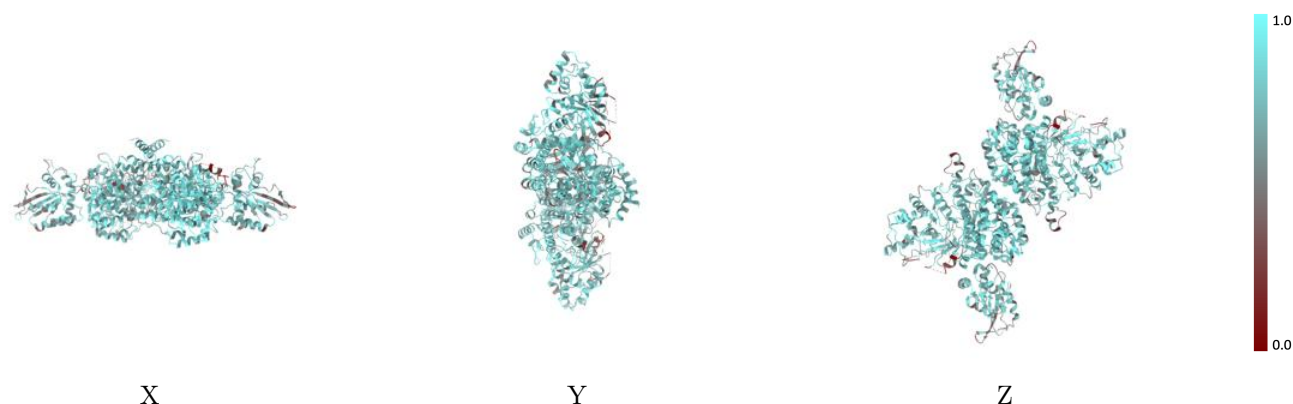
The images above show the 3D surface view of the map at the recommended contour level 10.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



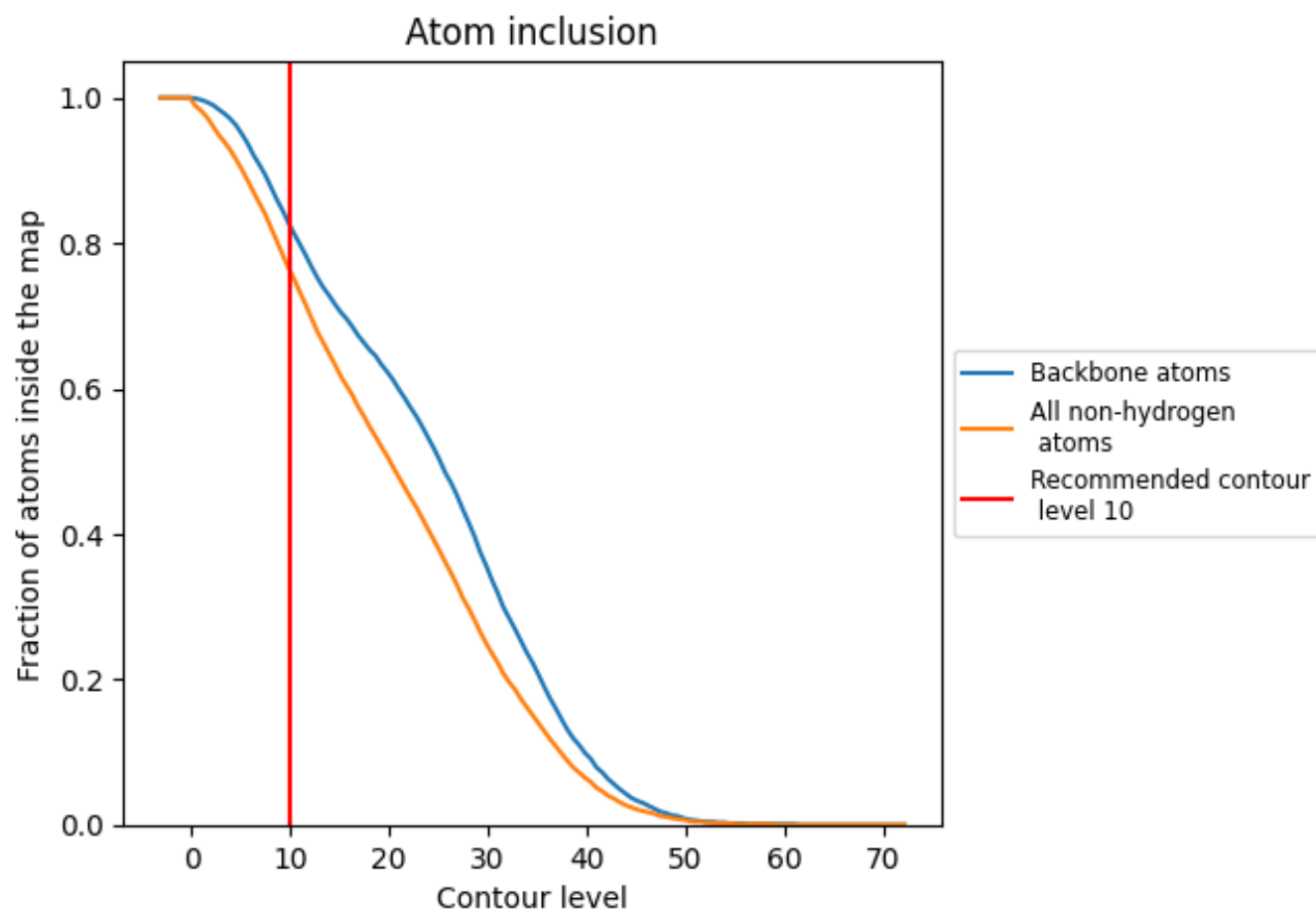
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (10).

9.4 Atom inclusion ⓘ



At the recommended contour level, 82% of all backbone atoms, 76% 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 (10) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.7620	<div></div> 0.5330
A	<div></div> 0.7790	<div></div> 0.5480
B	<div></div> 0.7000	<div></div> 0.4890
C	<div></div> 0.7820	<div></div> 0.5440
D	<div></div> 0.6980	<div></div> 0.4870

