



## Full wwPDB EM Validation Report ⓘ

Apr 22, 2024 – 10:25 AM JST

PDB ID : 8XJ8  
EMDB ID : EMD-38396  
Title : The Cryo-EM structure of MPXV E5 C-terminal in complex with DNA  
Authors : Zhang, W.; Liu, Y.; Gao, H.; Gan, J.  
Deposited on : 2023-12-20  
Resolution : 2.67 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

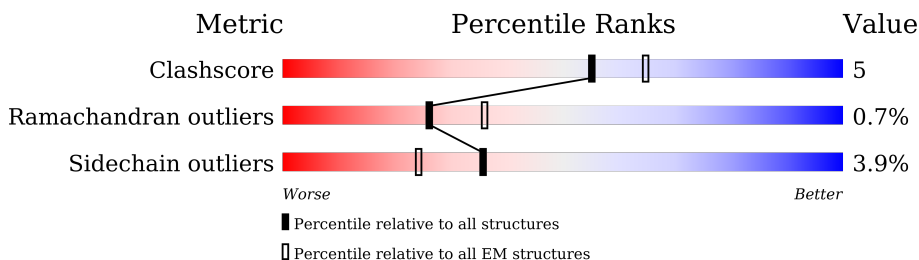
EMDB validation analysis : 0.0.1.dev92  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
MolProbity : 4.02b-467  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36.2

# 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 2.67 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion  $< 40\%$ ). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	X	70	
2	A	463	
2	B	463	
2	C	463	
2	D	463	
2	E	463	
2	F	463	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 15904 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.

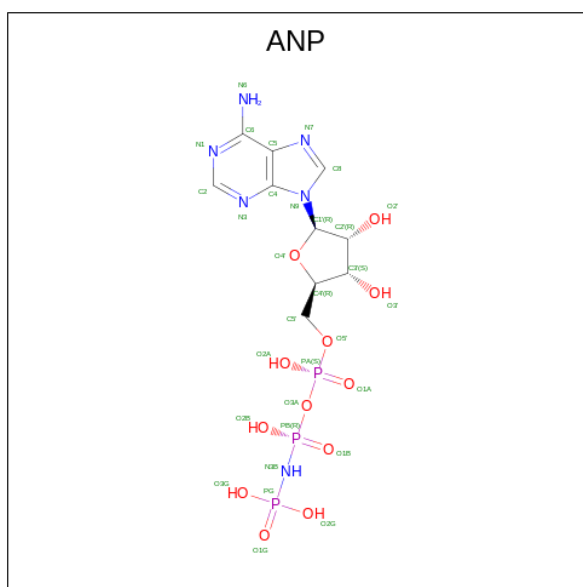
- Molecule 1 is a DNA chain called DNA (70-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	X	6	120	60	12	42	6	0	0

- Molecule 2 is a protein called Monkeypox virus E5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	A	422	3427	2196	581	634	16	0	0
2	B	372	2990	1902	512	560	16	0	0
2	C	379	3053	1950	519	568	16	0	0
2	D	353	2846	1823	484	524	15	0	0
2	E	123	979	629	167	179	4	0	0
2	F	308	2393	1537	398	447	11	0	0

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>6</sub>O<sub>12</sub>P<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
3	A	1	Total	C	N	O	P	0
			31	10	6	12	3	
3	B	1	Total	C	N	O	P	0
			31	10	6	12	3	
3	C	1	Total	C	N	O	P	0
			31	10	6	12	3	

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
			Total	Mg	
4	A	1	Total	Mg	0
			1	1	
4	B	1	Total	Mg	0
			1	1	
4	C	1	Total	Mg	0
			1	1	

### 3 Residue-property plots [i](#)


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.

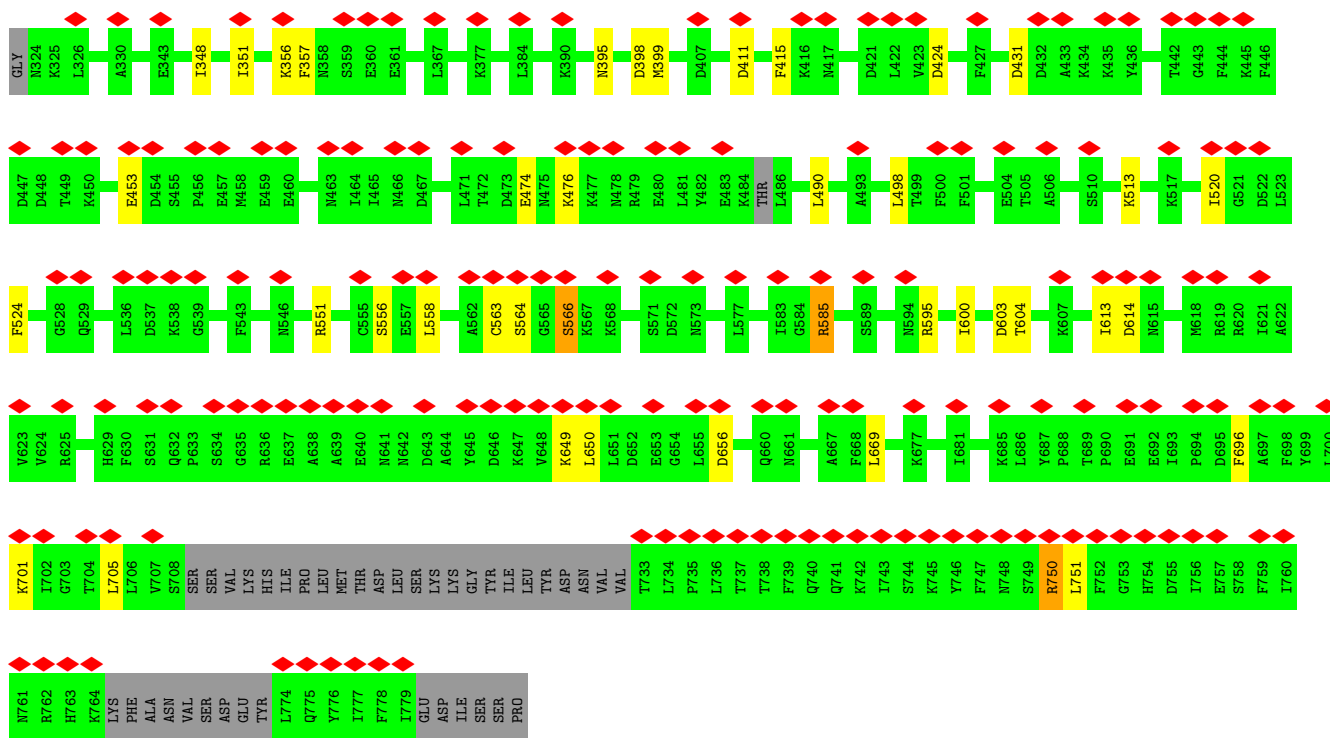
- Molecule 1: DNA (70-MER)

Chain X:  7% . 91%

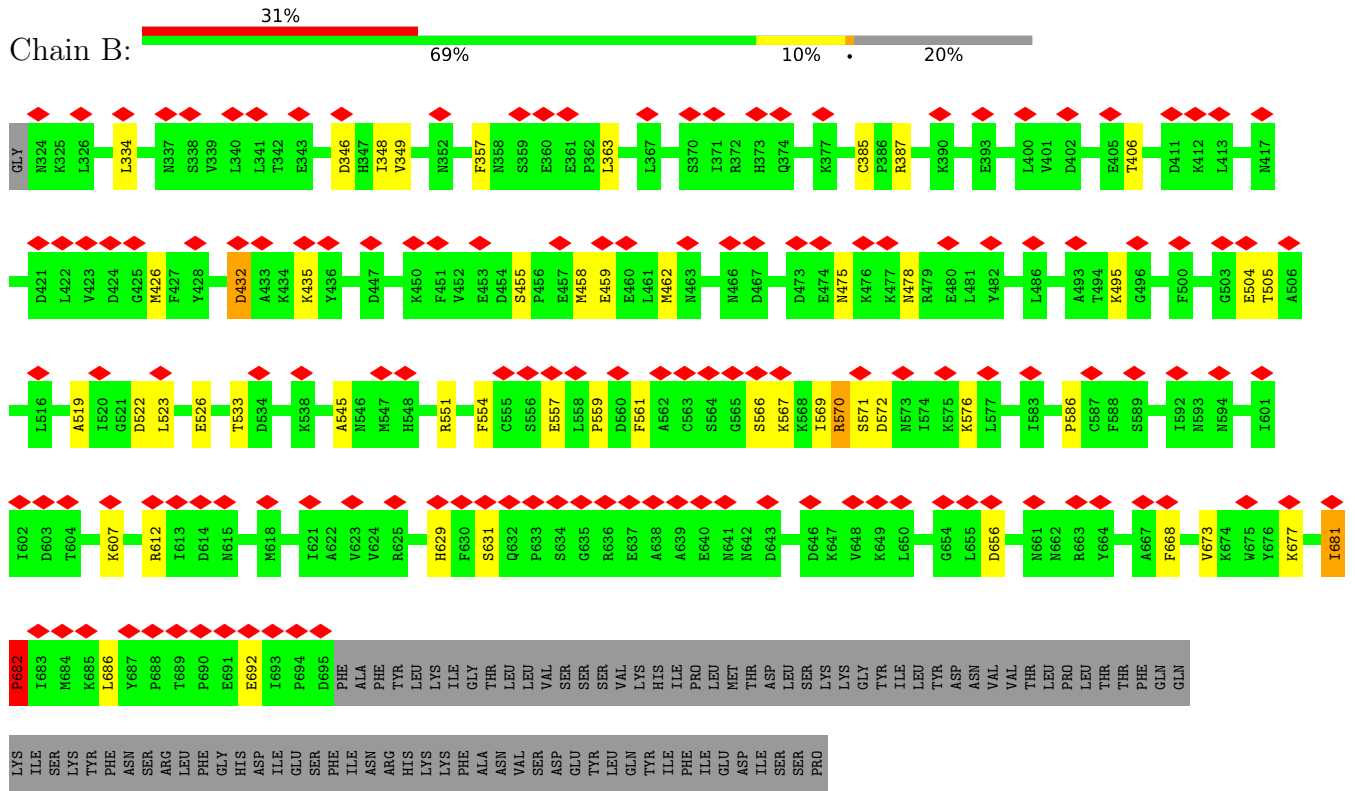


- Molecule 2: Monkeypox virus E5

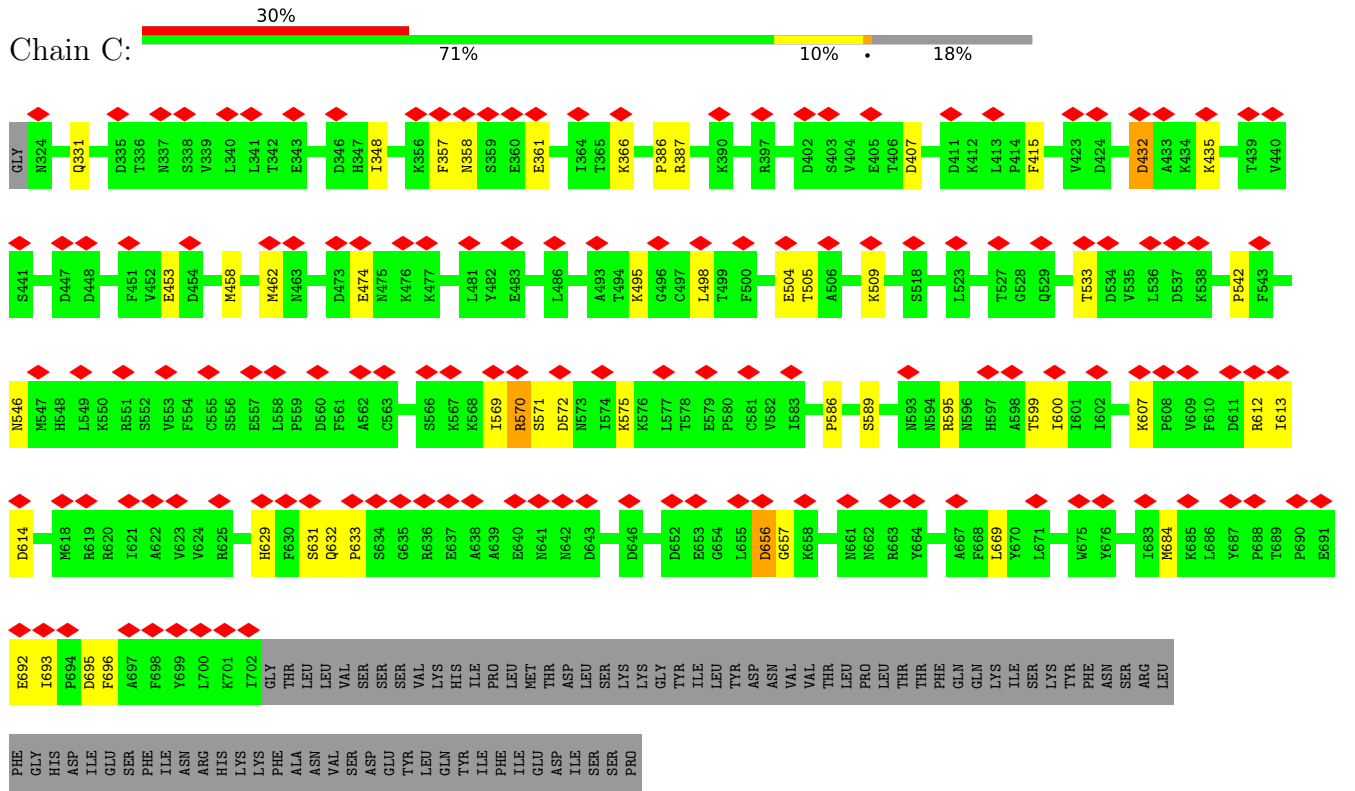
Chain A:  37% 82% 8% . 9%



- Molecule 2: Monkeypox virus E5

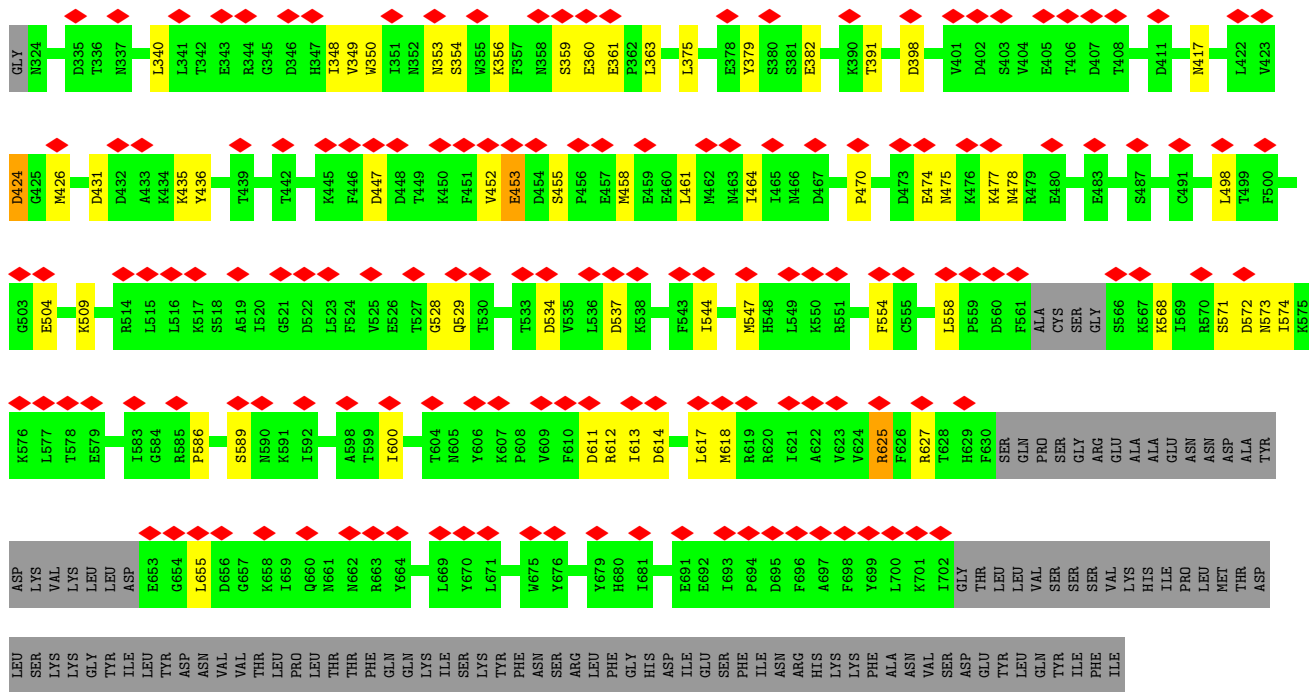


• Molecule 2: Monkeypox virus E5

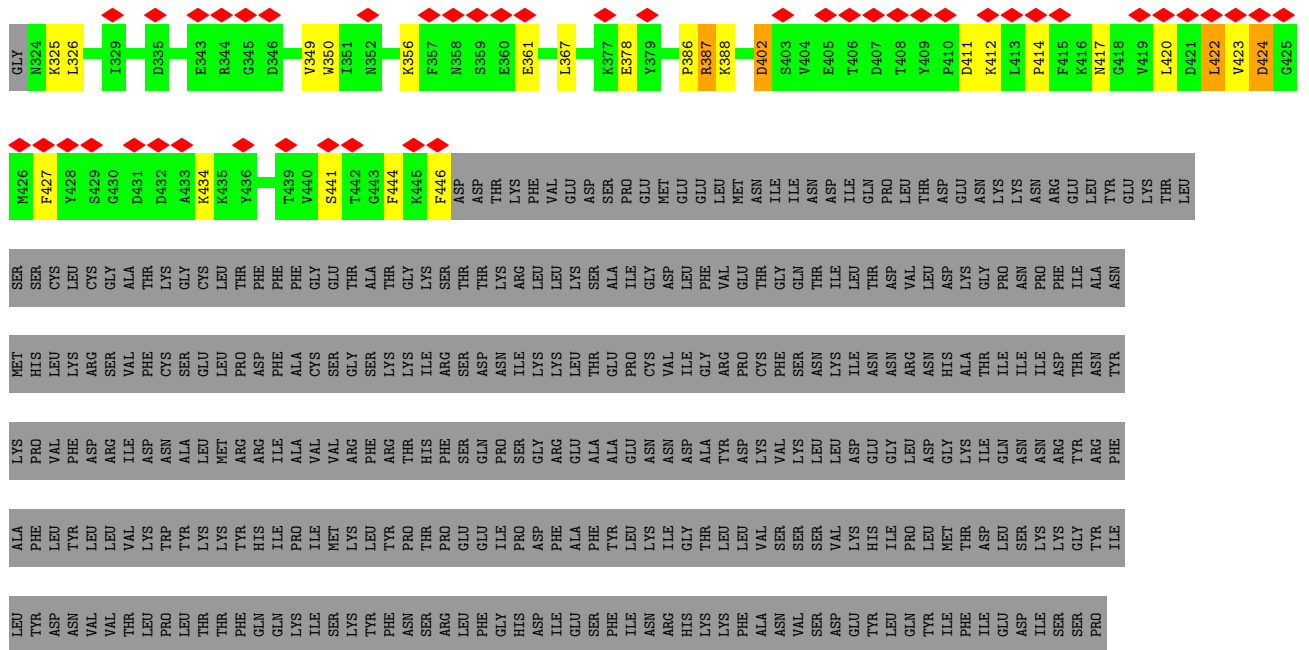


• Molecule 2: Monkeypox virus E5

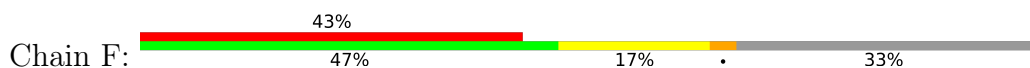




• Molecule 2: Monkeypox virus E5



• Molecule 2: Monkeypox virus E5



THR	LEU	PRO	LEU	LEU	THR	THR	PHE	GLN	GLN	LYS	SER	TYR	THR	PRO	THR	PRO	GLY	HIS	GLU	HIS	LYS	LYS	PHE	ALA	ALA	GLU	ASN	ASN	ASN	THR	LEU	VAL	VAL	VAL	ASP	GLU	GLU	TYR	TYR	GLN	TYR	ILE	ILE	ILE	ILE	GLU	THR	ILE	SER	PRO														
V673	K674	M675	V676	K677	K678	V679	H680	P682	I683	M684	LYS	LEU	TYR	ARG	LEU	THR	PRO	GLY	HIS	GLU	ASP	ILE	ILE	ASP	ASP	GLU	GLY	ASN	ASP	THR	LEU	VAL	VAL	ASP	GLU	GLU	TYR	TYR	GLN	TYR	ILE	ILE	ILE	ILE	GLU	THR	ILE	SER	PRO															
G491	G492	A493	T494	K495	G496	C497	L498	T499	F500	F501	F502	G503	E504	T505	A506	T507	G508	K509	S510	T511	T512	K513	R514	L515	L516	K517	S518	A519	I520	E461	M462	M463	I464	I465	I466	D467	I468	Q469	P470	L471	T472	D473	P474	E474	M475	K476	M477	M478	R479	E480	L481	Y482	E483	L486	S487	S488	C489	L490						
C491	G492	A493	T494	K495	G496	C497	L498	T499	F500	F501	F502	G503	E504	T505	A506	T507	G508	K509	S510	T511	T512	K513	R514	L515	L516	K517	S518	A519	I520	E461	M462	M463	I464	I465	I466	D467	I468	Q469	P470	L471	T472	D473	P474	E474	M475	K476	M477	M478	R479	E480	L481	Y482	E483	L486	S487	S488	C489	L490						
F554	C555	S556	E557	L558	P559	D560	F561	A562	SER	GLY	SER	K567	K568	I569	R570	S571	D572	N573	I574	K575	K576	L577	T578	E579	PRO	CYS	VAL	ILE	GLY	ARG	PRO	CYS	PHE	ASN	ASN	ILE	ASN	ASN	ARG	N596	H597	A598	T599	I600	I601	I602	D603	T604	N605	Y606	K607	F608	V609	F610	D611	H612								
I613	D614	N615	A616	L617	M618	R619	R620	I621	A622	V623	V624	R625	F626	R627	T628	H629	PHE	SER	GLN	PRO	PRO	SER	GLY	ALA	ARG	GLU	ALA	ALA	GLU	ASN	ASN	ASP	VAL	ALA	TYR	ASP	LYS	VAL	LYS	LEU	SER	HIS	ILE	PRO	MET	THR	ASP	LEU	SER	LYS	GLY	TYR	ILE	LEU	R663	Y664	R665	F666	A667	F668	L669	Y670	L671	L672
V673	K674	M675	V676	K677	K678	V679	H680	P682	I683	M684	LYS	LEU	TYR	ARG	LEU	THR	PRO	GLY	HIS	GLU	ASP	ILE	ILE	ASP	ASP	GLU	GLY	ASN	ASP	THR	LEU	VAL	VAL	ASP	GLU	GLU	TYR	TYR	GLN	TYR	ILE	ILE	ILE	ILE	GLU	THR	ILE	SER	PRO															
GLY	R324	I333	T336	N337	E343	R344	G345	D346	H347	I348	V349	M350	I351	N352	N353	S354	F357	N358	S359	E360	E361	P362	T365	K366	K377	R389	K390	T391	A394	S403	V404	E405	T406	D407	T408	D411	K412	L413	P414	F415	K416	L420	D421	L422	V423	D424	G425	M426																



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	247898	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	50.0	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	8.140	Depositor
Minimum map value	-3.638	Depositor
Average map value	-0.002	Depositor
Map value standard deviation	0.159	Depositor
Recommended contour level	0.733	Depositor
Map size ( $\text{\AA}$ )	302.68, 302.68, 302.68	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.081, 1.081, 1.081	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: ANP, MG

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	X	0.44	0/131	1.25	0/200
2	A	0.24	0/3498	0.46	0/4718
2	B	0.30	0/3050	0.53	2/4119 (0.0%)
2	C	0.24	0/3117	0.50	0/4210
2	D	0.25	0/2905	0.48	0/3924
2	E	0.24	0/999	0.48	0/1353
2	F	0.26	0/2437	0.52	0/3301
All	All	0.26	0/16137	0.51	2/21825 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	B	682	PRO	N-CA-CB	7.33	112.09	103.30
2	B	682	PRO	N-CA-C	-6.20	95.97	112.10

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	X	120	0	73	1	0
2	A	3427	0	3451	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2990	0	3000	27	0
2	C	3053	0	3070	28	0
2	D	2846	0	2857	33	0
2	E	979	0	982	14	0
2	F	2393	0	2322	54	0
3	A	31	0	13	0	0
3	B	31	0	13	0	0
3	C	31	0	13	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
All	All	15904	0	15794	168	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 5.

All (168) 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:F:343:GLU:N	2:F:343:GLU:OE1	2.19	0.75
2:D:504:GLU:O	2:D:509:LYS:NZ	2.28	0.67
2:D:361:GLU:N	2:D:361:GLU:OE1	2.28	0.66
2:F:453:GLU:HA	2:F:670:TYR:CE2	2.30	0.66
2:C:542:PRO:O	2:C:546:ASN:ND2	2.30	0.65
2:F:655:LEU:O	2:F:659:ILE:N	2.29	0.65
2:F:574:ILE:H	2:F:574:ILE:HD12	1.63	0.64
1:X:2:DT:OP1	2:A:585:ARG:NH2	2.31	0.63
2:E:378:GLU:OE1	2:E:378:GLU:N	2.29	0.62
2:C:386:PRO:HD2	2:C:387:ARG:HH21	1.65	0.61
2:F:459:GLU:N	2:F:459:GLU:OE2	2.33	0.61
2:A:513:LYS:NZ	2:A:556:SER:OG	2.33	0.61
2:F:348:ILE:HG22	2:F:357:PHE:HB2	1.83	0.61
2:F:432:ASP:N	2:F:432:ASP:OD1	2.32	0.60
2:B:571:SER:OG	2:B:612:ARG:O	2.21	0.58
2:D:349:VAL:HG21	2:D:363:LEU:HB3	1.85	0.58
2:C:656:ASP:OD1	2:C:657:GLY:N	2.35	0.58
2:E:402:ASP:OD1	2:E:402:ASP:N	2.33	0.58
2:E:387:ARG:H	2:E:387:ARG:HD2	1.69	0.57
2:F:479:ARG:O	2:F:483:GLU:HG3	2.05	0.57
2:F:664:TYR:O	2:F:666:PHE:N	2.37	0.57
2:D:359:SER:OG	2:D:360:GLU:OE1	2.18	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:476:LYS:O	2:F:480:GLU:HG2	2.04	0.56
2:E:414:PRO:HG2	2:E:441:SER:HA	1.88	0.56
2:C:632:GLN:HG3	2:C:633:PRO:HD2	1.88	0.56
2:F:459:GLU:HA	2:F:462:MET:HG2	1.87	0.56
2:B:673:VAL:HG12	2:B:677:LYS:HE3	1.88	0.56
2:D:498:LEU:HG	2:D:600:ILE:HB	1.88	0.56
2:F:571:SER:H	2:F:611:ASP:HB3	1.69	0.56
2:B:348:ILE:HG22	2:B:357:PHE:HB2	1.86	0.55
2:A:348:ILE:HG22	2:A:357:PHE:HB3	1.88	0.55
2:C:348:ILE:HG22	2:C:357:PHE:HB3	1.90	0.54
2:D:534:ASP:O	2:D:573:ASN:ND2	2.32	0.54
2:D:612:ARG:NH1	2:D:613:ILE:O	2.40	0.54
2:B:455:SER:O	2:B:459:GLU:OE2	2.26	0.54
2:F:562:ALA:HB2	2:F:609:VAL:HG21	1.91	0.53
2:C:474:GLU:N	2:C:474:GLU:OE2	2.41	0.53
2:B:523:LEU:HD13	2:B:551:ARG:HD2	1.91	0.53
2:B:607:LYS:NZ	2:B:692:GLU:O	2.40	0.52
2:D:361:GLU:HG2	2:D:363:LEU:HG	1.91	0.52
2:A:564:SER:OG	2:A:566:SER:OG	2.28	0.52
2:C:612:ARG:NH1	2:C:613:ILE:O	2.43	0.52
2:D:417:ASN:N	2:D:417:ASN:OD1	2.43	0.52
2:D:528:GLY:O	2:D:529:GLN:NE2	2.43	0.51
2:F:490:LEU:HD11	2:F:672:LEU:HB3	1.92	0.51
2:A:395:ASN:O	2:A:399:MET:HG3	2.11	0.51
2:B:504:GLU:HG3	2:B:505:THR:H	1.76	0.50
2:C:569:ILE:O	2:C:570:ARG:HB2	2.12	0.50
2:A:351:ILE:O	2:A:356:LYS:NZ	2.45	0.50
2:E:423:VAL:HG13	2:E:424:ASP:OD1	2.12	0.50
2:C:504:GLU:O	2:C:509:LYS:NZ	2.44	0.50
2:B:557:GLU:HG2	2:C:575:LYS:HB3	1.93	0.50
2:B:569:ILE:O	2:B:570:ARG:HB2	2.12	0.50
2:F:495:LYS:NZ	2:F:597:HIS:O	2.37	0.50
2:C:415:PHE:HZ	2:C:669:LEU:HD21	1.76	0.49
2:D:625:ARG:HD3	2:D:627:ARG:HH21	1.78	0.49
2:C:366:LYS:HG2	2:D:398:ASP:HA	1.95	0.49
2:F:467:ASP:OD1	2:F:468:ILE:HD12	2.13	0.49
2:F:486:LEU:HD21	2:F:668:PHE:CZ	2.48	0.49
2:A:398:ASP:OD2	2:F:389:ARG:NH2	2.45	0.49
2:D:478:ASN:HD21	2:D:625:ARG:H	1.61	0.49
2:F:495:LYS:HE3	2:F:599:THR:HG22	1.95	0.49
2:B:458:MET:O	2:B:462:MET:HG2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:354:SER:OG	2:D:356:LYS:NZ	2.44	0.48
2:D:375:LEU:HD23	2:D:379:TYR:HB3	1.94	0.48
2:C:498:LEU:HD22	2:C:600:ILE:HB	1.96	0.48
2:F:509:LYS:NZ	2:F:604:THR:O	2.47	0.48
2:F:670:TYR:CE1	2:F:674:LYS:HG3	2.49	0.48
2:D:571:SER:O	2:D:574:ILE:HG13	2.14	0.47
2:A:498:LEU:HG	2:A:600:ILE:HB	1.96	0.47
2:A:750:ARG:NH1	2:A:751:LEU:HB2	2.30	0.47
2:A:613:ILE:HD12	2:A:696:PHE:HE1	1.80	0.47
2:E:420:LEU:HA	2:E:427:PHE:HA	1.96	0.47
2:F:350:TRP:CD2	2:F:434:LYS:HG3	2.49	0.47
2:B:561:PHE:HD1	2:B:566:SER:HB2	1.80	0.47
2:F:487:SER:OG	2:F:680:HIS:NE2	2.46	0.47
2:F:525:VAL:HB	2:F:550:LYS:HD2	1.96	0.47
2:A:490:LEU:O	2:A:551:ARG:NH1	2.48	0.47
2:B:572:ASP:O	2:B:576:LYS:HG3	2.15	0.47
2:C:533:THR:HA	2:C:569:ILE:O	2.15	0.47
2:C:570:ARG:NH2	2:C:572:ASP:OD2	2.48	0.47
2:B:432:ASP:OD1	2:B:432:ASP:N	2.47	0.46
2:F:669:LEU:O	2:F:673:VAL:HG23	2.14	0.46
2:B:629:HIS:CE1	2:B:631:SER:HB3	2.51	0.46
2:B:561:PHE:CD1	2:B:566:SER:HB2	2.50	0.46
2:E:386:PRO:HG2	2:F:391:THR:HG23	1.97	0.46
2:D:464:ILE:HG12	2:D:655:LEU:HD21	1.97	0.46
2:F:513:LYS:NZ	2:F:603:ASP:OD1	2.49	0.46
2:E:349:VAL:HG13	2:E:367:LEU:HD22	1.98	0.45
2:F:522:ASP:O	2:F:550:LYS:NZ	2.27	0.45
2:D:614:ASP:O	2:D:618:MET:HG2	2.17	0.45
2:B:559:PRO:HD2	2:B:561:PHE:CZ	2.50	0.45
2:D:470:PRO:HD2	2:D:475:ASN:HD22	1.81	0.45
2:B:545:ALA:HB1	2:B:586:PRO:HG3	1.98	0.45
2:C:432:ASP:N	2:C:432:ASP:OD1	2.49	0.45
2:F:671:LEU:HD23	2:F:671:LEU:HA	1.82	0.45
2:A:520:ILE:HD11	2:A:524:PHE:HB2	1.97	0.45
2:C:458:MET:O	2:C:462:MET:HG2	2.17	0.45
2:F:498:LEU:HD13	2:F:574:ILE:HG23	1.97	0.45
2:F:527:THR:OG1	2:F:528:GLY:N	2.49	0.45
2:A:558:LEU:HB2	2:A:604:THR:HG22	1.99	0.45
2:C:607:LYS:NZ	2:C:692:GLU:O	2.33	0.45
2:F:574:ILE:HA	2:F:577:LEU:HB2	1.99	0.45
2:D:424:ASP:HB3	2:D:426:MET:HG2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:504:GLU:HG3	2:C:505:THR:H	1.82	0.44
2:F:346:ASP:HB3	2:F:357:PHE:HE1	1.81	0.44
2:C:614:ASP:N	2:C:614:ASP:OD1	2.50	0.44
2:F:677:LYS:HB2	2:F:677:LYS:HE2	1.76	0.44
2:D:452:VAL:O	2:D:453:GLU:HB2	2.17	0.44
2:A:701:LYS:O	2:A:705:LEU:HB2	2.18	0.44
2:B:681:ILE:O	2:B:682:PRO:CB	2.66	0.44
2:F:517:LYS:O	2:F:517:LYS:HD2	2.18	0.44
2:F:412:LYS:HE3	2:F:436:TYR:CE1	2.53	0.43
2:F:666:PHE:HA	2:F:669:LEU:HB2	2.00	0.43
2:A:415:PHE:HZ	2:A:669:LEU:HD21	1.83	0.43
2:B:533:THR:HA	2:B:569:ILE:O	2.17	0.43
2:F:377:LYS:HB3	2:F:377:LYS:HE3	1.80	0.43
2:E:326:LEU:HD11	2:E:388:LYS:HG2	1.99	0.43
2:B:567:LYS:HB3	2:B:567:LYS:HE2	1.62	0.43
2:C:571:SER:OG	2:C:612:ARG:O	2.33	0.43
2:D:350:TRP:CZ2	2:D:353:ASN:HA	2.53	0.43
2:F:664:TYR:HB3	2:F:667:ALA:HB3	1.99	0.43
2:E:422:LEU:HD22	2:E:422:LEU:HA	1.86	0.43
2:B:519:ALA:HB2	2:B:668:PHE:HD2	1.84	0.43
2:F:457:GLU:HA	2:F:460:GLU:HG3	2.00	0.43
2:F:513:LYS:HG2	2:F:554:PHE:CE2	2.54	0.43
2:F:577:LEU:HD23	2:F:577:LEU:HA	1.86	0.43
2:E:417:ASN:N	2:E:417:ASN:OD1	2.52	0.43
2:B:385:CYS:SG	2:B:387:ARG:NH2	2.92	0.43
2:B:495:LYS:HB3	2:B:686:LEU:HD12	2.00	0.43
2:F:510:SER:HA	2:F:513:LYS:HB2	2.01	0.43
2:C:407:ASP:O	2:C:595:ARG:NH2	2.50	0.42
2:F:461:LEU:HB3	2:F:664:TYR:CE2	2.54	0.42
2:F:462:MET:O	2:F:466:ASN:HB2	2.19	0.42
2:D:571:SER:HB3	2:D:617:LEU:HD13	2.01	0.42
2:F:600:ILE:HG22	2:F:602:ILE:HD11	2.01	0.42
2:C:693:ILE:HG22	2:C:696:PHE:H	1.84	0.42
2:B:349:VAL:HG21	2:B:363:LEU:HB3	2.01	0.42
2:B:475:ASN:OD1	2:B:478:ASN:HB3	2.20	0.42
2:D:475:ASN:OD1	2:D:478:ASN:HB3	2.19	0.42
2:D:568:LYS:HD2	2:D:611:ASP:OD1	2.20	0.42
2:E:417:ASN:ND2	2:E:446:PHE:O	2.52	0.42
2:D:544:ILE:HA	2:D:547:MET:HG3	2.00	0.42
2:D:435:LYS:HE3	2:D:436:TYR:CZ	2.55	0.41
2:F:507:THR:HG22	2:F:509:LYS:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:599:THR:O	2:F:599:THR:OG1	2.38	0.41
2:A:649:LYS:HG2	2:A:650:LEU:H	1.85	0.41
2:D:458:MET:HE3	2:D:461:LEU:HD23	2.02	0.41
2:C:358:ASN:HB3	2:C:361:GLU:O	2.20	0.41
2:E:411:ASP:O	2:E:412:LYS:NZ	2.38	0.41
2:D:477:LYS:HE3	2:D:477:LYS:HB3	1.75	0.41
2:D:586:PRO:HD2	2:D:589:SER:HB3	2.02	0.41
2:E:350:TRP:CD2	2:E:434:LYS:HG3	2.55	0.41
2:F:365:THR:HG23	2:F:389:ARG:HB3	2.02	0.41
2:C:386:PRO:HG2	2:D:391:THR:HG23	2.02	0.41
2:F:333:ILE:O	2:F:336:THR:OG1	2.37	0.41
2:F:362:PRO:O	2:F:366:LYS:HG3	2.21	0.41
2:F:530:THR:HA	2:F:533:THR:HG22	2.03	0.41
2:A:411:ASP:OD1	2:A:411:ASP:N	2.47	0.41
2:C:629:HIS:CE1	2:C:631:SER:HB2	2.56	0.41
2:F:450:LYS:HB2	2:F:450:LYS:HE2	1.79	0.41
2:A:563:CYS:HB2	2:B:612:ARG:HD3	2.02	0.40
2:B:435:LYS:HE2	2:B:435:LYS:HB3	1.66	0.40
2:D:340:LEU:HB2	2:D:348:ILE:HG13	2.02	0.40
2:F:453:GLU:HA	2:F:670:TYR:HE2	1.81	0.40
2:D:558:LEU:HD23	2:D:558:LEU:HA	1.96	0.40
2:C:495:LYS:HE3	2:C:599:THR:OG1	2.20	0.40
2:C:586:PRO:HD2	2:C:589:SER:HB3	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	414/463 (89%)	410 (99%)	4 (1%)	0	100	100
2	B	370/463 (80%)	354 (96%)	12 (3%)	4 (1%)	14	31

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	377/463 (81%)	365 (97%)	10 (3%)	2 (0%)	29	52
2	D	347/463 (75%)	328 (94%)	17 (5%)	2 (1%)	25	47
2	E	121/463 (26%)	113 (93%)	7 (6%)	1 (1%)	19	40
2	F	298/463 (64%)	266 (89%)	27 (9%)	5 (2%)	9	20
All	All	1927/2778 (69%)	1836 (95%)	77 (4%)	14 (1%)	26	44

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	570	ARG
2	B	681	ILE
2	B	682	PRO
2	C	453	GLU
2	C	570	ARG
2	F	665	ARG
2	D	453	GLU
2	F	424	ASP
2	F	679	TYR
2	B	406	THR
2	E	361	GLU
2	F	455	SER
2	F	361	GLU
2	D	455	SER

### 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	
2	A	387/425 (91%)	375 (97%)	12 (3%)	40	67
2	B	336/425 (79%)	328 (98%)	8 (2%)	49	75
2	C	343/425 (81%)	337 (98%)	6 (2%)	60	82
2	D	319/425 (75%)	310 (97%)	9 (3%)	43	70
2	E	109/425 (26%)	102 (94%)	7 (6%)	17	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	F	255/425 (60%)	229 (90%)	26 (10%)	7 15
All	All	1749/2550 (69%)	1681 (96%)	68 (4%)	36 58

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	424	ASP
2	A	431	ASP
2	A	453	GLU
2	A	474	GLU
2	A	476	LYS
2	A	566	SER
2	A	585	ARG
2	A	595	ARG
2	A	603	ASP
2	A	614	ASP
2	A	656	ASP
2	A	750	ARG
2	B	334	LEU
2	B	346	ASP
2	B	426	MET
2	B	432	ASP
2	B	522	ASP
2	B	526	GLU
2	B	554	PHE
2	B	656	ASP
2	C	331	GLN
2	C	432	ASP
2	C	435	LYS
2	C	656	ASP
2	C	684	MET
2	C	695	ASP
2	D	382	GLU
2	D	424	ASP
2	D	431	ASP
2	D	447	ASP
2	D	474	GLU
2	D	537	ASP
2	D	554	PHE
2	D	572	ASP
2	D	625	ARG
2	E	325	LYS

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Mol	Chain	Res	Type
2	E	356	LYS
2	E	387	ARG
2	E	402	ASP
2	E	422	LEU
2	E	424	ASP
2	E	444	PHE
2	F	405	GLU
2	F	411	ASP
2	F	432	ASP
2	F	436	TYR
2	F	447	ASP
2	F	461	LEU
2	F	471	LEU
2	F	483	GLU
2	F	495	LYS
2	F	497	CYS
2	F	514	ARG
2	F	516	LEU
2	F	522	ASP
2	F	546	ASN
2	F	554	PHE
2	F	555	CYS
2	F	556	SER
2	F	571	SER
2	F	603	ASP
2	F	604	THR
2	F	623	VAL
2	F	663	ARG
2	F	665	ARG
2	F	669	LEU
2	F	671	LEU
2	F	679	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	529	GLN
2	F	475	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](#)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	ANP	C	801	4	29,33,33	1.05	4 (13%)	31,52,52	1.09	2 (6%)
3	ANP	A	801	4	29,33,33	1.07	4 (13%)	31,52,52	1.06	2 (6%)
3	ANP	B	801	4	29,33,33	1.06	4 (13%)	31,52,52	1.05	2 (6%)

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
3	ANP	C	801	4	-	2/14/38/38	0/3/3/3
3	ANP	A	801	4	-	3/14/38/38	0/3/3/3
3	ANP	B	801	4	-	5/14/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	801	ANP	PG-N3B	2.51	1.69	1.63
3	C	801	ANP	PG-N3B	2.48	1.69	1.63
3	A	801	ANP	PG-N3B	2.46	1.69	1.63
3	A	801	ANP	PG-O1G	2.34	1.49	1.46
3	B	801	ANP	PG-O1G	2.28	1.49	1.46
3	C	801	ANP	PG-O1G	2.27	1.49	1.46
3	A	801	ANP	PB-O3A	-2.24	1.56	1.59
3	B	801	ANP	PB-O3A	-2.20	1.56	1.59
3	C	801	ANP	PB-O3A	-2.17	1.56	1.59
3	A	801	ANP	PB-O1B	2.16	1.49	1.46
3	C	801	ANP	PB-O1B	2.14	1.49	1.46
3	B	801	ANP	PB-O1B	2.11	1.49	1.46

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	801	ANP	PB-O3A-PA	-3.33	120.88	132.62
3	A	801	ANP	PB-O3A-PA	-3.19	121.38	132.62
3	B	801	ANP	PB-O3A-PA	-3.18	121.42	132.62
3	A	801	ANP	C5-C6-N6	2.30	123.84	120.35
3	B	801	ANP	C5-C6-N6	2.28	123.81	120.35
3	C	801	ANP	C5-C6-N6	2.27	123.80	120.35

There are no chirality outliers.

All (10) torsion outliers are listed below:

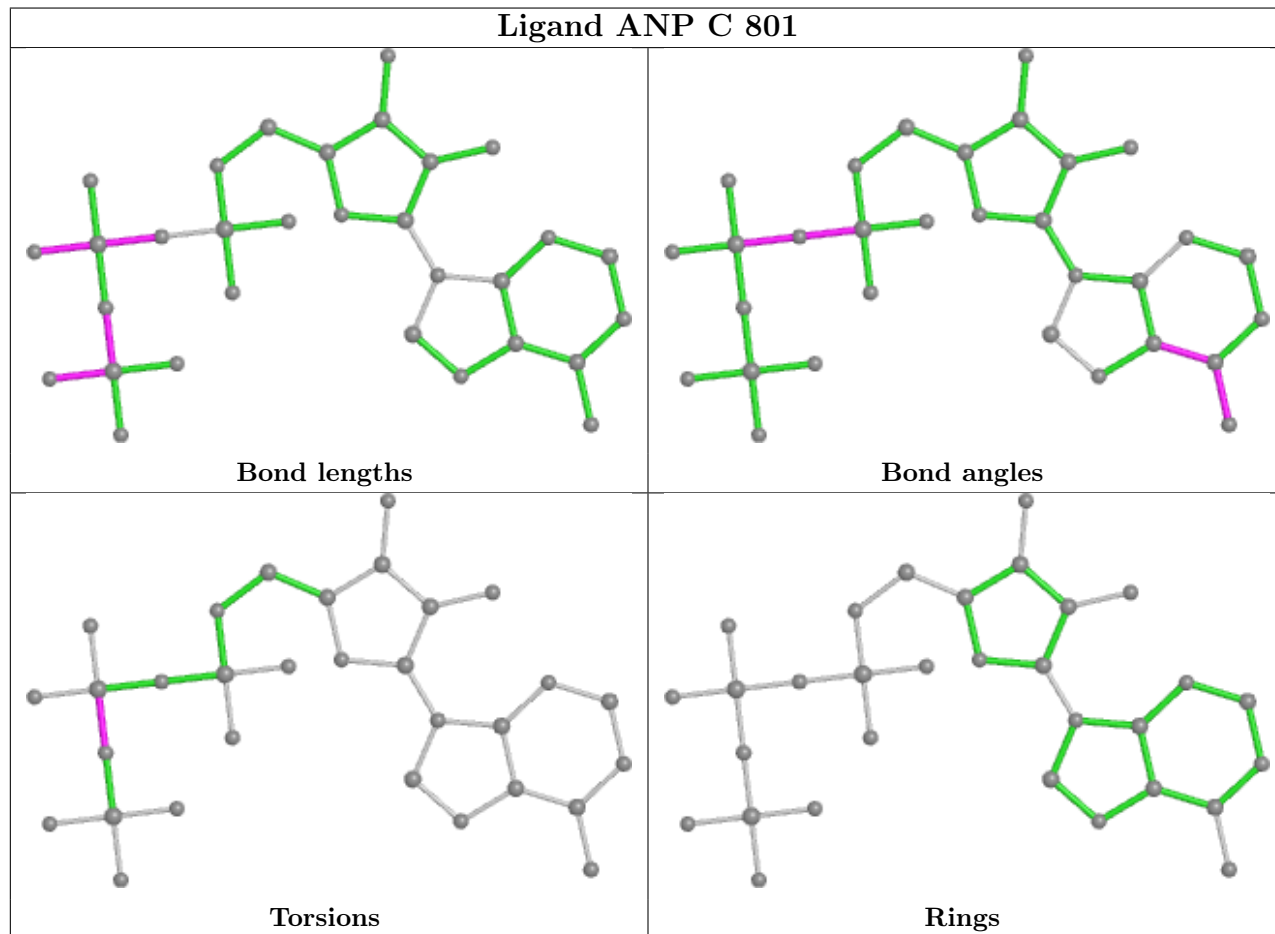
Mol	Chain	Res	Type	Atoms
3	A	801	ANP	PG-N3B-PB-O1B
3	A	801	ANP	PG-N3B-PB-O3A
3	B	801	ANP	PB-N3B-PG-O1G
3	B	801	ANP	PG-N3B-PB-O1B
3	B	801	ANP	PG-N3B-PB-O3A
3	C	801	ANP	PG-N3B-PB-O1B
3	C	801	ANP	PG-N3B-PB-O3A
3	B	801	ANP	C5'-O5'-PA-O3A
3	B	801	ANP	C5'-O5'-PA-O2A
3	A	801	ANP	C5'-O5'-PA-O3A

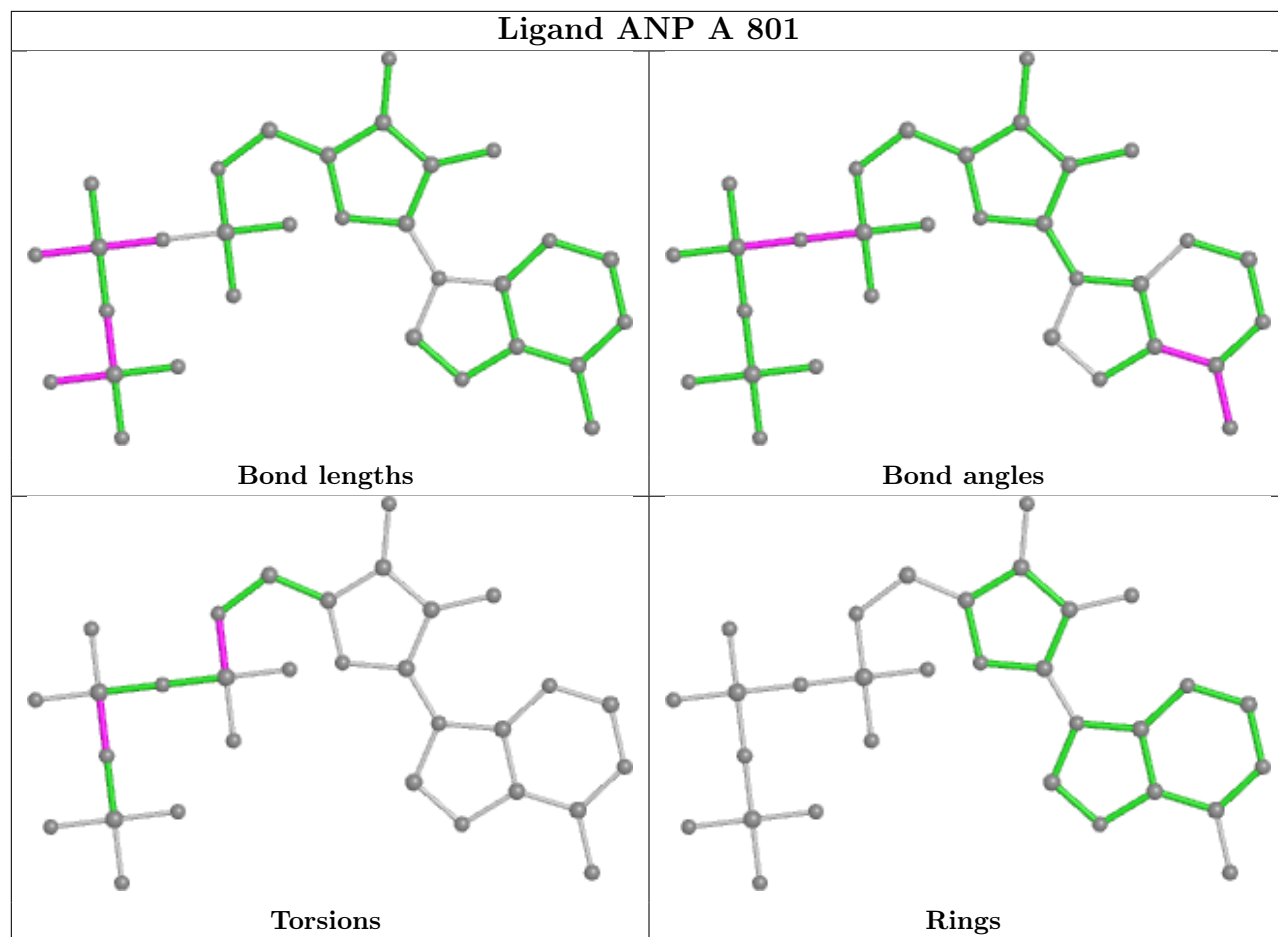
There are no ring outliers.

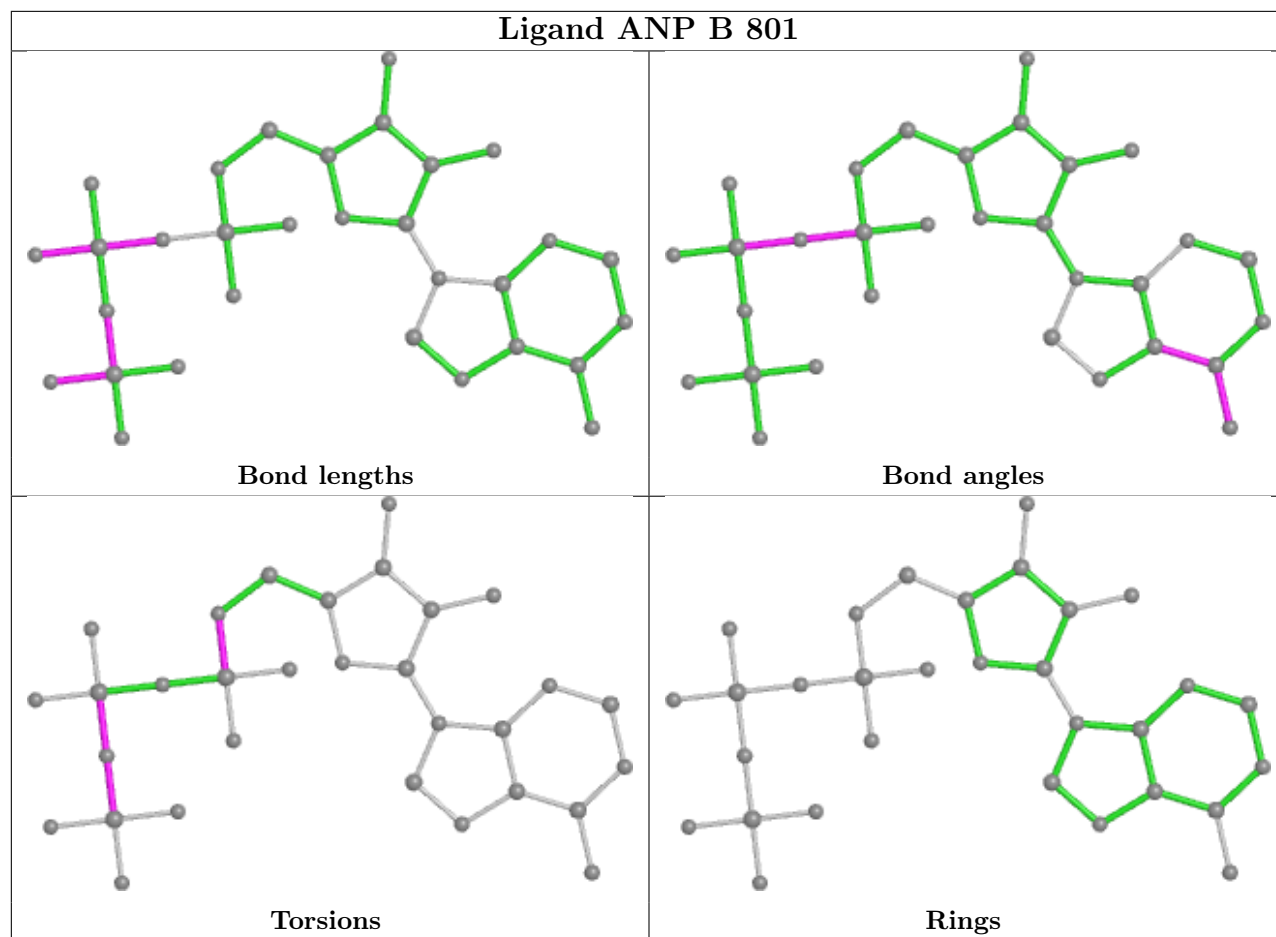
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

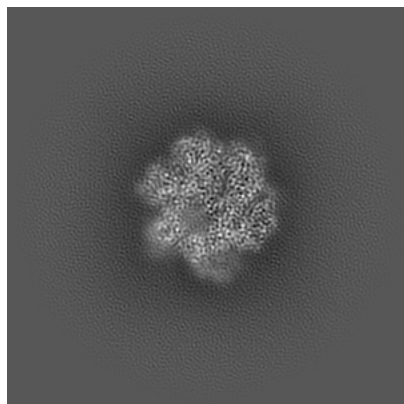
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-38396. These allow visual inspection of the internal detail of the map and identification of artifacts.

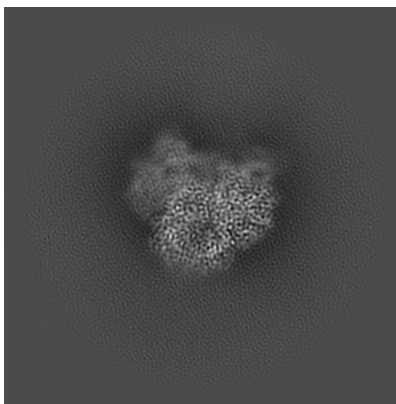
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

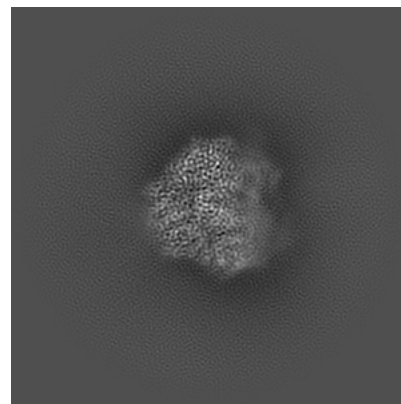
#### 6.1.1 Primary map



X

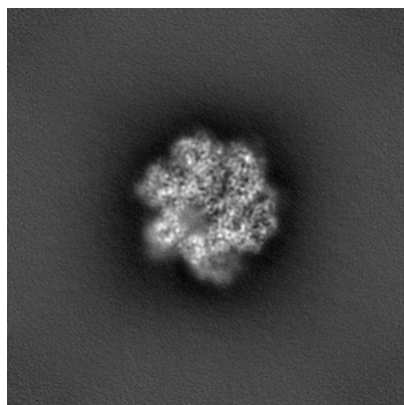


Y

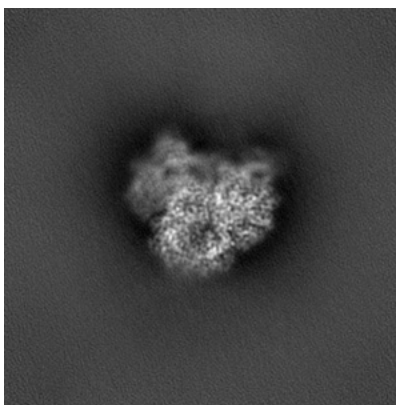


Z

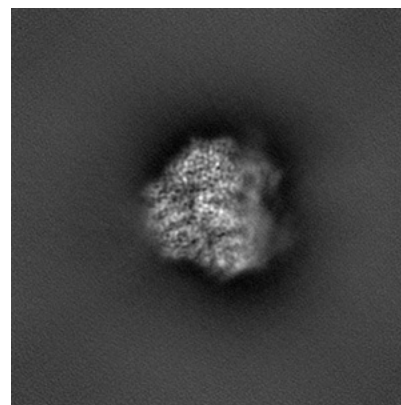
#### 6.1.2 Raw map



X



Y



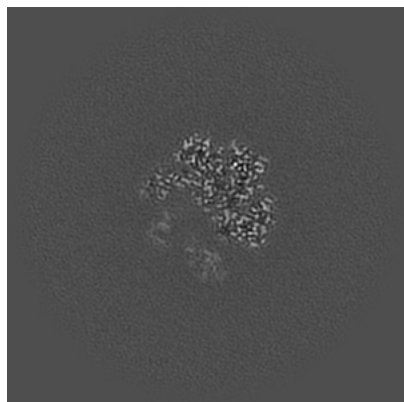
Z

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

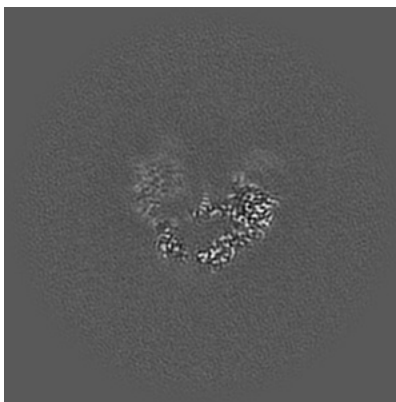


## 6.2 Central slices [i](#)

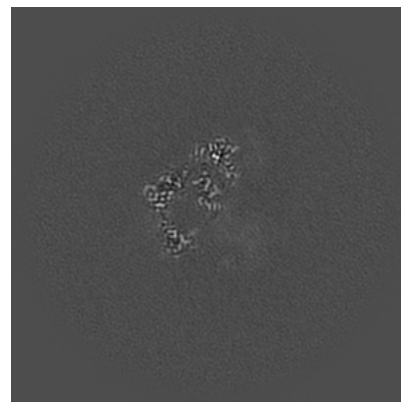
### 6.2.1 Primary map



X Index: 140

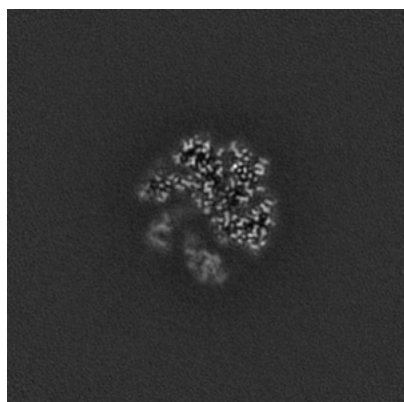


Y Index: 140



Z Index: 140

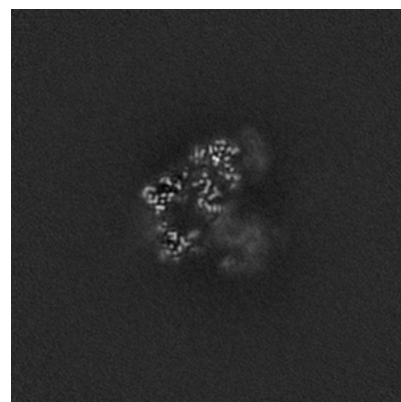
### 6.2.2 Raw map



X Index: 140



Y Index: 140

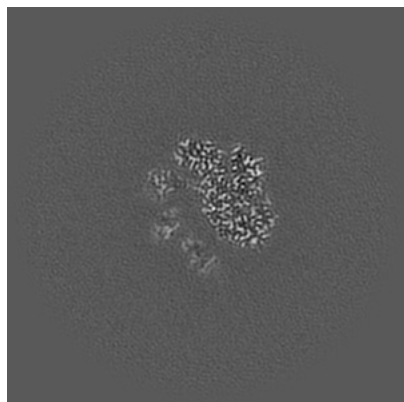


Z Index: 140

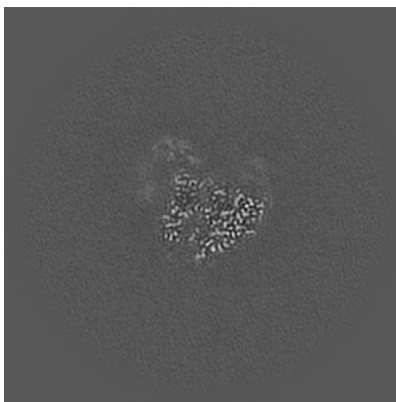
The images above show central slices of the map in three orthogonal directions.

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

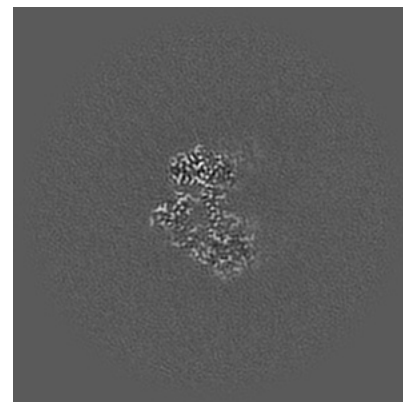
### 6.3.1 Primary map



X Index: 136

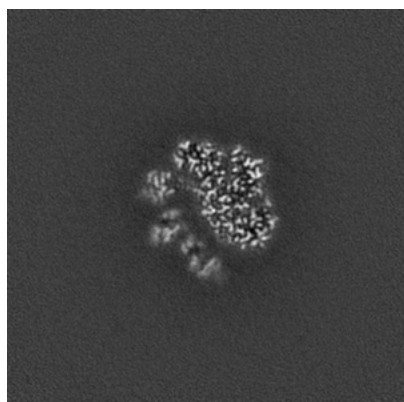


Y Index: 160

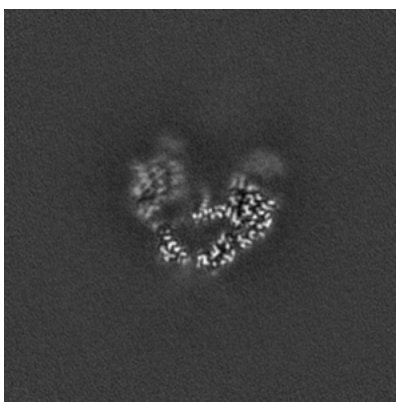


Z Index: 155

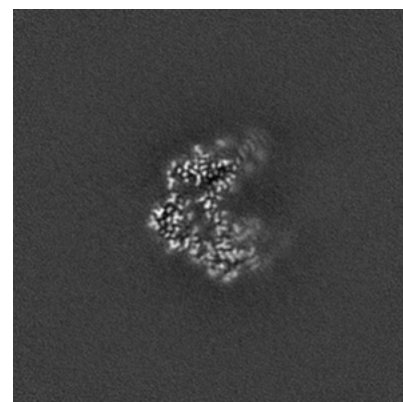
### 6.3.2 Raw map



X Index: 136



Y Index: 141

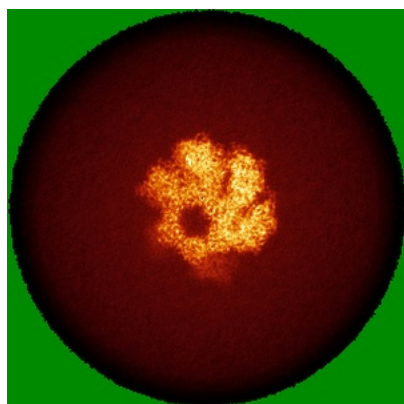


Z Index: 151

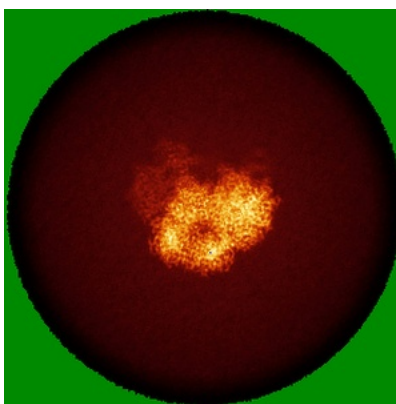
The images above show the largest variance slices of the map in three orthogonal directions.

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

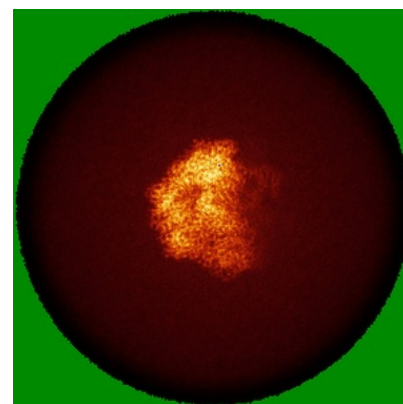
### 6.4.1 Primary map



X

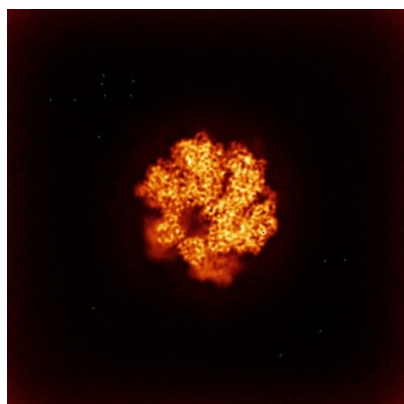


Y

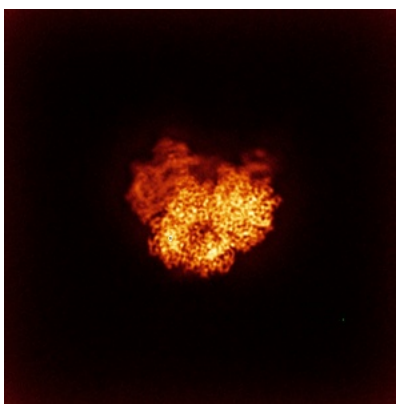


Z

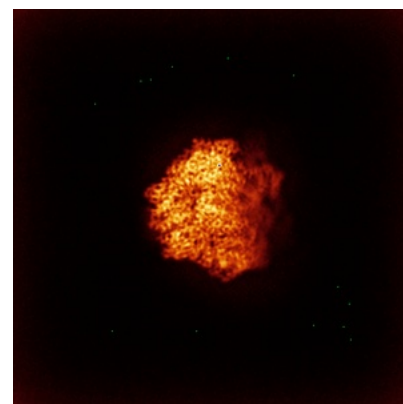
### 6.4.2 Raw map



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



X



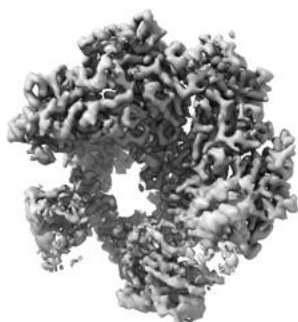
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.733. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

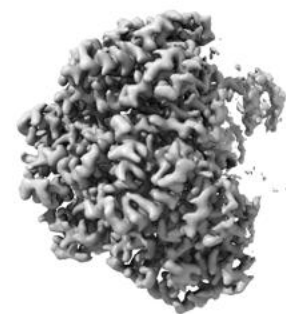
### 6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

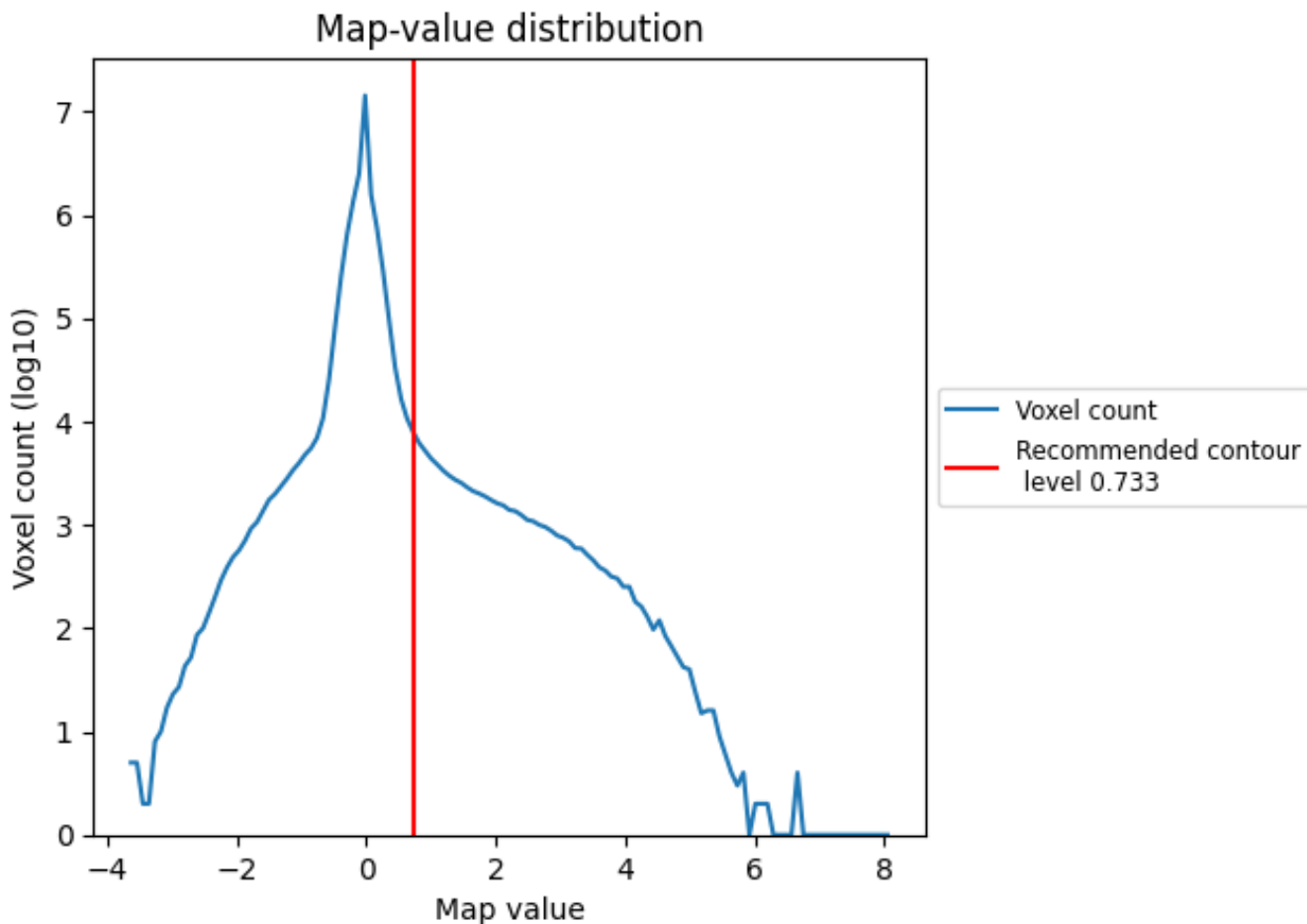
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

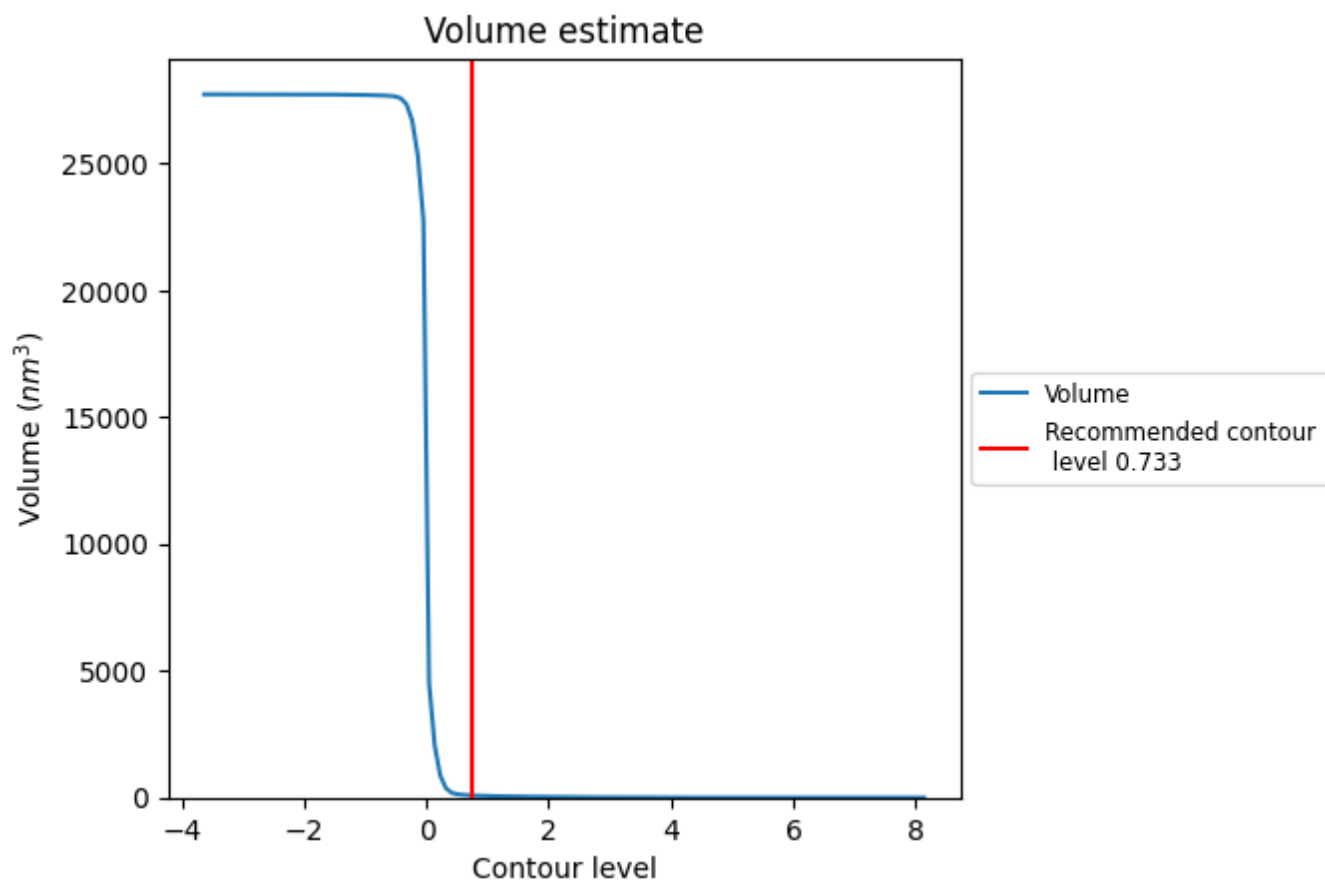
This section contains the results of statistical analysis of the map.

### 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.

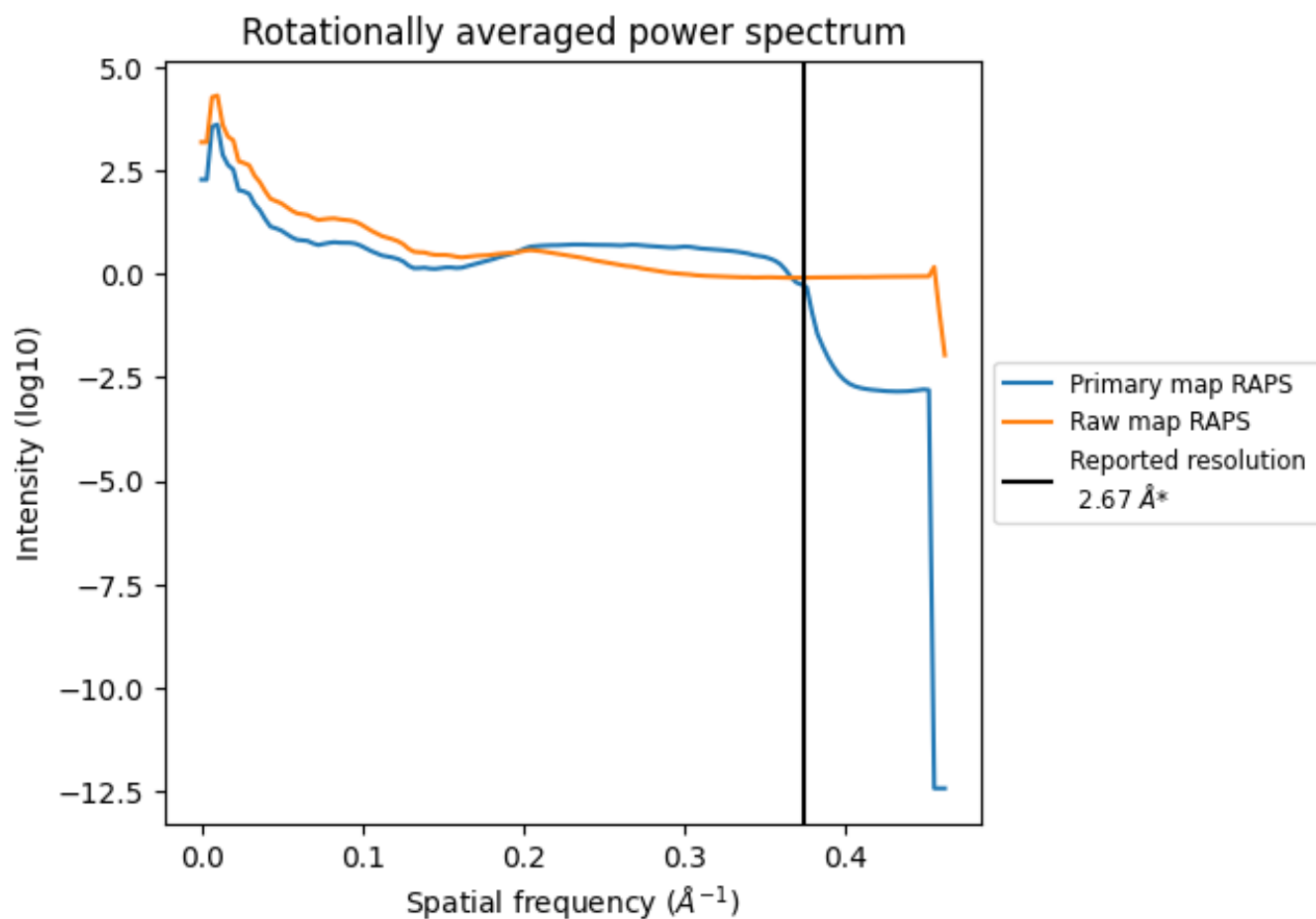
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 86 nm<sup>3</sup>; this corresponds to an approximate mass of 78 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](#)

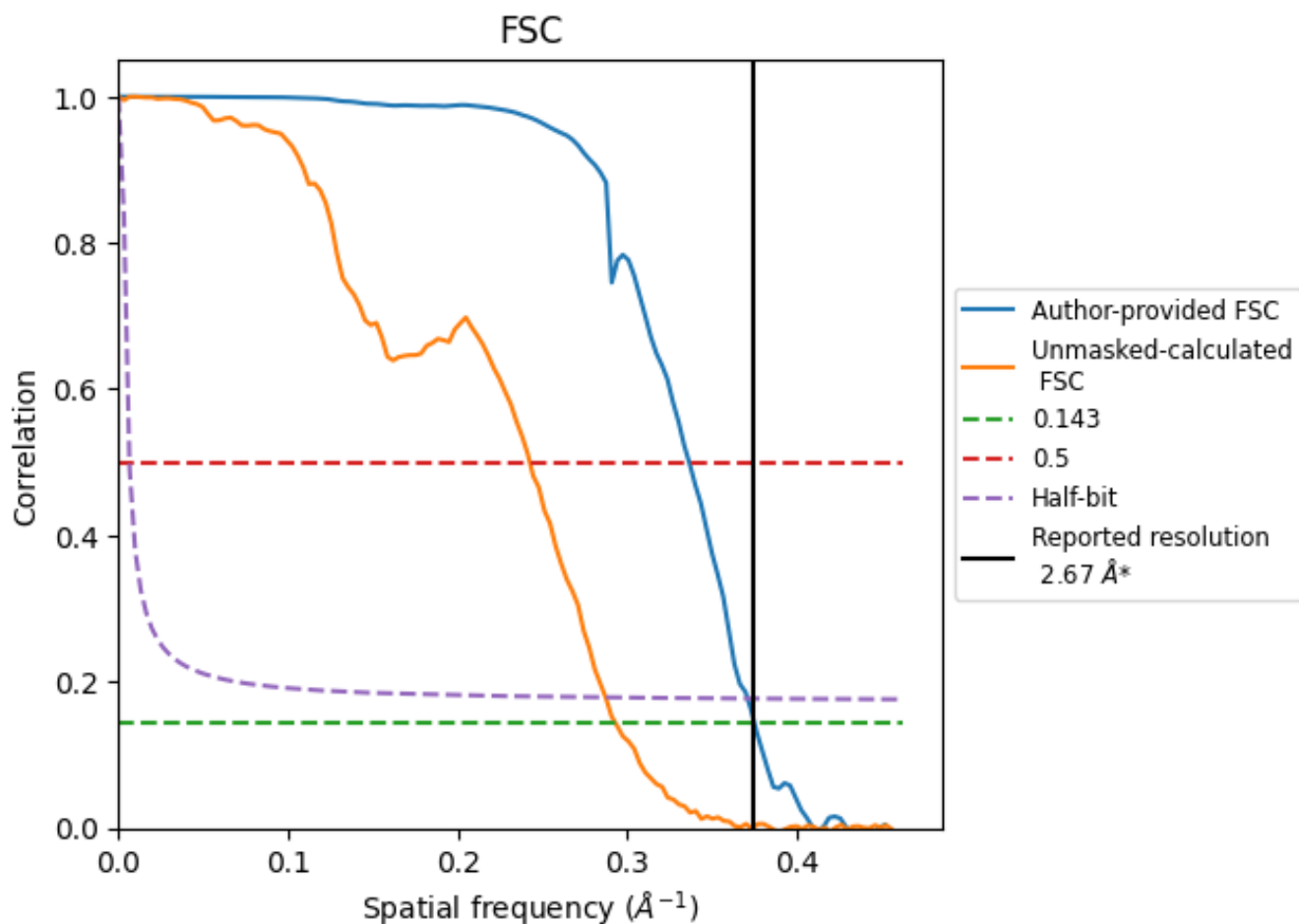


\*Reported resolution corresponds to spatial frequency of 0.375 Å<sup>-1</sup>

## 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 0.375 Å<sup>-1</sup>



## 8.2 Resolution estimates

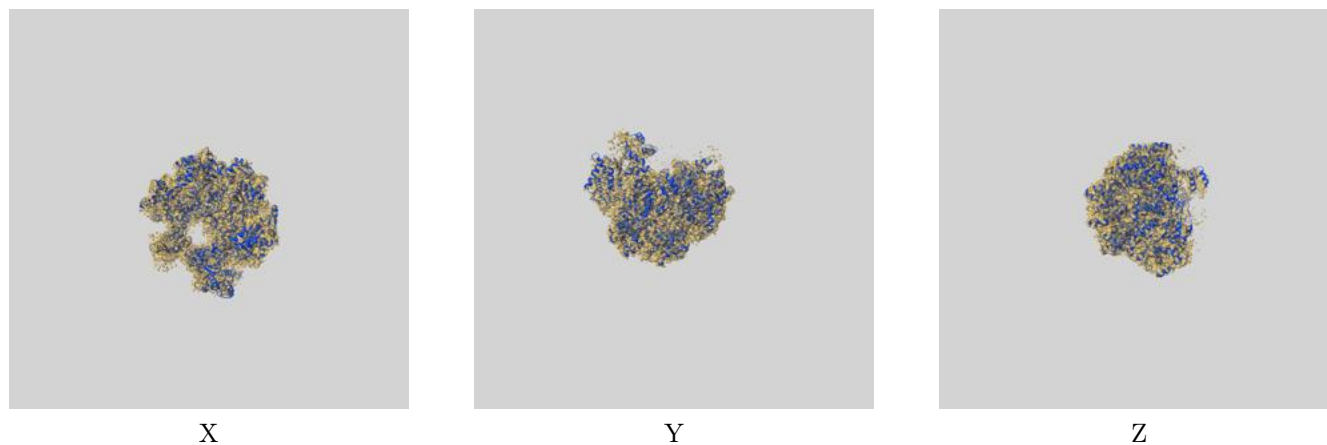
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.67	-	-
Author-provided FSC curve	2.67	2.97	2.70
Unmasked-calculated*	3.41	4.13	3.48

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.41 differs from the reported value 2.67 by more than 10 %

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

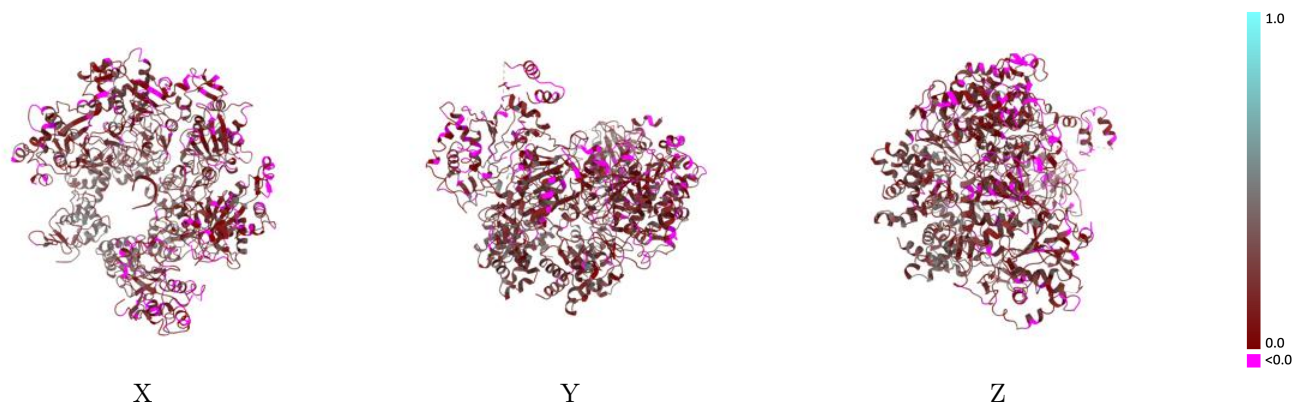
This section contains information regarding the fit between EMDB map EMD-38396 and PDB model 8XJ8. Per-residue inclusion information can be found in section 3 on page 5.

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



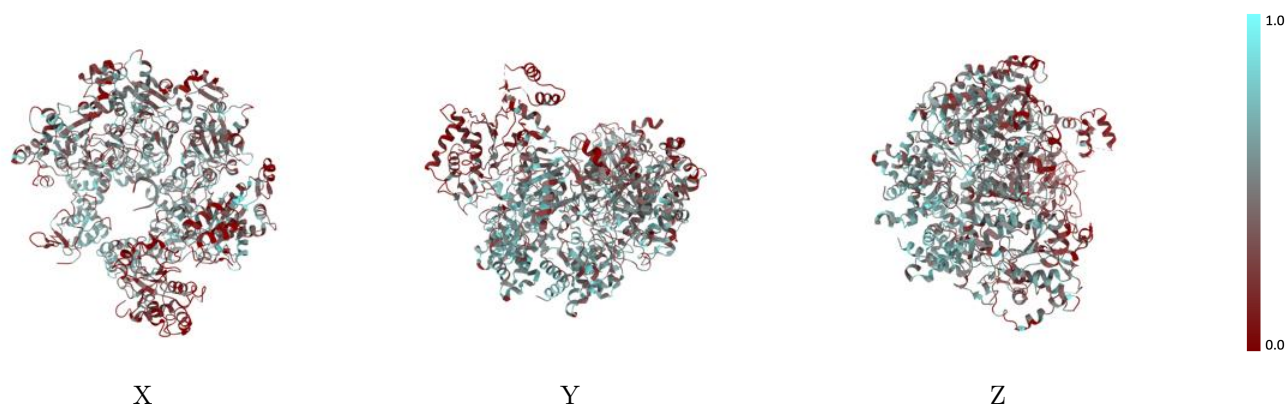
The images above show the 3D surface view of the map at the recommended contour level 0.733 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



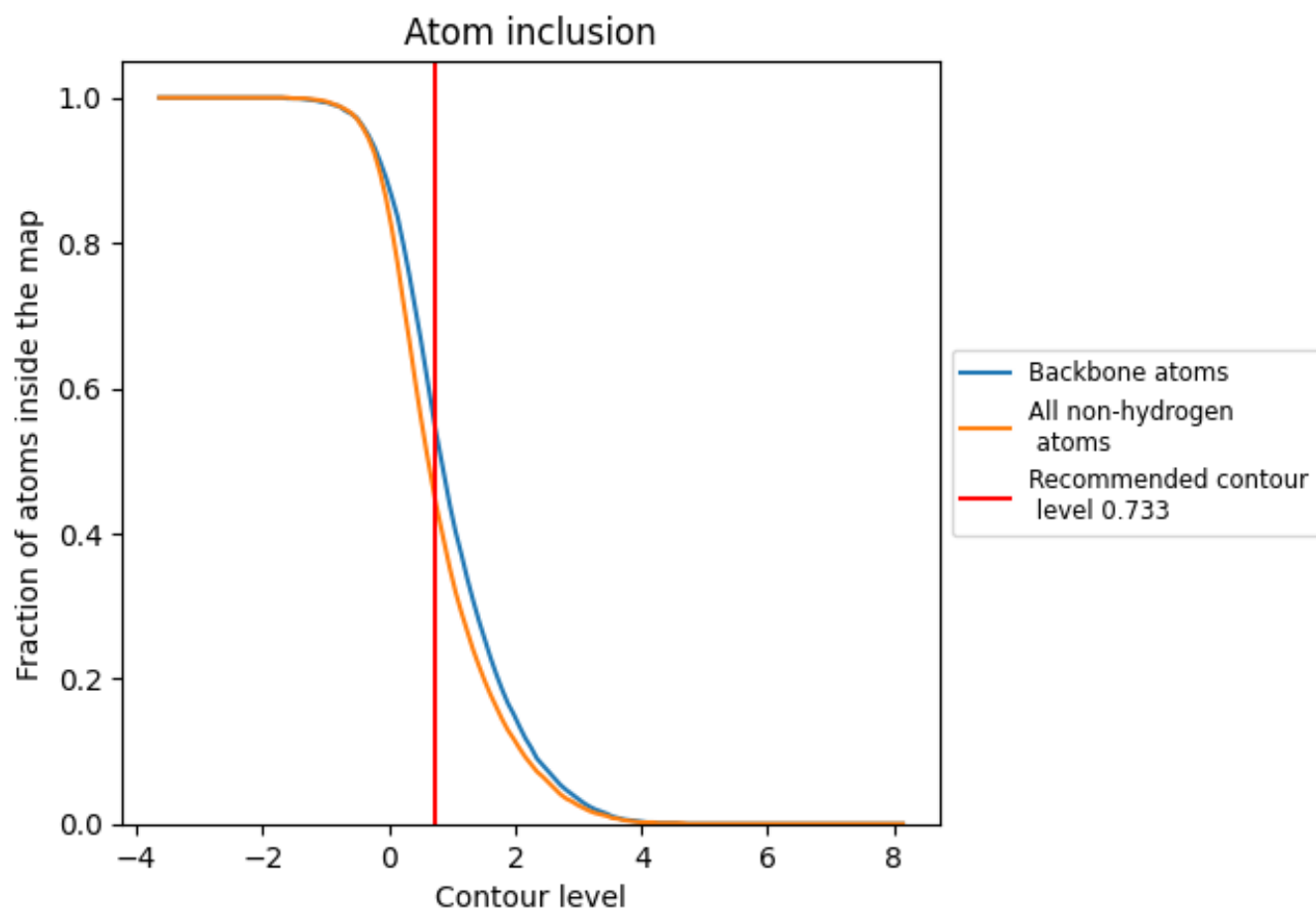
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.733).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 54% of all backbone atoms, 44% of all non-hydrogen atoms, are inside the map.

## 9.5 Map-model fit summary [i](#)

The table lists the average atom inclusion at the recommended contour level (0.733) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.4450	0.1890
A	0.4650	0.1800
B	0.4750	0.1620
C	0.4830	0.1770
D	0.4550	0.2030
E	0.4680	0.3140
F	0.3110	0.1890
X	0.3920	0.1190

