



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 02:44 PM UTC

PDB ID : 5KWL / pdb_00005kwl
EMDB ID : EMD-8277
Title : expanded poliovirus in complex with VHH 10E
Authors : Strauss, M.; Schotte, L.; Filman, D.J.; Hogle, J.M.
Deposited on : 2016-07-18
Resolution : 4.80 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:

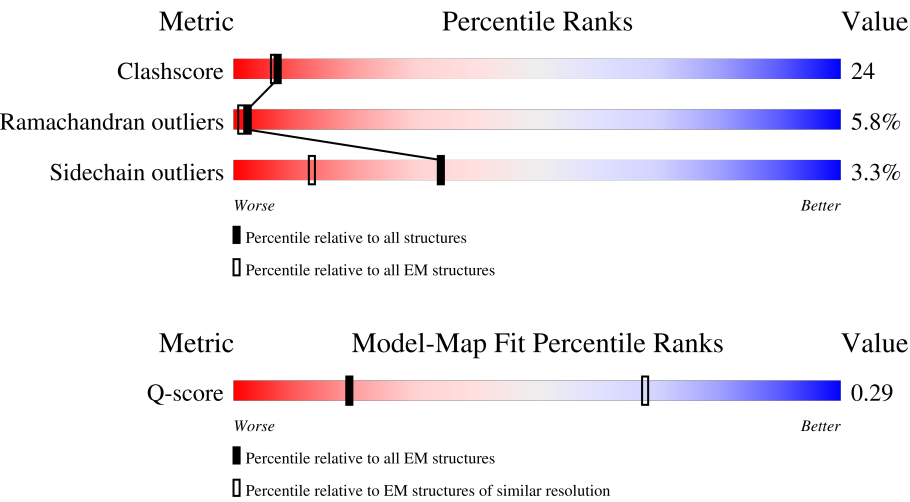
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 4.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	1575 (4.30 - 5.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1	209	5% 51%31%8%9%
2	2	268	54%32%10%
3	3	230	49%41%9%
4	7	124	39%46%44%10%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6124 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	190	Total	C	N	O	S	0	0
			1537	993	260	279	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	241	Total	C	N	O	S	0	0
			1863	1185	314	350	14		

- Molecule 3 is a protein called VP3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	230	Total	C	N	O	S	0	0
			1794	1144	292	341	17		

There is a discrepancy between the modelled and reference sequences:

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

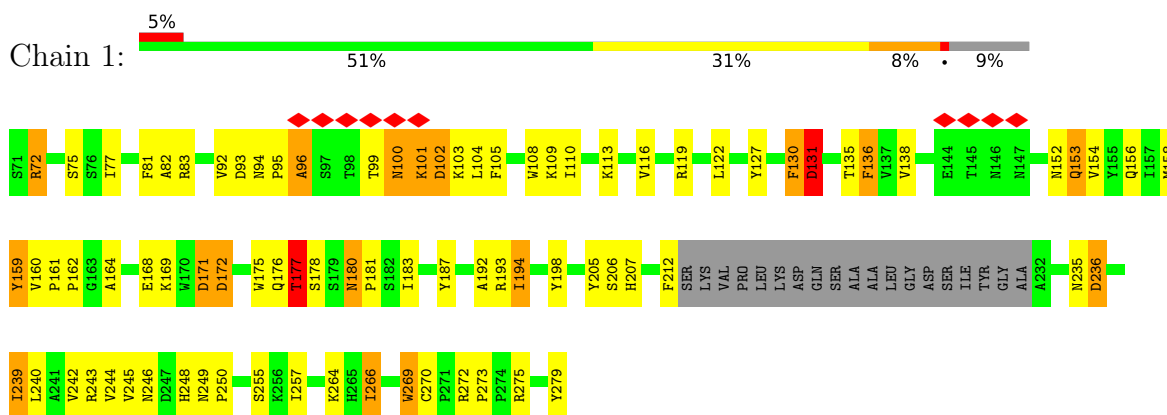
- Molecule 4 is a protein called VHH 10E.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	7	124	Total	C	N	O	S	0	0
			930	577	162	187	4		

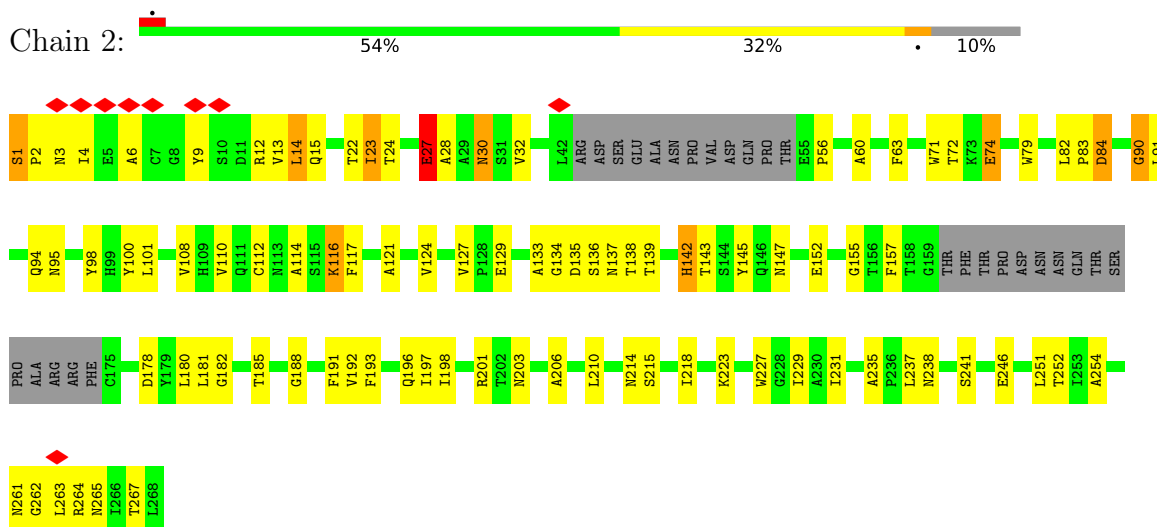
3 Residue-property plots

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.

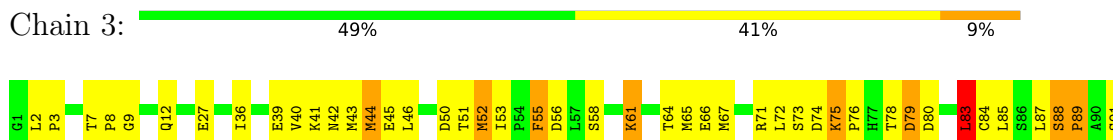
• Molecule 1: VP1

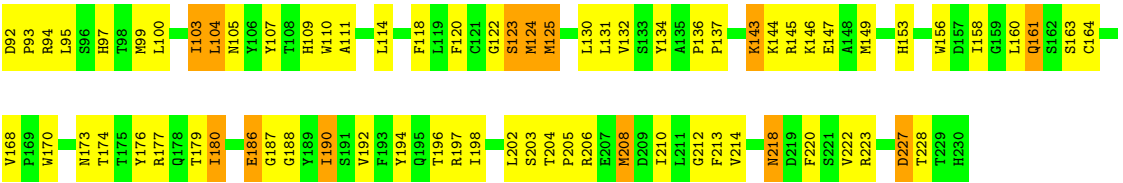


• Molecule 2: VP2

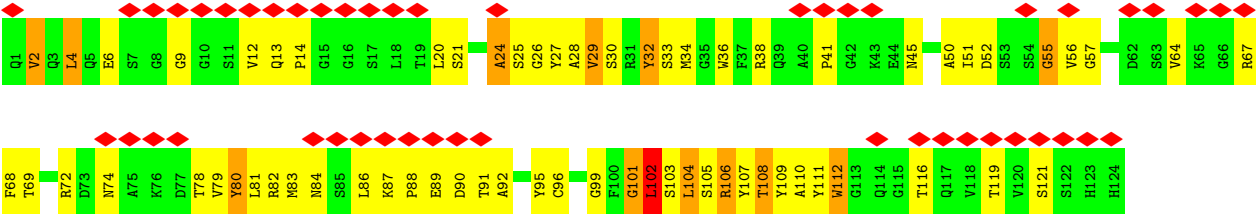


• Molecule 3: VP3





• Molecule 4: VHH 10E



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	13938	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 ($e^-/\text{\AA}^2$)	35	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	0.017	Depositor
Minimum map value	-0.004	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.00707	Depositor
Map size (Å)	504.832, 504.832, 504.832	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.986, 0.986, 0.986	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	1	0.70	0/1584	1.27	18/2161 (0.8%)
2	2	0.65	0/1912	1.19	14/2611 (0.5%)
3	3	0.68	0/1841	1.26	19/2509 (0.8%)
4	7	0.73	0/950	1.29	10/1286 (0.8%)
All	All	0.68	0/6287	1.24	61/8567 (0.7%)

There are no bond length outliers.

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	7	108	THR	N-CA-C	-9.34	101.18	112.58
1	1	102	ASP	N-CA-C	-9.21	101.60	113.12
3	3	103	ILE	N-CA-C	-8.03	102.25	111.00
4	7	101	GLY	CA-C-N	7.65	135.47	121.70
4	7	101	GLY	C-N-CA	7.65	135.47	121.70
3	3	88	SER	N-CA-C	-7.45	95.83	108.94
4	7	32	TYR	N-CA-C	-7.32	101.09	110.53
2	2	1	SER	CA-C-N	7.05	128.65	119.84
2	2	1	SER	C-N-CA	7.05	128.65	119.84
4	7	56	VAL	N-CA-C	6.60	116.97	108.12
3	3	71	ARG	N-CA-C	6.52	118.78	108.67
2	2	116	LYS	N-CA-C	-6.48	105.12	113.16
1	1	177	THR	N-CA-C	6.41	124.45	110.80
1	1	100	ASN	N-CA-C	6.41	120.56	112.87
1	1	193	ARG	N-CA-C	6.35	118.41	107.93
1	1	131	ASP	N-CA-C	-6.31	100.32	109.59
3	3	65	MET	N-CA-C	6.25	118.61	111.11
3	3	55	PHE	N-CA-C	6.13	117.96	111.28
4	7	80	TYR	N-CA-C	6.11	118.81	107.99
4	7	112	TRP	N-CA-C	6.07	113.76	108.78
3	3	227	ASP	N-CA-C	6.00	116.50	108.23
2	2	27	GLU	N-CA-C	-5.95	98.13	110.80
2	2	142	HIS	N-CA-C	5.94	123.45	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	96	ALA	N-CA-C	-5.92	105.37	113.30
2	2	223	LYS	N-CA-C	5.87	118.91	111.69
1	1	269	TRP	N-CA-C	5.83	116.92	108.34
1	1	244	VAL	N-CA-C	-5.83	100.19	108.58
2	2	152	GLU	N-CA-C	5.80	119.53	112.23
2	2	138	THR	N-CA-C	-5.76	105.42	113.37
1	1	159	TYR	N-CA-C	-5.73	99.39	108.73
1	1	266	ILE	N-CA-C	5.72	116.82	108.58
3	3	186	GLU	N-CA-C	5.62	117.55	108.34
1	1	130	PHE	N-CA-C	5.55	117.42	108.32
4	7	72	ARG	N-CA-C	5.54	117.81	109.23
4	7	24	ALA	N-CA-C	5.52	117.82	110.53
1	1	172	ASP	N-CA-C	5.49	122.49	110.80
1	1	116	VAL	N-CA-C	5.46	116.70	112.12
3	3	72	LEU	N-CA-C	-5.45	99.91	108.63
3	3	92	ASP	CA-C-N	5.45	125.80	119.47
3	3	92	ASP	C-N-CA	5.45	125.80	119.47
3	3	97	HIS	N-CA-C	5.43	119.53	113.02
1	1	136	PHE	N-CA-C	5.41	117.49	108.02
1	1	270	CYS	N-CA-C	5.41	117.27	109.48
2	2	23	ILE	CB-CA-C	-5.38	103.18	111.31
3	3	136	PRO	CA-C-N	5.37	126.56	119.84
3	3	136	PRO	C-N-CA	5.37	126.56	119.84
3	3	161	GLN	N-CA-C	5.36	117.25	108.52
1	1	194	ILE	N-CA-C	5.31	117.51	108.86
1	1	99	THR	N-CA-C	-5.31	102.27	109.54
1	1	153	GLN	N-CA-C	5.16	118.52	110.32
2	2	127	VAL	N-CA-C	5.15	113.60	107.73
3	3	143	LYS	N-CA-C	5.15	119.62	113.23
4	7	112	TRP	CA-C-O	5.14	120.92	117.94
3	3	61	LYS	N-CA-C	5.10	119.55	113.23
2	2	84	ASP	N-CA-C	5.08	121.62	110.80
3	3	203	SER	N-CA-C	5.08	117.32	110.06
2	2	117	PHE	N-CA-C	5.05	120.07	113.30
2	2	91	LEU	N-CA-C	5.03	118.19	111.75
2	2	22	THR	N-CA-C	5.02	115.61	107.23
3	3	27	GLU	N-CA-C	5.01	118.97	112.86
3	3	52	MET	N-CA-C	5.01	116.95	109.59

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	1537	0	1498	61	0
2	2	1863	0	1797	60	0
3	3	1794	0	1773	127	0
4	7	930	0	881	81	0
All	All	6124	0	5949	284	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 24.

All (284) 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)
3:3:110:TRP:CD1	3:3:179:THR:HG22	1.55	1.41
3:3:110:TRP:HD1	3:3:179:THR:CG2	1.46	1.27
2:2:238:ASN:HB3	3:3:205:PRO:HG2	1.35	1.06
4:7:24:ALA:HB2	4:7:29:VAL:HG13	1.33	1.05
2:2:143:THR:HG22	2:2:147:ASN:HB2	1.40	1.00
3:3:170:TRP:HH2	3:3:177:ARG:HB2	1.24	1.00
3:3:143:LYS:NZ	4:7:102:LEU:C	2.20	0.98
2:2:56:PRO:HD2	2:2:95:ASN:HD21	1.28	0.95
3:3:143:LYS:CE	4:7:101:GLY:H	1.78	0.95
3:3:61:LYS:HA	3:3:61:LYS:HE2	1.45	0.94
3:3:134:TYR:CE1	3:3:168:VAL:HG13	2.03	0.92
3:3:83:LEU:HA	4:7:107:TYR:HE1	1.37	0.90
4:7:34:MET:HG2	4:7:79:VAL:HG11	1.54	0.90
4:7:20:LEU:HD12	4:7:81:LEU:HD23	1.52	0.89
1:1:177:THR:OG1	1:1:180:ASN:HB2	1.72	0.89
4:7:4:LEU:HD22	4:7:24:ALA:HA	1.55	0.88
3:3:170:TRP:CH2	3:3:177:ARG:HB2	2.08	0.87
2:2:101:LEU:HD12	2:2:261:ASN:HB2	1.56	0.86
3:3:143:LYS:HZ3	4:7:102:LEU:C	1.83	0.84
3:3:9:GLY:O	3:3:12:GLN:HG2	1.78	0.84
3:3:7:THR:HB	3:3:8:PRO:HD2	1.60	0.83
1:1:109:LYS:HA	1:1:239:ILE:HG22	1.62	0.82
4:7:80:TYR:HB3	4:7:82:ARG:CZ	2.08	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:80:ASP:OD1	4:7:103:SER:HB2	1.80	0.80
1:1:212:PHE:HZ	2:2:143:THR:O	1.64	0.80
3:3:61:LYS:HD3	3:3:66:GLU:HB2	1.64	0.79
2:2:23:ILE:HG13	2:2:63:PHE:CE1	2.18	0.78
2:2:56:PRO:HD2	2:2:95:ASN:ND2	1.97	0.77
3:3:110:TRP:CD1	3:3:179:THR:CG2	2.37	0.77
3:3:87:LEU:HB3	3:3:188:GLY:H	1.50	0.77
3:3:110:TRP:HD1	3:3:179:THR:HG22	0.64	0.75
2:2:143:THR:HG22	2:2:147:ASN:CB	2.16	0.75
1:1:205:TYR:HA	1:1:235:ASN:ND2	2.01	0.75
3:3:84:CYS:HB2	4:7:108:THR:HG23	1.70	0.74
3:3:44:MET:SD	3:3:220:PHE:HD2	2.10	0.74
3:3:143:LYS:HZ2	4:7:102:LEU:HB3	1.52	0.73
1:1:95:PRO:O	1:1:248:HIS:ND1	2.19	0.72
3:3:87:LEU:O	3:3:187:GLY:HA2	1.89	0.72
2:2:188:GLY:HA2	3:3:52:MET:CE	2.19	0.71
3:3:87:LEU:HD11	3:3:114:LEU:HD12	1.73	0.71
3:3:85:LEU:HA	3:3:94:ARG:HH22	1.55	0.70
1:1:177:THR:OG1	1:1:180:ASN:CB	2.39	0.70
3:3:53:ILE:HD11	3:3:214:VAL:HB	1.74	0.69
2:2:98:TYR:O	2:2:264:ARG:HG2	1.92	0.69
1:1:269:TRP:CD1	3:3:36:ILE:HB	2.27	0.69
4:7:105:SER:O	4:7:107:TYR:N	2.22	0.69
2:2:188:GLY:HA2	3:3:52:MET:HE2	1.74	0.69
3:3:143:LYS:CE	4:7:101:GLY:N	2.55	0.68
1:1:100:ASN:HA	1:1:103:LYS:HB2	1.75	0.68
3:3:137:PRO:HA	3:3:186:GLU:HB3	1.74	0.68
4:7:105:SER:C	4:7:107:TYR:H	2.00	0.68
3:3:143:LYS:HZ2	4:7:102:LEU:CB	2.07	0.67
2:2:191:PHE:HE1	3:3:52:MET:HE2	1.59	0.67
1:1:212:PHE:CZ	2:2:143:THR:O	2.47	0.67
3:3:143:LYS:NZ	4:7:101:GLY:H	1.93	0.67
3:3:146:LYS:O	3:3:149:MET:HE3	1.95	0.66
3:3:44:MET:HA	3:3:44:MET:HE2	1.78	0.66
3:3:143:LYS:NZ	4:7:102:LEU:CA	2.58	0.66
3:3:137:PRO:CA	3:3:186:GLU:HB3	2.25	0.65
3:3:87:LEU:HB3	3:3:188:GLY:N	2.10	0.65
4:7:20:LEU:HB2	4:7:81:LEU:HB3	1.78	0.65
1:1:152:ASN:OD1	1:1:187:TYR:HB3	1.96	0.65
4:7:57:GLY:O	4:7:102:LEU:HD21	1.98	0.64
3:3:143:LYS:HE2	4:7:101:GLY:H	1.60	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:120:PHE:HA	3:3:210:ILE:HG22	1.79	0.63
2:2:72:THR:HG22	2:2:246:GLU:HB3	1.81	0.62
3:3:109:HIS:HB3	3:3:176:TYR:HB3	1.80	0.62
3:3:180:ILE:HD13	3:3:180:ILE:H	1.65	0.62
2:2:90:GLY:O	2:2:94:GLN:HG2	1.99	0.62
4:7:33:SER:OG	4:7:101:GLY:C	2.42	0.62
4:7:67:ARG:O	4:7:84:ASN:HB2	1.99	0.62
3:3:39:GLU:HG2	3:3:40:VAL:N	2.15	0.62
4:7:36:TRP:CG	4:7:81:LEU:HD22	2.34	0.62
4:7:13:GLN:HE21	4:7:14:PRO:HD2	1.65	0.62
2:2:56:PRO:O	2:2:60:ALA:HB3	2.01	0.61
3:3:83:LEU:HA	4:7:107:TYR:CE1	2.28	0.61
1:1:96:ALA:HA	1:1:248:HIS:HB3	1.83	0.61
3:3:83:LEU:CA	4:7:107:TYR:HE1	2.10	0.61
3:3:143:LYS:HZ1	4:7:102:LEU:C	2.06	0.61
4:7:21:SER:HA	4:7:79:VAL:O	2.00	0.60
4:7:2:VAL:HG22	4:7:111:TYR:CG	2.37	0.60
3:3:75:LYS:HB2	3:3:76:PRO:HD2	1.85	0.59
1:1:131:ASP:HB2	1:1:264:LYS:HG3	1.84	0.59
3:3:76:PRO:C	3:3:78:THR:H	2.09	0.59
2:2:30:ASN:HD22	2:2:203:ASN:ND2	2.01	0.59
2:2:201:ARG:NH1	3:3:124:MET:HA	2.17	0.58
3:3:61:LYS:HE2	3:3:61:LYS:CA	2.28	0.58
2:2:32:VAL:HG23	2:2:196:GLN:HE21	1.69	0.58
3:3:55:PHE:CE2	3:3:212:GLY:HA3	2.39	0.58
4:7:2:VAL:HB	4:7:25:SER:O	2.02	0.58
3:3:56:ASP:OD2	4:7:107:TYR:HD2	1.85	0.58
1:1:94:ASN:HA	1:1:104:LEU:HD23	1.85	0.58
4:7:87:LYS:HB3	4:7:88:PRO:HD2	1.85	0.58
4:7:9:GLY:HA3	4:7:116:THR:OG1	2.03	0.58
3:3:134:TYR:HE1	3:3:168:VAL:HG13	1.64	0.58
4:7:91:THR:HG23	4:7:119:THR:HA	1.85	0.58
3:3:132:VAL:HG22	3:3:192:VAL:HG22	1.86	0.58
2:2:110:VAL:HG22	2:2:251:LEU:HG	1.86	0.57
4:7:33:SER:OG	4:7:101:GLY:O	2.22	0.57
3:3:104:LEU:HD11	3:3:220:PHE:CE2	2.39	0.56
2:2:63:PHE:HB3	2:2:254:ALA:HB2	1.87	0.56
1:1:153:GLN:HA	1:1:246:ASN:HD21	1.71	0.56
4:7:69:THR:HB	4:7:82:ARG:HB2	1.86	0.56
1:1:138:VAL:HG22	1:1:257:ILE:HG13	1.88	0.56
3:3:100:LEU:O	3:3:104:LEU:HD13	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:208:MET:SD	3:3:208:MET:N	2.80	0.55
1:1:136:PHE:HB2	1:1:192:ALA:HB3	1.87	0.55
4:7:89:GLU:C	4:7:91:THR:H	2.14	0.55
4:7:33:SER:HG	4:7:101:GLY:C	2.13	0.55
1:1:177:THR:O	1:1:178:SER:C	2.48	0.55
3:3:44:MET:SD	3:3:220:PHE:CD2	2.97	0.55
3:3:120:PHE:CE2	3:3:122:GLY:HA3	2.42	0.55
1:1:110:ILE:HD11	1:1:240:LEU:HG	1.89	0.55
4:7:2:VAL:HG22	4:7:111:TYR:CD2	2.42	0.54
3:3:88:SER:HB3	3:3:91:SER:OG	2.07	0.54
3:3:149:MET:SD	3:3:149:MET:C	2.90	0.54
4:7:45:ASN:H	4:7:106:ARG:NH1	2.05	0.54
3:3:143:LYS:HZ1	4:7:101:GLY:H	1.54	0.54
4:7:13:GLN:NE2	4:7:14:PRO:HD2	2.22	0.54
4:7:41:PRO:HD3	4:7:92:ALA:HA	1.89	0.54
3:3:46:LEU:HB2	3:3:103:ILE:HD11	1.88	0.54
1:1:122:LEU:HB3	1:1:207:HIS:NE2	2.23	0.54
3:3:156:TRP:CD1	3:3:164:CYS:HB2	2.42	0.54
1:1:127:TYR:HB2	1:1:269:TRP:HB2	1.90	0.54
4:7:27:TYR:CE2	4:7:29:VAL:HA	2.42	0.54
3:3:87:LEU:HD22	3:3:168:VAL:HG11	1.90	0.54
4:7:14:PRO:HD3	4:7:121:SER:HA	1.89	0.53
3:3:85:LEU:HA	3:3:94:ARG:NH2	2.21	0.53
1:1:235:ASN:O	1:1:236:ASP:HB3	2.08	0.53
3:3:147:GLU:OE1	3:3:147:GLU:HA	2.08	0.53
1:1:101:LYS:H	1:1:101:LYS:HD2	1.73	0.53
4:7:34:MET:HB3	4:7:79:VAL:HG21	1.90	0.53
4:7:99:GLY:HA3	4:7:108:THR:O	2.09	0.53
3:3:64:THR:O	3:3:67:MET:HG2	2.09	0.53
2:2:116:LYS:HD3	3:3:125:MET:HE2	1.89	0.52
4:7:50:ALA:HB1	4:7:102:LEU:HA	1.91	0.52
2:2:72:THR:HG22	2:2:246:GLU:CB	2.39	0.52
4:7:51:ILE:HA	4:7:102:LEU:HD23	1.90	0.52
4:7:29:VAL:O	4:7:32:TYR:O	2.28	0.52
3:3:7:THR:HB	3:3:8:PRO:CD	2.36	0.52
3:3:130:LEU:HG	3:3:158:ILE:HD11	1.90	0.52
1:1:81:PHE:C	1:1:83:ARG:H	2.16	0.51
1:1:275:ARG:HG2	1:1:279:TYR:CE2	2.44	0.51
4:7:21:SER:HB3	4:7:78:THR:CG2	2.39	0.51
2:2:63:PHE:HB2	2:2:252:THR:HG22	1.92	0.51
3:3:107:TYR:CG	3:3:222:VAL:HB	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:112:CYS:HB3	2:2:198:ILE:HG21	1.92	0.51
2:2:188:GLY:HA2	3:3:52:MET:HE3	1.92	0.51
2:2:196:GLN:OE1	2:2:206:ALA:HA	2.11	0.51
4:7:6:GLU:HG2	4:7:96:CYS:SG	2.51	0.51
4:7:107:TYR:O	4:7:109:TYR:HD2	1.94	0.51
3:3:61:LYS:HD3	3:3:66:GLU:CB	2.38	0.50
4:7:83:MET:HB3	4:7:86:LEU:HD21	1.93	0.50
1:1:102:ASP:O	1:1:105:PHE:O	2.30	0.50
1:1:93:ASP:O	1:1:104:LEU:HD23	2.12	0.50
3:3:75:LYS:HB2	3:3:76:PRO:CD	2.41	0.49
3:3:52:MET:HA	3:3:212:GLY:O	2.11	0.49
4:7:52:ASP:HB3	4:7:55:GLY:O	2.11	0.49
1:1:273:PRO:HB2	2:2:185:THR:HB	1.93	0.49
2:2:79:TRP:CZ2	2:2:155:GLY:HA3	2.47	0.49
2:2:191:PHE:CE1	3:3:52:MET:HE2	2.42	0.49
3:3:87:LEU:CD2	3:3:168:VAL:HG11	2.42	0.49
1:1:273:PRO:HD3	2:2:193:PHE:CZ	2.47	0.49
4:7:101:GLY:HA3	4:7:102:LEU:HB3	1.95	0.49
3:3:173:ASN:CG	3:3:174:THR:H	2.21	0.49
3:3:190:ILE:N	3:3:190:ILE:HD12	2.28	0.49
2:2:121:ALA:N	2:2:235:ALA:HB3	2.28	0.48
4:7:105:SER:C	4:7:107:TYR:N	2.70	0.48
2:2:180:LEU:C	2:2:182:GLY:N	2.71	0.48
4:7:38:ARG:O	4:7:38:ARG:HG3	2.14	0.48
3:3:120:PHE:HE2	3:3:122:GLY:HA3	1.78	0.48
3:3:111:ALA:O	3:3:220:PHE:HA	2.14	0.48
3:3:79:ASP:O	3:3:80:ASP:C	2.57	0.48
3:3:198:ILE:HG21	3:3:208:MET:HE1	1.96	0.48
4:7:2:VAL:HG11	4:7:27:TYR:CD1	2.49	0.48
2:2:82:LEU:HA	2:2:83:PRO:C	2.38	0.47
2:2:2:PRO:HG2	2:2:197:ILE:HG13	1.97	0.47
2:2:74:GLU:H	2:2:74:GLU:CD	2.22	0.47
1:1:72:ARG:O	1:1:75:SER:OG	2.29	0.47
3:3:55:PHE:HE2	3:3:212:GLY:HA3	1.80	0.47
3:3:158:ILE:HD12	3:3:158:ILE:N	2.29	0.47
1:1:176:GLN:OE1	1:1:176:GLN:HA	2.15	0.47
1:1:177:THR:OG1	1:1:177:THR:O	2.26	0.47
1:1:183:ILE:HD11	1:1:194:ILE:CG2	2.45	0.47
3:3:56:ASP:OD2	4:7:107:TYR:CD2	2.66	0.47
1:1:249:ASN:HA	1:1:250:PRO:HD2	1.83	0.47
4:7:64:VAL:HB	4:7:68:PHE:CD1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:159:TYR:O	1:1:181:PRO:HD2	2.14	0.46
1:1:272:ARG:NH2	2:2:129:GLU:O	2.48	0.46
2:2:74:GLU:CD	2:2:74:GLU:N	2.73	0.46
3:3:158:ILE:HD12	3:3:158:ILE:H	1.81	0.46
2:2:14:LEU:HG	2:2:27:GLU:O	2.16	0.46
3:3:79:ASP:HB3	3:3:144:LYS:HA	1.97	0.46
2:2:56:PRO:HB2	2:2:60:ALA:HB2	1.98	0.46
3:3:41:LYS:HB2	3:3:45:GLU:OE2	2.15	0.46
2:2:56:PRO:CD	2:2:95:ASN:HD21	2.12	0.46
3:3:104:LEU:HD11	3:3:220:PHE:HE2	1.80	0.46
2:2:71:TRP:NE1	2:2:237:LEU:HD11	2.30	0.46
3:3:87:LEU:C	3:3:89:PRO:HD3	2.41	0.46
3:3:137:PRO:HB3	3:3:186:GLU:HB3	1.97	0.46
1:1:153:GLN:HA	1:1:246:ASN:ND2	2.31	0.46
2:2:1:SER:HA	3:3:50:ASP:HB2	1.98	0.46
4:7:52:ASP:HB2	4:7:102:LEU:HG	1.96	0.46
1:1:101:LYS:H	1:1:101:LYS:CD	2.28	0.46
2:2:100:TYR:HB2	2:2:264:ARG:HH21	1.81	0.46
1:1:154:VAL:HG23	1:1:245:VAL:HB	1.97	0.45
2:2:9:TYR:CE1	3:3:124:MET:HE3	2.51	0.45
3:3:42:ASN:OD1	3:3:43:MET:N	2.49	0.45
3:3:186:GLU:HA	3:3:187:GLY:HA3	1.36	0.45
4:7:30:SER:HB3	4:7:74:ASN:HB3	1.98	0.45
3:3:55:PHE:CZ	3:3:212:GLY:HA3	2.51	0.45
1:1:93:ASP:O	1:1:104:LEU:HA	2.17	0.45
1:1:113:LYS:HA	1:1:119:ARG:HD2	1.99	0.45
3:3:137:PRO:CB	3:3:186:GLU:HB3	2.47	0.45
2:2:121:ALA:H	2:2:235:ALA:HB3	1.82	0.45
4:7:36:TRP:HA	4:7:95:TYR:O	2.17	0.45
1:1:158:MET:HE1	1:1:175:TRP:HA	1.98	0.44
3:3:227:ASP:OD1	3:3:227:ASP:N	2.50	0.44
1:1:160:VAL:HG12	1:1:164:ALA:HB3	1.98	0.44
3:3:218:ASN:OD1	3:3:218:ASN:N	2.51	0.44
3:3:85:LEU:HD21	3:3:95:LEU:HD11	1.99	0.44
1:1:81:PHE:O	1:1:83:ARG:N	2.48	0.44
3:3:61:LYS:O	3:3:67:MET:HB3	2.18	0.44
4:7:82:ARG:HD2	4:7:82:ARG:N	2.32	0.44
4:7:2:VAL:HG11	4:7:27:TYR:HD1	1.83	0.44
3:3:2:LEU:HD12	3:3:3:PRO:HD2	2.00	0.44
2:2:261:ASN:HA	2:2:262:GLY:HA2	1.81	0.43
3:3:177:ARG:NH1	3:3:186:GLU:O	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:214:ASN:CG	2:2:215:SER:N	2.76	0.43
3:3:143:LYS:HE3	4:7:101:GLY:N	2.31	0.43
1:1:171:ASP:OD1	1:1:171:ASP:N	2.51	0.43
1:1:102:ASP:O	1:1:105:PHE:N	2.51	0.43
2:2:108:VAL:HG21	2:2:229:ILE:HD11	2.00	0.43
3:3:123:SER:O	3:3:124:MET:C	2.61	0.43
2:2:192:VAL:HG11	3:3:99:MET:HE3	2.01	0.43
3:3:51:THR:O	3:3:213:PHE:HA	2.19	0.43
4:7:2:VAL:HG12	4:7:26:GLY:HA3	2.01	0.42
4:7:28:ALA:O	4:7:29:VAL:C	2.62	0.42
4:7:38:ARG:HB2	4:7:92:ALA:HB3	2.01	0.42
2:2:15:GLN:HB3	2:2:24:THR:HG22	2.01	0.42
3:3:76:PRO:C	3:3:78:THR:N	2.76	0.42
3:3:143:LYS:NZ	4:7:102:LEU:O	2.33	0.42
1:1:168:GLU:HB2	1:1:169:LYS:H	1.37	0.42
1:1:119:ARG:NH2	2:2:137:ASN:OD1	2.51	0.42
3:3:103:ILE:C	3:3:105:ASN:H	2.28	0.42
4:7:67:ARG:NH1	4:7:68:PHE:HE1	2.17	0.42
1:1:130:PHE:HB3	1:1:266:ILE:HG13	2.01	0.42
3:3:194:TYR:HB3	3:3:196:THR:O	2.19	0.42
4:7:36:TRP:NE1	4:7:81:LEU:HB2	2.35	0.42
3:3:124:MET:C	3:3:124:MET:SD	3.02	0.42
1:1:183:ILE:HD11	1:1:194:ILE:HG21	2.00	0.42
3:3:177:ARG:HA	3:3:177:ARG:HD2	1.50	0.42
4:7:107:TYR:HD1	4:7:108:THR:N	2.17	0.42
1:1:102:ASP:O	1:1:103:LYS:C	2.62	0.42
2:2:1:SER:CA	3:3:50:ASP:HB2	2.50	0.42
2:2:210:LEU:HD13	2:2:227:TRP:CE3	2.54	0.42
3:3:131:LEU:HD21	3:3:153:HIS:HB2	2.02	0.42
2:2:101:LEU:HD22	2:2:218:ILE:HG12	2.02	0.41
3:3:204:THR:O	3:3:206:ARG:HG3	2.20	0.41
1:1:77:ILE:HD13	3:3:41:LYS:C	2.45	0.41
2:2:157:PHE:CD2	2:2:178:ASP:HA	2.56	0.41
1:1:104:LEU:O	1:1:243:ARG:HA	2.20	0.41
1:1:206:SER:H	1:1:235:ASN:HD21	1.68	0.41
3:3:143:LYS:HZ2	4:7:102:LEU:CA	2.29	0.41
2:2:180:LEU:C	2:2:182:GLY:H	2.28	0.41
2:2:79:TRP:CE2	2:2:155:GLY:HA3	2.55	0.41
1:1:108:TRP:CZ3	1:1:110:ILE:HA	2.56	0.41
4:7:107:TYR:CD1	4:7:108:THR:N	2.89	0.41
1:1:160:VAL:HA	1:1:161:PRO:HD2	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:93:PRO:HG3	4:7:110:ALA:HB1	2.03	0.41
1:1:135:THR:HA	1:1:192:ALA:O	2.20	0.41
1:1:159:TYR:HA	1:1:240:LEU:HD23	2.04	0.40
3:3:58:SER:HA	4:7:112:TRP:CE2	2.56	0.40
3:3:109:HIS:HD2	3:3:223:ARG:HG3	1.85	0.40
3:3:118:PHE:O	3:3:163:SER:HA	2.21	0.40
3:3:79:ASP:CG	3:3:145:ARG:H	2.29	0.40
1:1:156:GLN:O	1:1:242:VAL:HA	2.20	0.40
1:1:160:VAL:HB	1:1:239:ILE:CD1	2.52	0.40
1:1:205:TYR:HA	1:1:235:ASN:HD21	1.79	0.40
2:2:12:ARG:HH21	2:2:13:VAL:HG12	1.87	0.40
2:2:124:VAL:HG22	2:2:231:ILE:HG12	2.03	0.40
3:3:84:CYS:HB2	4:7:108:THR:CG2	2.47	0.40
3:3:93:PRO:CG	4:7:110:ALA:HB1	2.51	0.40
1:1:92:VAL:HG23	1:1:255:SER:HB2	2.02	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	1	186/209 (89%)	151 (81%)	28 (15%)	7 (4%)	2	19
2	2	235/268 (88%)	188 (80%)	28 (12%)	19 (8%)	1	9
3	3	228/230 (99%)	186 (82%)	30 (13%)	12 (5%)	1	15
4	7	122/124 (98%)	99 (81%)	16 (13%)	7 (6%)	1	14
All	All	771/831 (93%)	624 (81%)	102 (13%)	45 (6%)	2	14

All (45) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	172	ASP
1	1	177	THR
2	2	6	ALA
2	2	28	ALA
2	2	84	ASP
2	2	135	ASP
2	2	139	THR
2	2	142	HIS
2	2	241	SER
2	2	263	LEU
2	2	265	ASN
2	2	267	THR
3	3	83	LEU
3	3	202	LEU
1	1	82	ALA
1	1	236	ASP
2	2	30	ASN
2	2	90	GLY
2	2	133	ALA
3	3	73	SER
3	3	79	ASP
3	3	124	MET
3	3	228	THR
4	7	29	VAL
4	7	55	GLY
4	7	106	ARG
1	1	198	TYR
2	2	145	TYR
2	2	181	LEU
3	3	123	SER
4	7	12	VAL
4	7	90	ASP
4	7	102	LEU
4	7	104	LEU
1	1	162	PRO
3	3	125	MET
3	3	160	LEU
1	1	72	ARG
2	2	27	GLU
2	2	114	ALA
3	3	74	ASP
3	3	104	LEU
2	2	136	SER

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Mol	Chain	Res	Type
2	2	134	GLY
3	3	89	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	169/183 (92%)	164 (97%)	5 (3%)	36	57
2	2	203/228 (89%)	199 (98%)	4 (2%)	48	66
3	3	206/206 (100%)	197 (96%)	9 (4%)	25	47
4	7	97/97 (100%)	93 (96%)	4 (4%)	27	49
All	All	675/714 (94%)	653 (97%)	22 (3%)	34	55

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	101	LYS
1	1	131	ASP
1	1	171	ASP
1	1	180	ASN
1	1	239	ILE
2	2	3	ASN
2	2	4	ILE
2	2	14	LEU
2	2	74	GLU
3	3	44	MET
3	3	75	LYS
3	3	83	LEU
3	3	161	GLN
3	3	180	ILE
3	3	190	ILE
3	3	197	ARG
3	3	208	MET
3	3	218	ASN
4	7	2	VAL

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Mol	Chain	Res	Type
4	7	4	LEU
4	7	102	LEU
4	7	104	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	100	ASN
1	1	117	GLN
1	1	147	ASN
1	1	149	HIS
1	1	203	ASN
2	2	3	ASN
2	2	15	GLN
2	2	30	ASN
2	2	95	ASN
2	2	99	HIS
2	2	203	ASN
2	2	204	ASN
3	3	6	ASN
3	3	109	HIS
3	3	161	GLN
3	3	178	GLN
4	7	5	GLN
4	7	13	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

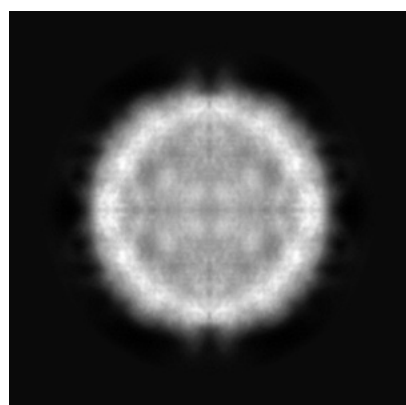
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8277. These allow visual inspection of the internal detail of the map and identification of artifacts.

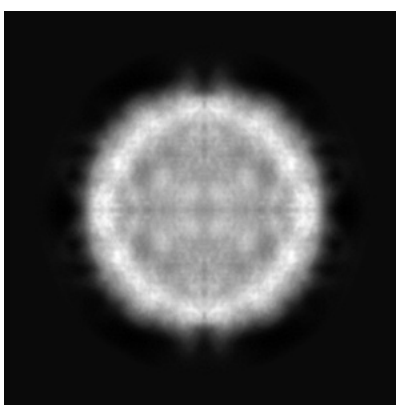
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

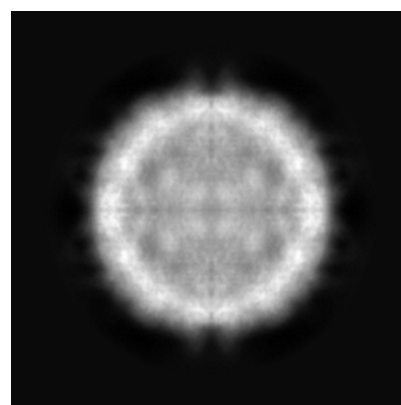
6.1.1 Primary map



X



Y

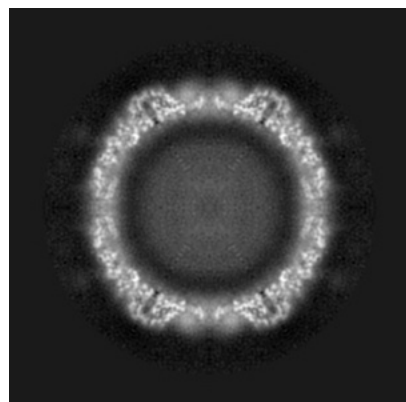


Z

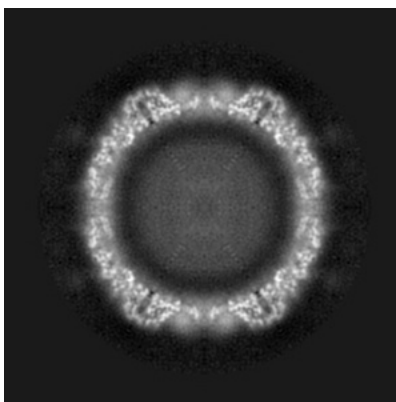
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

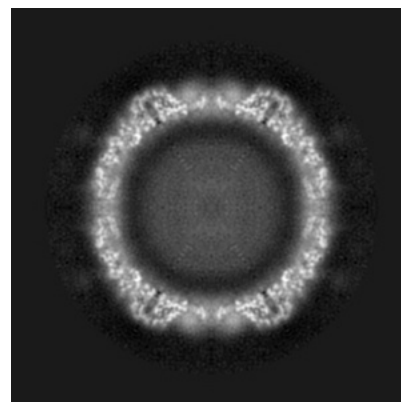
6.2.1 Primary map



X Index: 256



Y Index: 256

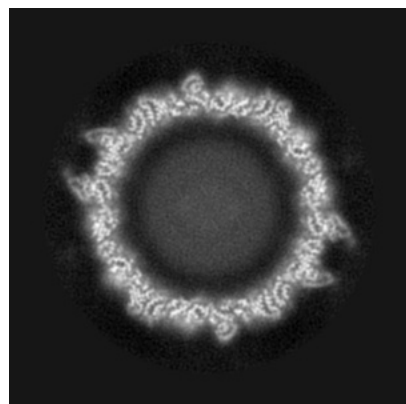


Z Index: 256

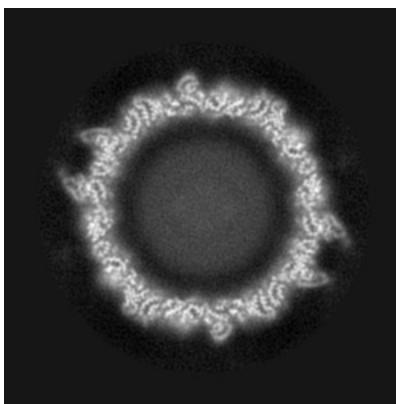
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

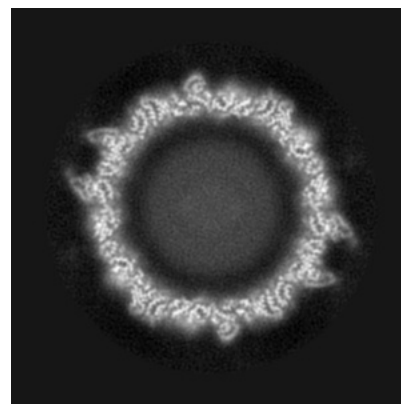
6.3.1 Primary map



X Index: 238



Y Index: 238

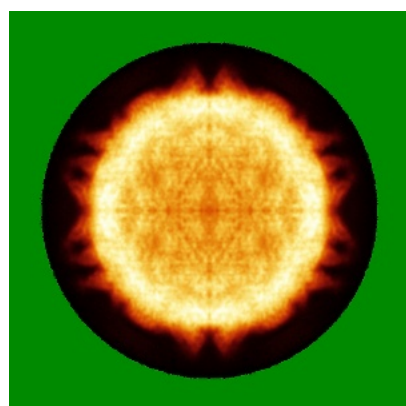


Z Index: 238

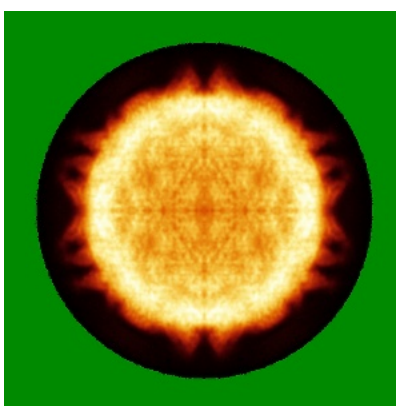
The images above show the largest variance slices of the map in three orthogonal directions.

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

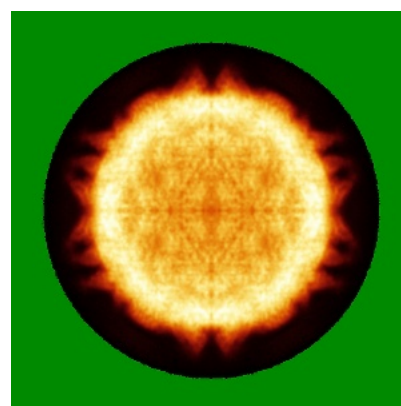
6.4.1 Primary map



X



Y

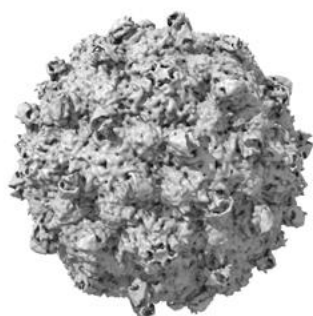


Z

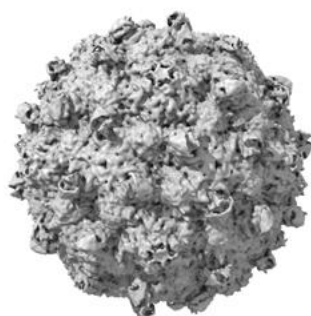
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

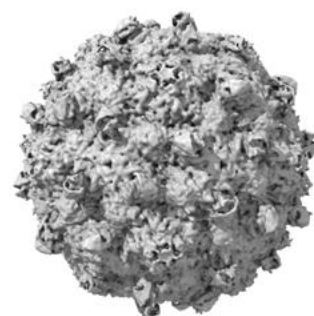
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.00707. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

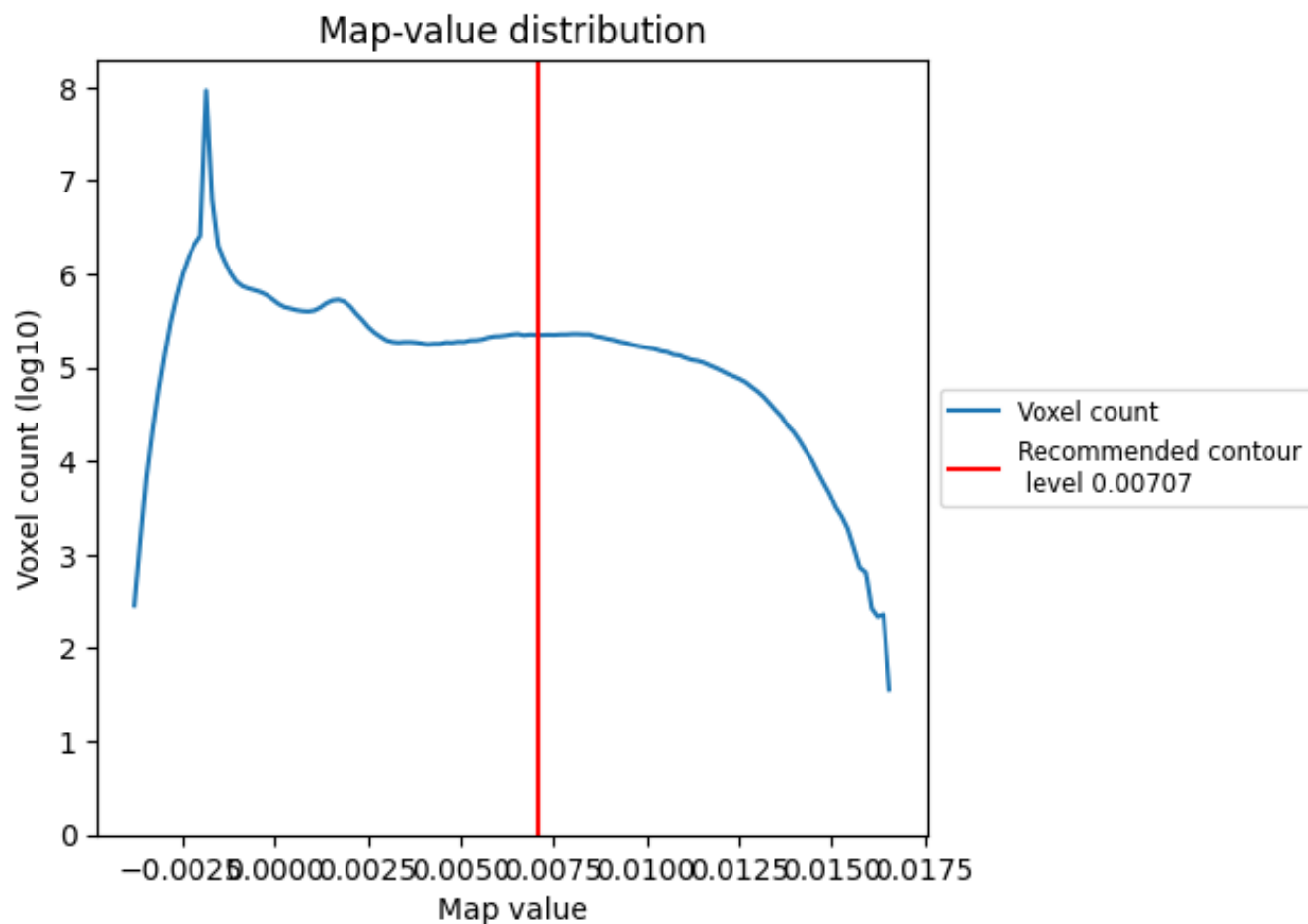
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

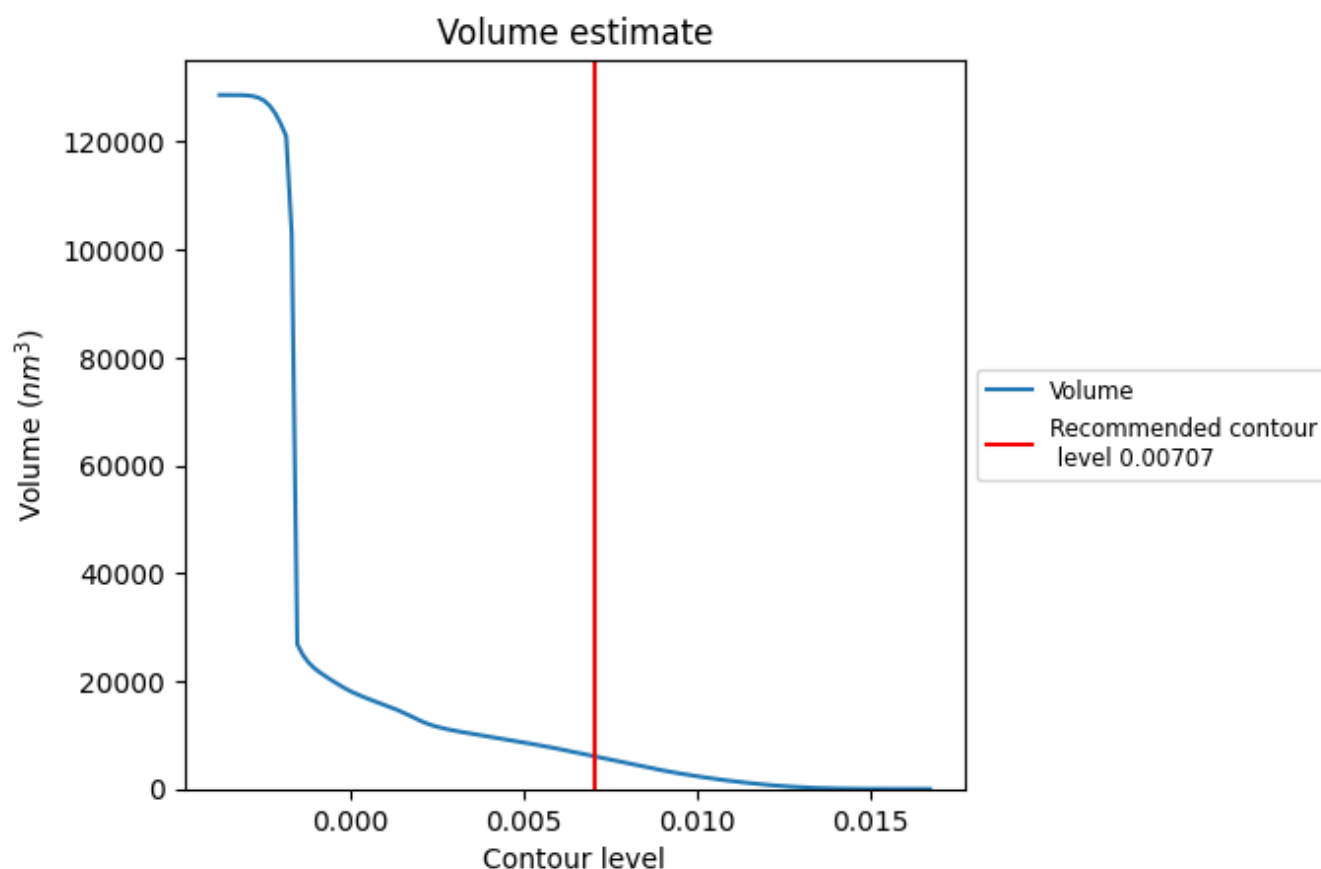
This section contains the results of statistical analysis of the map.

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.

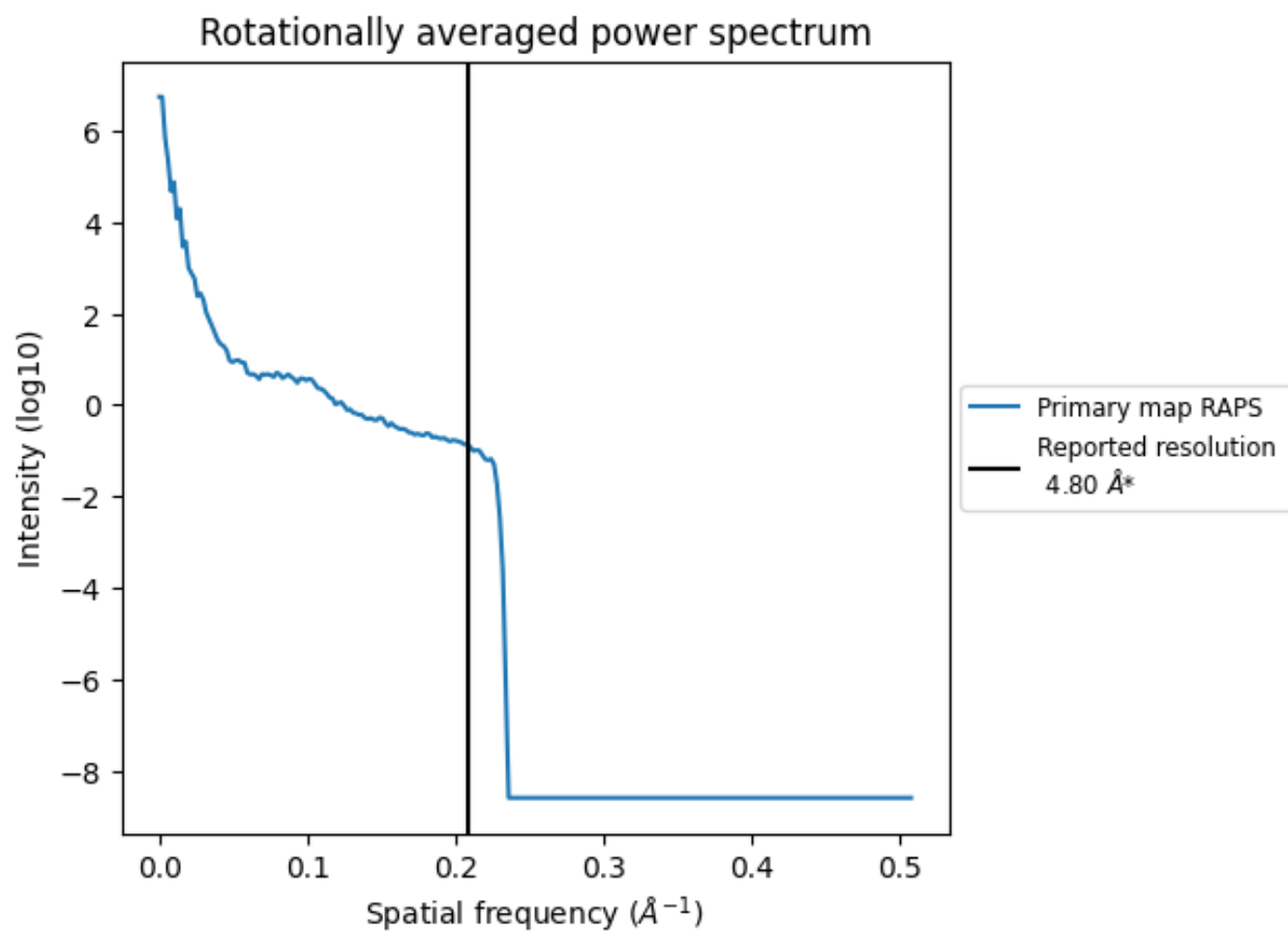
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 6013 nm³; this corresponds to an approximate mass of 5432 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 0.208 Å⁻¹

8 Fourier-Shell correlation

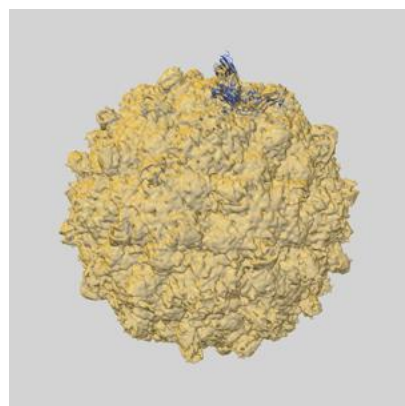
This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

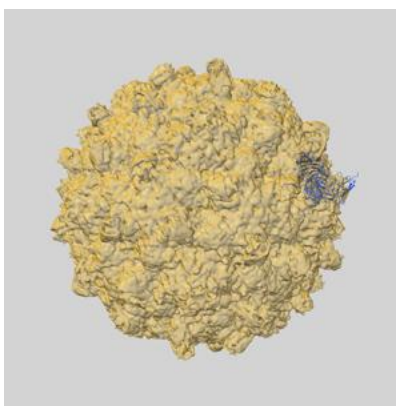
This section contains information regarding the fit between EMDB map EMD-8277 and PDB model 5KWL. Per-residue inclusion information can be found in section [3](#) on page [4](#).

9.1 Map-model overlays

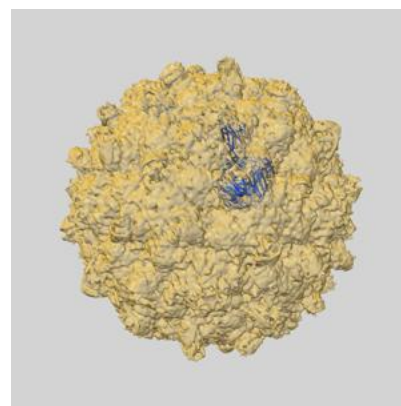
9.1.1 Map-model overlay [i](#)



X

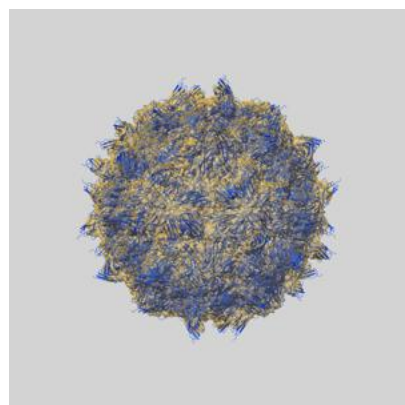


Y

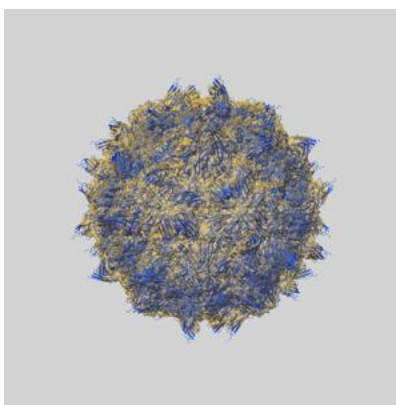


Z

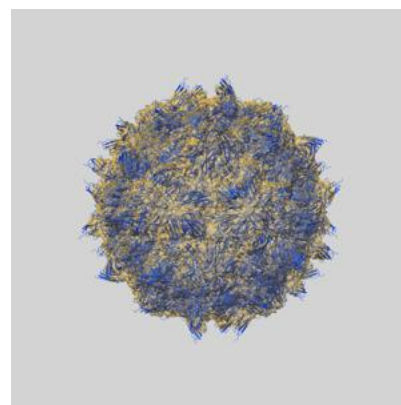
9.1.2 Map-model assembly overlay [i](#)



X



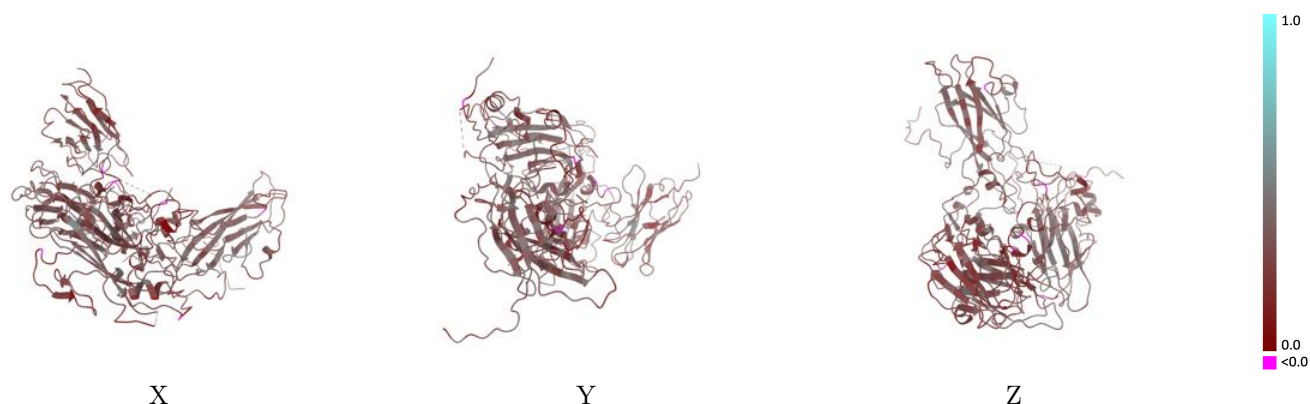
Y



Z

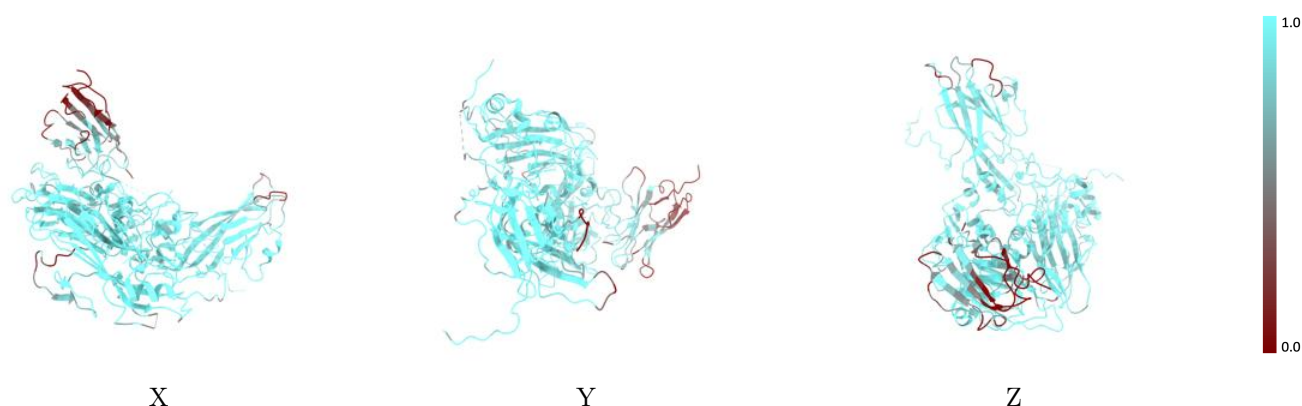
The images above show the 3D surface view of the map at the recommended contour level 0.00707 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



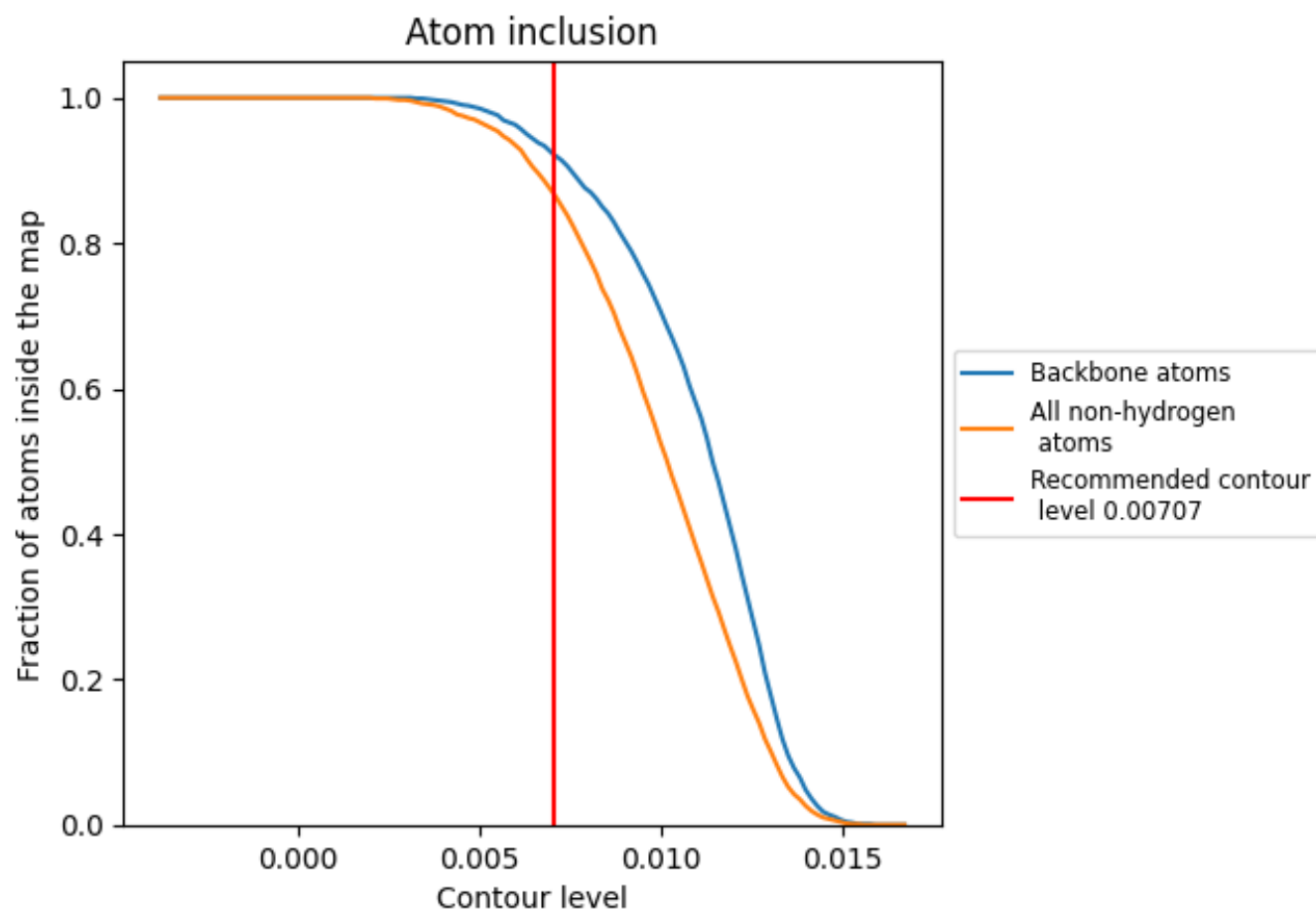
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.00707).

9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 87% 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 (0.00707) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8670	0.2900
1	0.9060	0.3140
2	0.9230	0.2850
3	0.9460	0.2920
7	0.5340	0.2590

