



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 12:12 PM UTC

PDB ID : 6Q0K / pdb_00006q0k
EMDB ID : EMD-20551
Title : Structure of a MAPK pathway complex
Authors : Park, E.; Rawson, S.; Jeon, H.; Eck, M.J.
Deposited on : 2019-08-01
Resolution : 6.80 Å(reported)
Based on initial models : 4MNE, 3NKX

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:

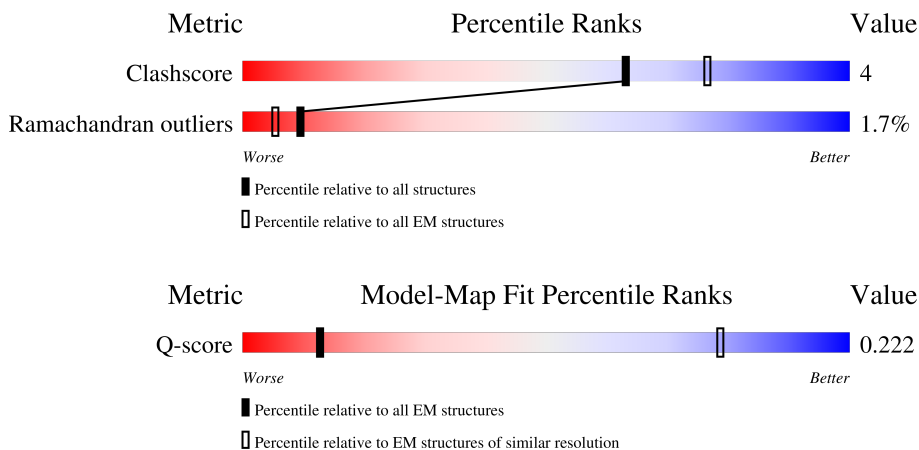
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
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 6.80 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	-
Q-score	-	25397	503 (6.30 - 7.30)

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	805	
1	B	805	
2	X	245	
2	Y	245	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5079 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Serine/threonine-protein kinase B-raf.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	285	Total	C	N	O	P	0	0
			1411	836	285	289	1		
1	B	281	Total	C	N	O	P	0	0
			1392	825	281	285	1		

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-26	MET	-	expression tag	UNP P15056
A	-25	SER	-	expression tag	UNP P15056
A	-24	TYR	-	expression tag	UNP P15056
A	-23	TYR	-	expression tag	UNP P15056
A	-22	HIS	-	expression tag	UNP P15056
A	-21	HIS	-	expression tag	UNP P15056
A	-20	HIS	-	expression tag	UNP P15056
A	-19	HIS	-	expression tag	UNP P15056
A	-18	HIS	-	expression tag	UNP P15056
A	-17	HIS	-	expression tag	UNP P15056
A	-16	HIS	-	expression tag	UNP P15056
A	-15	HIS	-	expression tag	UNP P15056
A	-14	ASP	-	expression tag	UNP P15056
A	-13	ILE	-	expression tag	UNP P15056
A	-12	PRO	-	expression tag	UNP P15056
A	-11	THR	-	expression tag	UNP P15056
A	-10	THR	-	expression tag	UNP P15056
A	-9	GLU	-	expression tag	UNP P15056
A	-8	ASN	-	expression tag	UNP P15056
A	-7	LEU	-	expression tag	UNP P15056
A	-6	TYR	-	expression tag	UNP P15056
A	-5	PHE	-	expression tag	UNP P15056
A	-4	GLN	-	expression tag	UNP P15056
A	-3	GLY	-	expression tag	UNP P15056
A	-2	ALA	-	expression tag	UNP P15056
A	-1	MET	-	expression tag	UNP P15056

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ASP	-	expression tag	UNP P15056
A	767	GLY	-	expression tag	UNP P15056
A	768	THR	-	expression tag	UNP P15056
A	769	SER	-	expression tag	UNP P15056
A	770	ALA	-	expression tag	UNP P15056
A	771	TRP	-	expression tag	UNP P15056
A	772	SER	-	expression tag	UNP P15056
A	773	HIS	-	expression tag	UNP P15056
A	774	PRO	-	expression tag	UNP P15056
A	775	GLN	-	expression tag	UNP P15056
A	776	PHE	-	expression tag	UNP P15056
A	777	GLU	-	expression tag	UNP P15056
A	778	LYS	-	expression tag	UNP P15056
B	-26	MET	-	expression tag	UNP P15056
B	-25	SER	-	expression tag	UNP P15056
B	-24	TYR	-	expression tag	UNP P15056
B	-23	TYR	-	expression tag	UNP P15056
B	-22	HIS	-	expression tag	UNP P15056
B	-21	HIS	-	expression tag	UNP P15056
B	-20	HIS	-	expression tag	UNP P15056
B	-19	HIS	-	expression tag	UNP P15056
B	-18	HIS	-	expression tag	UNP P15056
B	-17	HIS	-	expression tag	UNP P15056
B	-16	HIS	-	expression tag	UNP P15056
B	-15	HIS	-	expression tag	UNP P15056
B	-14	ASP	-	expression tag	UNP P15056
B	-13	ILE	-	expression tag	UNP P15056
B	-12	PRO	-	expression tag	UNP P15056
B	-11	THR	-	expression tag	UNP P15056
B	-10	THR	-	expression tag	UNP P15056
B	-9	GLU	-	expression tag	UNP P15056
B	-8	ASN	-	expression tag	UNP P15056
B	-7	LEU	-	expression tag	UNP P15056
B	-6	TYR	-	expression tag	UNP P15056
B	-5	PHE	-	expression tag	UNP P15056
B	-4	GLN	-	expression tag	UNP P15056
B	-3	GLY	-	expression tag	UNP P15056
B	-2	ALA	-	expression tag	UNP P15056
B	-1	MET	-	expression tag	UNP P15056
B	0	ASP	-	expression tag	UNP P15056
B	767	GLY	-	expression tag	UNP P15056
B	768	THR	-	expression tag	UNP P15056

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Chain	Residue	Modelled	Actual	Comment	Reference
B	769	SER	-	expression tag	UNP P15056
B	770	ALA	-	expression tag	UNP P15056
B	771	TRP	-	expression tag	UNP P15056
B	772	SER	-	expression tag	UNP P15056
B	773	HIS	-	expression tag	UNP P15056
B	774	PRO	-	expression tag	UNP P15056
B	775	GLN	-	expression tag	UNP P15056
B	776	PHE	-	expression tag	UNP P15056
B	777	GLU	-	expression tag	UNP P15056
B	778	LYS	-	expression tag	UNP P15056

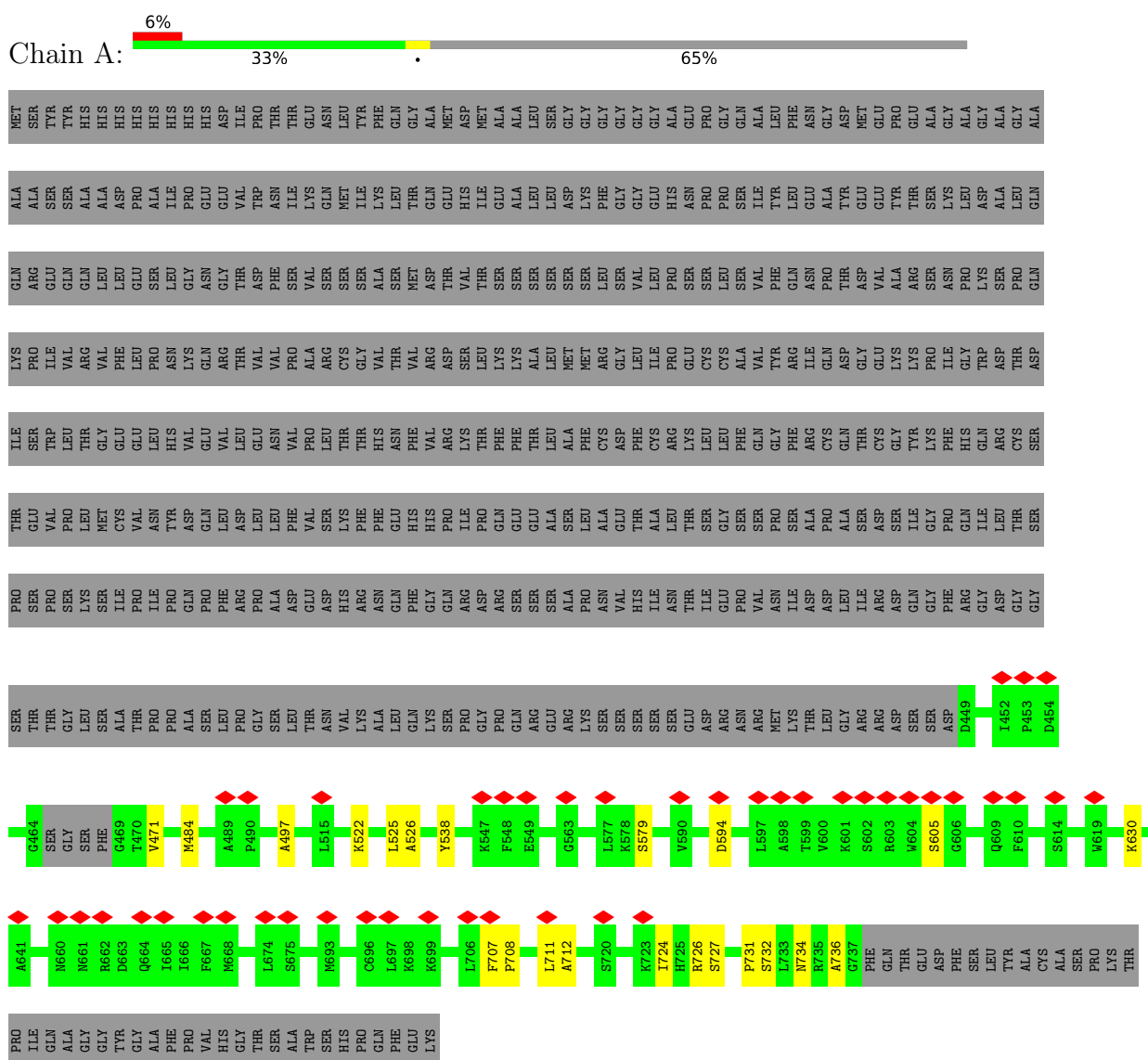
- Molecule 2 is a protein called 14-3-3 protein zeta/delta.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	X	229	Total	C	N	O	0	0
			1138	680	229	229		
2	Y	229	Total	C	N	O	0	0
			1138	680	229	229		

3 Residue-property plots

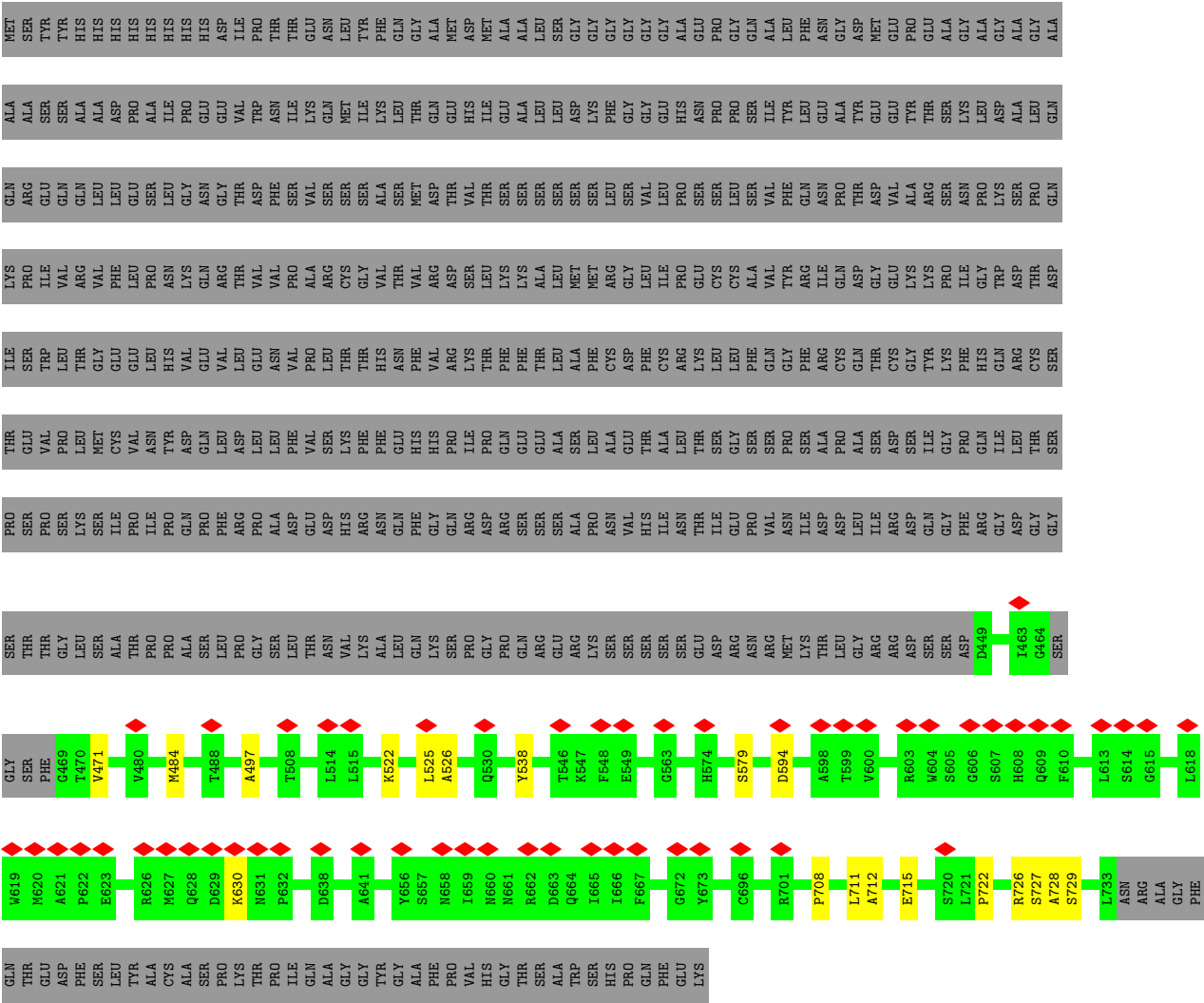
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

- Molecule 1: Serine/threonine-protein kinase B-raf

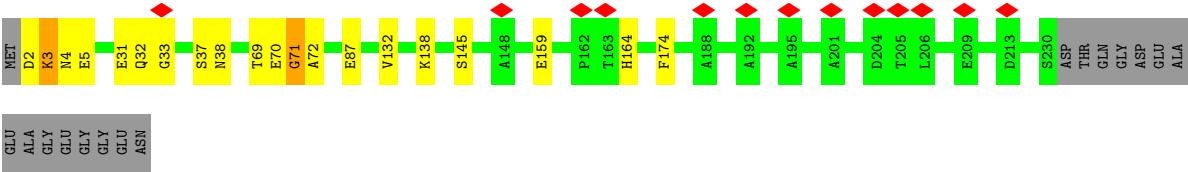
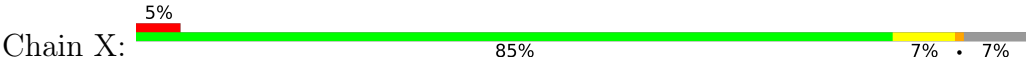


- Molecule 1: Serine/threonine-protein kinase B-raf

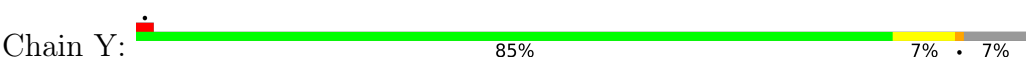




• Molecule 2: 14-3-3 protein zeta/delta



• Molecule 2: 14-3-3 protein zeta/delta



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	66215	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.128	Depositor
Minimum map value	-0.038	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.021	Depositor
Map size ($\text{\AA}$)	217.6, 217.6, 217.6	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.85, 0.85, 0.85	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SEP

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	0.80	0/1398	1.17	5/1941 (0.3%)
1	B	0.80	0/1379	1.14	3/1915 (0.2%)
2	X	1.56	1/1137 (0.1%)	1.51	10/1586 (0.6%)
2	Y	1.56	1/1137 (0.1%)	1.51	10/1586 (0.6%)
All	All	1.20	2/5051 (0.0%)	1.33	28/7028 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	X	87	GLU	CA-C	-5.54	1.45	1.52
2	Y	87	GLU	CA-C	-5.50	1.45	1.52

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Y	164	HIS	CA-C-N	8.48	128.99	119.32
2	Y	164	HIS	C-N-CA	8.48	128.99	119.32
2	X	164	HIS	CA-C-N	8.45	128.95	119.32
2	X	164	HIS	C-N-CA	8.45	128.95	119.32
2	Y	159	GLU	N-CA-C	6.70	121.74	113.50
2	X	159	GLU	N-CA-C	6.69	121.73	113.50
2	X	145	SER	N-CA-C	-6.33	103.67	111.40
2	Y	145	SER	N-CA-C	-6.31	103.70	111.40
2	X	38	ASN	CB-CA-C	6.08	120.22	109.65
2	Y	38	ASN	CB-CA-C	6.07	120.21	109.65
2	X	37	SER	CA-C-N	-5.98	111.92	122.26
2	X	37	SER	C-N-CA	-5.98	111.92	122.26
2	Y	37	SER	CA-C-N	-5.96	111.94	122.26
2	Y	37	SER	C-N-CA	-5.96	111.94	122.26
1	A	605	SER	N-CA-C	-5.70	101.86	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	X	132	VAL	N-CA-CB	-5.57	105.12	112.36
2	Y	132	VAL	N-CA-CB	-5.57	105.12	112.36
1	A	471	VAL	N-CA-C	5.53	115.86	108.11
1	B	471	VAL	N-CA-C	5.51	115.83	108.11
2	X	138	LYS	N-CA-C	5.51	117.28	111.28
2	Y	138	LYS	N-CA-C	5.46	117.23	111.28
1	B	630	LYS	N-CA-C	-5.32	105.40	111.14
1	A	731	PRO	N-CA-C	-5.29	101.58	112.47
1	A	630	LYS	N-CA-C	-5.27	105.45	111.14
1	A	497	ALA	N-CA-C	-5.20	105.30	110.97
1	B	497	ALA	N-CA-C	-5.18	105.32	110.97
2	Y	174	PHE	N-CA-C	-5.06	105.84	111.36
2	X	174	PHE	N-CA-C	-5.05	105.86	111.36

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1411	0	623	9	0
1	B	1392	0	611	7	0
2	X	1138	0	529	7	0
2	Y	1138	0	529	8	0
All	All	5079	0	2292	31	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 (31) 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:484:MET:CB	1:B:526:ALA:HB2	2.04	0.87
1:A:484:MET:CB	1:A:526:ALA:HB2	2.04	0.87
1:B:711:LEU:O	1:B:715:GLU:CB	2.23	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:X:71:GLY:HA2	2:X:72:ALA:C	2.17	0.70
2:Y:71:GLY:HA2	2:Y:72:ALA:C	2.17	0.69
2:X:2:ASP:O	2:X:4:ASN:N	2.29	0.66
2:Y:2:ASP:O	2:Y:4:ASN:N	2.29	0.65
2:X:31:GLU:C	2:X:33:GLY:H	2.09	0.60
1:A:538:TYR:N	1:A:579:SER:O	2.34	0.60
2:Y:31:GLU:C	2:Y:33:GLY:H	2.08	0.59
1:B:538:TYR:N	1:B:579:SER:O	2.37	0.57
1:A:484:MET:HA	1:A:526:ALA:HA	1.88	0.55
1:B:484:MET:HA	1:B:526:ALA:HA	1.88	0.54
2:Y:31:GLU:C	2:Y:33:GLY:N	2.65	0.53
1:A:734:ASN:C	1:A:736:ALA:H	2.16	0.53
2:X:31:GLU:C	2:X:33:GLY:N	2.65	0.53
1:B:484:MET:HA	1:B:525:LEU:O	2.10	0.51
1:A:484:MET:HA	1:A:525:LEU:O	2.11	0.50
2:X:69:THR:O	2:X:70:GLU:C	2.56	0.48
2:X:2:ASP:O	2:X:3:LYS:C	2.57	0.48
2:Y:2:ASP:O	2:Y:3:LYS:C	2.57	0.48
2:Y:69:THR:O	2:Y:70:GLU:C	2.56	0.47
2:Y:2:ASP:O	2:Y:5:GLU:N	2.46	0.47
1:B:727:SER:O	1:B:729:SEP:N	2.48	0.47
1:A:707:PHE:O	1:A:711:LEU:N	2.37	0.46
2:X:2:ASP:O	2:X:5:GLU:N	2.46	0.46
1:B:708:PRO:O	1:B:712:ALA:N	2.39	0.45
1:A:708:PRO:O	1:A:712:ALA:N	2.46	0.43
1:A:734:ASN:C	1:A:736:ALA:N	2.77	0.43
1:A:484:MET:CB	1:A:526:ALA:CB	2.89	0.41
2:Y:2:ASP:C	2:Y:4:ASN:N	2.78	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	280/805 (35%)	253 (90%)	21 (8%)	6 (2%)	5	30
1	B	276/805 (34%)	259 (94%)	12 (4%)	5 (2%)	6	34
2	X	227/245 (93%)	217 (96%)	7 (3%)	3 (1%)	9	42
2	Y	227/245 (93%)	217 (96%)	7 (3%)	3 (1%)	9	42
All	All	1010/2100 (48%)	946 (94%)	47 (5%)	17 (2%)	9	36

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	522	LYS
1	A	732	SER
1	B	522	LYS
2	X	3	LYS
2	Y	3	LYS
1	A	724	ILE
1	A	727	SER
1	B	722	PRO
1	B	726	ARG
1	B	728	ALA
2	X	71	GLY
2	Y	71	GLY
1	A	594	ASP
1	B	594	ASP
2	X	32	GLN
2	Y	32	GLN
1	A	726	ARG

5.3.2 Protein sidechains ⓘ

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	SEP	A	729	1	8,9,10	0.70	0	7,12,14	0.63	0
1	SEP	B	729	1	8,9,10	0.67	0	7,12,14	0.62	0

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	SEP	A	729	1	-	5/6/8/10	-
1	SEP	B	729	1	-	4/6/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	729	SEP	CA-CB-OG-P
1	A	729	SEP	CB-OG-P-O1P
1	A	729	SEP	CB-OG-P-O2P
1	A	729	SEP	CB-OG-P-O3P
1	B	729	SEP	CB-OG-P-O1P
1	B	729	SEP	CB-OG-P-O2P
1	B	729	SEP	CB-OG-P-O3P
1	A	729	SEP	N-CA-CB-OG
1	B	729	SEP	N-CA-CB-OG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	729	SEP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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-20551. These allow visual inspection of the internal detail of the map and identification of artifacts.

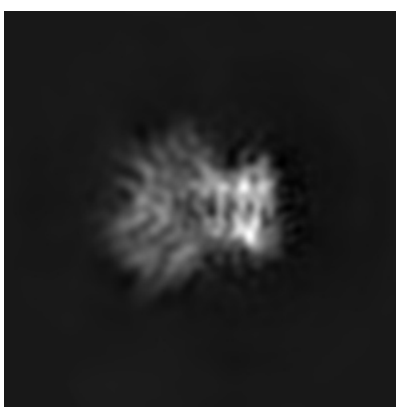
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

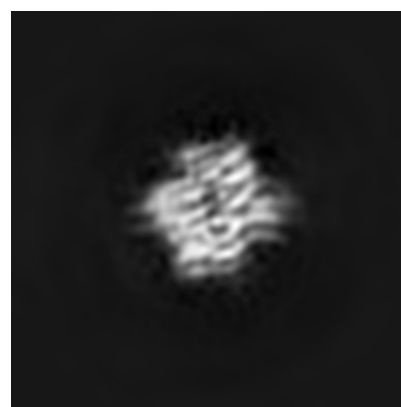
6.1.1 Primary map



X



Y



Z

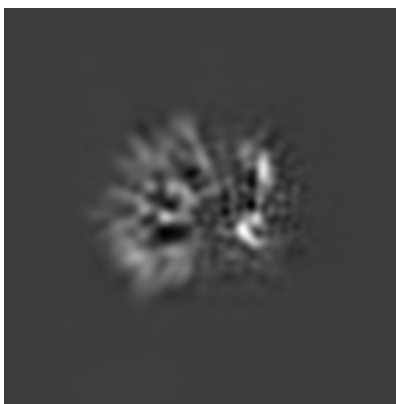
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

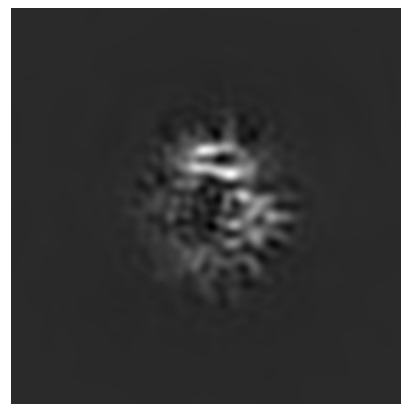
6.2.1 Primary map



X Index: 128



Y Index: 128

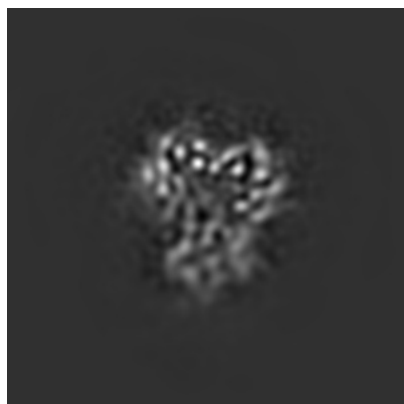


Z Index: 128

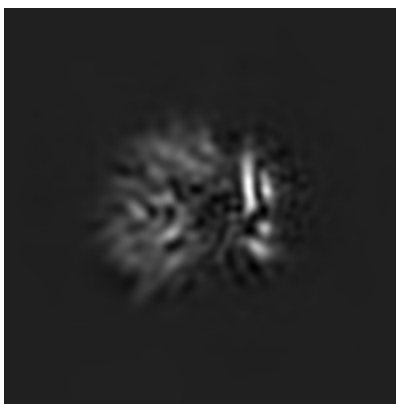
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 140



Y Index: 122



Z Index: 156

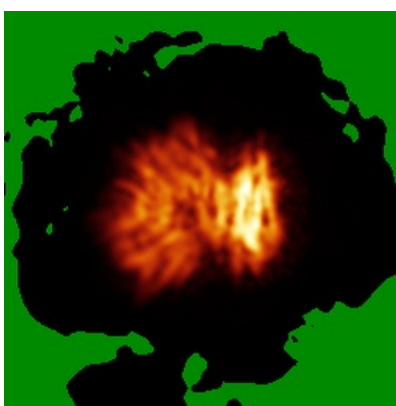
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



X



Y

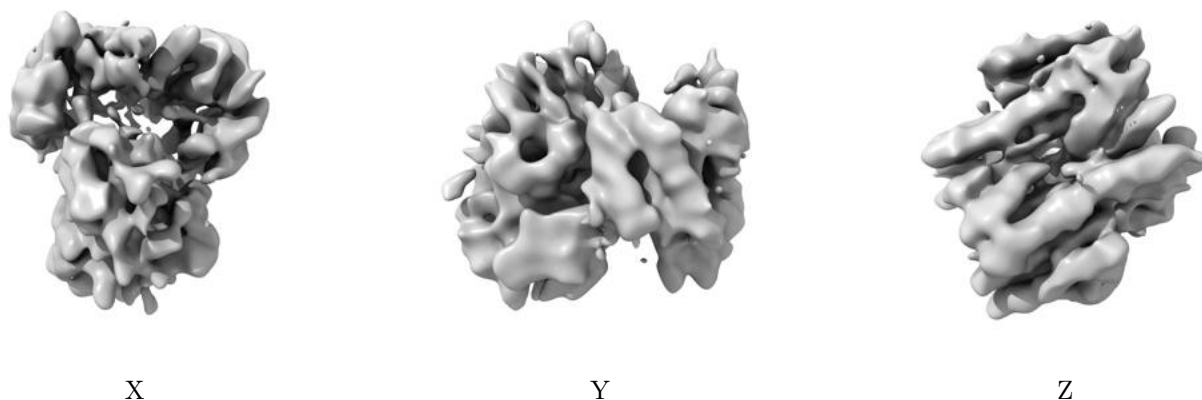


Z

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 0.021. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

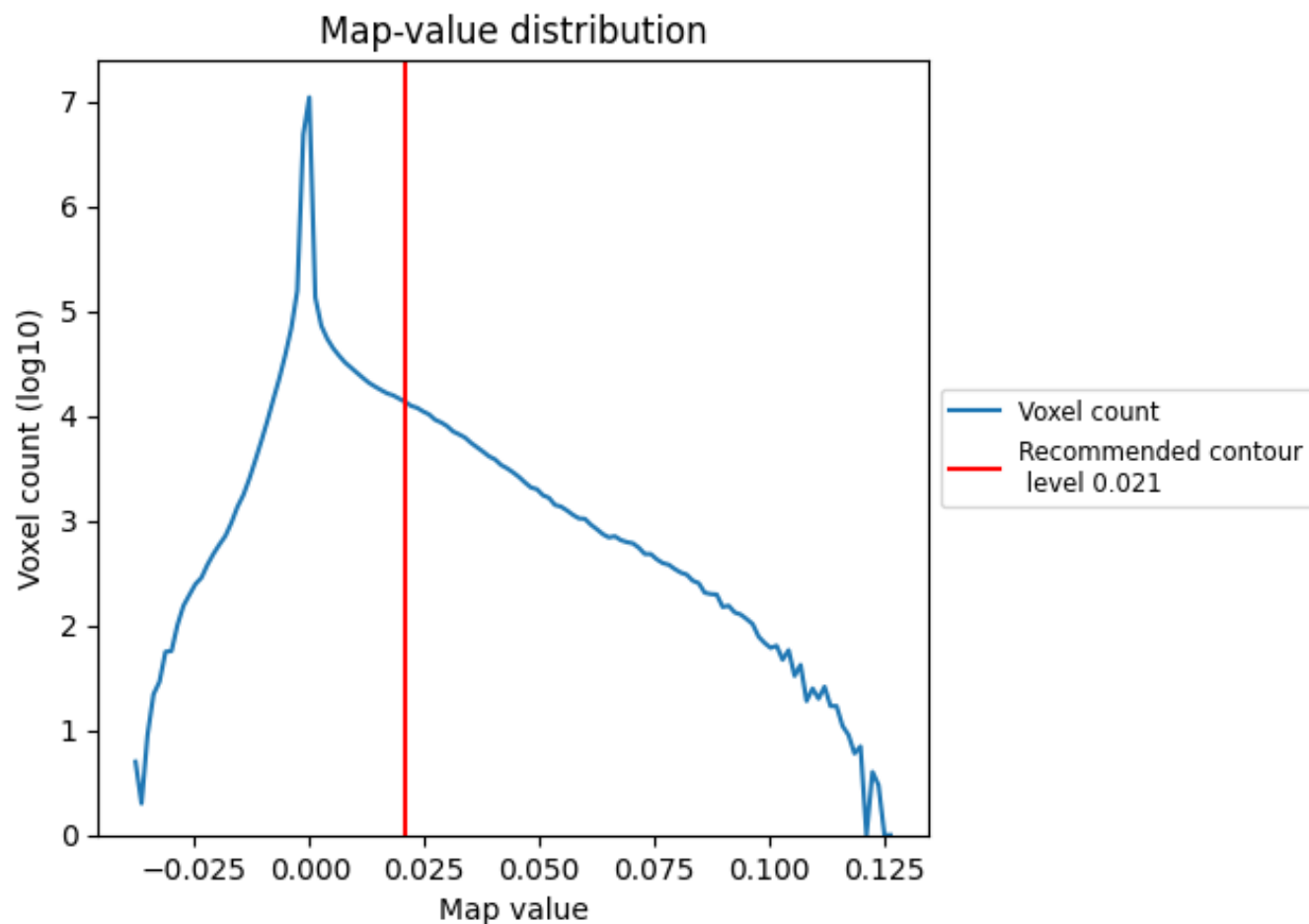
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

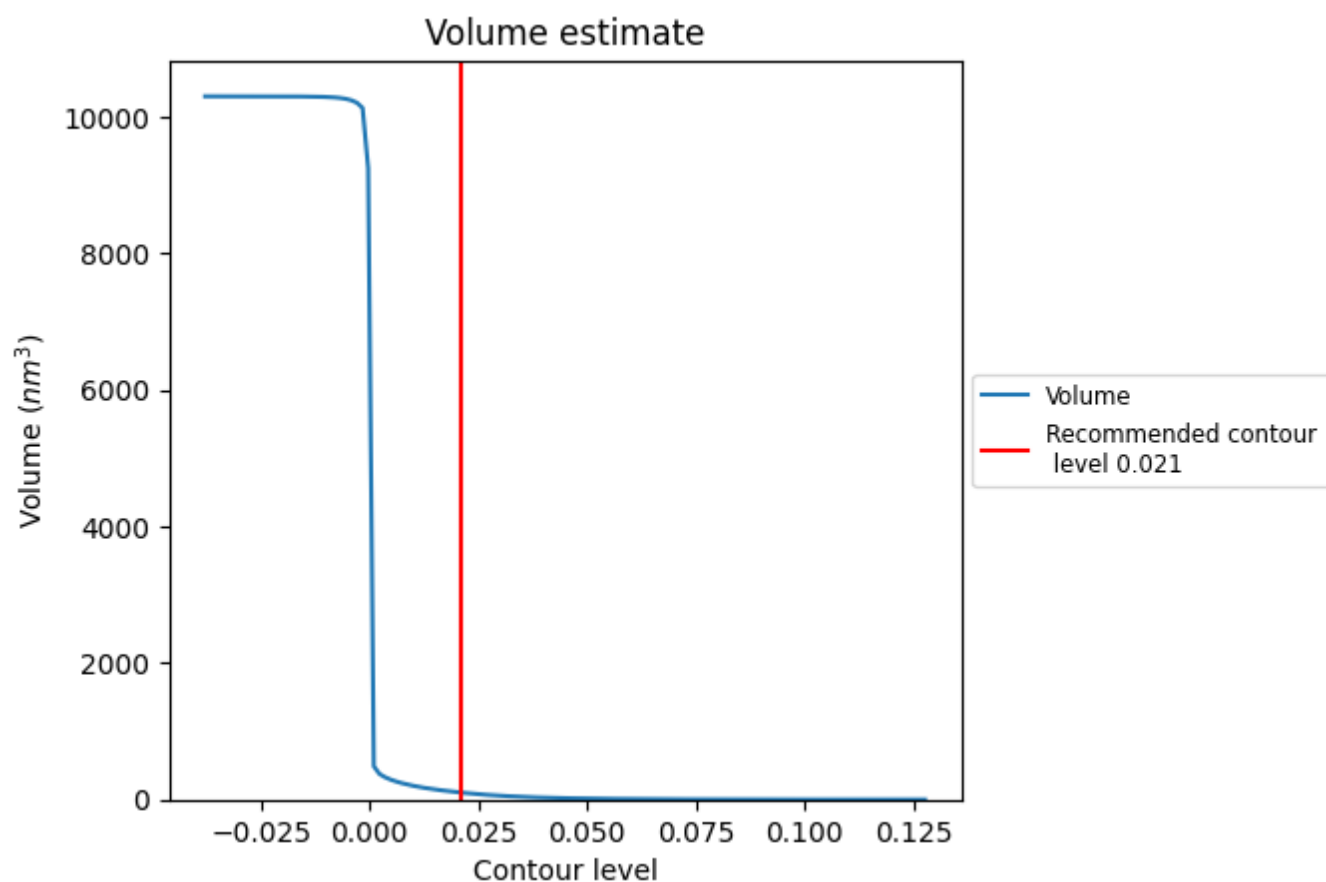
This section contains the results of statistical analysis of the map.

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.

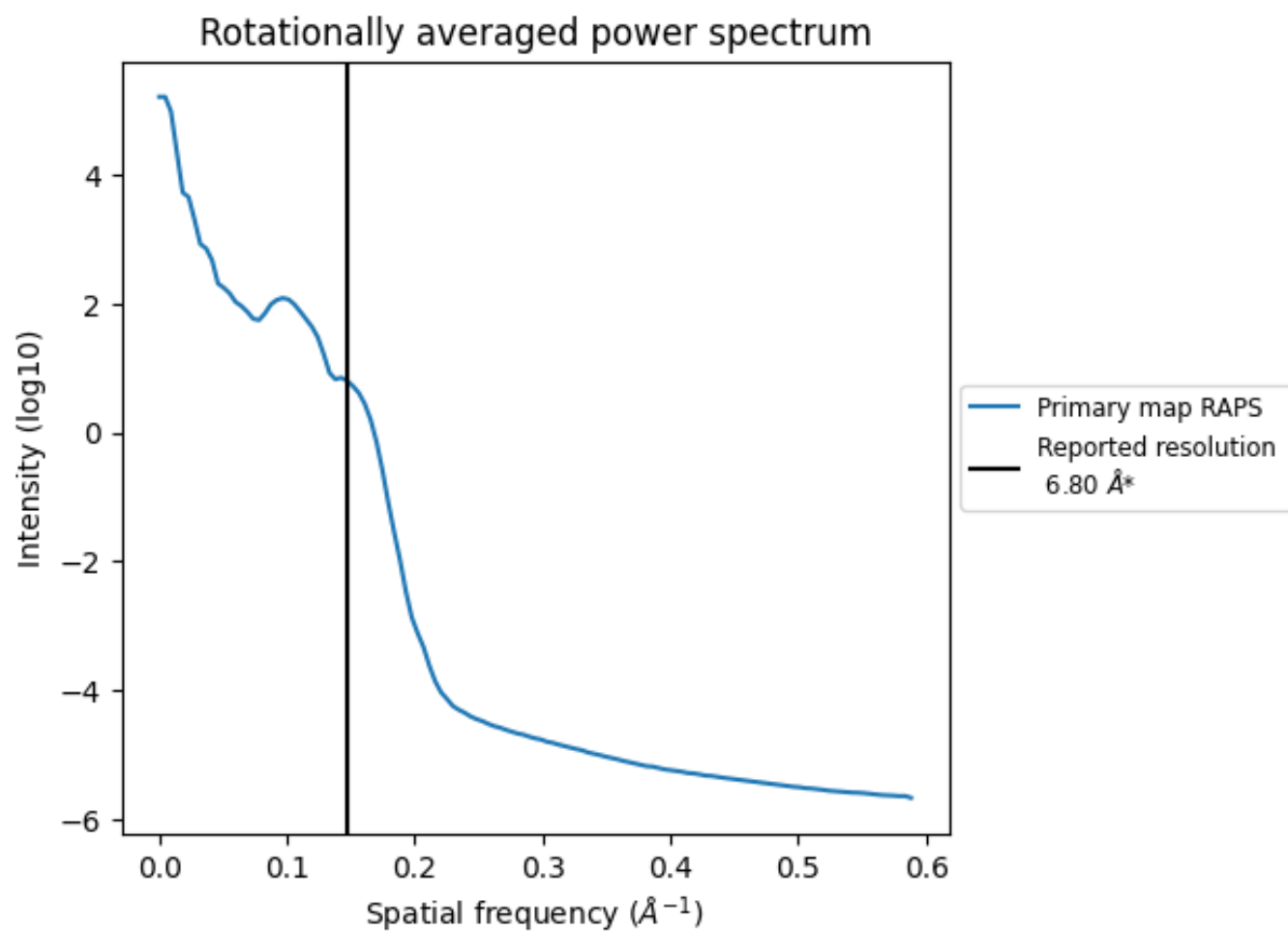
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 102 nm³; this corresponds to an approximate mass of 93 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 0.147 Å⁻¹

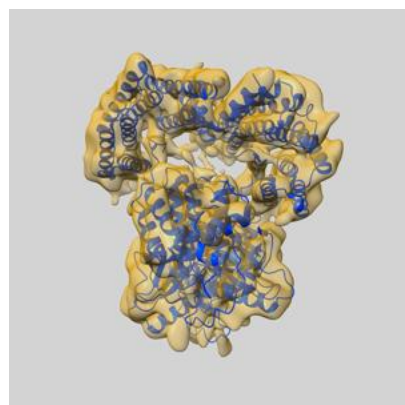
8 Fourier-Shell correlation

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

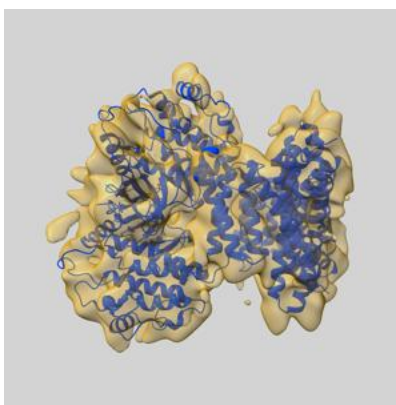
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-20551 and PDB model 6Q0K. Per-residue inclusion information can be found in section 3 on page 6.

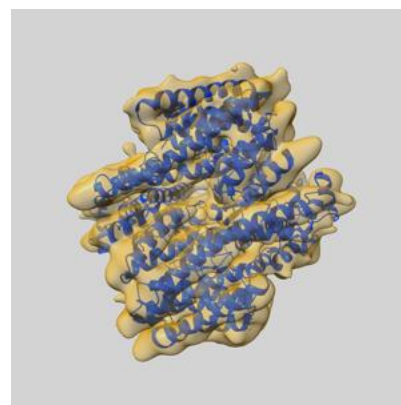
9.1 Map-model overlay [i](#)



X



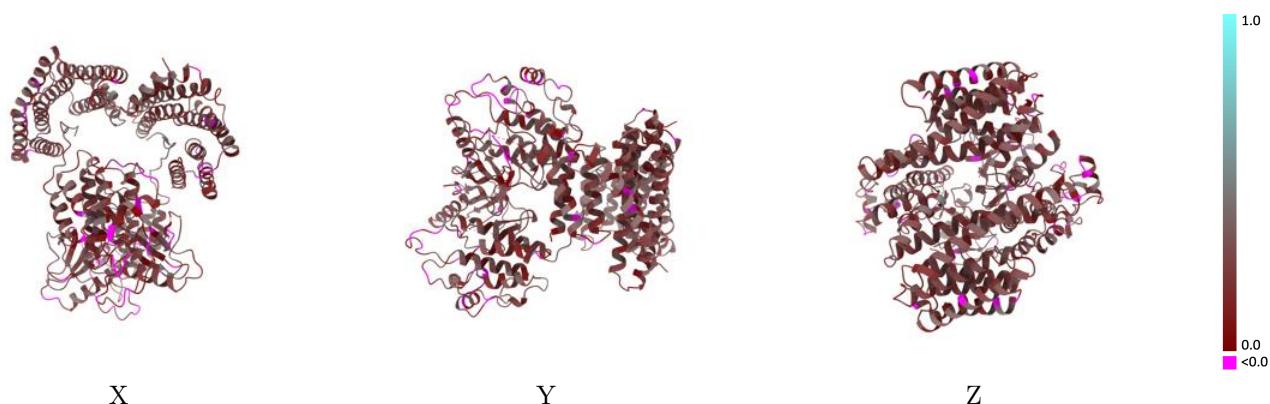
Y



Z

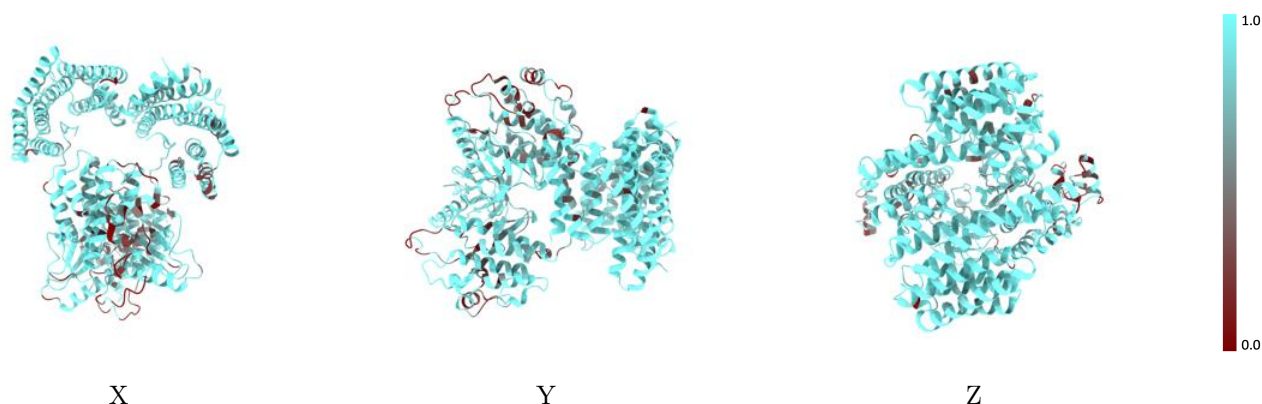
The images above show the 3D surface view of the map at the recommended contour level 0.021 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



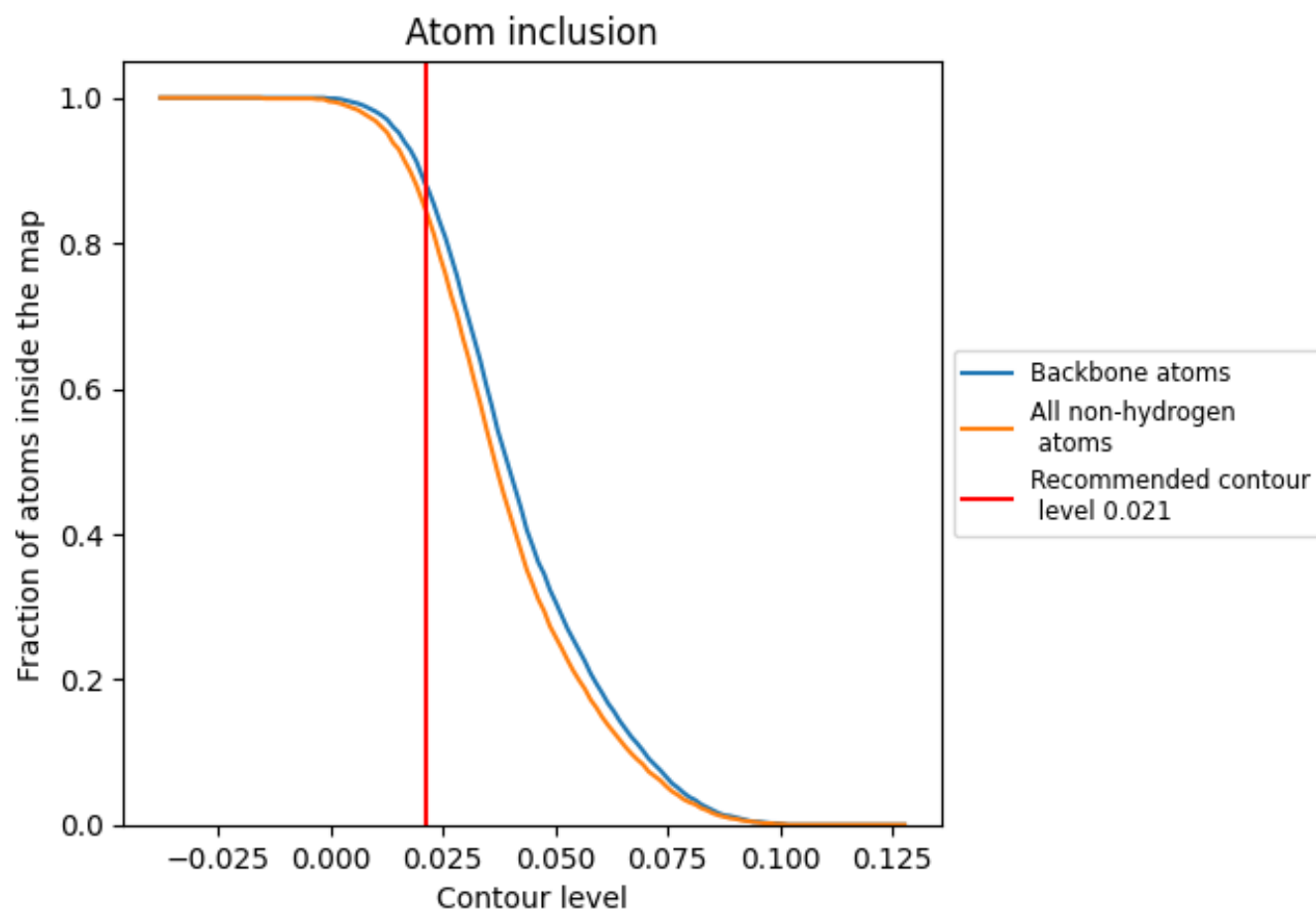
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.021).

9.4 Atom inclusion ⓘ



At the recommended contour level, 88% 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 (0.021) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8470	0.2220
A	0.8010	0.2160
B	0.7590	0.1990
X	0.9180	0.2340
Y	0.9420	0.2430

