



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 11:15 AM UTC

PDB ID : 8UYQ / pdb_00008uyq
EMDB ID : EMD-42817
Title : Consensus olfactory receptor consOR4 bound to 2-methylthiazoline and in complex with mini-Gs trimeric protein
Authors : Billesboelle, C.B.; Del Torrent, C.L.; Manglik, A.
Deposited on : 2023-11-13
Resolution : 3.50 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

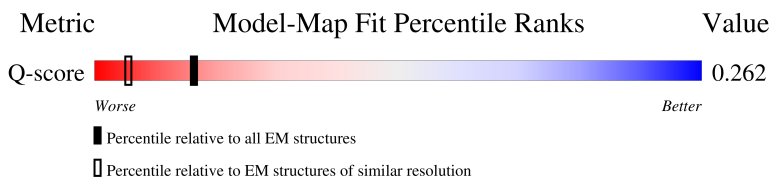
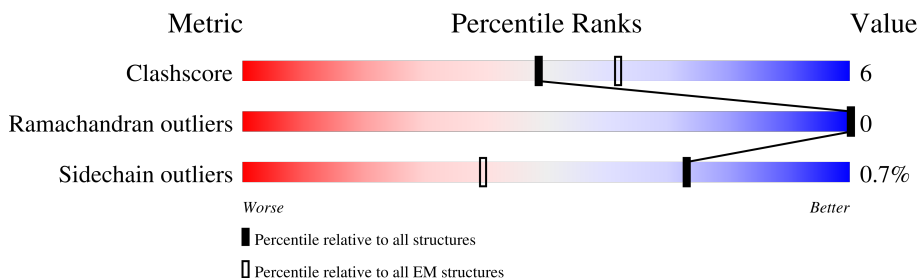
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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	13950 (3.00 - 4.00)

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	321	31% 88% 6% • 5%
2	N	145	20% 73% 14% 13%
3	X	261	25% 69% 13% • 17%
4	Y	370	23% 69% 22% 9%

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Mol	Chain	Length	Quality of chain
5	Z	71	63% 75% 21%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	XVT	A	401	X	-	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16486 atoms, of which 8246 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 consOR4.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	306	Total	C	H	N	O	S	0	0
			4974	1622	2548	382	403	19		

- Molecule 2 is a protein called Nanobody 35.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	N	126	Total	C	H	N	O	S	0	0
			1890	599	929	168	188	6		

- Molecule 3 is a protein called GNAS complex locus.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	X	216	Total	C	H	N	O	S	0	0
			3632	1153	1814	332	326	7		

There are 94 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	-7	GLY	-	expression tag	UNP A0A804HIH4
X	-6	GLY	-	expression tag	UNP A0A804HIH4
X	-5	SER	-	expression tag	UNP A0A804HIH4
X	-4	LEU	-	expression tag	UNP A0A804HIH4
X	-3	GLU	-	expression tag	UNP A0A804HIH4
X	-2	VAL	-	expression tag	UNP A0A804HIH4
X	-1	LEU	-	expression tag	UNP A0A804HIH4
X	0	PHE	-	expression tag	UNP A0A804HIH4
X	1	GLN	-	expression tag	UNP A0A804HIH4
X	2	GLY	-	expression tag	UNP A0A804HIH4
X	3	PRO	-	expression tag	UNP A0A804HIH4
X	4	SER	-	expression tag	UNP A0A804HIH4
X	5	GLY	-	expression tag	UNP A0A804HIH4
X	6	ASN	-	expression tag	UNP A0A804HIH4
X	7	SER	-	expression tag	UNP A0A804HIH4

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Chain	Residue	Modelled	Actual	Comment	Reference
X	8	LYS	-	expression tag	UNP A0A804HIH4
X	9	THR	-	expression tag	UNP A0A804HIH4
X	10	GLU	-	expression tag	UNP A0A804HIH4
X	11	ASP	-	expression tag	UNP A0A804HIH4
X	12	GLN	-	expression tag	UNP A0A804HIH4
X	13	ARG	-	expression tag	UNP A0A804HIH4
X	14	ASN	-	expression tag	UNP A0A804HIH4
X	15	GLU	-	expression tag	UNP A0A804HIH4
X	16	GLU	-	expression tag	UNP A0A804HIH4
X	17	LYS	-	expression tag	UNP A0A804HIH4
X	18	ALA	-	expression tag	UNP A0A804HIH4
X	19	GLN	-	expression tag	UNP A0A804HIH4
X	20	ARG	-	expression tag	UNP A0A804HIH4
X	21	GLU	-	expression tag	UNP A0A804HIH4
X	22	ALA	-	expression tag	UNP A0A804HIH4
X	23	ASN	-	expression tag	UNP A0A804HIH4
X	24	LYS	-	expression tag	UNP A0A804HIH4
X	25	LYS	-	expression tag	UNP A0A804HIH4
X	26	ILE	-	expression tag	UNP A0A804HIH4
X	27	GLU	-	expression tag	UNP A0A804HIH4
X	28	LYS	-	expression tag	UNP A0A804HIH4
X	29	GLN	-	expression tag	UNP A0A804HIH4
X	30	LEU	-	expression tag	UNP A0A804HIH4
X	31	GLN	-	expression tag	UNP A0A804HIH4
X	32	LYS	-	expression tag	UNP A0A804HIH4
X	33	ASP	-	expression tag	UNP A0A804HIH4
X	34	LYS	-	expression tag	UNP A0A804HIH4
X	35	GLN	-	expression tag	UNP A0A804HIH4
X	36	VAL	-	expression tag	UNP A0A804HIH4
X	37	TYR	-	expression tag	UNP A0A804HIH4
X	38	ARG	-	expression tag	UNP A0A804HIH4
X	39	ALA	-	expression tag	UNP A0A804HIH4
X	40	THR	-	expression tag	UNP A0A804HIH4
X	41	HIS	-	expression tag	UNP A0A804HIH4
X	42	ARG	-	expression tag	UNP A0A804HIH4
X	43	LEU	-	expression tag	UNP A0A804HIH4
X	44	LEU	-	expression tag	UNP A0A804HIH4
X	45	LEU	-	expression tag	UNP A0A804HIH4
X	46	LEU	-	expression tag	UNP A0A804HIH4
X	47	GLY	-	expression tag	UNP A0A804HIH4
X	48	ALA	-	expression tag	UNP A0A804HIH4
X	49	ASP	-	expression tag	UNP A0A804HIH4

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Chain	Residue	Modelled	Actual	Comment	Reference
X	50	ASN	-	expression tag	UNP A0A804HIH4
X	51	SER	-	expression tag	UNP A0A804HIH4
X	52	GLY	-	expression tag	UNP A0A804HIH4
X	53	LYS	-	expression tag	UNP A0A804HIH4
X	54	SER	-	expression tag	UNP A0A804HIH4
X	55	THR	-	expression tag	UNP A0A804HIH4
X	56	ILE	-	expression tag	UNP A0A804HIH4
X	57	VAL	-	expression tag	UNP A0A804HIH4
X	58	LYS	-	expression tag	UNP A0A804HIH4
X	59	GLN	-	expression tag	UNP A0A804HIH4
X	60	MET	-	expression tag	UNP A0A804HIH4
X	61	ARG	-	expression tag	UNP A0A804HIH4
X	193	ILE	-	expression tag	UNP A0A804HIH4
X	194	LEU	-	expression tag	UNP A0A804HIH4
X	195	HIS	-	expression tag	UNP A0A804HIH4
X	196	GLY	-	expression tag	UNP A0A804HIH4
X	197	GLY	-	expression tag	UNP A0A804HIH4
X	198	SER	-	expression tag	UNP A0A804HIH4
X	199	GLY	-	expression tag	UNP A0A804HIH4
X	200	GLY	-	expression tag	UNP A0A804HIH4
X	201	SER	-	expression tag	UNP A0A804HIH4
X	202	GLY	-	expression tag	UNP A0A804HIH4
X	203	GLY	-	expression tag	UNP A0A804HIH4
X	249	ASP	ALA	conflict	UNP A0A804HIH4
X	252	ASP	SER	conflict	UNP A0A804HIH4
X	?	-	MET	deletion	UNP A0A804HIH4
X	?	-	VAL	deletion	UNP A0A804HIH4
X	?	-	ILE	deletion	UNP A0A804HIH4
X	?	-	ARG	deletion	UNP A0A804HIH4
X	?	-	GLU	deletion	UNP A0A804HIH4
X	?	-	ASP	deletion	UNP A0A804HIH4
X	?	-	ASN	deletion	UNP A0A804HIH4
X	?	-	GLN	deletion	UNP A0A804HIH4
X	?	-	THR	deletion	UNP A0A804HIH4
X	?	-	ASN	deletion	UNP A0A804HIH4
X	372	ALA	ILE	conflict	UNP A0A804HIH4
X	375	ILE	VAL	conflict	UNP A0A804HIH4

- Molecule 4 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	Y	338	Total	C	H	N	O	S	0	0
			5106	1604	2506	467	508	21		

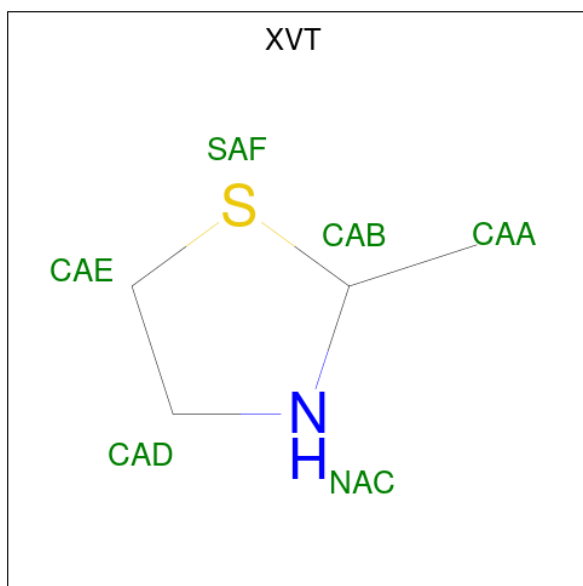
There are 31 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Y	-29	MET	-	initiating methionine	UNP P62873
Y	-28	HIS	-	expression tag	UNP P62873
Y	-27	HIS	-	expression tag	UNP P62873
Y	-26	HIS	-	expression tag	UNP P62873
Y	-25	HIS	-	expression tag	UNP P62873
Y	-24	HIS	-	expression tag	UNP P62873
Y	-23	HIS	-	expression tag	UNP P62873
Y	-22	LEU	-	expression tag	UNP P62873
Y	-21	GLU	-	expression tag	UNP P62873
Y	-20	VAL	-	expression tag	UNP P62873
Y	-19	LEU	-	expression tag	UNP P62873
Y	-18	PHE	-	expression tag	UNP P62873
Y	-17	GLN	-	expression tag	UNP P62873
Y	-16	GLY	-	expression tag	UNP P62873
Y	-15	PRO	-	expression tag	UNP P62873
Y	-14	GLU	-	expression tag	UNP P62873
Y	-13	ASP	-	expression tag	UNP P62873
Y	-12	GLN	-	expression tag	UNP P62873
Y	-11	VAL	-	expression tag	UNP P62873
Y	-10	ASP	-	expression tag	UNP P62873
Y	-9	PRO	-	expression tag	UNP P62873
Y	-8	ARG	-	expression tag	UNP P62873
Y	-7	LEU	-	expression tag	UNP P62873
Y	-6	ILE	-	expression tag	UNP P62873
Y	-5	ASP	-	expression tag	UNP P62873
Y	-4	GLY	-	expression tag	UNP P62873
Y	-3	LYS	-	expression tag	UNP P62873
Y	-2	GLY	-	expression tag	UNP P62873
Y	-1	SER	-	expression tag	UNP P62873
Y	0	SER	-	expression tag	UNP P62873
Y	1	GLY	-	expression tag	UNP P62873

- Molecule 5 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	Z	56	Total	C	H	N	O	S	0	0
			871	269	442	76	81	3		

- Molecule 6 is (2R)-2-methyl-1,3-thiazolidine (CCD ID: XVT) (formula: C₄H₉NS) (labeled as "Ligand of Interest" by depositor).

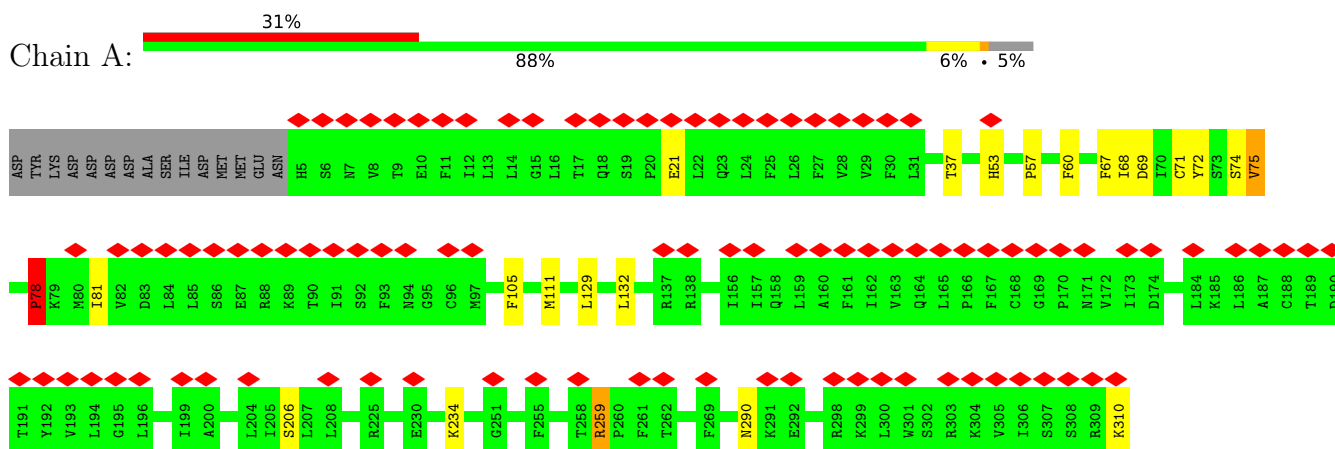


Mol	Chain	Residues	Atoms					AltConf
6	A	1	Total	C	H	N	S	0
			13	4	7	1	1	

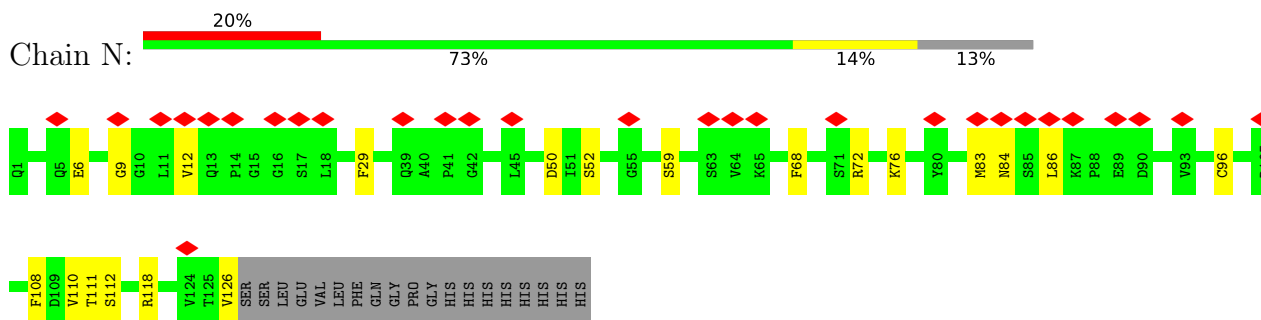
3 Residue-property plots [i](#)

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.

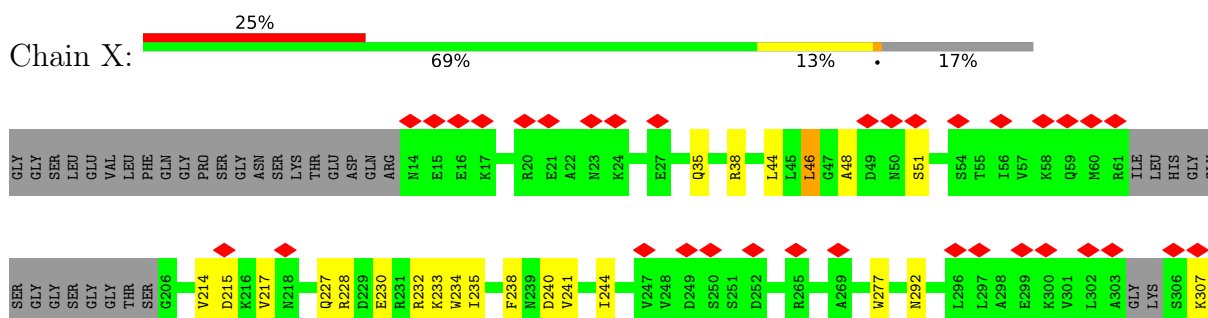
• Molecule 1: consOR4

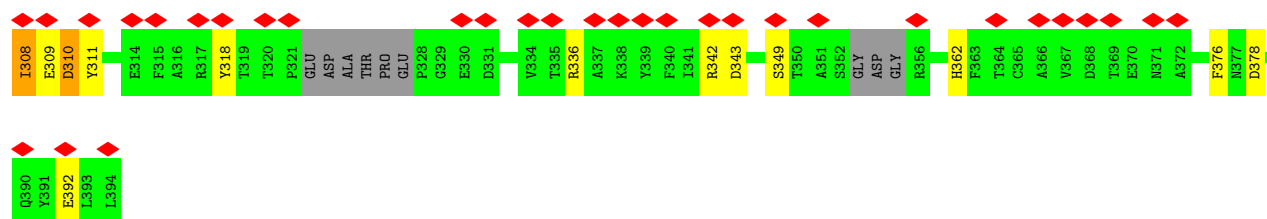


• Molecule 2: Nanobody 35

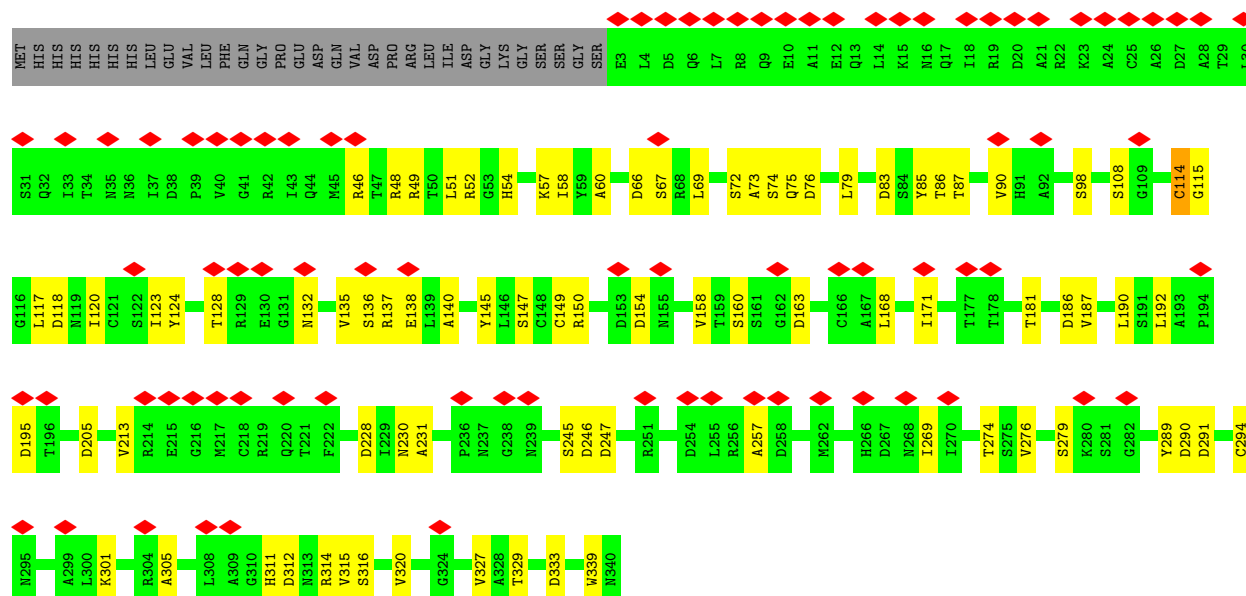


• Molecule 3: GNAS complex locus

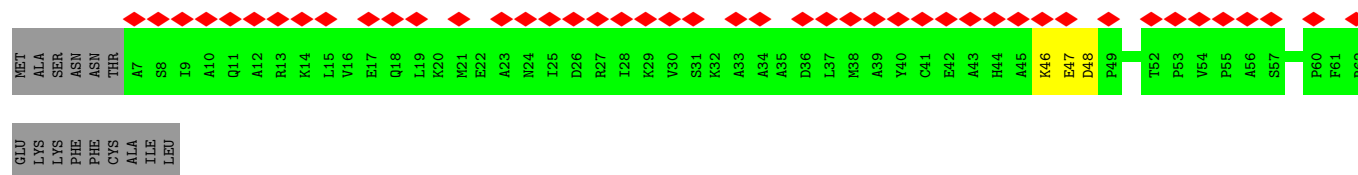
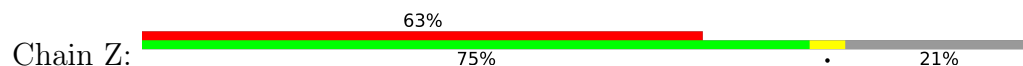




- Molecule 4: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1



- Molecule 5: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	318570	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.570	Depositor
Minimum map value	-1.058	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.029	Depositor
Recommended contour level	0.15	Depositor
Map size (Å)	246.528, 246.528, 246.528	wwPDB
Map dimensions	288, 288, 288	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.856, 0.856, 0.856	Depositor

5 Model quality

5.1 Standard geometry

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

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	1.11	11/2488 (0.4%)	1.38	9/3389 (0.3%)
2	N	0.16	0/981	0.74	8/1329 (0.6%)
3	X	0.25	0/1851	0.69	12/2483 (0.5%)
4	Y	0.17	0/2647	0.99	34/3589 (0.9%)
5	Z	0.16	0/435	0.33	0/587
All	All	0.63	11/8402 (0.1%)	1.02	63/11377 (0.6%)

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	X	1	0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	75	VAL	C-O	16.07	1.42	1.24
1	A	72	TYR	C-O	13.35	1.39	1.24
1	A	81	ILE	C-O	9.30	1.35	1.24
1	A	21	GLU	C-O	7.91	1.33	1.24
1	A	67	PHE	C-O	7.48	1.32	1.24
1	A	68	ILE	C-O	5.96	1.31	1.24
1	A	78	PRO	C-O	5.60	1.31	1.24
1	A	69	ASP	C-O	5.57	1.30	1.24
1	A	57	PRO	C-O	5.23	1.30	1.24
1	A	37	THR	C-O	5.18	1.30	1.24
1	A	60	PHE	C-O	5.10	1.30	1.24

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	Y	291	ASP	CA-C-N	11.57	143.64	124.31
4	Y	291	ASP	C-N-CA	11.57	143.64	124.31
4	Y	333	ASP	CA-C-N	11.53	141.68	123.93
4	Y	333	ASP	C-N-CA	11.53	141.68	123.93
4	Y	52	ARG	CA-C-N	10.58	142.15	121.41
4	Y	52	ARG	C-N-CA	10.58	142.15	121.41
4	Y	181	THR	CA-C-N	10.31	141.62	121.41
4	Y	181	THR	C-N-CA	10.31	141.62	121.41
1	A	78	PRO	CA-N-CD	-10.22	97.69	112.00
4	Y	140	ALA	CA-C-N	10.16	141.33	121.41
4	Y	140	ALA	C-N-CA	10.16	141.33	121.41
4	Y	115	GLY	CA-C-N	10.12	141.24	121.41
4	Y	115	GLY	C-N-CA	10.12	141.24	121.41
3	X	215	ASP	CA-C-N	9.22	137.52	123.05
3	X	215	ASP	C-N-CA	9.22	137.52	123.05
4	Y	301	LYS	CA-C-N	9.16	139.12	123.91
4	Y	301	LYS	C-N-CA	9.16	139.12	123.91
4	Y	66	ASP	CA-C-N	9.06	137.28	123.05
4	Y	66	ASP	C-N-CA	9.06	137.28	123.05
4	Y	257	ALA	CA-C-N	8.96	135.54	122.63
4	Y	257	ALA	C-N-CA	8.96	135.54	122.63
4	Y	118	ASP	CA-C-N	8.82	136.90	123.05
4	Y	118	ASP	C-N-CA	8.82	136.90	123.05
4	Y	163	ASP	CA-C-N	8.78	135.14	122.40
4	Y	163	ASP	C-N-CA	8.78	135.14	122.40
4	Y	247	ASP	CA-C-N	8.70	135.02	122.40
4	Y	247	ASP	C-N-CA	8.70	135.02	122.40
2	N	9	GLY	CA-C-N	8.67	132.88	120.56
2	N	9	GLY	C-N-CA	8.67	132.88	120.56
4	Y	305	ALA	CA-C-N	8.58	134.51	121.04
4	Y	305	ALA	C-N-CA	8.58	134.51	121.04
2	N	108	PHE	CA-C-N	8.58	136.52	123.05
2	N	108	PHE	C-N-CA	8.58	136.52	123.05
3	X	292	ASN	CA-C-N	8.47	134.14	122.19
3	X	292	ASN	C-N-CA	8.47	134.14	122.19
3	X	214	VAL	CA-C-N	8.33	133.98	122.07
3	X	214	VAL	C-N-CA	8.33	133.98	122.07
4	Y	114	CYS	CA-C-N	8.26	134.24	120.86
4	Y	114	CYS	C-N-CA	8.26	134.24	120.86
4	Y	205	ASP	CA-C-N	8.23	133.80	122.19
4	Y	205	ASP	C-N-CA	8.23	133.80	122.19
4	Y	195	ASP	CA-C-N	8.16	134.24	122.40
4	Y	195	ASP	C-N-CA	8.16	134.24	122.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	84	ASN	CA-C-N	8.00	134.44	122.68
2	N	84	ASN	C-N-CA	8.00	134.44	122.68
2	N	76	LYS	CA-C-N	7.54	134.32	122.93
2	N	76	LYS	C-N-CA	7.54	134.32	122.93
4	Y	86	THR	CA-C-N	7.25	134.43	123.05
4	Y	86	THR	C-N-CA	7.25	134.43	123.05
3	X	46	LEU	CA-C-N	7.09	132.35	120.86
3	X	46	LEU	C-N-CA	7.09	132.35	120.86
1	A	72	TYR	N-CA-CB	-6.86	100.07	110.01
3	X	308	ILE	CA-C-O	-5.84	114.32	120.57
1	A	67	PHE	O-C-N	-5.81	115.96	122.12
1	A	259	ARG	NE-CZ-NH2	5.54	124.19	119.20
1	A	105	PHE	CA-CB-CG	5.53	119.33	113.80
3	X	308	ILE	N-CA-C	-5.32	104.40	111.05
3	X	308	ILE	CA-CB-CG1	5.29	119.39	110.40
3	X	310	ASP	CB-CA-C	5.19	119.41	110.79
1	A	78	PRO	CA-C-N	-5.08	113.83	120.44
1	A	78	PRO	C-N-CA	-5.08	113.83	120.44
1	A	72	TYR	CA-CB-CG	5.07	123.03	113.90
1	A	71	CYS	N-CA-C	5.05	116.78	111.28

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	X	308	ILE	CB

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	2426	2548	2544	27	0
2	N	961	929	932	9	0
3	X	1818	1814	1810	40	0
4	Y	2600	2506	2505	47	0
5	Z	429	442	441	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	6	7	0	0	0
All	All	8240	8246	8232	95	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 6.

All (95) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:LEU:HD22	3:X:38:ARG:HH12	1.30	0.96
1:A:290:ASN:O	3:X:392:GLU:CD	2.11	0.93
1:A:129:LEU:CD1	3:X:376:PHE:CE1	2.58	0.87
1:A:290:ASN:HA	3:X:392:GLU:HG3	1.58	0.85
1:A:132:LEU:HD22	3:X:38:ARG:NH1	1.92	0.84
4:Y:320:VAL:HG22	4:Y:327:VAL:HG22	1.71	0.73
3:X:228:ARG:NH1	4:Y:186:ASP:OD1	2.22	0.72
4:Y:147:SER:OG	4:Y:187:VAL:O	2.07	0.70
3:X:240:ASP:OD1	4:Y:57:LYS:NZ	2.26	0.69
4:Y:274:THR:OG1	4:Y:315:VAL:O	2.08	0.68
4:Y:137:ARG:NH1	4:Y:171:ILE:O	2.28	0.67
4:Y:74:SER:OG	4:Y:76:ASP:OD1	2.10	0.67
3:X:310:ASP:OD1	3:X:311:TYR:CD2	2.48	0.67
4:Y:79:LEU:HD11	4:Y:114:CYS:HB3	1.77	0.67
1:A:290:ASN:CA	3:X:392:GLU:HG3	2.25	0.66
4:Y:160:SER:HB3	4:Y:190:LEU:HD23	1.80	0.64
1:A:132:LEU:HD13	3:X:38:ARG:HH11	1.63	0.64
3:X:318:TYR:OH	3:X:343:ASP:OD2	2.13	0.63
1:A:129:LEU:HD12	3:X:376:PHE:CZ	2.33	0.63
1:A:129:LEU:CD1	3:X:376:PHE:CZ	2.82	0.62
1:A:290:ASN:O	3:X:392:GLU:OE2	2.17	0.61
3:X:234:TRP:HB2	4:Y:117:LEU:HD21	1.82	0.61
1:A:290:ASN:O	3:X:392:GLU:CG	2.49	0.61
4:Y:294:CYS:SG	4:Y:315:VAL:HG11	2.41	0.61
1:A:310:LYS:HE2	4:Y:48:ARG:HH21	1.67	0.60
1:A:129:LEU:HD13	3:X:217:VAL:HG21	1.84	0.59
4:Y:311:HIS:NE2	4:Y:329:THR:O	2.36	0.58
5:Z:46:LYS:NZ	5:Z:47:GLU:OE2	2.36	0.58
3:X:227:GLN:OE1	4:Y:145:TYR:N	2.37	0.57
2:N:52:SER:O	2:N:72:ARG:NH1	2.38	0.57
3:X:44:LEU:HD22	3:X:238:PHE:CB	2.36	0.56
4:Y:128:THR:OG1	4:Y:132:ASN:O	2.17	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:LEU:HD13	3:X:376:PHE:CE1	2.41	0.55
4:Y:54:HIS:HE2	4:Y:72:SER:HG	1.52	0.55
4:Y:51:LEU:HD23	4:Y:87:THR:HG23	1.88	0.55
1:A:111:MET:HE1	1:A:206:SER:OG	2.07	0.54
1:A:290:ASN:O	3:X:392:GLU:HG3	2.08	0.54
4:Y:60:ALA:HB3	4:Y:73:ALA:HB3	1.90	0.54
1:A:129:LEU:HD12	3:X:376:PHE:CE1	2.43	0.54
2:N:29:PHE:O	2:N:72:ARG:NH2	2.40	0.53
4:Y:75:GLN:O	4:Y:98:SER:OG	2.26	0.53
4:Y:329:THR:OG1	4:Y:339:TRP:NE1	2.41	0.53
3:X:44:LEU:HD22	3:X:238:PHE:HB2	1.91	0.52
3:X:230:GLU:OE1	3:X:232:ARG:NH1	2.42	0.52
4:Y:54:HIS:NE2	4:Y:72:SER:OG	2.41	0.52
3:X:318:TYR:O	3:X:336:ARG:NH1	2.43	0.52
4:Y:108:SER:OG	4:Y:154:ASP:OD1	2.28	0.52
2:N:83:MET:HE2	2:N:86:LEU:HD21	1.92	0.52
3:X:342:ARG:NH2	3:X:343:ASP:OD1	2.43	0.51
1:A:310:LYS:HD3	4:Y:46:ARG:HH11	1.75	0.51
4:Y:58:ILE:O	4:Y:316:SER:OG	2.29	0.51
1:A:290:ASN:C	3:X:392:GLU:HG3	2.36	0.50
3:X:48:ALA:N	3:X:51:SER:OG	2.44	0.50
3:X:46:LEU:HD11	3:X:235:ILE:HD11	1.93	0.50
1:A:310:LYS:HE2	4:Y:48:ARG:NH2	2.27	0.50
3:X:362:HIS:NE2	3:X:378:ASP:OD2	2.37	0.48
4:Y:69:LEU:HD22	4:Y:90:VAL:HG21	1.94	0.48
1:A:132:LEU:HD13	3:X:38:ARG:NH1	2.27	0.48
1:A:129:LEU:HD11	3:X:376:PHE:CE1	2.46	0.47
4:Y:269:ILE:HG21	4:Y:289:TYR:CE2	2.49	0.47
4:Y:51:LEU:CD2	4:Y:87:THR:HG23	2.44	0.46
2:N:68:PHE:CZ	2:N:83:MET:HE3	2.50	0.46
3:X:35:GLN:OE1	3:X:38:ARG:NH2	2.49	0.46
3:X:44:LEU:HD23	3:X:244:ILE:HG12	1.96	0.46
4:Y:67:SER:O	4:Y:67:SER:OG	2.26	0.46
4:Y:228:ASP:C	4:Y:245:SER:HG	2.18	0.46
4:Y:290:ASP:OD1	4:Y:314:ARG:NE	2.47	0.46
2:N:50:ASP:OD1	2:N:59:SER:OG	2.33	0.45
1:A:310:LYS:HD3	4:Y:46:ARG:NH1	2.31	0.45
1:A:74:SER:O	1:A:78:PRO:HD2	2.17	0.45
4:Y:124:TYR:CE2	4:Y:135:VAL:HG22	2.52	0.45
1:A:74:SER:O	1:A:75:VAL:C	2.58	0.44
2:N:110:VAL:HG23	2:N:111:THR:HG23	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:Y:123:ILE:O	4:Y:136:SER:N	2.50	0.44
4:Y:168:LEU:HD22	4:Y:213:VAL:HG13	2.00	0.44
3:X:307:LYS:O	3:X:310:ASP:OD1	2.37	0.43
3:X:362:HIS:HE2	3:X:378:ASP:CG	2.22	0.43
4:Y:279:SER:HG	5:Z:48:ASP:CG	2.26	0.43
3:X:277:TRP:NE1	3:X:349:SER:O	2.53	0.42
4:Y:149:CYS:O	4:Y:150:ARG:NH1	2.50	0.42
4:Y:158:VAL:CG1	4:Y:192:LEU:HD21	2.50	0.42
3:X:238:PHE:O	3:X:241:VAL:HG22	2.20	0.42
1:A:310:LYS:NZ	4:Y:46:ARG:NE	2.63	0.41
4:Y:83:ASP:O	4:Y:87:THR:N	2.52	0.41
4:Y:312:ASP:OD1	4:Y:312:ASP:N	2.52	0.41
4:Y:245:SER:OG	4:Y:246:ASP:N	2.53	0.41
2:N:6:GLU:OE2	2:N:96:CYS:N	2.49	0.41
4:Y:69:LEU:CD2	4:Y:90:VAL:HG21	2.51	0.41
2:N:112:SER:O	2:N:118:ARG:NH1	2.52	0.41
4:Y:231:ALA:HB1	4:Y:276:VAL:HG22	2.03	0.40
4:Y:49:ARG:NH1	4:Y:85:TYR:O	2.53	0.40
2:N:12:VAL:O	2:N:126:VAL:HG12	2.21	0.40
4:Y:120:ILE:HG21	4:Y:138:GLU:OE1	2.22	0.40
1:A:129:LEU:HD13	3:X:376:PHE:HE1	1.82	0.40
3:X:233:LYS:NZ	4:Y:230:ASN:OD1	2.25	0.40

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	
1	A	304/321 (95%)	290 (95%)	14 (5%)	0	100	100
2	N	124/145 (86%)	116 (94%)	8 (6%)	0	100	100
3	X	206/261 (79%)	204 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	Y	336/370 (91%)	323 (96%)	13 (4%)	0	100	100
5	Z	54/71 (76%)	53 (98%)	1 (2%)	0	100	100
All	All	1024/1168 (88%)	986 (96%)	38 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	279/293 (95%)	275 (99%)	4 (1%)	59	71
2	N	104/121 (86%)	104 (100%)	0	100	100
3	X	197/228 (86%)	195 (99%)	2 (1%)	68	75
4	Y	281/309 (91%)	281 (100%)	0	100	100
5	Z	45/58 (78%)	45 (100%)	0	100	100
All	All	906/1009 (90%)	900 (99%)	6 (1%)	73	78

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	HIS
1	A	78	PRO
1	A	234	LYS
1	A	259	ARG
3	X	308	ILE
3	X	309	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	171	ASN
3	X	387	HIS

5.3.3 RNA [i](#)

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](#)

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	$\# Z > 2$
6	XVT	A	401	-	4,6,6	2.45	1 (25%)	3,7,7	6.02	2 (66%)

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
6	XVT	A	401	-	1/1/1/2	-	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	401	XVT	CAB-SAF	-4.82	1.74	1.84

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	401	XVT	CAA-CAB-NAC	9.49	124.87	111.81
6	A	401	XVT	CAD-CAE-SAF	-4.32	98.99	106.16

All (1) chirality outliers are listed below:

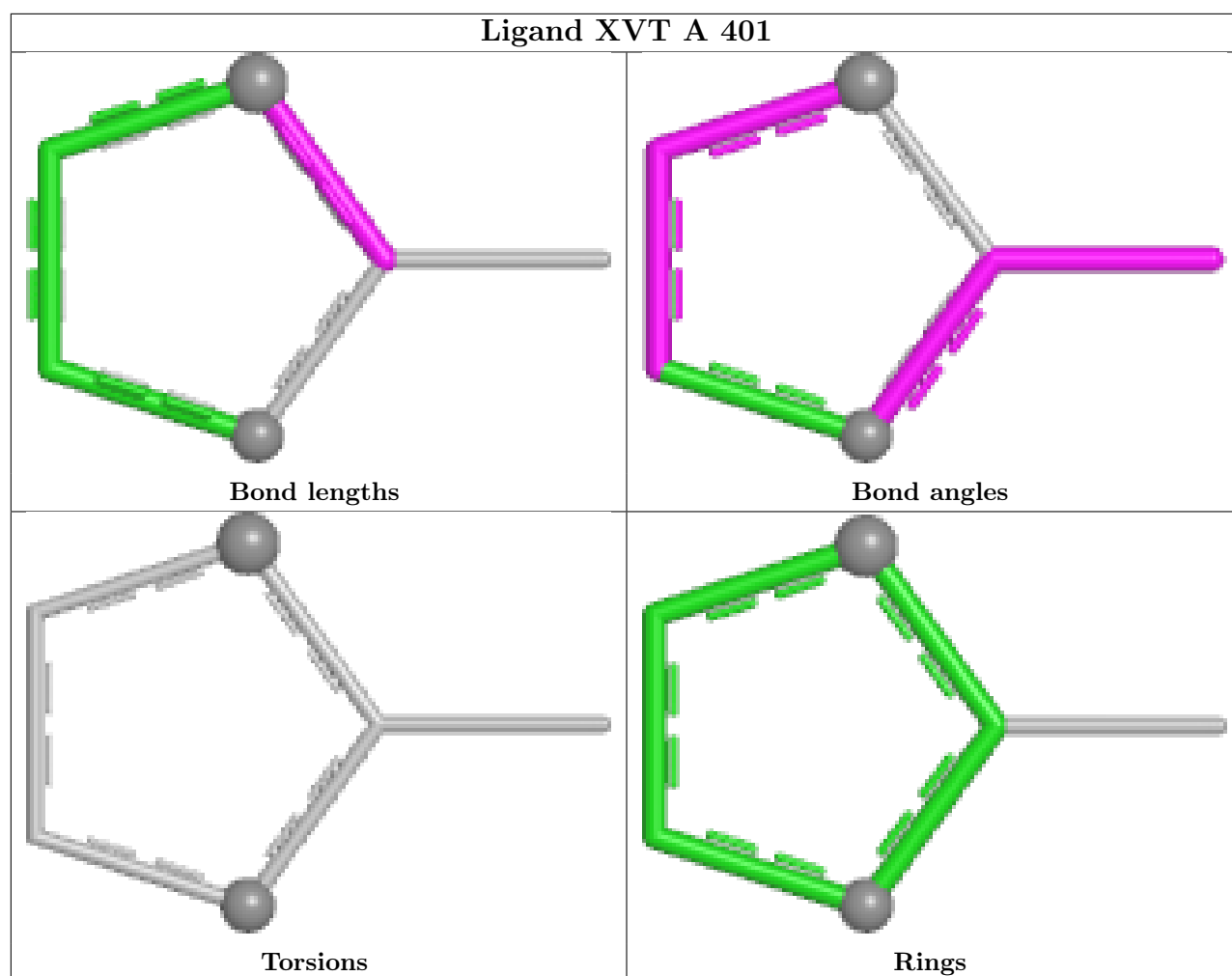
Mol	Chain	Res	Type	Atom
6	A	401	XVT	CAB

There are no torsion outliers.

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 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.

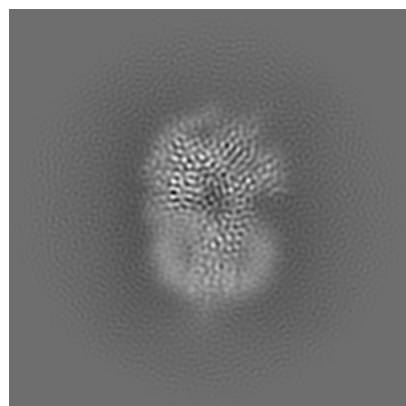
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-42817. These allow visual inspection of the internal detail of the map and identification of artifacts.

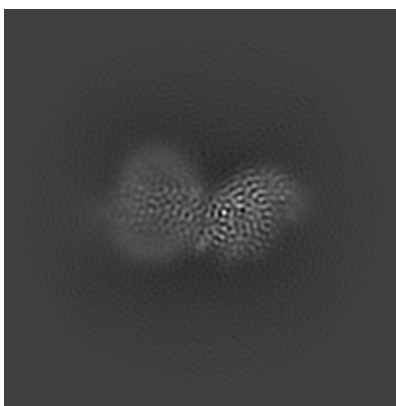
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

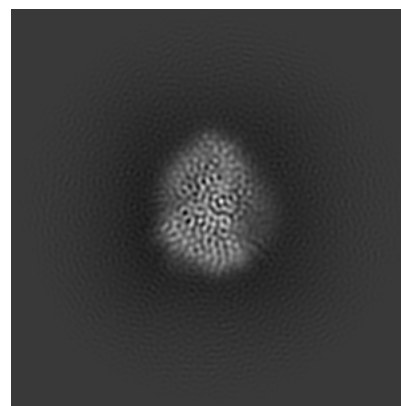
6.1.1 Primary map



X

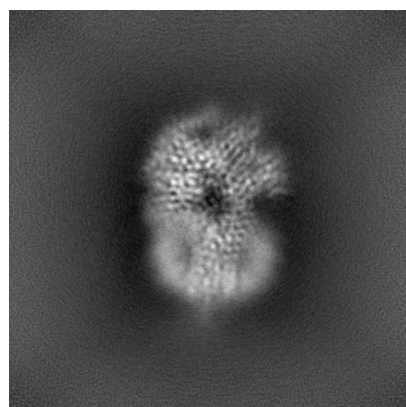


Y

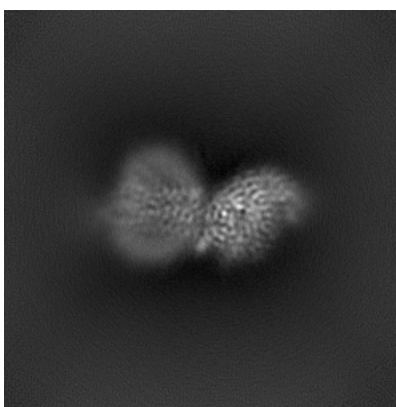


Z

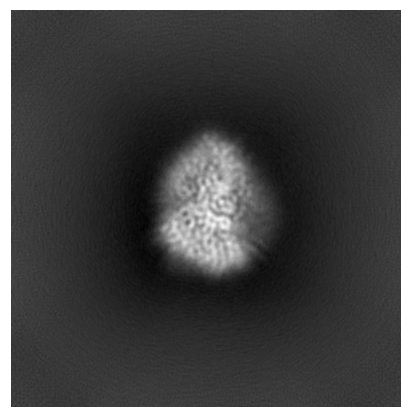
6.1.2 Raw map



X



Y

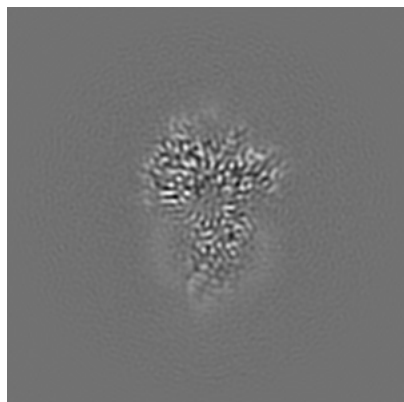


Z

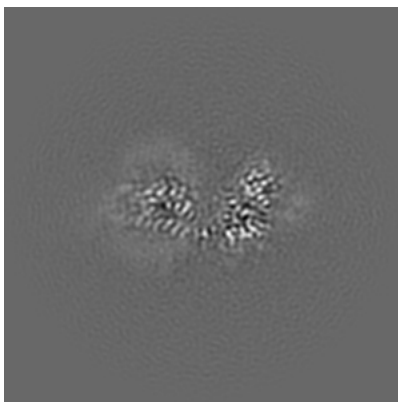
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

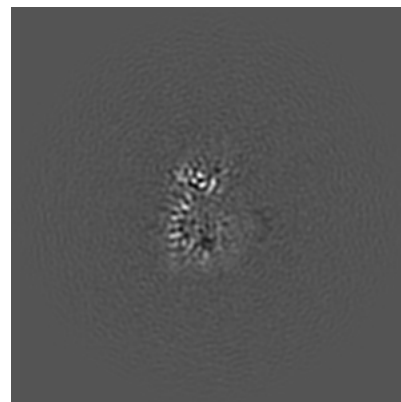
6.2.1 Primary map



X Index: 144

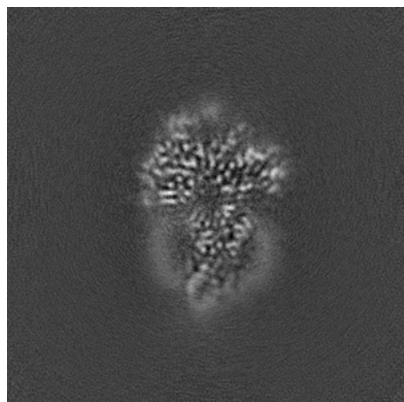


Y Index: 144

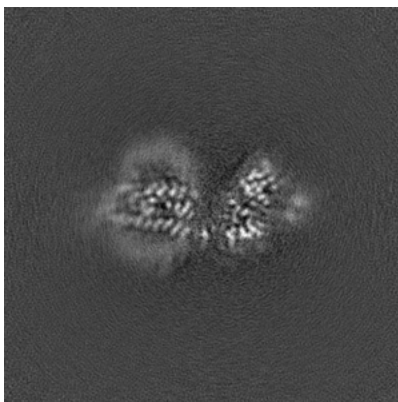


Z Index: 144

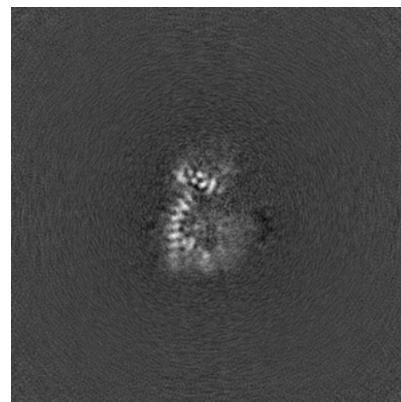
6.2.2 Raw map



X Index: 144



Y Index: 144

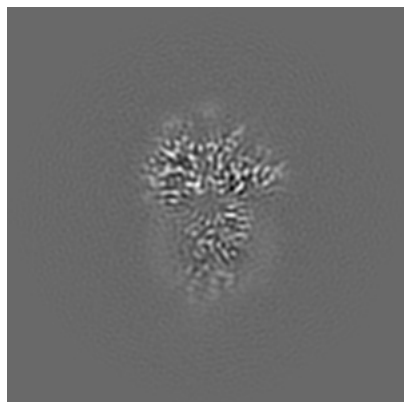


Z Index: 144

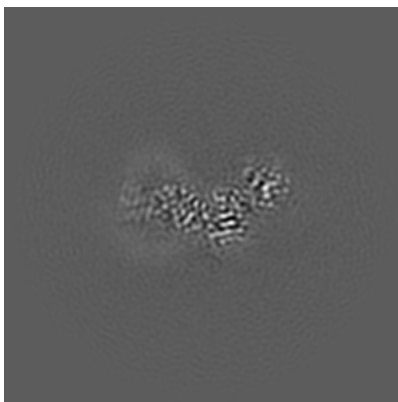
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

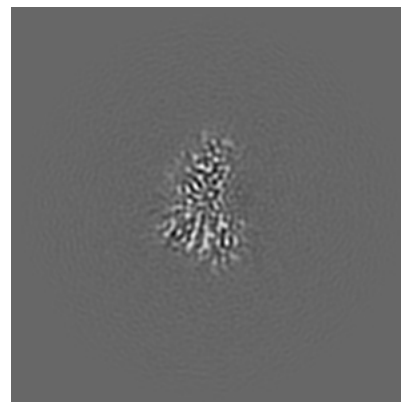
6.3.1 Primary map



X Index: 146

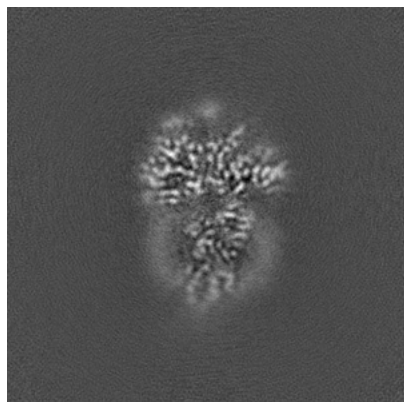


Y Index: 159

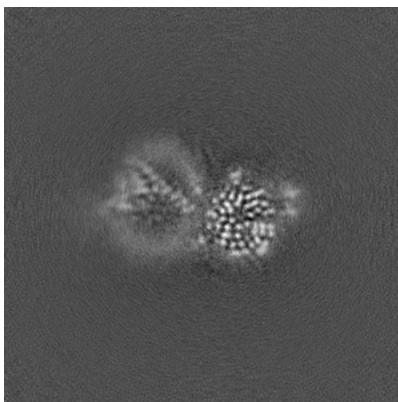


Z Index: 164

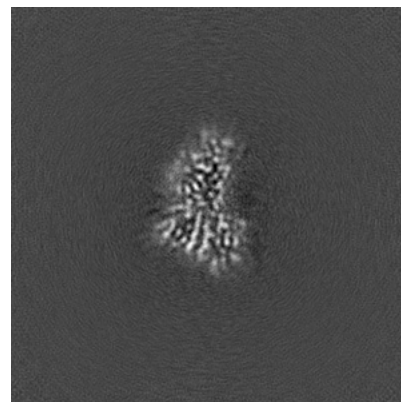
6.3.2 Raw map



X Index: 146



Y Index: 132

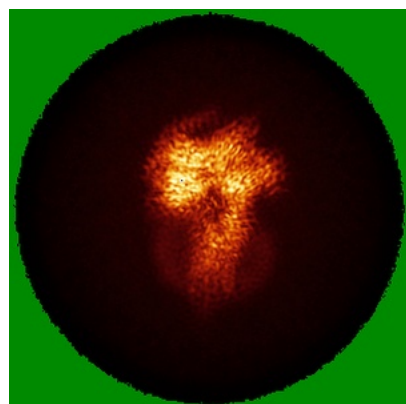


Z Index: 164

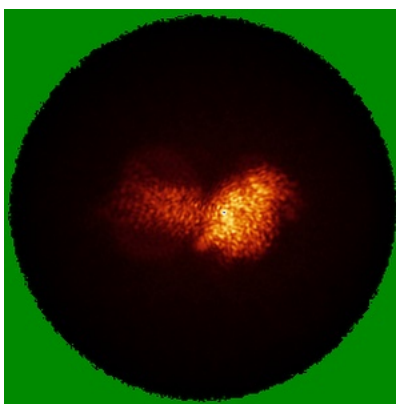
The images above show the largest variance slices of the map in three orthogonal directions.

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

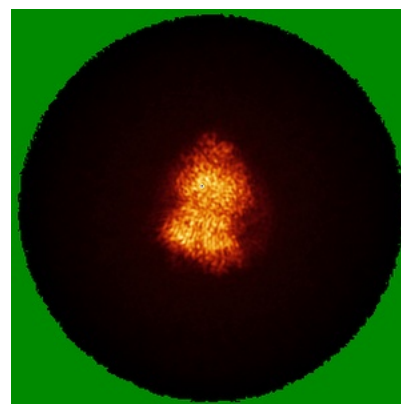
6.4.1 Primary map



X

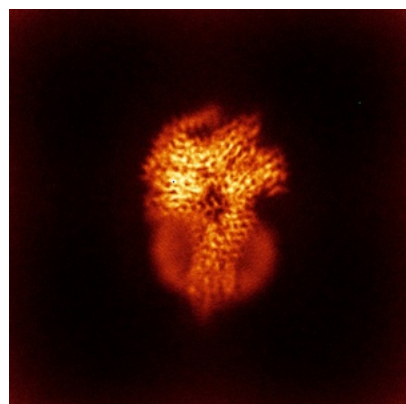


Y

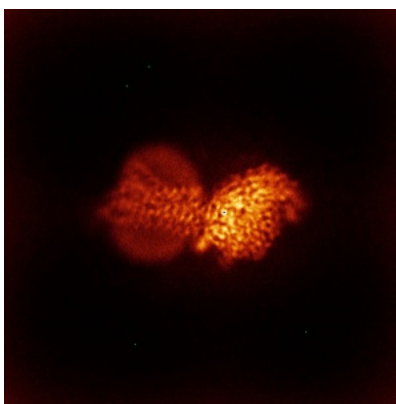


Z

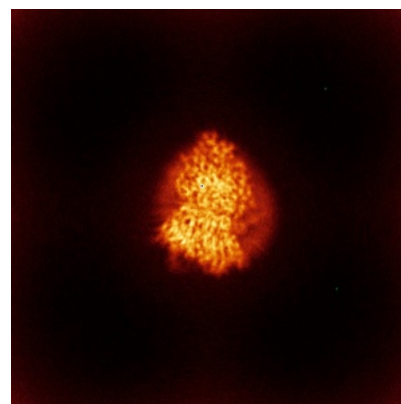
6.4.2 Raw map



X



Y

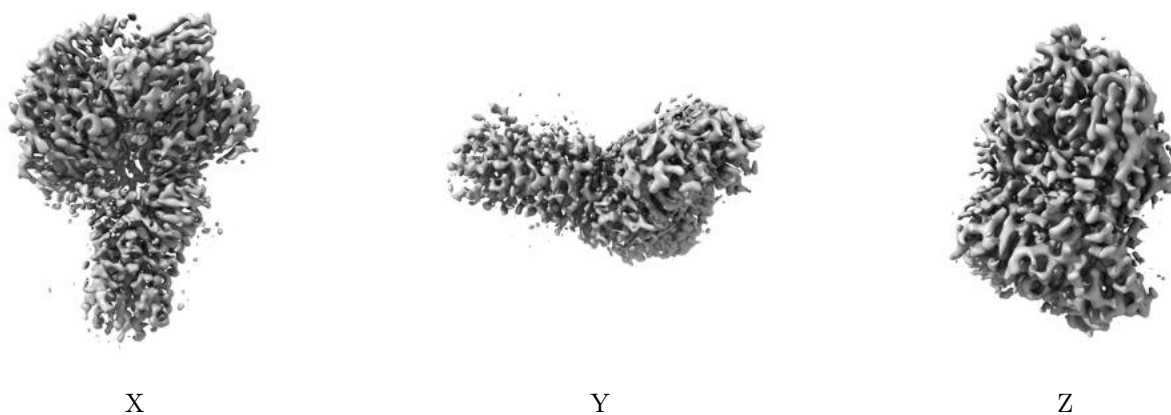


Z

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

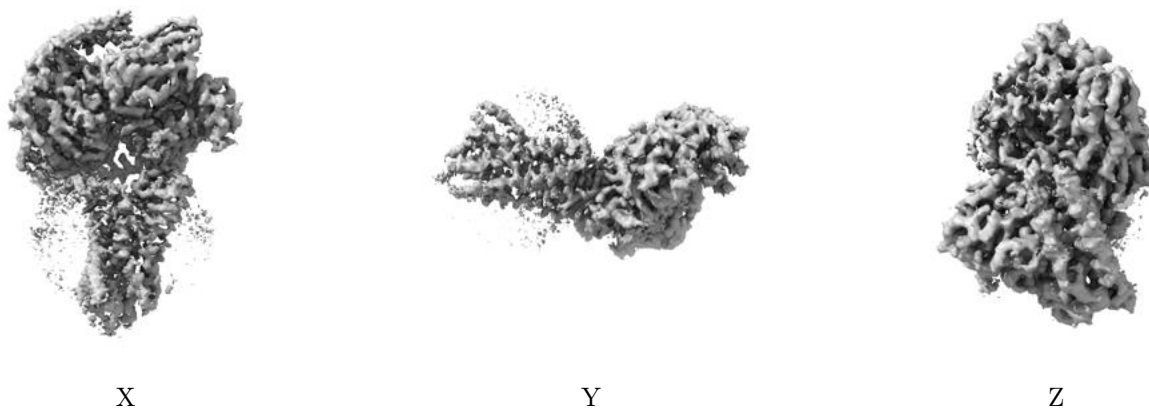
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.15. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

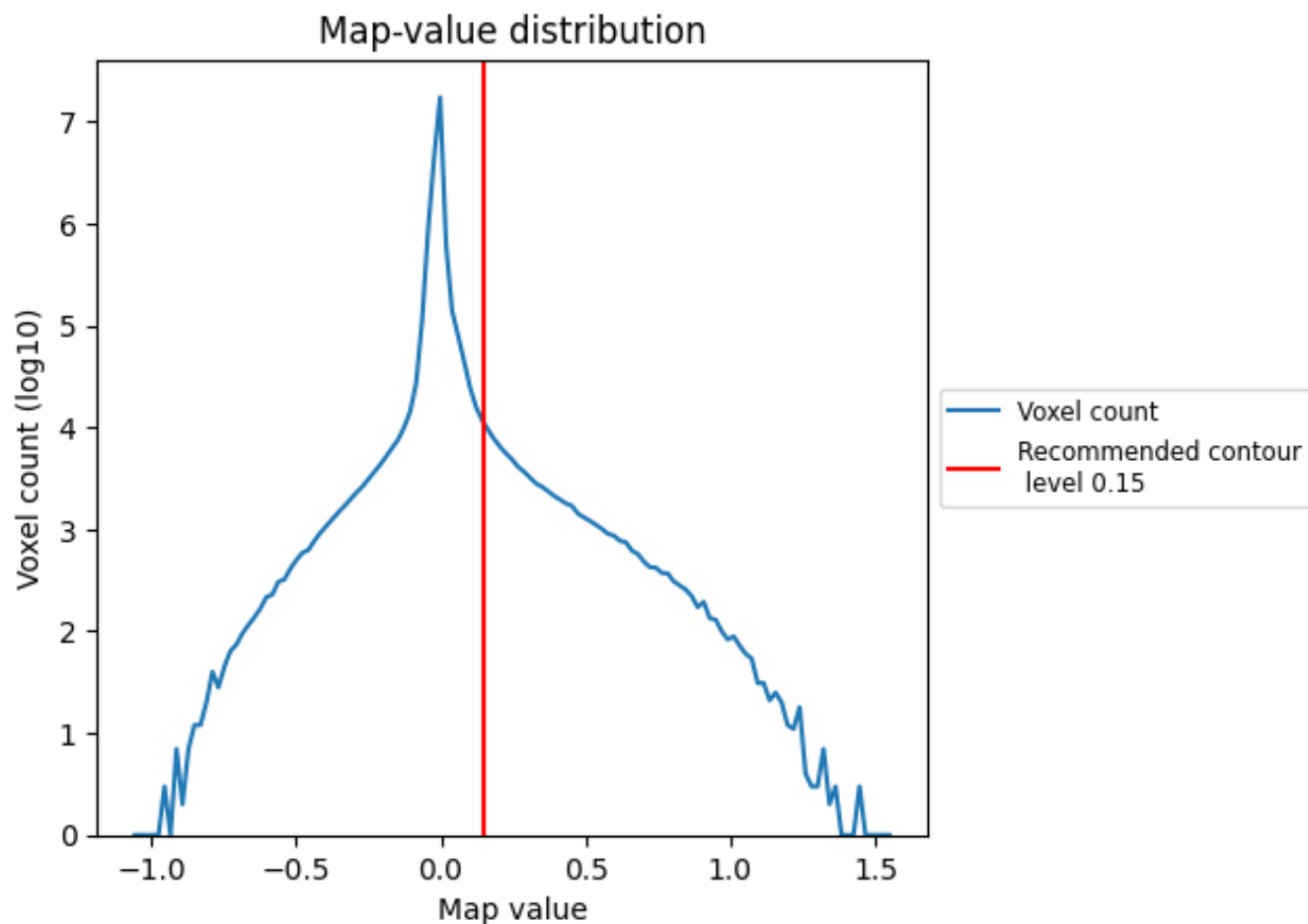
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

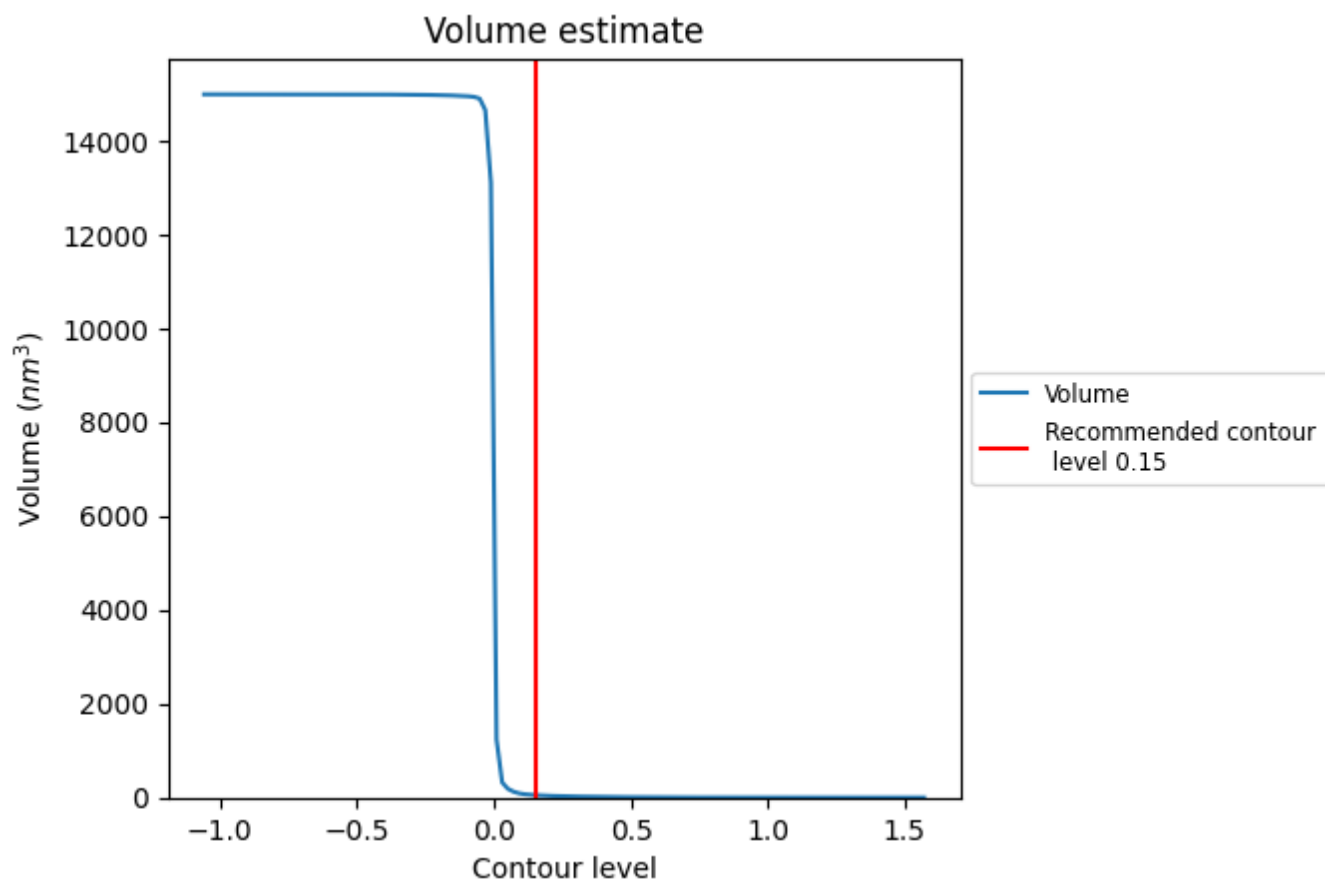
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

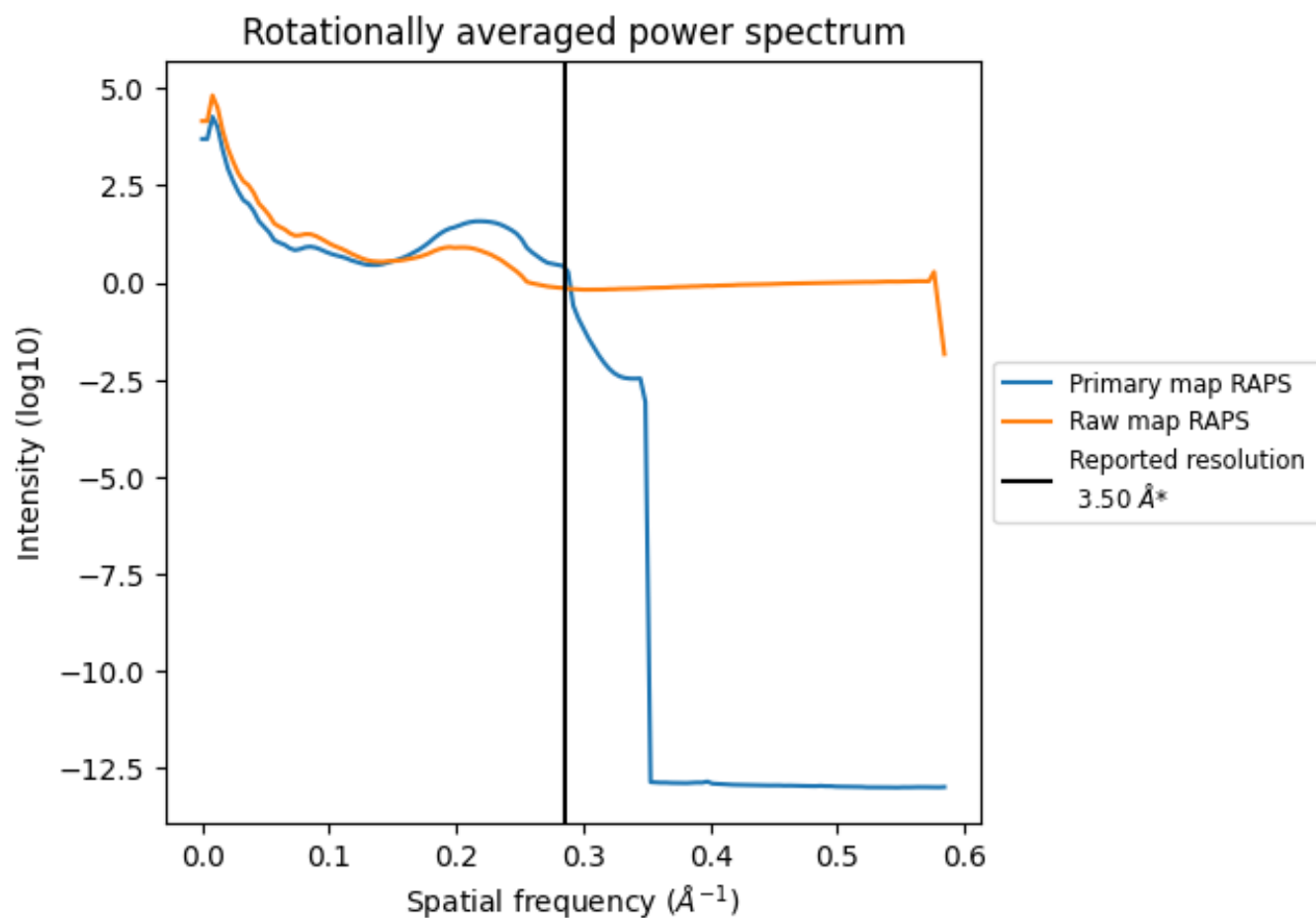
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 52 nm³; this corresponds to an approximate mass of 47 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 ⓘ

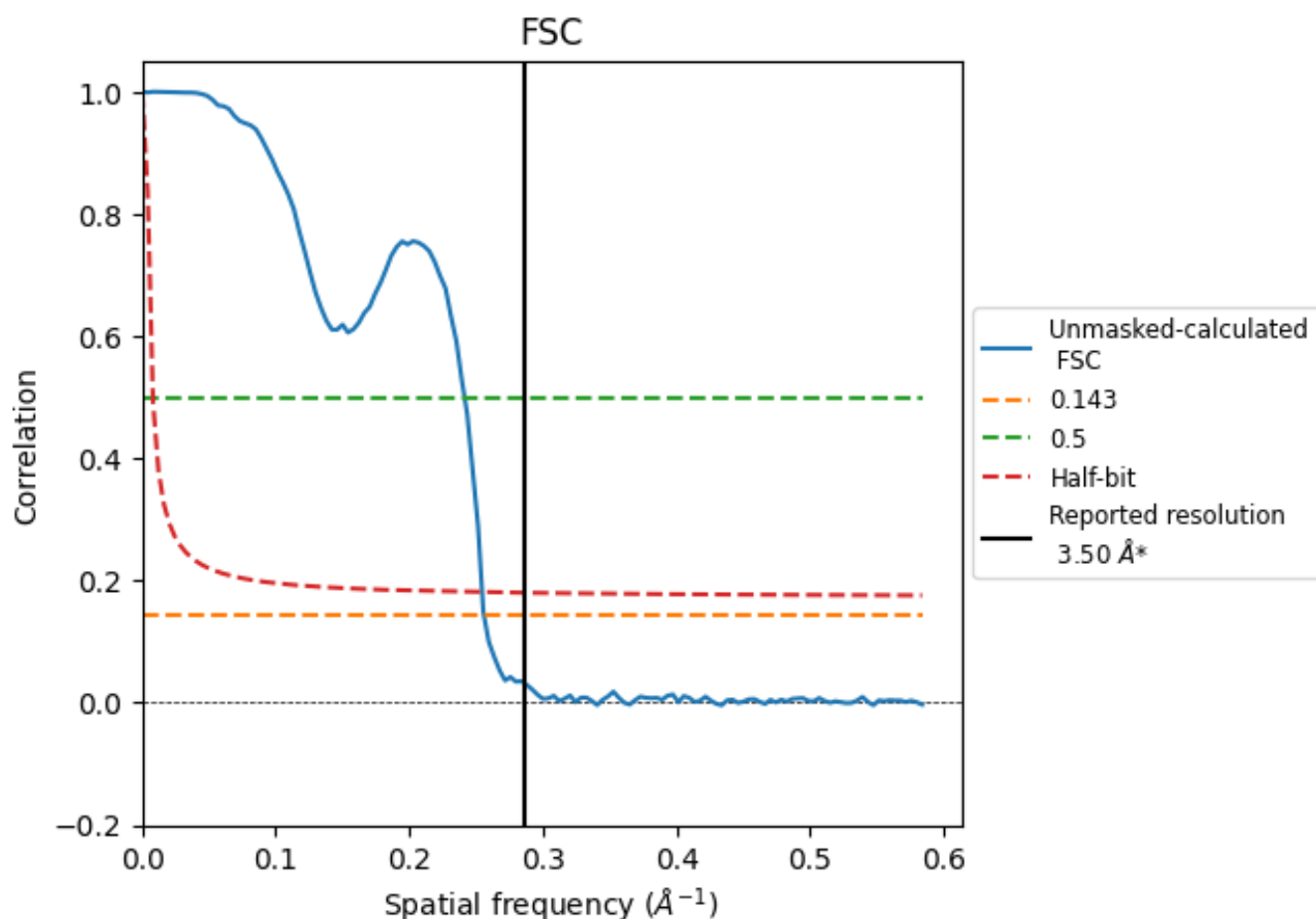


*Reported resolution corresponds to spatial frequency of 0.286 $\text{\AA}^{-1}$

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 0.286 Å⁻¹

8.2 Resolution estimates [i](#)

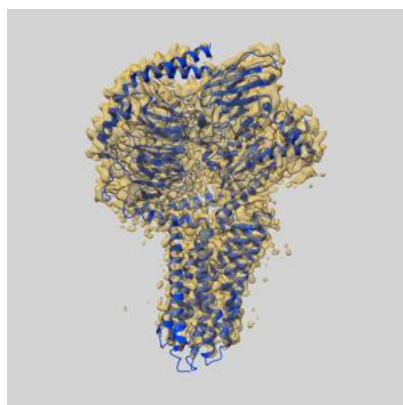
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.91	4.14	3.93

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.91 differs from the reported value 3.5 by more than 10 %

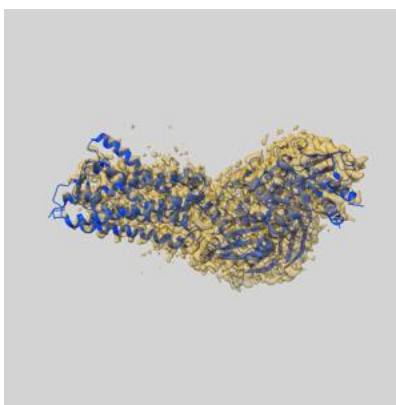
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-42817 and PDB model 8UYQ. Per-residue inclusion information can be found in [section 3](#) on [page 9](#).

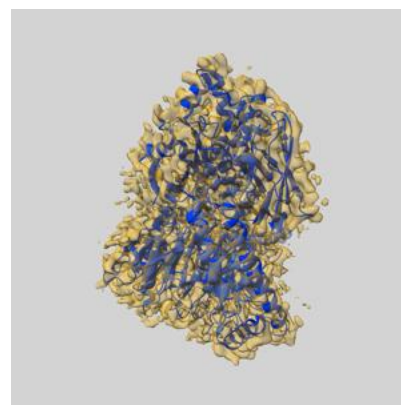
9.1 Map-model overlay [i](#)



X



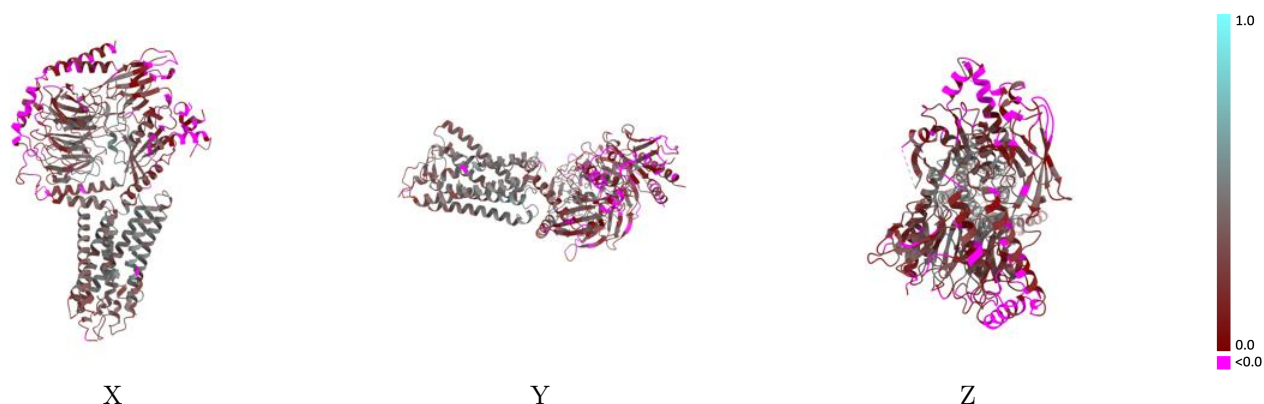
Y



Z

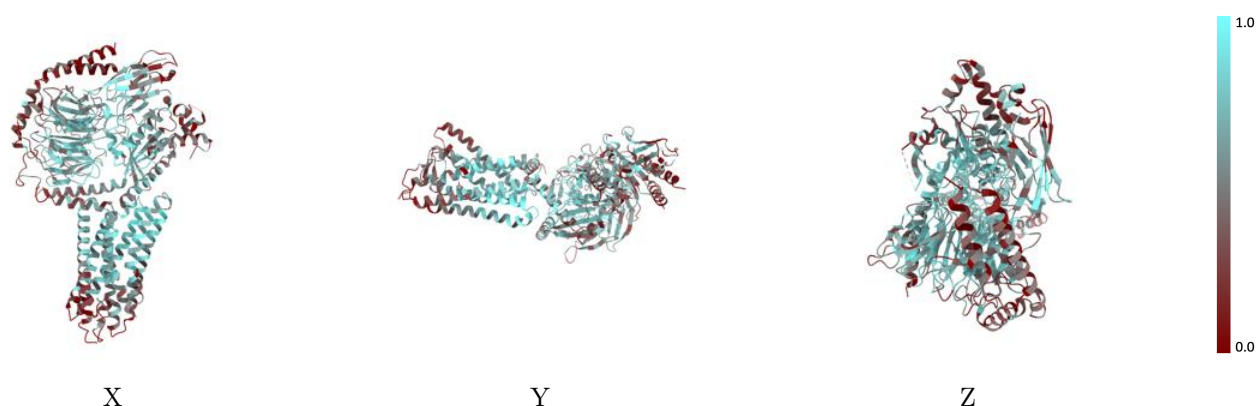
The images above show the 3D surface view of the map at the recommended contour level 0.15 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



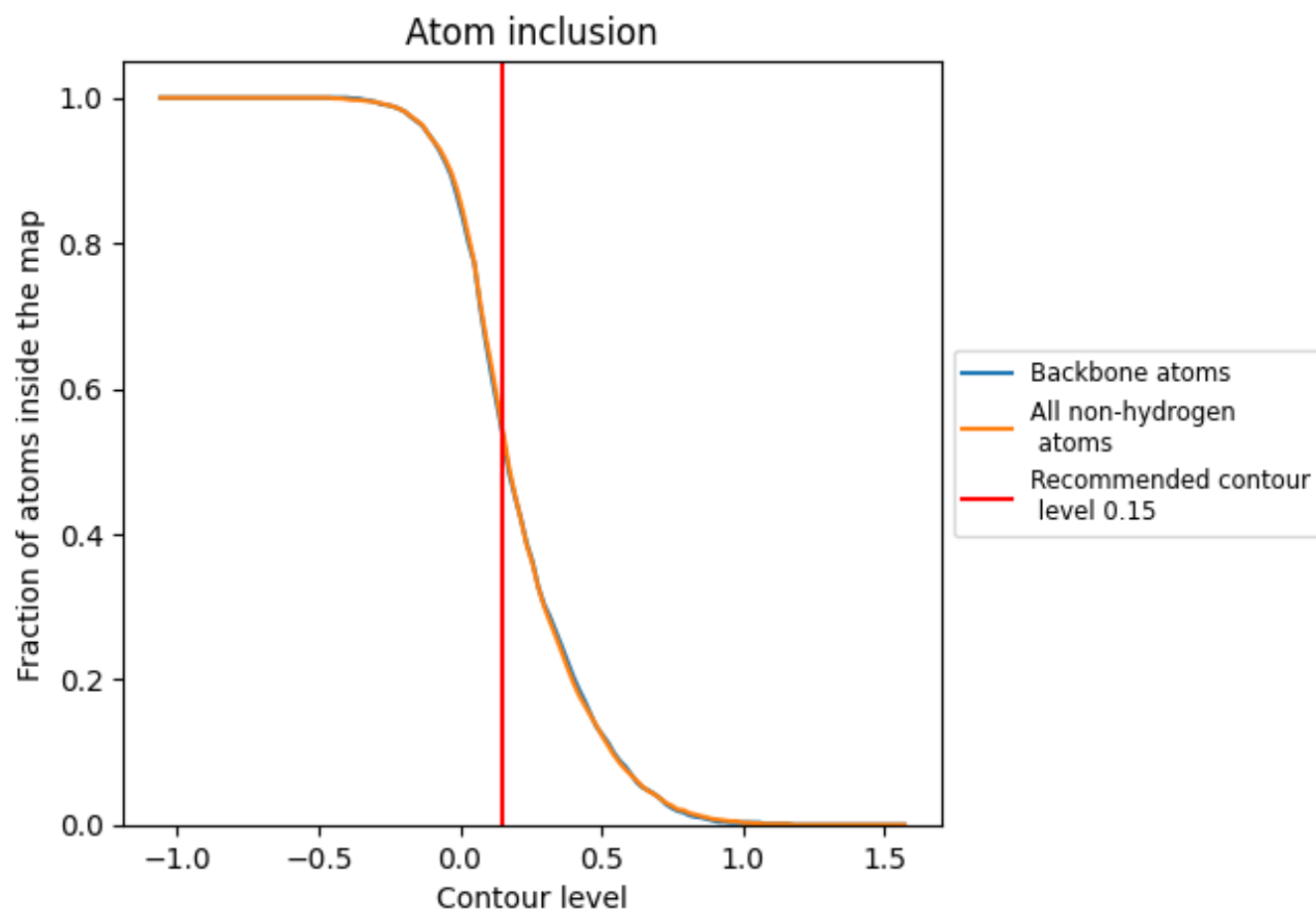
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.15).

9.4 Atom inclusion [i](#)



At the recommended contour level, 54% of all backbone atoms, 54% 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 (0.15) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5430	0.2620
A	0.5440	0.3890
N	0.5600	0.2020
X	0.5440	0.2030
Y	0.5910	0.2440
Z	0.2750	0.0350

