



Full wwPDB EM Validation Report ⓘ

Oct 12, 2024 – 08:37 pm BST

PDB ID : 7ZR0
EMDB ID : EMD-14875
Title : CryoEM structure of HSP90-CDC37-BRAF(V600E) complex.
Authors : Oberoi, J.; Pearl, L.H.
Deposited on : 2022-05-03
Resolution : 3.40 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
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
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

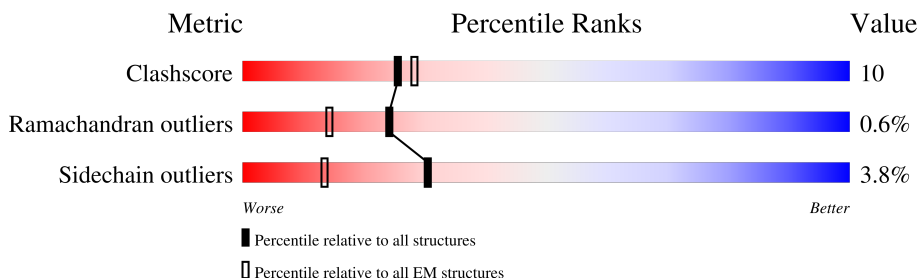
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.40 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	749	
1	B	749	
2	C	396	
3	K	828	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14052 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 Heat shock protein HSP 90-beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	633	Total	C	N	O	S	0	0
			5132	3251	862	996	23		
1	B	631	Total	C	N	O	S	0	0
			5112	3240	857	992	23		

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	initiating methionine	UNP P08238
A	-23	ARG	-	expression tag	UNP P08238
A	-22	GLY	-	expression tag	UNP P08238
A	-21	SER	-	expression tag	UNP P08238
A	-20	HIS	-	expression tag	UNP P08238
A	-19	HIS	-	expression tag	UNP P08238
A	-18	HIS	-	expression tag	UNP P08238
A	-17	HIS	-	expression tag	UNP P08238
A	-16	HIS	-	expression tag	UNP P08238
A	-15	HIS	-	expression tag	UNP P08238
A	-14	HIS	-	expression tag	UNP P08238
A	-13	HIS	-	expression tag	UNP P08238
A	-12	GLY	-	expression tag	UNP P08238
A	-11	MET	-	expression tag	UNP P08238
A	-10	ALA	-	expression tag	UNP P08238
A	-9	LEU	-	expression tag	UNP P08238
A	-8	GLU	-	expression tag	UNP P08238
A	-7	VAL	-	expression tag	UNP P08238
A	-6	LEU	-	expression tag	UNP P08238
A	-5	PHE	-	expression tag	UNP P08238
A	-4	GLN	-	expression tag	UNP P08238
A	-3	GLY	-	expression tag	UNP P08238
A	-2	PRO	-	expression tag	UNP P08238
A	-1	SER	-	expression tag	UNP P08238
A	0	ALA	-	expression tag	UNP P08238
B	-24	MET	-	initiating methionine	UNP P08238

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-23	ARG	-	expression tag	UNP P08238
B	-22	GLY	-	expression tag	UNP P08238
B	-21	SER	-	expression tag	UNP P08238
B	-20	HIS	-	expression tag	UNP P08238
B	-19	HIS	-	expression tag	UNP P08238
B	-18	HIS	-	expression tag	UNP P08238
B	-17	HIS	-	expression tag	UNP P08238
B	-16	HIS	-	expression tag	UNP P08238
B	-15	HIS	-	expression tag	UNP P08238
B	-14	HIS	-	expression tag	UNP P08238
B	-13	HIS	-	expression tag	UNP P08238
B	-12	GLY	-	expression tag	UNP P08238
B	-11	MET	-	expression tag	UNP P08238
B	-10	ALA	-	expression tag	UNP P08238
B	-9	LEU	-	expression tag	UNP P08238
B	-8	GLU	-	expression tag	UNP P08238
B	-7	VAL	-	expression tag	UNP P08238
B	-6	LEU	-	expression tag	UNP P08238
B	-5	PHE	-	expression tag	UNP P08238
B	-4	GLN	-	expression tag	UNP P08238
B	-3	GLY	-	expression tag	UNP P08238
B	-2	PRO	-	expression tag	UNP P08238
B	-1	SER	-	expression tag	UNP P08238
B	0	ALA	-	expression tag	UNP P08238

- Molecule 2 is a protein called Hsp90 co-chaperone Cdc37.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	C	266	Total	C	N	O	P	S	0	0
			2243	1393	399	434	1	16		

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	379	LEU	-	expression tag	UNP Q16543
C	380	GLU	-	expression tag	UNP Q16543
C	381	VAL	-	expression tag	UNP Q16543
C	382	LEU	-	expression tag	UNP Q16543
C	383	PHE	-	expression tag	UNP Q16543
C	384	GLN	-	expression tag	UNP Q16543
C	385	GLY	-	expression tag	UNP Q16543
C	386	PRO	-	expression tag	UNP Q16543

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Chain	Residue	Modelled	Actual	Comment	Reference
C	387	LEU	-	expression tag	UNP Q16543
C	388	GLU	-	expression tag	UNP Q16543
C	389	HIS	-	expression tag	UNP Q16543
C	390	HIS	-	expression tag	UNP Q16543
C	391	HIS	-	expression tag	UNP Q16543
C	392	HIS	-	expression tag	UNP Q16543
C	393	HIS	-	expression tag	UNP Q16543
C	394	HIS	-	expression tag	UNP Q16543
C	395	HIS	-	expression tag	UNP Q16543
C	396	HIS	-	expression tag	UNP Q16543

- Molecule 3 is a protein called Serine/threonine-protein kinase B-raf.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	K	186	Total	C	N	O	S	0	0
			1503	966	260	266	11		

There are 63 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	-61	MET	-	initiating methionine	UNP P15056
K	-60	GLY	-	expression tag	UNP P15056
K	-59	GLY	-	expression tag	UNP P15056
K	-58	SER	-	expression tag	UNP P15056
K	-57	HIS	-	expression tag	UNP P15056
K	-56	HIS	-	expression tag	UNP P15056
K	-55	HIS	-	expression tag	UNP P15056
K	-54	HIS	-	expression tag	UNP P15056
K	-53	HIS	-	expression tag	UNP P15056
K	-52	HIS	-	expression tag	UNP P15056
K	-51	HIS	-	expression tag	UNP P15056
K	-50	HIS	-	expression tag	UNP P15056
K	-49	GLY	-	expression tag	UNP P15056
K	-48	GLY	-	expression tag	UNP P15056
K	-47	SER	-	expression tag	UNP P15056
K	-46	TRP	-	expression tag	UNP P15056
K	-45	SER	-	expression tag	UNP P15056
K	-44	HIS	-	expression tag	UNP P15056
K	-43	PRO	-	expression tag	UNP P15056
K	-42	GLN	-	expression tag	UNP P15056
K	-41	PHE	-	expression tag	UNP P15056
K	-40	GLU	-	expression tag	UNP P15056

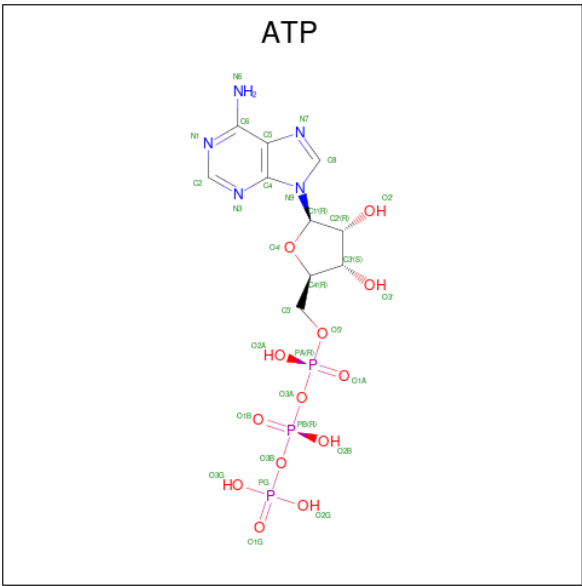
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Chain	Residue	Modelled	Actual	Comment	Reference
K	-39	LYS	-	expression tag	UNP P15056
K	-38	GLY	-	expression tag	UNP P15056
K	-37	GLY	-	expression tag	UNP P15056
K	-36	GLY	-	expression tag	UNP P15056
K	-35	SER	-	expression tag	UNP P15056
K	-34	GLY	-	expression tag	UNP P15056
K	-33	GLY	-	expression tag	UNP P15056
K	-32	GLY	-	expression tag	UNP P15056
K	-31	SER	-	expression tag	UNP P15056
K	-30	GLY	-	expression tag	UNP P15056
K	-29	GLY	-	expression tag	UNP P15056
K	-28	GLY	-	expression tag	UNP P15056
K	-27	SER	-	expression tag	UNP P15056
K	-26	TRP	-	expression tag	UNP P15056
K	-25	SER	-	expression tag	UNP P15056
K	-24	HIS	-	expression tag	UNP P15056
K	-23	PRO	-	expression tag	UNP P15056
K	-22	GLN	-	expression tag	UNP P15056
K	-21	PHE	-	expression tag	UNP P15056
K	-20	GLU	-	expression tag	UNP P15056
K	-19	LYS	-	expression tag	UNP P15056
K	-18	GLY	-	expression tag	UNP P15056
K	-17	ALA	-	expression tag	UNP P15056
K	-16	GLU	-	expression tag	UNP P15056
K	-15	THR	-	expression tag	UNP P15056
K	-14	ALA	-	expression tag	UNP P15056
K	-13	VAL	-	expression tag	UNP P15056
K	-12	PRO	-	expression tag	UNP P15056
K	-11	ASN	-	expression tag	UNP P15056
K	-10	SER	-	expression tag	UNP P15056
K	-9	LEU	-	expression tag	UNP P15056
K	-8	GLU	-	expression tag	UNP P15056
K	-7	VAL	-	expression tag	UNP P15056
K	-6	LEU	-	expression tag	UNP P15056
K	-5	PHE	-	expression tag	UNP P15056
K	-4	GLN	-	expression tag	UNP P15056
K	-3	GLY	-	expression tag	UNP P15056
K	-2	PRO	-	expression tag	UNP P15056
K	-1	SER	-	expression tag	UNP P15056
K	0	ALA	-	expression tag	UNP P15056
K	600	GLU	VAL	variant	UNP P15056

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula:

C₁₀H₁₆N₅O₁₃P₃).

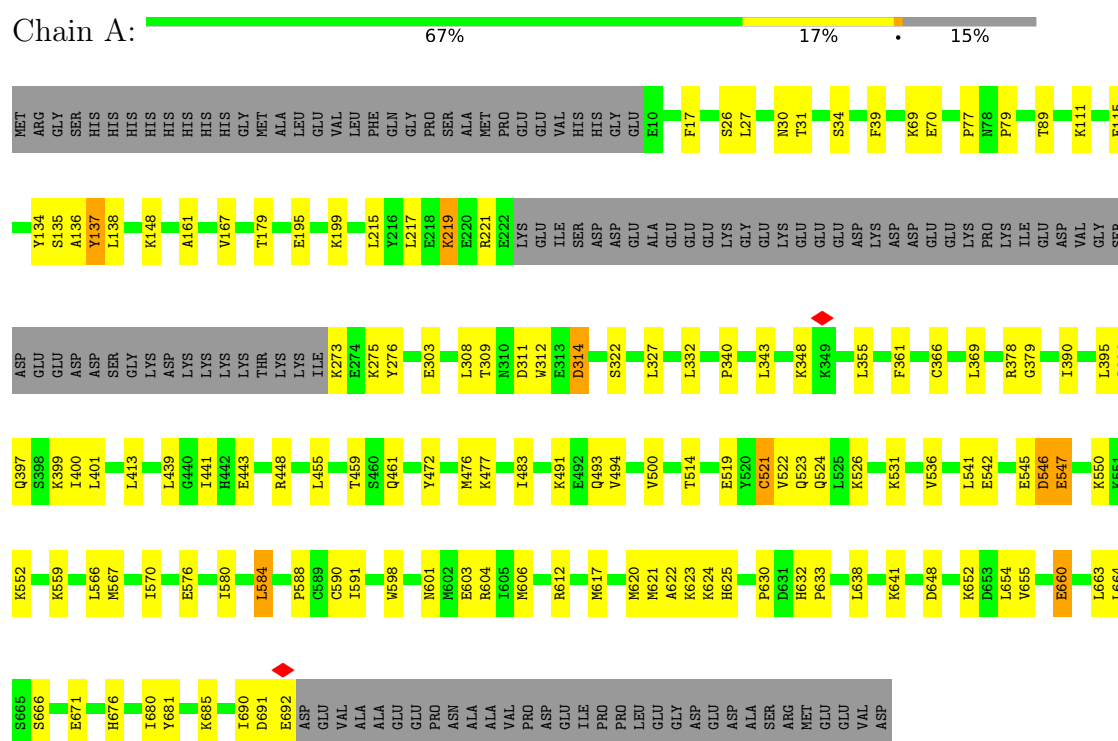


Mol	Chain	Residues	Atoms					AltConf
4	A	1	Total	C	N	O	P	0
			31	10	5	13	3	
4	B	1	Total	C	N	O	P	0
			31	10	5	13	3	

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

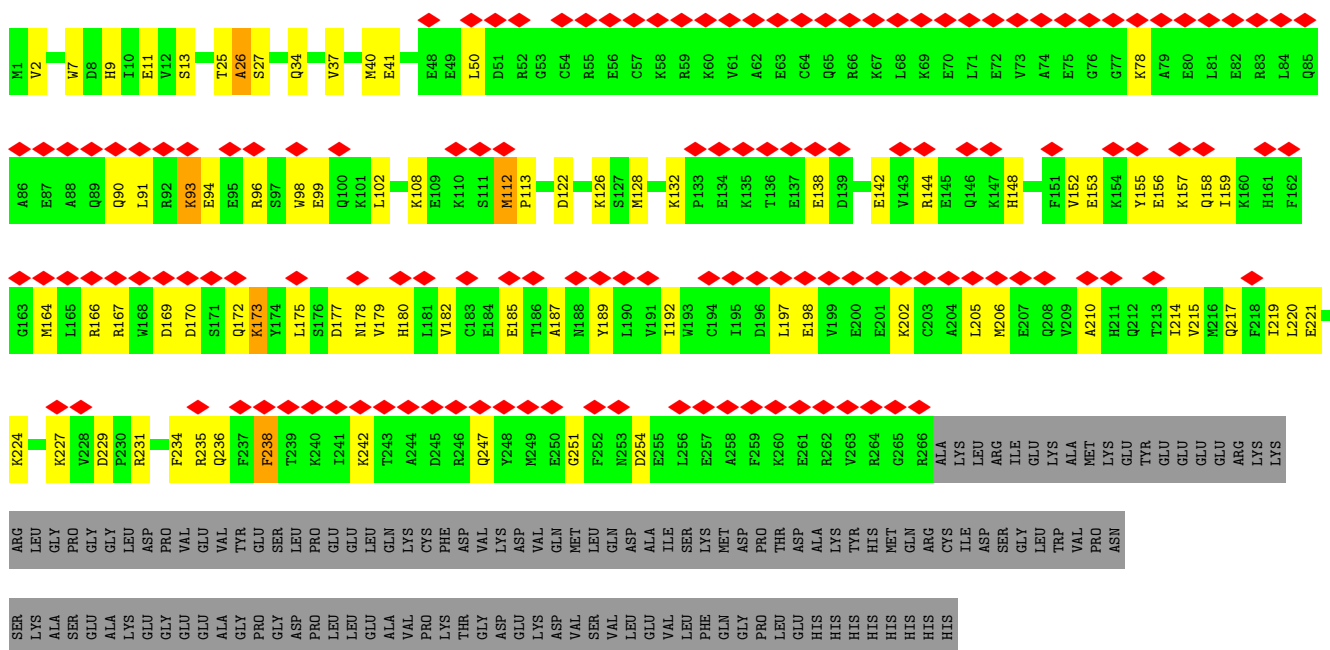
• Molecule 1: Heat shock protein HSP 90-beta



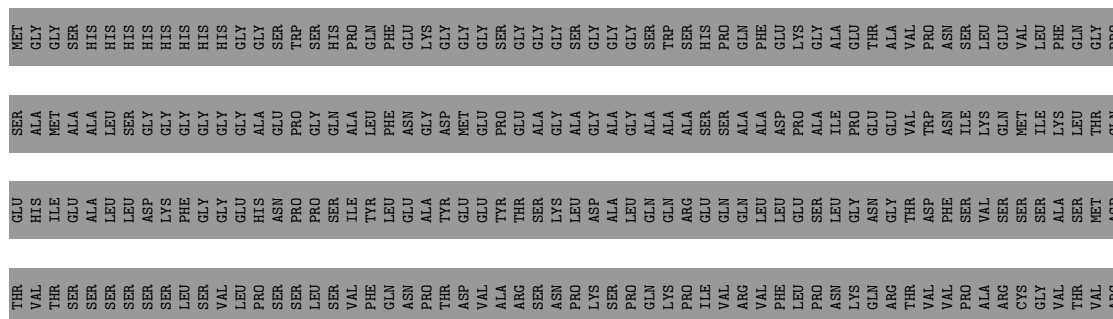
• Molecule 1: Heat shock protein HSP 90-beta



- Molecule 2: Hsp90 co-chaperone Cdc37



- Molecule 3: Serine/threonine-protein kinase B-raf





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	400624	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 ($e^-/\text{\AA}^2$)	45	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	1.844	Depositor
Minimum map value	-0.002	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.026	Depositor
Recommended contour level	0.02	Depositor
Map size (\AA)	282.08002, 282.08002, 282.08002	wwPDB
Map dimensions	328, 328, 328	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.8600001, 0.8600001, 0.8600001	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: SEP, ATP

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	0.38	3/5216 (0.1%)	0.64	4/7010 (0.1%)
1	B	0.32	0/5196	0.56	3/6984 (0.0%)
2	C	0.26	0/2270	0.52	0/3034
3	K	0.27	0/1534	0.54	0/2066
All	All	0.33	3/14216 (0.0%)	0.59	7/19094 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
2	C	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	79	PRO	CB-CG	-12.74	0.86	1.50
1	A	79	PRO	CG-CD	-7.43	1.26	1.50
1	A	77	PRO	N-CD	-5.36	1.40	1.47

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	79	PRO	CB-CG-CD	18.83	179.95	106.50
1	A	79	PRO	N-CD-CG	-16.20	78.91	103.20
1	A	79	PRO	CA-CB-CG	-14.17	77.07	104.00
1	B	59	LEU	CA-CB-CG	6.77	130.88	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	688	LEU	CA-CB-CG	5.84	128.73	115.30
1	A	79	PRO	CA-N-CD	-5.17	104.27	111.50
1	B	566	LEU	CA-CB-CG	5.02	126.84	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	309	THR	Peptide
2	C	112	MET	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5132	0	5162	103	0
1	B	5112	0	5143	102	0
2	C	2243	0	2192	63	0
3	K	1503	0	1543	33	0
4	A	31	0	12	2	0
4	B	31	0	12	0	0
All	All	14052	0	14064	284	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (284) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:MET:HG3	1:A:621:MET:CE	1.53	1.36
1:A:620:MET:CG	1:A:621:MET:HE1	1.61	1.29
2:C:148:HIS:O	2:C:152:VAL:HB	1.53	1.07
1:A:620:MET:HG3	1:A:621:MET:HE1	1.07	1.05
1:A:620:MET:CG	1:A:621:MET:CE	2.25	0.99
1:A:308:LEU:HD23	1:A:309:THR:HG23	1.47	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:MET:HG3	1:A:621:MET:HE2	1.51	0.93
1:A:620:MET:HG2	1:A:621:MET:HE1	1.52	0.90
1:A:308:LEU:HD23	1:A:309:THR:CG2	2.06	0.86
1:B:195:GLU:HG2	1:B:217:LEU:HD11	1.61	0.82
1:B:660:GLU:OE1	1:B:676:HIS:HB2	1.82	0.79
3:K:650:MET:SD	3:K:689:MET:SD	2.81	0.78
1:A:580:ILE:HD11	1:A:630:PRO:HG3	1.64	0.77
2:C:11:GLU:HB2	2:C:113:PRO:HG2	1.67	0.77
1:B:94:THR:OG1	1:B:152:ASP:OD1	2.01	0.75
1:A:542:GLU:N	1:A:542:GLU:OE1	2.20	0.75
3:K:650:MET:CE	3:K:689:MET:SD	2.75	0.75
1:A:179:THR:HG21	4:A:801:ATP:HN62	1.52	0.74
1:A:459:THR:HG23	1:A:461:GLN:H	1.53	0.74
1:B:195:GLU:CD	1:B:217:LEU:HD21	2.09	0.73
2:C:179:VAL:HB	2:C:231:ARG:HA	1.70	0.72
1:A:576:GLU:N	1:A:576:GLU:OE1	2.21	0.72
1:B:70:GLU:N	1:B:70:GLU:OE1	2.23	0.72
1:B:361:PHE:HZ	1:B:364:ASP:HB3	1.55	0.71
1:A:523:GLN:HE22	3:K:523:PRO:HB3	1.56	0.70
2:C:25:THR:O	2:C:27:SER:N	2.24	0.70
2:C:164:MET:SD	2:C:164:MET:N	2.65	0.70
2:C:178:ASN:O	2:C:182:VAL:N	2.23	0.70
1:B:632:HIS:CD2	1:B:633:PRO:HD2	2.26	0.70
1:A:638:LEU:HD12	1:A:654:LEU:HD23	1.74	0.70
1:A:397:GLN:HE22	1:A:399:LYS:HB3	1.56	0.69
2:C:108:LYS:HD3	2:C:112:MET:HG3	1.74	0.69
1:A:524:GLN:O	1:A:526:LYS:NZ	2.26	0.69
1:A:691:ASP:C	1:A:692:GLU:OE1	2.31	0.69
1:A:70:GLU:OE1	1:A:70:GLU:N	2.26	0.68
1:A:493:GLN:C	1:A:493:GLN:OE1	2.31	0.68
1:A:690:ILE:O	1:A:692:GLU:OE1	2.12	0.67
1:A:620:MET:CG	1:A:621:MET:HE2	2.15	0.67
2:C:187:ALA:HB1	2:C:220:LEU:HD11	1.76	0.67
2:C:108:LYS:O	2:C:112:MET:N	2.18	0.67
1:A:27:LEU:O	1:A:31:THR:OG1	2.13	0.67
1:B:219:LYS:HA	1:B:219:LYS:HE3	1.77	0.66
1:A:441:ILE:O	1:A:448:ARG:NH1	2.29	0.66
1:A:134:TYR:HA	1:A:137:TYR:HB2	1.77	0.65
1:B:87:VAL:HG22	1:B:180:LYS:HG2	1.78	0.65
1:B:560:PHE:HB2	1:B:580:ILE:HD11	1.79	0.65
1:B:353:ILE:HD12	1:B:370:ILE:HD13	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:570:ILE:HD13	1:A:655:VAL:HG11	1.80	0.64
2:C:50:LEU:HD11	2:C:98:TRP:HB3	1.78	0.64
1:B:663:LEU:HD22	1:B:670:LEU:HD13	1.78	0.64
1:A:632:HIS:HD2	1:A:633:PRO:HD2	1.61	0.64
1:A:632:HIS:CD2	1:A:633:PRO:HD2	2.33	0.63
2:C:155:TYR:HA	2:C:158:GLN:HB2	1.80	0.63
1:B:341:PHE:HZ	2:C:236:GLN:HB3	1.65	0.61
1:B:97:ASP:CG	1:B:101:ASN:HD21	2.04	0.61
1:B:160:SER:O	1:B:160:SER:OG	2.18	0.60
1:A:395:LEU:HD13	1:A:401:LEU:HD21	1.84	0.60
2:C:25:THR:OG1	2:C:26:ALA:N	2.34	0.60
1:B:67:SER:O	1:B:177:ARG:NH1	2.34	0.60
1:B:139:VAL:HG23	1:B:139:VAL:O	2.00	0.60
1:B:616:THR:HG21	3:K:586:GLU:HG3	1.84	0.60
1:A:526:LYS:HE3	2:C:247:GLN:HG2	1.84	0.59
1:A:612:ARG:HH21	1:B:516:PRO:HD3	1.67	0.59
1:A:308:LEU:CD2	1:A:309:THR:HG23	2.26	0.59
1:B:36:LYS:HG2	1:B:138:LEU:HD22	1.85	0.59
1:A:598:TRP:CD2	1:A:621:MET:HG3	2.38	0.58
1:A:523:GLN:NE2	3:K:523:PRO:HB3	2.18	0.58
1:A:493:GLN:OE1	1:A:494:VAL:N	2.36	0.58
1:B:51:LEU:HD22	1:B:71:LEU:HB3	1.85	0.58
3:K:564:MET:HE1	3:K:577:LEU:HD13	1.86	0.58
1:B:632:HIS:HD2	1:B:633:PRO:HD2	1.67	0.58
1:A:500:VAL:HG23	1:A:500:VAL:O	2.04	0.58
1:B:314:ASP:OD1	1:B:314:ASP:N	2.36	0.57
1:A:493:GLN:C	1:A:493:GLN:CD	2.63	0.57
1:A:570:ILE:HD11	1:A:652:LYS:HG3	1.85	0.56
2:C:215:VAL:O	2:C:219:ILE:HG12	2.04	0.56
1:A:179:THR:HG21	4:A:801:ATP:N6	2.20	0.56
1:A:355:LEU:HD11	1:A:369:LEU:HD11	1.87	0.56
1:B:581:SER:HB3	1:B:627:GLU:HB3	1.87	0.56
1:A:396:GLN:O	1:A:397:GLN:HG3	2.04	0.56
2:C:144:ARG:HD3	2:C:185:GLU:HG3	1.86	0.56
1:B:353:ILE:HB	1:B:366:CYS:SG	2.45	0.56
1:B:381:VAL:HG12	1:B:381:VAL:O	2.06	0.56
1:A:167:VAL:HG12	1:A:167:VAL:O	2.05	0.56
1:B:341:PHE:CZ	2:C:236:GLN:HB3	2.41	0.55
1:B:443:GLU:O	1:B:443:GLU:HG2	2.07	0.55
2:C:108:LYS:NZ	2:C:112:MET:SD	2.78	0.55
1:B:597:GLY:HA3	1:B:622:ALA:HB2	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:ASN:ND2	2:C:2:VAL:O	2.37	0.55
1:A:195:GLU:OE2	1:A:199:LYS:NZ	2.40	0.55
1:B:190:THR:O	1:B:190:THR:HG23	2.07	0.55
2:C:197:LEU:HD22	2:C:202:LYS:HD2	1.88	0.55
1:B:539:GLU:O	1:B:539:GLU:HG3	2.07	0.55
1:B:297:THR:O	1:B:299:GLU:N	2.40	0.55
1:A:309:THR:HG21	1:A:378:ARG:HH11	1.70	0.54
1:B:501:GLU:HG3	1:B:585:VAL:HG12	1.87	0.54
2:C:189:TYR:HA	2:C:192:ILE:HG22	1.89	0.54
1:B:589:CYS:SG	1:B:662:ALA:HB2	2.47	0.54
1:B:27:LEU:O	1:B:31:THR:HB	2.07	0.54
1:A:217:LEU:HD11	1:A:219:LYS:NZ	2.22	0.54
1:B:185:LEU:HD12	1:B:193:LEU:HD21	1.88	0.54
1:A:311:ASP:OD1	1:A:312:TRP:N	2.40	0.54
2:C:185:GLU:N	2:C:185:GLU:OE1	2.40	0.53
1:B:59:LEU:O	1:B:59:LEU:HD12	2.09	0.53
1:B:195:GLU:HG2	1:B:217:LEU:CD1	2.35	0.53
1:A:111:LYS:O	1:A:115:GLU:HG2	2.08	0.53
1:B:564:CYS:SG	1:B:578:VAL:HG12	2.49	0.53
1:B:688:LEU:HD12	1:B:688:LEU:O	2.09	0.53
3:K:575:ARG:HD2	3:K:633:TYR:CD2	2.44	0.52
1:A:620:MET:CB	1:A:621:MET:HE2	2.39	0.52
3:K:703:GLU:N	3:K:703:GLU:OE1	2.41	0.52
1:A:221:ARG:H	1:A:273:LYS:N	2.07	0.52
1:A:519:GLU:OE2	1:A:601:ASN:ND2	2.38	0.52
2:C:90:GLN:O	2:C:93:LYS:HG3	2.10	0.52
1:B:691:ASP:OD1	1:B:691:ASP:N	2.40	0.52
1:A:588:PRO:HD3	1:A:632:HIS:CE1	2.46	0.51
1:B:97:ASP:OD1	1:B:101:ASN:ND2	2.28	0.51
1:B:475:ARG:HG3	1:B:475:ARG:O	2.10	0.51
1:B:633:PRO:O	1:B:637:THR:HG23	2.10	0.51
1:B:576:GLU:OE1	1:B:625:HIS:NE2	2.44	0.51
3:K:692:LEU:HD11	3:K:710:ILE:HG23	1.92	0.51
1:A:39:PHE:HB2	1:A:136:ALA:HA	1.92	0.51
3:K:568:HIS:HD1	3:K:635:PHE:HD1	1.59	0.51
1:A:545:GLU:HG2	1:A:550:LYS:HE3	1.93	0.51
3:K:552:LYS:HE2	3:K:588:LEU:HD13	1.93	0.51
1:B:545:GLU:OE1	1:B:545:GLU:N	2.43	0.51
2:C:217:GLN:O	2:C:221:GLU:HG2	2.11	0.51
1:B:311:ASP:OD1	1:B:313:GLU:N	2.37	0.50
1:B:447:ASN:OD1	1:B:447:ASN:N	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:175:LEU:HA	2:C:178:ASN:HB2	1.93	0.50
2:C:198:GLU:HB2	2:C:206:MET:CE	2.40	0.50
1:A:390:ILE:HD13	1:B:27:LEU:HD11	1.93	0.50
3:K:629:ASP:O	3:K:630:LYS:HG2	2.12	0.50
1:A:521:CYS:SG	1:A:522:VAL:N	2.85	0.50
3:K:564:MET:HG3	3:K:568:HIS:HD2	1.76	0.50
2:C:9:HIS:O	2:C:113:PRO:HB3	2.12	0.50
2:C:198:GLU:HB2	2:C:206:MET:HE1	1.93	0.50
1:B:195:GLU:CG	1:B:217:LEU:HD21	2.41	0.49
2:C:167:ARG:HB3	2:C:170:ASP:HB3	1.93	0.49
2:C:202:LYS:HB2	2:C:205:LEU:HB2	1.94	0.49
1:A:215:LEU:HD21	1:A:217:LEU:HD23	1.93	0.49
2:C:251:GLY:HA2	2:C:254:ASP:OD2	2.11	0.49
1:A:135:SER:O	1:A:138:LEU:HB2	2.13	0.49
1:B:195:GLU:OE2	1:B:217:LEU:HD21	2.12	0.49
1:A:327:LEU:HD21	1:A:400:ILE:HD12	1.94	0.49
1:A:439:LEU:O	1:A:443:GLU:HG3	2.12	0.49
1:A:576:GLU:HG2	1:A:623:LYS:HD2	1.95	0.49
2:C:122:ASP:OD1	2:C:122:ASP:N	2.45	0.49
1:B:648:ASP:OD1	1:B:649:LYS:N	2.46	0.48
1:A:17:PHE:CE1	1:B:165:PHE:HD1	2.31	0.48
1:A:598:TRP:CE3	1:A:621:MET:HG3	2.48	0.48
1:B:301:TYR:HE2	2:C:128:MET:HE1	1.77	0.48
3:K:621:ALA:HA	3:K:640:TYR:CE1	2.48	0.48
2:C:152:VAL:O	2:C:156:GLU:HB3	2.13	0.48
2:C:153:GLU:O	2:C:157:LYS:HG2	2.14	0.48
2:C:50:LEU:HD13	2:C:102:LEU:HG	1.96	0.48
2:C:152:VAL:O	2:C:156:GLU:CB	2.62	0.48
1:B:542:GLU:OE1	1:B:543:LEU:O	2.32	0.48
1:A:340:PRO:HG2	1:A:343:LEU:HB3	1.95	0.47
1:B:51:LEU:HD23	1:B:90:GLY:HA2	1.95	0.47
1:B:219:LYS:O	1:B:275:LYS:HG2	2.14	0.47
1:B:391:SER:O	1:B:393:GLU:N	2.47	0.47
1:B:80:GLN:OE1	1:B:81:GLU:OE1	2.31	0.47
1:A:491:LYS:HA	1:A:494:VAL:HG12	1.96	0.47
1:A:567:MET:HA	1:A:570:ILE:HG22	1.97	0.47
2:C:91:LEU:O	2:C:94:GLU:HG3	2.14	0.47
3:K:651:THR:HG22	3:K:653:GLN:HG2	1.95	0.47
1:B:368:GLU:HG3	1:B:401:LEU:HD23	1.97	0.47
2:C:210:ALA:O	2:C:214:ILE:HG23	2.15	0.47
1:A:217:LEU:HD11	1:A:219:LYS:HZ3	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:441:ILE:HD11	1:A:455:LEU:HD11	1.97	0.47
1:A:348:LYS:HA	1:A:348:LYS:HD3	1.66	0.46
1:B:452:SER:O	1:B:452:SER:OG	2.32	0.46
3:K:537:LEU:HD12	3:K:584:LEU:HG	1.98	0.46
1:A:476:MET:HE1	1:A:531:LYS:HE2	1.98	0.46
1:B:82:ARG:NH2	1:B:187:GLU:OE2	2.49	0.46
1:B:346:ASN:O	1:B:348:LYS:N	2.48	0.46
1:B:301:TYR:CE2	2:C:128:MET:HE1	2.51	0.46
2:C:78:LYS:HD2	2:C:78:LYS:HA	1.74	0.46
1:A:584:LEU:HD21	1:A:590:CYS:HB3	1.98	0.46
3:K:662:ARG:O	3:K:666:ILE:HG22	2.16	0.46
1:B:188:ASP:OD1	1:B:189:GLN:HG2	2.16	0.46
1:B:352:ASN:OD1	1:B:352:ASN:N	2.42	0.46
1:B:367:ASP:OD2	1:B:367:ASP:N	2.39	0.45
2:C:156:GLU:HA	2:C:159:ILE:HG22	1.98	0.45
2:C:34:GLN:O	2:C:37:VAL:HG12	2.16	0.45
2:C:214:ILE:HA	2:C:217:GLN:HG3	1.99	0.45
1:A:691:ASP:O	1:A:692:GLU:OE1	2.34	0.45
3:K:625:ILE:HG21	3:K:666:ILE:HG13	1.97	0.45
1:B:566:LEU:O	1:B:570:ILE:HD12	2.17	0.45
1:A:591:ILE:HD12	1:A:663:LEU:HD13	1.99	0.45
1:A:601:ASN:OD1	1:A:604:ARG:NH2	2.50	0.45
1:A:526:LYS:HB2	1:A:526:LYS:HE2	1.57	0.44
1:A:566:LEU:O	1:A:570:ILE:HG22	2.16	0.44
1:A:680:ILE:HD11	1:B:680:ILE:HD11	1.99	0.44
1:B:361:PHE:CZ	1:B:364:ASP:HB3	2.42	0.44
2:C:214:ILE:HG13	2:C:215:VAL:N	2.31	0.44
2:C:227:LYS:HD2	2:C:227:LYS:HA	1.70	0.44
2:C:138:GLU:OE2	2:C:142:GLU:HB3	2.17	0.44
2:C:220:LEU:O	2:C:224:LYS:HG2	2.18	0.44
1:A:591:ILE:HG23	1:A:625:HIS:O	2.18	0.44
1:A:89:THR:O	1:A:89:THR:OG1	2.35	0.44
1:A:413:LEU:HD23	1:A:413:LEU:HA	1.81	0.44
3:K:691:ARG:HD2	3:K:691:ARG:HA	1.81	0.44
1:A:69:LYS:HA	1:A:69:LYS:HD3	1.81	0.44
1:B:195:GLU:HG2	1:B:217:LEU:HD21	2.00	0.44
2:C:9:HIS:C	2:C:113:PRO:HB3	2.38	0.44
1:A:547:GLU:HA	1:A:550:LYS:HB2	1.99	0.44
1:A:552:LYS:HB2	1:A:552:LYS:HE3	1.69	0.44
1:A:660:GLU:OE1	1:A:676:HIS:HB2	2.18	0.44
1:B:481:LYS:O	1:B:482:SER:OG	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:621:ALA:HB3	3:K:624:VAL:HG23	2.00	0.44
1:A:397:GLN:NE2	1:A:399:LYS:HB3	2.29	0.43
1:B:25:MET:O	1:B:29:ILE:HG23	2.17	0.43
1:B:539:GLU:OE2	1:B:595:THR:OG1	2.29	0.43
1:A:681:TYR:CZ	1:B:588:PRO:HG2	2.52	0.43
3:K:640:TYR:HE2	3:K:697:LEU:HD13	1.83	0.43
3:K:650:MET:HE1	3:K:689:MET:SD	2.57	0.43
1:A:332:LEU:O	1:A:379:GLY:HA3	2.19	0.43
1:B:53:LYS:HE2	1:B:125:MET:HE1	2.00	0.43
1:B:118:GLN:HB3	1:B:119:ALA:H	1.43	0.43
1:B:341:PHE:CD1	1:B:341:PHE:C	2.92	0.43
2:C:175:LEU:HA	2:C:175:LEU:HD12	1.80	0.43
2:C:221:GLU:HA	2:C:224:LYS:HG2	2.01	0.43
1:A:546:ASP:OD1	1:A:547:GLU:N	2.52	0.43
1:B:340:PRO:HG2	1:B:343:LEU:HD13	2.00	0.43
1:B:293:PRO:HB2	2:C:126:LYS:HD2	2.00	0.43
3:K:640:TYR:HB2	3:K:704:ARG:HH11	1.84	0.42
1:B:20:GLU:O	1:B:24:LEU:HB2	2.19	0.42
1:A:308:LEU:HD23	1:A:309:THR:HG22	1.93	0.42
1:B:297:THR:O	1:B:297:THR:OG1	2.28	0.42
1:B:188:ASP:OD1	1:B:188:ASP:C	2.57	0.42
1:A:664:LEU:HD11	1:B:664:LEU:HD11	2.02	0.42
1:A:685:LYS:HG2	1:A:690:ILE:HB	2.01	0.42
1:B:202:VAL:HG11	1:B:282:LEU:HD12	2.01	0.42
1:B:311:ASP:OD1	1:B:312:TRP:N	2.52	0.42
1:B:394:MET:HG2	1:B:395:LEU:H	1.83	0.42
2:C:229:ASP:OD1	2:C:229:ASP:N	2.52	0.42
1:A:459:THR:HG21	1:A:461:GLN:HE21	1.85	0.42
1:A:641:LYS:HB2	1:A:641:LYS:HE3	1.87	0.42
2:C:156:GLU:HG3	2:C:189:TYR:CE2	2.55	0.42
1:A:493:GLN:CD	1:A:493:GLN:O	2.59	0.42
3:K:548:PHE:HD2	3:K:553:LEU:HD21	1.85	0.42
3:K:626:ARG:HD2	3:K:626:ARG:HA	1.77	0.42
3:K:677:ASP:HB3	3:K:680:LYS:HE2	2.02	0.42
2:C:177:ASP:HA	2:C:180:HIS:HB2	2.02	0.41
1:B:383:SER:OG	1:B:384:GLU:N	2.52	0.41
3:K:568:HIS:ND1	3:K:635:PHE:HD1	2.16	0.41
1:A:477:LYS:H	1:A:477:LYS:HG3	1.73	0.41
1:B:281:GLU:HA	1:B:284:LYS:HD2	2.03	0.41
2:C:202:LYS:HD3	2:C:205:LEU:HD22	2.02	0.41
2:C:238:PHE:O	2:C:242:LYS:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:GLU:O	1:B:173:GLU:HG2	2.21	0.41
1:B:580:ILE:HG23	1:B:630:PRO:HG3	2.03	0.41
2:C:96:ARG:NH1	2:C:99:GLU:OE1	2.51	0.41
1:A:666:SER:O	1:A:666:SER:OG	2.33	0.41
1:B:571:LEU:HD23	1:B:571:LEU:HA	1.84	0.41
1:B:97:ASP:O	1:B:101:ASN:ND2	2.54	0.41
2:C:169:ASP:OD1	2:C:170:ASP:N	2.53	0.41
2:C:170:ASP:O	2:C:173:LYS:HG3	2.21	0.41
2:C:172:GLN:HG3	2:C:238:PHE:CD2	2.56	0.41
3:K:530:GLN:O	3:K:531:TRP:HB3	2.20	0.41
3:K:668:MET:O	3:K:673:TYR:HA	2.21	0.41
1:A:514:THR:O	1:A:514:THR:OG1	2.34	0.41
1:A:576:GLU:OE1	1:A:624:LYS:O	2.39	0.41
1:B:322:SER:HB2	1:B:330:ARG:HG2	2.02	0.41
3:K:573:ILE:HG22	3:K:574:HIS:H	1.86	0.41
1:A:472:TYR:HE1	1:A:483:ILE:HG12	1.84	0.41
1:B:394:MET:HG2	1:B:395:LEU:N	2.36	0.41
1:B:397:GLN:HG3	2:C:7:TRP:O	2.21	0.41
2:C:156:GLU:HG3	2:C:189:TYR:HE2	1.86	0.41
3:K:529:THR:OG1	3:K:530:GLN:N	2.54	0.40
3:K:650:MET:HE3	3:K:689:MET:SD	2.58	0.40
1:B:539:GLU:HB3	1:B:593:THR:O	2.21	0.40
1:A:314:ASP:OD1	1:A:314:ASP:N	2.55	0.40
1:A:603:GLU:OE1	1:A:622:ALA:HB2	2.21	0.40
1:B:55:ARG:HG2	1:B:55:ARG:HH11	1.86	0.40
1:B:456:ARG:HA	1:B:456:ARG:HD3	1.83	0.40
1:A:536:VAL:HA	1:A:541:LEU:HD13	2.03	0.40
1:B:487:THR:HG22	1:B:513:MET:HB2	2.03	0.40
1:B:567:MET:HE1	1:B:628:ILE:HD12	2.04	0.40
1:B:623:LYS:H	1:B:623:LYS:HG2	1.55	0.40
1:A:308:LEU:CD2	1:A:309:THR:CG2	2.89	0.40
3:K:578:LYS:HA	3:K:619:TRP:NE1	2.37	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	629/749 (84%)	560 (89%)	66 (10%)	3 (0%)	25	54
1	B	627/749 (84%)	556 (89%)	66 (10%)	5 (1%)	16	44
2	C	263/396 (66%)	247 (94%)	15 (6%)	1 (0%)	30	60
3	K	182/828 (22%)	157 (86%)	24 (13%)	1 (0%)	25	54
All	All	1701/2722 (62%)	1520 (89%)	171 (10%)	10 (1%)	24	50

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	118	GLN
1	B	298	GLN
1	B	574	LYS
2	C	26	ALA
3	K	523	PRO
1	B	275	LYS
1	B	347	LYS
1	A	161	ALA
1	A	366	CYS
1	A	276	TYR

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	572/672 (85%)	552 (96%)	20 (4%)	31	56
1	B	570/672 (85%)	548 (96%)	22 (4%)	27	53
2	C	244/357 (68%)	235 (96%)	9 (4%)	29	54
3	K	167/702 (24%)	159 (95%)	8 (5%)	21	48
All	All	1553/2403 (65%)	1494 (96%)	59 (4%)	30	54

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	SER
1	A	34	SER
1	A	137	TYR
1	A	148	LYS
1	A	219	LYS
1	A	275	LYS
1	A	303	GLU
1	A	314	ASP
1	A	322	SER
1	A	361	PHE
1	A	521	CYS
1	A	546	ASP
1	A	547	GLU
1	A	559	LYS
1	A	584	LEU
1	A	606	MET
1	A	617	MET
1	A	648	ASP
1	A	660	GLU
1	A	671	GLU
1	B	52	ASP
1	B	66	ASP
1	B	74	ASP
1	B	115	GLU
1	B	171	HIS
1	B	188	ASP
1	B	216	TYR
1	B	313	GLU
1	B	337	ARG
1	B	338	ARG
1	B	341	PHE
1	B	417	SER
1	B	444	ASP
1	B	468	SER
1	B	489	GLU
1	B	521	CYS
1	B	539	GLU
1	B	548	GLU
1	B	581	SER
1	B	589	CYS
1	B	620	MET
1	B	641	LYS

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Mol	Chain	Res	Type
2	C	40	MET
2	C	41	GLU
2	C	93	LYS
2	C	132	LYS
2	C	166	ARG
2	C	173	LYS
2	C	234	PHE
2	C	235	ARG
2	C	238	PHE
3	K	548	PHE
3	K	577	LEU
3	K	595	PHE
3	K	637	SER
3	K	657	SER
3	K	682	ARG
3	K	685	CYS
3	K	722	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	397	GLN
1	A	461	GLN
1	A	523	GLN
1	A	632	HIS
1	B	101	ASN
1	B	632	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SEP	C	13	2	8,9,10	1.49	1 (12%)	8,12,14	1.61	2 (25%)

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
2	SEP	C	13	2	-	0/5/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	13	SEP	P-O1P	3.27	1.61	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	13	SEP	P-OG-CB	-2.96	110.14	118.30
2	C	13	SEP	OG-CB-CA	2.87	110.94	108.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	ATP	A	801	-	26,33,33	0.62	0	31,52,52	0.79	2 (6%)
4	ATP	B	801	-	26,33,33	0.62	0	31,52,52	0.75	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
4	ATP	A	801	-	-	5/18/38/38	0/3/3/3
4	ATP	B	801	-	-	11/18/38/38	0/3/3/3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	801	ATP	C5-C6-N6	2.30	123.85	120.35
4	A	801	ATP	C5-C6-N6	2.23	123.73	120.35
4	A	801	ATP	PB-O3B-PG	2.07	139.93	132.83
4	B	801	ATP	PB-O3B-PG	2.06	139.91	132.83

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	801	ATP	PB-O3B-PG-O2G
4	B	801	ATP	C5'-O5'-PA-O1A
4	B	801	ATP	C5'-O5'-PA-O2A
4	B	801	ATP	C5'-O5'-PA-O3A
4	A	801	ATP	PA-O3A-PB-O1B
4	A	801	ATP	PB-O3A-PA-O5'
4	B	801	ATP	O4'-C4'-C5'-O5'
4	B	801	ATP	C3'-C4'-C5'-O5'
4	B	801	ATP	PB-O3A-PA-O1A

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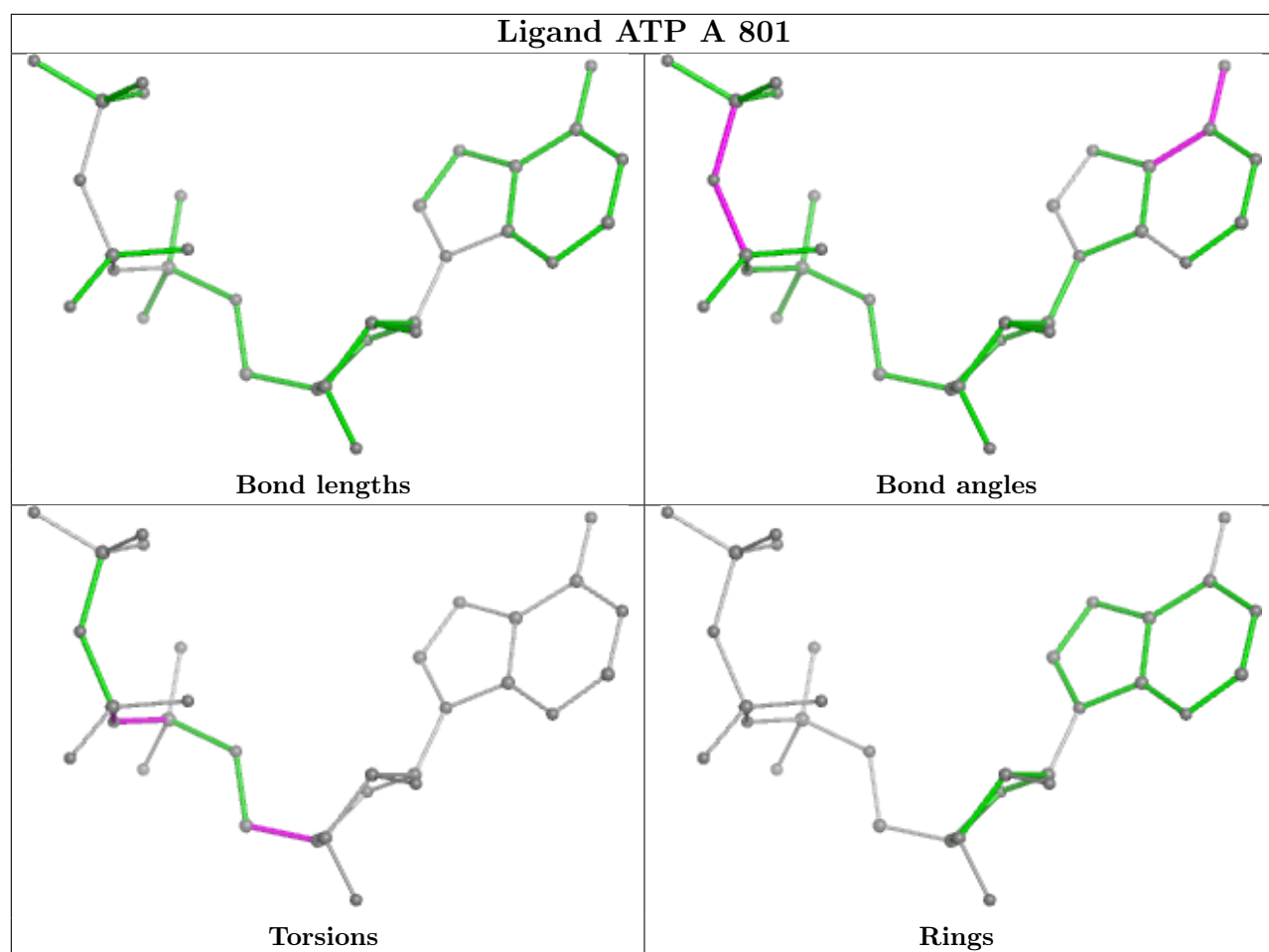
Mol	Chain	Res	Type	Atoms
4	B	801	ATP	PB-O3A-PA-O2A
4	A	801	ATP	O4'-C4'-C5'-O5'
4	A	801	ATP	C3'-C4'-C5'-O5'
4	B	801	ATP	PB-O3B-PG-O1G
4	B	801	ATP	PB-O3B-PG-O3G
4	A	801	ATP	PA-O3A-PB-O2B
4	B	801	ATP	PG-O3B-PB-O2B

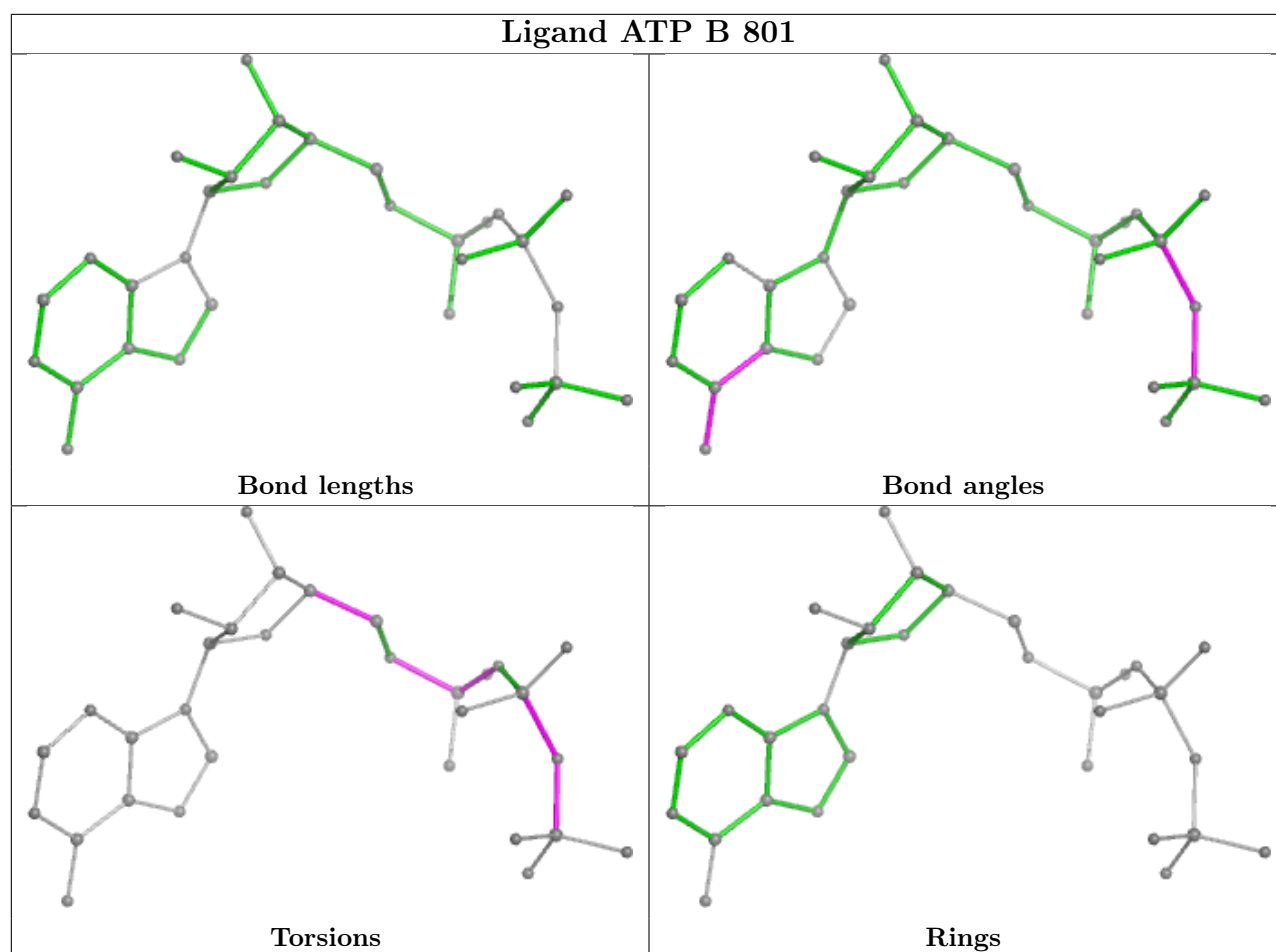
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	801	ATP	2	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

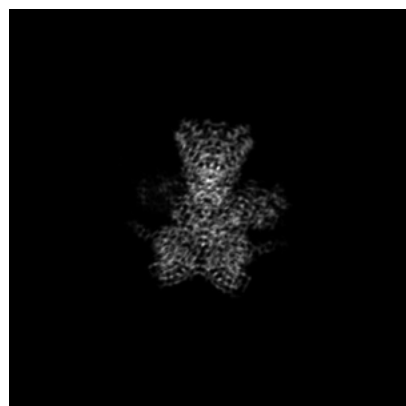
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-14875. These allow visual inspection of the internal detail of the map and identification of artifacts.

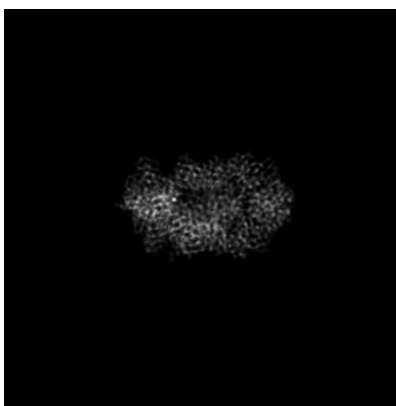
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

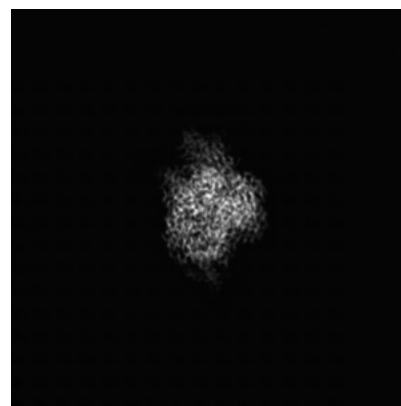
6.1.1 Primary map



X

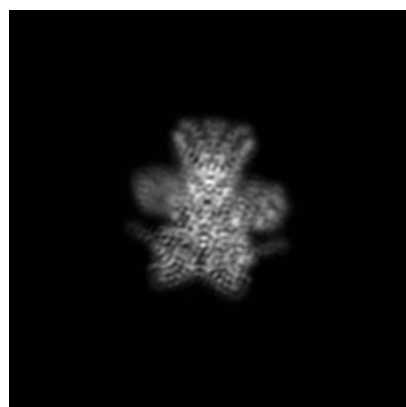


Y

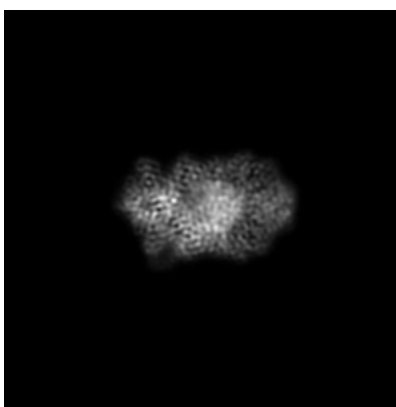


Z

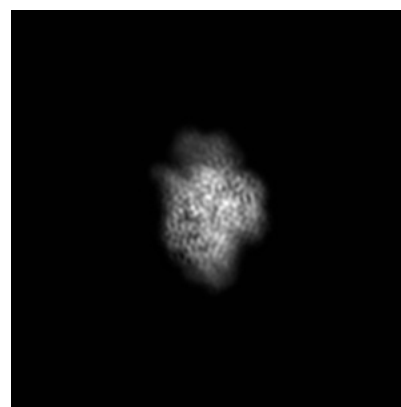
6.1.2 Raw map



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: 164

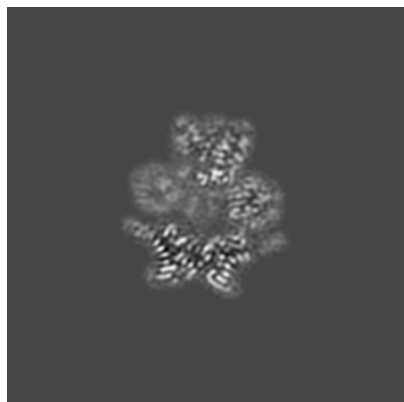


Y Index: 164

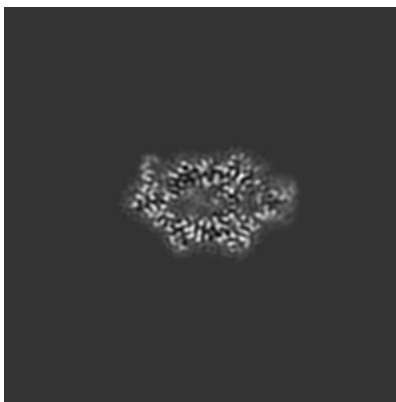


Z Index: 164

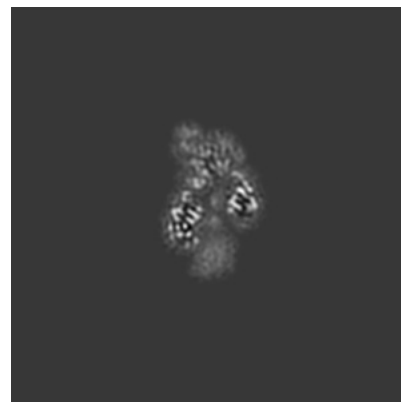
6.2.2 Raw map



X Index: 164



Y Index: 164



Z Index: 164

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: 162

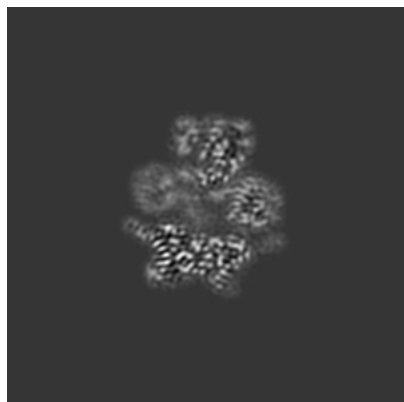


Y Index: 171

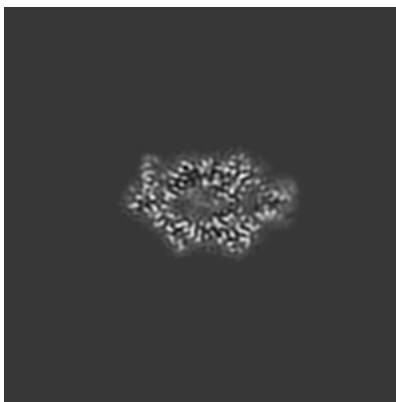


Z Index: 122

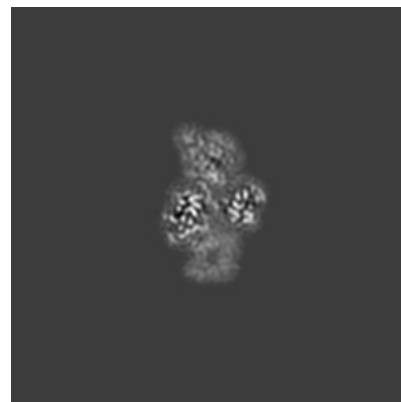
6.3.2 Raw map



X Index: 162



Y Index: 165

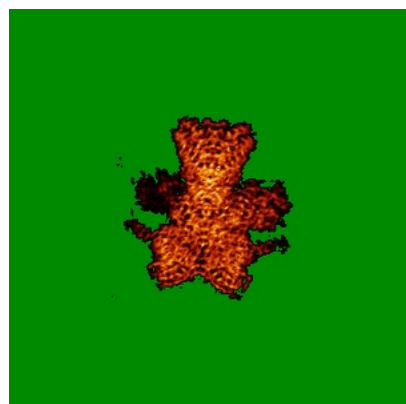


Z Index: 173

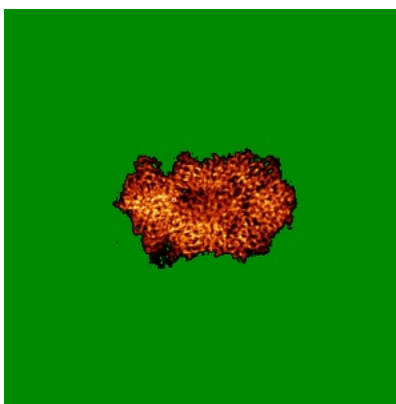
The images above show the largest variance slices of the map in three orthogonal directions.

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

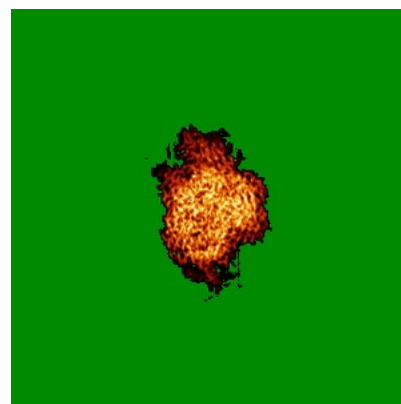
6.4.1 Primary map



X

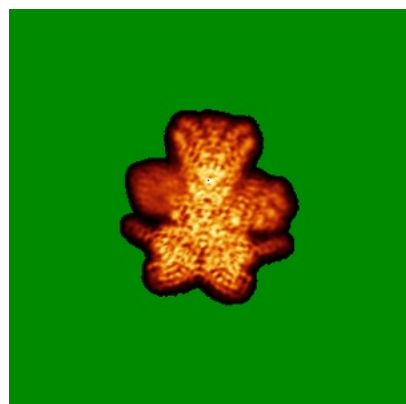


Y

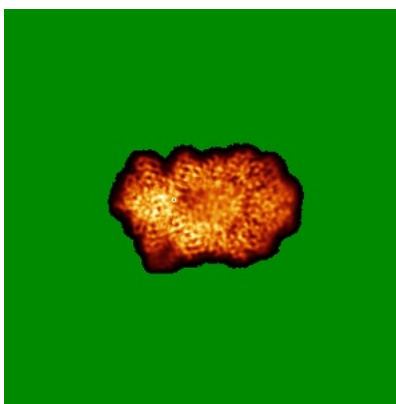


Z

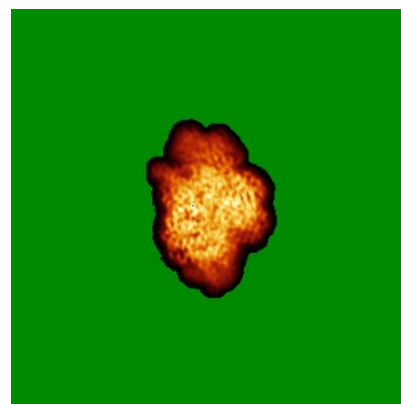
6.4.2 Raw map



X



Y

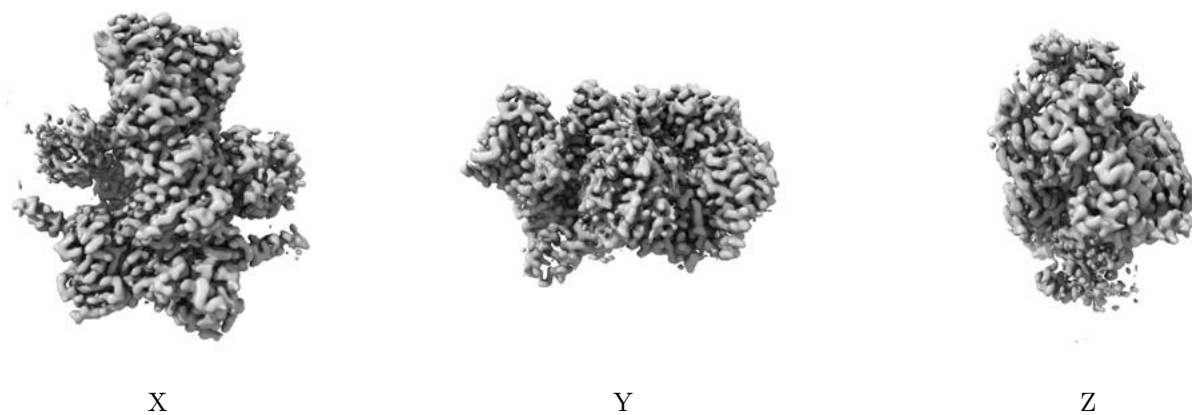


Z

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.

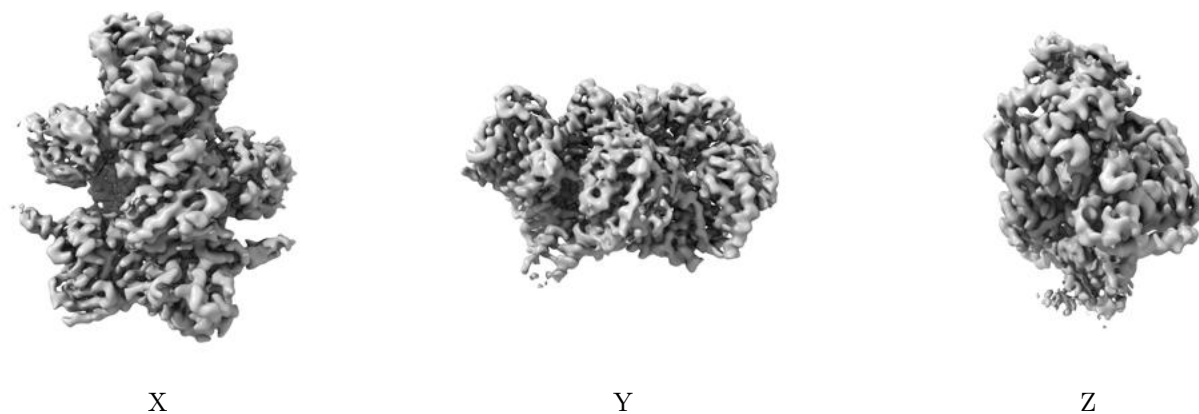
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.02. 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



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.

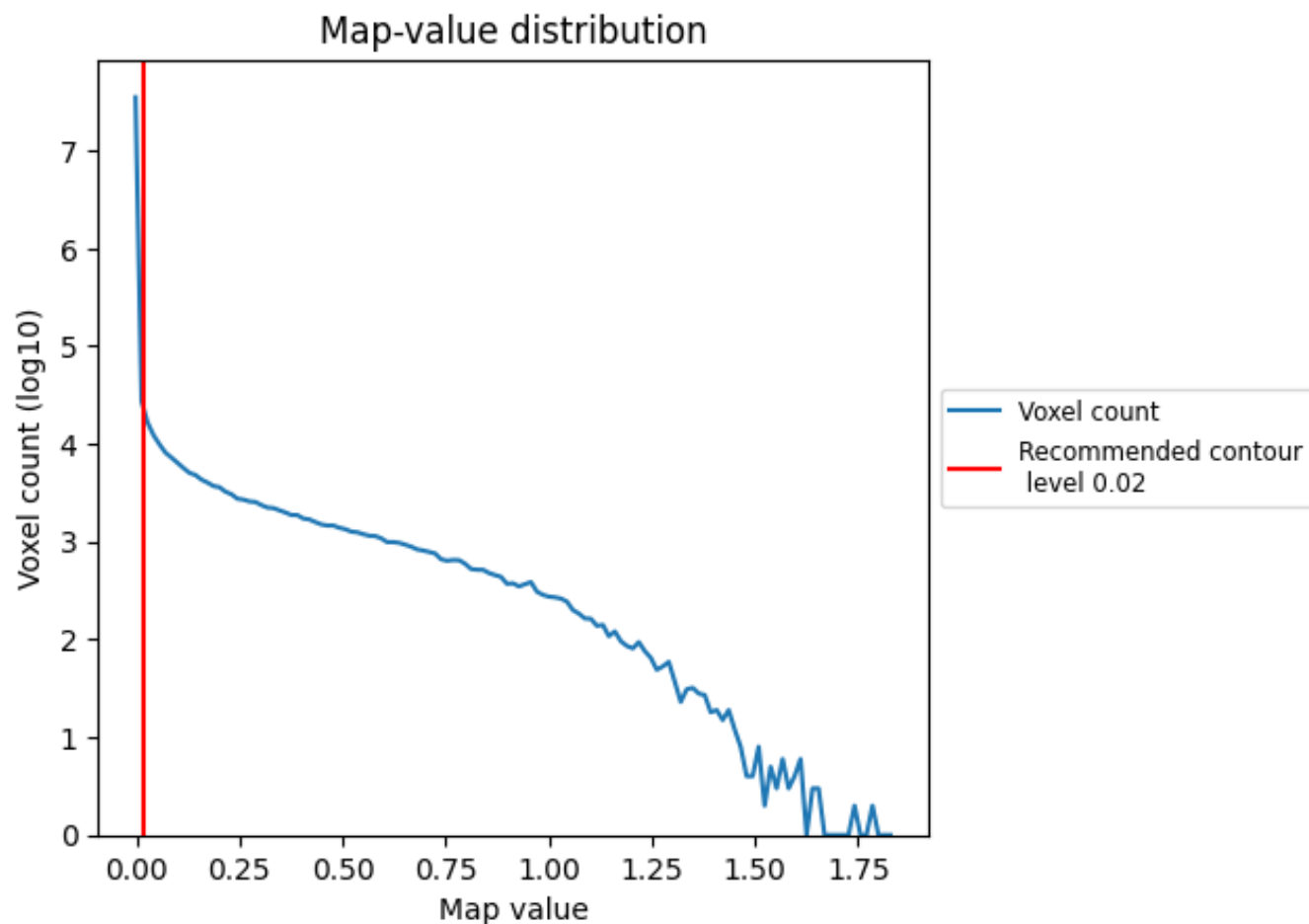
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

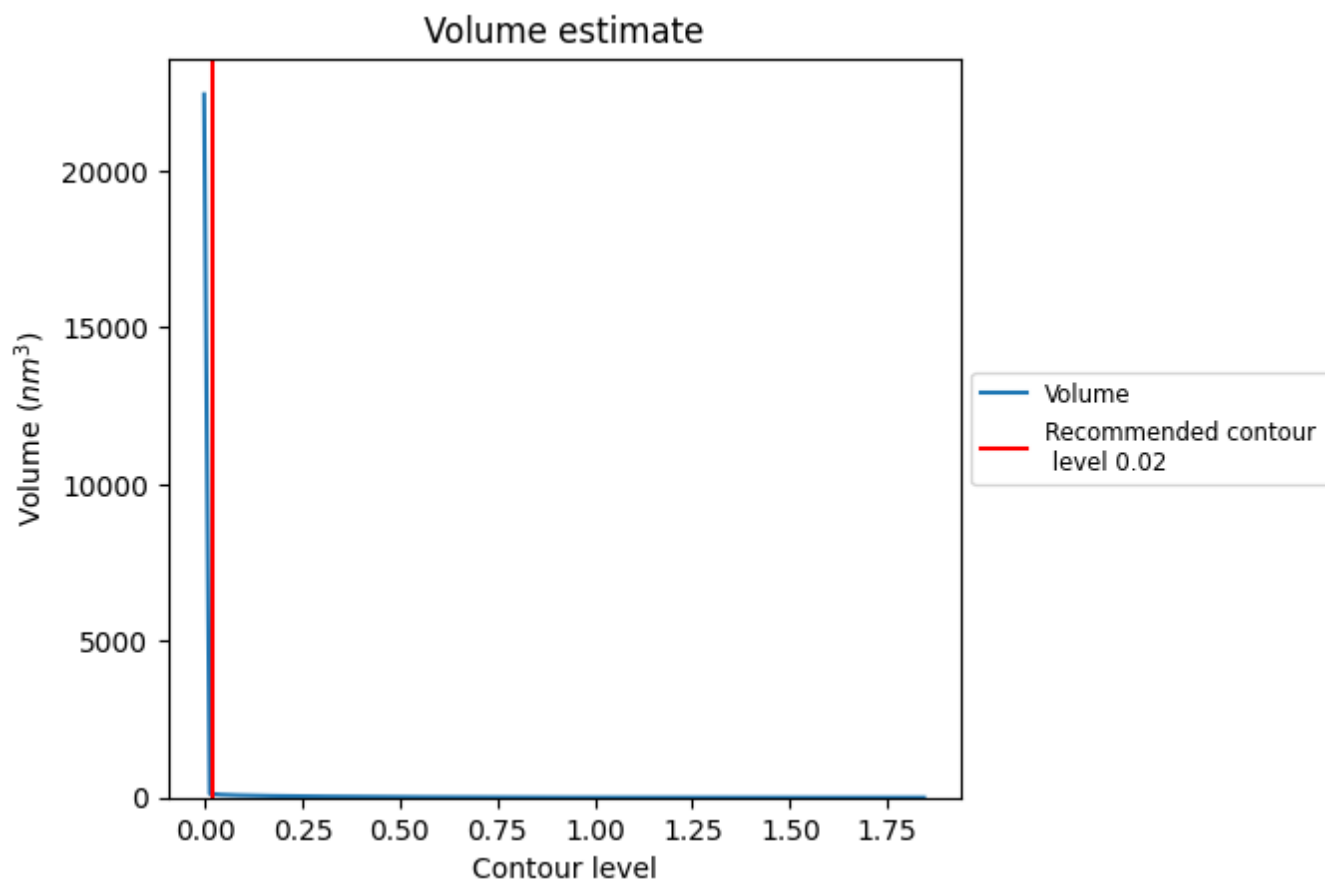
This section contains the results of statistical analysis of the map.

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.

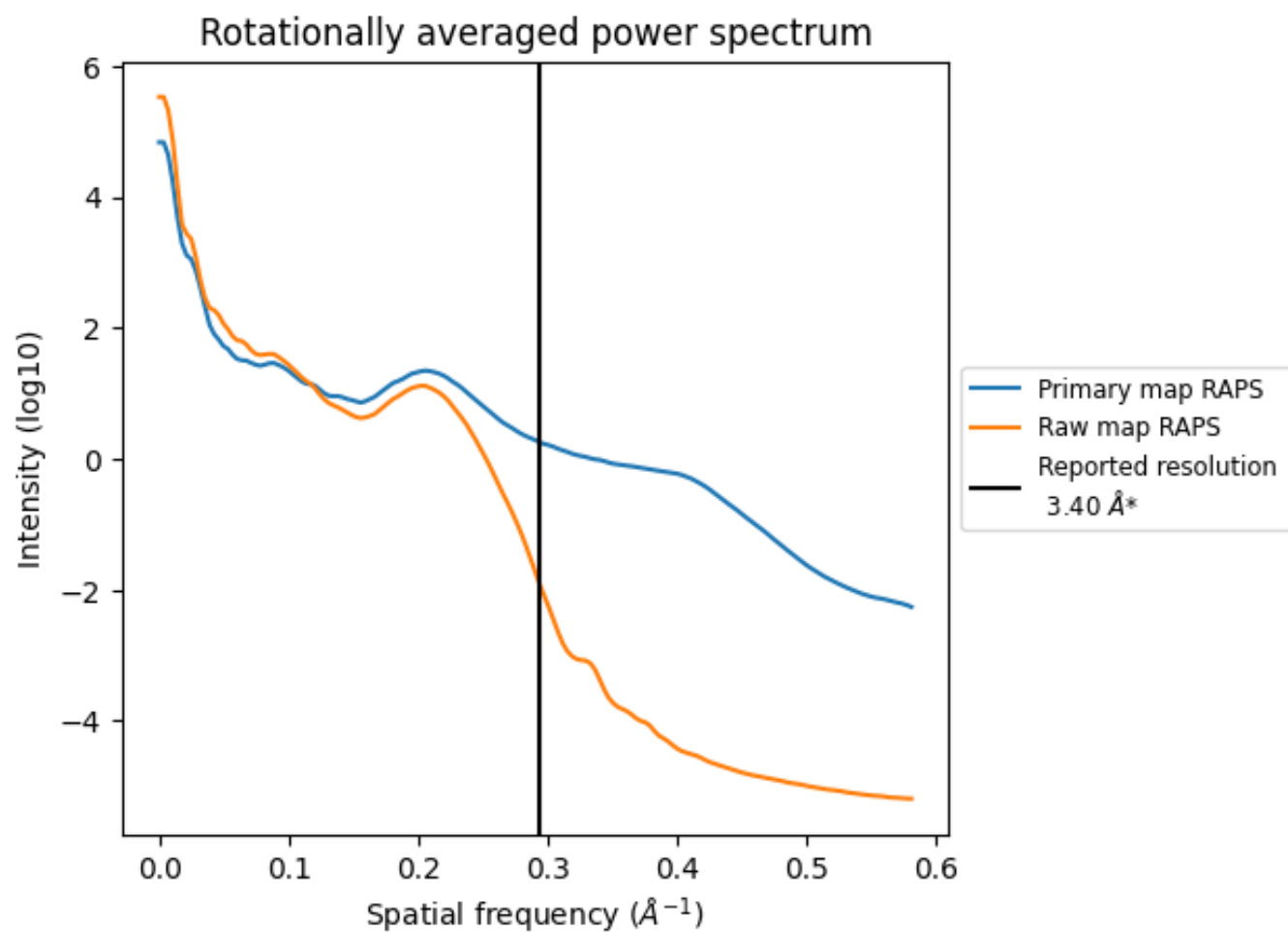
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 112 nm³; this corresponds to an approximate mass of 101 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 ⓘ

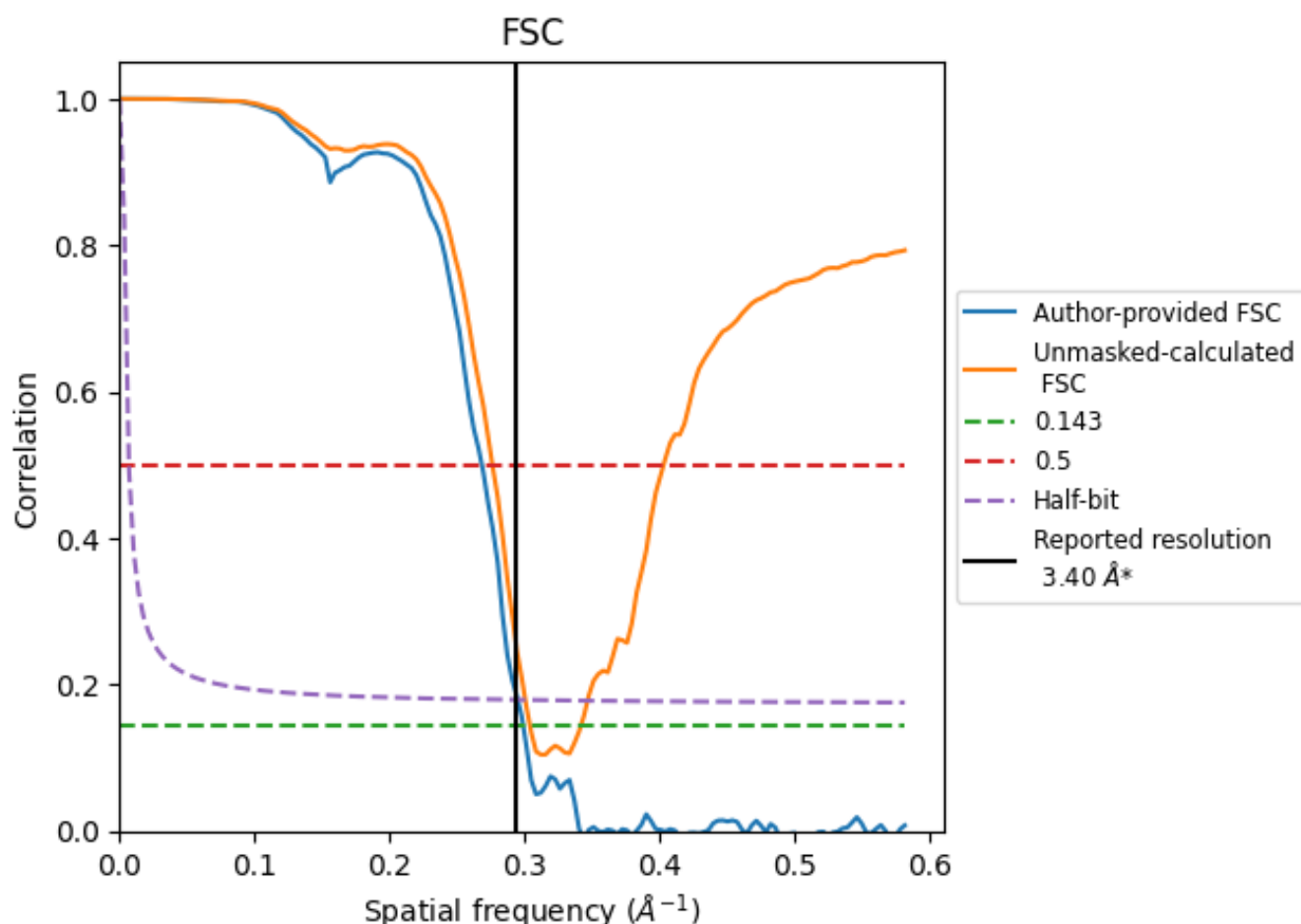


*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.294 \AA^{-1}

8.2 Resolution estimates [i](#)

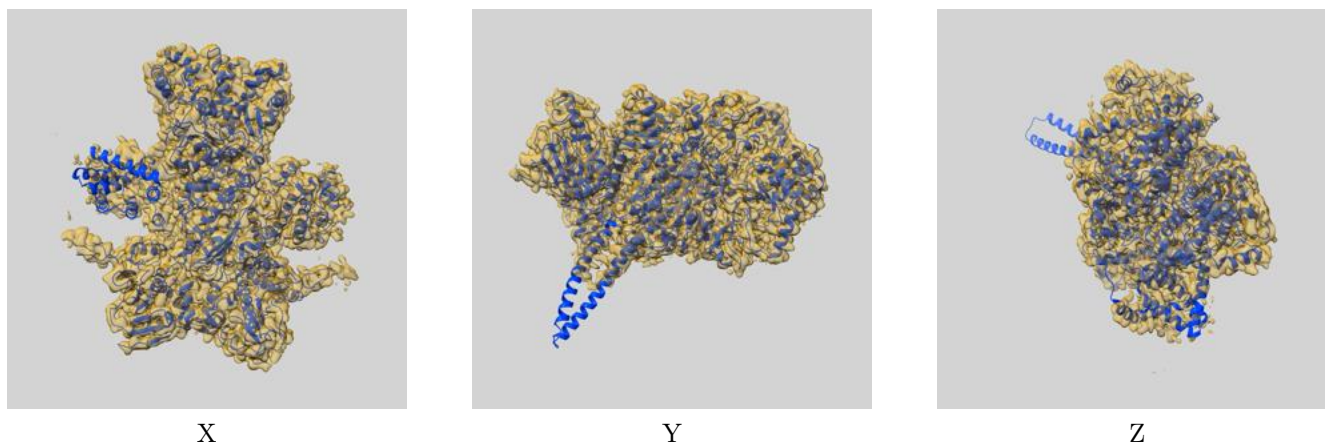
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	3.35	3.73	3.39
Unmasked-calculated*	3.29	3.62	3.33

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

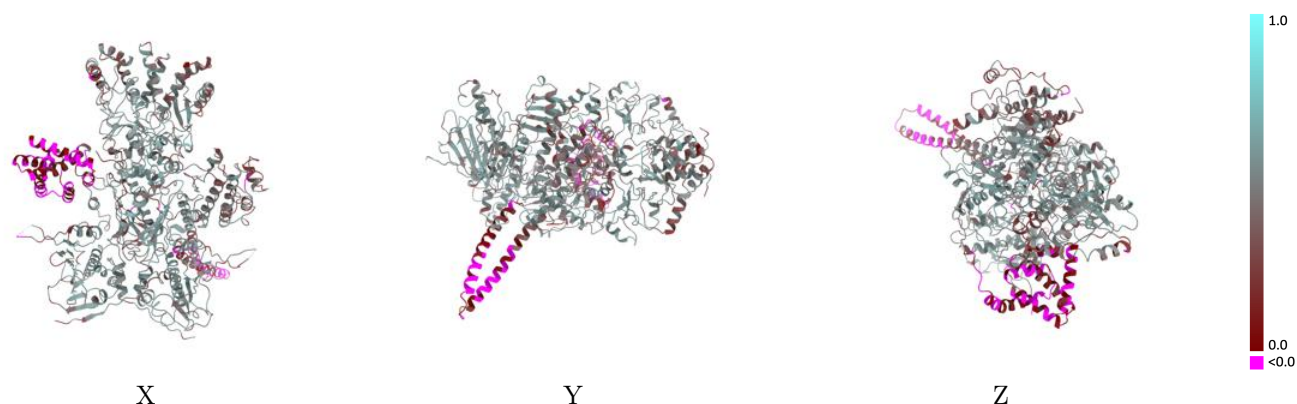
This section contains information regarding the fit between EMDB map EMD-14875 and PDB model 7ZR0. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



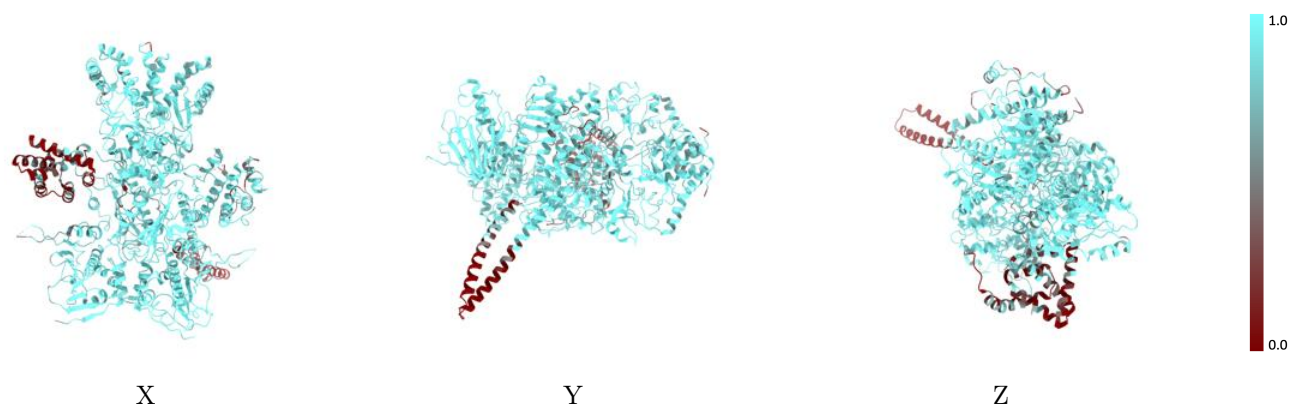
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



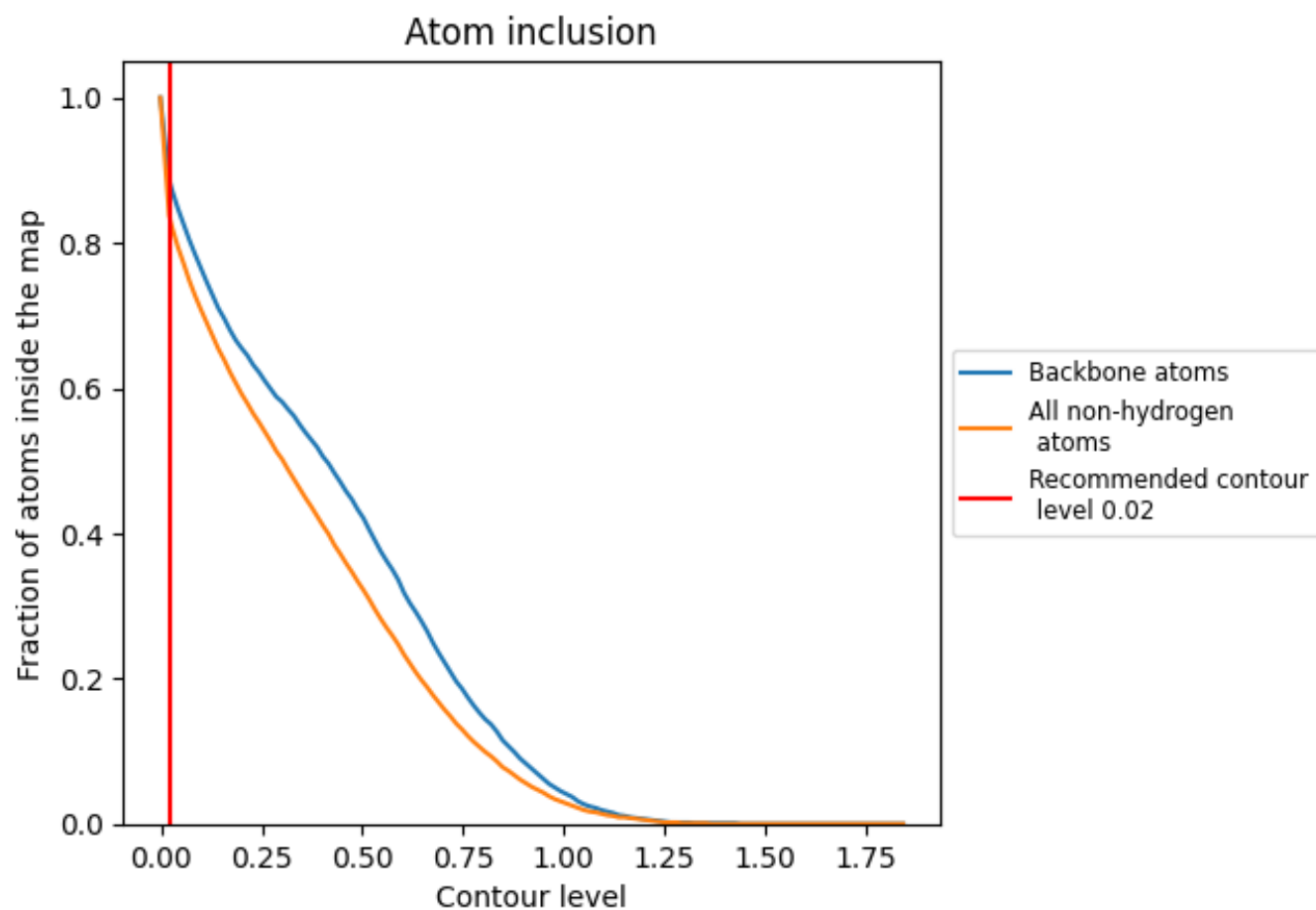
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.02).

9.4 Atom inclusion ⓘ



At the recommended contour level, 89% of all backbone atoms, 84% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8380	<div></div> 0.4270
A	<div></div> 0.9230	<div></div> 0.4890
B	<div></div> 0.9260	<div></div> 0.5030
C	<div></div> 0.4360	<div></div> 0.1340
K	<div></div> 0.8360	<div></div> 0.3960

