



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 03:16 PM UTC

PDB ID : 5KU2 / pdb_00005ku2
EMDB ID : EMD-8286
Title : expanded poliovirus in complex with VHH 7A
Authors : Strauss, M.; Schotte, L.; Filman, D.J.; Hogle, J.M.
Deposited on : 2016-07-12
Resolution : 5.30 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

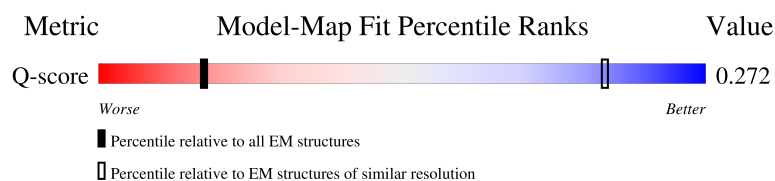
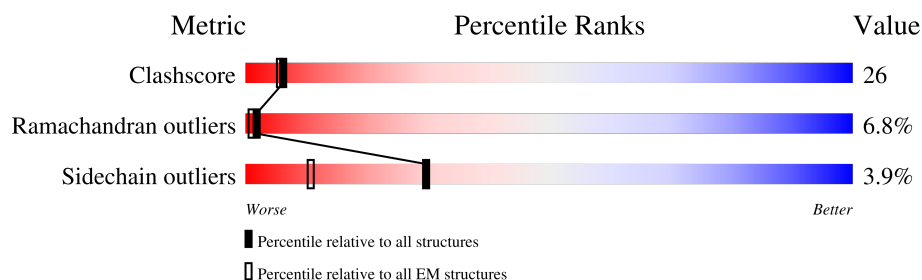
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 5.30 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	625 (4.80 - 5.80)

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	1	209	
2	2	268	
3	3	230	
4	7	125	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5994 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 VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	191	Total	C	N	O	S	0	0
			1543	996	261	281	5		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	228	ILE	LEU	conflict	UNP P03300

- Molecule 2 is a protein called VP2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	2	235	Total	C	N	O	S	0	0
			1822	1160	310	339	13		

- Molecule 3 is a protein called VP3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	221	Total	C	N	O	S	0	0
			1712	1093	279	323	17		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
3	123	SER	PHE	conflict	UNP P03300

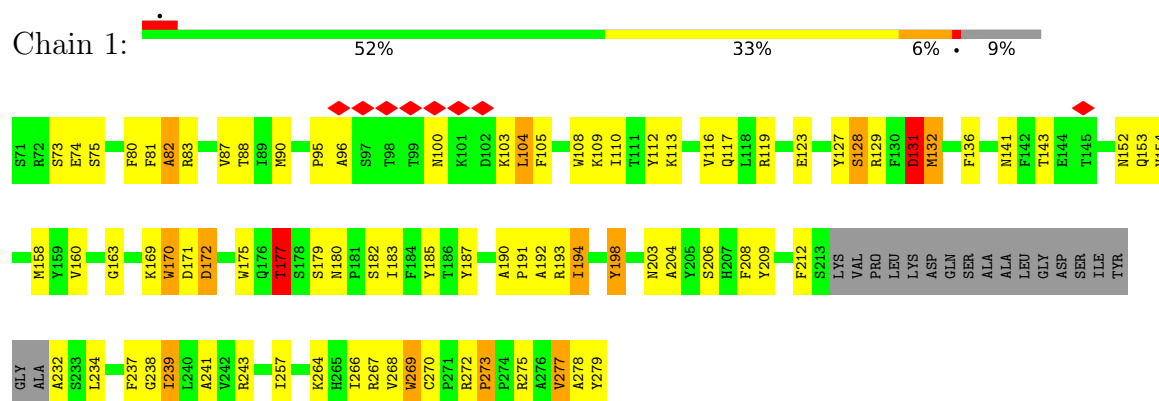
- Molecule 4 is a protein called VHH 7A.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	7	124	Total	C	N	O	S	0	0
			917	567	159	186	5		

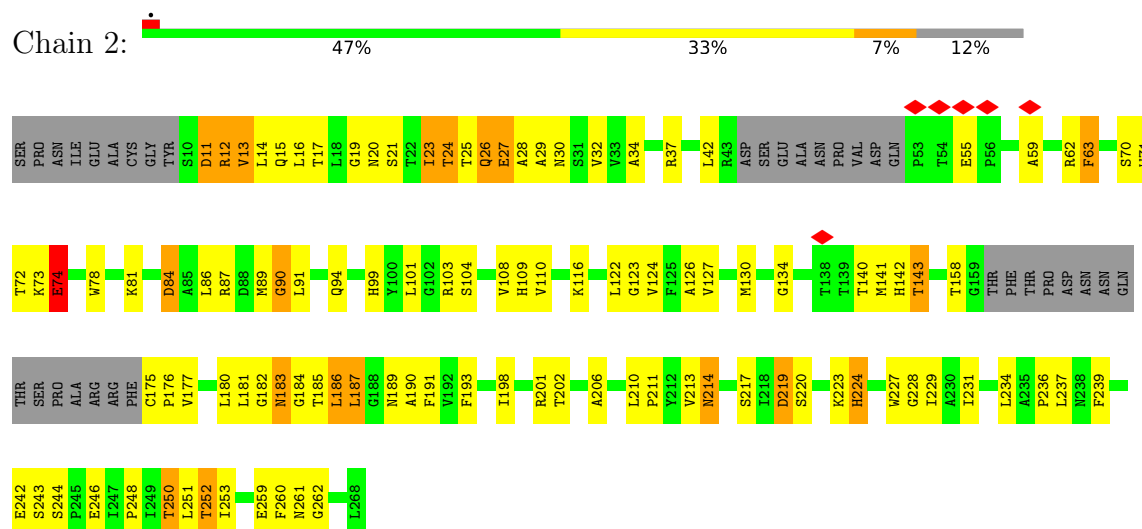
3 Residue-property plots

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.

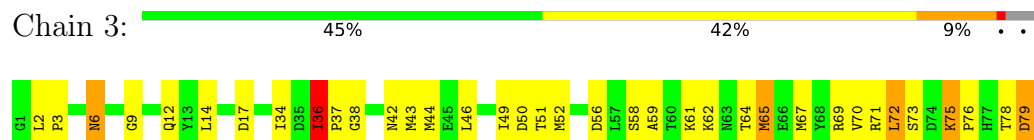
• Molecule 1: VP1

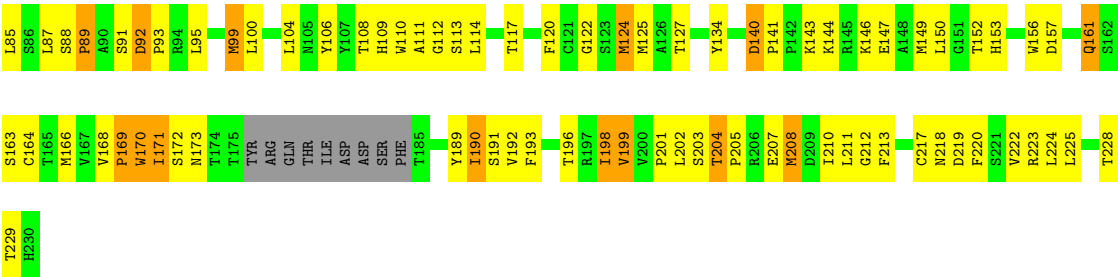


• Molecule 2: VP2

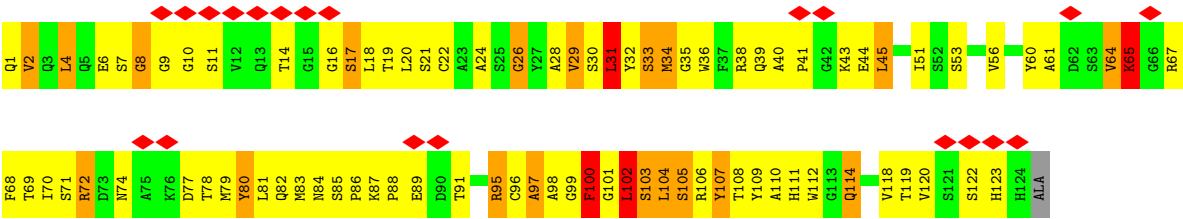


• Molecule 3: VP3





● Molecule 4: VHH 7A



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	17654	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	30	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.014	Depositor
Minimum map value	-0.004	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.00481	Depositor
Map size (Å)	422.144, 422.144, 422.144	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8245, 0.8245, 0.8245	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	1	0.78	0/1590	1.35	19/2169 (0.9%)
2	2	0.86	1/1870 (0.1%)	1.36	16/2553 (0.6%)
3	3	0.89	2/1756 (0.1%)	1.53	32/2392 (1.3%)
4	7	1.13	1/935 (0.1%)	1.71	28/1265 (2.2%)
All	All	0.90	4/6151 (0.1%)	1.46	95/8379 (1.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	3	196	THR	CA-C	6.64	1.56	1.52
2	2	202	THR	CA-C	6.43	1.55	1.52
3	3	172	SER	N-CA	5.23	1.52	1.46
4	7	85	SER	CA-CB	5.14	1.56	1.52

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	3	36	ILE	N-CA-C	-14.21	96.41	108.63
3	3	170	TRP	N-CA-C	10.08	124.68	109.25
4	7	101	GLY	N-CA-C	9.91	125.94	111.76
3	3	143	LYS	N-CA-C	9.54	122.68	111.71
3	3	80	ASP	CA-C-N	9.52	131.75	119.84
3	3	80	ASP	C-N-CA	9.52	131.75	119.84
4	7	78	THR	N-CA-C	9.16	123.16	110.24
3	3	82	ILE	N-CA-C	-8.77	102.29	110.53
4	7	64	VAL	N-CA-C	8.59	124.16	113.00
4	7	29	VAL	N-CA-C	-8.39	100.53	112.35
2	2	223	LYS	N-CA-C	8.35	120.46	111.36
1	1	268	VAL	N-CA-C	7.41	119.50	108.46
4	7	123	HIS	N-CA-C	7.04	120.15	108.96
1	1	269	TRP	N-CA-C	6.98	119.78	108.41
1	1	112	TYR	N-CA-C	-6.92	104.98	113.50
3	3	65	MET	N-CA-C	6.92	121.33	112.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	3	99	MET	N-CA-C	-6.75	103.85	111.07
3	3	173	ASN	N-CA-C	6.72	120.84	110.42
2	2	26	GLN	N-CA-C	-6.70	102.39	112.04
1	1	80	PHE	N-CA-C	-6.70	104.13	112.90
3	3	92	ASP	CA-C-N	6.67	128.18	119.84
3	3	92	ASP	C-N-CA	6.67	128.18	119.84
4	7	9	GLY	N-CA-C	6.63	124.41	112.54
4	7	31	LEU	CA-C-N	6.62	130.91	122.16
4	7	31	LEU	C-N-CA	6.62	130.91	122.16
3	3	198	ILE	N-CA-C	6.59	118.30	108.54
3	3	108	THR	N-CA-C	-6.58	105.78	113.88
4	7	26	GLY	CA-C-N	6.57	130.67	120.82
4	7	26	GLY	C-N-CA	6.57	130.67	120.82
1	1	177	THR	N-CA-C	6.56	124.76	110.80
1	1	131	ASP	N-CA-C	-6.52	100.00	109.59
3	3	225	LEU	N-CA-C	6.41	118.60	108.67
4	7	35	GLY	N-CA-C	6.35	120.08	110.18
1	1	128	SER	N-CA-C	6.34	118.49	108.79
2	2	63	PHE	N-CA-C	6.33	119.55	110.24
3	3	222	VAL	N-CA-C	6.28	118.17	108.44
3	3	140	ASP	CA-C-N	-6.22	113.97	120.38
3	3	140	ASP	C-N-CA	-6.22	113.97	120.38
2	2	224	HIS	N-CA-C	6.22	119.08	109.07
1	1	272	ARG	CA-C-N	-6.10	114.10	120.38
1	1	272	ARG	C-N-CA	-6.10	114.10	120.38
4	7	4	LEU	N-CA-C	6.09	119.56	109.46
1	1	143	THR	N-CA-C	6.03	117.52	111.07
3	3	124	MET	N-CA-C	-6.02	104.80	111.36
4	7	18	LEU	N-CA-C	5.99	118.39	109.59
1	1	190	ALA	CA-C-N	5.98	127.31	119.84
1	1	190	ALA	C-N-CA	5.98	127.31	119.84
1	1	132	MET	N-CA-C	5.89	118.71	107.75
4	7	10	GLY	N-CA-C	5.87	117.90	110.38
2	2	219	ASP	N-CA-C	5.87	117.99	109.07
1	1	153	GLN	N-CA-C	5.80	118.31	109.95
1	1	237	PHE	N-CA-C	-5.79	105.88	113.17
1	1	193	ARG	N-CA-C	5.76	118.85	108.48
2	2	104	SER	CA-C-N	5.72	126.63	121.82
2	2	104	SER	C-N-CA	5.72	126.63	121.82
4	7	74	ASN	CA-C-N	5.70	127.92	120.28
4	7	74	ASN	C-N-CA	5.70	127.92	120.28
4	7	44	GLU	N-CA-C	5.70	118.54	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	3	163	SER	N-CA-C	5.70	118.08	107.99
3	3	204	THR	CA-C-N	5.69	126.95	119.84
3	3	204	THR	C-N-CA	5.69	126.95	119.84
3	3	106	TYR	N-CA-C	5.65	119.16	112.72
2	2	252	THR	N-CA-C	-5.63	99.72	108.90
4	7	34	MET	N-CA-C	5.62	116.49	108.74
3	3	157	ASP	CA-C-N	5.61	127.75	120.56
3	3	157	ASP	C-N-CA	5.61	127.75	120.56
4	7	98	ALA	N-CA-C	5.61	119.70	112.41
2	2	32	VAL	N-CA-C	5.59	116.18	108.12
2	2	183	ASN	N-CA-C	5.58	119.74	111.87
3	3	111	ALA	N-CA-C	5.58	117.59	108.55
4	7	56	VAL	N-CA-C	5.58	116.80	112.12
3	3	191	SER	N-CA-C	5.57	117.87	109.23
1	1	273	PRO	N-CA-C	-5.57	103.91	110.70
1	1	172	ASP	N-CA-C	5.53	122.57	110.80
2	2	250	THR	N-CA-C	5.51	118.08	108.76
3	3	172	SER	N-CA-C	5.47	117.42	108.55
2	2	70	SER	N-CA-C	5.45	117.52	108.20
3	3	79	ASP	N-CA-C	5.42	119.38	112.87
1	1	96	ALA	N-CA-C	-5.42	106.51	113.23
3	3	80	ASP	N-CA-C	-5.40	98.74	109.60
4	7	8	GLY	N-CA-C	5.36	121.89	114.92
4	7	80	TYR	N-CA-C	5.34	117.33	108.20
4	7	104	LEU	N-CA-C	5.32	118.23	109.72
3	3	72	LEU	N-CA-C	-5.28	97.92	107.75
2	2	74	GLU	N-CA-C	5.25	118.95	112.54
4	7	65	LYS	N-CA-CB	5.25	118.27	110.29
3	3	78	THR	N-CA-C	-5.25	99.96	108.41
2	2	185	THR	N-CA-C	5.23	117.92	107.62
3	3	223	ARG	N-CA-C	5.20	117.16	108.99
2	2	23	ILE	N-CA-C	5.18	120.12	109.34
4	7	30	SER	N-CA-C	-5.18	107.02	113.19
4	7	105	SER	N-CA-C	-5.09	107.03	113.20
2	2	24	THR	N-CA-CB	5.04	119.02	110.49
4	7	6	GLU	N-CA-C	5.04	117.52	109.96
4	7	31	LEU	CA-CB-CG	5.03	133.89	116.30

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	1	1543	0	1503	59	0
2	2	1822	0	1762	87	0
3	3	1712	0	1697	101	0
4	7	917	0	872	100	0
All	All	5994	0	5834	306	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 26.

All (306) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:198:ILE:HG21	3:3:208:MET:HE1	1.31	1.13
1:1:209:TYR:CE1	2:2:224:HIS:HD2	1.66	1.13
3:3:56:ASP:OD2	4:7:106:ARG:HD3	1.50	1.09
1:1:209:TYR:HE1	2:2:224:HIS:CD2	1.78	1.01
1:1:109:LYS:HA	1:1:239:ILE:HG22	1.44	0.99
3:3:125:MET:HB3	3:3:201:PRO:HG3	1.49	0.94
1:1:209:TYR:CE1	2:2:224:HIS:CD2	2.55	0.93
1:1:209:TYR:HE1	2:2:224:HIS:HD2	0.91	0.88
4:7:14:THR:H	4:7:122:SER:HB3	1.41	0.85
2:2:201:ARG:HH11	3:3:124:MET:HA	1.43	0.84
3:3:84:CYS:SG	4:7:107:TYR:CD1	2.71	0.83
2:2:101:LEU:HB2	2:2:261:ASN:HB2	1.60	0.83
2:2:201:ARG:NH1	3:3:124:MET:HA	1.95	0.82
2:2:158:THR:HG22	2:2:176:PRO:HD2	1.62	0.81
4:7:102:LEU:HD13	4:7:103:SER:H	1.44	0.81
1:1:136:PHE:HB2	1:1:192:ALA:HB3	1.62	0.81
4:7:11:SER:HA	4:7:119:THR:O	1.80	0.80
2:2:122:LEU:HB2	2:2:198:ILE:HB	1.64	0.80
4:7:87:LYS:HB3	4:7:88:PRO:HD2	1.63	0.79
3:3:198:ILE:CG2	3:3:208:MET:HE1	2.12	0.78
4:7:16:GLY:H	4:7:86:PRO:HD2	1.47	0.78
1:1:269:TRP:HA	3:3:38:GLY:O	1.83	0.77
4:7:20:LEU:HD12	4:7:81:LEU:HD23	1.64	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:15:GLN:HB3	2:2:24:THR:HG23	1.67	0.76
2:2:74:GLU:H	2:2:74:GLU:CD	1.90	0.76
3:3:70:VAL:HG13	4:7:106:ARG:HB3	1.66	0.75
3:3:50:ASP:HB3	3:3:213:PHE:HB3	1.69	0.74
4:7:14:THR:N	4:7:122:SER:HB3	2.02	0.74
4:7:14:THR:HG23	4:7:87:LYS:HA	1.67	0.74
2:2:116:LYS:O	3:3:125:MET:HG2	1.89	0.73
3:3:125:MET:CB	3:3:201:PRO:HG3	2.19	0.72
2:2:110:VAL:HG22	2:2:251:LEU:HG	1.72	0.72
2:2:23:ILE:HD11	2:2:63:PHE:CE1	2.24	0.72
1:1:183:ILE:HD11	1:1:194:ILE:CG2	2.20	0.71
2:2:213:VAL:HG22	3:3:37:PRO:HD2	1.72	0.71
3:3:51:THR:HG21	3:3:99:MET:HB2	1.72	0.71
4:7:31:LEU:CD2	4:7:32:TYR:H	2.03	0.71
4:7:109:TYR:CE2	4:7:111:HIS:CE1	2.78	0.71
3:3:88:SER:HB3	3:3:91:SER:OG	1.91	0.70
2:2:180:LEU:HD13	2:2:184:GLY:HA2	1.72	0.70
4:7:38:ARG:O	4:7:45:LEU:HA	1.92	0.69
3:3:100:LEU:O	3:3:104:LEU:HD13	1.92	0.69
1:1:232:ALA:HB2	2:2:140:THR:OG1	1.94	0.67
4:7:32:TYR:CD2	4:7:100:PHE:HB2	2.28	0.67
4:7:29:VAL:HA	4:7:31:LEU:HD13	1.77	0.66
4:7:72:ARG:HA	4:7:79:MET:HA	1.78	0.66
3:3:58:SER:HB3	4:7:105:SER:OG	1.96	0.66
2:2:37:ARG:HG3	3:3:37:PRO:HB3	1.77	0.66
3:3:146:LYS:O	3:3:150:LEU:HG	1.96	0.66
2:2:189:ASN:C	2:2:191:PHE:H	2.03	0.66
4:7:107:TYR:HD1	4:7:108:THR:H	1.45	0.65
1:1:183:ILE:HD11	1:1:194:ILE:HG23	1.78	0.65
3:3:79:ASP:HB2	3:3:144:LYS:HD2	1.78	0.65
4:7:39:GLN:HB3	4:7:45:LEU:HB3	1.77	0.65
1:1:116:VAL:HB	3:3:229:THR:CG2	2.27	0.64
4:7:67:ARG:HG3	4:7:68:PHE:HD1	1.62	0.64
3:3:125:MET:HB3	3:3:201:PRO:CG	2.26	0.64
1:1:131:ASP:HB2	1:1:264:LYS:HG3	1.80	0.63
3:3:189:TYR:CD1	4:7:100:PHE:HE2	2.16	0.63
2:2:55:GLU:HG2	2:2:259:GLU:HB3	1.79	0.63
1:1:116:VAL:HB	3:3:229:THR:HG22	1.79	0.63
4:7:103:SER:HB2	4:7:107:TYR:OH	1.99	0.62
4:7:105:SER:C	4:7:107:TYR:H	2.07	0.62
2:2:11:ASP:CG	2:2:12:ARG:H	2.06	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:7:102:LEU:CD1	4:7:103:SER:H	2.11	0.62
4:7:20:LEU:HB2	4:7:81:LEU:HB3	1.82	0.61
4:7:31:LEU:HD23	4:7:32:TYR:H	1.64	0.61
2:2:234:LEU:HB3	3:3:211:LEU:HD11	1.83	0.60
2:2:42:LEU:HB2	2:2:103:ARG:HD3	1.83	0.60
1:1:177:THR:OG1	1:1:180:ASN:N	2.35	0.60
2:2:20:ASN:HB3	2:2:62:ARG:NE	2.17	0.60
4:7:33:SER:O	4:7:99:GLY:HA2	2.02	0.59
2:2:103:ARG:HD2	2:2:217:SER:O	2.02	0.59
4:7:99:GLY:N	4:7:109:TYR:HB2	2.17	0.59
2:2:13:VAL:HA	2:2:27:GLU:HA	1.85	0.59
3:3:14:LEU:HB3	3:3:17:ASP:HB3	1.84	0.59
4:7:51:ILE:HD12	4:7:71:SER:HA	1.84	0.59
3:3:84:CYS:HB3	4:7:100:PHE:CE1	2.37	0.59
3:3:89:PRO:HA	3:3:95:LEU:HB3	1.85	0.59
2:2:116:LYS:CB	3:3:125:MET:HE2	2.33	0.59
1:1:212:PHE:CE2	2:2:141:MET:HB2	2.38	0.58
3:3:110:TRP:HB2	3:3:220:PHE:CZ	2.38	0.58
4:7:95:ARG:HH11	4:7:95:ARG:CG	2.15	0.58
3:3:140:ASP:HB3	3:3:141:PRO:HD2	1.84	0.58
4:7:91:THR:HG23	4:7:119:THR:HA	1.84	0.58
2:2:181:LEU:HG	2:2:183:ASN:HD21	1.69	0.58
2:2:42:LEU:HD11	2:2:217:SER:HB3	1.86	0.58
2:2:17:THR:HA	2:2:21:SER:O	2.04	0.58
2:2:110:VAL:HB	2:2:206:ALA:HB3	1.86	0.58
3:3:79:ASP:HB2	3:3:144:LYS:CD	2.34	0.57
1:1:82:ALA:C	1:1:83:ARG:HD2	2.29	0.57
4:7:107:TYR:CD1	4:7:108:THR:N	2.72	0.57
3:3:100:LEU:HD21	3:3:114:LEU:HD22	1.86	0.57
3:3:189:TYR:CD1	4:7:100:PHE:CE2	2.92	0.57
3:3:84:CYS:HB3	4:7:100:PHE:CD1	2.40	0.57
4:7:34:MET:HE2	4:7:99:GLY:N	2.20	0.57
1:1:152:ASN:OD1	1:1:187:TYR:HB3	2.05	0.56
4:7:97:ALA:HB1	4:7:111:HIS:O	2.05	0.56
2:2:108:VAL:HA	2:2:253:ILE:HG12	1.87	0.56
3:3:85:LEU:HD22	3:3:190:ILE:HD13	1.87	0.56
2:2:187:LEU:HD23	3:3:65:MET:SD	2.46	0.56
3:3:127:THR:OG1	3:3:199:VAL:HG22	2.06	0.56
2:2:116:LYS:HB2	3:3:125:MET:HE2	1.88	0.56
2:2:189:ASN:O	2:2:191:PHE:N	2.39	0.56
3:3:201:PRO:HD2	3:3:204:THR:HG21	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:7:95:ARG:NH1	4:7:114:GLN:O	2.39	0.55
2:2:42:LEU:HD13	2:2:103:ARG:HH11	1.71	0.55
3:3:147:GLU:HA	3:3:150:LEU:HD12	1.87	0.55
4:7:32:TYR:O	4:7:53:SER:HA	2.06	0.55
2:2:16:LEU:HB2	2:2:23:ILE:HD12	1.87	0.55
2:2:72:THR:HG22	2:2:246:GLU:HG2	1.88	0.54
2:2:177:VAL:HG23	2:2:180:LEU:HB2	1.89	0.54
4:7:29:VAL:CA	4:7:31:LEU:HD13	2.37	0.54
4:7:28:ALA:O	4:7:31:LEU:HB3	2.07	0.54
4:7:34:MET:HG3	4:7:79:MET:HE2	1.90	0.54
2:2:219:ASP:OD1	2:2:220:SER:N	2.39	0.54
2:2:189:ASN:C	2:2:191:PHE:N	2.65	0.54
3:3:52:MET:HA	3:3:212:GLY:O	2.08	0.54
1:1:154:VAL:HA	1:1:185:TYR:O	2.07	0.54
4:7:45:LEU:HD23	4:7:45:LEU:H	1.72	0.54
4:7:2:VAL:HA	4:7:26:GLY:HA3	1.90	0.54
1:1:103:LYS:O	1:1:105:PHE:N	2.41	0.54
2:2:81:LYS:HB2	2:2:84:ASP:HB3	1.89	0.54
3:3:59:ALA:HB2	4:7:95:ARG:HG2	1.89	0.53
3:3:79:ASP:O	3:3:193:PHE:HE2	1.90	0.53
3:3:83:LEU:HG	3:3:192:VAL:HB	1.90	0.53
4:7:2:VAL:HG12	4:7:26:GLY:HA3	1.91	0.53
4:7:60:TYR:HB2	4:7:65:LYS:HG3	1.91	0.53
2:2:213:VAL:O	2:2:214:ASN:HB2	2.09	0.53
3:3:58:SER:HB2	4:7:45:LEU:HD11	1.90	0.53
1:1:158:MET:HE1	1:1:175:TRP:O	2.09	0.53
4:7:8:GLY:HA3	4:7:20:LEU:HA	1.89	0.53
1:1:160:VAL:O	1:1:238:GLY:HA3	2.09	0.52
1:1:212:PHE:HZ	2:2:143:THR:O	1.91	0.52
3:3:87:LEU:HD21	3:3:171:ILE:HD13	1.92	0.52
4:7:61:ALA:HB3	4:7:64:VAL:HG22	1.90	0.52
4:7:40:ALA:HB3	4:7:43:LYS:HB2	1.91	0.52
4:7:97:ALA:HB3	4:7:112:TRP:HA	1.91	0.51
3:3:89:PRO:HD2	3:3:110:TRP:CZ2	2.45	0.51
2:2:20:ASN:HB3	2:2:62:ARG:HE	1.75	0.51
1:1:103:LYS:HD3	1:1:170:TRP:CD1	2.45	0.51
3:3:51:THR:HG21	3:3:99:MET:CB	2.40	0.51
2:2:90:GLY:O	2:2:94:GLN:HG2	2.11	0.51
3:3:6:ASN:N	3:3:6:ASN:HD22	2.09	0.51
4:7:28:ALA:O	4:7:31:LEU:CD1	2.59	0.51
2:2:177:VAL:N	2:2:182:GLY:HA2	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:273:PRO:HB3	2:2:189:ASN:HB2	1.93	0.51
3:3:189:TYR:HD1	4:7:100:PHE:HE2	1.56	0.51
4:7:31:LEU:HG	4:7:109:TYR:CZ	2.46	0.51
1:1:100:ASN:HA	1:1:103:LYS:HB2	1.93	0.50
1:1:158:MET:HB2	1:1:241:ALA:HB3	1.93	0.50
3:3:110:TRP:CE3	3:3:220:PHE:HE1	2.29	0.50
2:2:244:SER:C	2:2:246:GLU:H	2.20	0.50
3:3:72:LEU:HD21	3:3:210:ILE:HG23	1.92	0.50
3:3:89:PRO:HD2	3:3:110:TRP:HZ2	1.76	0.50
3:3:149:MET:O	3:3:153:HIS:CE1	2.64	0.50
3:3:202:LEU:HG	3:3:203:SER:H	1.77	0.50
4:7:95:ARG:HH11	4:7:95:ARG:HG3	1.77	0.50
2:2:34:ALA:HB3	2:2:211:PRO:HD2	1.94	0.50
2:2:59:ALA:HB1	2:2:91:LEU:HD13	1.94	0.49
3:3:44:MET:HA	3:3:44:MET:HE2	1.93	0.49
3:3:46:LEU:HB3	3:3:99:MET:HE3	1.94	0.49
2:2:11:ASP:O	2:2:13:VAL:N	2.45	0.49
3:3:64:THR:O	3:3:67:MET:HG2	2.12	0.49
1:1:169:LYS:C	1:1:171:ASP:H	2.20	0.49
1:1:273:PRO:HB3	2:2:189:ASN:CB	2.43	0.49
2:2:210:LEU:HD11	2:2:229:ILE:HD11	1.94	0.49
4:7:1:GLN:HG2	4:7:2:VAL:N	2.27	0.49
2:2:13:VAL:HB	2:2:26:GLN:O	2.12	0.49
1:1:87:VAL:HG23	1:1:88:THR:HG22	1.94	0.49
2:2:110:VAL:HA	2:2:250:THR:O	2.13	0.49
3:3:112:GLY:HA2	3:3:170:TRP:CZ3	2.47	0.49
2:2:71:TRP:NE1	2:2:237:LEU:HD11	2.28	0.49
4:7:31:LEU:HB2	4:7:32:TYR:HD1	1.78	0.49
1:1:73:SER:C	1:1:75:SER:N	2.70	0.48
4:7:99:GLY:O	4:7:100:PHE:CB	2.61	0.48
3:3:120:PHE:CE2	3:3:122:GLY:HA3	2.47	0.48
3:3:141:PRO:HB3	4:7:100:PHE:CZ	2.48	0.48
3:3:109:HIS:HE1	3:3:224:LEU:HD13	1.77	0.48
4:7:17:SER:OG	4:7:82:GLN:NE2	2.47	0.48
3:3:127:THR:OG1	3:3:199:VAL:CG2	2.61	0.48
1:1:103:LYS:HD3	1:1:170:TRP:HD1	1.79	0.48
3:3:87:LEU:HD11	3:3:114:LEU:HD12	1.94	0.48
4:7:87:LYS:HB3	4:7:88:PRO:CD	2.41	0.48
2:2:127:VAL:HG13	2:2:193:PHE:CE1	2.49	0.47
3:3:82:ILE:HD13	3:3:83:LEU:N	2.29	0.47
2:2:210:LEU:HD13	2:2:227:TRP:CD2	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:81:PHE:C	1:1:83:ARG:H	2.21	0.47
1:1:116:VAL:HG23	1:1:117:GLN:H	1.79	0.47
1:1:158:MET:HE1	1:1:175:TRP:HA	1.95	0.47
3:3:9:GLY:O	3:3:12:GLN:HG2	2.14	0.47
3:3:75:LYS:HB2	3:3:76:PRO:CD	2.45	0.47
4:7:32:TYR:HB2	4:7:99:GLY:HA3	1.96	0.47
3:3:42:ASN:OD1	3:3:43:MET:N	2.47	0.47
3:3:61:LYS:HZ2	4:7:106:ARG:HD2	1.80	0.47
3:3:110:TRP:CE3	3:3:220:PHE:CE1	3.03	0.47
2:2:130:MET:HB2	2:2:228:GLY:HA3	1.97	0.47
1:1:104:LEU:HD12	1:1:170:TRP:CE2	2.50	0.47
4:7:51:ILE:HD13	4:7:72:ARG:H	1.79	0.47
3:3:46:LEU:HB3	3:3:99:MET:CE	2.45	0.46
4:7:33:SER:C	4:7:99:GLY:HA2	2.40	0.46
4:7:29:VAL:C	4:7:31:LEU:H	2.23	0.46
4:7:51:ILE:CD1	4:7:71:SER:HA	2.45	0.46
1:1:127:TYR:HB2	1:1:269:TRP:HB2	1.96	0.46
2:2:34:ALA:HB2	2:2:210:LEU:HD23	1.97	0.46
3:3:85:LEU:HD21	3:3:95:LEU:HD11	1.97	0.46
3:3:156:TRP:HE1	3:3:161:GLN:HE22	1.64	0.46
1:1:177:THR:C	1:1:179:SER:H	2.23	0.46
4:7:107:TYR:HD1	4:7:108:THR:N	2.10	0.46
1:1:183:ILE:HD11	1:1:194:ILE:HG21	1.98	0.46
3:3:73:SER:HA	3:3:207:GLU:HG2	1.97	0.46
1:1:198:TYR:OH	1:1:204:ALA:N	2.48	0.46
2:2:86:LEU:HA	2:2:89:MET:HG2	1.97	0.46
3:3:156:TRP:CD1	3:3:164:CYS:HB2	2.51	0.46
3:3:134:TYR:CE2	3:3:169:PRO:HD2	2.51	0.45
3:3:190:ILE:N	3:3:190:ILE:HD12	2.30	0.45
2:2:71:TRP:HA	2:2:78:TRP:HH2	1.80	0.45
3:3:113:SER:HB3	3:3:168:VAL:O	2.16	0.45
2:2:63:PHE:HD2	2:2:252:THR:HG22	1.81	0.45
2:2:126:ALA:O	2:2:193:PHE:HB3	2.17	0.45
4:7:70:ILE:HA	4:7:80:TYR:O	2.17	0.45
3:3:2:LEU:HD12	3:3:3:PRO:HD2	1.99	0.45
1:1:90:MET:SD	1:1:108:TRP:HB3	2.57	0.45
2:2:158:THR:HG23	2:2:175:CYS:SG	2.57	0.45
2:2:186:LEU:HD22	3:3:67:MET:HE1	1.99	0.45
4:7:34:MET:HE2	4:7:34:MET:HA	1.99	0.45
3:3:82:ILE:O	4:7:107:TYR:CD2	2.70	0.45
4:7:34:MET:CE	4:7:99:GLY:H	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:7:36:TRP:CG	4:7:81:LEU:HD22	2.52	0.44
3:3:46:LEU:HA	3:3:49:ILE:HD12	1.99	0.44
3:3:117:THR:HA	3:3:164:CYS:O	2.17	0.44
1:1:128:SER:HA	1:1:267:ARG:O	2.16	0.44
2:2:21:SER:HB2	2:2:63:PHE:HB2	2.00	0.44
3:3:44:MET:SD	3:3:220:PHE:HD2	2.41	0.44
3:3:152:THR:HG22	3:3:166:MET:HE3	1.99	0.44
1:1:73:SER:C	1:1:75:SER:H	2.25	0.44
1:1:206:SER:HB3	1:1:208:PHE:O	2.17	0.44
3:3:82:ILE:HD12	3:3:192:VAL:CG1	2.48	0.44
4:7:102:LEU:HD13	4:7:103:SER:N	2.21	0.44
4:7:4:LEU:HD22	4:7:22:CYS:SG	2.58	0.44
1:1:116:VAL:HB	3:3:229:THR:HG21	1.99	0.44
4:7:16:GLY:O	4:7:17:SER:HB2	2.17	0.43
4:7:79:MET:SD	4:7:79:MET:N	2.90	0.43
4:7:96:CYS:O	4:7:97:ALA:CB	2.65	0.43
1:1:203:ASN:ND2	3:3:34:ILE:HG23	2.32	0.43
2:2:84:ASP:O	2:2:87:ARG:HG2	2.18	0.43
3:3:89:PRO:HA	3:3:95:LEU:CB	2.48	0.43
4:7:97:ALA:CB	4:7:111:HIS:O	2.65	0.43
3:3:17:ASP:OD1	3:3:17:ASP:C	2.62	0.43
4:7:64:VAL:HB	4:7:68:PHE:HB2	2.01	0.43
1:1:108:TRP:CZ3	1:1:110:ILE:HA	2.53	0.43
1:1:129:ARG:O	1:1:266:ILE:HA	2.18	0.43
2:2:236:PRO:HB3	3:3:69:ARG:NH1	2.33	0.43
4:7:32:TYR:HB3	4:7:100:PHE:N	2.34	0.43
1:1:269:TRP:CD2	3:3:36:ILE:HD12	2.53	0.43
2:2:15:GLN:HA	2:2:24:THR:H	1.84	0.43
4:7:19:THR:HA	4:7:82:GLN:HA	2.01	0.43
2:2:181:LEU:HB3	2:2:183:ASN:OD1	2.18	0.43
3:3:92:ASP:HA	3:3:93:PRO:HD2	1.72	0.43
4:7:60:TYR:HB3	4:7:64:VAL:HG23	2.00	0.42
4:7:69:THR:HB	4:7:82:GLN:HB3	2.01	0.42
4:7:67:ARG:HD2	4:7:84:ASN:O	2.19	0.42
1:1:158:MET:HG2	1:1:182:SER:OG	2.18	0.42
3:3:71:ARG:HD3	3:3:207:GLU:HB3	2.00	0.42
2:2:71:TRP:CG	2:2:237:LEU:HD21	2.55	0.42
2:2:124:VAL:HG13	2:2:231:ILE:HG12	2.01	0.42
4:7:32:TYR:HB3	4:7:99:GLY:C	2.44	0.42
4:7:83:MET:HE1	4:7:118:VAL:HG11	2.02	0.42
4:7:89:GLU:C	4:7:91:THR:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:71:TRP:HA	2:2:78:TRP:CH2	2.54	0.42
2:2:109:HIS:HA	2:2:206:ALA:O	2.19	0.42
1:1:103:LYS:HG2	1:1:104:LEU:N	2.35	0.42
2:2:239:PHE:HE1	2:2:242:GLU:HA	1.85	0.42
4:7:111:HIS:CD2	4:7:111:HIS:H	2.37	0.42
4:7:105:SER:C	4:7:107:TYR:N	2.75	0.41
1:1:83:ARG:HD2	1:1:83:ARG:N	2.35	0.41
1:1:104:LEU:HD12	1:1:170:TRP:CD2	2.55	0.41
1:1:132:MET:HE2	1:1:132:MET:HB3	1.92	0.41
3:3:58:SER:O	3:3:62:LYS:N	2.50	0.41
3:3:82:ILE:HD13	3:3:82:ILE:C	2.44	0.41
3:3:208:MET:SD	3:3:208:MET:N	2.93	0.41
4:7:2:VAL:HG12	4:7:26:GLY:C	2.45	0.41
2:2:191:PHE:HE1	3:3:52:MET:HE2	1.85	0.41
1:1:73:SER:O	1:1:75:SER:N	2.53	0.41
1:1:113:LYS:HG3	1:1:119:ARG:NH1	2.36	0.41
2:2:99:HIS:CB	2:2:260:PHE:HB3	2.50	0.41
2:2:130:MET:CB	2:2:228:GLY:HA3	2.49	0.41
4:7:64:VAL:HB	4:7:68:PHE:CG	2.56	0.41
2:2:14:LEU:HD23	2:2:14:LEU:H	1.85	0.41
2:2:72:THR:O	2:2:237:LEU:HD13	2.20	0.41
2:2:110:VAL:O	2:2:206:ALA:N	2.54	0.41
2:2:123:GLY:O	2:2:231:ILE:HA	2.20	0.41
3:3:217:CYS:HB2	3:3:219:ASP:OD1	2.21	0.41
2:2:124:VAL:HG22	2:2:231:ILE:HG12	2.03	0.41
1:1:95:PRO:HD2	1:1:104:LEU:CD2	2.51	0.40
3:3:83:LEU:HD12	3:3:83:LEU:C	2.47	0.40
2:2:73:LYS:HB2	2:2:74:GLU:OE1	2.21	0.40
4:7:4:LEU:HB3	4:7:22:CYS:SG	2.62	0.40
4:7:14:THR:HA	4:7:120:VAL:CG1	2.52	0.40
4:7:67:ARG:HB2	4:7:84:ASN:O	2.21	0.40
1:1:277:VAL:HB	1:1:278:ALA:H	1.50	0.40
4:7:4:LEU:HD21	4:7:24:ALA:HB2	2.02	0.40
4:7:100:PHE:O	4:7:108:THR:HA	2.20	0.40
1:1:170:TRP:HE3	1:1:243:ARG:HH11	1.68	0.40
1:1:275:ARG:HG2	1:1:279:TYR:CZ	2.57	0.40
4:7:7:SER:OG	4:7:21:SER:HB2	2.21	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	1	187/209 (90%)	154 (82%)	19 (10%)	14 (8%)	1	9
2	2	229/268 (85%)	180 (79%)	29 (13%)	20 (9%)	0	8
3	3	217/230 (94%)	182 (84%)	30 (14%)	5 (2%)	5	26
4	7	122/125 (98%)	90 (74%)	20 (16%)	12 (10%)	0	7
All	All	755/832 (91%)	606 (80%)	98 (13%)	51 (7%)	2	11

All (51) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	104	LEU
1	1	172	ASP
1	1	177	THR
1	1	198	TYR
1	1	234	LEU
2	2	12	ARG
2	2	25	THR
2	2	27	GLU
2	2	30	ASN
2	2	142	HIS
3	3	81	PRO
4	7	17	SER
4	7	31	LEU
4	7	77	ASP
4	7	97	ALA
4	7	100	PHE
4	7	103	SER
4	7	110	ALA
2	2	11	ASP
2	2	29	ALA
2	2	90	GLY
2	2	134	GLY

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Mol	Chain	Res	Type
2	2	186	LEU
2	2	187	LEU
2	2	190	ALA
2	2	214	ASN
3	3	169	PRO
4	7	107	TYR
1	1	74	GLU
1	1	123	GLU
1	1	170	TRP
2	2	84	ASP
3	3	228	THR
4	7	114	GLN
1	1	82	ALA
1	1	141	ASN
1	1	191	PRO
1	1	270	CYS
2	2	28	ALA
2	2	243	SER
4	7	33	SER
1	1	277	VAL
2	2	19	GLY
2	2	262	GLY
3	3	89	PRO
3	3	205	PRO
4	7	102	LEU
1	1	163	GLY
2	2	143	THR
2	2	248	PRO
4	7	41	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	1	170/183 (93%)	166 (98%)	4 (2%)	43	62
2	2	198/228 (87%)	196 (99%)	2 (1%)	68	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	3	196/206 (95%)	185 (94%)	11 (6%)	19	40
4	7	96/97 (99%)	87 (91%)	9 (9%)	8	25
All	All	660/714 (92%)	634 (96%)	26 (4%)	30	49

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	131	ASP
1	1	194	ILE
1	1	239	ILE
1	1	257	ILE
2	2	13	VAL
2	2	74	GLU
3	3	6	ASN
3	3	36	ILE
3	3	75	LYS
3	3	82	ILE
3	3	83	LEU
3	3	161	GLN
3	3	171	ILE
3	3	190	ILE
3	3	199	VAL
3	3	208	MET
3	3	218	ASN
4	7	2	VAL
4	7	31	LEU
4	7	45	LEU
4	7	65	LYS
4	7	72	ARG
4	7	95	ARG
4	7	100	PHE
4	7	102	LEU
4	7	104	LEU

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

Mol	Chain	Res	Type
1	1	117	GLN
1	1	147	ASN
1	1	153	GLN

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Mol	Chain	Res	Type
1	1	265	HIS
2	2	111	GLN
2	2	137	ASN
2	2	199	ASN
2	2	224	HIS
3	3	6	ASN
3	3	97	HIS
3	3	161	GLN
4	7	13	GLN
4	7	82	GLN
4	7	111	HIS
4	7	117	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.

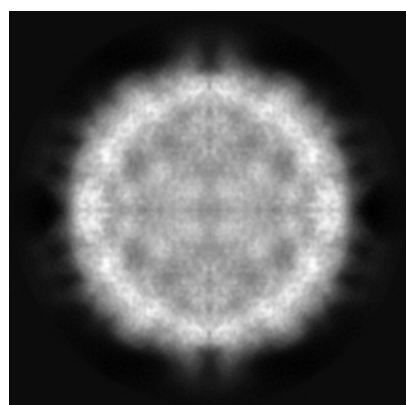
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8286. These allow visual inspection of the internal detail of the map and identification of artifacts.

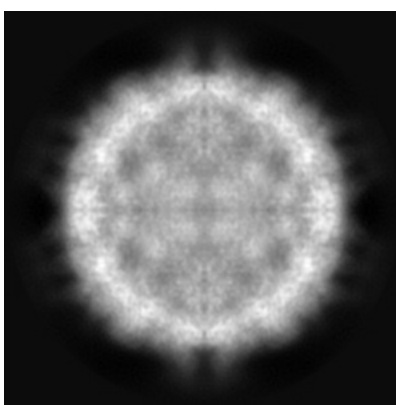
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

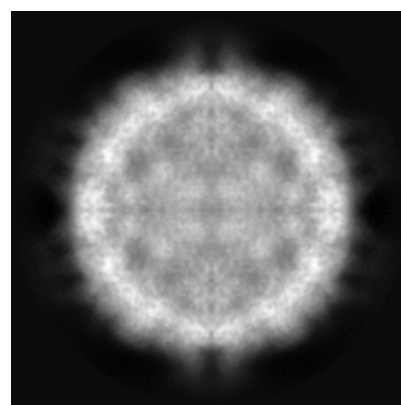
6.1.1 Primary map



X



Y

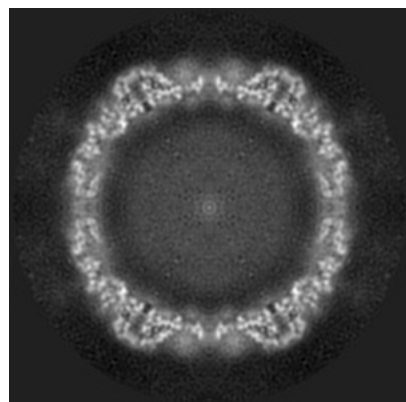


Z

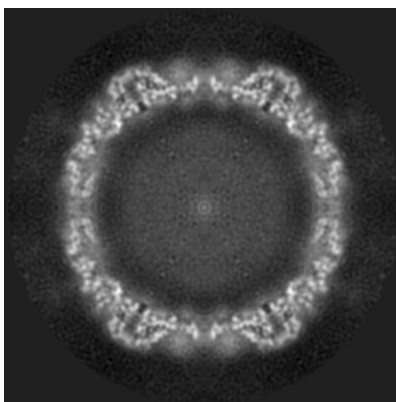
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

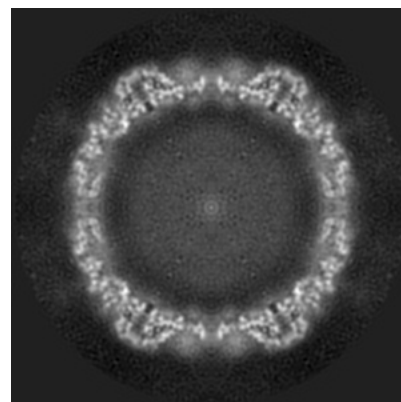
6.2.1 Primary map



X Index: 256



Y Index: 256

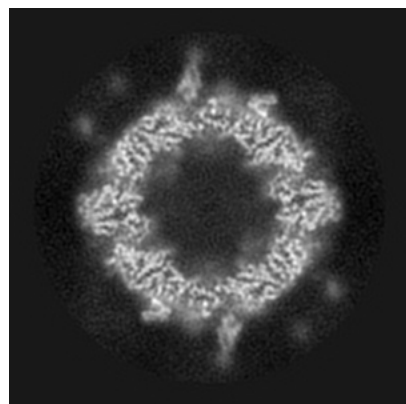


Z Index: 256

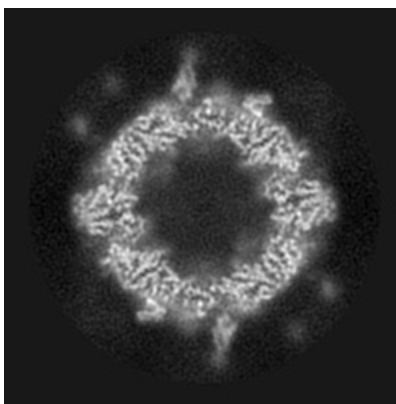
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

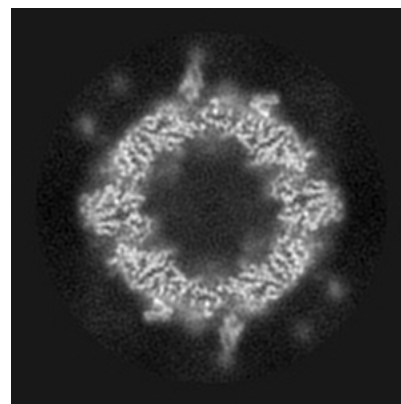
6.3.1 Primary map



X Index: 141



Y Index: 141

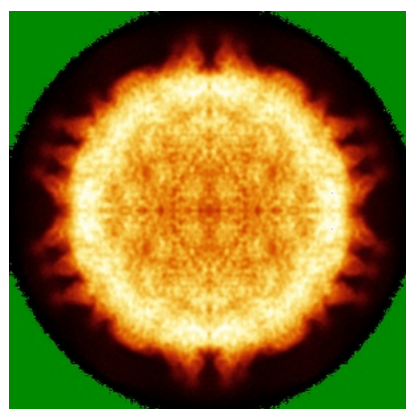


Z Index: 141

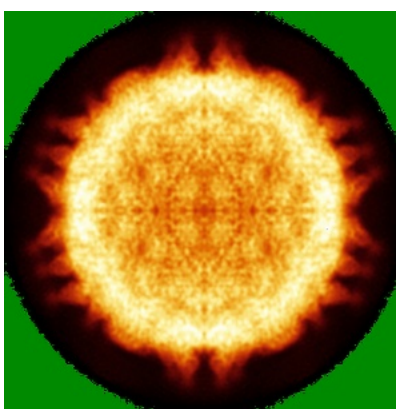
The images above show the largest variance slices of the map in three orthogonal directions.

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

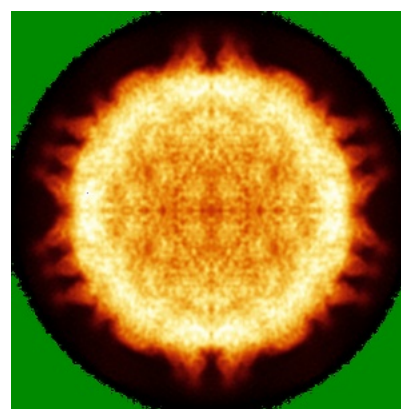
6.4.1 Primary map



X



Y

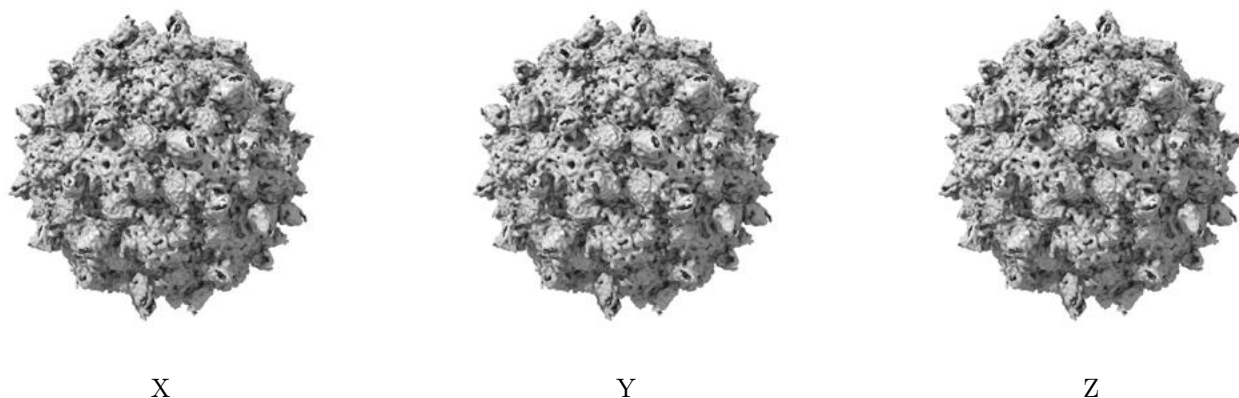


Z

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 0.00481. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

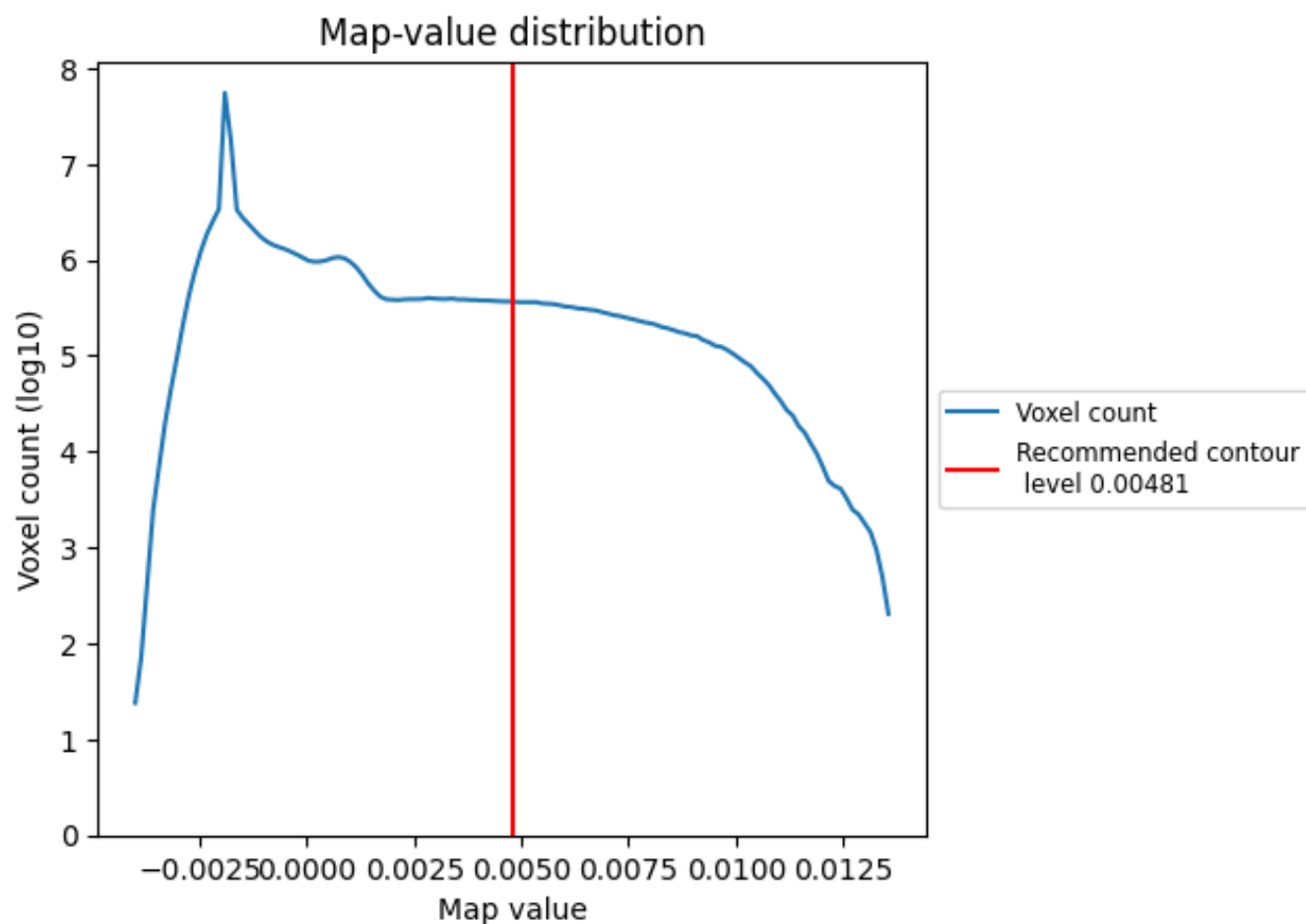
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

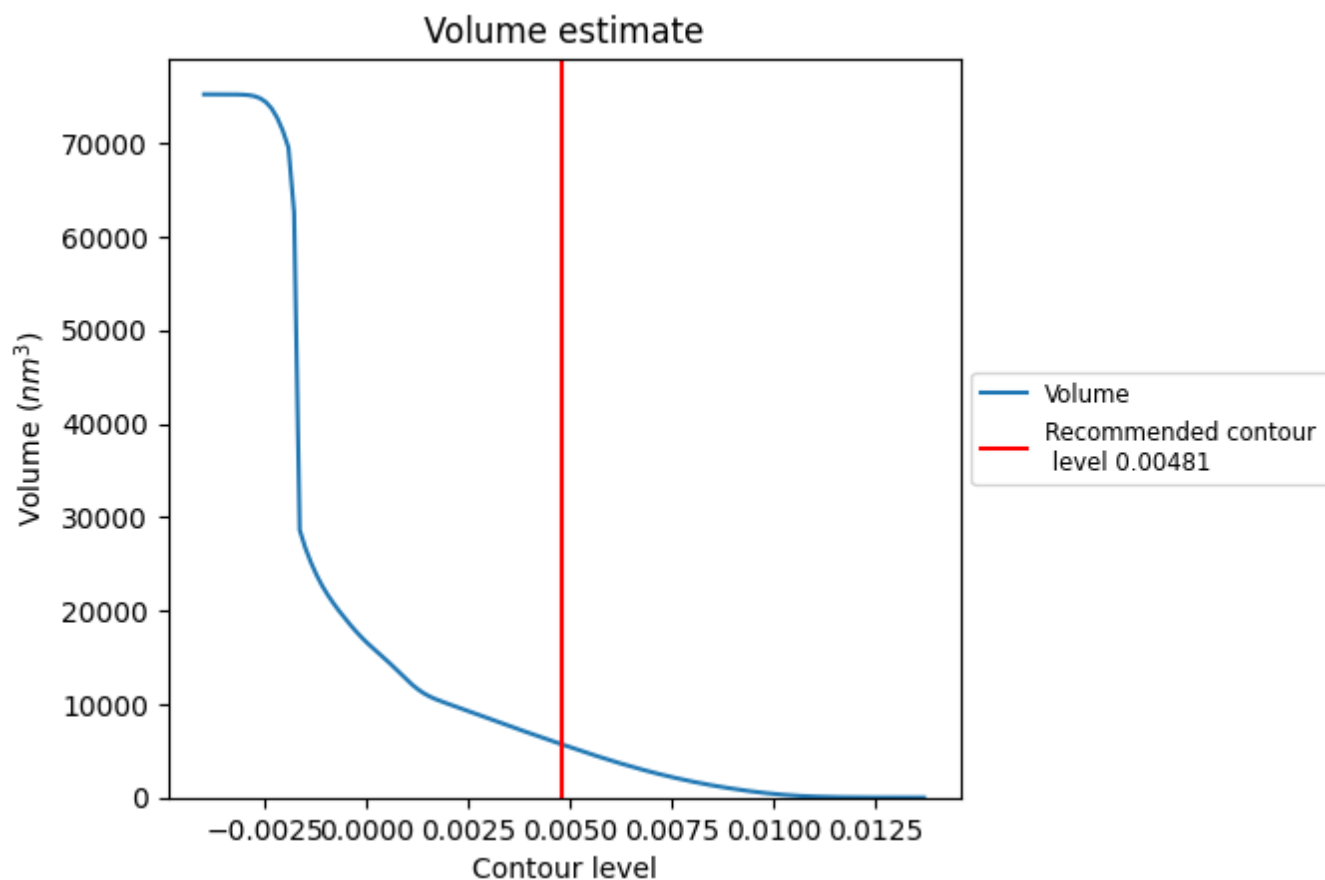
This section contains the results of statistical analysis of the map.

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.

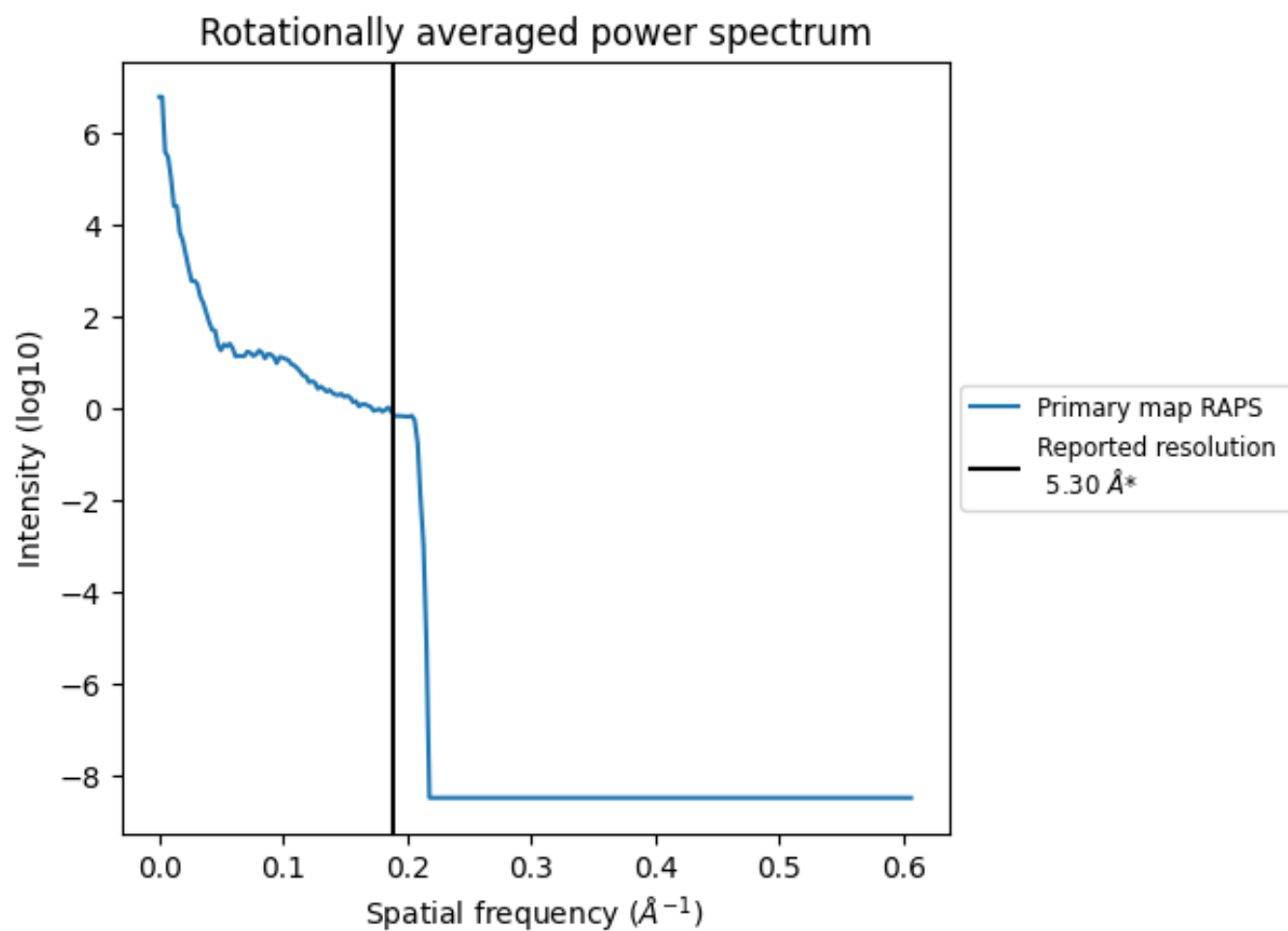
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 5673 nm³; this corresponds to an approximate mass of 5124 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 0.189 Å⁻¹

8 Fourier-Shell correlation ⓘ

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

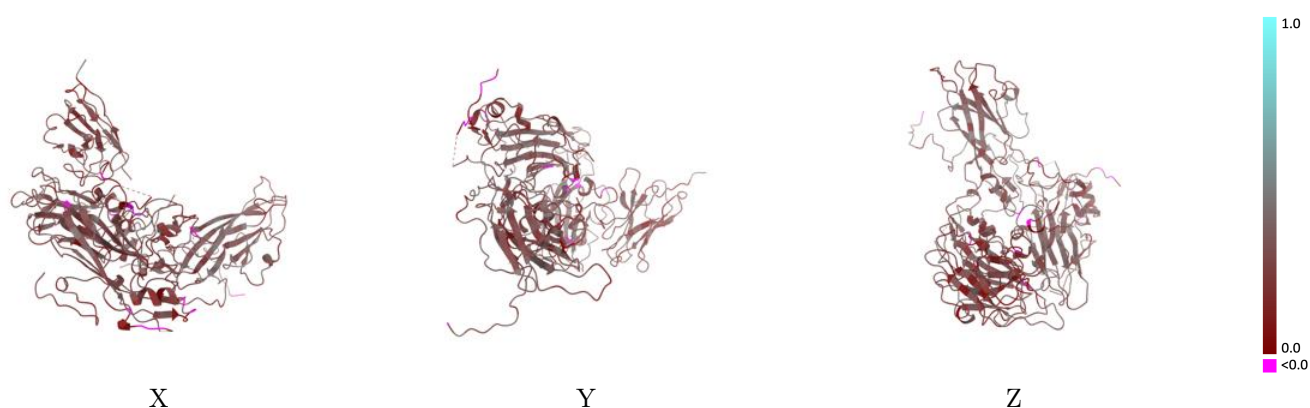
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-8286 and PDB model 5KU2. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlay [i](#)

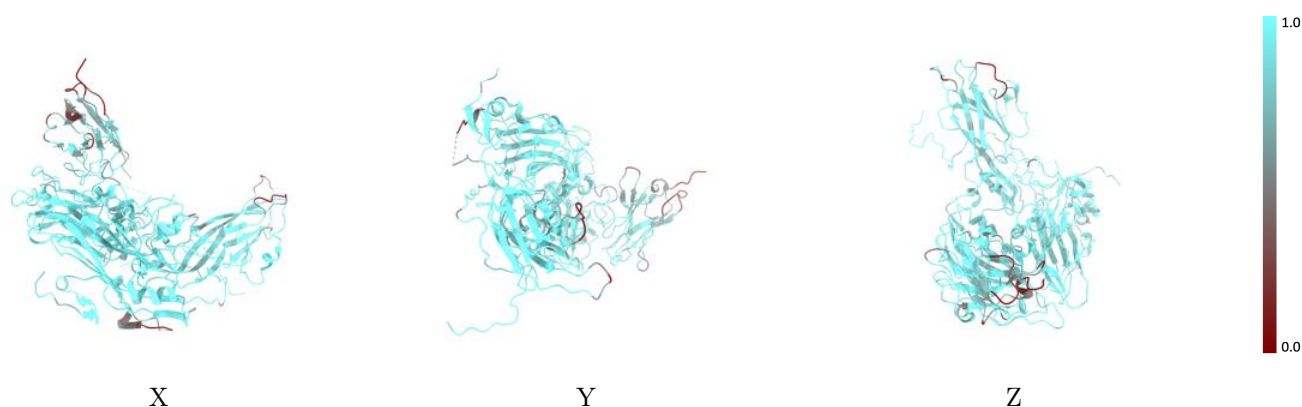
This section was not generated.

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



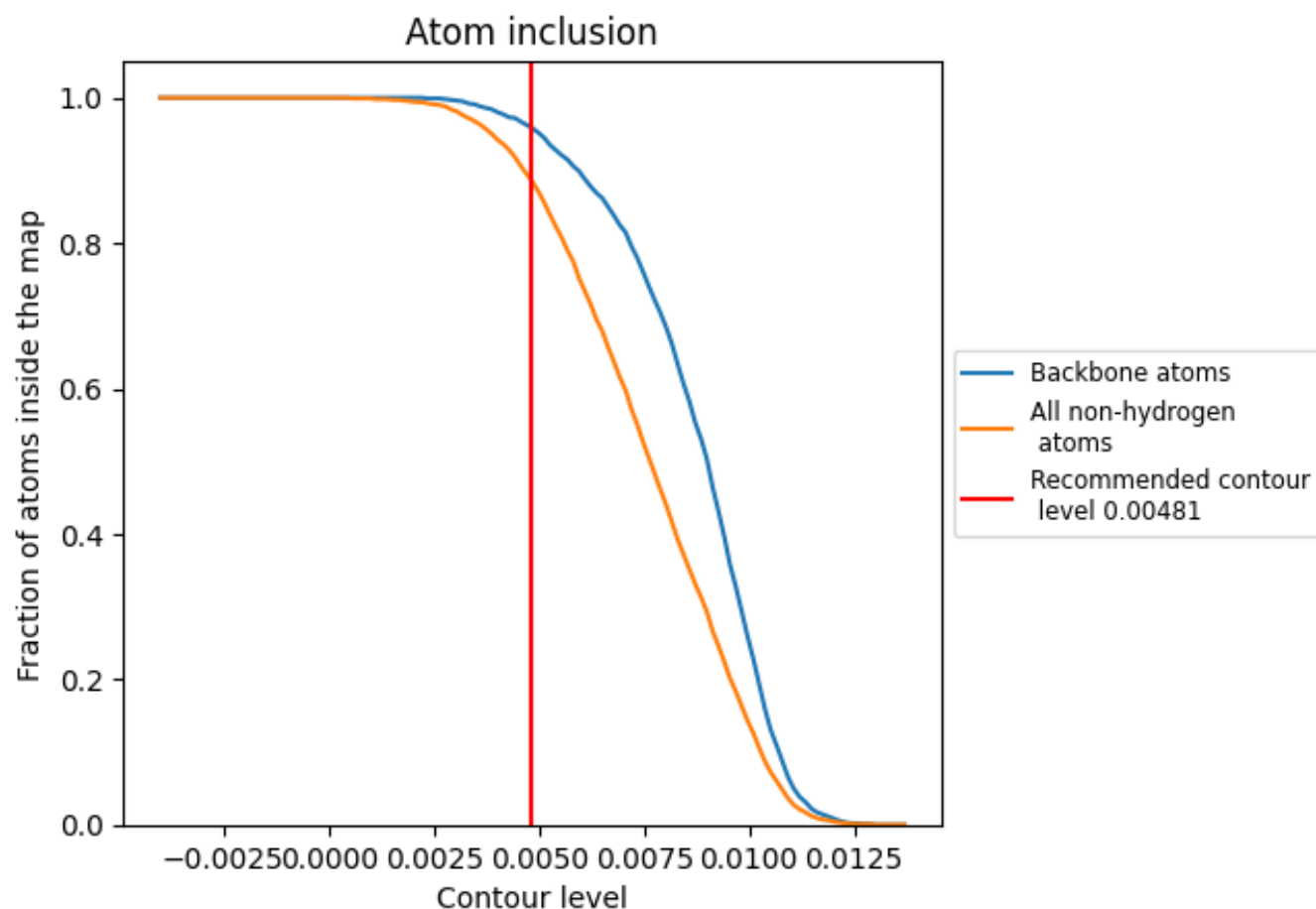
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.00481).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 89% 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 (0.00481) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8870	0.2720
1	0.8910	0.2950
2	0.9280	0.2620
3	0.9250	0.2800
7	0.7250	0.2340

