



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 05:53 PM UTC

PDB ID : 5FWL / pdb_00005fwl
EMDB ID : EMD-3341
Title : Atomic cryoEM structure of Hsp90-Cdc37-Cdk4 complex
Authors : Verba, K.A.; Wang, R.Y.R.; Arakawa, A.; Liu, Y.; Yokoyama, S.; Agard, D.A.
Deposited on : 2016-02-18
Resolution : 9.00 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.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

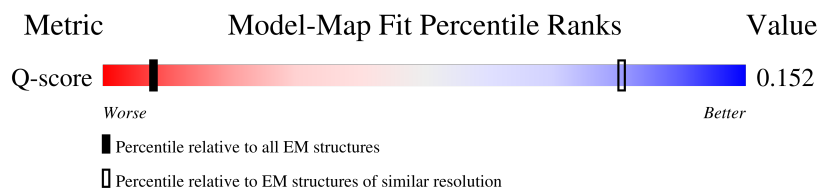
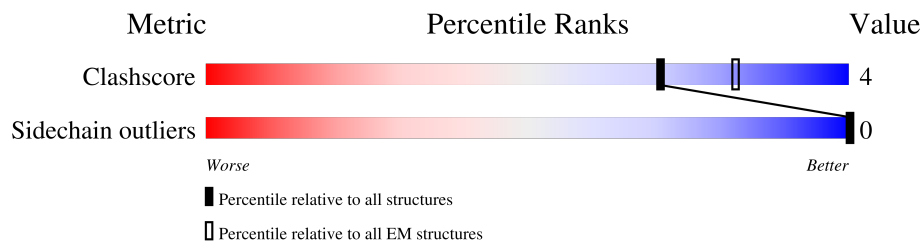
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 9.00 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	257 (8.50 - 9.50)

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	727	
1	B	727	
2	E	378	
3	K	310	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 29596 atoms, of which 14789 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	638	Total	C	H	N	O	S	0	0	
			10399	3280	5226	870	1000	23			
1	B	630	Total	C	H	N	O	S	0	0	
			10242	3235	5139	856	989	23			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P08238
A	-1	GLY	-	expression tag	UNP P08238
A	0	PHE	-	expression tag	UNP P08238
B	-2	GLY	-	expression tag	UNP P08238
B	-1	GLY	-	expression tag	UNP P08238
B	0	PHE	-	expression tag	UNP P08238

- Molecule 2 is a protein called HSP90 CO-CHAPERONE CDC37.

Mol	Chain	Residues	Atoms							AltConf	Trace
2	E	259	Total	C	H	N	O	P	S	0	0
			4307	1358	2125	383	425	1	15		

- Molecule 3 is a protein called CYCLIN-DEPENDENT KINASE 4.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	K	291	Total	C	H	N	O	S	0	0
			4584	1463	2299	403	408	11		

There are 7 discrepancies between the modelled and reference sequences:

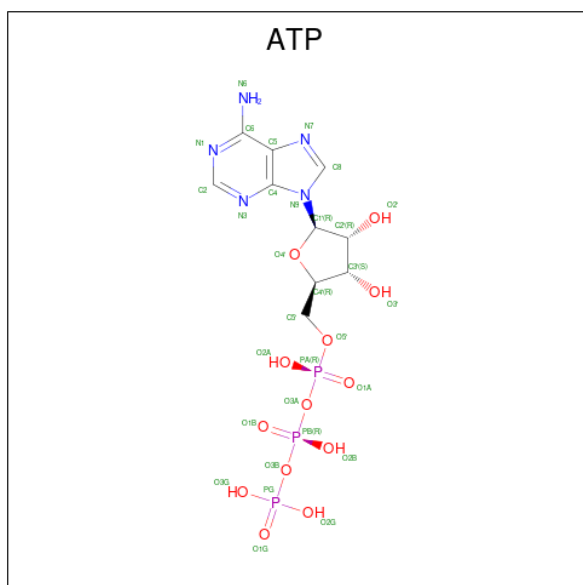
Chain	Residue	Modelled	Actual	Comment	Reference
K	-6	GLY	-	expression tag	UNP P11802
K	-5	ALA	-	expression tag	UNP P11802
K	-4	MET	-	expression tag	UNP P11802

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Chain	Residue	Modelled	Actual	Comment	Reference
K	-3	ASP	-	expression tag	UNP P11802
K	-2	PRO	-	expression tag	UNP P11802
K	-1	GLU	-	expression tag	UNP P11802
K	0	PHE	-	expression tag	UNP P11802

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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	

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
5	A	1	Total	Mg	0
			1	1	
5	B	1	Total	Mg	0
			1	1	

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	45974	Depositor
Resolution determination method	Not provided	
CTF correction method	Not provided	
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	44	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	3800	Depositor
Magnification	22500	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.085	Depositor
Minimum map value	-0.024	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.02	Depositor
Map size ($\text{\AA}$)	336.64, 336.64, 336.64	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.315, 1.315, 1.315	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 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	0.95	0/5257	0.93	27/7062 (0.4%)
1	B	0.94	1/5187 (0.0%)	0.95	33/6972 (0.5%)
2	E	0.92	3/2209 (0.1%)	0.83	6/2955 (0.2%)
3	K	0.96	2/2342 (0.1%)	1.20	39/3181 (1.2%)
All	All	0.94	6/14995 (0.0%)	0.97	105/20170 (0.5%)

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
3	K	0	8

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	K	85	ARG	CA-C	6.46	1.61	1.52
2	E	132	LYS	CA-C	5.70	1.60	1.52
1	B	629	ASN	CA-CB	-5.61	1.46	1.53
2	E	237	PHE	CA-C	-5.29	1.45	1.52
2	E	131	THR	CA-C	-5.13	1.46	1.52
3	K	59	LEU	CA-C	-5.08	1.46	1.52

All (105) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	42	GLY	CA-C-O	-10.05	115.69	122.22
3	K	250	PRO	CA-C-N	9.75	129.83	119.89
3	K	250	PRO	C-N-CA	9.75	129.83	119.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	587	SER	CA-C-N	9.72	129.38	119.56
1	B	587	SER	C-N-CA	9.72	129.38	119.56
1	A	629	ASN	CA-C-N	9.29	129.04	119.56
1	A	629	ASN	C-N-CA	9.29	129.04	119.56
3	K	253	GLY	CA-C-N	9.20	129.32	120.31
3	K	253	GLY	C-N-CA	9.20	129.32	120.31
3	K	108	PRO	CA-C-N	9.16	129.81	120.38
3	K	108	PRO	C-N-CA	9.16	129.81	120.38
1	B	629	ASN	CA-C-N	8.95	128.59	119.82
1	B	629	ASN	C-N-CA	8.95	128.59	119.82
1	B	78	ASN	CA-C-N	8.93	129.10	119.28
1	B	78	ASN	C-N-CA	8.93	129.10	119.28
1	A	61	ASP	CA-C-N	8.75	129.62	119.47
1	A	61	ASP	C-N-CA	8.75	129.62	119.47
1	A	587	SER	CA-C-N	8.71	128.44	119.64
1	A	587	SER	C-N-CA	8.71	128.44	119.64
3	K	261	VAL	CA-C-N	8.59	129.12	119.93
3	K	261	VAL	C-N-CA	8.59	129.12	119.93
1	B	632	HIS	CA-C-N	8.52	128.66	119.28
1	B	632	HIS	C-N-CA	8.52	128.66	119.28
1	B	672	ASP	CA-C-N	8.44	129.26	119.47
1	B	672	ASP	C-N-CA	8.44	129.26	119.47
1	A	335	ILE	CA-C-N	8.39	128.42	119.78
1	A	335	ILE	C-N-CA	8.39	128.42	119.78
1	B	370	ILE	CA-C-N	8.37	128.40	119.78
1	B	370	ILE	C-N-CA	8.37	128.40	119.78
1	B	76	ILE	CA-C-N	8.34	128.37	119.78
1	B	76	ILE	C-N-CA	8.34	128.37	119.78
1	A	672	ASP	CA-C-N	8.33	129.13	119.47
1	A	672	ASP	C-N-CA	8.33	129.13	119.47
1	A	78	ASN	CA-C-N	8.17	128.95	119.47
1	A	78	ASN	C-N-CA	8.17	128.95	119.47
1	A	632	HIS	CA-C-N	8.12	128.58	119.32
1	A	632	HIS	C-N-CA	8.12	128.58	119.32
3	K	107	ALA	CA-C-N	8.12	128.74	120.38
3	K	107	ALA	C-N-CA	8.12	128.74	120.38
2	E	229	ASP	CA-C-N	8.07	128.51	119.32
2	E	229	ASP	C-N-CA	8.07	128.51	119.32
1	B	292	ASN	CA-C-N	8.06	128.82	119.47
1	B	292	ASN	C-N-CA	8.06	128.82	119.47
3	K	279	ASN	CA-C-N	8.00	127.72	119.56
3	K	279	ASN	C-N-CA	8.00	127.72	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	286	LYS	CA-C-N	7.93	127.93	119.76
1	B	286	LYS	C-N-CA	7.93	127.93	119.76
1	B	335	ILE	CA-C-N	7.88	127.81	119.85
1	B	335	ILE	C-N-CA	7.88	127.81	119.85
1	A	339	ALA	CA-C-N	7.67	127.82	120.31
1	A	339	ALA	C-N-CA	7.67	127.82	120.31
1	B	61	ASP	CA-C-N	7.66	129.42	119.84
1	B	61	ASP	C-N-CA	7.66	129.42	119.84
3	K	238	TRP	CA-C-N	7.64	127.60	120.03
3	K	238	TRP	C-N-CA	7.64	127.60	120.03
2	E	20	HIS	CA-C-N	7.55	127.58	119.28
2	E	20	HIS	C-N-CA	7.55	127.58	119.28
2	E	112	MET	CA-C-N	7.42	127.40	119.76
2	E	112	MET	C-N-CA	7.42	127.40	119.76
3	K	233	PRO	CA-C-N	7.17	128.80	119.84
3	K	233	PRO	C-N-CA	7.17	128.80	119.84
3	K	109	PRO	CA-C-N	6.85	128.40	119.84
3	K	109	PRO	C-N-CA	6.85	128.40	119.84
1	A	292	ASN	CA-C-N	6.82	128.37	119.84
1	A	292	ASN	C-N-CA	6.82	128.37	119.84
1	A	370	ILE	CA-C-N	6.74	126.65	119.85
1	A	370	ILE	C-N-CA	6.74	126.65	119.85
1	A	286	LYS	CA-C-N	6.68	127.05	119.83
1	A	286	LYS	C-N-CA	6.68	127.05	119.83
1	B	211	TYR	CA-C-N	6.67	127.03	119.83
1	B	211	TYR	C-N-CA	6.67	127.03	119.83
3	K	172	THR	CA-C-N	6.66	126.36	119.56
3	K	172	THR	C-N-CA	6.66	126.36	119.56
1	A	211	TYR	CA-C-N	6.63	126.86	119.90
1	A	211	TYR	C-N-CA	6.63	126.86	119.90
1	B	515	GLU	CA-C-N	6.45	126.03	119.19
1	B	515	GLU	C-N-CA	6.45	126.03	119.19
3	K	68	HIS	CA-C-N	6.34	127.77	119.84
3	K	68	HIS	C-N-CA	6.34	127.77	119.84
3	K	193	THR	CA-C-N	6.24	125.80	119.19
3	K	193	THR	C-N-CA	6.24	125.80	119.19
3	K	182	ALA	CA-C-N	6.07	126.24	119.32
3	K	182	ALA	C-N-CA	6.07	126.24	119.32
1	B	339	ALA	CA-C-N	5.96	126.46	120.14
1	B	339	ALA	C-N-CA	5.96	126.46	120.14
3	K	249	PHE	CA-C-N	5.95	126.51	120.38
3	K	249	PHE	C-N-CA	5.95	126.51	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	244	LEU	CA-C-N	5.66	126.91	119.84
3	K	244	LEU	C-N-CA	5.66	126.91	119.84
3	K	25	ASP	CA-C-N	5.62	126.87	119.84
3	K	25	ASP	C-N-CA	5.62	126.87	119.84
1	B	173	GLU	CA-C-N	5.61	127.58	120.51
1	B	173	GLU	C-N-CA	5.61	127.58	120.51
3	K	7	GLU	CA-C-N	5.61	126.49	119.98
3	K	7	GLU	C-N-CA	5.61	126.49	119.98
1	B	543	LEU	CA-C-N	5.60	126.85	119.84
1	B	543	LEU	C-N-CA	5.60	126.85	119.84
1	A	76	ILE	CA-C-N	5.48	125.95	120.14
1	A	76	ILE	C-N-CA	5.48	125.95	120.14
3	K	142	LYS	CA-C-N	5.34	126.51	119.84
3	K	142	LYS	C-N-CA	5.34	126.51	119.84
1	A	54	ILE	CB-CA-C	-5.23	105.27	111.97
1	B	54	ILE	CB-CA-C	-5.12	105.42	111.97
3	K	211	LYS	CA-C-N	5.10	126.21	119.84
3	K	211	LYS	C-N-CA	5.10	126.21	119.84

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	K	42	GLY	Mainchain,Peptide
3	K	81	SER	Mainchain,Peptide
3	K	85	ARG	Mainchain,Peptide
3	K	86	GLU	Mainchain,Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5173	5226	5224	28	0
1	B	5103	5139	5137	32	0
2	E	2182	2125	2124	11	0
3	K	2285	2299	2297	35	0
4	A	31	0	12	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	31	0	12	2	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
All	All	14807	14789	14806	104	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 4.

All (104) 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:B:609:GLN:HB3	1:B:612:ARG:CG	1.76	1.15
1:B:609:GLN:HB3	1:B:612:ARG:HG2	1.44	0.98
1:B:609:GLN:CB	1:B:612:ARG:HG2	1.96	0.95
1:B:609:GLN:CG	1:B:612:ARG:HG2	1.96	0.94
1:B:609:GLN:HB3	1:B:612:ARG:HG3	1.52	0.89
4:A:1691:ATP:O2G	4:A:1691:ATP:O2A	2.00	0.79
3:K:87:ILE:O	3:K:87:ILE:HG22	1.90	0.71
2:E:132:LYS:HB3	2:E:133:PRO:HA	1.71	0.71
4:B:1692:ATP:O2G	4:B:1692:ATP:O2B	2.03	0.71
1:B:609:GLN:CB	1:B:612:ARG:CG	2.56	0.68
3:K:41:ASN:O	3:K:41:ASN:OD1	2.15	0.65
1:B:609:GLN:HG3	1:B:612:ARG:HG2	1.81	0.61
1:B:425:ASN:ND2	2:E:131:THR:OG1	2.35	0.60
1:A:279:GLN:O	1:A:279:GLN:HG3	2.01	0.60
1:B:80:GLN:N	1:B:80:GLN:OE1	2.35	0.59
3:K:85:ARG:HB3	3:K:86:GLU:HA	1.84	0.58
2:E:132:LYS:HB3	2:E:133:PRO:CA	2.33	0.58
1:B:609:GLN:HG2	1:B:612:ARG:CD	2.33	0.57
1:A:221:ARG:HB2	1:A:273:LYS:HB2	1.86	0.57
1:B:609:GLN:CG	1:B:612:ARG:CG	2.77	0.57
2:E:131:THR:O	2:E:132:LYS:HB2	2.05	0.56
3:K:85:ARG:HB3	3:K:86:GLU:CA	2.36	0.56
1:B:298:GLN:O	1:B:298:GLN:NE2	2.40	0.54
1:A:402:LYS:HG2	3:K:55:ARG:HB3	1.89	0.54
1:A:38:ILE:HD12	1:A:38:ILE:C	2.34	0.53
3:K:50:PRO:CD	3:K:50:PRO:O	2.56	0.53
3:K:255:ARG:O	3:K:256:PRO:C	2.52	0.53
1:A:128:GLN:O	1:A:128:GLN:CD	2.52	0.52
2:E:132:LYS:CB	2:E:133:PRO:HA	2.38	0.52
3:K:189:SER:OG	3:K:190:THR:N	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1691:ATP:N3	4:A:1691:ATP:H2'	2.24	0.51
1:A:402:LYS:HD3	3:K:59:LEU:HB2	1.93	0.51
1:B:93:MET:SD	4:B:1692:ATP:H1'	2.51	0.51
1:B:366:CYS:HA	1:B:369:LEU:HB3	1.93	0.50
1:A:486:ILE:O	1:A:486:ILE:HG23	2.12	0.50
1:A:124:SER:OG	1:A:358:ARG:NH1	2.45	0.49
2:E:131:THR:O	2:E:132:LYS:CB	2.59	0.49
4:A:1691:ATP:O1G	4:A:1691:ATP:O3A	2.30	0.49
1:B:220:GLU:HG2	1:B:220:GLU:O	2.13	0.49
3:K:206:GLU:O	3:K:210:ARG:N	2.46	0.49
1:A:128:GLN:O	1:A:128:GLN:OE1	2.30	0.49
1:A:399:LYS:HA	1:A:402:LYS:HD2	1.94	0.49
1:B:609:GLN:HG2	1:B:612:ARG:HD3	1.93	0.49
3:K:71:VAL:HG23	3:K:71:VAL:O	2.13	0.48
3:K:75:MET:O	3:K:76:ASP:CB	2.61	0.48
3:K:187:LEU:HD22	3:K:242:VAL:HG13	1.96	0.48
3:K:210:ARG:HG2	3:K:210:ARG:O	2.14	0.48
3:K:142:LYS:HB2	3:K:143:PRO:CD	2.43	0.48
1:B:157:TRP:O	1:B:157:TRP:HE3	1.97	0.47
2:E:222:LEU:HB2	2:E:237:PHE:CE2	2.49	0.47
3:K:142:LYS:HB2	3:K:143:PRO:HD2	1.96	0.47
3:K:81:SER:HA	3:K:82:ARG:HB2	1.96	0.47
3:K:77:VAL:HG22	3:K:77:VAL:O	2.14	0.47
1:A:473:VAL:HA	1:A:476:MET:SD	2.55	0.46
3:K:84:ASP:O	3:K:85:ARG:HB2	2.16	0.46
1:B:207:GLN:OE1	1:B:207:GLN:N	2.32	0.46
3:K:85:ARG:HB3	3:K:86:GLU:CB	2.46	0.46
1:B:219:LYS:O	1:B:220:GLU:C	2.59	0.46
1:A:347:LYS:HA	3:K:44:GLY:HA3	1.98	0.46
1:A:430:TYR:O	1:A:434:SER:N	2.48	0.46
3:K:244:LEU:HB3	3:K:245:PRO:HD2	1.98	0.46
3:K:34:LEU:HG	3:K:82:ARG:HD2	1.99	0.45
1:A:79:PRO:HA	1:A:193:LEU:HD13	1.99	0.45
1:A:456:ARG:HB3	1:A:466:MET:HG2	1.98	0.45
1:A:594:SER:OG	1:A:596:TYR:O	2.34	0.44
3:K:50:PRO:O	3:K:50:PRO:HD2	2.17	0.44
1:A:67:SER:OG	1:A:177:ARG:NH2	2.50	0.44
1:A:597:GLY:HA2	1:A:621:MET:SD	2.56	0.44
2:E:185:GLU:OE1	2:E:185:GLU:N	2.39	0.44
3:K:183:PRO:O	3:K:187:LEU:HG	2.18	0.44
1:B:486:ILE:HG23	1:B:486:ILE:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:260:LYS:HG2	2:E:260:LYS:O	2.17	0.43
1:B:473:VAL:HA	1:B:476:MET:SD	2.58	0.43
1:A:629:ASN:HA	1:A:630:PRO:HD2	1.87	0.43
3:K:161:LEU:HA	3:K:164:ILE:HG12	2.01	0.43
1:B:342:ASP:OD1	1:B:342:ASP:C	2.61	0.43
3:K:223:LEU:HB3	3:K:227:PHE:CE2	2.53	0.43
3:K:63:LEU:HD13	3:K:71:VAL:HB	2.01	0.43
1:A:331:ALA:HA	1:A:380:VAL:O	2.19	0.43
1:B:593:THR:HG21	1:B:599:THR:HA	2.00	0.43
1:A:82:ARG:HG2	1:A:82:ARG:O	2.19	0.42
2:E:133:PRO:O	2:E:134:GLU:HB2	2.19	0.42
3:K:48:GLY:O	3:K:49:LEU:C	2.62	0.42
1:A:316:LEU:HB2	1:A:337:ARG:HA	2.02	0.42
1:B:148:LYS:HD3	1:B:148:LYS:C	2.44	0.42
2:E:25:THR:HA	2:E:28:LEU:HB3	2.01	0.42
1:B:472:TYR:CD1	1:B:472:TYR:C	2.98	0.42
1:B:78:ASN:OD1	1:B:78:ASN:C	2.63	0.41
1:B:157:TRP:O	1:B:157:TRP:CE3	2.73	0.41
3:K:198:TRP:CH2	3:K:202:CYS:SG	3.10	0.41
1:A:141:GLU:N	1:A:184:HIS:O	2.49	0.41
1:B:83:THR:HG22	1:B:184:HIS:HA	2.02	0.41
1:B:347:LYS:HA	3:K:41:ASN:CB	2.50	0.41
1:B:345:GLU:HB2	3:K:88:LYS:HD3	2.01	0.41
1:A:389:ASN:OD1	1:A:389:ASN:C	2.63	0.41
1:A:65:LEU:HB3	1:A:69:LYS:HG2	2.03	0.41
1:B:74:ASP:HB2	1:B:87:VAL:HB	2.02	0.41
1:A:362:ILE:HG22	1:A:363:MET:HG3	2.03	0.41
1:A:74:ASP:HB2	1:A:87:VAL:HB	2.02	0.40
3:K:35:LYS:O	3:K:82:ARG:NH1	2.54	0.40
3:K:21:TYR:HB2	3:K:34:LEU:HB3	2.03	0.40
1:A:564:CYS:SG	1:A:628:ILE:HD12	2.62	0.40
1:B:292:ASN:HA	1:B:293:PRO:HD3	1.89	0.40
3:K:51:ILE:HG13	3:K:55:ARG:HG2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

There are no protein backbone outliers to report in this entry.

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	
1	A	577/653 (88%)	577 (100%)	0	100	100
1	B	569/653 (87%)	569 (100%)	0	100	100
2	E	238/340 (70%)	238 (100%)	0	100	100
3	K	247/262 (94%)	247 (100%)	0	100	100
All	All	1631/1908 (86%)	1631 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	461	GLN
1	B	184	HIS
1	B	396	GLN
1	B	674	GLN
2	E	148	HIS
2	E	158	GLN
3	K	98	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](#)

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	E	13	2	8,9,10	1.48	1 (12%)	7,12,14	1.25	1 (14%)

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	E	13	2	-	0/6/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	13	SEP	P-O1P	3.39	1.61	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	13	SEP	OG-CB-CA	3.01	111.07	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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
4	ATP	B	1692	5	32,33,33	1.55	7 (21%)	48,52,52	1.77	8 (16%)
4	ATP	A	1691	5	32,33,33	1.47	6 (18%)	48,52,52	1.72	6 (12%)

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	B	1692	5	-	8/22/38/38	0/3/3/3
4	ATP	A	1691	5	-	9/22/38/38	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1692	ATP	C5-C4	4.87	1.47	1.39
4	A	1691	ATP	C5-C4	4.80	1.47	1.39
4	B	1692	ATP	PB-O3A	-3.05	1.56	1.59
4	B	1692	ATP	C5-N7	-2.79	1.34	1.39
4	A	1691	ATP	PB-O3B	-2.55	1.56	1.59
4	A	1691	ATP	C5-C6	2.53	1.48	1.41
4	B	1692	ATP	PA-O3A	-2.43	1.56	1.59
4	B	1692	ATP	C5-C6	2.40	1.47	1.41
4	B	1692	ATP	C4-N9	-2.33	1.32	1.37
4	B	1692	ATP	C8-N7	2.31	1.36	1.31
4	A	1691	ATP	C5-N7	-2.28	1.34	1.39
4	A	1691	ATP	C8-N7	2.18	1.35	1.31
4	A	1691	ATP	C4-N9	-2.08	1.33	1.37

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1692	ATP	C5-C4-N3	-5.88	118.61	126.72
4	A	1691	ATP	C5-C4-N3	-5.79	118.75	126.72
4	B	1692	ATP	N3-C4-N9	4.52	134.85	127.17
4	A	1691	ATP	N3-C4-N9	4.52	134.85	127.17
4	B	1692	ATP	C2-N3-C4	3.75	121.00	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1691	ATP	C2-N3-C4	3.74	120.96	111.83
4	B	1692	ATP	N3-C2-N1	-3.36	123.50	128.58
4	A	1691	ATP	N3-C2-N1	-3.35	123.51	128.58
4	B	1692	ATP	C4-C5-N7	-3.15	106.98	110.58
4	A	1691	ATP	C4-C5-N7	-3.06	107.09	110.58
4	B	1692	ATP	C3'-C2'-C1'	2.42	106.04	101.46
4	A	1691	ATP	O2B-PB-O1B	2.16	122.50	112.44
4	B	1692	ATP	O2G-PG-O1G	2.05	118.83	110.83
4	B	1692	ATP	O2B-PB-O1B	2.05	121.96	112.44

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1691	ATP	O4'-C4'-C5'-O5'
4	B	1692	ATP	C5'-O5'-PA-O2A
4	B	1692	ATP	C5'-O5'-PA-O3A
4	A	1691	ATP	C3'-C4'-C5'-O5'
4	B	1692	ATP	O4'-C4'-C5'-O5'
4	B	1692	ATP	C3'-C4'-C5'-O5'
4	A	1691	ATP	PA-O3A-PB-O2B
4	B	1692	ATP	PG-O3B-PB-O2B
4	B	1692	ATP	PB-O3A-PA-O2A
4	A	1691	ATP	C5'-O5'-PA-O1A
4	A	1691	ATP	C5'-O5'-PA-O2A
4	A	1691	ATP	C5'-O5'-PA-O3A
4	A	1691	ATP	C2'-C1'-N9-C8
4	A	1691	ATP	C2'-C1'-N9-C4
4	B	1692	ATP	PB-O3A-PA-O1A
4	A	1691	ATP	PA-O3A-PB-O1B
4	B	1692	ATP	PG-O3B-PB-O1B

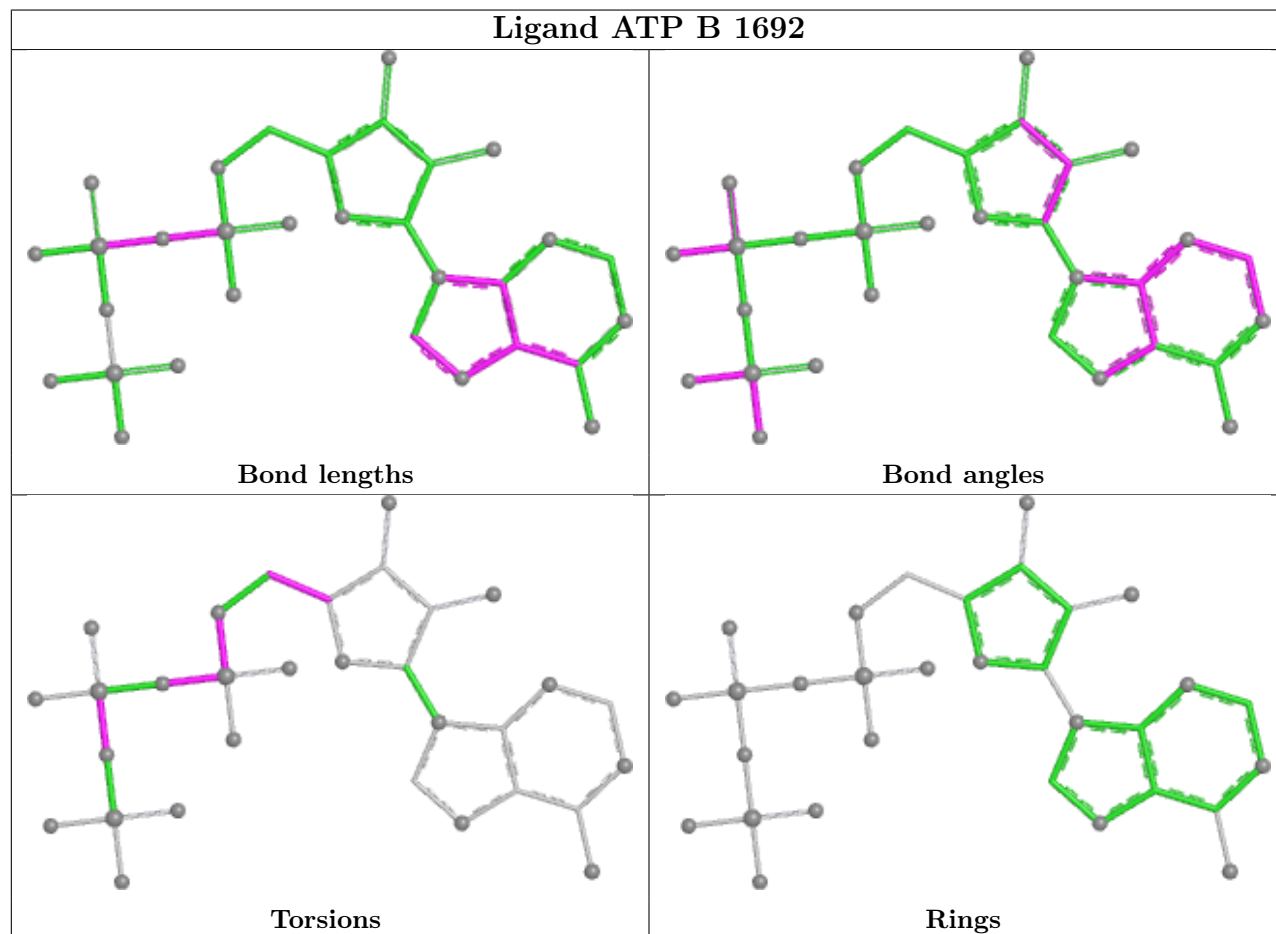
There are no ring outliers.

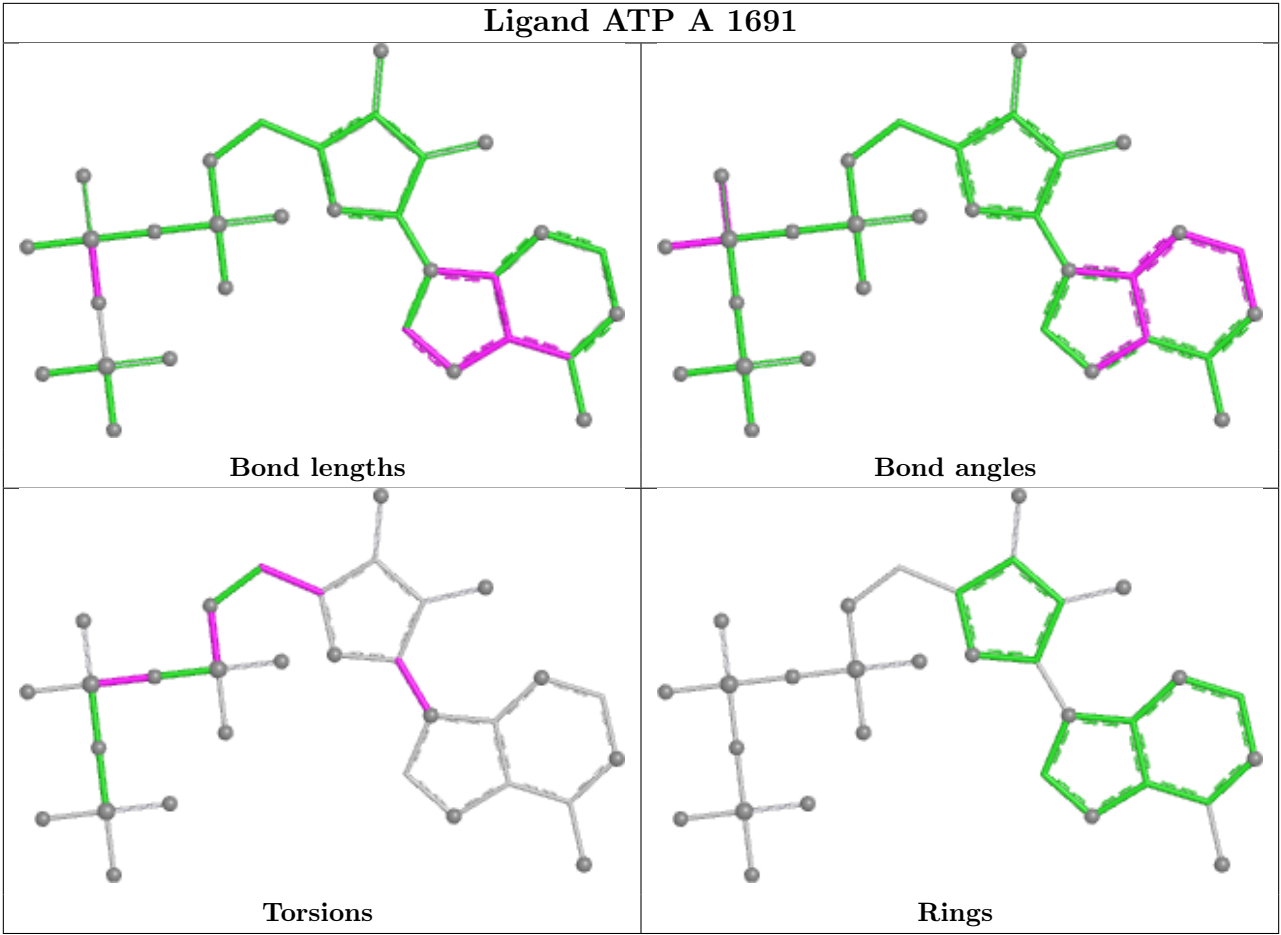
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1692	ATP	2	0
4	A	1691	ATP	3	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	K	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	K	49:LEU	C	50:PRO	N	5.53
1	K	67:GLU	C	68:HIS	N	3.83

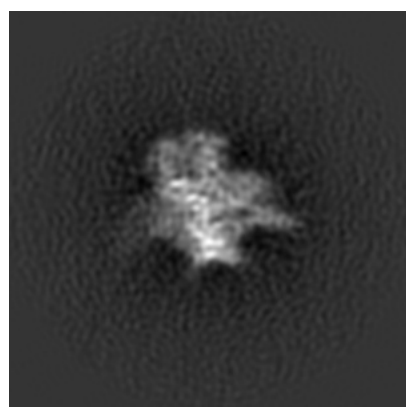
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-3341. These allow visual inspection of the internal detail of the map and identification of artifacts.

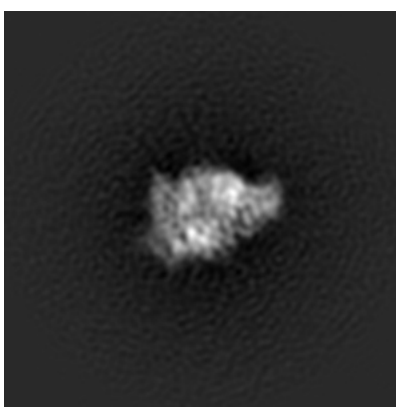
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

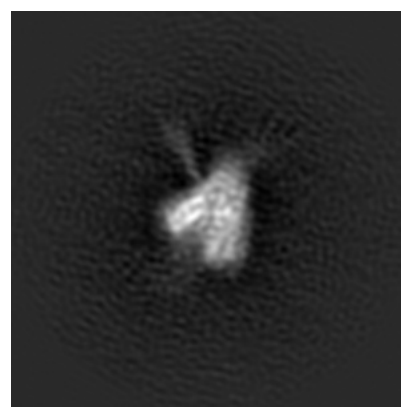
6.1.1 Primary map



X



Y

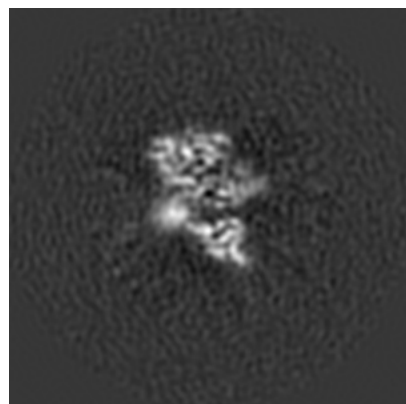


Z

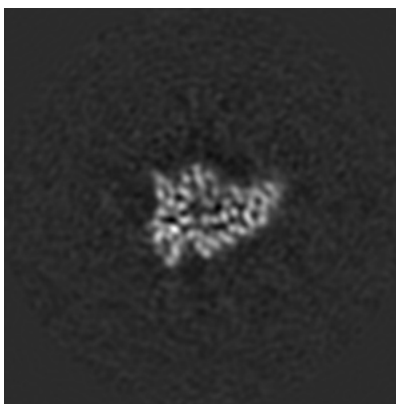
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

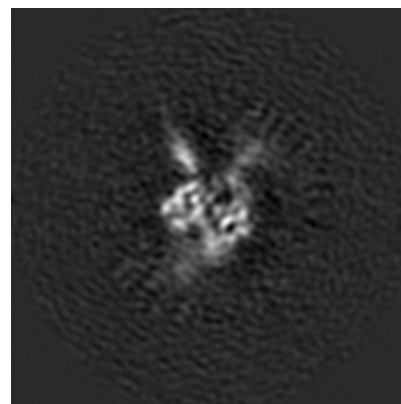
6.2.1 Primary map



X Index: 128



Y Index: 128

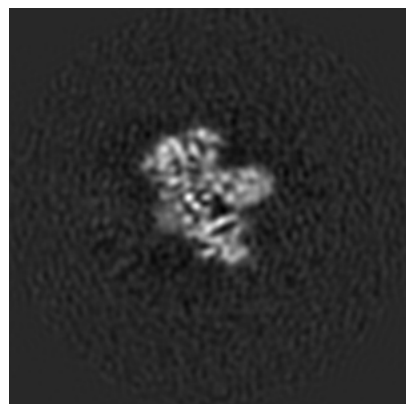


Z Index: 128

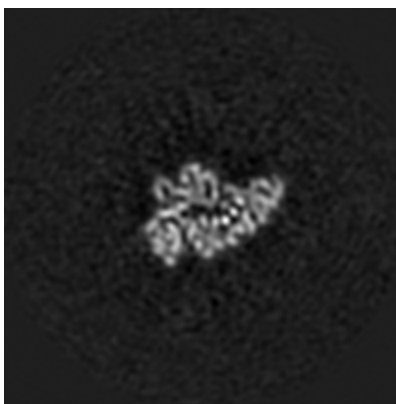
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

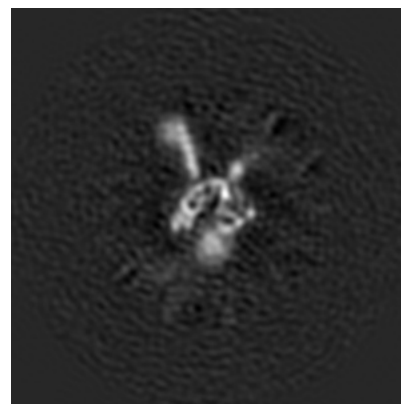
6.3.1 Primary map



X Index: 138



Y Index: 125

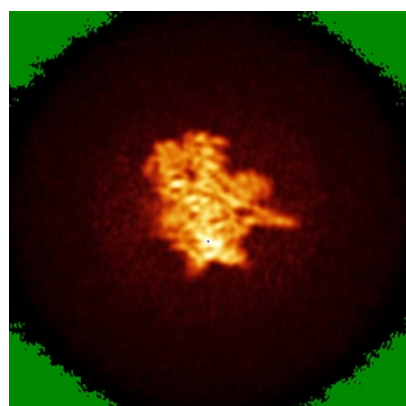


Z Index: 120

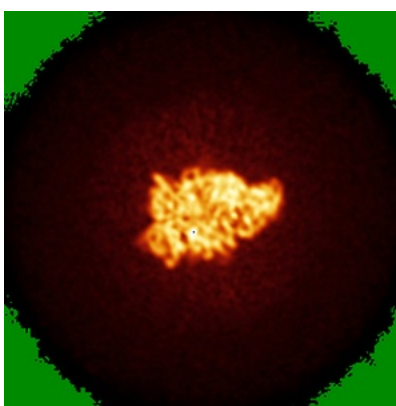
The images above show the largest variance slices of the map in three orthogonal directions.

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

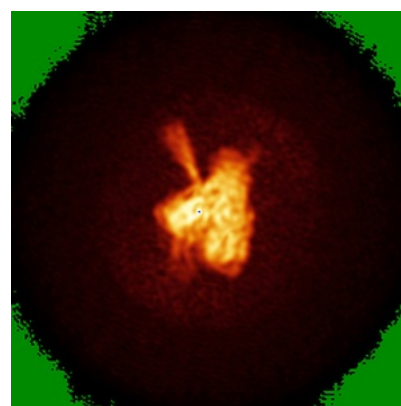
6.4.1 Primary map



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

This section was not generated.

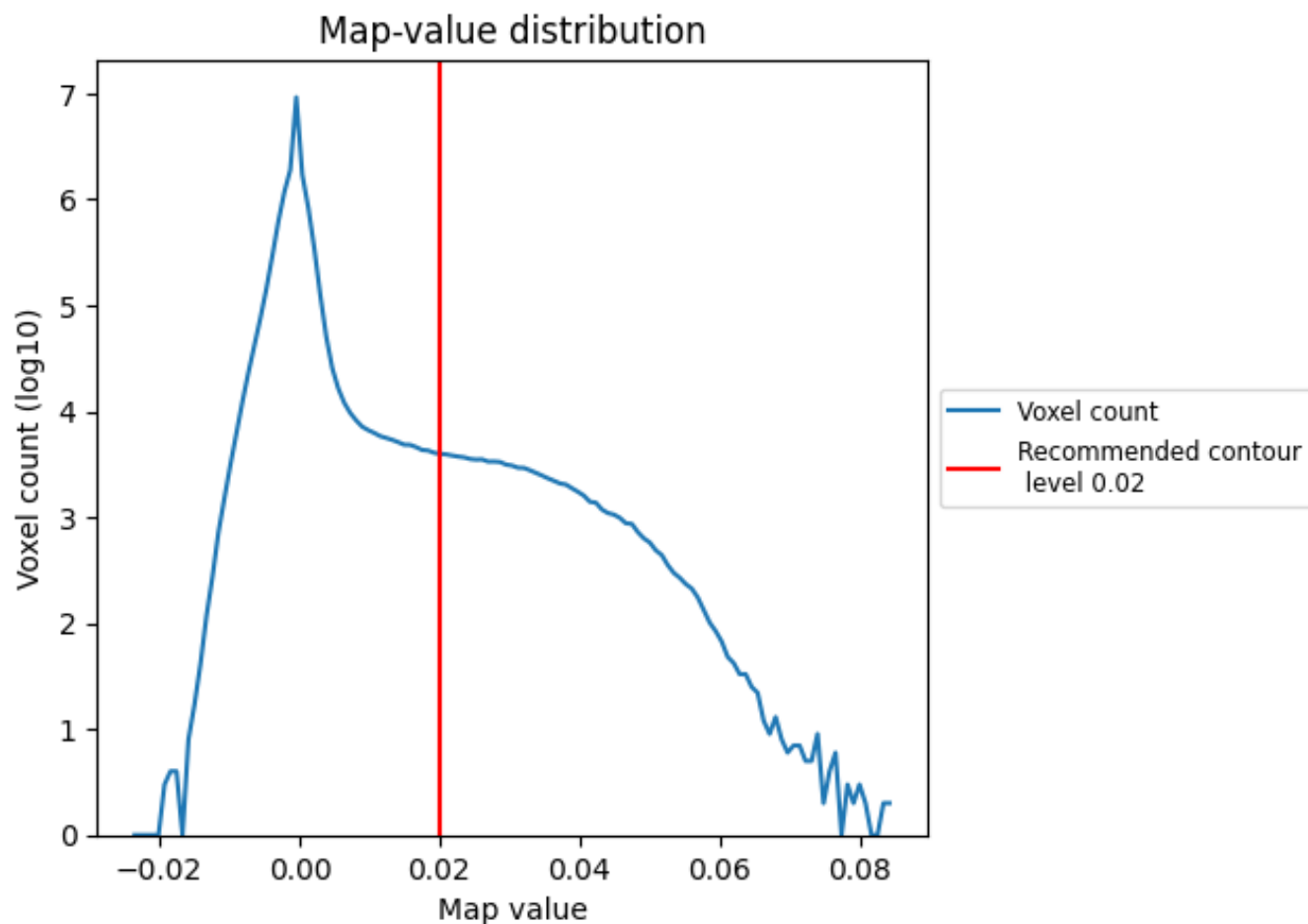
6.6 Mask visualisation

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

7 Map analysis [i](#)

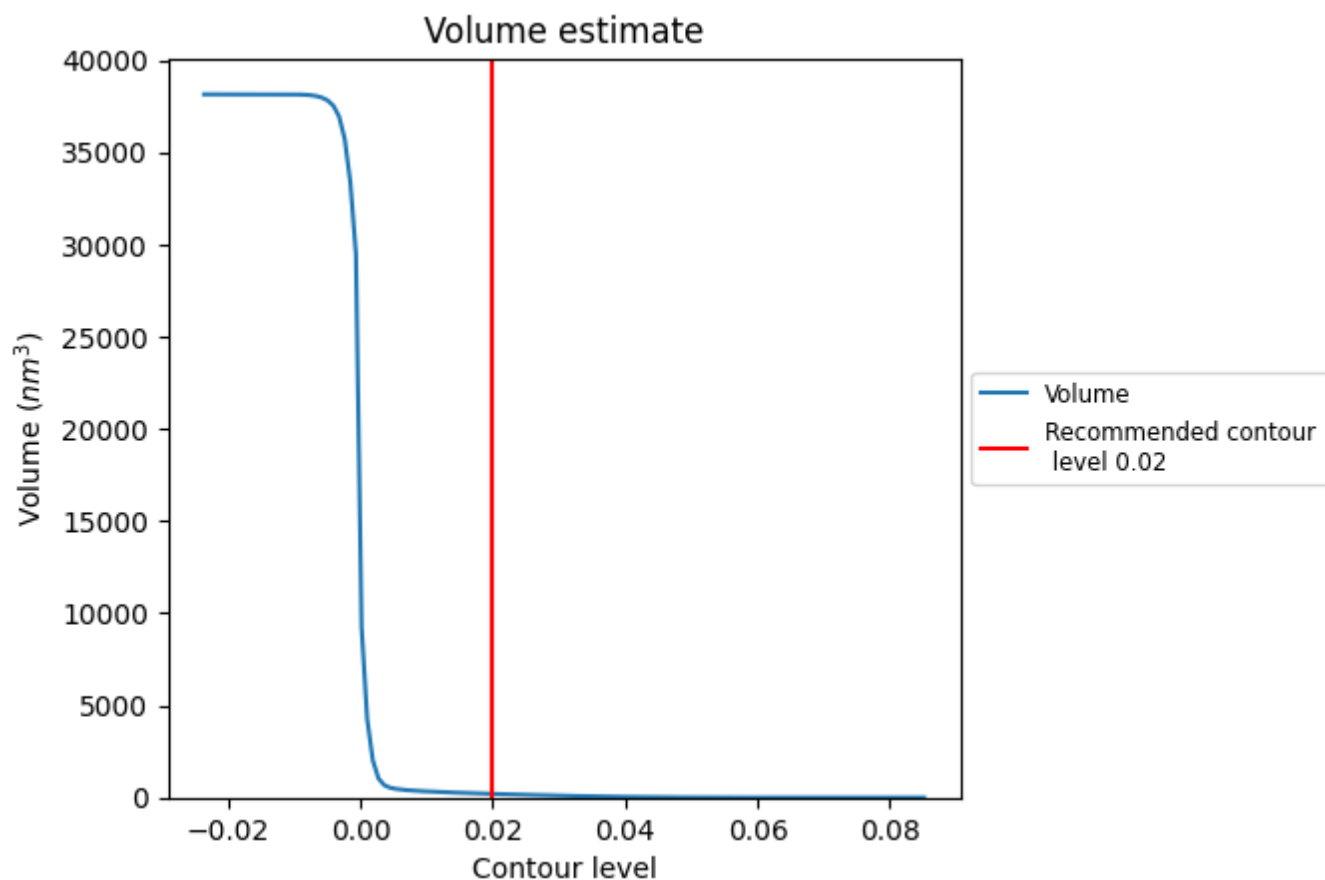
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

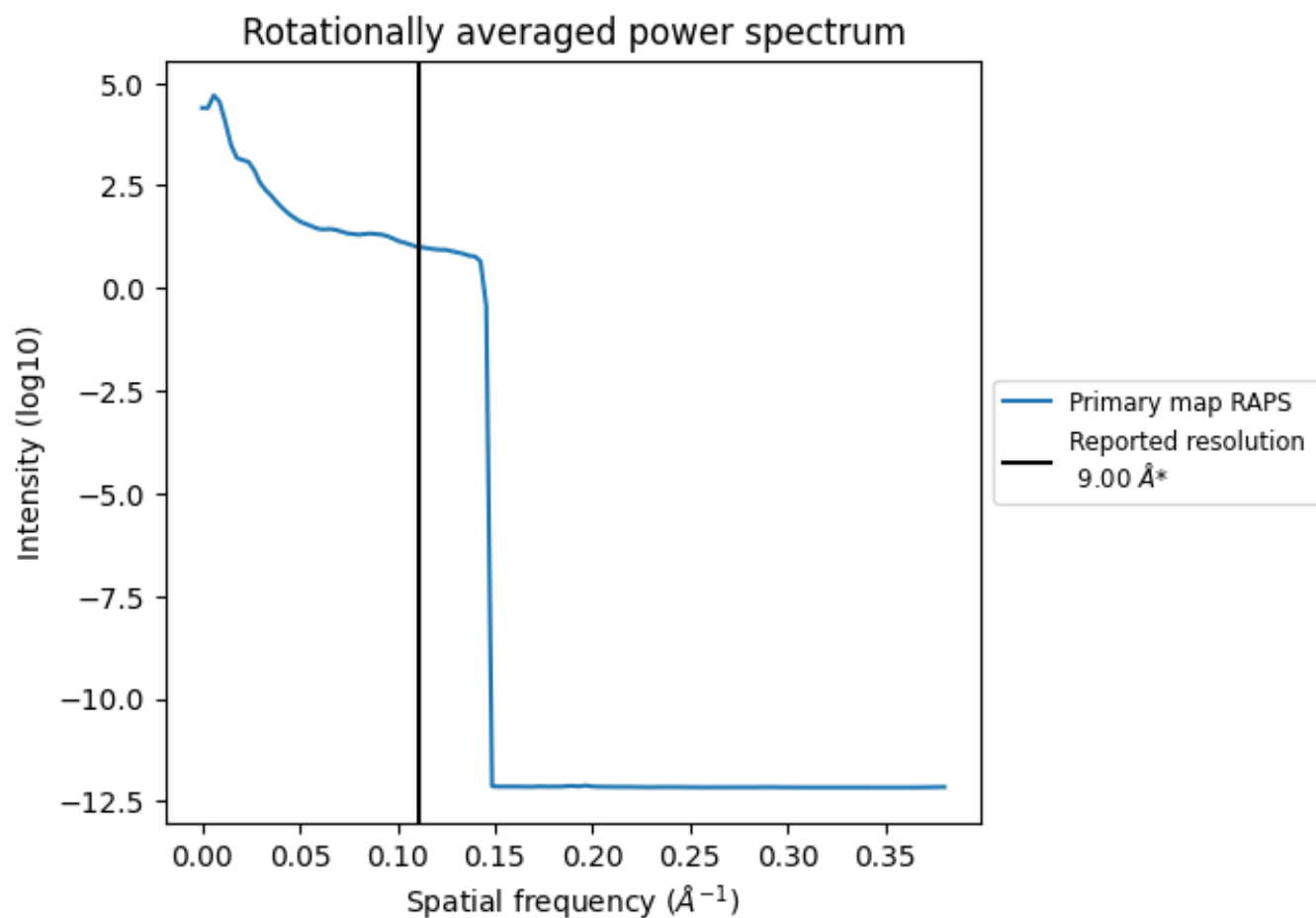
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 200 nm³; this corresponds to an approximate mass of 180 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 ⓘ

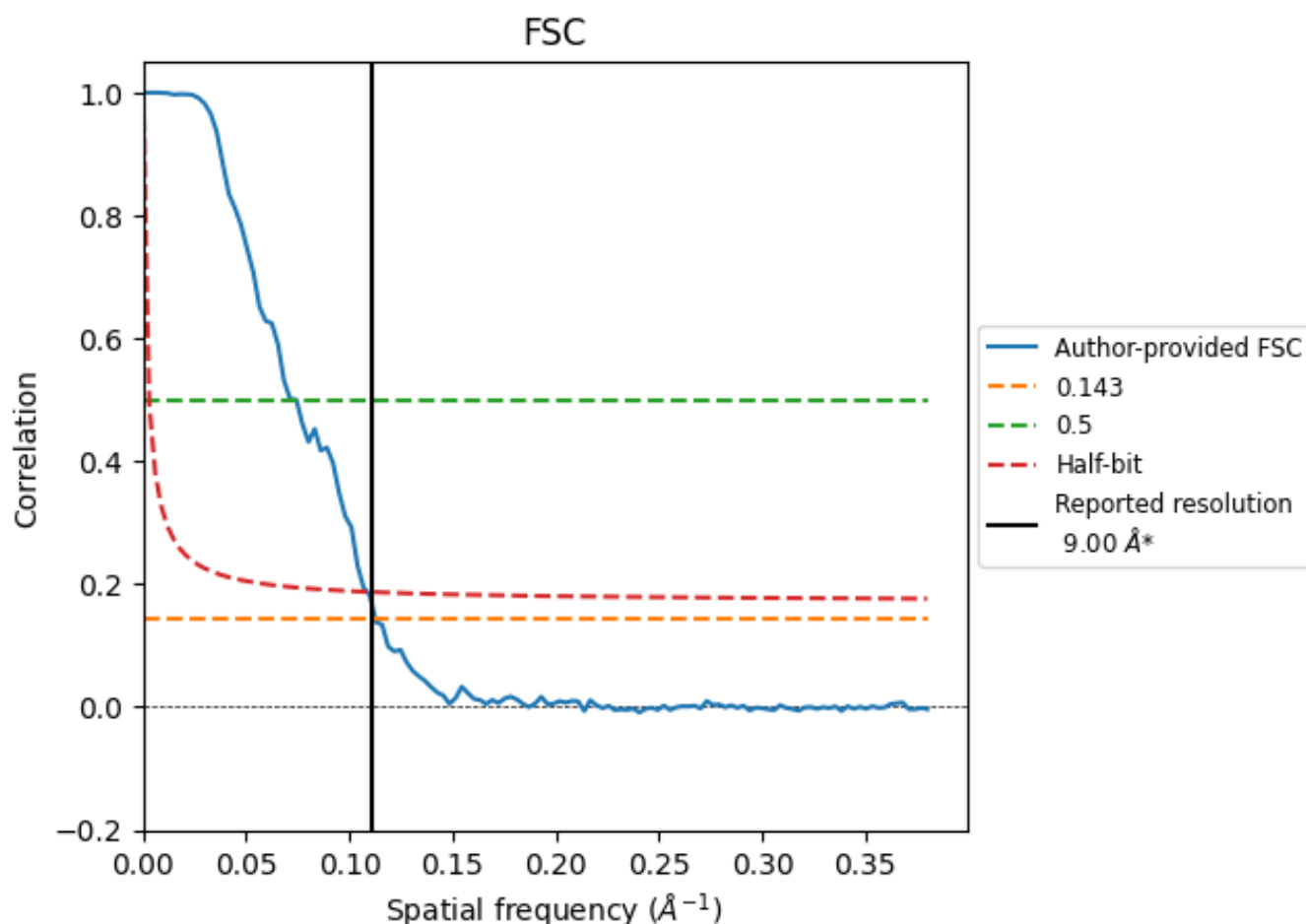


*Reported resolution corresponds to spatial frequency of 0.111 Å⁻¹

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 0.111 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	-	-	-
Author-provided FSC curve	8.89	13.55	9.21
Unmasked-calculated*	-	-	-

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

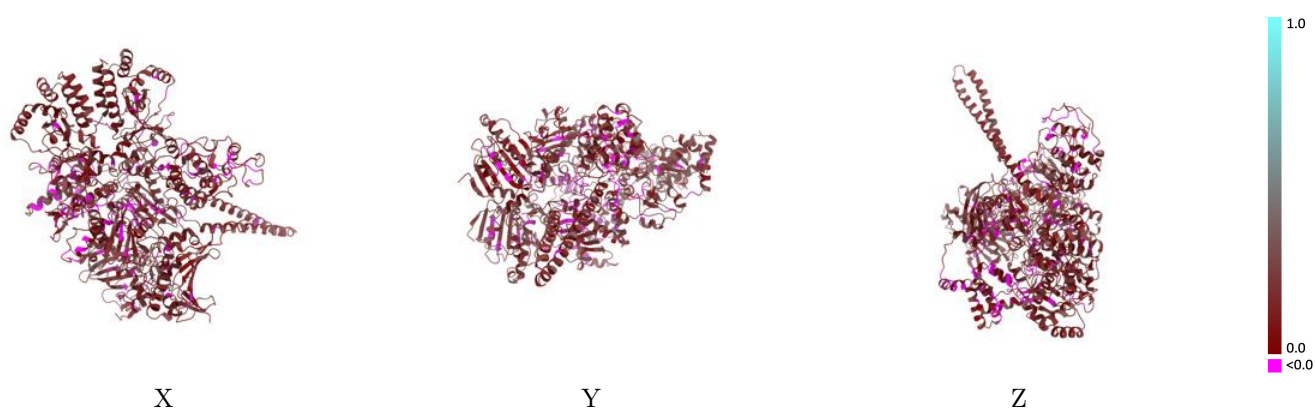
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-3341 and PDB model 5FWL. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)

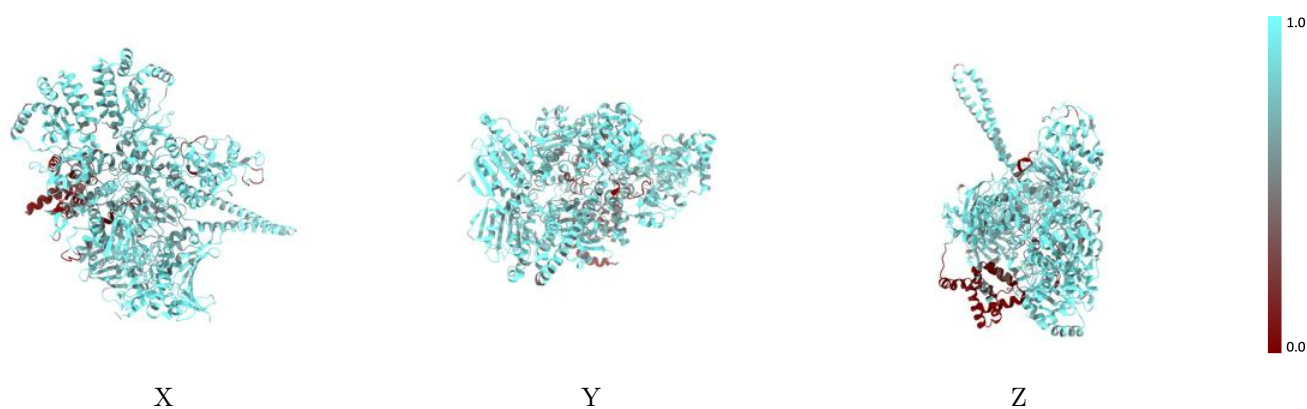
This section was not generated.

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



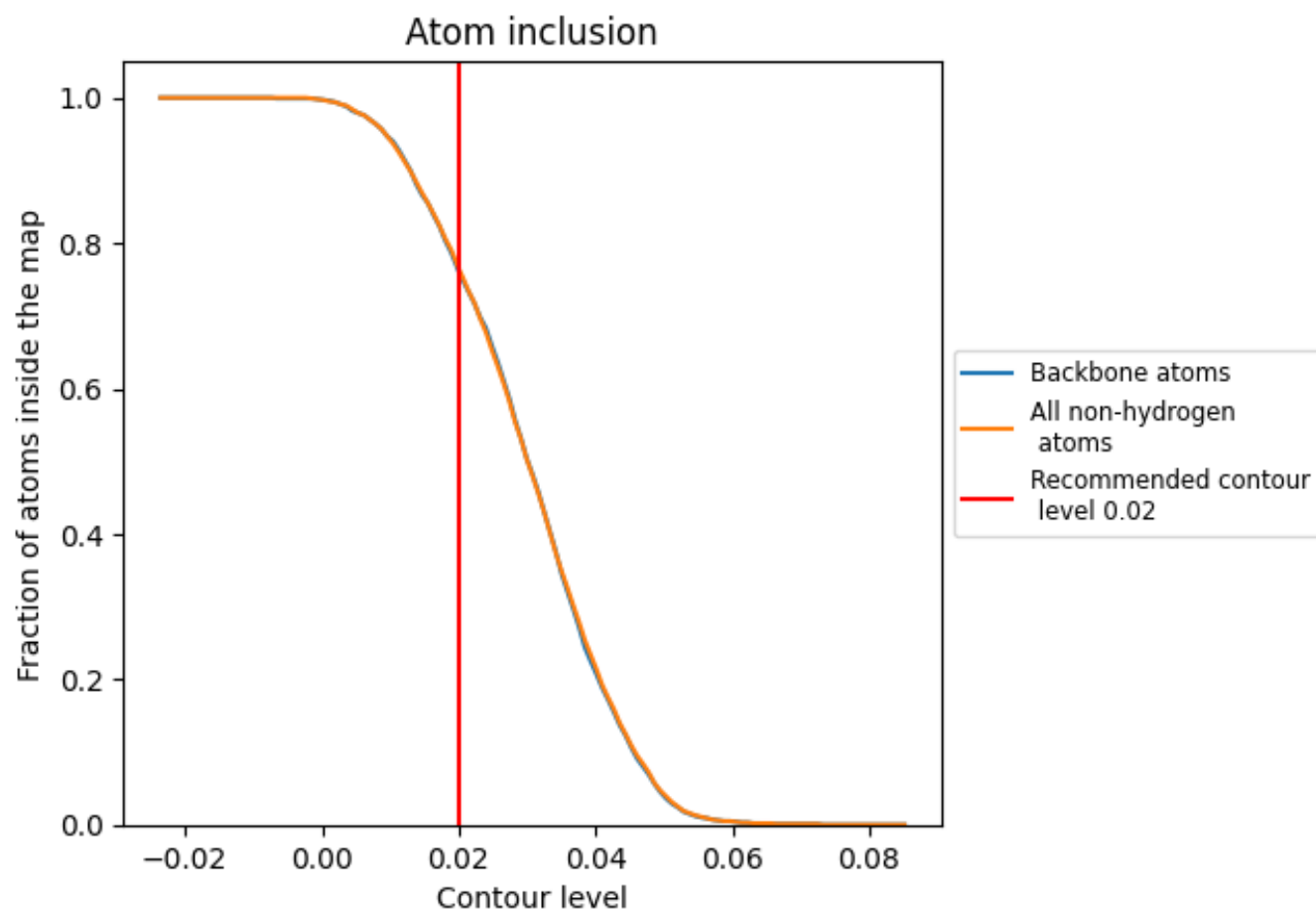
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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



At the recommended contour level, 76% 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 (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7650	0.1520
A	0.8320	0.1660
B	0.8360	0.1650
E	0.4670	0.1260
K	0.7600	0.1150

