wwPDB EM Validation Summary Report (i)

Sep 16, 2026 – 05:52 PM EDT

PDB ID	:	6PCH / pdb_00006pch
EMDB ID	:	EMD-20300
Title	:	E. coli 50S ribosome bound to compound 21
Authors	:	Pellegrino, J.; Lee, D.J.; Fraser, J.S.; Seiple, I.B.
Deposited on	:	2019-06-17
Resolution	:	2.90 Å(reported)

This is a wwPDB EM Validation Summary 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 (i) 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 (i)) were used in the production of this report:

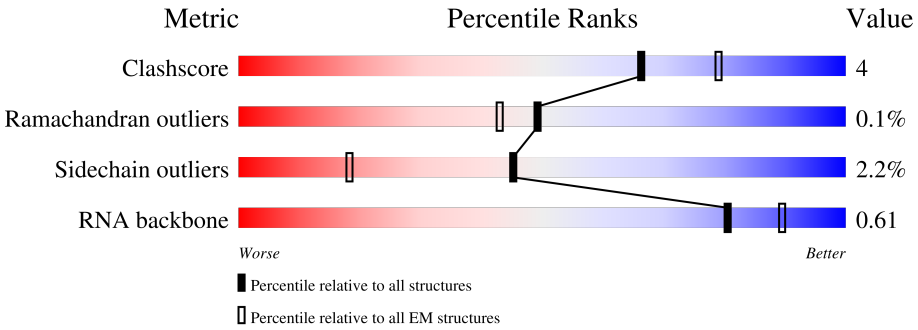
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4-5-2 with Phenix2.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data 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)	:	2.50

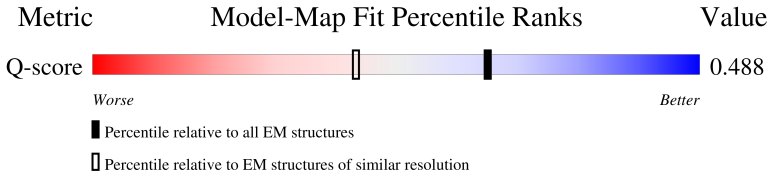
1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.90 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	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	13054 (2.40 - 3.40)

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 <=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	I	2904	5% 75% 21%
2	J	118	86% 13%
3	K	271	78% 20%

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Mol	Chain	Length	Quality of chain
4	L	144	
5	M	201	
6	N	209	
7	O	142	

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 72134 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 RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	I	2897	62206	27757	11443	20109	2897	0	0

- Molecule 2 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	J	118	2529	1126	464	821	118	0	0

- Molecule 3 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	K	271	2083	1288	423	365	7	0	0

- Molecule 4 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	L	144	1053	654	207	190	2	0	0

- Molecule 5 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	M	201	1552	974	283	290	5	0	0

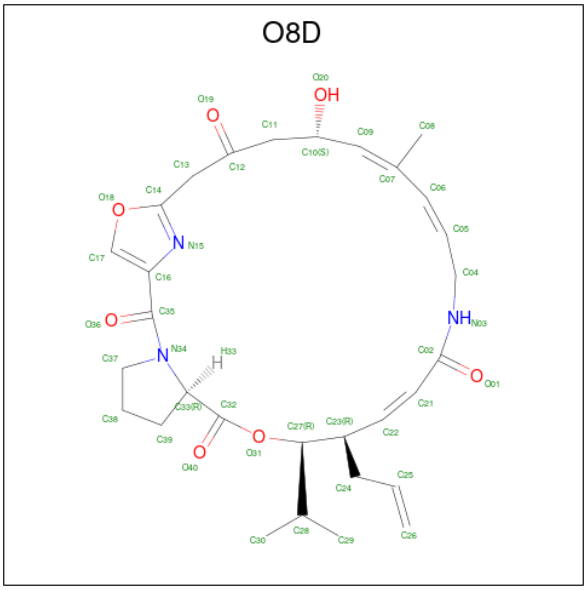
- Molecule 6 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	N	206	1542	965	284	289	4	0	0

- Molecule 7 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	O	142	1129	714	212	199	4	0	0

- Molecule 8 is (3R,4R,5E,10E,12E,14S,26aR)-14-hydroxy-12-methyl-3-(propan-2-yl)-4-(prop-2-en-1-yl)-8,9,14,15,24,25,26,26a-octahydro-1H,3H,22H-21,18-(azeno)pyrrolo[2,1-c][1,8,4,19]dioxadiazacyclotetracosine-1,7,16,22(4H,17H)-tetrone (CCD ID: O8D) (formula: $C_{30}H_{39}N_3O_7$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
8	I	1	40	30	3	7	0

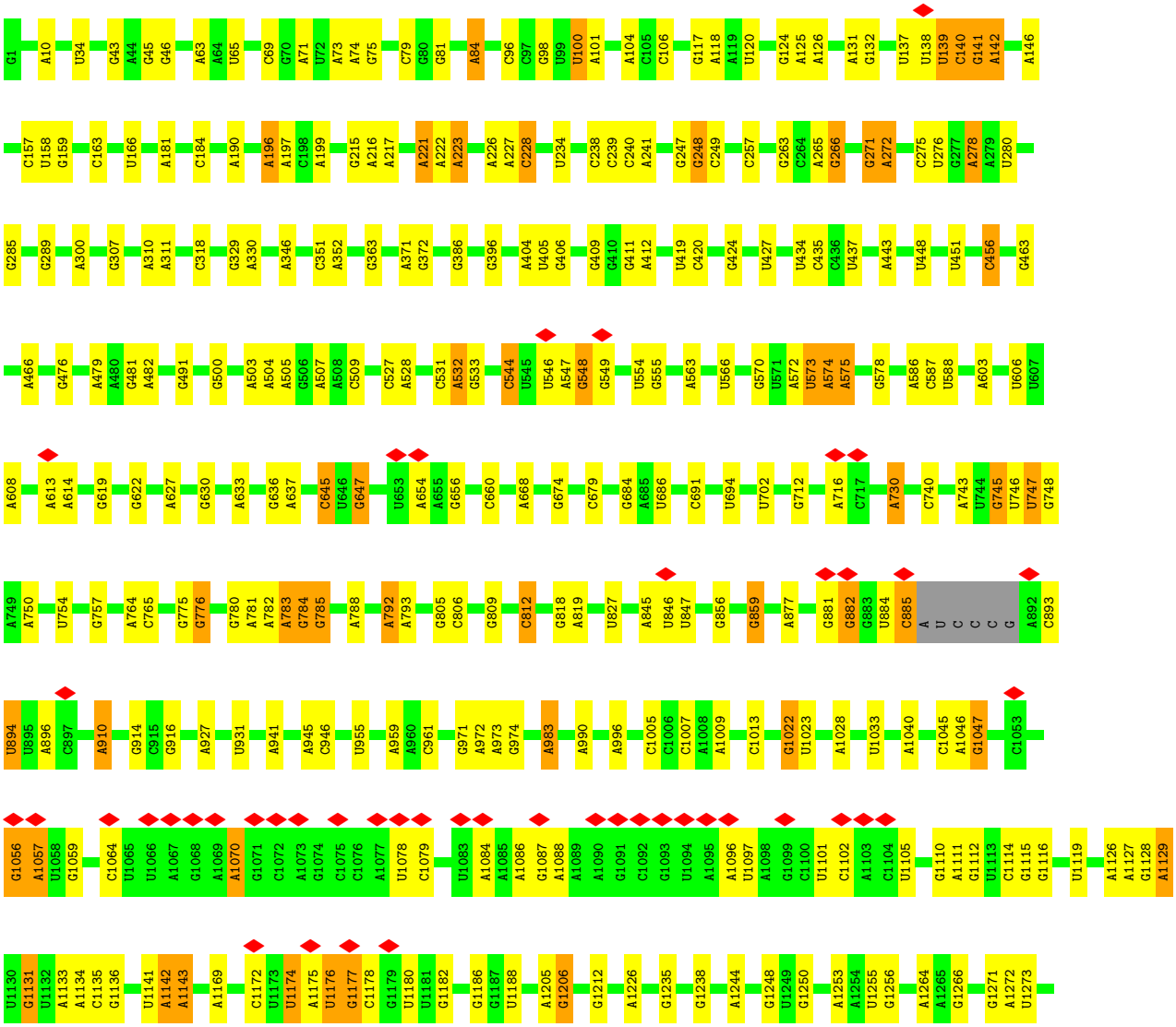
3 Residue-property plots (i)

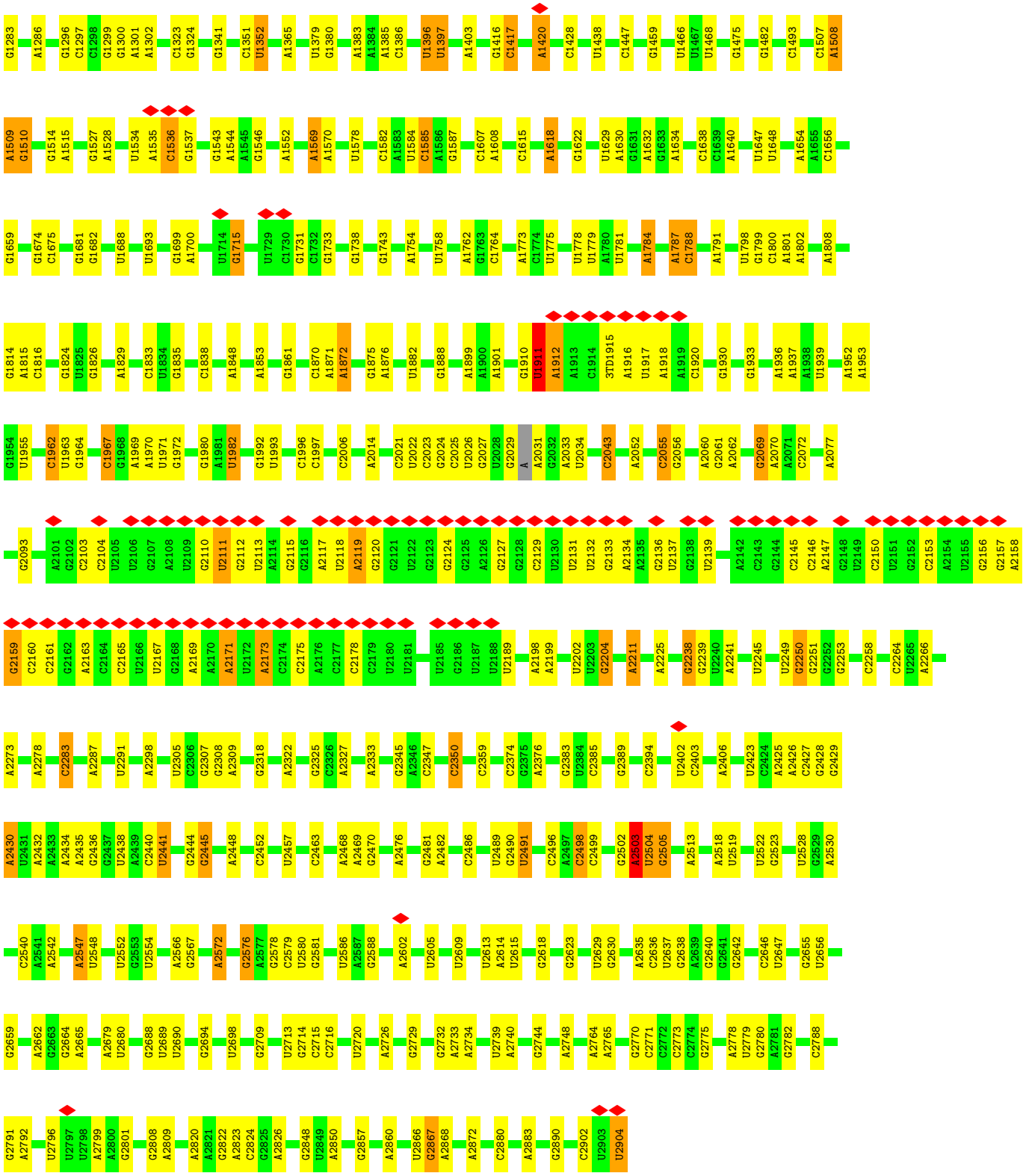
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: 23S ribosomal RNA

Chain I:



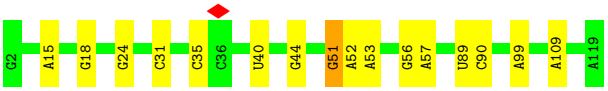




• Molecule 2: 5S ribosomal RNA

Chain J:

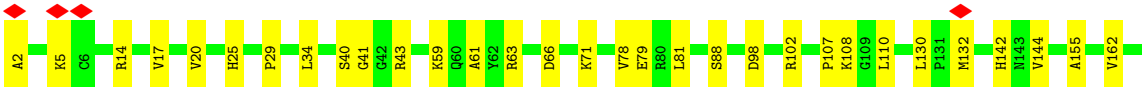


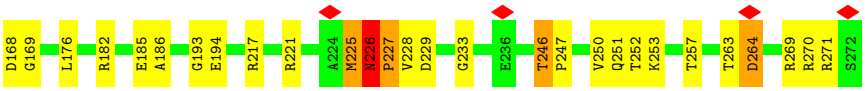


• Molecule 3: 50S ribosomal protein L2

Chain K:



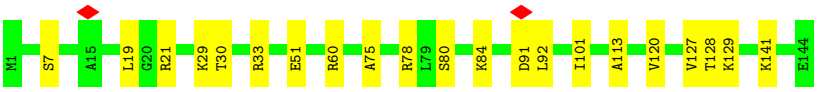




• Molecule 4: 50S ribosomal protein L15

Chain L:

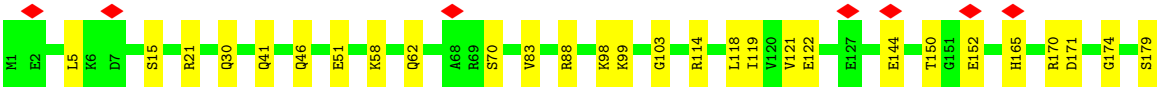


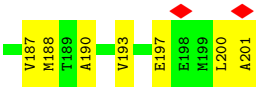


• Molecule 5: 50S ribosomal protein L4

Chain M:





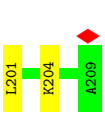


• Molecule 6: 50S ribosomal protein L3

Chain N:



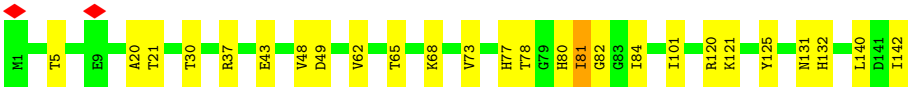




• Molecule 7: 50S ribosomal protein L13

Chain O:




M1
T5
E9
A20
T21
T30
R37
E43
V48
D49
V62
T65
K68
V73
H77
T78
G79
H80
I81
G82
G83
I84
I101
R120
K121
Y125
N131
H132
L140
D141
I142

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	23048	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	75.2	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	27.817	Depositor
Minimum map value	-17.460	Depositor
Average map value	-0.017	Depositor
Map value standard deviation	1.289	Depositor
Recommended contour level	4	Depositor
Map size (Å)	496.98, 496.98, 496.98	wwPDB
Map dimensions	600, 600, 600	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8283, 0.8283, 0.8283	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 6MZ, OMU, 5MU, 1MG, G7M, 5MC, 2MA, OMG, OMC, 2MG, O8D, PSU, 3TD

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	I	0.10	0/69165	0.16	0/107896
2	J	0.08	0/2828	0.14	0/4410
3	K	0.38	3/2122 (0.1%)	0.41	1/2852 (0.0%)
4	L	0.15	0/1062	0.34	0/1413
5	M	0.13	0/1571	0.29	0/2113
6	N	0.15	0/1561	0.28	0/2097
7	O	0.15	0/1152	0.27	0/1551
All	All	0.12	3/79461 (0.0%)	0.18	1/122332 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	K	227	PRO	C-O	-6.81	1.15	1.24
3	K	226	ASN	CA-C	-6.21	1.45	1.52
3	K	228	VAL	C-O	-5.37	1.17	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	K	225	MET	O-C-N	5.51	128.82	121.99

There are no chirality outliers.

There are no planarity outliers.

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	I	62206	0	31305	266	0
2	J	2529	0	1281	4	0
3	K	2083	0	2154	44	0
4	L	1053	0	1129	15	0
5	M	1552	0	1618	23	0
6	N	1542	0	1593	19	0
7	O	1129	0	1162	20	0
8	I	40	0	0	0	0
All	All	72134	0	40242	346	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.

The worst 5 of 346 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)
3:K:225:MET:O	3:K:226:ASN:O	1.70	1.07
1:I:1798:U:OP2	3:K:271:ARG:NH2	2.06	0.89
1:I:2576:G:O2'	1:I:2579:C:OP2	1.91	0.87
1:I:1788:C:OP2	3:K:221:ARG:NH1	2.08	0.86
1:I:1174:U:OP1	1:I:1177:G:N2	2.08	0.85

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
3	K	269/271 (99%)	234 (87%)	34 (13%)	1 (0%)	30 59
4	L	142/144 (99%)	131 (92%)	11 (8%)	0	100 100
5	M	199/201 (99%)	187 (94%)	12 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
6	N	202/209 (97%)	187 (93%)	15 (7%)	0	100 100
7	O	140/142 (99%)	137 (98%)	3 (2%)	0	100 100
All	All	952/967 (98%)	876 (92%)	75 (8%)	1 (0%)	49 77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	K	226	ASN

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
3	K	216/216 (100%)	209 (97%)	7 (3%)	34 68
4	L	103/103 (100%)	103 (100%)	0	100 100
5	M	165/165 (100%)	162 (98%)	3 (2%)	51 80
6	N	161/164 (98%)	157 (98%)	4 (2%)	42 74
7	O	116/116 (100%)	113 (97%)	3 (3%)	40 73
All	All	761/764 (100%)	744 (98%)	17 (2%)	45 77

5 of 17 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
7	O	5	THR
7	O	81	ILE
5	M	144	GLU
5	M	150	THR
5	M	200	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
6	N	126	ASN
7	O	128	ASN
7	O	138	GLN
7	O	135	GLN
3	K	243	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	I	2894/2904 (99%)	375 (12%)	9 (0%)
2	J	117/118 (99%)	9 (7%)	0
All	All	3011/3022 (99%)	384 (12%)	9 (0%)

5 of 384 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	I	10	A
1	I	34	U
1	I	43	G
1	I	46	G
1	I	63	A

5 of 9 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	I	2425	A
1	I	2481	G
1	I	1056	G
1	I	1787	A
1	I	1875	G

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

21 non-standard protein/DNA/RNA residues 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	$\#\|Z\|>2$
1	PSU	I	1917	1	18,21,22	0.86	0	21,30,33	1.97	4 (19%)
1	2MA	I	2503	1	22,25,26	1.35	3 (13%)	32,37,40	1.95	7 (21%)
1	PSU	I	1911	1	18,21,22	0.81	0	21,30,33	1.97	4 (19%)
1	PSU	I	746	1	18,21,22	0.97	1 (5%)	21,30,33	1.84	4 (19%)
1	PSU	I	955	1	18,21,22	0.91	0	21,30,33	1.96	4 (19%)
1	2MG	I	2445	1	23,26,27	0.78	0	33,38,41	2.14	8 (24%)
1	PSU	I	2457	1	18,21,22	0.91	0	21,30,33	1.94	4 (19%)
1	OMC	I	2498	1	19,22,23	0.87	1 (5%)	25,31,34	1.00	1 (4%)
1	5MC	I	1962	1	19,22,23	1.23	2 (10%)	26,32,35	1.12	3 (11%)
1	3TD	I	1915	1	19,22,23	1.07	2 (10%)	23,32,35	1.81	3 (13%)
1	PSU	I	2580	1	18,21,22	0.90	0	21,30,33	1.94	4 (19%)
1	5MU	I	747	1	19,22,23	1.02	3 (15%)	27,32,35	2.10	6 (22%)
1	PSU	I	2605	1	18,21,22	0.88	0	21,30,33	1.96	4 (19%)
1	G7M	I	2069	1	23,26,27	0.51	0	34,39,42	0.91	1 (2%)
1	OMU	I	2552	1	19,22,23	0.93	1 (5%)	25,31,34	1.93	6 (24%)
1	6MZ	I	1618	1	22,25,26	1.09	1 (4%)	29,36,39	2.17	9 (31%)
1	PSU	I	2504	1	18,21,22	0.90	0	21,30,33	1.95	4 (19%)
1	2MG	I	1835	1	23,26,27	0.75	0	33,38,41	2.16	9 (27%)
1	1MG	I	745	1	23,26,27	0.91	1 (4%)	33,39,42	1.73	6 (18%)
1	5MU	I	1939	1	19,22,23	1.02	2 (10%)	27,32,35	2.09	6 (22%)
1	OMG	I	2251	1	23,26,27	0.71	0	32,38,41	1.94	9 (28%)

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
1	PSU	I	1917	1	-	0/7/25/26	0/2/2/2
1	2MA	I	2503	1	-	4/7/25/26	0/3/3/3
1	PSU	I	1911	1	-	2/7/25/26	0/2/2/2
1	PSU	I	746	1	-	1/7/25/26	0/2/2/2
1	PSU	I	955	1	-	0/7/25/26	0/2/2/2
1	2MG	I	2445	1	-	2/9/27/28	0/3/3/3
1	PSU	I	2457	1	-	1/7/25/26	0/2/2/2
1	OMC	I	2498	1	-	2/9/27/28	0/2/2/2
1	5MC	I	1962	1	-	0/7/25/26	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	3TD	I	1915	1	-	3/7/25/26	0/2/2/2
1	PSU	I	2580	1	-	0/7/25/26	0/2/2/2
1	5MU	I	747	1	-	0/7/25/26	0/2/2/2
1	PSU	I	2605	1	-	0/7/25/26	0/2/2/2
1	G7M	I	2069	1	-	2/7/25/26	0/3/3/3
1	OMU	I	2552	1	-	0/9/27/28	0/2/2/2
1	6MZ	I	1618	1	-	2/9/27/28	0/3/3/3
1	PSU	I	2504	1	-	2/7/25/26	0/2/2/2
1	2MG	I	1835	1	-	0/9/27/28	0/3/3/3
1	1MG	I	745	1	-	0/7/25/26	0/3/3/3
1	5MU	I	1939	1	-	0/7/25/26	0/2/2/2
1	OMG	I	2251	1	-	2/9/27/28	0/3/3/3

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	1962	5MC	C5-C4	-3.99	1.41	1.44
1	I	2503	2MA	C6-N1	3.94	1.40	1.35
1	I	1618	6MZ	C6-N6	3.86	1.38	1.34
1	I	2503	2MA	C5-N7	-2.93	1.33	1.39
1	I	2498	OMC	C2-N1	-2.58	1.34	1.40

The worst 5 of 106 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	I	1835	2MG	C2-N3-C4	7.21	121.02	112.00
1	I	2445	2MG	C2-N3-C4	7.12	120.90	112.00
1	I	1915	3TD	N1-C2-N3	6.43	120.80	116.13
1	I	2503	2MA	C5-C4-N3	-5.76	121.11	127.18
1	I	1911	PSU	N1-C2-N3	5.66	121.13	115.17

There are no chirality outliers.

5 of 23 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	I	1618	6MZ	C5-C6-N6-C9
1	I	1618	6MZ	N1-C6-N6-C9
1	I	1915	3TD	O4'-C1'-C5-C4
1	I	1915	3TD	O4'-C1'-C5-C6
1	I	2445	2MG	C3'-C4'-C5'-O5'

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	I	2503	2MA	1	0
1	I	1911	PSU	4	0
1	I	1962	5MC	1	0
1	I	1618	6MZ	1	0
1	I	745	1MG	1	0

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand 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	$\#\|Z\| > 2$
8	O8D	I	3001	-	41,42,42	1.35	5 (12%)	48,57,57	0.87	2 (4%)

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
8	O8D	I	3001	-	-	3/51/61/61	0/2/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	I	3001	O8D	C17-C16	4.61	1.39	1.34
8	I	3001	O8D	C16-C35	-3.04	1.45	1.49

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	I	3001	O8D	C16-N15	-2.73	1.32	1.38
8	I	3001	O8D	C14-N15	2.15	1.31	1.29
8	I	3001	O8D	C21-C02	-2.00	1.44	1.48

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
8	I	3001	O8D	C32-C33-N34	2.66	117.32	112.08
8	I	3001	O8D	C27-O31-C32	2.14	121.32	117.76

There are no chirality outliers.

All (3) torsion outliers are listed below:

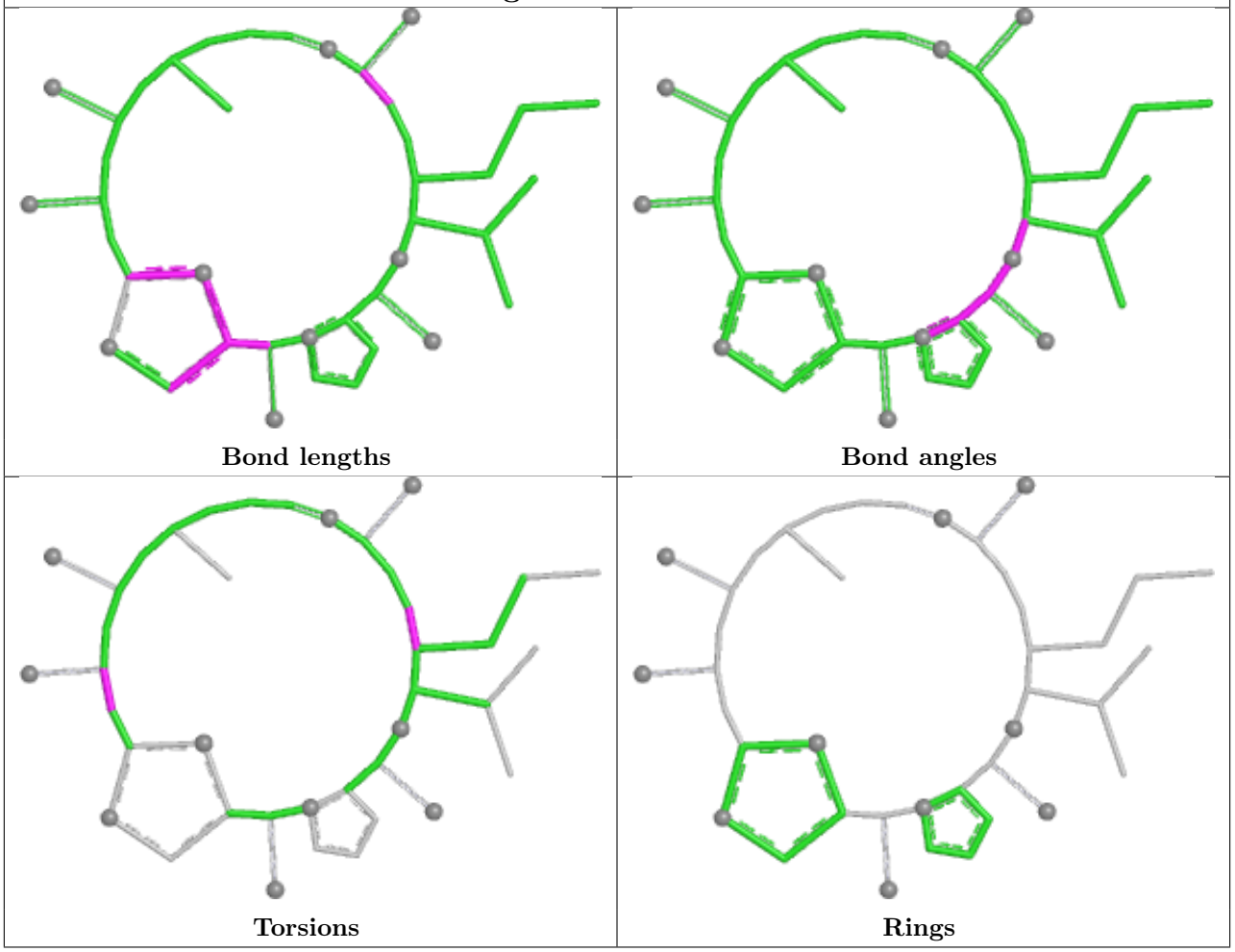
Mol	Chain	Res	Type	Atoms
8	I	3001	O8D	C21-C22-C23-C24
8	I	3001	O8D	O19-C12-C13-C14
8	I	3001	O8D	C11-C12-C13-C14

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 then 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.

Ligand O8D I 3001



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.

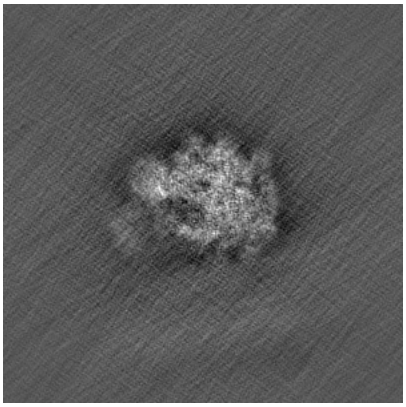
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-20300. These allow visual inspection of the internal detail of the map and identification of artifacts.

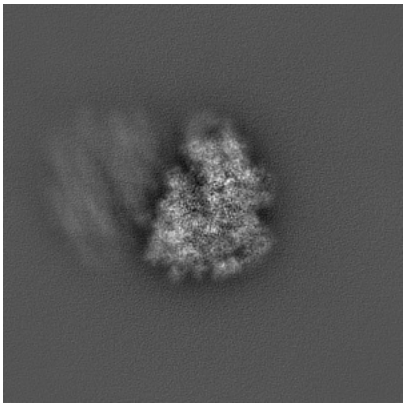
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

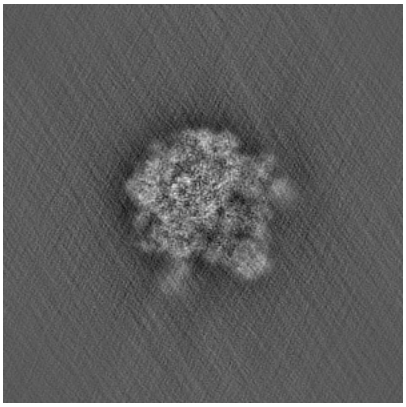
6.1.1 Primary map



X



Y

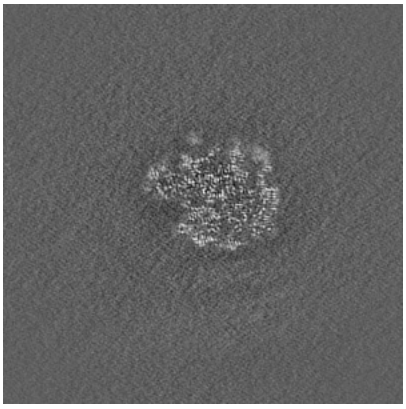


Z

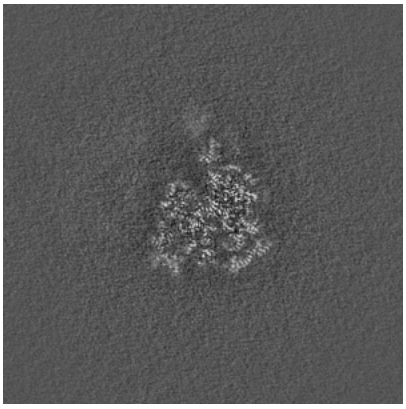
The images above show the map projected in three orthogonal directions.

6.2 Central slices

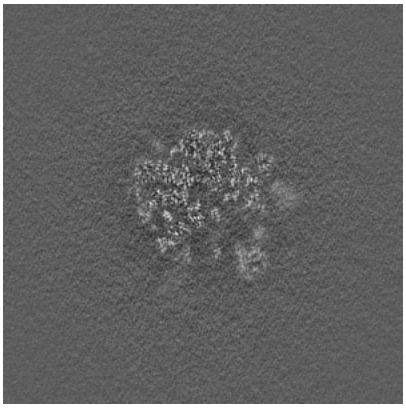
6.2.1 Primary map



X Index: 300



Y Index: 300

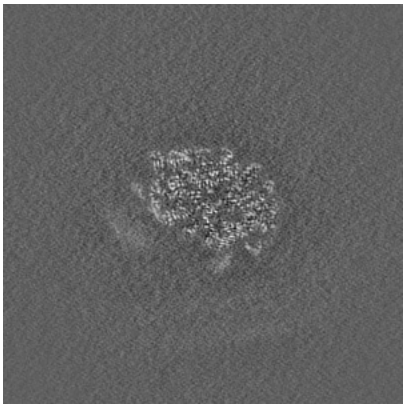


Z Index: 300

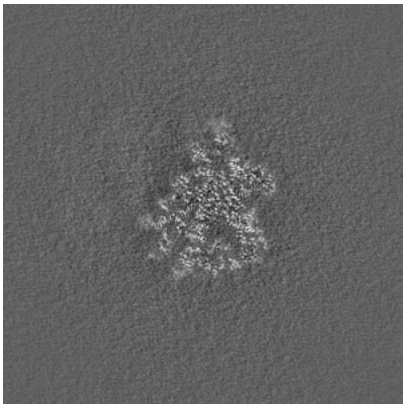
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

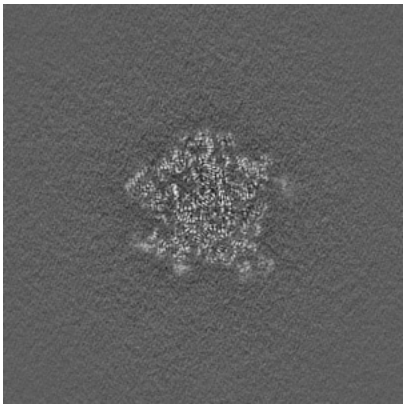
6.3.1 Primary map



X Index: 274



Y Index: 336

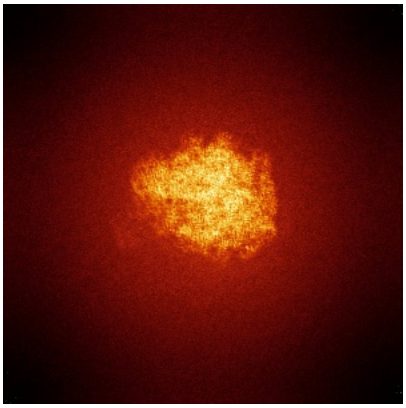


Z Index: 317

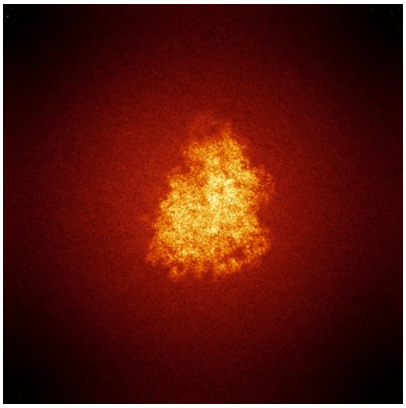
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

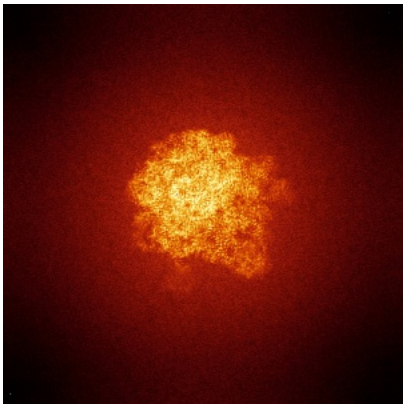
6.4.1 Primary map



X



Y

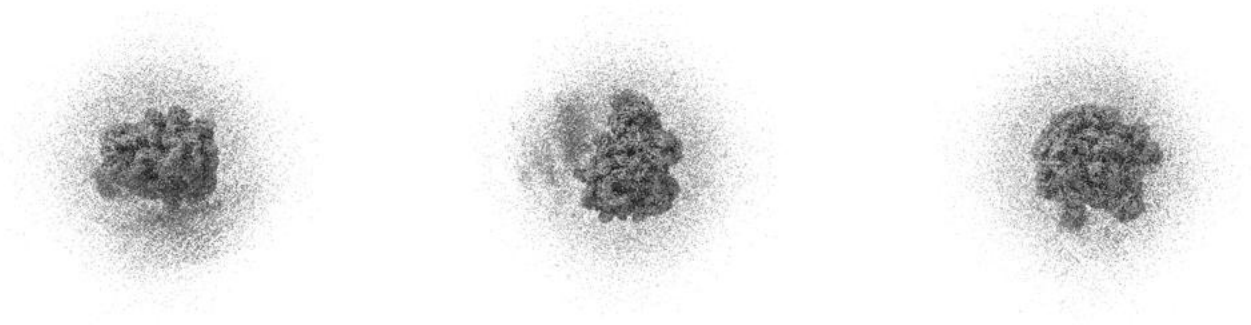


Z

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

6.5 Orthogonal surface views (i)

6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 4.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

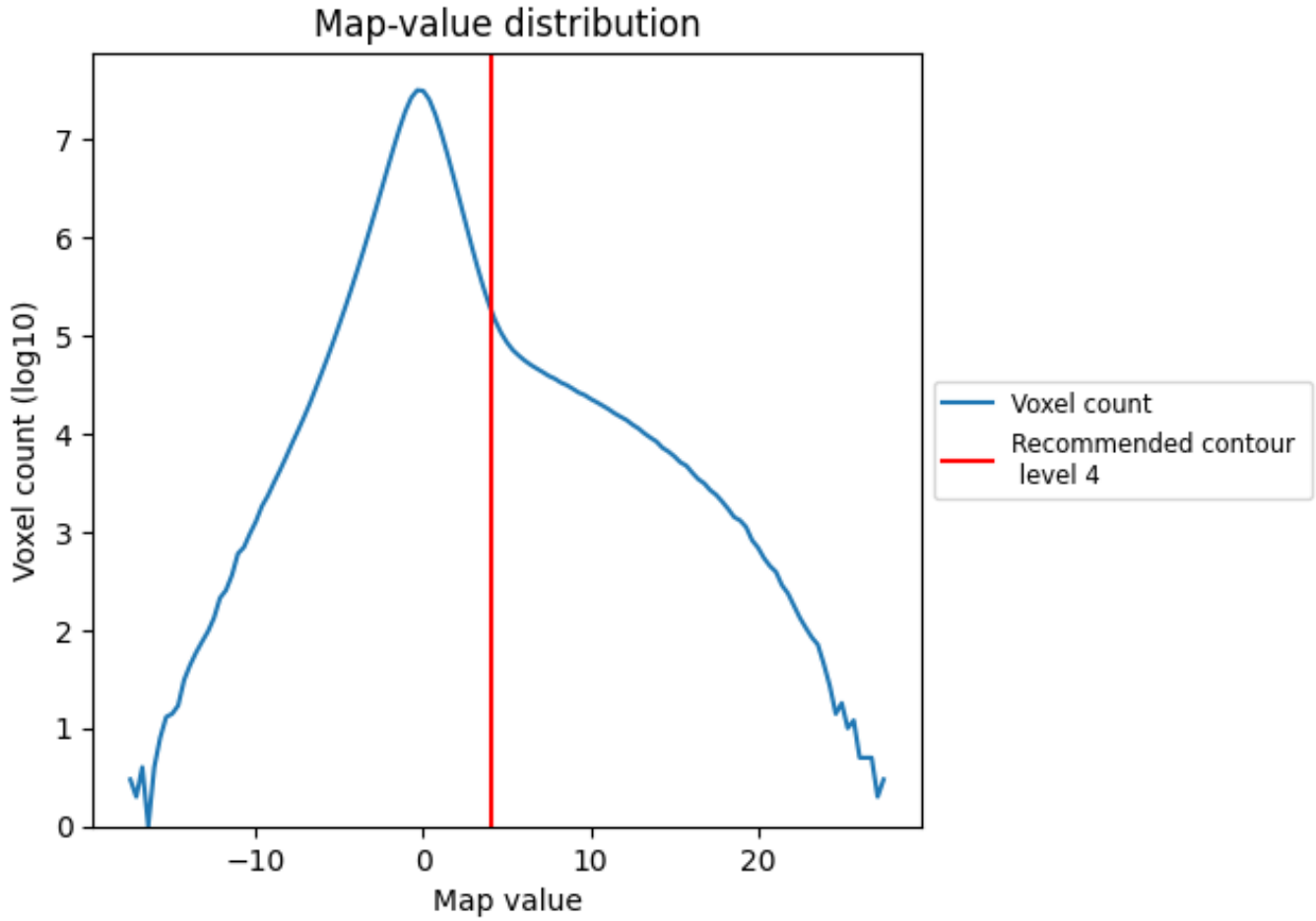
6.6 Mask visualisation (i)

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

7 Map analysis

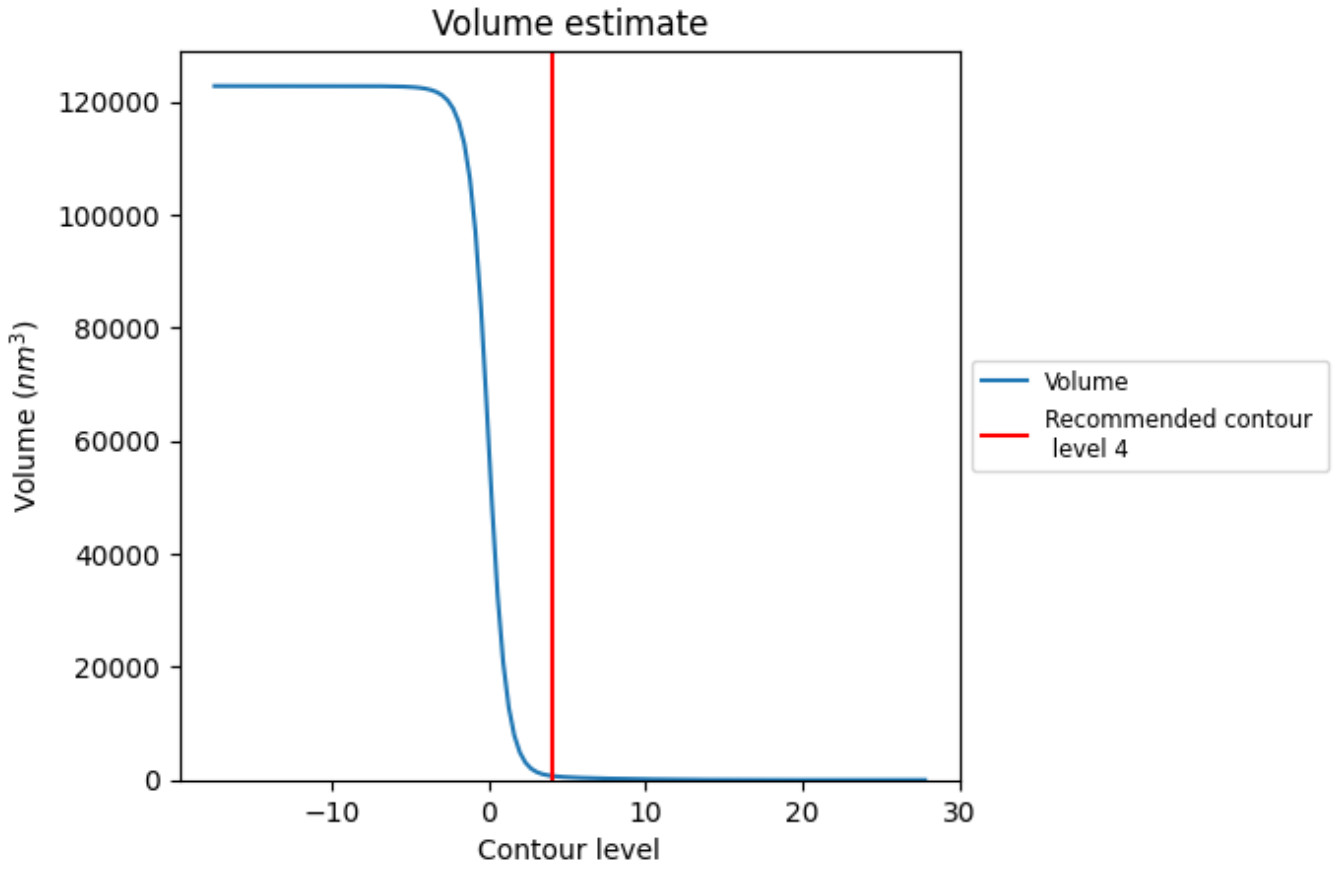
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

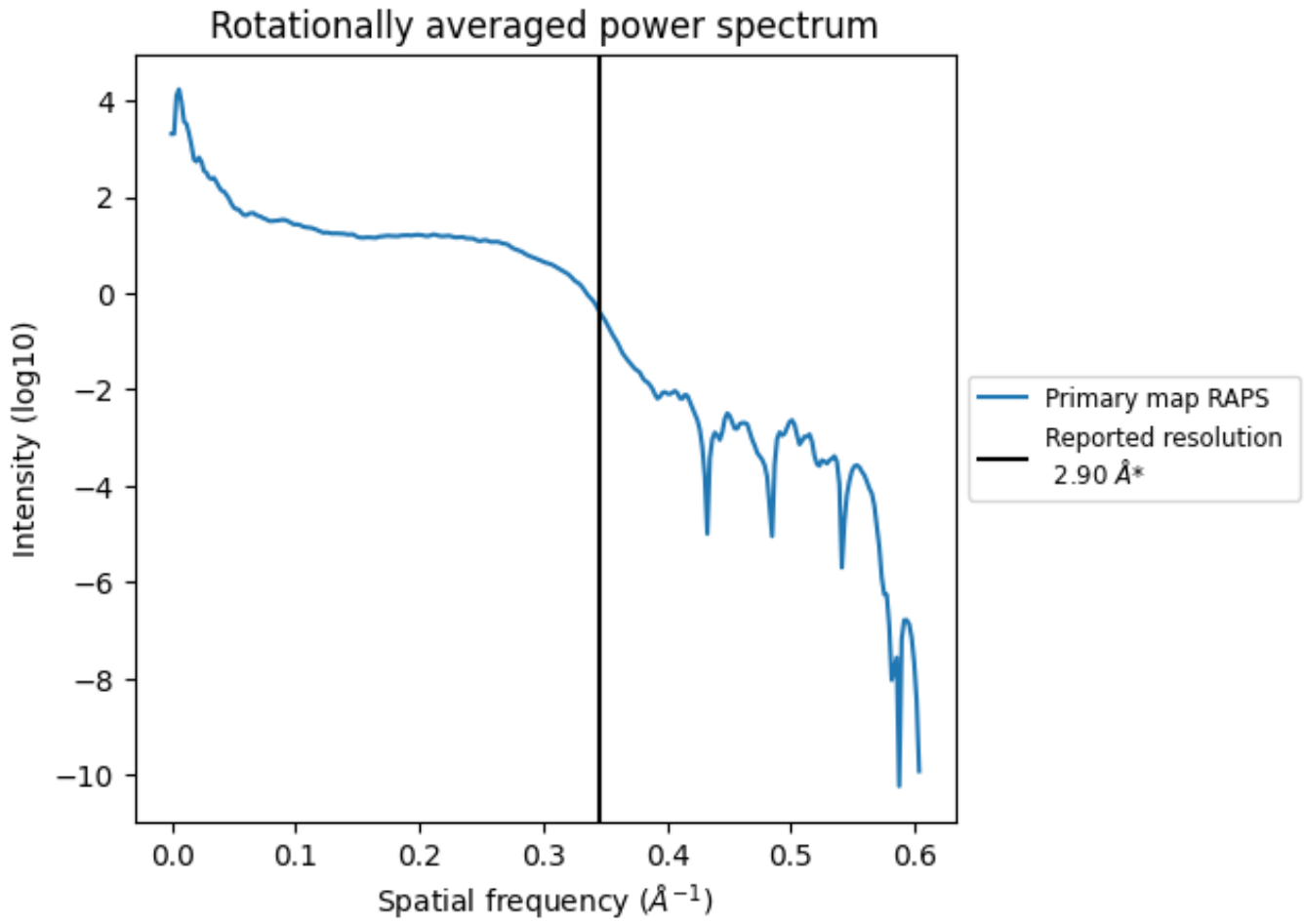
7.2 Volume estimate (i)



The volume at the recommended contour level is 731 nm^3; this corresponds to an approximate mass of 661 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 0.345 Å^{-1}

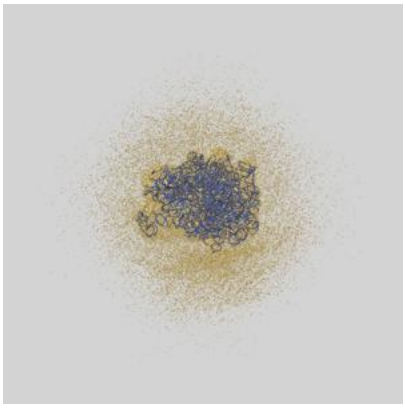
8 Fourier-Shell correlation

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

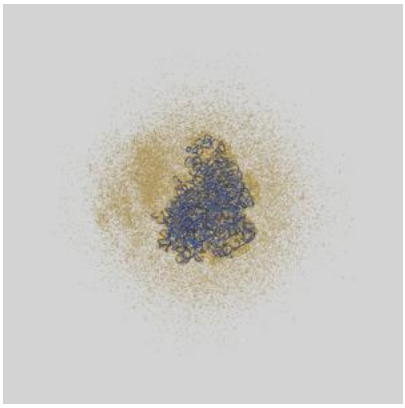
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-20300 and PDB model 6PCH. Per-residue inclusion information can be found in section 3 on page 6.

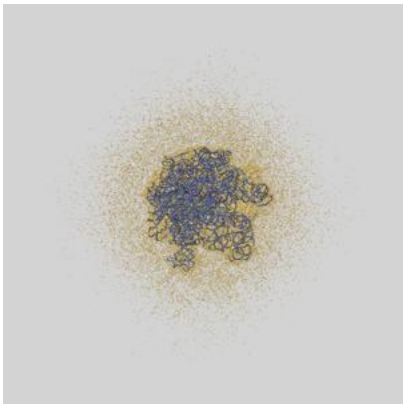
9.1 Map-model overlay



X



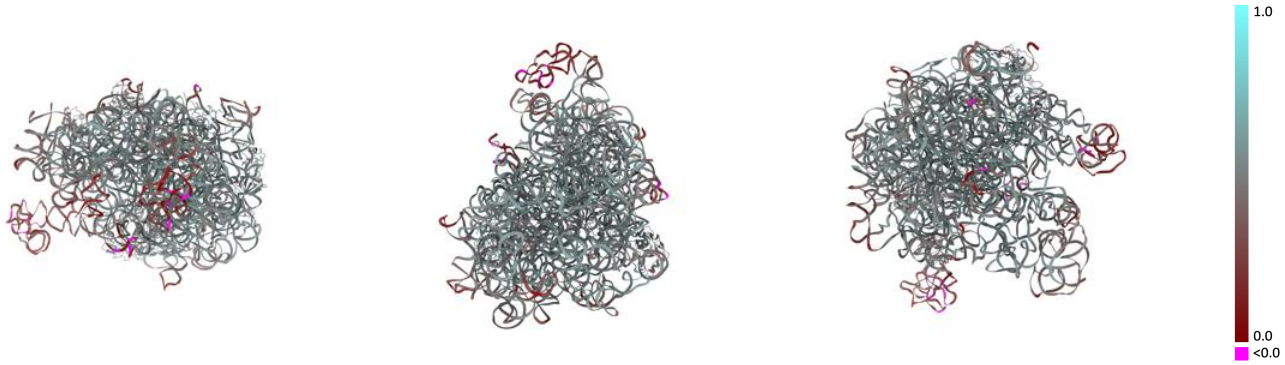
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 4.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

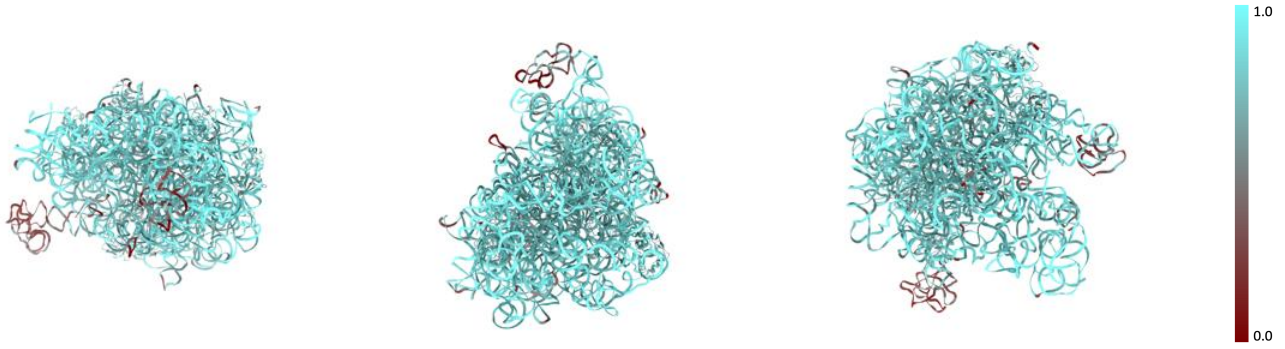
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

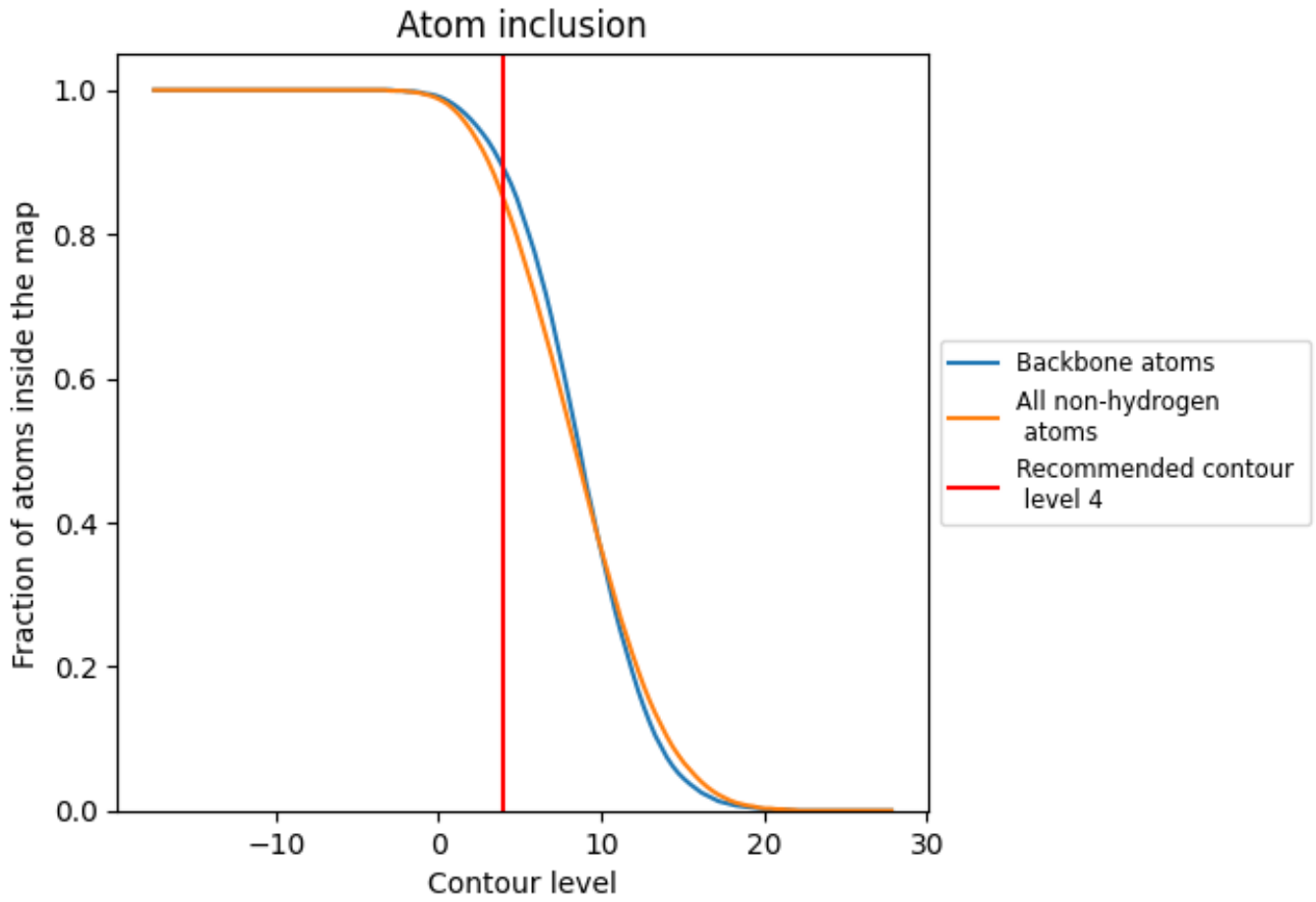
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (4).

9.4 Atom inclusion (i)



At the recommended contour level, 89% of all backbone atoms, 85% 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 (4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8490	0.4880
I	0.8570	0.4870
J	0.8810	0.4570
K	0.7760	0.5090
L	0.7540	0.4960
M	0.7470	0.4960
N	0.7730	0.5160
O	0.8020	0.5230

