



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 08:32 PM UTC

PDB ID : 6H3L / pdb_00006h3l
EMDB ID : EMD-0135
Title : Structure of VgrG1 in the Type VI secretion "pre-firing" VgrG1-Tse6-EagT6-EF-Tu-Tsi6 complex
Authors : Quentin, D.; Raunser, S.
Deposited on : 2018-07-19
Resolution : 4.20 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
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

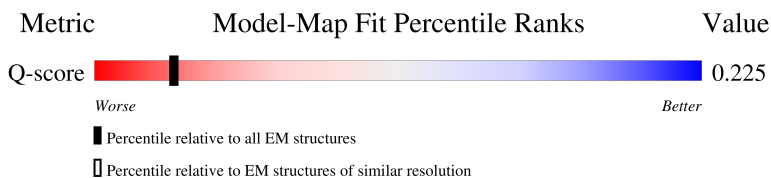
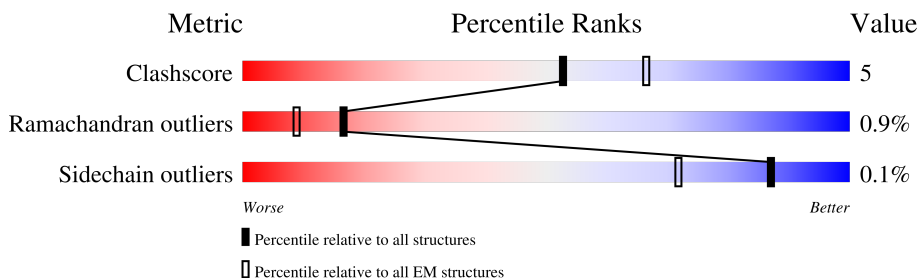
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.20 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	5410 (3.70 - 4.70)

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	643	49% 84% 12% ..
1	B	643	47% 83% 13% ..
1	C	643	51% 84% 12% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 29319 atoms, of which 14388 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.

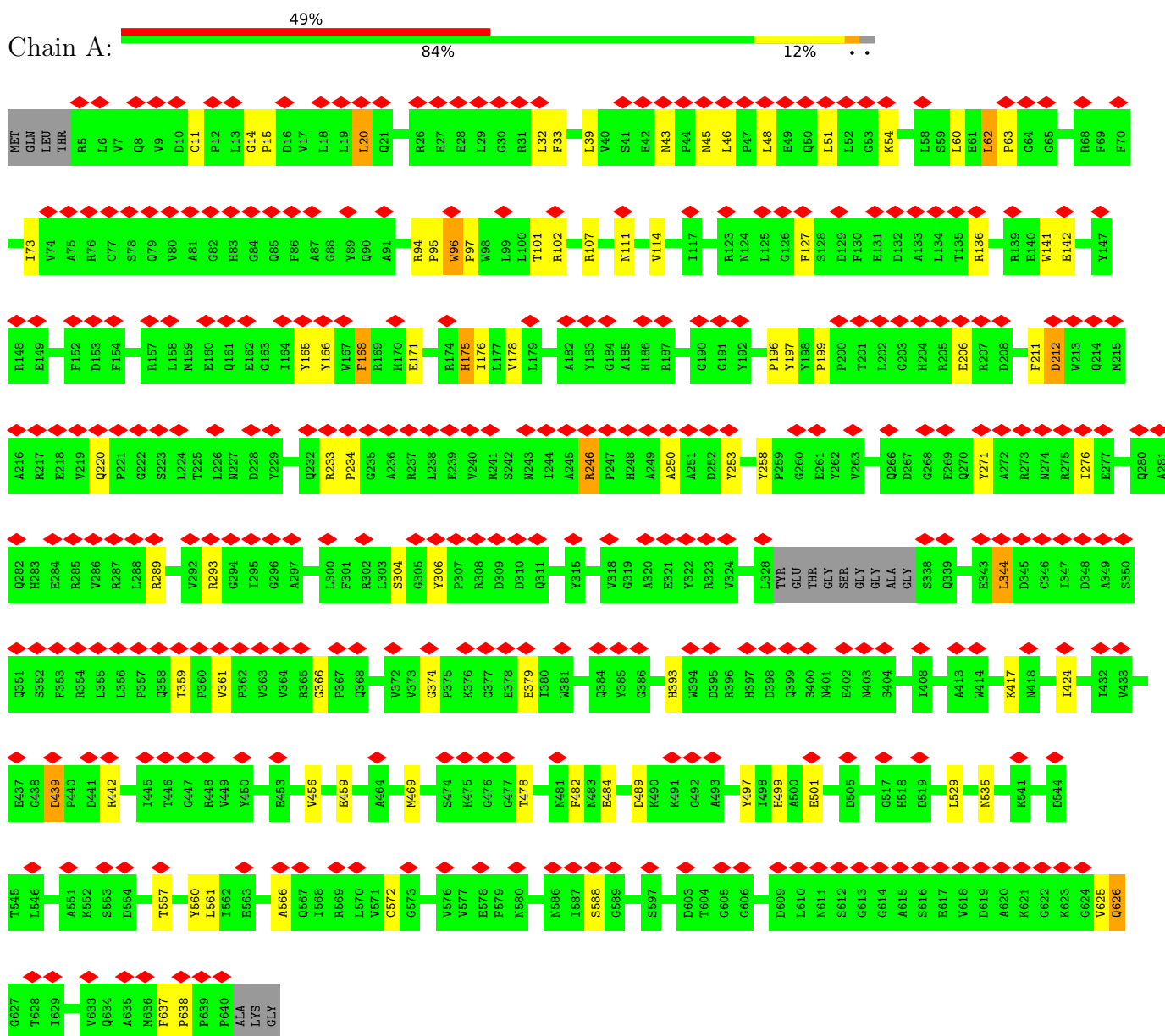
- Molecule 1 is a protein called VgrG1.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	627	Total	C	H	N	O	S	0	0
			9773	3113	4796	910	939	15		
1	B	627	Total	C	H	N	O	S	0	0
			9773	3113	4796	910	939	15		
1	C	627	Total	C	H	N	O	S	0	0
			9773	3113	4796	910	939	15		

3 Residue-property plots

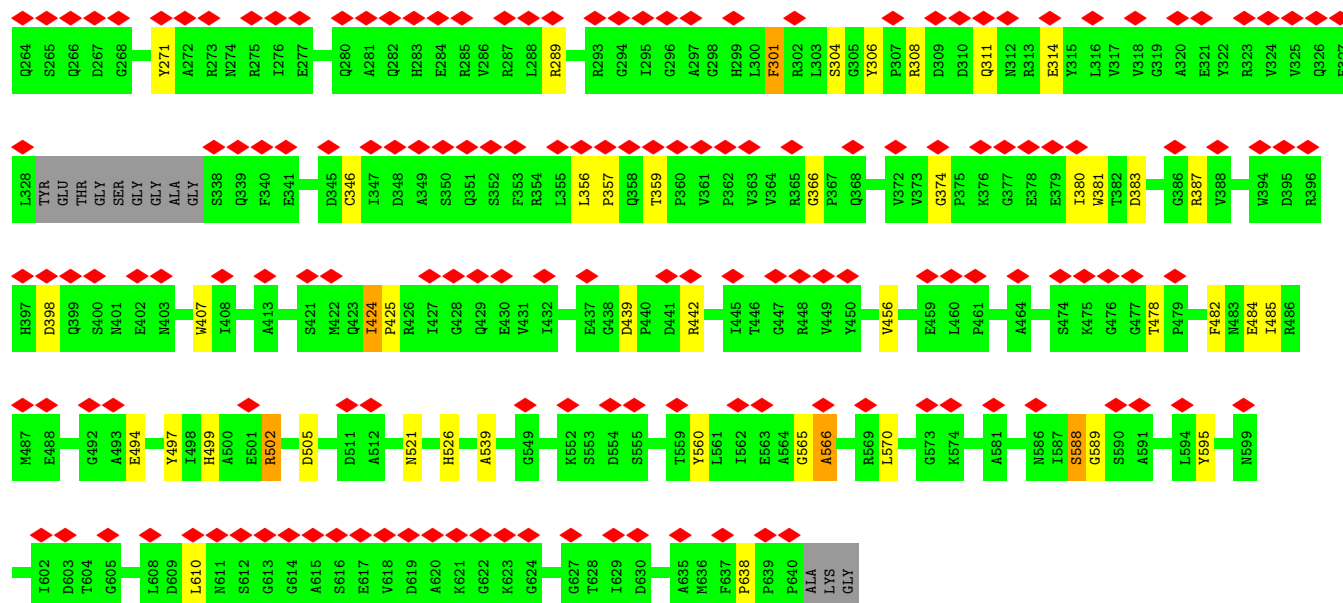
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: VgrG1



• Molecule 1: VgrG1





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	55000	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1700	Depositor
Maximum defocus (nm)	4200	Depositor
Magnification	59000	Depositor
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.115	Depositor
Minimum map value	-0.044	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.053	Depositor
Map size (Å)	410.4, 410.4, 410.4	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.14, 1.14, 1.14	Depositor

5 Model quality

5.1 Standard geometry

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	0.98	0/5095	1.27	70/6905 (1.0%)
1	B	0.96	1/5095 (0.0%)	1.29	70/6905 (1.0%)
1	C	0.95	0/5095	1.28	68/6905 (1.0%)
All	All	0.97	1/15285 (0.0%)	1.28	208/20715 (1.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	625	VAL	CA-C	-5.31	1.48	1.53

All (208) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	638	PRO	CA-C-N	10.44	127.18	119.66
1	B	638	PRO	C-N-CA	10.44	127.18	119.66
1	A	11	CYS	N-CA-C	9.99	116.96	108.07
1	C	638	PRO	CA-C-N	9.88	126.77	119.66
1	C	638	PRO	C-N-CA	9.88	126.77	119.66
1	A	637	PHE	CA-C-N	9.67	126.62	119.66
1	A	637	PHE	C-N-CA	9.67	126.62	119.66
1	B	199	PRO	CA-C-N	9.13	128.78	119.56
1	B	199	PRO	C-N-CA	9.13	128.78	119.56
1	B	175	HIS	CA-CB-CG	-8.91	104.89	113.80
1	A	456	VAL	CA-C-N	8.74	128.90	119.28
1	A	456	VAL	C-N-CA	8.74	128.90	119.28
1	B	424	ILE	CA-C-N	8.60	128.54	119.85
1	B	424	ILE	C-N-CA	8.60	128.54	119.85
1	C	43	ASN	CA-C-N	8.54	128.19	119.82
1	C	43	ASN	C-N-CA	8.54	128.19	119.82
1	C	199	PRO	CA-C-N	8.48	128.21	119.56
1	C	199	PRO	C-N-CA	8.48	128.21	119.56
1	B	366	GLY	CA-C-N	8.43	128.46	119.78
1	B	366	GLY	C-N-CA	8.43	128.46	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	246	ARG	CA-C-N	8.42	129.24	119.47
1	A	246	ARG	C-N-CA	8.42	129.24	119.47
1	A	199	PRO	CA-C-N	8.41	128.14	119.56
1	A	199	PRO	C-N-CA	8.41	128.14	119.56
1	A	43	ASN	CA-C-N	8.37	128.03	119.82
1	A	43	ASN	C-N-CA	8.37	128.03	119.82
1	C	94	ARG	CA-C-N	8.22	128.25	119.78
1	C	94	ARG	C-N-CA	8.22	128.25	119.78
1	B	356	LEU	CA-C-N	8.19	130.08	119.84
1	B	356	LEU	C-N-CA	8.19	130.08	119.84
1	B	54	LYS	CA-C-N	8.15	127.91	119.76
1	B	54	LYS	C-N-CA	8.15	127.91	119.76
1	B	246	ARG	CA-C-N	8.15	128.92	119.47
1	B	246	ARG	C-N-CA	8.15	128.92	119.47
1	B	456	VAL	CA-C-N	8.03	128.47	119.32
1	B	456	VAL	C-N-CA	8.03	128.47	119.32
1	A	54	LYS	CA-C-N	8.01	127.77	119.76
1	A	54	LYS	C-N-CA	8.01	127.77	119.76
1	B	359	THR	CA-C-N	8.01	127.96	120.03
1	B	359	THR	C-N-CA	8.01	127.96	120.03
1	B	220	GLN	CA-C-N	8.01	128.49	119.92
1	B	220	GLN	C-N-CA	8.01	128.49	119.92
1	C	442	ARG	CA-C-N	7.93	127.86	119.85
1	C	442	ARG	C-N-CA	7.93	127.86	119.85
1	C	301	PHE	CA-CB-CG	-7.88	105.92	113.80
1	A	442	ARG	CA-C-N	7.84	127.76	119.85
1	A	442	ARG	C-N-CA	7.84	127.76	119.85
1	C	220	GLN	CA-C-N	7.83	127.85	119.78
1	C	220	GLN	C-N-CA	7.83	127.85	119.78
1	B	188	SER	CA-C-N	7.73	127.68	120.03
1	B	188	SER	C-N-CA	7.73	127.68	120.03
1	A	175	HIS	CA-CB-CG	-7.73	106.07	113.80
1	A	253	TYR	CA-C-N	7.68	128.29	119.99
1	A	253	TYR	C-N-CA	7.68	128.29	119.99
1	C	246	ARG	CA-C-N	7.65	128.34	119.47
1	C	246	ARG	C-N-CA	7.65	128.34	119.47
1	B	253	TYR	CA-C-N	7.64	128.24	120.14
1	B	253	TYR	C-N-CA	7.64	128.24	120.14
1	B	198	TYR	CA-C-N	7.62	128.23	120.38
1	B	198	TYR	C-N-CA	7.62	128.23	120.38
1	A	11	CYS	CA-C-N	7.62	127.62	119.78
1	A	11	CYS	C-N-CA	7.62	127.62	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	94	ARG	CA-C-N	7.62	127.57	120.03
1	B	94	ARG	C-N-CA	7.62	127.57	120.03
1	A	359	THR	CA-C-N	7.58	127.58	119.78
1	A	359	THR	C-N-CA	7.58	127.58	119.78
1	C	356	LEU	CA-C-N	7.53	129.25	119.84
1	C	356	LEU	C-N-CA	7.53	129.25	119.84
1	B	43	ASN	CA-C-N	7.44	128.38	120.12
1	B	43	ASN	C-N-CA	7.44	128.38	120.12
1	C	54	LYS	CA-C-N	7.37	127.13	119.76
1	C	54	LYS	C-N-CA	7.37	127.13	119.76
1	A	366	GLY	CA-C-N	7.35	127.35	119.78
1	A	366	GLY	C-N-CA	7.35	127.35	119.78
1	A	638	PRO	CA-C-N	7.35	127.95	120.38
1	A	638	PRO	C-N-CA	7.35	127.95	120.38
1	C	195	VAL	CA-C-N	7.33	127.50	120.31
1	C	195	VAL	C-N-CA	7.33	127.50	120.31
1	A	220	GLN	CA-C-N	7.29	128.95	119.84
1	A	220	GLN	C-N-CA	7.29	128.95	119.84
1	B	478	THR	CA-C-N	7.29	126.99	119.56
1	B	478	THR	C-N-CA	7.29	126.99	119.56
1	A	276	ILE	N-CA-C	7.27	117.40	110.42
1	A	374	GLY	CA-C-N	7.18	127.17	119.78
1	A	374	GLY	C-N-CA	7.18	127.17	119.78
1	B	69	PHE	CA-CB-CG	-7.15	106.65	113.80
1	C	424	ILE	CA-C-N	7.11	128.72	119.84
1	C	424	ILE	C-N-CA	7.11	128.72	119.84
1	C	366	GLY	CA-C-N	7.10	127.02	119.85
1	C	366	GLY	C-N-CA	7.10	127.02	119.85
1	C	359	THR	CA-C-N	7.05	127.02	119.76
1	C	359	THR	C-N-CA	7.05	127.02	119.76
1	C	478	THR	CA-C-N	7.04	126.72	119.82
1	C	478	THR	C-N-CA	7.04	126.72	119.82
1	C	136	ARG	CA-C-N	7.00	126.92	119.85
1	C	136	ARG	C-N-CA	7.00	126.92	119.85
1	C	253	TYR	CA-C-N	6.98	127.54	120.14
1	C	253	TYR	C-N-CA	6.98	127.54	120.14
1	B	145	VAL	N-CA-C	6.98	118.77	108.45
1	A	439	ASP	CA-C-N	6.95	128.53	119.84
1	A	439	ASP	C-N-CA	6.95	128.53	119.84
1	B	374	GLY	CA-C-N	6.87	126.85	119.78
1	B	374	GLY	C-N-CA	6.87	126.85	119.78
1	B	306	TYR	CA-C-N	6.79	127.06	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	306	TYR	C-N-CA	6.79	127.06	119.32
1	C	374	GLY	CA-C-N	6.75	126.73	119.78
1	C	374	GLY	C-N-CA	6.75	126.73	119.78
1	C	456	VAL	CA-C-N	6.74	128.26	119.84
1	C	456	VAL	C-N-CA	6.74	128.26	119.84
1	C	439	ASP	CA-C-N	6.72	128.24	119.84
1	C	439	ASP	C-N-CA	6.72	128.24	119.84
1	B	442	ARG	CA-C-N	6.69	126.61	119.85
1	B	442	ARG	C-N-CA	6.69	126.61	119.85
1	A	96	TRP	CA-C-N	6.59	127.11	119.47
1	A	96	TRP	C-N-CA	6.59	127.11	119.47
1	B	195	VAL	CA-C-N	6.53	126.50	120.03
1	B	195	VAL	C-N-CA	6.53	126.50	120.03
1	A	136	ARG	CA-C-N	6.51	127.05	120.14
1	A	136	ARG	C-N-CA	6.51	127.05	120.14
1	C	526	HIS	CB-CA-C	-6.45	107.13	116.54
1	A	306	TYR	CA-C-N	6.39	126.61	119.32
1	A	306	TYR	C-N-CA	6.39	126.61	119.32
1	C	176	ILE	N-CA-C	6.39	117.70	107.73
1	B	176	ILE	N-CA-C	6.38	117.11	108.17
1	B	233	ARG	CA-C-N	6.34	127.77	119.84
1	B	233	ARG	C-N-CA	6.34	127.77	119.84
1	B	11	CYS	CA-C-N	6.34	126.31	119.78
1	B	11	CYS	C-N-CA	6.34	126.31	119.78
1	B	439	ASP	CA-C-N	6.20	127.59	119.84
1	B	439	ASP	C-N-CA	6.20	127.59	119.84
1	B	136	ARG	CA-C-N	6.19	126.70	120.14
1	B	136	ARG	C-N-CA	6.19	126.70	120.14
1	C	114	VAL	CA-C-N	6.18	125.67	119.24
1	C	114	VAL	C-N-CA	6.18	125.67	119.24
1	C	198	TYR	CA-C-N	6.14	126.71	120.38
1	C	198	TYR	C-N-CA	6.14	126.71	120.38
1	C	306	TYR	CA-C-N	6.14	126.32	119.32
1	C	306	TYR	C-N-CA	6.14	126.32	119.32
1	A	176	ILE	N-CA-C	6.13	117.13	108.36
1	A	233	ARG	CA-C-N	6.13	127.50	119.84
1	A	233	ARG	C-N-CA	6.13	127.50	119.84
1	B	46	LEU	CA-C-N	6.13	127.50	119.84
1	B	46	LEU	C-N-CA	6.13	127.50	119.84
1	A	114	VAL	CA-C-N	6.13	126.30	119.32
1	A	114	VAL	C-N-CA	6.13	126.30	119.32
1	A	478	THR	CA-C-N	6.09	125.77	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	478	THR	C-N-CA	6.09	125.77	119.56
1	B	62	LEU	CA-C-N	6.09	126.11	119.90
1	B	62	LEU	C-N-CA	6.09	126.11	119.90
1	C	588	SER	N-CA-C	6.03	116.52	108.38
1	A	46	LEU	CA-C-N	6.00	125.96	119.78
1	A	46	LEU	C-N-CA	6.00	125.96	119.78
1	C	188	SER	CA-C-N	6.00	125.94	119.76
1	C	188	SER	C-N-CA	6.00	125.94	119.76
1	C	311	GLN	N-CA-C	-5.99	106.11	113.41
1	A	588	SER	N-CA-C	5.93	117.26	108.60
1	A	51	LEU	N-CA-C	-5.93	106.08	113.55
1	B	186	HIS	CA-CB-CG	-5.92	107.88	113.80
1	A	94	ARG	CA-C-N	5.89	126.08	120.31
1	A	94	ARG	C-N-CA	5.89	126.08	120.31
1	C	186	HIS	CA-CB-CG	-5.83	107.97	113.80
1	A	168	PHE	N-CA-C	5.82	118.03	109.24
1	B	301	PHE	CA-CB-CG	-5.82	107.98	113.80
1	B	526	HIS	CA-CB-CG	5.77	119.57	113.80
1	B	114	VAL	CA-C-N	5.67	125.13	119.24
1	B	114	VAL	C-N-CA	5.67	125.13	119.24
1	C	346	CYS	N-CA-C	5.67	117.58	109.14
1	C	62	LEU	CA-C-N	5.62	125.55	119.76
1	C	62	LEU	C-N-CA	5.62	125.55	119.76
1	A	379	GLU	N-CA-C	-5.54	107.06	113.88
1	C	526	HIS	CA-CB-CG	5.50	119.31	113.80
1	C	168	PHE	CA-CB-CG	5.48	119.28	113.80
1	C	38	HIS	CA-CB-CG	-5.45	108.35	113.80
1	C	11	CYS	CA-C-N	5.45	125.35	119.85
1	C	11	CYS	C-N-CA	5.45	125.35	119.85
1	A	344	LEU	N-CA-C	5.44	117.34	109.07
1	C	595	TYR	N-CA-C	5.44	118.31	109.06
1	C	46	LEU	CA-C-N	5.42	126.62	119.84
1	C	46	LEU	C-N-CA	5.42	126.62	119.84
1	A	361	VAL	CA-C-N	5.37	126.55	119.84
1	A	361	VAL	C-N-CA	5.37	126.55	119.84
1	C	69	PHE	CA-CB-CG	-5.35	108.45	113.80
1	A	48	LEU	N-CA-C	-5.31	106.48	113.12
1	A	171	GLU	N-CA-C	-5.29	103.70	110.53
1	B	11	CYS	N-CA-C	5.27	114.20	108.14
1	B	361	VAL	CA-C-N	5.26	125.02	119.76
1	B	361	VAL	C-N-CA	5.26	125.02	119.76
1	A	393	HIS	CA-CB-CG	-5.25	108.55	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	LEU	CA-C-N	5.25	125.17	119.76
1	A	62	LEU	C-N-CA	5.25	125.17	119.76
1	B	414	TRP	N-CA-C	5.25	115.98	107.32
1	B	14	GLY	CA-C-N	5.20	126.34	119.84
1	B	14	GLY	C-N-CA	5.20	126.34	119.84
1	C	233	ARG	CA-C-N	5.16	126.29	119.84
1	C	233	ARG	C-N-CA	5.16	126.29	119.84
1	A	293	ARG	N-CA-C	5.16	118.73	112.23
1	A	212	ASP	CA-CB-CG	5.16	117.76	112.60
1	A	14	GLY	CA-C-N	5.12	126.24	119.84
1	A	14	GLY	C-N-CA	5.12	126.24	119.84
1	A	424	ILE	CA-C-N	5.12	124.88	119.76
1	A	424	ILE	C-N-CA	5.12	124.88	119.76
1	B	637	PHE	CA-C-N	5.09	125.63	120.38
1	B	637	PHE	C-N-CA	5.09	125.63	120.38
1	A	211	PHE	CA-CB-CG	-5.05	108.75	113.80
1	C	484	GLU	N-CA-C	5.05	115.97	108.60
1	B	126	GLY	N-CA-C	-5.02	108.35	115.43
1	C	271	TYR	CA-CB-CG	-5.01	104.88	113.90
1	A	276	ILE	CB-CA-C	-5.01	105.56	111.97

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4977	4796	4794	50	0
1	B	4977	4796	4794	61	0
1	C	4977	4796	4794	48	0
All	All	14931	14388	14382	155	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 5.

All (155) 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:C:192:TYR:CD2	1:C:192:TYR:O	2.12	1.02
1:B:177:LEU:C	1:B:177:LEU:HD23	2.05	0.81
1:A:20:LEU:HD23	1:A:20:LEU:O	1.82	0.78
1:A:96:TRP:C	1:A:96:TRP:CD1	2.63	0.75
1:C:192:TYR:O	1:C:192:TYR:HD2	1.67	0.74
1:B:409:ARG:N	1:B:409:ARG:HD3	2.02	0.73
1:A:96:TRP:CD1	1:A:96:TRP:O	2.45	0.70
1:B:96:TRP:C	1:B:96:TRP:CD1	2.69	0.70
1:A:96:TRP:O	1:A:96:TRP:HD1	1.74	0.69
1:A:20:LEU:HD23	1:A:20:LEU:C	2.18	0.69
1:B:409:ARG:HD3	1:B:409:ARG:H	1.56	0.69
1:B:123:ARG:NE	1:B:123:ARG:HA	2.11	0.66
1:C:494:GLU:OE1	1:C:494:GLU:N	2.24	0.65
1:B:96:TRP:CD1	1:B:96:TRP:O	2.51	0.64
1:B:414:TRP:C	1:B:414:TRP:CD1	2.76	0.64
1:A:96:TRP:C	1:A:96:TRP:HD1	2.04	0.64
1:B:422:MET:SD	1:B:422:MET:O	2.55	0.64
1:B:96:TRP:O	1:B:96:TRP:HD1	1.81	0.63
1:A:20:LEU:O	1:A:20:LEU:CD2	2.46	0.62
1:A:101:THR:OG1	1:A:102:ARG:NH1	2.33	0.61
1:A:501:GLU:O	1:A:501:GLU:HG2	2.00	0.61
1:B:494:GLU:N	1:B:494:GLU:OE1	2.33	0.60
1:B:96:TRP:C	1:B:96:TRP:HD1	2.09	0.59
1:C:134:LEU:C	1:C:134:LEU:HD12	2.26	0.59
1:B:258:TYR:CD2	1:B:258:TYR:C	2.79	0.59
1:A:344:LEU:C	1:A:344:LEU:HD12	2.28	0.58
1:B:177:LEU:C	1:B:177:LEU:CD2	2.76	0.58
1:B:422:MET:SD	1:B:422:MET:C	2.88	0.57
1:C:502:ARG:HG2	1:C:502:ARG:O	2.04	0.57
1:A:212:ASP:HB3	1:A:289:ARG:HB3	1.87	0.57
1:C:195:VAL:O	1:C:195:VAL:HG13	2.05	0.57
1:A:206:GLU:H	1:A:206:GLU:CD	2.14	0.56
1:B:527:ASP:OD1	1:B:527:ASP:C	2.49	0.55
1:B:496:LEU:HD13	1:B:496:LEU:C	2.31	0.55
1:A:497:TYR:CD2	1:A:497:TYR:C	2.85	0.55
1:B:414:TRP:CD1	1:B:414:TRP:O	2.60	0.54
1:C:96:TRP:CD1	1:C:96:TRP:C	2.85	0.54
1:C:258:TYR:C	1:C:258:TYR:CD2	2.85	0.54
1:A:258:TYR:C	1:A:258:TYR:CD2	2.85	0.54
1:B:142:GLU:OE1	1:B:142:GLU:N	2.41	0.54
1:A:39:LEU:C	1:A:39:LEU:HD12	2.34	0.53
1:B:540:VAL:O	1:B:540:VAL:HG13	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:535:ASN:OD1	1:A:535:ASN:C	2.52	0.52
1:A:142:GLU:OE1	1:A:142:GLU:N	2.30	0.52
1:B:177:LEU:HD23	1:B:177:LEU:O	2.09	0.51
1:A:95:PRO:HB2	1:A:97:PRO:HD2	1.92	0.51
1:A:482:PHE:N	1:A:482:PHE:CD2	2.76	0.51
1:A:484:GLU:HG2	1:A:499:HIS:HB3	1.93	0.51
1:A:489:ASP:OD1	1:A:489:ASP:C	2.53	0.50
1:B:206:GLU:OE1	1:B:206:GLU:N	2.34	0.50
1:A:165:TYR:CD1	1:A:165:TYR:N	2.80	0.50
1:C:565:GLY:O	1:C:566:ALA:HB3	2.12	0.50
1:A:246:ARG:NE	1:A:246:ARG:HA	2.27	0.50
1:C:499:HIS:CD2	1:C:499:HIS:C	2.90	0.50
1:C:381:TRP:HA	1:C:381:TRP:CE3	2.47	0.49
1:A:459:GLU:OE1	1:A:459:GLU:N	2.33	0.49
1:B:177:LEU:CD2	1:B:177:LEU:O	2.61	0.49
1:C:610:LEU:HD12	1:C:610:LEU:C	2.37	0.49
1:C:588:SER:OG	1:C:589:GLY:N	2.45	0.49
1:B:40:VAL:HB	1:B:86:PHE:HB3	1.95	0.49
1:C:482:PHE:C	1:C:482:PHE:CD1	2.91	0.49
1:A:127:PHE:CD1	1:A:127:PHE:C	2.89	0.49
1:B:132:ASP:OD1	1:B:132:ASP:C	2.56	0.48
1:A:625:VAL:O	1:A:626:GLN:CB	2.62	0.48
1:C:132:ASP:C	1:C:132:ASP:OD1	2.56	0.48
1:B:505:ASP:OD1	1:B:505:ASP:C	2.56	0.48
1:A:33:PHE:CD1	1:A:33:PHE:C	2.92	0.47
1:A:32:LEU:HD12	1:A:32:LEU:N	2.29	0.47
1:C:521:ASN:OD1	1:C:521:ASN:C	2.56	0.47
1:A:197:TYR:C	1:A:197:TYR:CD2	2.93	0.47
1:B:10:ASP:OD1	1:B:10:ASP:C	2.57	0.47
1:A:469:MET:SD	1:A:469:MET:C	2.98	0.47
1:B:636:MET:HG2	1:C:539:ALA:HB3	1.97	0.47
1:B:48:LEU:HB3	1:B:77:CYS:SG	2.55	0.47
1:B:414:TRP:O	1:B:414:TRP:HD1	1.97	0.47
1:C:48:LEU:HB3	1:C:77:CYS:SG	2.55	0.46
1:A:102:ARG:NE	1:A:102:ARG:HA	2.31	0.46
1:B:570:LEU:C	1:B:570:LEU:HD23	2.41	0.46
1:C:62:LEU:HB3	1:C:63:PRO:HD2	1.97	0.46
1:C:502:ARG:O	1:C:502:ARG:CG	2.64	0.46
1:A:73:ILE:O	1:A:73:ILE:HG13	2.15	0.46
1:B:181:ASP:OD1	1:B:181:ASP:C	2.59	0.46
1:B:561:LEU:C	1:B:561:LEU:HD23	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:TYR:C	1:B:166:TYR:CD1	2.93	0.46
1:C:570:LEU:C	1:C:570:LEU:HD23	2.40	0.46
1:A:45:ASN:OD1	1:C:308:ARG:NH1	2.49	0.45
1:A:60:LEU:C	1:A:60:LEU:HD23	2.41	0.45
1:C:142:GLU:OE1	1:C:142:GLU:N	2.35	0.45
1:B:62:LEU:HB3	1:B:63:PRO:HD2	1.98	0.45
1:A:417:LYS:HA	1:B:454:GLN:HA	1.98	0.45
1:A:168:PHE:HB3	1:A:175:HIS:HE2	1.81	0.45
1:B:395:ASP:OD1	1:B:395:ASP:C	2.60	0.45
1:B:546:LEU:C	1:B:546:LEU:HD23	2.42	0.45
1:A:561:LEU:C	1:A:561:LEU:HD23	2.42	0.44
1:B:123:ARG:NE	1:B:123:ARG:CA	2.80	0.44
1:C:56:MET:SD	1:C:56:MET:C	3.00	0.44
1:B:96:TRP:HB3	1:B:97:PRO:HD3	1.98	0.44
1:B:544:ASP:OD1	1:B:544:ASP:C	2.60	0.44
1:B:233:ARG:O	1:B:235:GLY:N	2.49	0.44
1:C:301:PHE:CD1	1:C:301:PHE:C	2.90	0.44
1:C:96:TRP:HB3	1:C:97:PRO:HD3	2.00	0.44
1:B:246:ARG:NE	1:B:246:ARG:HA	2.32	0.44
1:C:424:ILE:HA	1:C:425:PRO:HD3	1.85	0.43
1:A:178:VAL:O	1:A:178:VAL:HG12	2.17	0.43
1:B:197:TYR:CD2	1:B:197:TYR:C	2.96	0.43
1:A:271:TYR:N	1:A:271:TYR:CD1	2.84	0.43
1:B:560:TYR:C	1:B:560:TYR:CD1	2.94	0.43
1:B:176:ILE:HD12	1:B:176:ILE:N	2.34	0.43
1:C:35:TYR:HB2	1:C:93:LEU:HB3	2.01	0.43
1:B:302:ARG:HA	1:B:314:GLU:HA	2.00	0.43
1:A:32:LEU:N	1:A:32:LEU:CD1	2.82	0.42
1:C:192:TYR:O	1:C:192:TYR:CG	2.68	0.42
1:C:497:TYR:CE2	1:C:499:HIS:HB2	2.53	0.42
1:C:383:ASP:OD1	1:C:383:ASP:C	2.61	0.42
1:B:33:PHE:CD1	1:B:33:PHE:C	2.97	0.42
1:A:62:LEU:HB3	1:A:63:PRO:HD2	2.01	0.42
1:C:6:LEU:C	1:C:6:LEU:HD12	2.44	0.42
1:C:114:VAL:HB	1:C:115:PRO:HD3	2.00	0.42
1:B:345:ASP:OD1	1:B:345:ASP:C	2.63	0.42
1:C:485:ILE:HG23	1:C:485:ILE:O	2.19	0.42
1:A:196:PRO:HA	1:A:304:SER:O	2.20	0.42
1:B:32:LEU:HD22	1:B:95:PRO:HG2	2.01	0.42
1:B:167:TRP:CD1	1:B:167:TRP:C	2.97	0.42
1:C:168:PHE:HA	1:C:176:ILE:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:439:ASP:OD1	1:A:439:ASP:C	2.63	0.41
1:A:529:LEU:C	1:A:529:LEU:HD23	2.45	0.41
1:B:134:LEU:HD11	1:B:179:LEU:HB3	2.02	0.41
1:C:80:VAL:HB	1:C:88:GLY:HA3	2.02	0.41
1:B:207:ARG:O	1:B:208:ASP:C	2.62	0.41
1:C:380:ILE:O	1:C:380:ILE:HG22	2.20	0.41
1:C:560:TYR:C	1:C:560:TYR:CD1	2.98	0.41
1:B:383:ASP:OD1	1:B:383:ASP:C	2.63	0.41
1:A:33:PHE:CE2	1:A:95:PRO:HB3	2.56	0.41
1:A:560:TYR:C	1:A:560:TYR:CD1	2.97	0.41
1:A:107:ARG:CZ	1:A:107:ARG:HB2	2.50	0.41
1:A:111:ASN:N	1:A:141:TRP:O	2.54	0.41
1:B:94:ARG:HB3	1:B:95:PRO:HD2	2.02	0.41
1:C:179:LEU:C	1:C:179:LEU:HD23	2.46	0.41
1:A:572:CYS:SG	1:B:567:GLN:HA	2.60	0.41
1:C:197:TYR:N	1:C:304:SER:O	2.53	0.41
1:B:38:HIS:CD2	1:B:38:HIS:C	2.99	0.41
1:B:301:PHE:O	1:B:301:PHE:CG	2.73	0.41
1:C:387:ARG:HB3	1:C:407:TRP:HB3	2.03	0.41
1:C:505:ASP:C	1:C:505:ASP:OD1	2.61	0.41
1:C:212:ASP:HB3	1:C:289:ARG:HB3	2.02	0.41
1:B:100:LEU:CD1	1:B:100:LEU:N	2.84	0.40
1:C:301:PHE:O	1:C:301:PHE:CG	2.69	0.40
1:C:398:ASP:OD1	1:C:398:ASP:C	2.63	0.40
1:B:301:PHE:CD1	1:B:301:PHE:C	2.97	0.40
1:A:166:TYR:CD1	1:A:166:TYR:O	2.74	0.40
1:B:100:LEU:N	1:B:100:LEU:HD12	2.36	0.40
1:B:114:VAL:HB	1:B:115:PRO:HD3	2.03	0.40
1:C:134:LEU:HD12	1:C:134:LEU:O	2.21	0.40
1:C:301:PHE:O	1:C:314:GLU:HA	2.21	0.40
1:C:381:TRP:HA	1:C:381:TRP:HE3	1.87	0.40

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	623/643 (97%)	576 (92%)	41 (7%)	6 (1%)	12	47
1	B	623/643 (97%)	569 (91%)	50 (8%)	4 (1%)	21	58
1	C	623/643 (97%)	564 (90%)	52 (8%)	7 (1%)	11	45
All	All	1869/1929 (97%)	1709 (91%)	143 (8%)	17 (1%)	16	49

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	626	GLN
1	B	234	PRO
1	B	236	ALA
1	C	566	ALA
1	A	15	PRO
1	A	20	LEU
1	A	250	ALA
1	A	566	ALA
1	C	199	PRO
1	C	502	ARG
1	B	15	PRO
1	C	15	PRO
1	C	47	PRO
1	C	74	VAL
1	C	357	PRO
1	A	234	PRO
1	B	74	VAL

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	526/535 (98%)	525 (100%)	1 (0%)	87	85
1	B	526/535 (98%)	526 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	526/535 (98%)	526 (100%)	0	100	100
All	All	1578/1605 (98%)	1577 (100%)	1 (0%)	87	88

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	557	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	110	GLN
1	A	120	GLN
1	A	220	GLN
1	A	232	GLN
1	A	601	ASN
1	A	626	GLN
1	B	232	GLN
1	B	351	GLN
1	B	506	ASN
1	C	21	GLN
1	C	110	GLN
1	C	146	GLN
1	C	270	GLN
1	C	423	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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](#)

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.

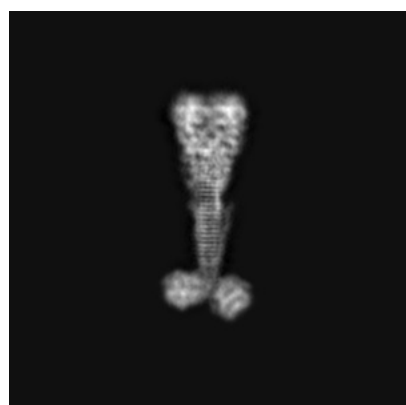
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-0135. These allow visual inspection of the internal detail of the map and identification of artifacts.

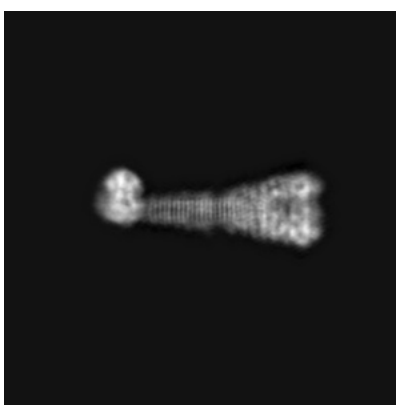
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

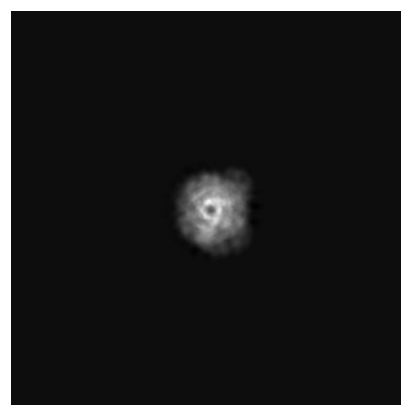
6.1.1 Primary map



X



Y

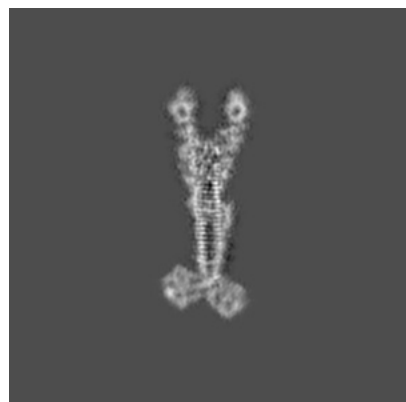


Z

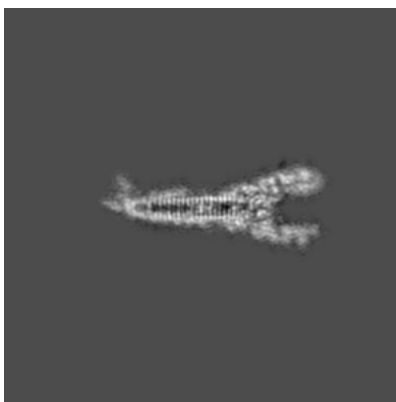
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

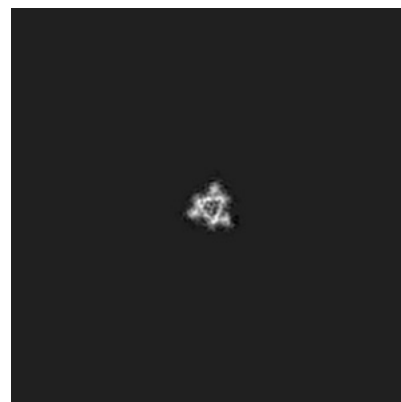
6.2.1 Primary map



X Index: 180



Y Index: 180

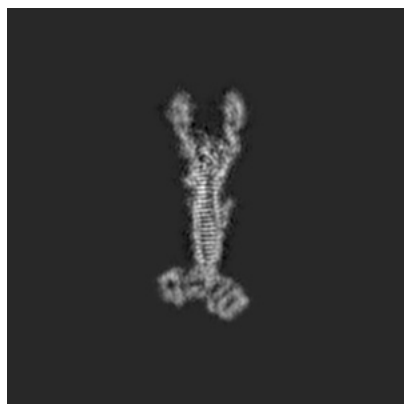


Z Index: 180

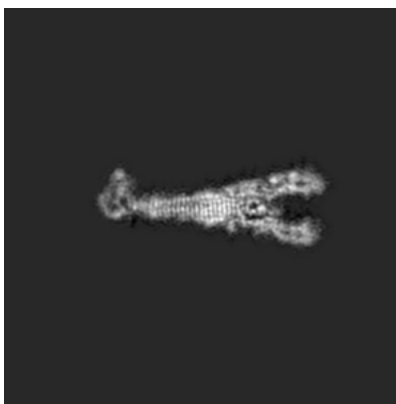
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

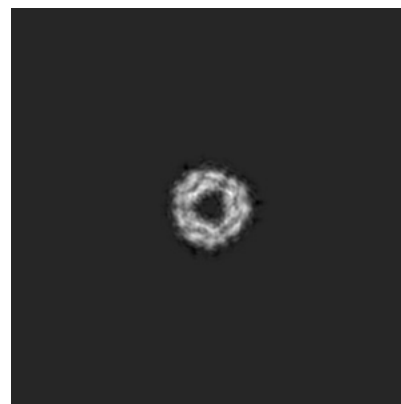
6.3.1 Primary map



X Index: 185



Y Index: 187

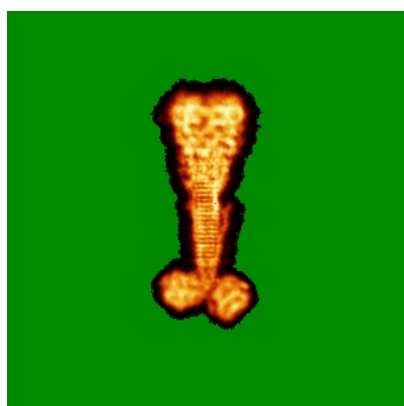


Z Index: 271

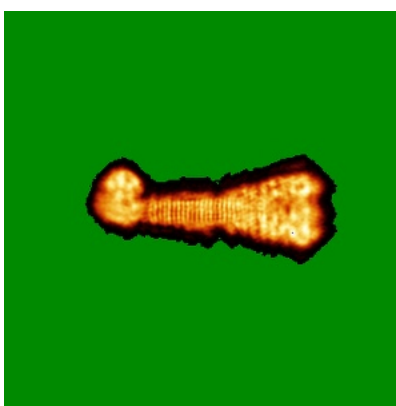
The images above show the largest variance slices of the map in three orthogonal directions.

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

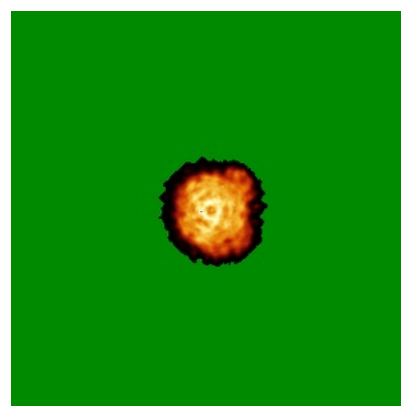
6.4.1 Primary map



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

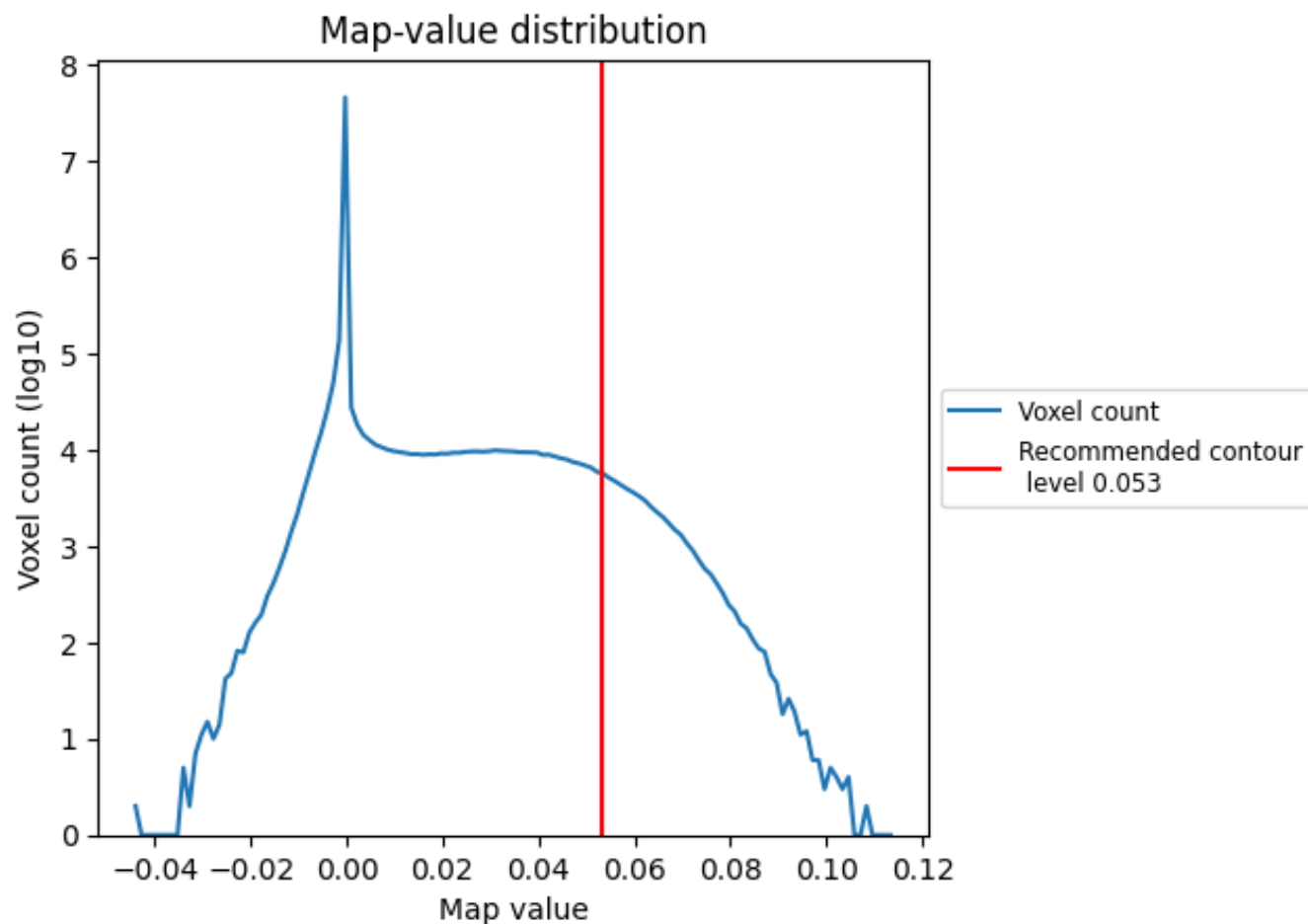
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

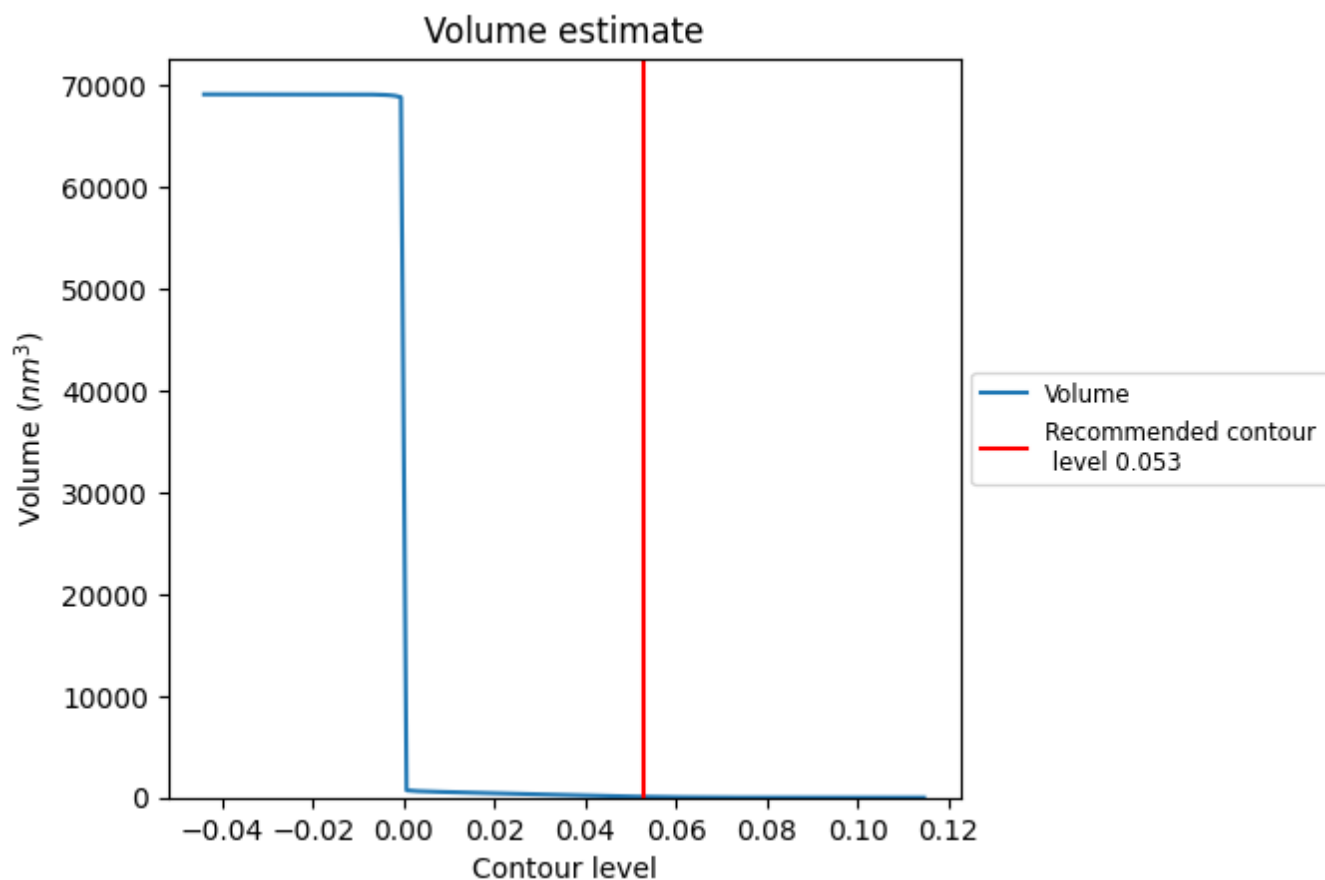
This section contains the results of statistical analysis of the map.

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.

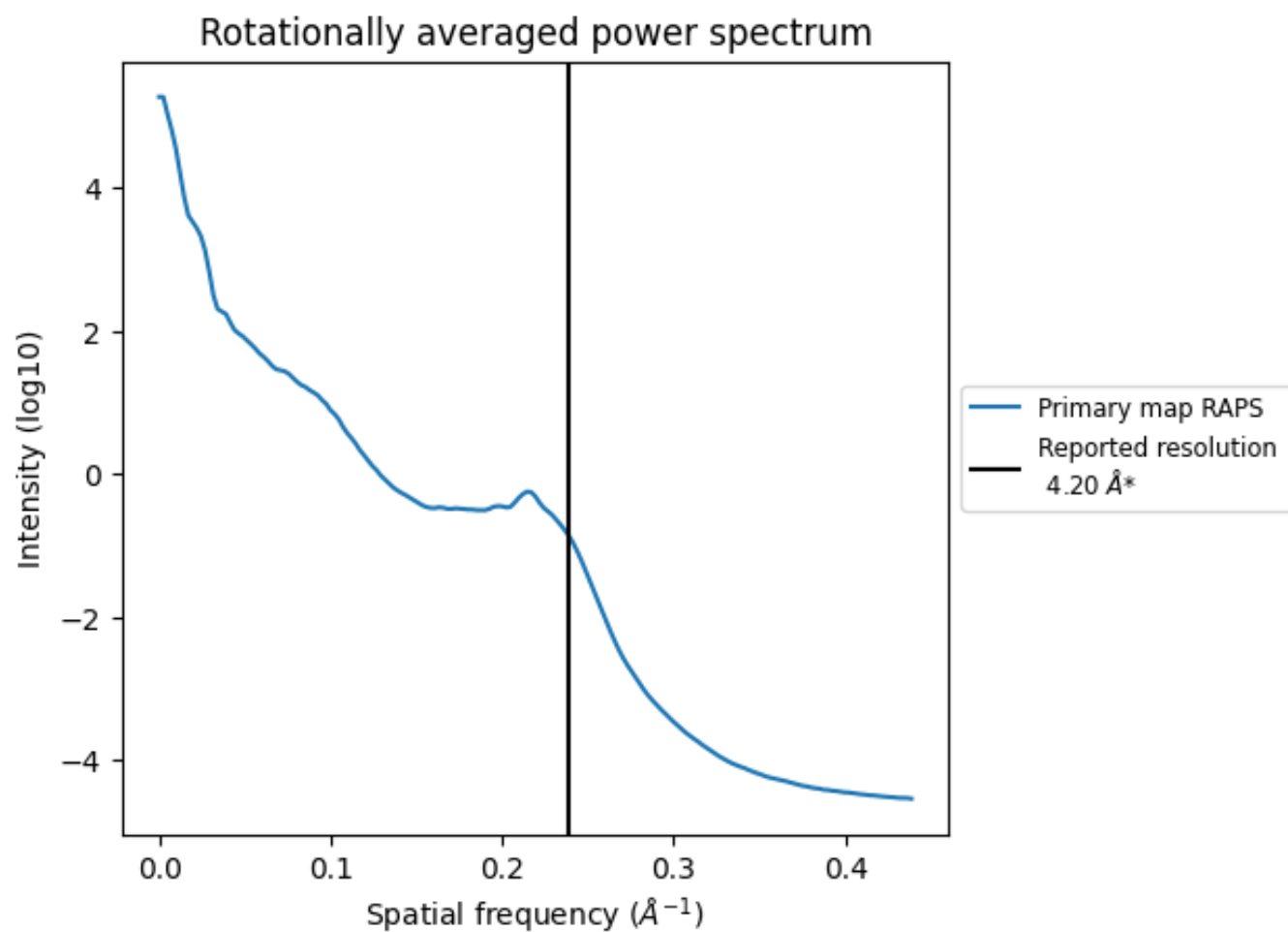
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 78 nm^3 ; this corresponds to an approximate mass of 71 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 0.238 Å⁻¹

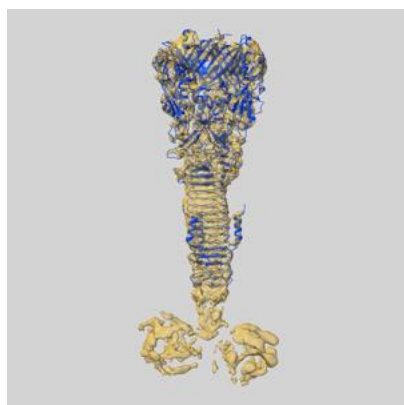
8 Fourier-Shell correlation ⓘ

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

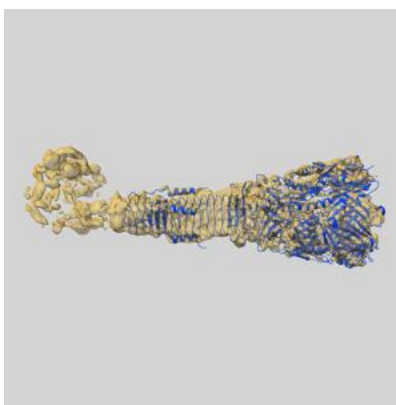
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-0135 and PDB model 6H3L. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

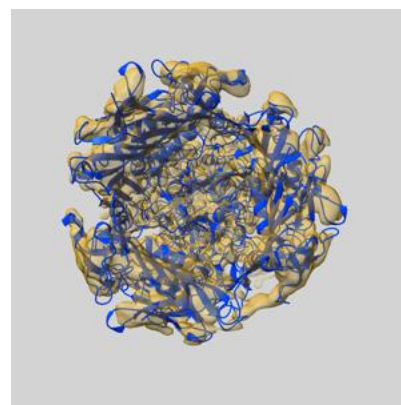
9.1 Map-model overlay [i](#)



X



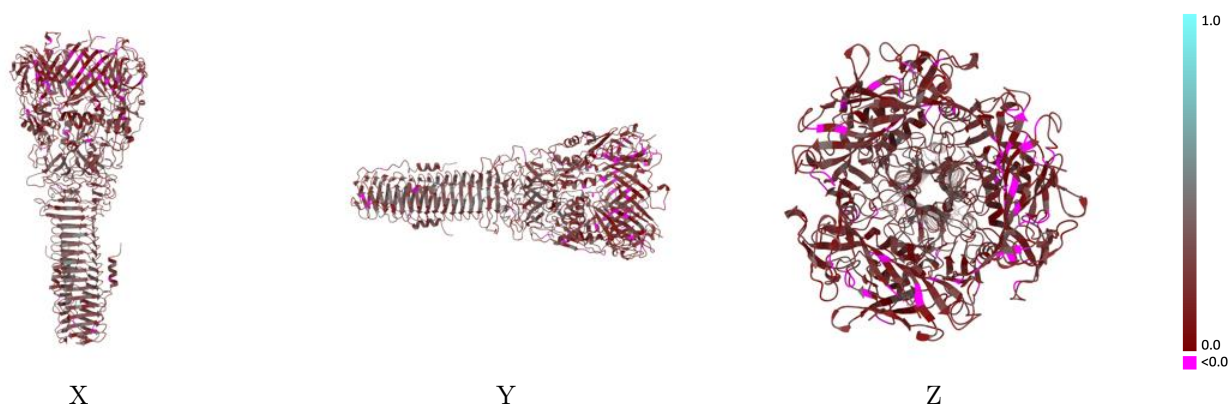
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.053 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](#)

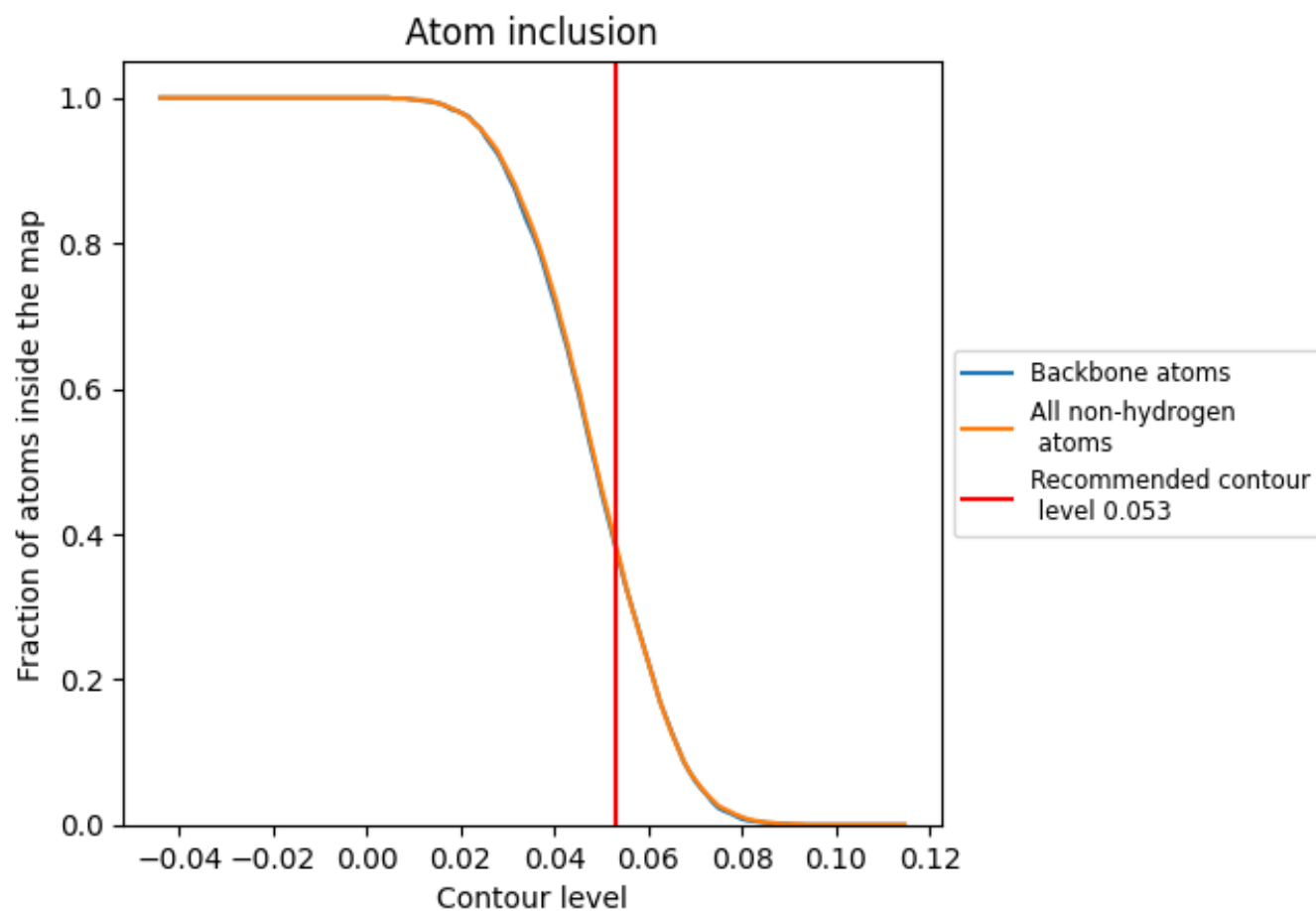


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 38% of all backbone atoms, 39% 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 (0.053) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.3860	0.2250
A	0.3970	0.2260
B	0.4070	0.2340
C	0.3780	0.2160

