



wwPDB EM Validation Summary Report ⓘ

Mar 6, 2026 – 09:38 AM UTC

PDB ID : 6PSR / pdb_00006psr
EMDB ID : EMD-20461
Title : Escherichia coli RNA polymerase promoter unwinding intermediate (TRPi1)
with TraR and rpsT P2 promoter
Authors : Chen, J.; Chiu, C.E.; Campbell, E.A.; Darst, S.A.
Deposited on : 2019-07-13
Resolution : 3.40 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 ⓘ 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:

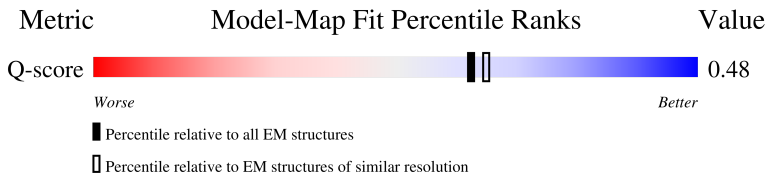
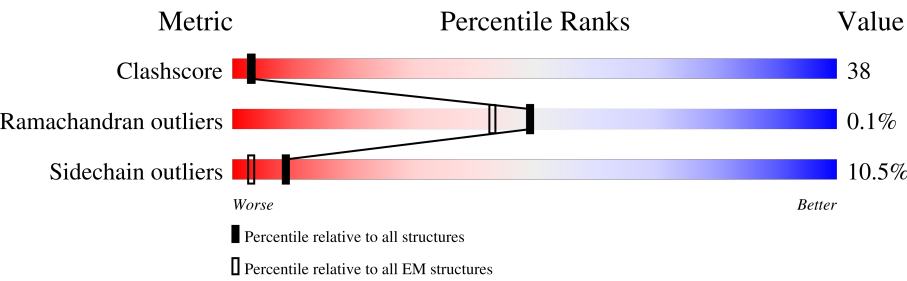
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 i

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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14717 (2.90 - 3.90)

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	G	329	34%32%5%30%
1	H	329	27%35%.33%
1	M	329	9% 5%16%.78%
2	I	1342	41%52%6%

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Mol	Chain	Length	Quality of chain
3	J	1430	10% 38% 50% 6% 6%
4	K	91	31% 46% 5% 18%
5	L	616	13% 32% 53% • 11%
6	N	72	51% 44% •
7	O	85	9% 33% 58%
8	P	85	• 39% 59%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 32342 atoms, of which 156 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 DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	G	230	Total	C	N	O	S	0	0
			1771	1106	314	345	6		
1	H	219	Total	C	N	O	S	0	0
			1678	1048	295	329	6		
1	M	73	Total	C	N	O	S	0	0
			572	362	100	108	2		

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	I	1340	Total	C	N	O	S	0	0
			10559	6627	1841	2048	43		

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	J	1345	Total	C	N	O	S	0	0
			10466	6577	1867	1972	50		

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	1	VAL	-	expression tag	UNP P0A8T7
J	1408	LEU	-	expression tag	UNP P0A8T7
J	1409	GLU	-	expression tag	UNP P0A8T7
J	1410	LEU	-	expression tag	UNP P0A8T7
J	1411	GLU	-	expression tag	UNP P0A8T7
J	1412	VAL	-	expression tag	UNP P0A8T7
J	1413	LEU	-	expression tag	UNP P0A8T7
J	1414	PHE	-	expression tag	UNP P0A8T7
J	1415	GLN	-	expression tag	UNP P0A8T7
J	1416	GLY	-	expression tag	UNP P0A8T7
J	1417	PRO	-	expression tag	UNP P0A8T7

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Chain	Residue	Modelled	Actual	Comment	Reference
J	1418	SER	-	expression tag	UNP P0A8T7
J	1419	SER	-	expression tag	UNP P0A8T7
J	1420	GLY	-	expression tag	UNP P0A8T7
J	1421	HIS	-	expression tag	UNP P0A8T7
J	1422	HIS	-	expression tag	UNP P0A8T7
J	1423	HIS	-	expression tag	UNP P0A8T7
J	1424	HIS	-	expression tag	UNP P0A8T7
J	1425	HIS	-	expression tag	UNP P0A8T7
J	1426	HIS	-	expression tag	UNP P0A8T7
J	1427	HIS	-	expression tag	UNP P0A8T7
J	1428	HIS	-	expression tag	UNP P0A8T7
J	1429	HIS	-	expression tag	UNP P0A8T7
J	1430	HIS	-	expression tag	UNP P0A8T7

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	K	75	Total	C	N	O	S	0	0
			600	365	114	120	1		

- Molecule 5 is a protein called RNA polymerase sigma factor RpoD.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	L	548	Total	C	N	O	S	0	0
			4407	2754	771	855	27		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	-2	SER	-	expression tag	UNP Q0P6L9
L	-1	GLU	-	expression tag	UNP Q0P6L9
L	0	PHE	-	expression tag	UNP Q0P6L9

- Molecule 6 is a protein called Protein TraR.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	N	72	Total	C	N	O	S	0	0
			565	350	102	108	5		

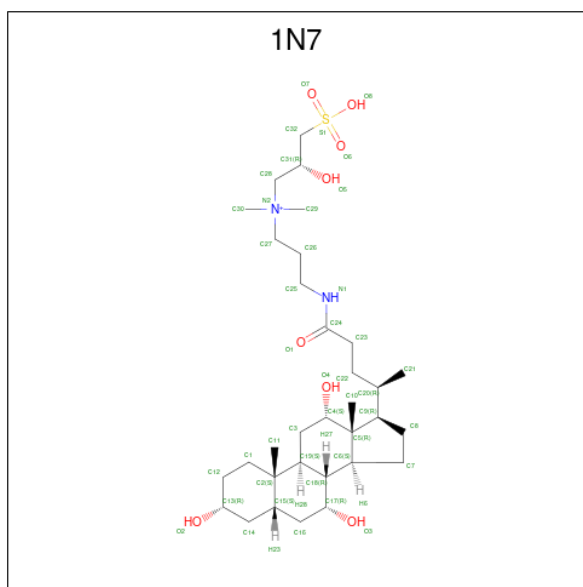
- Molecule 7 is a DNA chain called DNA (85-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
7	O	36	Total 746	C 353	N 154	O 203	P 36	0	0

- Molecule 8 is a DNA chain called DNA (85-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
8	P	35	Total 710	C 342	N 111	O 222	P 35	0	0

- Molecule 9 is CHAPSO (CCD ID: 1N7) (formula: $\text{C}_{32}\text{H}_{59}\text{N}_2\text{O}_8\text{S}$).



Mol	Chain	Residues	Atoms				AltConf
9	I	1	Total 66	C 24	H 39	O 3	0
9	J	1	Total 66	C 24	H 39	O 3	0
9	J	1	Total 66	C 24	H 39	O 3	0
9	L	1	Total 66	C 24	H 39	O 3	0

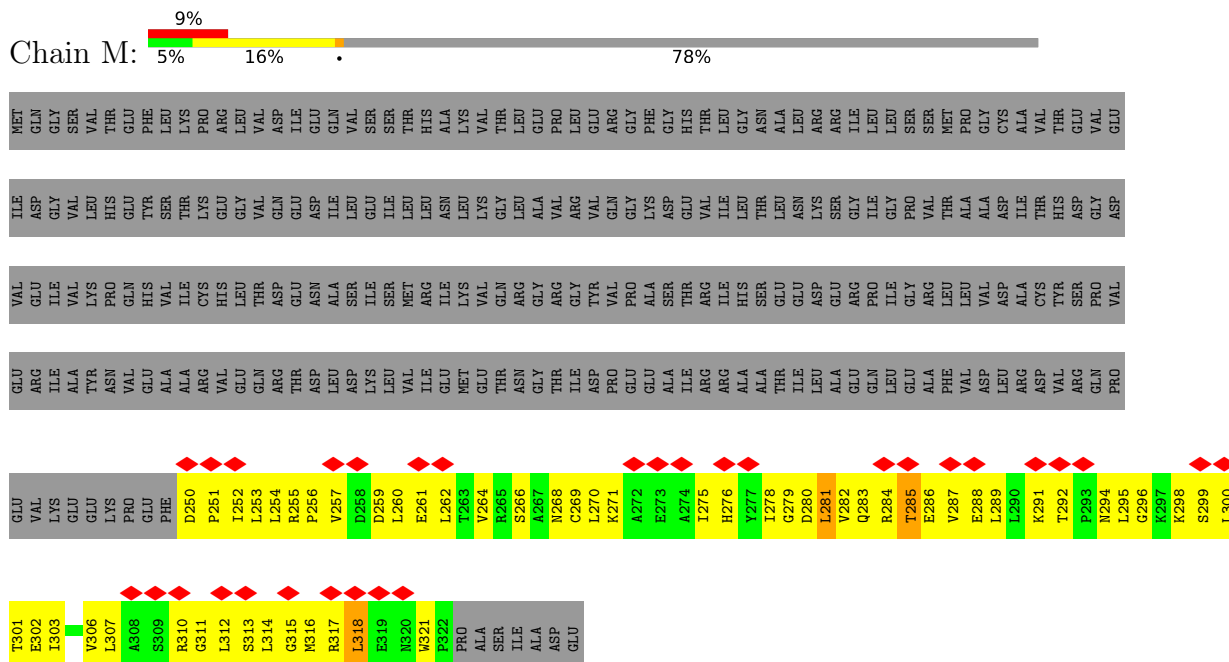
- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
10	J	1	Total	Mg	0
			1	1	

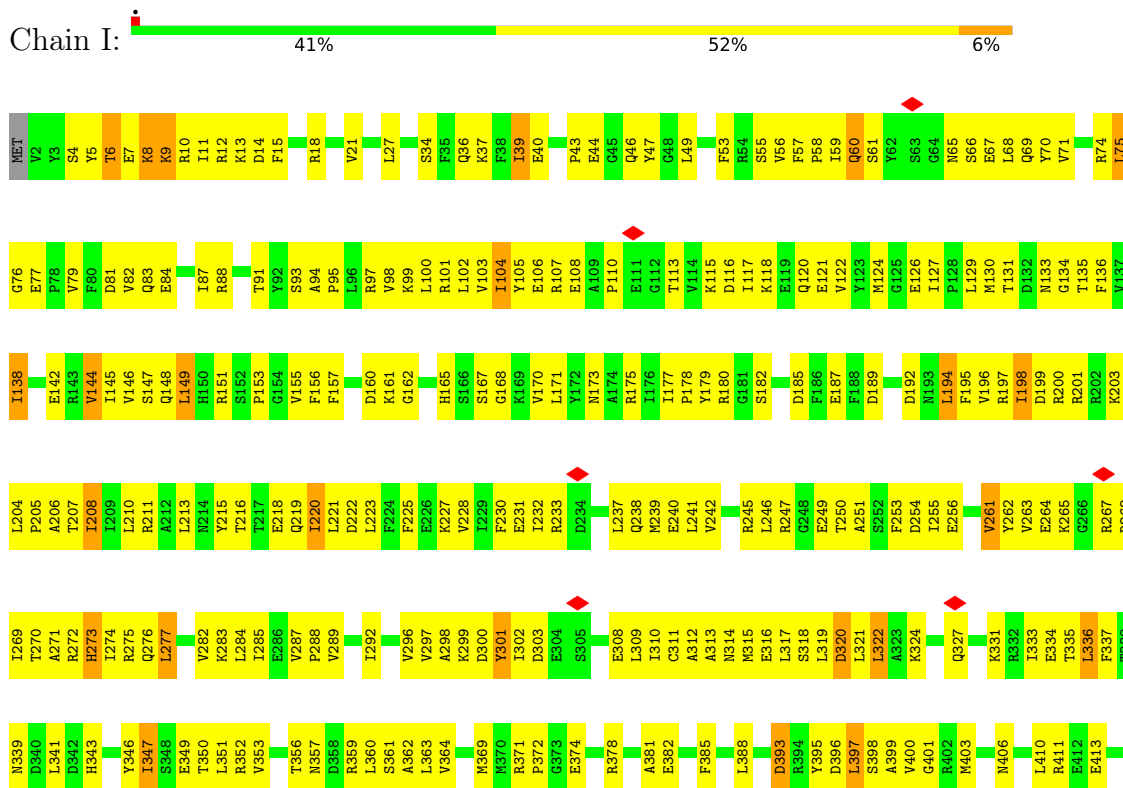
- Molecule 11 is ZINC ION (CCD ID: ZN) (formula: Zn).

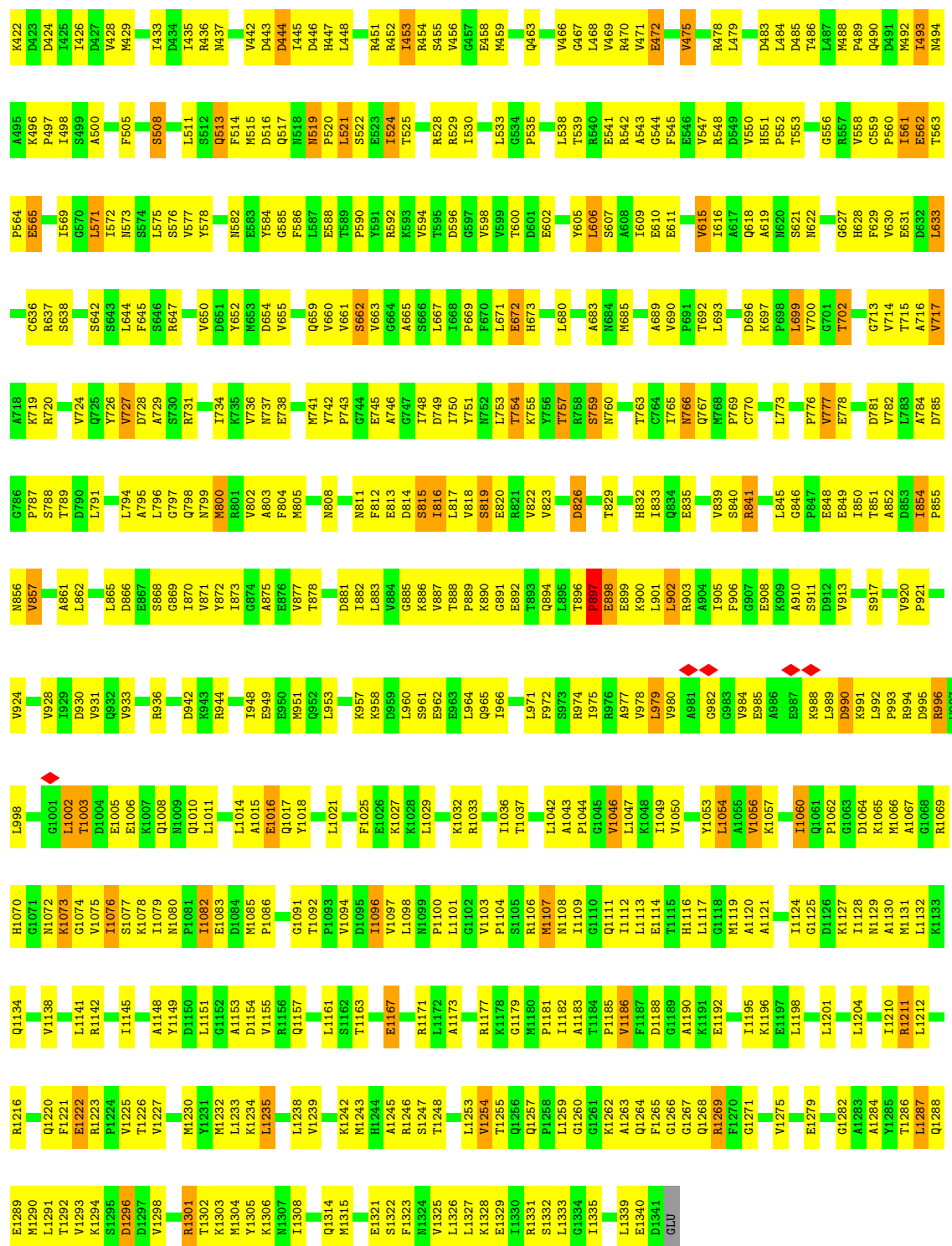
Mol	Chain	Residues	Atoms		AltConf
11	J	2	Total 2	Zn 2	0
11	N	1	Total 1	Zn 1	0

- Molecule 1: DNA-directed RNA polymerase subunit alpha



- Molecule 2: DNA-directed RNA polymerase subunit beta

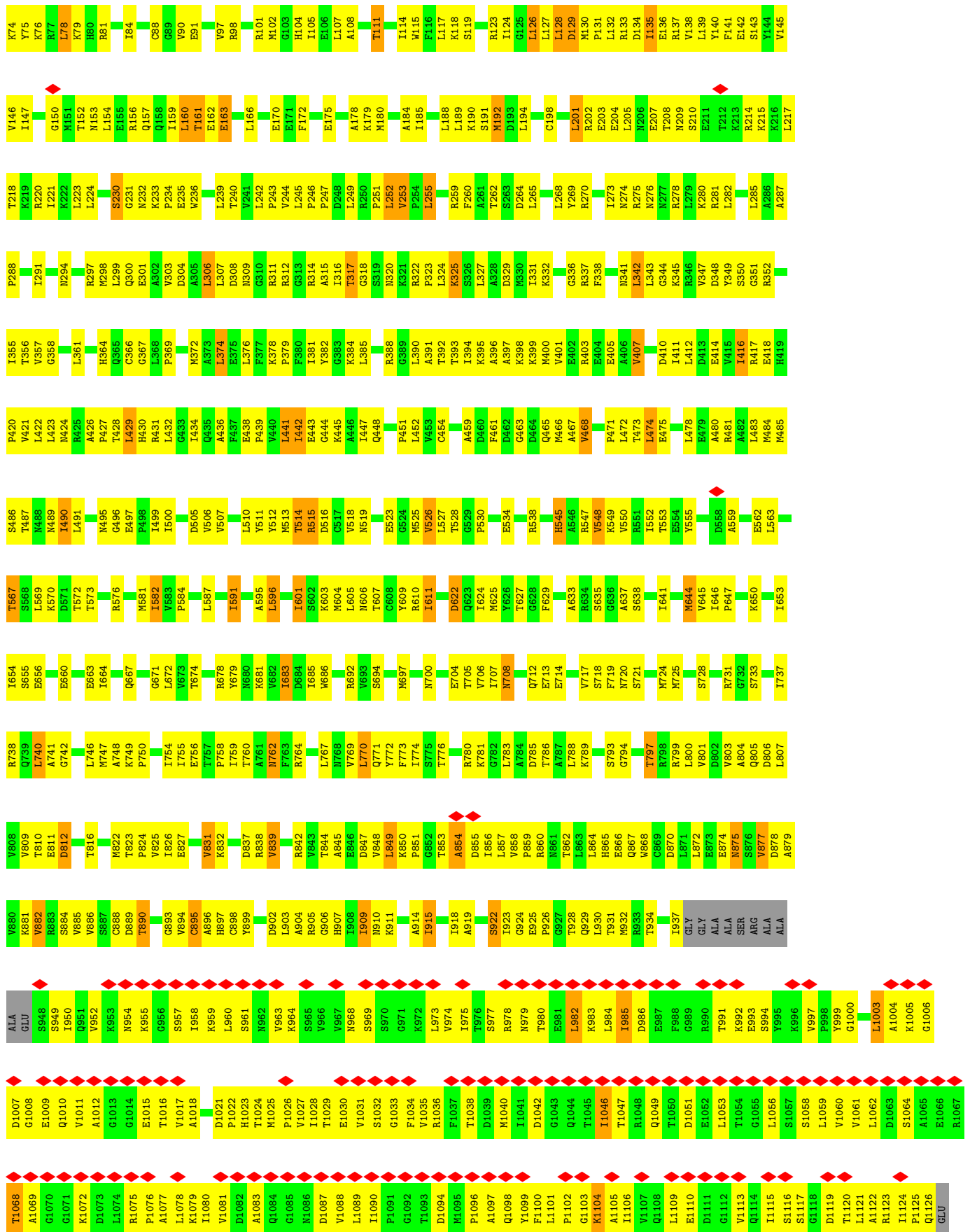


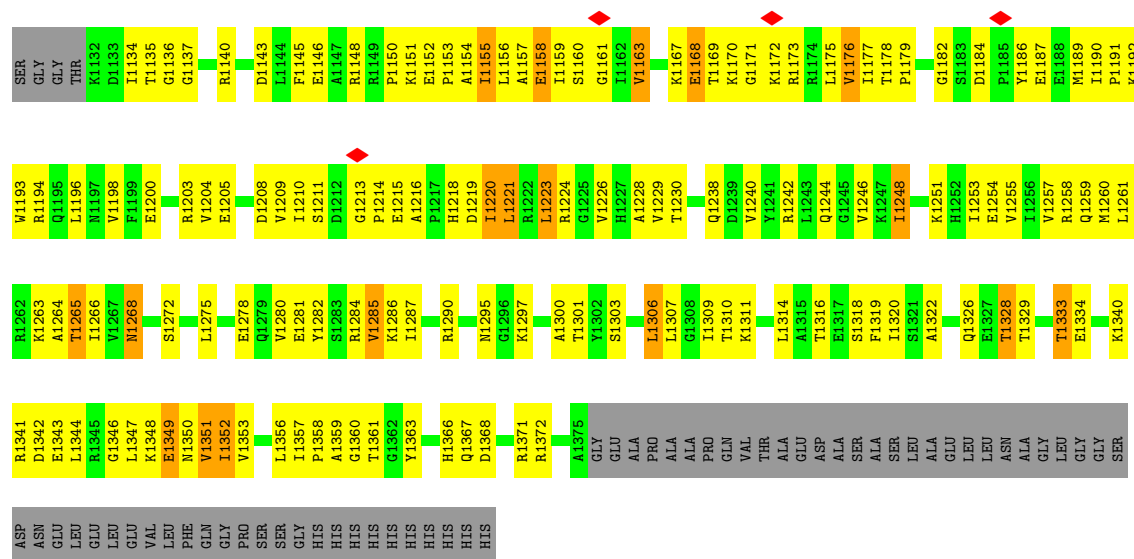


• Molecule 3: DNA-directed RNA polymerase subunit beta'



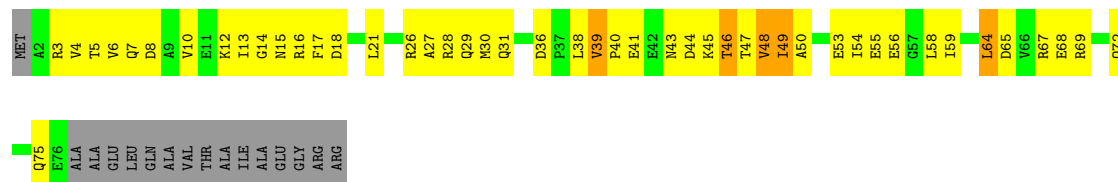
VAL	E16	A19	L20	K21	L22	A23	L24	A25	S26	M29	I30	S34	F35	K40	P41	E42	T43	L44	N45	Y46	R47	T48	R53	D54	G55	L56	F57	C58	I61	F62	V65	K66	E69	C70	L71	C72	C73														
LYS	GLU	ASP	LEU	LEU	LYS	PHE	LEU	LYS	ALA	GLN	THR	LYS	THR	GLU	E16	A19	L20	K21	L22	A23	L24	A25	S26	M29	I30	S34	F35	K40	P41	E42	T43	L44	N45	Y46	R47	T48	R53	D54	G55	L56	F57	C58	I61	F62	V65	K66	E69	C70	L71	C72	C73





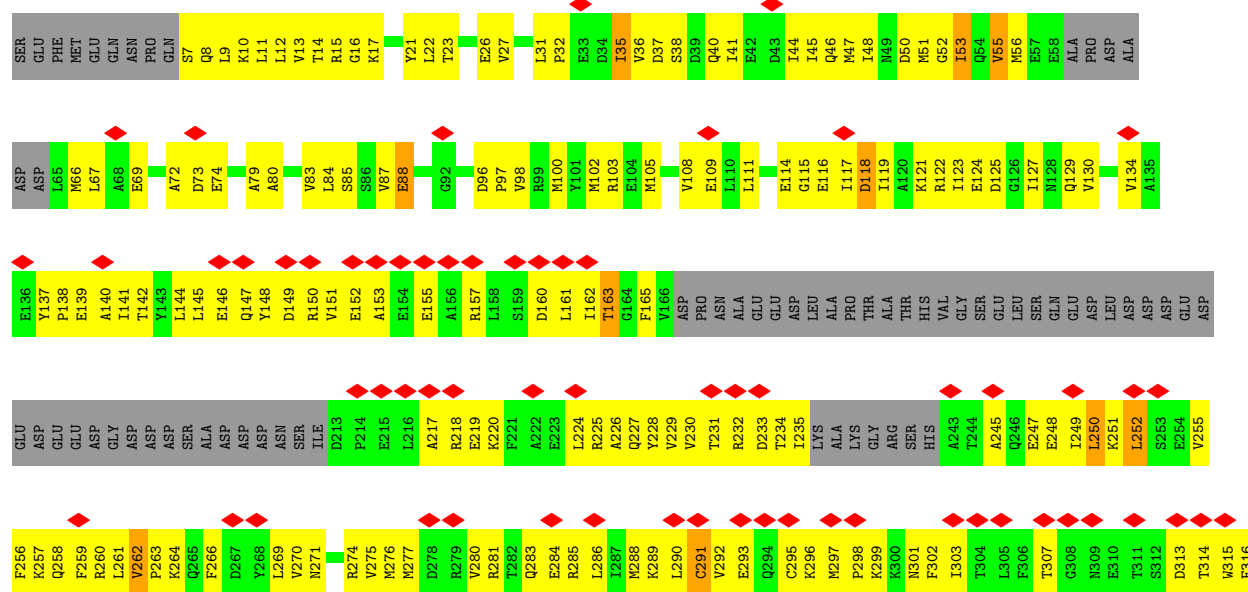
• Molecule 4: DNA-directed RNA polymerase subunit omega

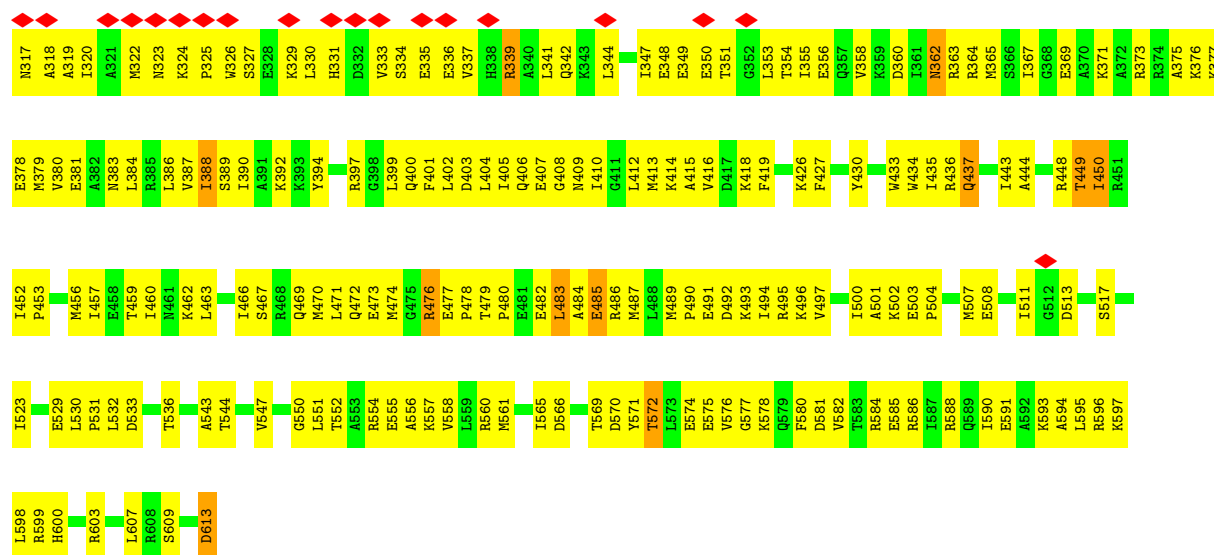
Chain K: 31% 46% 5% 18%



• Molecule 5: RNA polymerase sigma factor RpoD

Chain L: 13% 32% 53% 11%





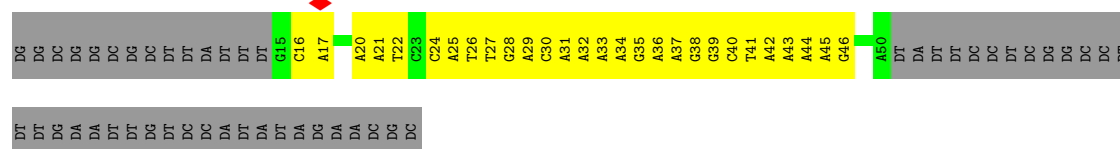
• Molecule 6: Protein TraR

Chain N: 51% 44%



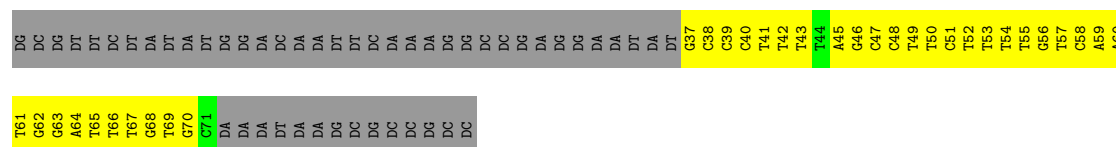
• Molecule 7: DNA (85-MER)

Chain O: 9% 33% 58%



• Molecule 8: DNA (85-MER)

Chain P: 39% 59%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	98324	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	80	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	0.223	Depositor
Minimum map value	-0.124	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.02	Depositor
Map size (Å)	332.8, 332.8, 332.8	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.3, 1.3, 1.3	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MG, 1N7

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	G	0.38	0/1793	0.55	0/2432
1	H	0.35	0/1697	0.54	0/2301
1	M	0.17	0/579	0.48	0/784
2	I	0.38	0/10728	0.55	2/14477 (0.0%)
3	J	0.36	0/10625	0.55	3/14345 (0.0%)
4	K	0.28	0/602	0.51	0/810
5	L	0.23	0/4461	0.51	0/6004
6	N	0.32	0/575	0.48	0/778
7	O	0.25	0/842	0.41	0/1297
8	P	0.25	0/790	0.47	0/1217
All	All	0.34	0/32692	0.53	5/44445 (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
2	I	0	2
3	J	0	1
All	All	0	3

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	9	LYS	N-CA-C	-6.20	104.78	112.72
3	J	53	ARG	CA-C-N	6.00	133.00	121.54
3	J	53	ARG	C-N-CA	6.00	133.00	121.54
3	J	839	VAL	N-CA-C	-5.91	107.37	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	I	162	GLY	N-CA-C	5.71	116.93	111.67

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	I	519	ASN	Peptide
2	I	897	PRO	Peptide
3	J	53	ARG	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	G	1771	0	1799	112	0
1	H	1678	0	1698	129	0
1	M	572	0	602	85	0
2	I	10559	0	10577	784	0
3	J	10466	0	10689	886	0
4	K	600	0	607	51	0
5	L	4407	0	4432	427	0
6	N	565	0	545	35	0
7	O	746	0	401	42	0
8	P	710	0	402	52	0
9	I	27	39	39	4	0
9	J	54	78	75	9	0
9	L	27	39	38	4	0
10	J	1	0	0	0	0
11	J	2	0	0	0	0
11	N	1	0	0	0	0
All	All	32186	156	31904	2439	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 38.

The worst 5 of 2439 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)
9:L:701:1N7:C19	9:L:701:1N7:C3	1.82	1.58
9:I:1401:1N7:C3	9:I:1401:1N7:C19	1.82	1.56
9:J:1505:1N7:C3	9:J:1505:1N7:C19	1.81	1.54
9:J:1504:1N7:C19	9:J:1504:1N7:C3	1.84	1.50
3:J:1179:PRO:HD2	3:J:1184:ASP:HA	1.34	1.08

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	G	228/329 (69%)	205 (90%)	23 (10%)	0	100	100
1	H	215/329 (65%)	189 (88%)	26 (12%)	0	100	100
1	M	71/329 (22%)	68 (96%)	3 (4%)	0	100	100
2	I	1338/1342 (100%)	1179 (88%)	157 (12%)	2 (0%)	48	78
3	J	1339/1430 (94%)	1197 (89%)	141 (10%)	1 (0%)	48	78
4	K	73/91 (80%)	68 (93%)	5 (7%)	0	100	100
5	L	540/616 (88%)	500 (93%)	40 (7%)	0	100	100
6	N	70/72 (97%)	65 (93%)	5 (7%)	0	100	100
All	All	3874/4538 (85%)	3471 (90%)	400 (10%)	3 (0%)	49	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	I	898	GLU
3	J	854	ALA
2	I	897	PRO

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	G	195/286 (68%)	168 (86%)	27 (14%)	3	14
1	H	184/286 (64%)	161 (88%)	23 (12%)	4	18
1	M	65/286 (23%)	62 (95%)	3 (5%)	24	50
2	I	1153/1157 (100%)	1020 (88%)	133 (12%)	5	20
3	J	1128/1189 (95%)	1010 (90%)	118 (10%)	6	23
4	K	65/75 (87%)	58 (89%)	7 (11%)	6	22
5	L	479/543 (88%)	448 (94%)	31 (6%)	15	42
6	N	60/61 (98%)	52 (87%)	8 (13%)	4	15
All	All	3329/3883 (86%)	2979 (90%)	350 (10%)	9	23

5 of 350 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	J	515	ARG
3	J	1220	ILE
3	J	591	ILE
3	J	837	ASP
3	J	1352	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 56 such sidechains are listed below:

Mol	Chain	Res	Type
3	J	340	GLN
6	N	68	GLN
3	J	792	ASN
1	M	276	HIS
5	L	258	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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$
9	1N7	J	1505	-	30,30,46	4.93	15 (50%)	47,48,72	2.12	15 (31%)
9	1N7	L	701	-	30,30,46	5.00	15 (50%)	47,48,72	2.48	20 (42%)
9	1N7	J	1504	-	30,30,46	5.20	16 (53%)	47,48,72	2.39	17 (36%)
9	1N7	I	1401	-	30,30,46	5.07	16 (53%)	47,48,72	2.09	10 (21%)

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
9	1N7	J	1505	-	-	0/7/72/92	0/4/4/4
9	1N7	L	701	-	-	5/7/72/92	0/4/4/4
9	1N7	J	1504	-	-	7/7/72/92	0/4/4/4
9	1N7	I	1401	-	-	5/7/72/92	0/4/4/4

The worst 5 of 62 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	J	1504	1N7	C3-C19	18.71	1.84	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	I	1401	1N7	C3-C19	17.97	1.82	1.53
9	L	701	1N7	C3-C19	17.48	1.82	1.53
9	J	1505	1N7	C3-C19	17.39	1.81	1.53
9	J	1504	1N7	C3-C4	12.37	1.73	1.53

The worst 5 of 62 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	L	701	1N7	C9-C5-C4	-7.86	110.59	117.67
9	J	1504	1N7	C9-C5-C4	-7.73	110.71	117.67
9	I	1401	1N7	C9-C5-C4	-6.76	111.58	117.67
9	L	701	1N7	C3-C19-C18	-6.29	101.58	110.89
9	J	1505	1N7	C19-C3-C4	-4.99	107.77	114.29

There are no chirality outliers.

5 of 17 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	J	1504	1N7	C21-C20-C9-C5
9	J	1504	1N7	C21-C20-C9-C8
9	I	1401	1N7	C21-C20-C9-C5
9	I	1401	1N7	C21-C20-C9-C8
9	I	1401	1N7	C22-C20-C9-C5

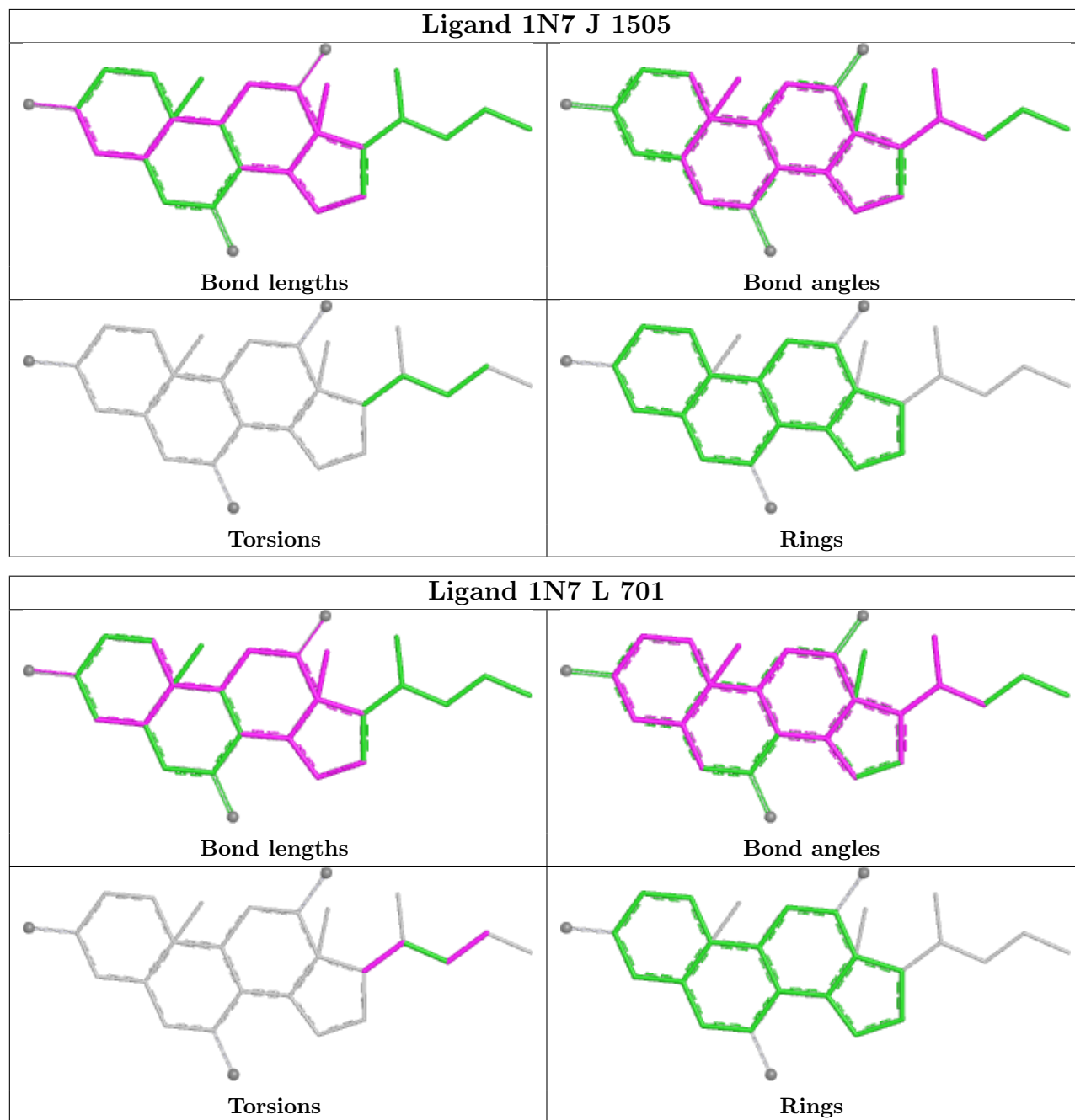
There are no ring outliers.

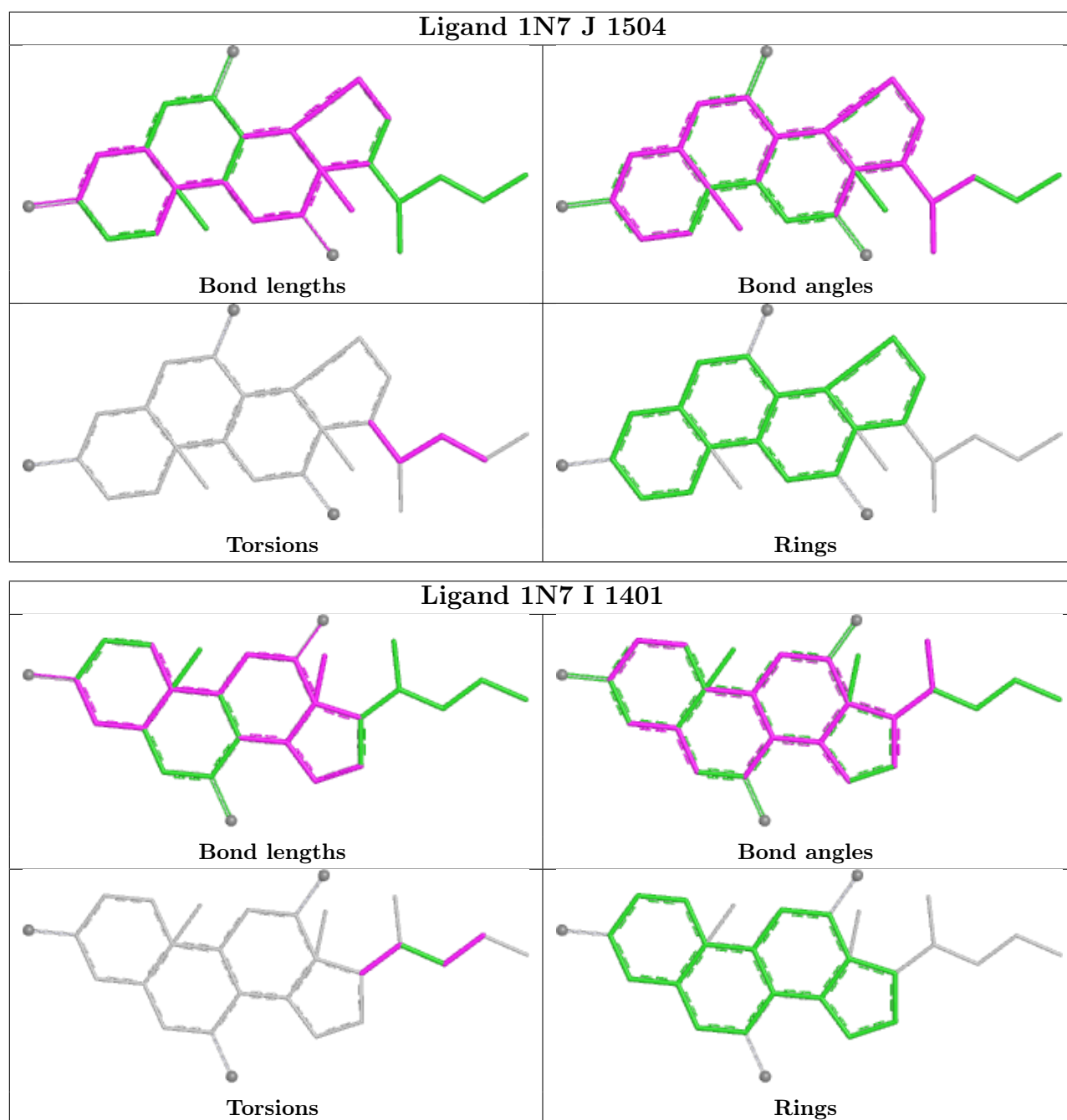
4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	J	1505	1N7	5	0
9	L	701	1N7	4	0
9	J	1504	1N7	4	0
9	I	1401	1N7	4	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.

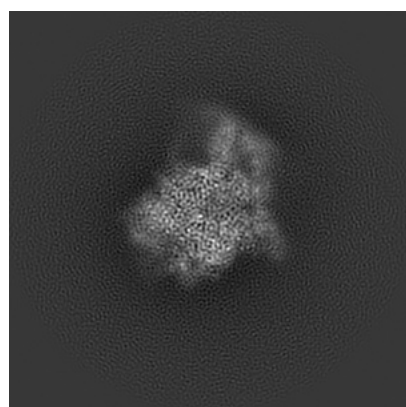
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20461. These allow visual inspection of the internal detail of the map and identification of artifacts.

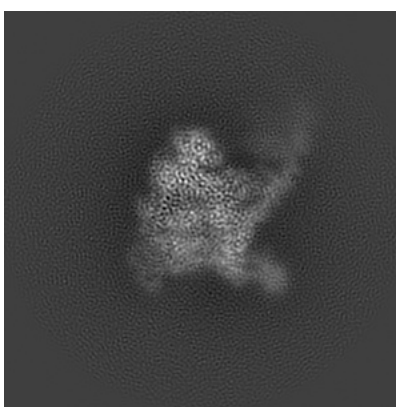
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

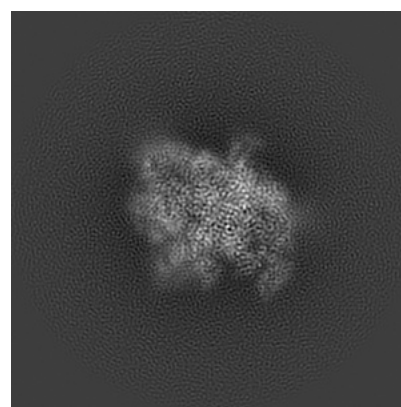
6.1.1 Primary map



X



Y

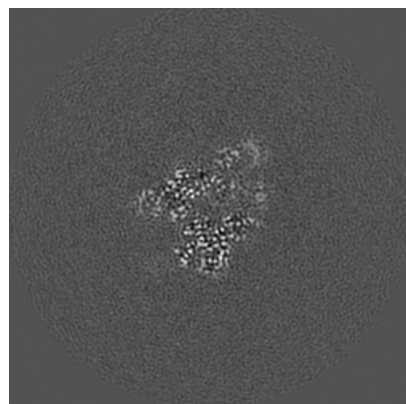


Z

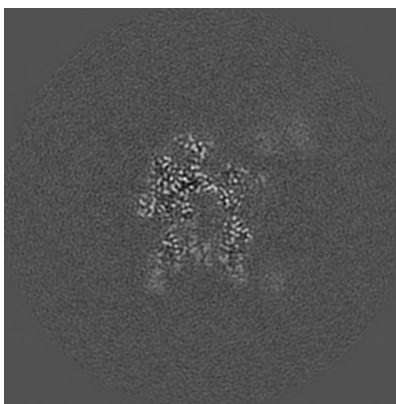
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

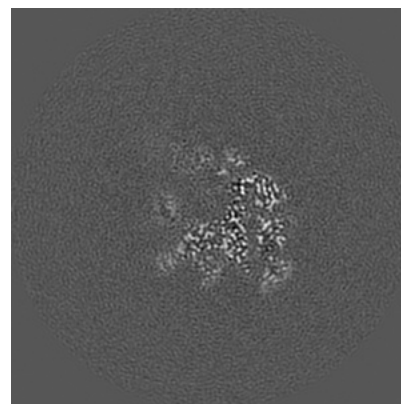
6.2.1 Primary map



X Index: 128



Y Index: 128

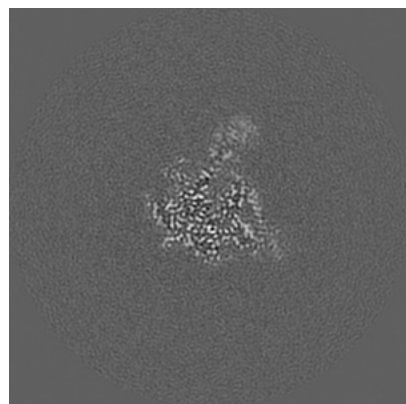


Z Index: 128

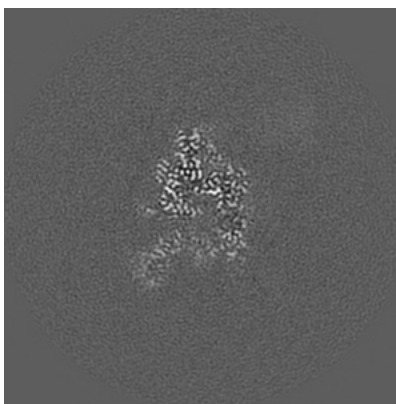
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

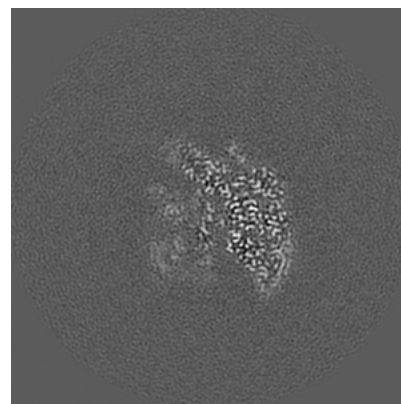
6.3.1 Primary map



X Index: 146



Y Index: 121

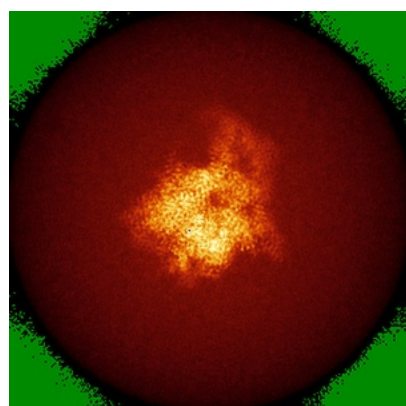


Z Index: 120

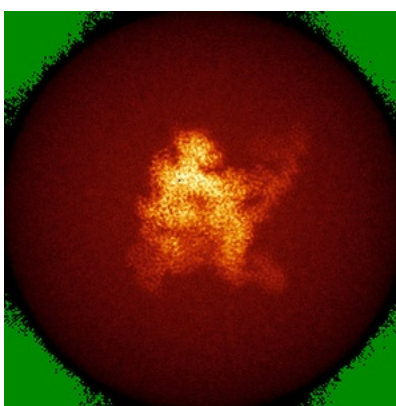
The images above show the largest variance slices of the map in three orthogonal directions.

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

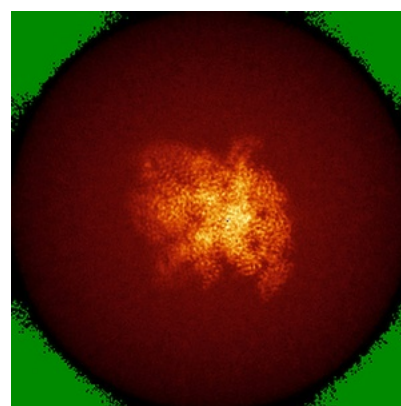
6.4.1 Primary map



X



Y

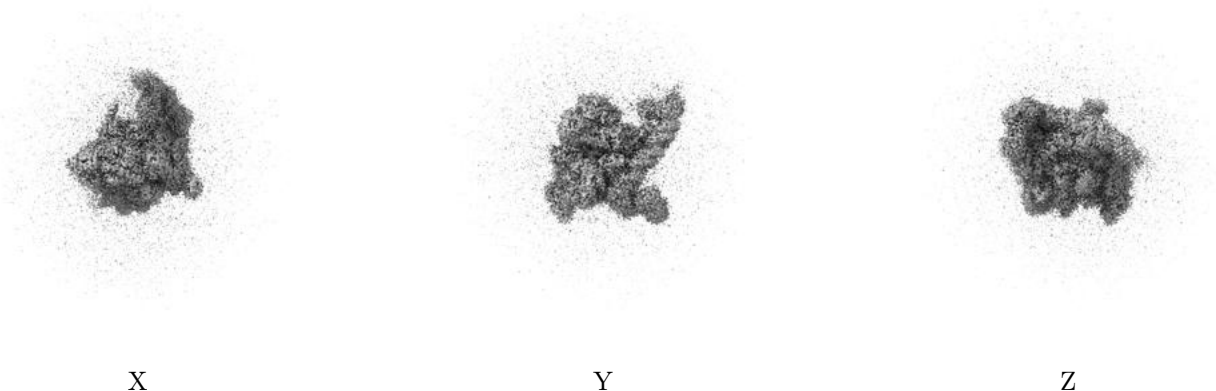


Z

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 [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.

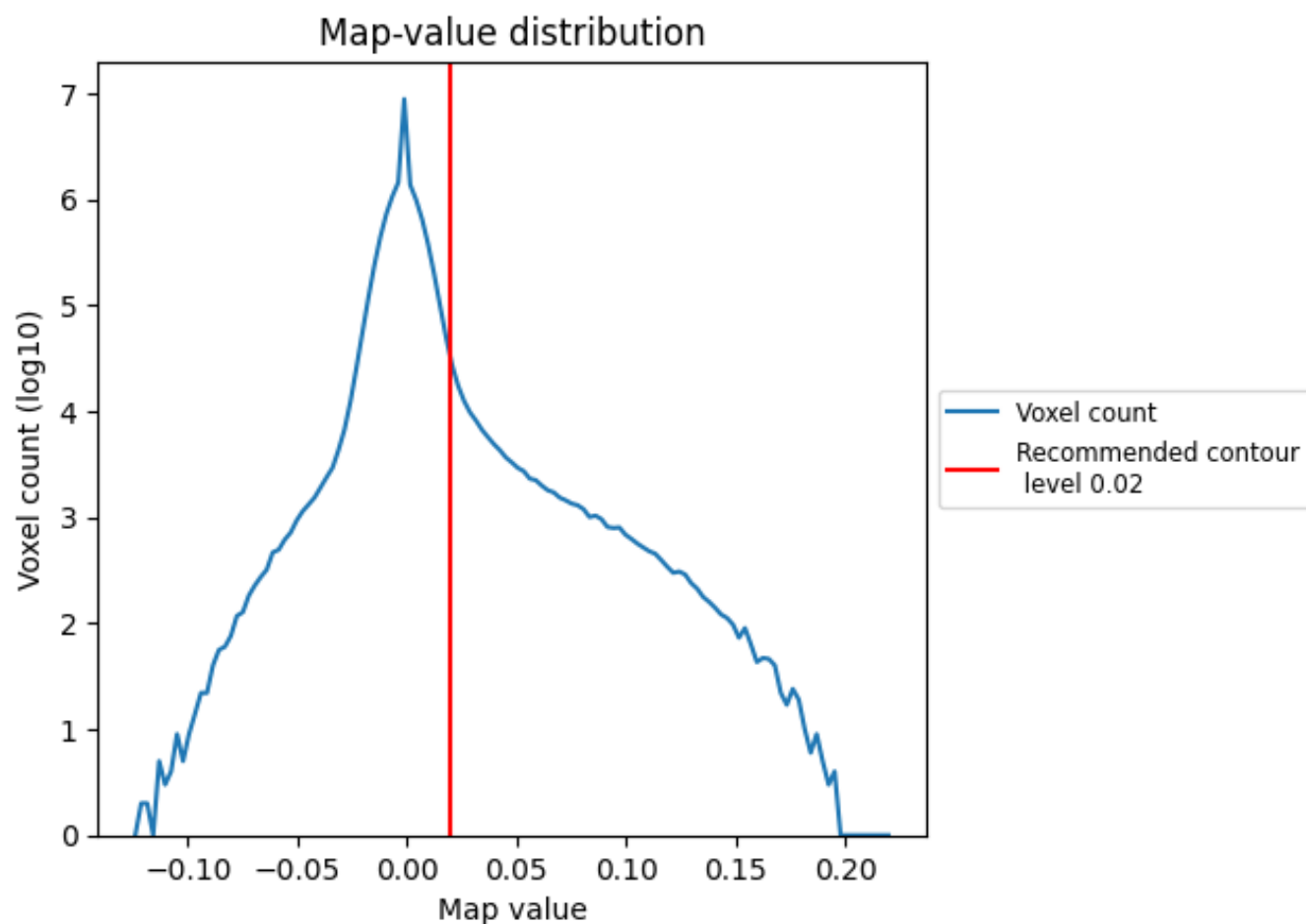
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

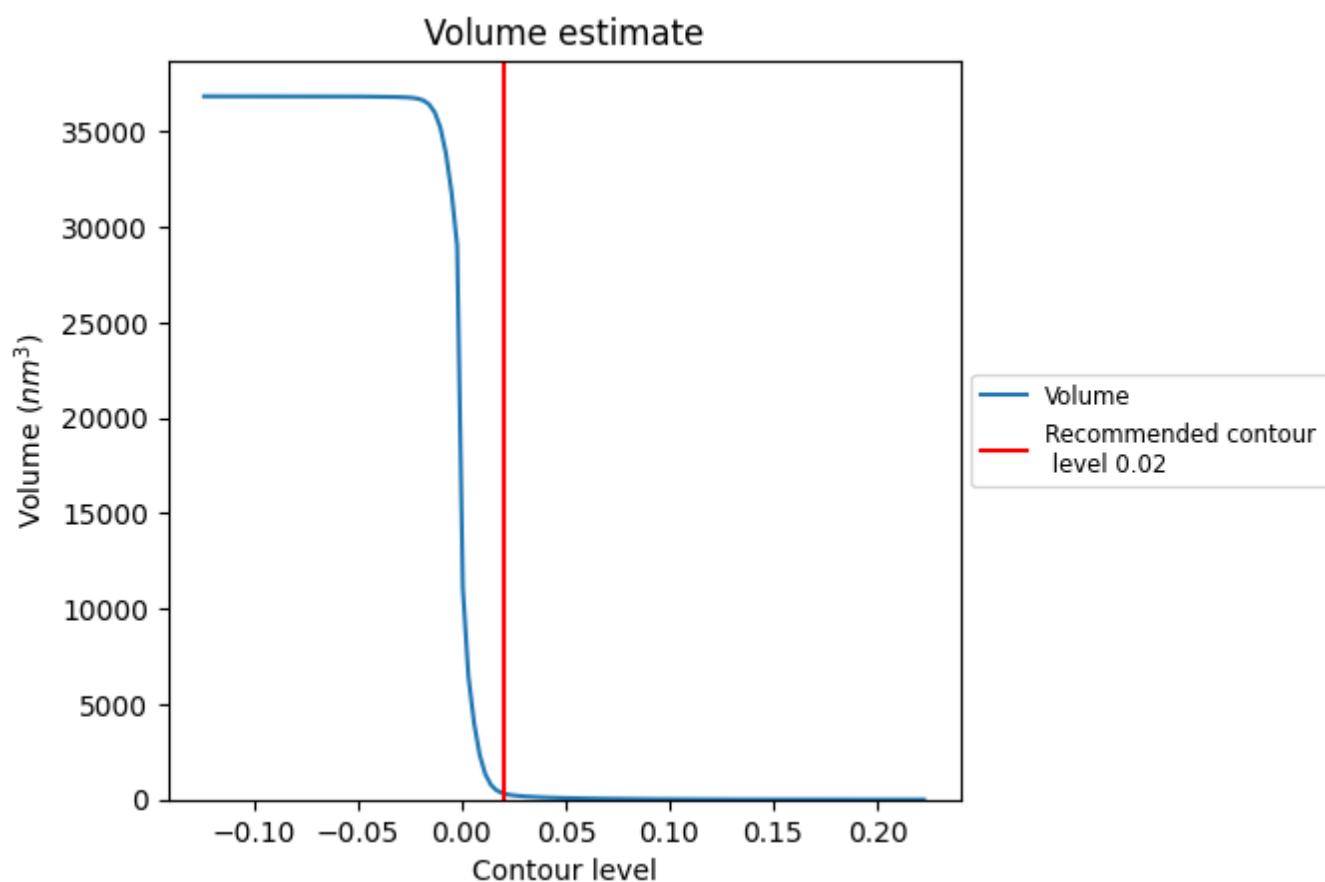
This section contains the results of statistical analysis of the map.

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.

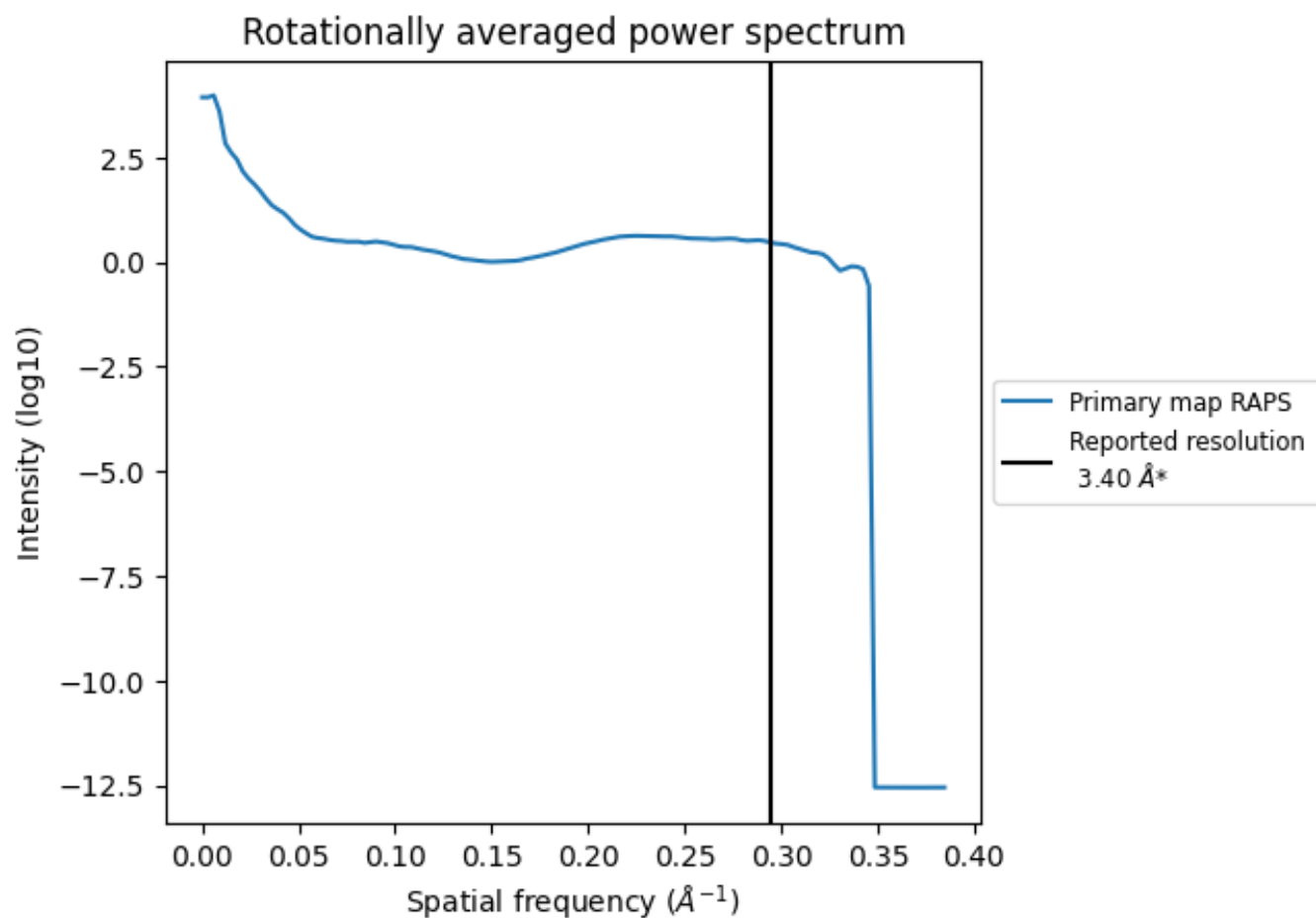
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 325 nm³; this corresponds to an approximate mass of 294 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 ⓘ

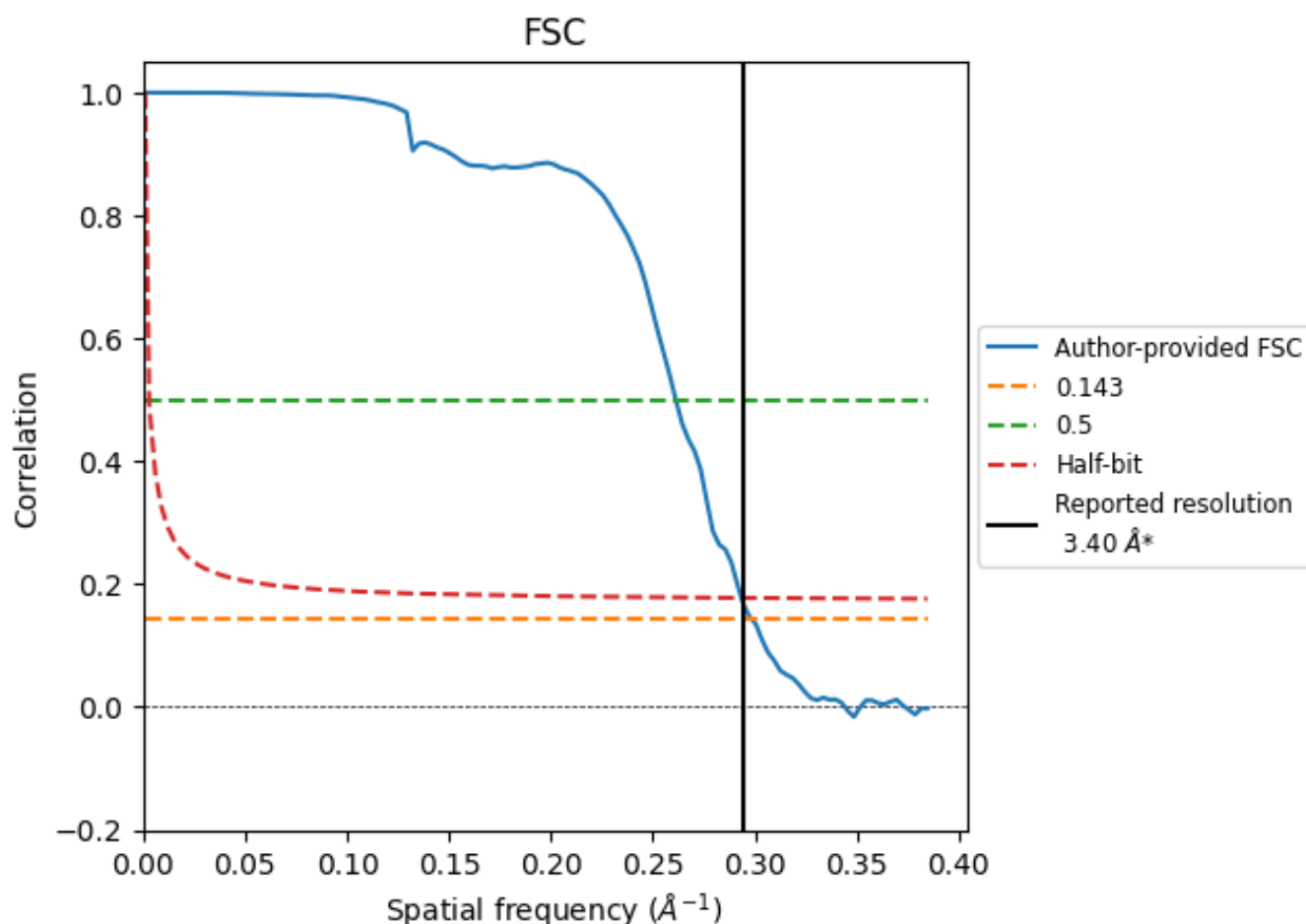


*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

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 Å⁻¹

8.2 Resolution estimates [i](#)

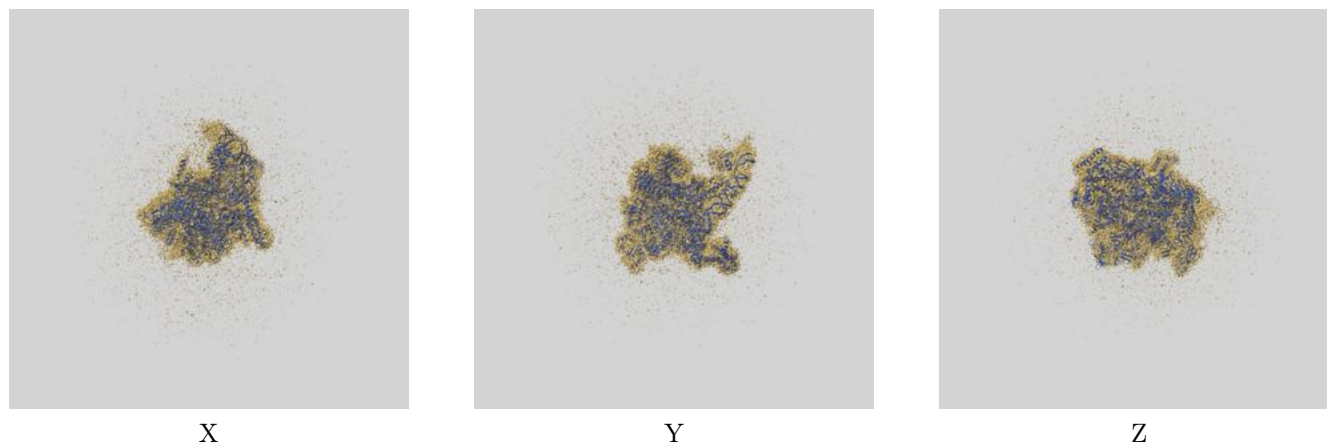
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	3.35	3.83	3.41
Unmasked-calculated*	-	-	-

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

9 Map-model fit [i](#)

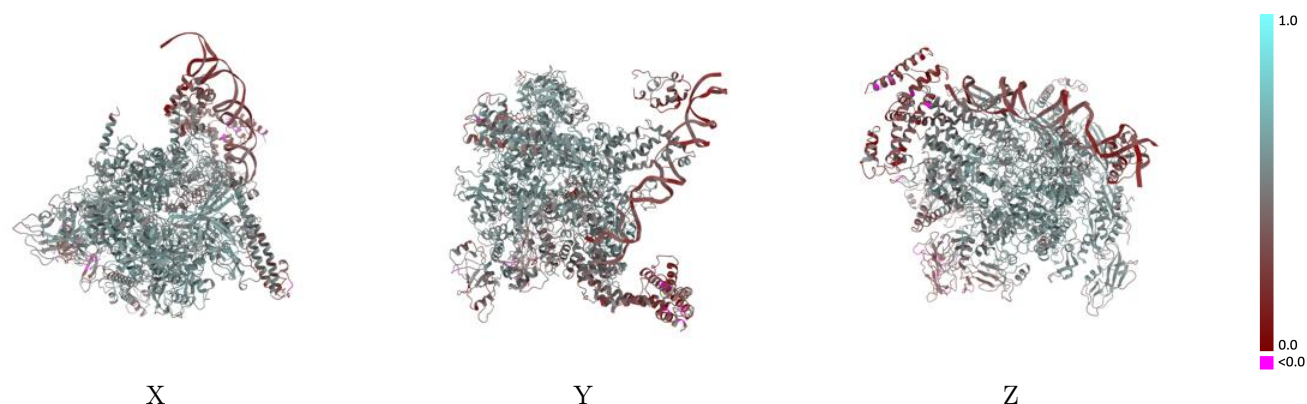
This section contains information regarding the fit between EMDB map EMD-20461 and PDB model 6PSR. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

9.1 Map-model overlay [i](#)



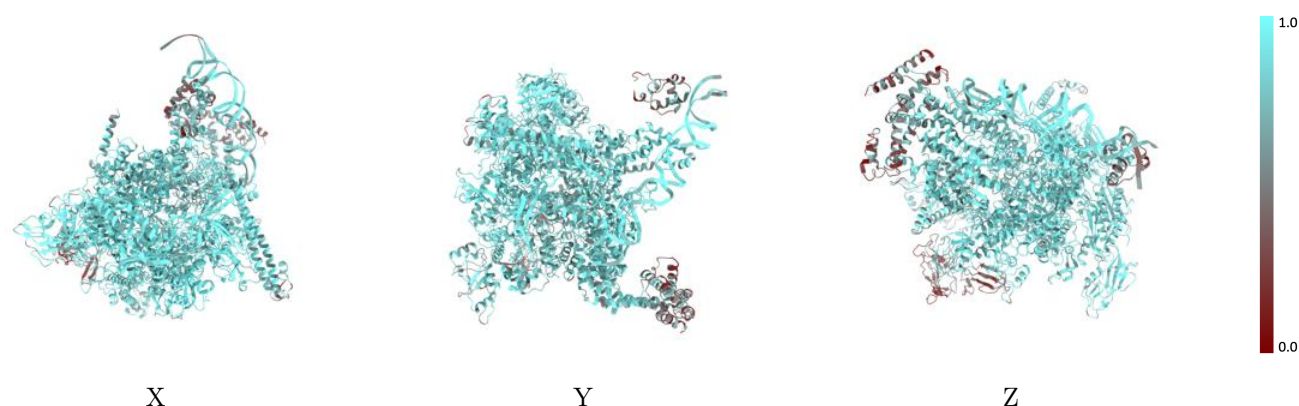
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



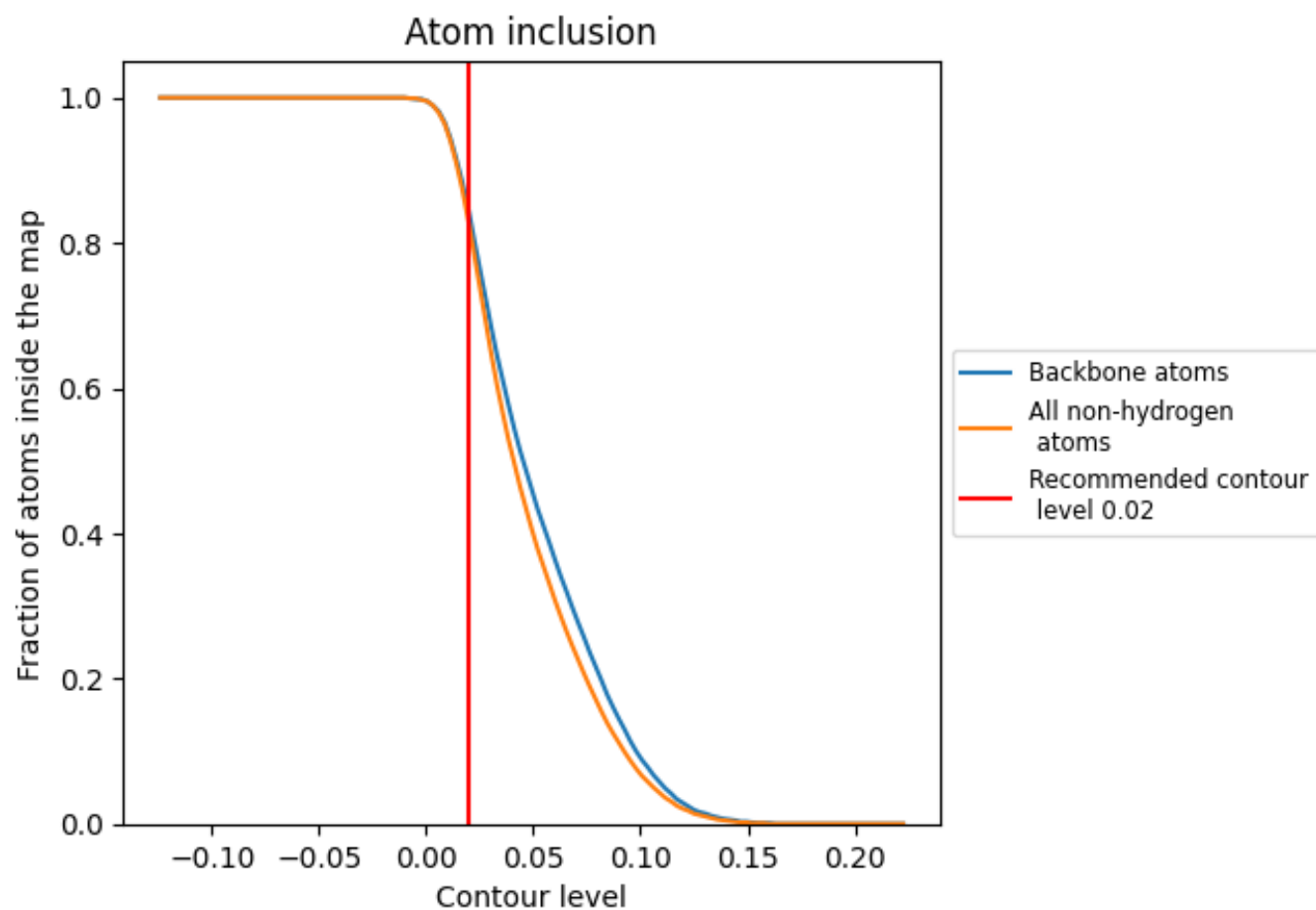
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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



At the recommended contour level, 84% of all backbone atoms, 82% 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	0.8250	0.4800
G	0.9000	0.5390
H	0.8870	0.5150
I	0.8860	0.5260
J	0.8120	0.4940
K	0.8240	0.5280
L	0.7080	0.3850
M	0.4550	0.3030
N	0.9150	0.5220
O	0.8060	0.2790
P	0.8110	0.2510

1.0
0.0
<0.0