



Full wwPDB EM Validation Report ⓘ

Mar 24, 2026 – 12:51 PM UTC

PDB ID : 7EO0 / pdb_00007eo0
EMDB ID : EMD-31223
Title : FOOT AND MOUTH DISEASE VIRUS O/TIBET/99-BOUND THE SINGLE CHAIN FRAGMENT ANTIBODY C4
Authors : He, Y.; Li, K.
Deposited on : 2021-04-21
Resolution : 3.75 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:

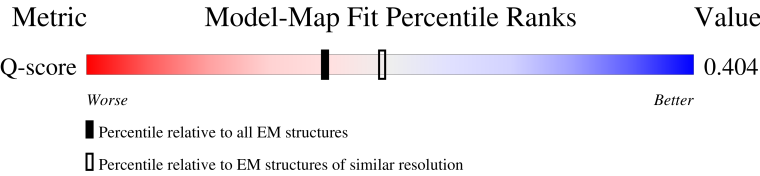
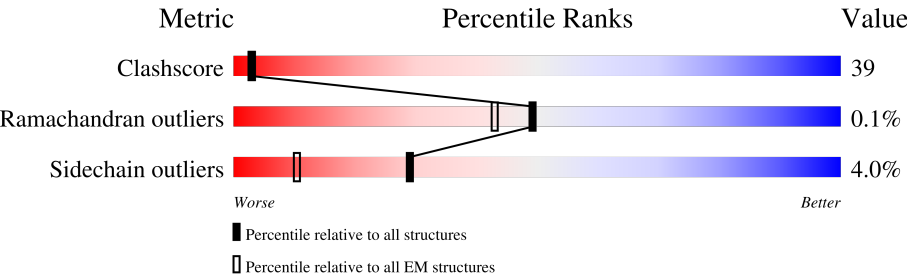
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.75 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	10301 (3.25 - 4.25)

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	213	5% 51% 27% 7% 15%
2	2	218	5% 50% 35% 8% 6%
3	3	220	5% 55% 38% 6%
4	4	85	13% 38% 11% 48%

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Mol	Chain	Length	Quality of chain
5	H	129	32% 46% 48% 5%
6	L	122	43% 45% 38% 6% 11%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6802 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 O/TIBET/99 VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	182	Total	C	N	O	S	0	0
			1421	901	252	265	3		

- Molecule 2 is a protein called O/TIBET/99 VP2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	2	205	Total	C	N	O	S	0	0
			1607	1026	275	299	7		

- Molecule 3 is a protein called O/TIBET/99 VP3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	219	Total	C	N	O	S	0	0
			1673	1073	272	319	9		

- Molecule 4 is a protein called O/TIBET/99 VP4.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	4	44	Total	C	N	O	S	0	0
			337	214	54	67	2		

- Molecule 5 is a protein called Ig heavy chain variable region.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	H	129	Total	C	N	O	S	0	0
			975	609	165	196	5		

- Molecule 6 is a protein called Ig lamda chain variable region.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	L	108	Total	C	N	O	S	0	0
			789	487	136	164	2		

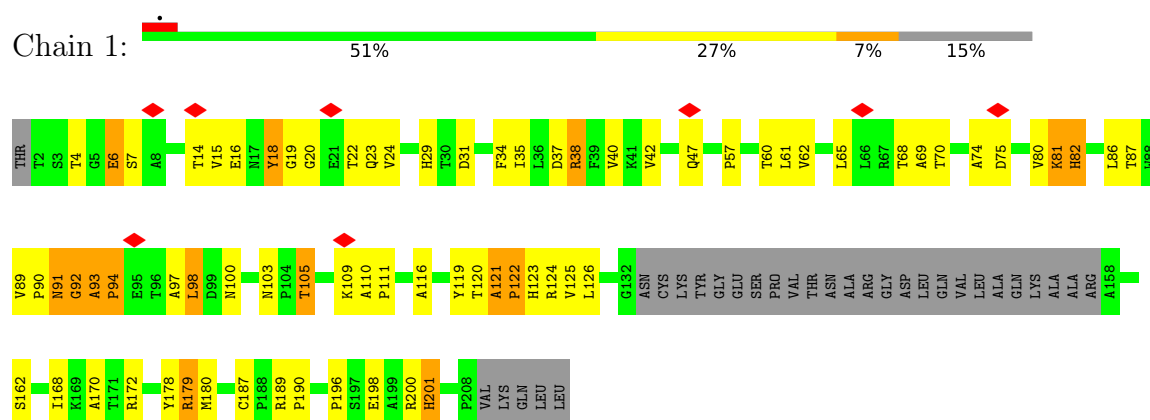
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	113	ASP	-	expression tag	UNP A0A6B9SCH7
L	114	TYR	-	expression tag	UNP A0A6B9SCH7
L	115	LYS	-	expression tag	UNP A0A6B9SCH7
L	116	ASP	-	expression tag	UNP A0A6B9SCH7
L	117	ASP	-	expression tag	UNP A0A6B9SCH7
L	118	ASP	-	expression tag	UNP A0A6B9SCH7
L	119	ASP	-	expression tag	UNP A0A6B9SCH7
L	120	LYS	-	expression tag	UNP A0A6B9SCH7
L	121	GLY	-	expression tag	UNP A0A6B9SCH7
L	122	GLY	-	expression tag	UNP A0A6B9SCH7

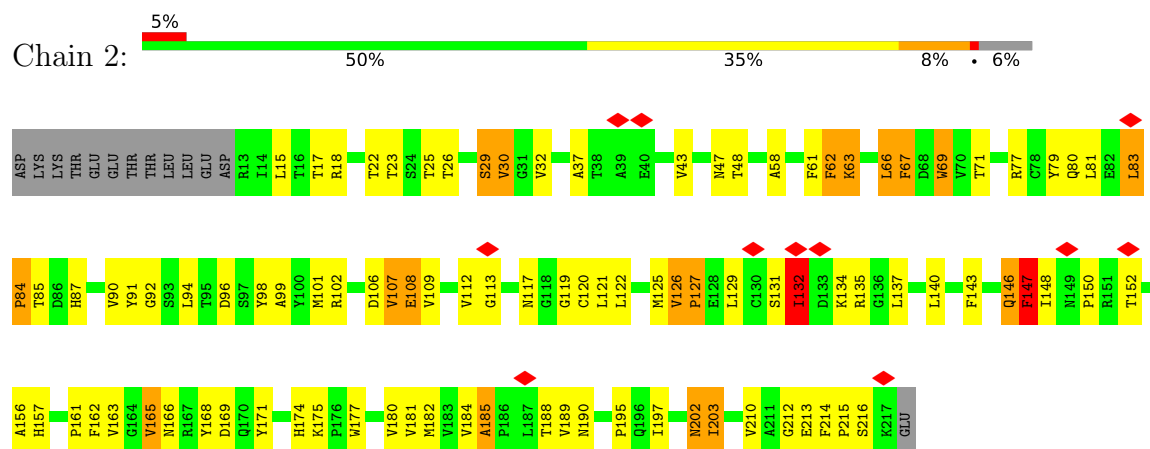
3 Residue-property plots

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.

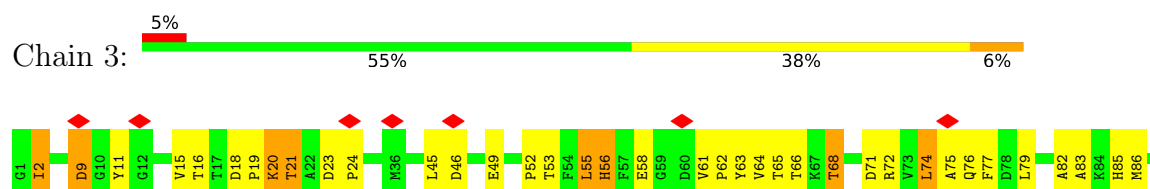
• Molecule 1: O/TIBET/99 VP1



• Molecule 2: O/TIBET/99 VP2



• Molecule 3: O/TIBET/99 VP3



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	14458	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.130	Depositor
Minimum map value	-0.086	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.023	Depositor
Map size ($\text{\AA}$)	446.4, 446.4, 446.4	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.93, 0.93, 0.93	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	0.97	4/1455 (0.3%)	1.59	30/1991 (1.5%)
2	2	0.93	5/1651 (0.3%)	1.42	32/2257 (1.4%)
3	3	0.70	1/1723 (0.1%)	1.62	40/2355 (1.7%)
4	4	0.59	0/343	1.67	13/459 (2.8%)
5	H	0.56	0/995	1.23	18/1353 (1.3%)
6	L	0.61	0/803	1.13	12/1093 (1.1%)
All	All	0.79	10/6970 (0.1%)	1.47	145/9508 (1.5%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1	94	PRO	N-CA	18.70	1.70	1.47
1	1	122	PRO	N-CA	17.99	1.70	1.47
2	2	127	PRO	N-CA	17.98	1.70	1.47
3	3	23	ASP	C-N	14.10	1.50	1.33
2	2	84	PRO	N-CA	13.06	1.69	1.47
1	1	93	ALA	C-N	8.94	1.45	1.33
2	2	216	SER	C-N	-8.46	1.21	1.33
2	2	83	LEU	C-N	6.02	1.45	1.34
1	1	121	ALA	C-N	5.06	1.45	1.33
2	2	126	VAL	C-N	5.01	1.45	1.33

All (145) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	3	74	LEU	N-CA-C	23.25	139.77	111.40
1	1	18	TYR	N-CA-C	-21.47	78.69	109.18
1	1	93	ALA	CA-C-N	18.65	138.99	119.78
1	1	93	ALA	C-N-CA	18.65	138.99	119.78
3	3	23	ASP	CA-C-N	16.11	136.81	119.90
3	3	23	ASP	C-N-CA	16.11	136.81	119.90
3	3	20	LYS	N-CA-C	-16.02	84.28	108.79
1	1	81	LYS	N-CA-C	-15.85	84.76	109.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	2	126	VAL	CA-C-N	15.48	139.20	119.84
2	2	126	VAL	C-N-CA	15.48	139.20	119.84
1	1	121	ALA	CA-C-N	15.18	138.81	119.84
1	1	121	ALA	C-N-CA	15.18	138.81	119.84
3	3	209	PHE	N-CA-C	-14.87	83.62	108.76
3	3	75	ALA	N-CA-C	-14.00	85.86	108.55
4	4	18	THR	N-CA-C	13.68	130.07	110.23
3	3	21	THR	N-CA-C	-13.40	84.97	108.69
3	3	24	PRO	CB-CA-C	-13.01	95.72	111.46
5	H	64	LYS	N-CA-C	-12.95	89.38	109.07
2	2	106	ASP	N-CA-C	-12.91	87.68	109.24
1	1	19	GLY	N-CA-C	11.26	139.85	113.18
2	2	146	GLN	N-CA-C	-11.09	88.52	108.48
3	3	83	ALA	N-CA-C	-10.95	90.03	108.75
4	4	83	LEU	N-CA-C	-10.84	91.76	109.96
4	4	83	LEU	CB-CA-C	10.73	127.23	109.53
3	3	19	PRO	N-CA-C	-10.67	93.75	110.50
1	1	81	LYS	CB-CA-C	10.65	128.85	109.37
2	2	84	PRO	CA-N-CD	-10.50	96.80	111.50
1	1	105	THR	N-CA-C	-10.31	93.74	109.41
1	1	82	HIS	N-CA-CB	10.29	127.82	111.56
3	3	158	SER	N-CA-C	-10.16	89.94	108.02
2	2	147	PHE	N-CA-C	-9.96	92.17	108.41
5	H	66	ARG	CB-CA-C	-9.94	96.77	112.07
2	2	147	PHE	N-CA-CB	9.66	125.72	110.65
3	3	20	LYS	CB-CA-C	9.62	128.51	110.62
2	2	62	PHE	CB-CA-C	9.61	130.90	109.56
3	3	75	ALA	N-CA-CB	9.61	127.99	110.99
6	L	47	LEU	CB-CA-C	-9.59	94.54	109.89
4	4	84	LEU	N-CA-C	-9.44	95.06	109.41
3	3	210	GLU	N-CA-C	-9.13	93.99	109.24
4	4	75	ALA	N-CA-CB	9.01	127.44	111.13
5	H	23	THR	N-CA-C	-8.98	96.50	109.96
2	2	62	PHE	N-CA-C	-8.92	95.40	109.23
3	3	76	GLN	N-CA-C	-8.82	94.07	109.06
3	3	83	ALA	N-CA-CB	8.69	126.85	111.13
1	1	92	GLY	N-CA-C	-8.52	92.98	113.18
3	3	141	HIS	N-CA-C	8.42	120.08	111.07
1	1	94	PRO	CA-N-CD	-8.40	100.23	112.00
3	3	46	ASP	N-CA-C	-8.27	103.33	113.41
2	2	216	SER	CB-CA-C	-8.12	94.52	109.46
4	4	75	ALA	N-CA-C	-8.12	94.86	108.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	179	ARG	N-CA-C	-8.07	96.32	108.99
3	3	21	THR	N-CA-CB	8.03	124.31	110.81
3	3	175	ALA	N-CA-C	7.93	120.93	111.33
2	2	127	PRO	CA-N-CD	-7.92	100.92	112.00
6	L	18	GLN	N-CA-C	-7.90	97.02	109.50
1	1	122	PRO	CA-N-CD	-7.87	100.98	112.00
2	2	216	SER	O-C-N	-7.86	114.08	122.96
5	H	23	THR	CB-CA-C	7.86	122.47	109.89
5	H	67	LEU	N-CA-CB	7.80	124.79	110.99
2	2	106	ASP	CB-CA-C	7.75	124.69	109.35
5	H	12	VAL	N-CA-C	7.71	121.96	109.78
3	3	24	PRO	N-CA-C	7.68	122.40	110.80
1	1	105	THR	CB-CA-C	-7.59	95.27	110.30
2	2	203	ILE	N-CA-C	-7.51	97.47	108.89
2	2	146	GLN	CB-CA-C	7.36	122.32	109.80
3	3	209	PHE	CB-CA-C	7.36	123.02	109.70
3	3	157	PHE	N-CA-C	-7.32	97.60	108.79
1	1	179	ARG	CB-CA-C	7.31	125.52	109.94
5	H	64	LYS	N-CA-CB	7.29	123.83	111.66
2	2	185	ALA	N-CA-C	-7.16	101.05	110.07
5	H	63	LEU	N-CA-C	-7.15	91.10	107.48
3	3	129	GLY	N-CA-C	7.11	124.53	113.86
3	3	76	GLN	CB-CA-C	7.10	123.80	109.38
2	2	202	ASN	N-CA-C	-7.08	96.09	108.20
6	L	80	SER	N-CA-C	-7.08	97.88	108.99
3	3	157	PHE	CB-CA-C	7.04	123.71	110.62
6	L	81	LEU	N-CA-C	-7.01	99.27	110.14
6	L	48	ARG	N-CA-CB	-6.88	100.73	111.05
4	4	33	SER	CB-CA-C	-6.84	97.82	109.51
2	2	107	VAL	N-CA-C	-6.81	97.77	108.81
1	1	82	HIS	N-CA-C	-6.80	98.82	108.96
5	H	97	LYS	N-CA-CB	-6.77	99.94	110.57
3	3	55	LEU	CB-CA-C	-6.73	97.60	109.71
2	2	107	VAL	N-CA-CB	6.72	124.48	111.92
1	1	105	THR	CA-C-O	-6.71	113.79	121.58
6	L	4	ALA	CA-C-N	-6.71	109.89	121.97
6	L	4	ALA	C-N-CA	-6.71	109.89	121.97
2	2	63	LYS	N-CA-C	-6.71	98.79	109.59
6	L	80	SER	CB-CA-C	6.71	124.23	109.94
2	2	67	PHE	N-CA-C	-6.68	101.71	110.33
3	3	74	LEU	CB-CA-C	-6.66	99.59	110.79
1	1	91	ASN	N-CA-C	6.66	118.19	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	31	SER	N-CA-C	6.64	121.38	113.28
1	1	74	ALA	CB-CA-C	6.60	122.74	109.68
5	H	96	ALA	CB-CA-C	-6.59	94.94	109.56
4	4	74	SER	N-CA-C	-6.56	94.84	107.57
1	1	93	ALA	N-CA-CB	6.55	122.31	111.17
1	1	116	ALA	N-CA-C	-6.50	97.09	108.20
6	L	48	ARG	CB-CA-C	6.47	122.41	110.24
3	3	130	MET	N-CA-CB	6.34	121.57	111.56
4	4	18	THR	N-CA-CB	-6.34	100.26	109.83
3	3	82	ALA	N-CA-C	-6.30	96.03	107.75
5	H	24	VAL	N-CA-C	-6.22	99.52	108.53
3	3	56	HIS	N-CA-C	6.19	118.83	108.73
3	3	77	PHE	N-CA-CB	6.13	122.12	111.39
4	4	84	LEU	N-CA-CB	6.06	121.06	111.43
3	3	188	GLN	N-CA-C	-6.04	97.87	108.20
2	2	132	ILE	N-CA-C	6.03	117.20	109.30
2	2	83	LEU	CA-C-N	5.96	141.31	127.00
2	2	83	LEU	C-N-CA	5.96	141.31	127.00
3	3	19	PRO	CB-CA-C	5.96	118.05	111.56
4	4	17	ASN	N-CA-C	5.96	120.09	112.34
1	1	69	ALA	N-CA-CB	-5.95	101.22	110.57
3	3	146	GLU	N-CA-C	-5.91	100.35	109.52
1	1	172	ARG	CB-CA-C	5.81	120.80	110.16
5	H	82	VAL	N-CA-CB	5.77	119.67	111.82
3	3	9	ASP	CB-CA-C	5.74	121.25	110.63
2	2	66	LEU	CA-C-N	5.74	131.16	120.95
2	2	66	LEU	C-N-CA	5.74	131.16	120.95
3	3	158	SER	N-CA-CB	5.72	119.53	110.84
6	L	107	THR	N-CA-C	-5.71	101.60	110.10
2	2	108	GLU	N-CA-C	-5.66	99.07	108.75
2	2	108	GLU	CB-CA-C	5.61	120.78	109.68
3	3	158	SER	CB-CA-C	5.50	118.92	110.62
4	4	34	MET	N-CA-CB	-5.48	101.49	110.42
2	2	15	LEU	N-CA-C	-5.45	99.40	108.34
2	2	132	ILE	CB-CA-C	-5.43	103.14	111.71
2	2	30	VAL	N-CA-C	5.39	121.20	113.16
1	1	180	MET	N-CA-C	-5.36	96.93	107.69
6	L	47	LEU	N-CA-C	5.32	117.94	109.96
1	1	122	PRO	N-CD-CG	5.32	111.18	103.20
6	L	107	THR	CB-CA-C	5.31	118.19	109.90
1	1	75	ASP	N-CA-C	-5.29	101.08	109.76
4	4	34	MET	N-CA-C	5.25	117.29	108.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	82	VAL	N-CA-C	-5.23	100.03	107.77
1	1	116	ALA	CB-CA-C	5.22	119.13	110.78
5	H	28	SER	N-CA-C	5.20	117.23	109.59
5	H	106	CYS	CA-C-N	5.17	126.60	120.14
5	H	106	CYS	C-N-CA	5.17	126.60	120.14
2	2	29	SER	CB-CA-C	-5.15	99.23	109.79
3	3	140	ALA	CB-CA-C	5.14	120.14	110.63
1	1	105	THR	CA-C-N	5.10	131.15	122.37
1	1	105	THR	C-N-CA	5.10	131.15	122.37
3	3	9	ASP	N-CA-C	-5.03	105.92	111.71
5	H	16	GLN	N-CA-CB	-5.02	102.24	111.37

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1	1421	0	1423	121	0
2	2	1607	0	1562	160	0
3	3	1673	0	1600	114	0
4	4	337	0	314	8	0
5	H	975	0	952	109	0
6	L	789	0	766	91	0
All	All	6802	0	6617	520	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 39.

All (520) 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)
3:3:58:GLU:OE2	5:H:59:TYR:CE2	1.64	1.49
2:2:84:PRO:N	2:2:84:PRO:CA	1.69	1.46
1:1:94:PRO:N	1:1:94:PRO:CA	1.70	1.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:122:PRO:N	1:1:122:PRO:CA	1.70	1.38
2:2:84:PRO:CD	2:2:175:LYS:HG2	1.53	1.37
5:H:98:TRP:CE2	5:H:108:VAL:HG23	1.59	1.37
2:2:63:LYS:NZ	2:2:202:ASN:HD21	1.20	1.36
1:1:89:VAL:CG1	1:1:98:LEU:HD12	1.55	1.34
1:1:196:PRO:HB3	1:1:201:HIS:CB	1.58	1.33
2:2:127:PRO:N	2:2:127:PRO:CA	1.70	1.32
1:1:100:ASN:ND2	3:3:216:ASP:OD2	1.64	1.28
2:2:81:LEU:HD11	2:2:85:THR:CG2	1.63	1.26
2:2:63:LYS:NZ	2:2:202:ASN:ND2	1.85	1.25
3:3:65:THR:CG2	3:3:196:GLY:HA2	1.66	1.24
2:2:94:LEU:HD22	2:2:101:MET:SD	1.78	1.23
3:3:58:GLU:OE2	5:H:59:TYR:HE2	0.93	1.23
5:H:58:GLY:HA3	5:H:105:ASP:OD1	1.37	1.22
6:L:5:VAL:CG2	6:L:101:VAL:CG1	2.18	1.21
1:1:125:VAL:O	2:2:165:VAL:HG23	1.39	1.20
2:2:94:LEU:CD2	2:2:101:MET:SD	2.30	1.19
2:2:63:LYS:HZ2	2:2:202:ASN:ND2	1.39	1.16
5:H:107:GLY:HA2	6:L:94:TYR:CD2	1.80	1.16
5:H:98:TRP:CH2	5:H:114:SER:HB2	1.83	1.14
1:1:34:PHE:HE1	1:1:38:ARG:NH1	1.45	1.14
1:1:196:PRO:HB3	1:1:201:HIS:HB2	1.24	1.13
6:L:5:VAL:CG2	6:L:101:VAL:HG12	1.78	1.13
2:2:83:LEU:CD2	2:2:125:MET:HE1	1.80	1.12
5:H:51:ILE:HA	5:H:57:THR:HG22	1.30	1.12
1:1:89:VAL:CG1	1:1:98:LEU:CD1	2.27	1.10
2:2:84:PRO:HD2	2:2:175:LYS:HG2	1.16	1.09
1:1:89:VAL:CG2	1:1:98:LEU:HD11	1.83	1.09
1:1:89:VAL:HG22	1:1:98:LEU:HD11	1.29	1.09
1:1:125:VAL:O	2:2:165:VAL:CG2	2.00	1.08
1:1:196:PRO:HB3	1:1:201:HIS:HB3	1.30	1.07
2:2:83:LEU:O	2:2:175:LYS:HB3	1.53	1.07
5:H:109:HIS:CD2	6:L:102:PHE:HZ	1.73	1.06
1:1:87:THR:HG22	1:1:105:THR:HG22	1.32	1.05
1:1:92:GLY:O	3:3:99:THR:HB	1.54	1.05
5:H:98:TRP:CE2	5:H:108:VAL:CG2	2.38	1.05
2:2:113:GLY:HA3	2:2:150:PRO:HB3	1.35	1.05
2:2:81:LEU:CD1	2:2:85:THR:HG21	1.86	1.04
3:3:45:LEU:O	3:3:49:GLU:HG3	1.57	1.04
5:H:98:TRP:CD2	5:H:108:VAL:HG23	1.93	1.03
5:H:35:ASN:ND2	5:H:103:GLY:O	1.93	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:84:PRO:HD3	2:2:175:LYS:HG2	1.39	1.01
4:4:79:LEU:HD23	4:4:79:LEU:H	1.26	1.01
2:2:126:VAL:CG2	2:2:129:LEU:HB2	1.91	1.00
1:1:89:VAL:HG13	1:1:98:LEU:HD12	1.03	0.99
6:L:5:VAL:HG22	6:L:101:VAL:HG12	1.43	0.99
6:L:29:SER:HB3	6:L:96:ILE:HG22	1.45	0.99
1:1:34:PHE:CE1	1:1:38:ARG:NH1	2.31	0.98
3:3:49:GLU:HG2	3:3:205:ALA:HB3	1.44	0.97
2:2:83:LEU:HD23	2:2:125:MET:HE1	1.47	0.96
6:L:32:GLY:HA2	6:L:69:LYS:HE3	1.46	0.96
1:1:123:HIS:HB3	2:2:165:VAL:HG22	1.48	0.95
1:1:89:VAL:CG2	1:1:98:LEU:CD1	2.44	0.94
3:3:65:THR:HG23	3:3:196:GLY:HA2	1.48	0.94
2:2:126:VAL:HG23	2:2:129:LEU:HB2	1.47	0.94
2:2:84:PRO:CD	2:2:175:LYS:CG	2.46	0.94
1:1:89:VAL:HG13	1:1:98:LEU:CD1	1.93	0.93
1:1:123:HIS:HE1	1:1:162:SER:OG	1.51	0.92
1:1:35:ILE:HD11	1:1:61:LEU:HD11	1.50	0.92
1:1:103:ASN:HD21	3:3:216:ASP:HB3	1.34	0.92
1:1:92:GLY:O	3:3:99:THR:CB	2.17	0.91
5:H:109:HIS:CD2	6:L:102:PHE:CZ	2.59	0.91
5:H:98:TRP:NE1	5:H:108:VAL:CG2	2.34	0.90
1:1:89:VAL:HG11	1:1:98:LEU:CD1	2.00	0.90
2:2:81:LEU:HD11	2:2:85:THR:HG21	0.90	0.90
1:1:92:GLY:O	3:3:99:THR:CG2	2.20	0.90
1:1:123:HIS:CE1	1:1:162:SER:OG	2.25	0.90
2:2:63:LYS:CE	2:2:202:ASN:HD21	1.84	0.89
3:3:119:ALA:O	3:3:149:THR:HG23	1.73	0.89
6:L:29:SER:HB3	6:L:96:ILE:CG2	2.02	0.89
6:L:5:VAL:CG2	6:L:101:VAL:HG11	2.02	0.88
6:L:32:GLY:HA2	6:L:69:LYS:CE	2.02	0.88
2:2:126:VAL:HG21	2:2:129:LEU:HD22	1.53	0.88
1:1:196:PRO:CB	1:1:201:HIS:HB2	2.03	0.88
6:L:78:ILE:HG22	6:L:80:SER:O	1.74	0.88
2:2:129:LEU:HD22	2:2:180:VAL:HG21	1.56	0.87
1:1:123:HIS:HB3	2:2:165:VAL:CG2	2.04	0.87
1:1:87:THR:CG2	1:1:105:THR:HG22	2.04	0.87
3:3:45:LEU:HD22	3:3:209:PHE:HD2	1.40	0.86
6:L:47:LEU:HD12	6:L:47:LEU:O	1.74	0.86
1:1:196:PRO:CB	1:1:201:HIS:CB	2.51	0.85
2:2:83:LEU:CD2	2:2:125:MET:CE	2.54	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:87:THR:HA	1:1:105:THR:HA	1.58	0.84
2:2:81:LEU:CD1	2:2:85:THR:CG2	2.51	0.84
3:3:100:GLN:NE2	3:3:213:LEU:HD12	1.91	0.84
5:H:98:TRP:CD1	5:H:108:VAL:CG2	2.60	0.84
5:H:98:TRP:CD1	5:H:116:LEU:HD12	2.12	0.84
1:1:34:PHE:HE1	1:1:38:ARG:HH12	0.86	0.84
2:2:190:ASN:HD21	2:2:195:PRO:HA	1.43	0.84
5:H:107:GLY:HA2	6:L:94:TYR:CE2	2.13	0.84
6:L:5:VAL:HG23	6:L:101:VAL:CG1	2.08	0.83
1:1:105:THR:O	3:3:15:VAL:O	1.96	0.83
1:1:89:VAL:HG11	1:1:98:LEU:HD12	1.54	0.83
6:L:33:ARG:HG2	6:L:33:ARG:HH21	1.44	0.83
2:2:63:LYS:HZ3	2:2:202:ASN:ND2	1.76	0.83
6:L:5:VAL:HG21	6:L:101:VAL:CG1	2.09	0.83
1:1:38:ARG:HG2	1:1:38:ARG:HH21	1.44	0.83
2:2:83:LEU:CB	2:2:177:TRP:O	2.27	0.82
5:H:51:ILE:HA	5:H:57:THR:CG2	2.08	0.82
5:H:60:ASN:O	5:H:63:LEU:O	1.98	0.82
3:3:65:THR:HG22	3:3:196:GLY:HA2	1.60	0.82
3:3:100:GLN:HE21	3:3:212:ARG:HE	1.25	0.82
5:H:45:LEU:HB2	6:L:102:PHE:CD2	2.14	0.82
3:3:207:LYS:NZ	3:3:207:LYS:HB3	1.95	0.82
1:1:68:THR:O	1:1:189:ARG:N	2.12	0.82
2:2:84:PRO:HD2	2:2:175:LYS:CG	2.07	0.81
5:H:66:ARG:O	5:H:67:LEU:HD23	1.80	0.81
2:2:84:PRO:HD3	2:2:175:LYS:CG	2.08	0.81
5:H:98:TRP:CZ2	5:H:114:SER:HB2	2.15	0.81
2:2:119:GLY:N	2:2:189:VAL:HG22	1.95	0.81
1:1:90:PRO:HD2	3:3:215:VAL:HG23	1.63	0.80
3:3:45:LEU:HD22	3:3:209:PHE:CD2	2.15	0.80
2:2:119:GLY:H	2:2:189:VAL:HG22	1.45	0.79
2:2:129:LEU:CD2	2:2:180:VAL:HG21	2.12	0.79
1:1:103:ASN:ND2	3:3:216:ASP:HB3	1.97	0.78
2:2:129:LEU:CD2	2:2:180:VAL:CG2	2.61	0.78
1:1:16:GLU:C	1:1:18:TYR:O	2.26	0.78
5:H:98:TRP:CH2	5:H:114:SER:CB	2.66	0.78
2:2:102:ARG:HB2	2:2:168:TYR:HD1	1.49	0.78
6:L:5:VAL:HG21	6:L:101:VAL:HG12	1.64	0.78
5:H:51:ILE:CA	5:H:57:THR:HG22	2.10	0.77
5:H:98:TRP:CD2	5:H:108:VAL:CG2	2.63	0.77
2:2:63:LYS:CD	2:2:202:ASN:HD21	1.96	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:6:GLU:OE2	2:2:152:THR:HG21	1.84	0.77
3:3:103:GLY:HA3	3:3:209:PHE:HA	1.65	0.77
1:1:38:ARG:HG2	1:1:38:ARG:NH2	1.99	0.77
2:2:22:THR:HG21	2:2:61:PHE:H	1.50	0.77
2:2:126:VAL:HG21	2:2:129:LEU:HB2	1.67	0.76
3:3:61:VAL:HG11	5:H:55:GLY:HA3	1.66	0.76
3:3:108:HIS:HB2	3:3:202:LEU:HB2	1.67	0.76
1:1:120:THR:OG1	3:3:213:LEU:CD1	2.34	0.76
2:2:63:LYS:CD	2:2:202:ASN:ND2	2.49	0.76
6:L:5:VAL:HG22	6:L:101:VAL:CG1	2.03	0.76
4:4:79:LEU:H	4:4:79:LEU:CD2	1.98	0.76
1:1:86:LEU:O	1:1:105:THR:HA	1.87	0.75
2:2:63:LYS:HD3	2:2:202:ASN:ND2	2.02	0.75
2:2:63:LYS:HZ2	2:2:202:ASN:HD21	0.76	0.75
1:1:70:THR:HG22	1:1:189:ARG:HE	1.50	0.75
5:H:98:TRP:NE1	5:H:108:VAL:HG22	2.01	0.74
3:3:18:ASP:OD2	3:3:20:LYS:HD2	1.88	0.74
5:H:50:GLY:N	5:H:104:TYR:O	2.20	0.74
1:1:89:VAL:HG21	1:1:98:LEU:CD1	2.18	0.74
3:3:45:LEU:O	3:3:49:GLU:CG	2.35	0.73
3:3:102:SER:HB3	3:3:210:GLU:HG3	1.70	0.73
2:2:58:ALA:HB2	2:2:90:VAL:HG12	1.69	0.73
3:3:108:HIS:HD2	3:3:156:THR:HG22	1.54	0.73
3:3:65:THR:HG23	3:3:196:GLY:CA	2.19	0.73
3:3:102:SER:HB3	3:3:210:GLU:CG	2.19	0.73
2:2:58:ALA:HB2	2:2:90:VAL:CG1	2.19	0.72
2:2:120:CYS:H	2:2:185:ALA:HB3	1.54	0.72
5:H:98:TRP:CZ2	5:H:108:VAL:HG23	2.20	0.72
1:1:196:PRO:CB	1:1:201:HIS:HB3	2.13	0.72
1:1:31:ASP:O	1:1:35:ILE:HG22	1.88	0.72
2:2:32:VAL:HG23	2:2:157:HIS:ND1	2.05	0.72
6:L:5:VAL:HG23	6:L:101:VAL:HG11	1.69	0.72
2:2:58:ALA:CB	2:2:90:VAL:HG12	2.19	0.72
1:1:200:ARG:O	1:1:200:ARG:HG2	1.90	0.72
2:2:71:THR:HA	2:2:188:THR:HG22	1.72	0.71
4:4:79:LEU:HD23	4:4:79:LEU:N	2.04	0.70
1:1:80:VAL:HG12	1:1:81:LYS:O	1.91	0.70
1:1:87:THR:HG22	1:1:105:THR:CG2	2.19	0.70
1:1:120:THR:OG1	3:3:213:LEU:HD13	1.92	0.70
2:2:32:VAL:HG23	2:2:157:HIS:HD1	1.56	0.70
1:1:198:GLU:N	1:1:198:GLU:OE1	2.23	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:132:ILE:HD12	2:2:140:LEU:CD1	2.21	0.70
2:2:98:TYR:CD2	2:2:210:VAL:HB	2.27	0.70
2:2:83:LEU:HB3	2:2:177:TRP:O	1.92	0.69
5:H:13:LYS:HD3	5:H:14:PRO:HD2	1.73	0.69
2:2:129:LEU:HD22	2:2:180:VAL:CG2	2.21	0.69
5:H:51:ILE:HG13	5:H:57:THR:CG2	2.22	0.69
1:1:89:VAL:HG22	1:1:98:LEU:CD1	2.12	0.68
2:2:77:ARG:HH22	2:2:79:TYR:HB2	1.58	0.68
3:3:120:ARG:NH2	3:3:148:ASP:OD1	2.24	0.68
5:H:110:SER:HA	6:L:37:THR:HG23	1.76	0.68
5:H:107:GLY:HA2	6:L:94:TYR:HD2	1.45	0.68
2:2:22:THR:HG23	2:2:61:PHE:HD1	1.56	0.68
1:1:123:HIS:HD2	2:2:166:ASN:HD21	1.39	0.68
3:3:74:LEU:HD11	3:3:188:GLN:HE21	1.59	0.68
5:H:24:VAL:HB	5:H:76:SER:O	1.93	0.68
6:L:9:PRO:HD3	6:L:23:THR:O	1.93	0.68
3:3:66:THR:HG23	3:3:188:GLN:CD	2.19	0.67
6:L:5:VAL:CG2	6:L:101:VAL:CB	2.72	0.67
1:1:7:SER:HA	2:2:146:GLN:NE2	2.09	0.67
5:H:57:THR:HA	5:H:104:TYR:HB2	1.77	0.67
3:3:151:LEU:O	3:3:151:LEU:HD23	1.95	0.67
1:1:189:ARG:NH2	2:2:127:PRO:HG2	2.10	0.67
3:3:65:THR:HG1	5:H:54:SER:HG	1.43	0.67
2:2:140:LEU:HD23	2:2:140:LEU:O	1.94	0.67
5:H:98:TRP:NE1	5:H:116:LEU:HD12	2.09	0.66
5:H:98:TRP:CZ2	5:H:114:SER:CB	2.80	0.66
5:H:53:THR:HG23	5:H:101:ARG:O	1.97	0.65
1:1:189:ARG:HG3	1:1:190:PRO:HD2	1.78	0.65
2:2:22:THR:CG2	2:2:61:PHE:H	2.09	0.65
2:2:120:CYS:HB3	2:2:185:ALA:HB2	1.78	0.65
6:L:33:ARG:HH21	6:L:33:ARG:CG	2.08	0.65
2:2:94:LEU:HD21	2:2:101:MET:SD	2.33	0.65
3:3:102:SER:O	3:3:102:SER:OG	2.13	0.65
3:3:65:THR:CG2	3:3:196:GLY:CA	2.59	0.65
3:3:207:LYS:HB3	3:3:207:LYS:HZ3	1.61	0.65
6:L:32:GLY:O	6:L:69:LYS:NZ	2.25	0.65
1:1:189:ARG:HH21	2:2:127:PRO:HG2	1.60	0.65
3:3:58:GLU:OE2	5:H:59:TYR:CD2	2.44	0.65
1:1:37:ASP:O	1:1:179:ARG:HD2	1.96	0.64
5:H:108:VAL:HG22	5:H:116:LEU:HD11	1.78	0.64
6:L:32:GLY:CA	6:L:69:LYS:HE3	2.24	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:107:GLY:CA	6:L:94:TYR:CE2	2.80	0.64
5:H:98:TRP:CG	5:H:108:VAL:CG2	2.80	0.64
2:2:132:ILE:CD1	2:2:140:LEU:CD1	2.75	0.64
5:H:98:TRP:CG	5:H:108:VAL:HG21	2.32	0.64
6:L:5:VAL:HG23	6:L:5:VAL:O	1.97	0.64
2:2:69:TRP:HE1	2:2:188:THR:HB	1.63	0.63
2:2:80:GLN:HE22	2:2:131:SER:HA	1.61	0.63
1:1:120:THR:OG1	3:3:213:LEU:HD11	1.98	0.63
3:3:101:TYR:HA	3:3:211:LEU:HD12	1.80	0.63
1:1:89:VAL:HG11	1:1:98:LEU:HD13	1.77	0.63
4:4:17:ASN:OD1	4:4:17:ASN:O	2.16	0.63
6:L:7:THR:HB	6:L:25:SER:HB2	1.79	0.63
1:1:35:ILE:HG12	1:1:65:LEU:HD22	1.81	0.63
5:H:109:HIS:NE2	6:L:102:PHE:HZ	1.97	0.62
5:H:61:PRO:HD2	6:L:99:ASN:HB3	1.80	0.62
2:2:126:VAL:HG21	2:2:129:LEU:CD2	2.29	0.62
1:1:92:GLY:O	3:3:99:THR:HG22	1.99	0.62
6:L:36:ALA:HB2	6:L:93:ALA:HB2	1.82	0.62
1:1:4:THR:HG21	4:4:76:PHE:CD2	2.35	0.62
2:2:83:LEU:HB2	2:2:177:TRP:O	2.00	0.61
5:H:98:TRP:CD1	5:H:108:VAL:HG21	2.35	0.61
2:2:137:LEU:HD23	2:2:182:MET:SD	2.39	0.61
1:1:103:ASN:OD1	3:3:216:ASP:N	2.34	0.61
2:2:102:ARG:HB2	2:2:168:TYR:CD1	2.33	0.61
5:H:110:SER:OG	6:L:94:TYR:CB	2.49	0.61
1:1:123:HIS:CB	2:2:165:VAL:CG2	2.77	0.61
3:3:159:ILE:HD13	3:3:184:VAL:HG11	1.83	0.61
3:3:151:LEU:HD23	3:3:151:LEU:C	2.26	0.60
5:H:39:GLN:OE1	6:L:41:GLN:NE2	2.34	0.60
5:H:98:TRP:CD1	5:H:115:TYR:O	2.54	0.60
6:L:11:SER:HB3	6:L:107:THR:O	2.01	0.60
6:L:47:LEU:O	6:L:47:LEU:CD1	2.49	0.60
2:2:83:LEU:HD22	2:2:125:MET:CE	2.30	0.60
2:2:22:THR:HG23	2:2:61:PHE:CD1	2.35	0.60
2:2:135:ARG:HH11	2:2:135:ARG:HG2	1.65	0.60
3:3:101:TYR:HD2	3:3:170:THR:HG22	1.66	0.60
5:H:110:SER:OG	6:L:94:TYR:HB2	2.00	0.60
3:3:66:THR:HG23	3:3:188:GLN:NE2	2.17	0.60
1:1:86:LEU:O	1:1:105:THR:CA	2.50	0.60
2:2:108:GLU:OE1	2:2:157:HIS:CD2	2.55	0.60
1:1:123:HIS:HD2	2:2:166:ASN:ND2	2.00	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:15:VAL:HG12	1:1:15:VAL:O	2.02	0.59
1:1:16:GLU:O	1:1:18:TYR:O	2.21	0.59
1:1:24:VAL:HG12	1:1:24:VAL:O	2.01	0.59
2:2:62:PHE:CD2	2:2:87:HIS:HE1	2.20	0.59
5:H:98:TRP:CD1	5:H:116:LEU:CD1	2.84	0.59
1:1:89:VAL:HG21	1:1:98:LEU:HD11	1.77	0.59
2:2:214:PHE:CZ	3:3:129:GLY:HA3	2.38	0.59
3:3:72:ARG:NH2	6:L:34:TYR:CD2	2.71	0.58
2:2:62:PHE:CD2	2:2:87:HIS:CE1	2.91	0.58
5:H:110:SER:HB3	6:L:93:ALA:HA	1.84	0.58
6:L:5:VAL:CG2	6:L:101:VAL:HB	2.34	0.58
3:3:98:TYR:HD2	3:3:211:LEU:HG	1.69	0.58
2:2:132:ILE:CD1	2:2:140:LEU:HD11	2.33	0.58
1:1:7:SER:HA	2:2:146:GLN:HE21	1.68	0.58
2:2:22:THR:HG21	2:2:61:PHE:N	2.17	0.58
5:H:119:TRP:CG	6:L:47:LEU:O	2.57	0.58
2:2:37:ALA:HB3	2:2:161:PRO:HG3	1.85	0.57
3:3:71:ASP:OD1	3:3:136:PRO:HG2	2.05	0.57
2:2:117:ASN:HA	2:2:189:VAL:O	2.04	0.57
2:2:215:PRO:HB3	3:3:133:PRO:HB3	1.87	0.57
2:2:113:GLY:CA	2:2:150:PRO:HB3	2.24	0.57
3:3:135:THR:HG23	3:3:138:ALA:H	1.70	0.56
6:L:5:VAL:HG23	6:L:101:VAL:CB	2.34	0.56
2:2:113:GLY:HA2	2:2:197:ILE:HD12	1.87	0.56
5:H:57:THR:HA	5:H:104:TYR:CB	2.34	0.56
2:2:120:CYS:H	2:2:185:ALA:CB	2.19	0.56
6:L:98:THR:HG22	6:L:98:THR:O	2.05	0.56
1:1:82:HIS:CE1	1:1:86:LEU:HD22	2.41	0.56
2:2:109:VAL:HG22	2:2:156:ALA:HB3	1.88	0.56
1:1:38:ARG:HH21	1:1:38:ARG:CG	2.17	0.56
2:2:22:THR:CG2	2:2:61:PHE:HD1	2.19	0.56
1:1:16:GLU:HB2	1:1:20:GLY:O	2.06	0.56
1:1:40:VAL:HB	1:1:62:VAL:HG11	1.86	0.56
1:1:93:ALA:CB	3:3:215:VAL:HG21	2.37	0.55
1:1:94:PRO:N	1:1:94:PRO:C	2.61	0.55
2:2:58:ALA:HB1	2:2:91:TYR:HB2	1.88	0.55
3:3:79:LEU:HD22	3:3:125:TYR:HE1	1.71	0.55
1:1:82:HIS:ND1	1:1:86:LEU:HD22	2.22	0.55
2:2:83:LEU:C	2:2:175:LYS:HB3	2.28	0.55
6:L:5:VAL:HG22	6:L:101:VAL:CB	2.34	0.55
6:L:5:VAL:HG23	6:L:101:VAL:HB	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:65:THR:HG21	3:3:196:GLY:HA2	1.79	0.55
2:2:83:LEU:HD23	2:2:125:MET:CE	2.28	0.55
6:L:12:VAL:O	6:L:108:LEU:HA	2.07	0.55
5:H:115:TYR:CE2	6:L:58:PRO:HB3	2.41	0.55
1:1:123:HIS:CD2	2:2:166:ASN:HD21	2.23	0.55
2:2:126:VAL:CG2	2:2:129:LEU:HD22	2.33	0.55
6:L:81:LEU:HD21	6:L:108:LEU:HD21	1.88	0.55
3:3:45:LEU:HD11	3:3:211:LEU:HB2	1.89	0.54
6:L:6:LEU:CD2	6:L:26:GLY:HA2	2.38	0.54
2:2:215:PRO:HG2	3:3:133:PRO:HD3	1.90	0.54
5:H:98:TRP:CZ3	5:H:114:SER:CB	2.90	0.54
2:2:163:VAL:HG23	2:2:163:VAL:O	2.08	0.54
1:1:68:THR:O	1:1:189:ARG:HB3	2.07	0.54
3:3:207:LYS:HB3	3:3:207:LYS:HZ2	1.72	0.54
5:H:119:TRP:CD1	6:L:47:LEU:O	2.61	0.54
3:3:63:TYR:HB3	3:3:200:VAL:HG12	1.90	0.54
5:H:115:TYR:CE2	5:H:117:ASP:HB3	2.43	0.54
3:3:49:GLU:HG2	3:3:205:ALA:CB	2.29	0.54
1:1:123:HIS:HE1	1:1:162:SER:CB	2.20	0.53
5:H:98:TRP:CZ3	5:H:114:SER:HB2	2.41	0.53
5:H:98:TRP:HE1	5:H:116:LEU:HD12	1.74	0.53
2:2:132:ILE:HD11	2:2:180:VAL:HG11	1.91	0.53
3:3:65:THR:HG23	3:3:196:GLY:C	2.34	0.53
5:H:51:ILE:CG1	5:H:57:THR:CG2	2.85	0.53
6:L:5:VAL:HG21	6:L:101:VAL:HG11	1.81	0.53
3:3:58:GLU:HB3	5:H:57:THR:O	2.09	0.53
5:H:47:CYS:O	5:H:60:ASN:ND2	2.36	0.53
5:H:45:LEU:CB	6:L:102:PHE:CD2	2.91	0.53
2:2:47:ASN:OD1	2:2:48:THR:N	2.42	0.52
3:3:109:PHE:CD2	3:3:201:VAL:HG12	2.45	0.52
2:2:101:MET:HG3	2:2:210:VAL:HG12	1.91	0.52
5:H:51:ILE:HG13	5:H:57:THR:HG22	1.89	0.52
6:L:10:SER:O	6:L:106:THR:HA	2.09	0.52
5:H:107:GLY:CA	6:L:94:TYR:CD2	2.73	0.52
6:L:30:ASN:HA	6:L:95:ASP:HA	1.92	0.52
1:1:89:VAL:CB	1:1:98:LEU:CD1	2.86	0.52
2:2:169:ASP:OD1	2:2:174:HIS:NE2	2.42	0.52
3:3:195:ASP:O	5:H:101:ARG:HB2	2.09	0.52
5:H:6:GLU:OE2	5:H:95:CYS:N	2.41	0.52
5:H:98:TRP:CE2	5:H:114:SER:HB3	2.45	0.52
3:3:55:LEU:HB2	3:3:201:VAL:HG23	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:35:ILE:HD11	1:1:61:LEU:HD21	1.91	0.51
2:2:63:LYS:CE	2:2:202:ASN:ND2	2.59	0.51
3:3:64:VAL:HG21	3:3:74:LEU:HB3	1.92	0.51
1:1:110:ALA:HB1	3:3:9:ASP:O	2.09	0.51
3:3:86:MET:O	3:3:89:THR:HG22	2.09	0.51
3:3:53:THR:OG1	3:3:89:THR:OG1	2.18	0.51
3:3:109:PHE:HD2	3:3:201:VAL:HG12	1.76	0.51
3:3:111:PHE:H	3:3:153:SER:HB2	1.74	0.51
5:H:51:ILE:HD13	5:H:71:LYS:HB2	1.91	0.51
3:3:89:THR:HG23	3:3:92:ALA:H	1.74	0.51
3:3:98:TYR:CD2	3:3:211:LEU:HG	2.45	0.51
2:2:94:LEU:C	2:2:94:LEU:HD23	2.35	0.51
3:3:100:GLN:NE2	3:3:212:ARG:HE	2.03	0.51
1:1:4:THR:CG2	4:4:76:PHE:CE2	2.94	0.51
1:1:103:ASN:ND2	3:3:216:ASP:CB	2.73	0.51
2:2:109:VAL:CG2	2:2:156:ALA:HB3	2.40	0.51
2:2:126:VAL:HG21	2:2:129:LEU:CB	2.41	0.51
1:1:35:ILE:CD1	1:1:61:LEU:HD21	2.41	0.50
2:2:83:LEU:HD21	2:2:125:MET:HE1	1.83	0.50
6:L:36:ALA:CB	6:L:93:ALA:HB2	2.41	0.50
1:1:35:ILE:HD11	1:1:61:LEU:CD1	2.34	0.50
2:2:120:CYS:CB	2:2:185:ALA:HB2	2.41	0.50
2:2:129:LEU:HD11	2:2:132:ILE:HG23	1.93	0.50
2:2:135:ARG:HG2	2:2:135:ARG:NH1	2.27	0.50
1:1:87:THR:OG1	1:1:98:LEU:O	2.22	0.50
1:1:91:ASN:HD22	1:1:121:ALA:HA	1.76	0.50
1:1:89:VAL:HG21	1:1:98:LEU:HD13	1.92	0.50
2:2:47:ASN:ND2	3:3:163:SER:O	2.45	0.50
3:3:61:VAL:HG11	5:H:55:GLY:CA	2.38	0.50
1:1:119:TYR:HE1	1:1:126:LEU:HD23	1.77	0.50
2:2:99:ALA:N	2:2:212:GLY:O	2.44	0.49
2:2:112:VAL:HG23	2:2:112:VAL:O	2.12	0.49
2:2:94:LEU:HD21	2:2:210:VAL:CG1	2.41	0.49
5:H:66:ARG:O	5:H:67:LEU:CD2	2.57	0.49
5:H:110:SER:OG	6:L:94:TYR:N	2.45	0.49
6:L:8:GLN:NE2	6:L:8:GLN:N	2.60	0.49
6:L:6:LEU:HD22	6:L:26:GLY:HA2	1.93	0.49
3:3:79:LEU:HD11	3:3:159:ILE:HG21	1.95	0.49
3:3:133:PRO:HB2	3:3:139:ALA:HB2	1.94	0.49
2:2:148:ILE:HG13	2:2:148:ILE:O	2.12	0.49
6:L:35:GLY:C	6:L:54:SER:HB3	2.37	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:67:PHE:CE2	2:2:69:TRP:HB3	2.48	0.48
5:H:18:LEU:O	5:H:81:SER:HA	2.13	0.48
1:1:68:THR:O	1:1:189:ARG:CB	2.61	0.48
1:1:109:LYS:HD3	1:1:110:ALA:HB2	1.94	0.48
2:2:94:LEU:HD23	2:2:101:MET:SD	2.46	0.48
3:3:141:HIS:H	3:3:141:HIS:CD2	2.30	0.48
5:H:12:VAL:HG23	5:H:127:VAL:HG12	1.94	0.48
3:3:117:ALA:HB1	3:3:194:ALA:HB2	1.95	0.48
3:3:151:LEU:C	3:3:151:LEU:CD2	2.86	0.48
1:1:87:THR:O	1:1:168:ILE:HA	2.13	0.48
5:H:50:GLY:CA	5:H:104:TYR:O	2.62	0.48
2:2:71:THR:HA	2:2:188:THR:CG2	2.42	0.48
2:2:98:TYR:CE2	2:2:210:VAL:HG21	2.48	0.48
1:1:35:ILE:CG1	1:1:65:LEU:HD22	2.44	0.48
2:2:108:GLU:OE1	2:2:157:HIS:HD2	1.96	0.47
6:L:32:GLY:C	6:L:69:LYS:HZ1	2.16	0.47
3:3:45:LEU:O	3:3:49:GLU:CD	2.58	0.47
6:L:33:ARG:CG	6:L:33:ARG:NH2	2.72	0.47
2:2:17:THR:HG22	2:2:18:ARG:H	1.79	0.47
6:L:69:LYS:HG3	6:L:74:VAL:HG12	1.97	0.47
2:2:127:PRO:N	2:2:127:PRO:C	2.64	0.47
3:3:71:ASP:OD1	3:3:136:PRO:CG	2.62	0.47
5:H:34:VAL:HA	5:H:96:ALA:O	2.14	0.47
1:1:190:PRO:HD3	2:2:143:PHE:CZ	2.50	0.47
2:2:63:LYS:HZ2	2:2:202:ASN:CG	2.16	0.47
2:2:122:LEU:HA	2:2:147:PHE:HA	1.96	0.47
3:3:61:VAL:HG21	5:H:56:SER:HB2	1.96	0.47
6:L:47:LEU:HD12	6:L:47:LEU:C	2.39	0.47
2:2:98:TYR:CD2	2:2:210:VAL:CG2	2.98	0.46
3:3:102:SER:HB2	3:3:167:TYR:CD1	2.49	0.46
2:2:107:VAL:HG23	2:2:203:ILE:HG13	1.95	0.46
3:3:121:TYR:CD2	3:3:199:LEU:HD22	2.51	0.46
5:H:14:PRO:HA	5:H:85:VAL:HG23	1.96	0.46
5:H:35:ASN:HD22	5:H:50:GLY:HA3	1.78	0.46
5:H:50:GLY:C	5:H:104:TYR:HA	2.40	0.46
5:H:51:ILE:CB	5:H:57:THR:CG2	2.94	0.46
5:H:110:SER:HA	6:L:37:THR:CG2	2.43	0.46
5:H:109:HIS:HD2	6:L:102:PHE:CZ	2.27	0.46
5:H:115:TYR:HB3	6:L:52:TYR:CB	2.46	0.46
6:L:47:LEU:O	6:L:47:LEU:CG	2.63	0.46
1:1:93:ALA:HB1	3:3:215:VAL:HG21	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:38:ARG:HD3	5:H:48:LEU:HD21	1.97	0.46
2:2:87:HIS:HB3	2:2:92:GLY:CA	2.45	0.46
2:2:98:TYR:CE2	2:2:210:VAL:CG2	2.99	0.46
1:1:91:ASN:ND2	1:1:121:ALA:HA	2.30	0.46
2:2:62:PHE:CG	2:2:87:HIS:CE1	3.03	0.46
6:L:6:LEU:HD22	6:L:26:GLY:CA	2.46	0.46
6:L:8:GLN:CD	6:L:8:GLN:H	2.24	0.46
2:2:96:ASP:O	2:2:214:PHE:HD2	1.98	0.46
5:H:60:ASN:O	5:H:63:LEU:C	2.59	0.46
5:H:37:VAL:HG23	5:H:47:CYS:HA	1.97	0.45
2:2:98:TYR:CD2	2:2:210:VAL:CB	2.98	0.45
1:1:110:ALA:HB1	1:1:111:PRO:HD2	1.98	0.45
2:2:213:GLU:OE1	2:2:213:GLU:N	2.49	0.45
1:1:57:PRO:HG2	1:1:60:THR:OG1	2.16	0.45
3:3:11:TYR:O	3:3:11:TYR:CD1	2.70	0.45
2:2:69:TRP:CZ3	2:2:121:LEU:HD22	2.51	0.45
3:3:72:ARG:HH22	6:L:34:TYR:HD2	1.64	0.45
5:H:98:TRP:HZ2	5:H:111:SER:HB3	1.80	0.45
1:1:125:VAL:O	2:2:165:VAL:HG21	2.06	0.45
6:L:32:GLY:HA2	6:L:69:LYS:NZ	2.31	0.45
1:1:47:GLN:HG3	1:1:170:ALA:O	2.17	0.45
1:1:87:THR:CB	1:1:105:THR:HG22	2.45	0.45
5:H:70:THR:HG22	5:H:79:SER:HB2	1.98	0.45
2:2:29:SER:OG	2:2:30:VAL:N	2.50	0.45
2:2:117:ASN:ND2	2:2:189:VAL:O	2.50	0.44
3:3:65:THR:OG1	5:H:54:SER:OG	2.18	0.44
2:2:126:VAL:CG2	2:2:129:LEU:CB	2.81	0.44
3:3:52:PRO:HB2	3:3:202:LEU:HD12	1.99	0.44
3:3:120:ARG:HG2	3:3:189:ILE:HD11	1.99	0.44
2:2:83:LEU:HD12	2:2:83:LEU:HA	1.76	0.44
2:2:126:VAL:HG12	2:2:143:PHE:CD2	2.52	0.44
5:H:39:GLN:CD	6:L:41:GLN:HE22	2.26	0.44
6:L:8:GLN:HG3	6:L:104:SER:HB3	1.98	0.44
5:H:75:LYS:HB3	5:H:75:LYS:HE3	1.75	0.44
5:H:98:TRP:CD2	5:H:114:SER:HB3	2.53	0.44
1:1:70:THR:HG22	1:1:189:ARG:NE	2.27	0.44
2:2:91:TYR:O	2:2:91:TYR:CD1	2.70	0.44
5:H:51:ILE:HB	5:H:57:THR:HG21	2.00	0.44
5:H:111:SER:O	6:L:53:GLY:HA2	2.18	0.44
3:3:117:ALA:CB	3:3:194:ALA:HB2	2.48	0.44
3:3:121:TYR:CE2	3:3:199:LEU:HD22	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:124:ARG:HB3	2:2:165:VAL:O	2.18	0.43
5:H:51:ILE:CB	5:H:57:THR:HG21	2.47	0.43
6:L:27:SER:HB3	6:L:30:ASN:ND2	2.33	0.43
6:L:27:SER:HB3	6:L:30:ASN:HD21	1.82	0.43
2:2:87:HIS:HB3	2:2:92:GLY:HA2	2.00	0.43
3:3:68:THR:HG22	3:3:195:ASP:HB2	2.00	0.43
5:H:4:LEU:HD22	5:H:24:VAL:HG22	2.01	0.43
3:3:61:VAL:CG1	5:H:55:GLY:HA3	2.43	0.43
4:4:18:THR:H	4:4:18:THR:HG23	1.41	0.43
5:H:8:GLY:HA3	5:H:20:LEU:HD23	1.99	0.43
6:L:35:GLY:CA	6:L:54:SER:HB3	2.48	0.43
1:1:40:VAL:O	1:1:40:VAL:HG13	2.18	0.43
1:1:100:ASN:HD22	3:3:216:ASP:HB3	1.84	0.43
3:3:2:ILE:H	3:3:2:ILE:HG12	1.52	0.43
2:2:62:PHE:CG	2:2:87:HIS:HE1	2.36	0.43
2:2:17:THR:HG22	2:2:18:ARG:N	2.34	0.43
2:2:43:VAL:HG22	2:2:102:ARG:NH2	2.34	0.43
3:3:72:ARG:NH2	6:L:34:TYR:HD2	2.14	0.43
4:4:76:PHE:CD2	4:4:76:PHE:O	2.71	0.43
5:H:112:ASP:HA	6:L:53:GLY:HA2	2.00	0.43
1:1:97:ALA:O	1:1:103:ASN:ND2	2.51	0.43
2:2:163:VAL:HG22	2:2:177:TRP:CZ2	2.54	0.43
6:L:81:LEU:HD23	6:L:81:LEU:HA	1.79	0.43
2:2:58:ALA:CA	2:2:90:VAL:HG12	2.49	0.43
1:1:70:THR:HG23	1:1:187:CYS:O	2.18	0.42
1:1:196:PRO:HB2	1:1:198:GLU:O	2.18	0.42
3:3:55:LEU:O	3:3:62:PRO:HA	2.19	0.42
5:H:34:VAL:HG11	5:H:78:VAL:HG11	2.01	0.42
5:H:115:TYR:HE2	5:H:117:ASP:HB3	1.83	0.42
2:2:96:ASP:O	2:2:214:PHE:CD2	2.72	0.42
5:H:58:GLY:CA	5:H:105:ASP:OD1	2.32	0.42
6:L:8:GLN:N	6:L:8:GLN:CD	2.77	0.42
5:H:98:TRP:NE1	5:H:116:LEU:CD1	2.82	0.42
6:L:8:GLN:HG3	6:L:104:SER:H	1.84	0.42
5:H:65:SER:O	5:H:65:SER:OG	2.34	0.42
2:2:134:LYS:HA	2:2:137:LEU:HD12	2.02	0.42
3:3:56:HIS:O	3:3:85:HIS:HA	2.20	0.42
3:3:122:MET:HG3	3:3:187:PHE:HB2	2.02	0.42
1:1:42:VAL:HG21	1:1:178:TYR:CD2	2.55	0.42
2:2:66:LEU:HD23	2:2:181:VAL:HG11	2.02	0.42
3:3:212:ARG:CD	3:3:213:LEU:HG	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:L:34:TYR:CD1	6:L:34:TYR:N	2.87	0.42
1:1:120:THR:OG1	1:1:120:THR:O	2.31	0.41
1:1:35:ILE:CD1	1:1:61:LEU:HD11	2.37	0.41
1:1:93:ALA:HB2	3:3:215:VAL:HG21	2.01	0.41
1:1:122:PRO:N	1:1:122:PRO:C	2.66	0.41
3:3:63:TYR:CE1	5:H:54:SER:HB2	2.55	0.41
5:H:27:PHE:CD2	5:H:27:PHE:C	2.98	0.41
6:L:29:SER:HB3	6:L:96:ILE:HG21	1.94	0.41
6:L:29:SER:CB	6:L:96:ILE:CG2	2.88	0.41
3:3:135:THR:OG1	3:3:136:PRO:HD2	2.21	0.41
3:3:208:ASP:OD1	3:3:208:ASP:N	2.52	0.41
1:1:125:VAL:O	2:2:165:VAL:CB	2.67	0.41
2:2:77:ARG:NH2	2:2:79:TYR:HB2	2.32	0.41
1:1:91:ASN:OD1	1:1:91:ASN:C	2.64	0.41
3:3:136:PRO:HB3	3:3:189:ILE:HG22	2.03	0.41
2:2:99:ALA:HB2	2:2:214:PHE:CE1	2.56	0.40
2:2:148:ILE:CG2	2:2:156:ALA:HB2	2.51	0.40
3:3:163:SER:HB2	3:3:181:GLN:HE22	1.86	0.40
6:L:16:LEU:HG	6:L:110:LEU:HD21	2.02	0.40
2:2:162:PHE:CZ	2:2:169:ASP:HB2	2.56	0.40
5:H:86:THR:OG1	5:H:88:GLU:HG3	2.22	0.40
2:2:84:PRO:HG2	2:2:171:TYR:CD2	2.57	0.40
3:3:66:THR:HG23	3:3:188:GLN:OE1	2.21	0.40
6:L:64:ARG:NH2	6:L:84:GLU:OE2	2.55	0.40
6:L:67:GLY:HA2	6:L:75:THR:O	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	178/213 (84%)	164 (92%)	13 (7%)	1 (1%)	21	52

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	2	203/218 (93%)	188 (93%)	15 (7%)	0	100	100
3	3	217/220 (99%)	212 (98%)	5 (2%)	0	100	100
4	4	40/85 (47%)	37 (92%)	3 (8%)	0	100	100
5	H	127/129 (98%)	122 (96%)	5 (4%)	0	100	100
6	L	106/122 (87%)	98 (92%)	8 (8%)	0	100	100
All	All	871/987 (88%)	821 (94%)	49 (6%)	1 (0%)	49	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	29	HIS

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	153/178 (86%)	146 (95%)	7 (5%)	24	47
2	2	177/190 (93%)	169 (96%)	8 (4%)	24	48
3	3	174/175 (99%)	169 (97%)	5 (3%)	37	56
4	4	35/67 (52%)	32 (91%)	3 (9%)	10	33
5	H	113/113 (100%)	110 (97%)	3 (3%)	39	58
6	L	89/99 (90%)	85 (96%)	4 (4%)	24	48
All	All	741/822 (90%)	711 (96%)	30 (4%)	29	51

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	6	GLU
1	1	14	THR
1	1	22	THR
1	1	23	GLN
1	1	38	ARG

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Mol	Chain	Res	Type
1	1	98	LEU
1	1	201	HIS
2	2	23	THR
2	2	25	THR
2	2	26	THR
2	2	69	TRP
2	2	132	ILE
2	2	147	PHE
2	2	165	VAL
2	2	184	VAL
3	3	2	ILE
3	3	16	THR
3	3	21	THR
3	3	68	THR
3	3	212	ARG
4	4	35	ASP
4	4	37	GLN
4	4	79	LEU
5	H	24	VAL
5	H	98	TRP
5	H	109	HIS
6	L	8	GLN
6	L	31	ILE
6	L	33	ARG
6	L	34	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	29	HIS
1	1	85	ASN
1	1	100	ASN
1	1	108	HIS
1	1	123	HIS
1	1	195	HIS
1	1	201	HIS
2	2	27	GLN
2	2	87	HIS
2	2	117	ASN
2	2	149	ASN
2	2	166	ASN
2	2	190	ASN

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Mol	Chain	Res	Type
2	2	202	ASN
3	3	85	HIS
3	3	88	ASN
3	3	100	GLN
3	3	108	HIS
3	3	141	HIS
3	3	179	ASN
3	3	181	GLN
3	3	188	GLN
4	4	17	ASN
4	4	23	ASN
4	4	24	ASN
5	H	35	ASN
5	H	39	GLN
5	H	109	HIS
6	L	8	GLN
6	L	41	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 ⓘ

There are no chain breaks in this entry.

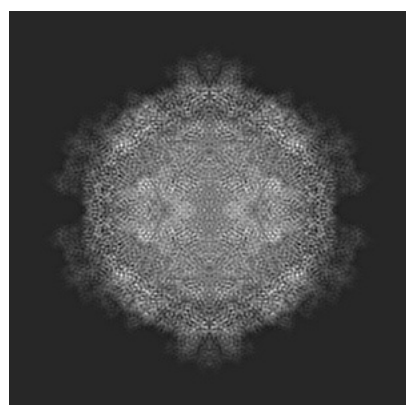
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31223. These allow visual inspection of the internal detail of the map and identification of artifacts.

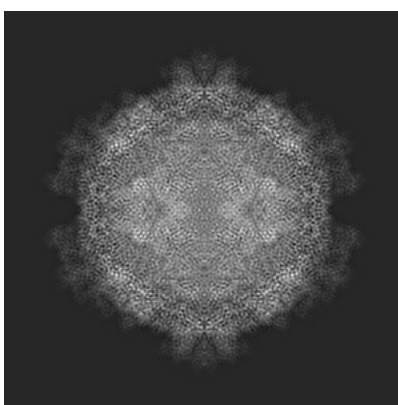
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

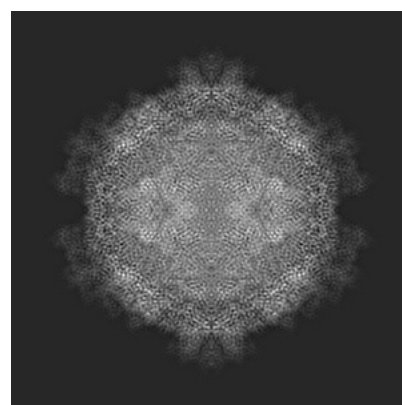
6.1.1 Primary map



X



Y

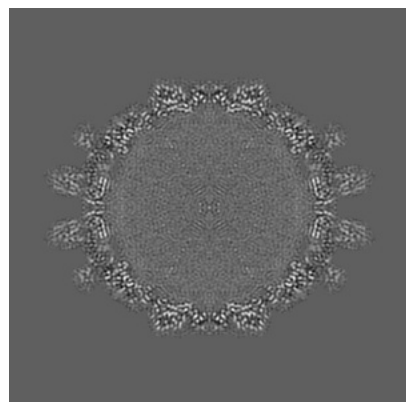


Z

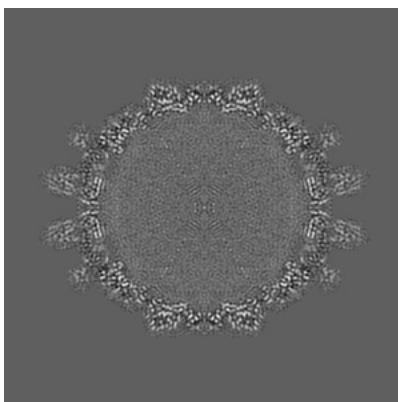
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

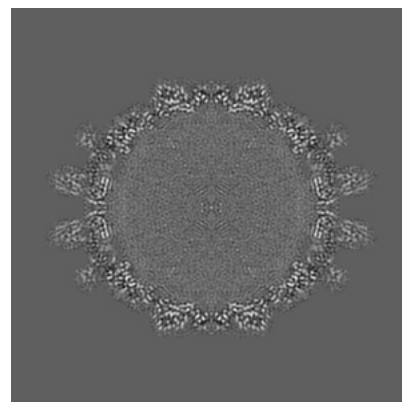
6.2.1 Primary map



X Index: 240



Y Index: 240

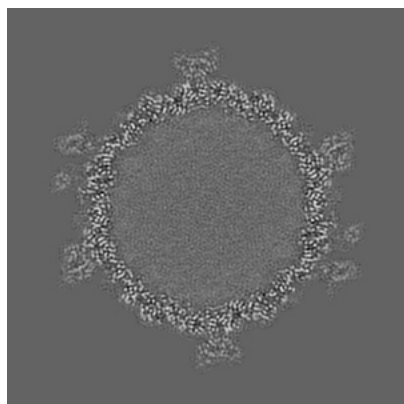


Z Index: 240

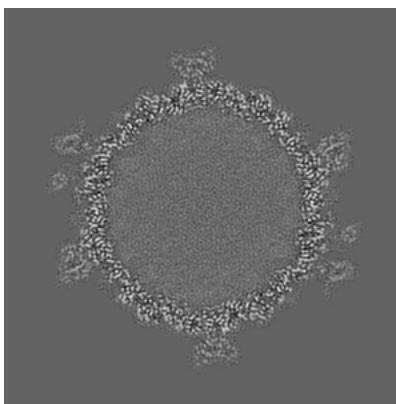
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

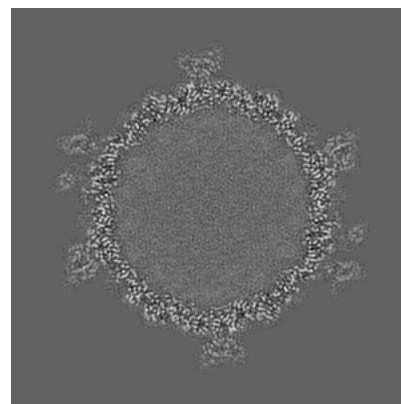
6.3.1 Primary map



X Index: 209



Y Index: 209

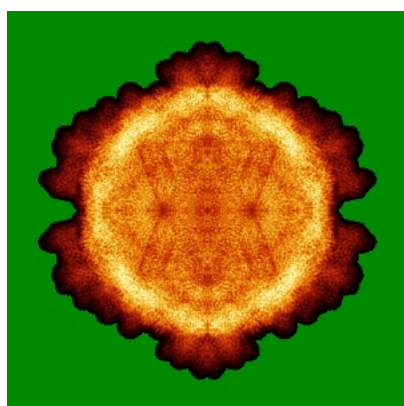


Z Index: 209

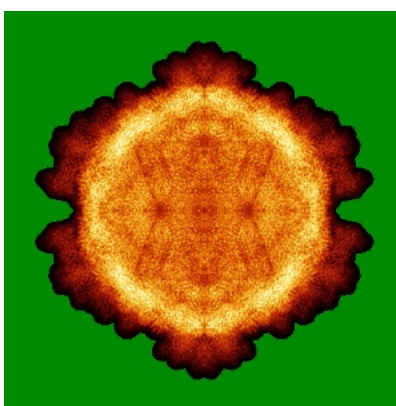
The images above show the largest variance slices of the map in three orthogonal directions.

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

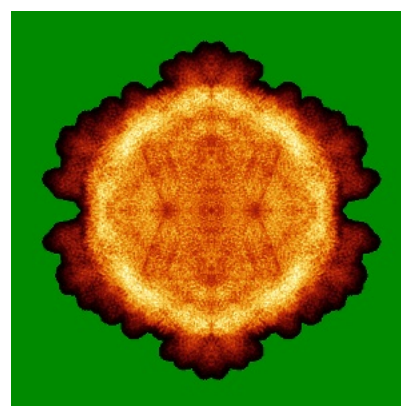
6.4.1 Primary map



X



Y

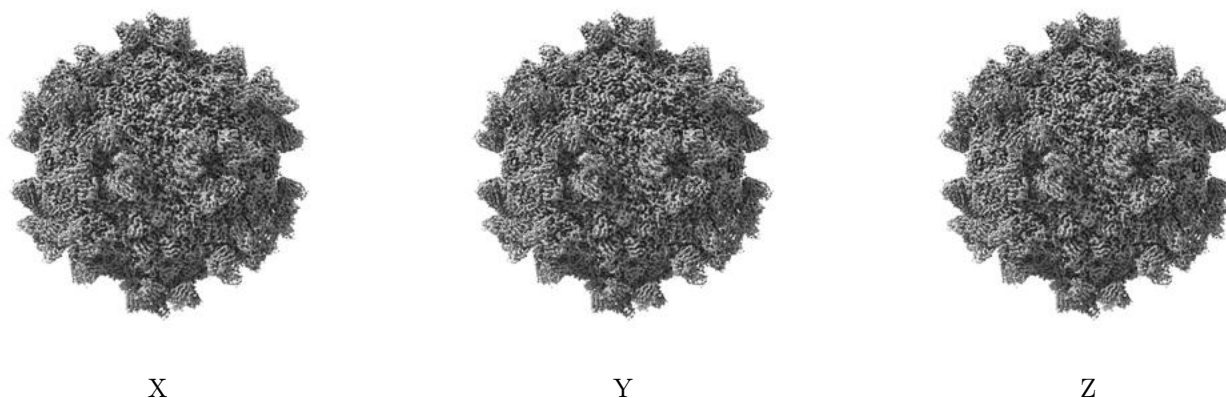


Z

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

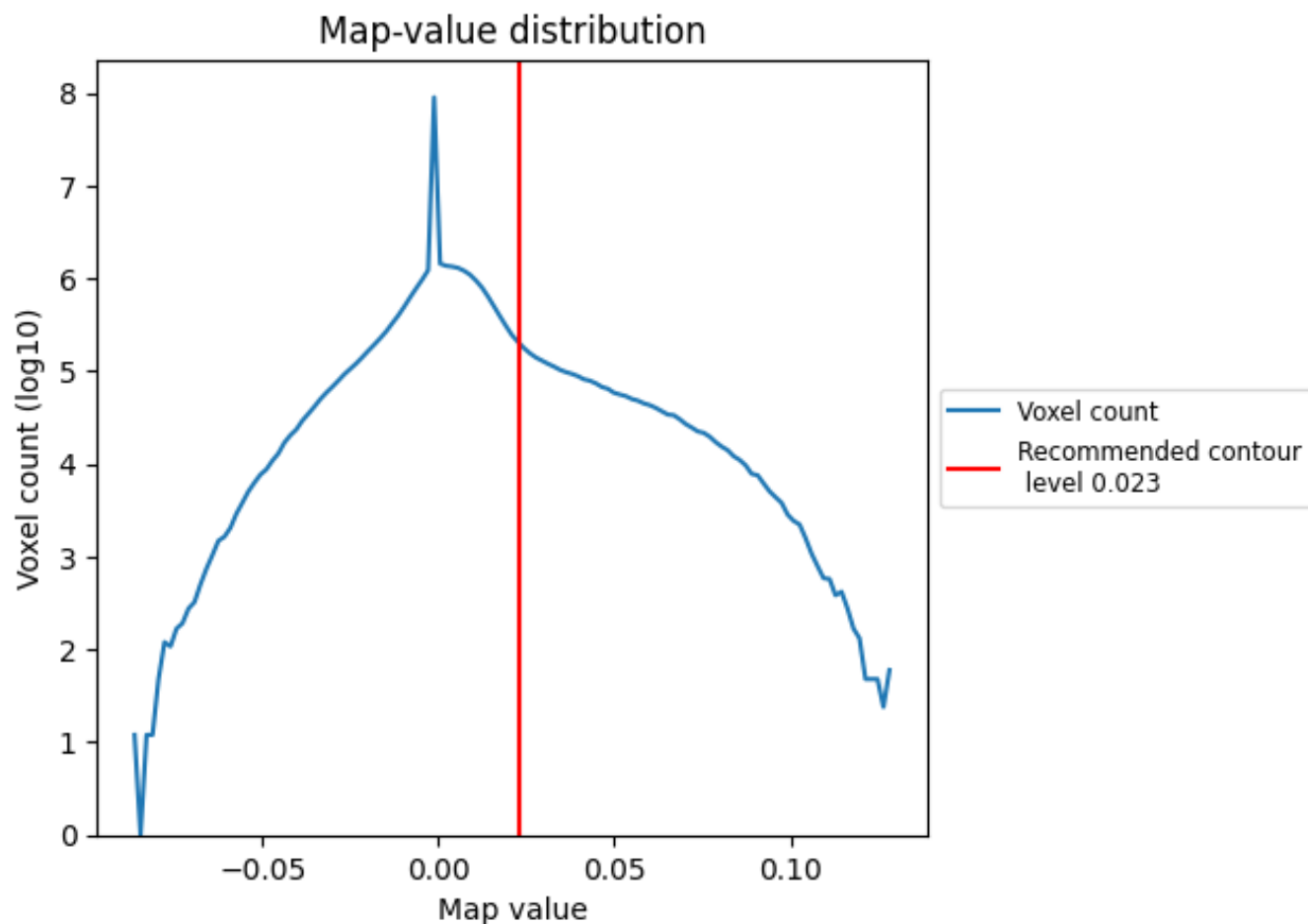
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

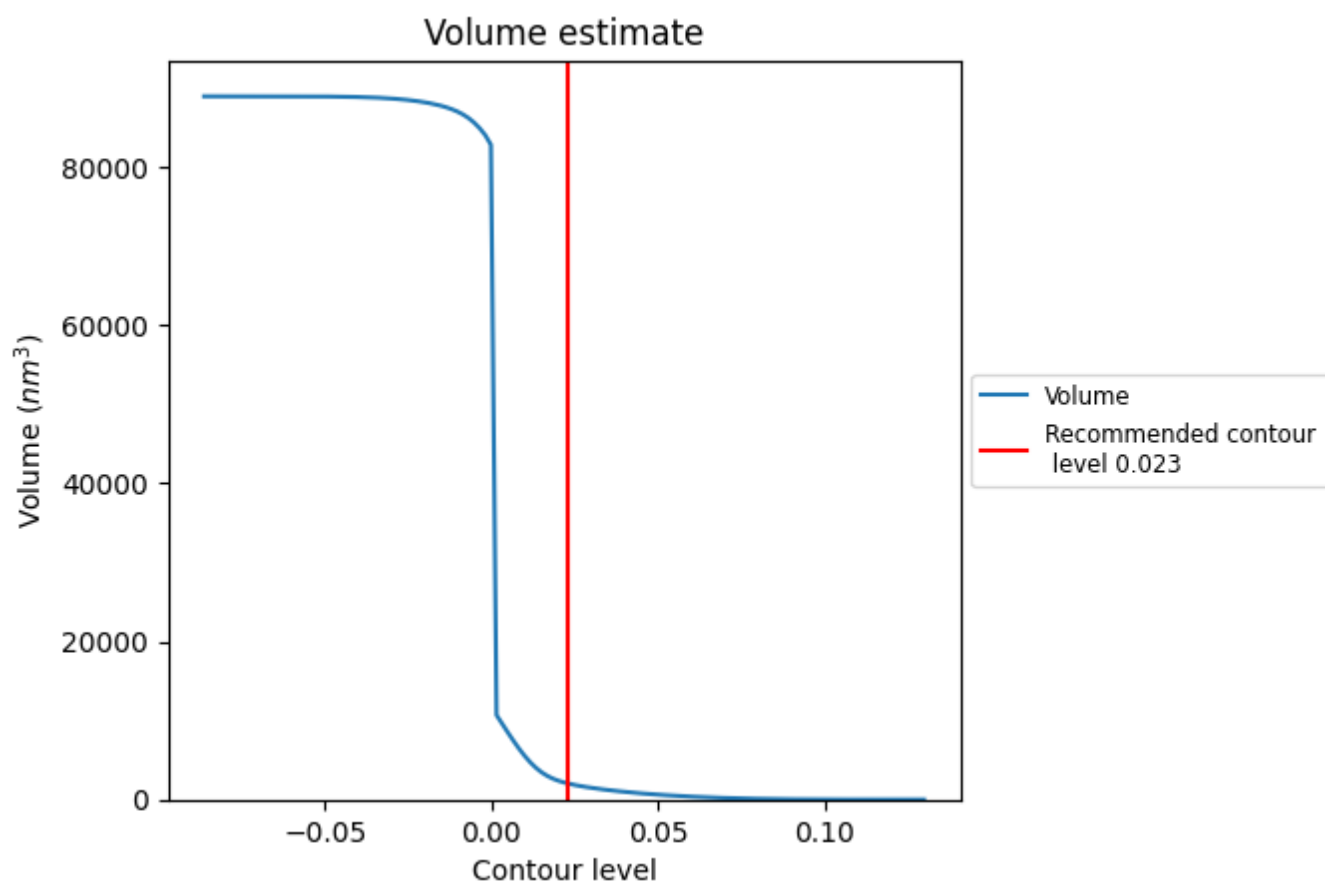
This section contains the results of statistical analysis of the map.

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.

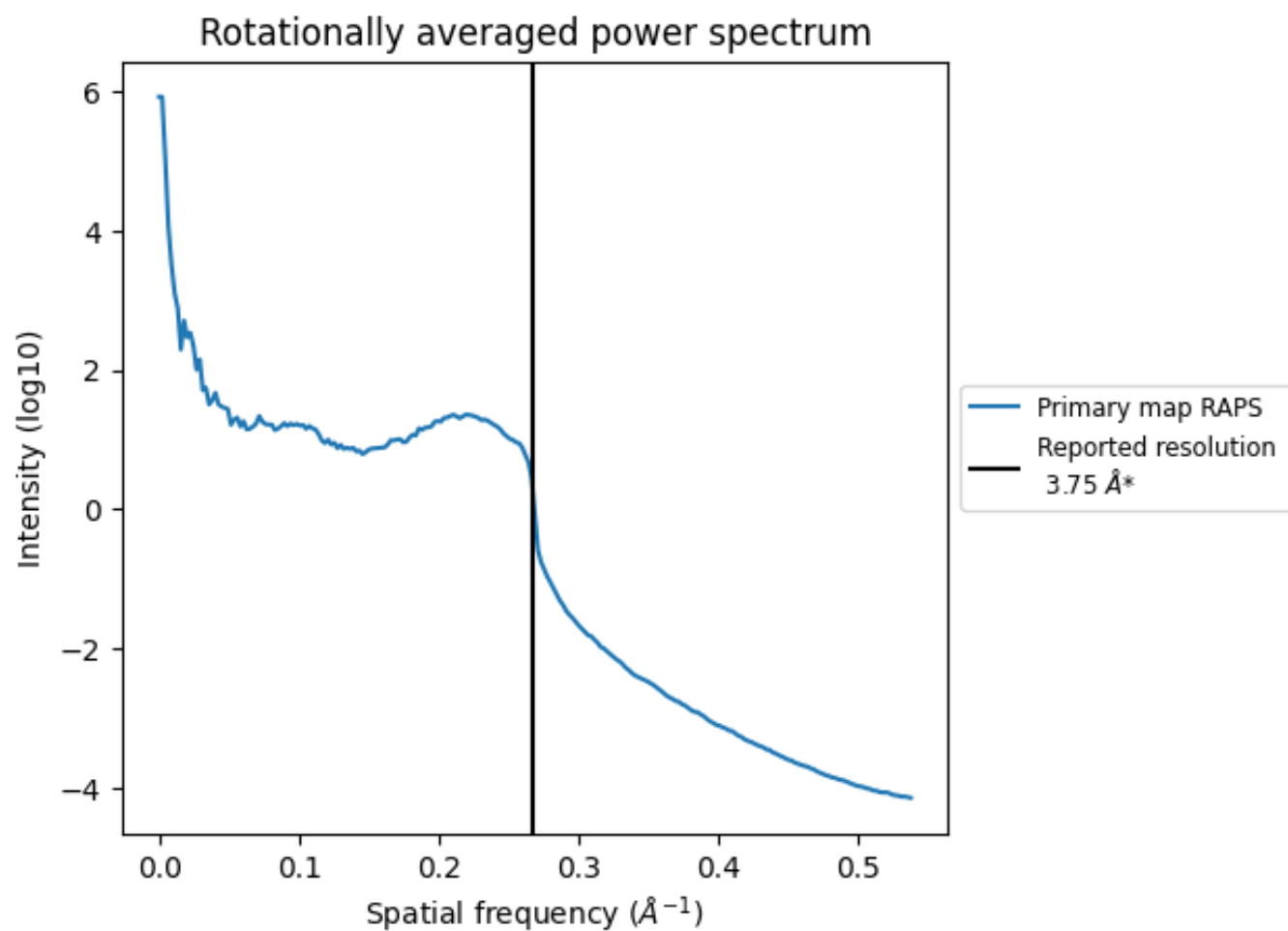
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2032 nm³; this corresponds to an approximate mass of 1836 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 0.267 Å⁻¹

8 Fourier-Shell correlation ⓘ

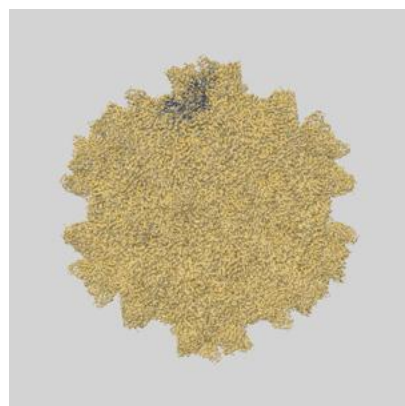
This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

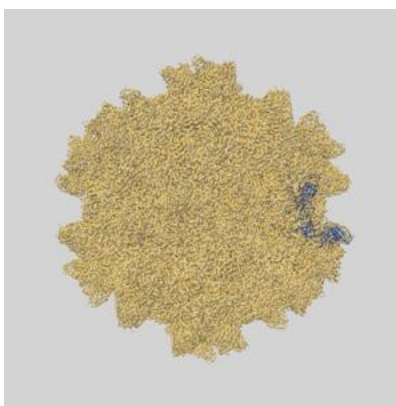
This section contains information regarding the fit between EMDB map EMD-31223 and PDB model 7EO0. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlays

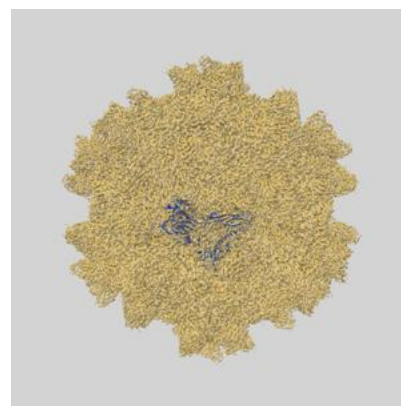
9.1.1 Map-model overlay [i](#)



X

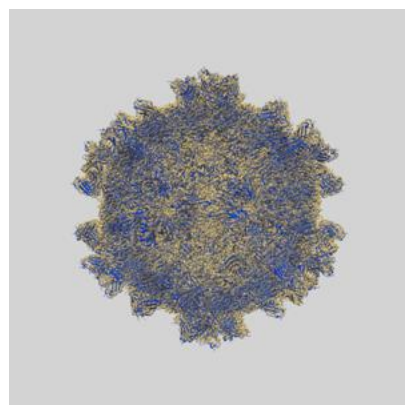


Y

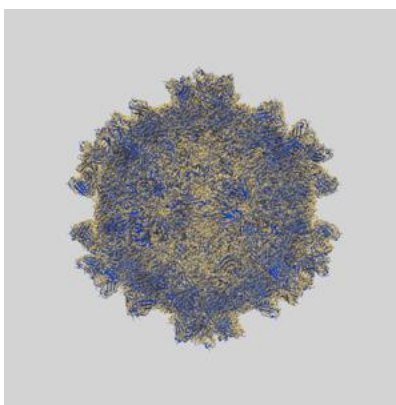


Z

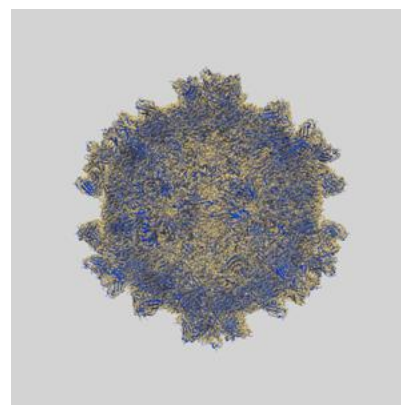
9.1.2 Map-model assembly overlay [i](#)



X



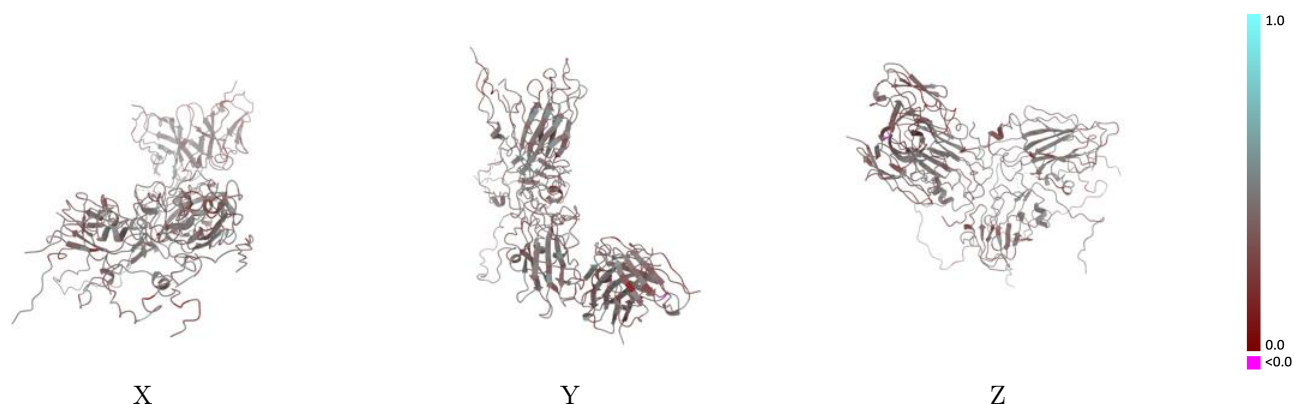
Y



Z

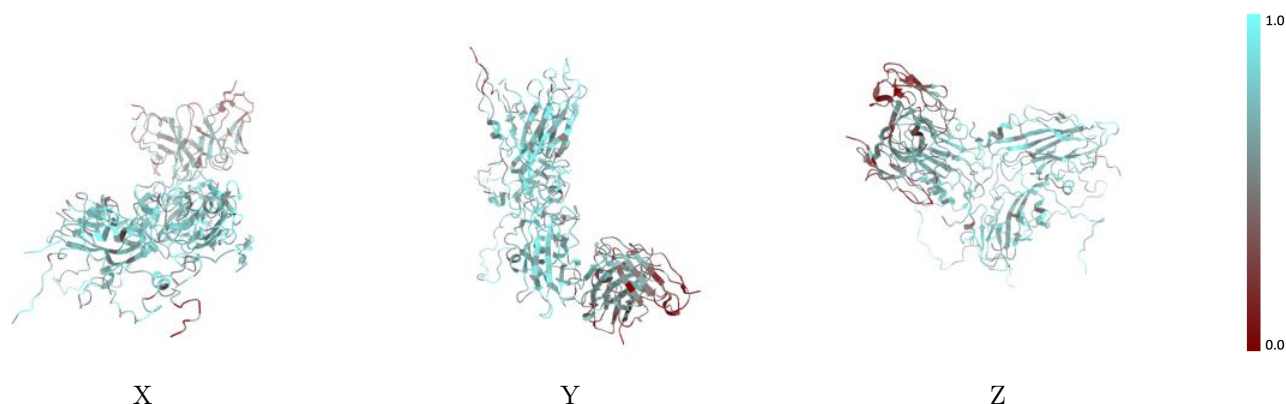
The images above show the 3D surface view of the map at the recommended contour level 0.023 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



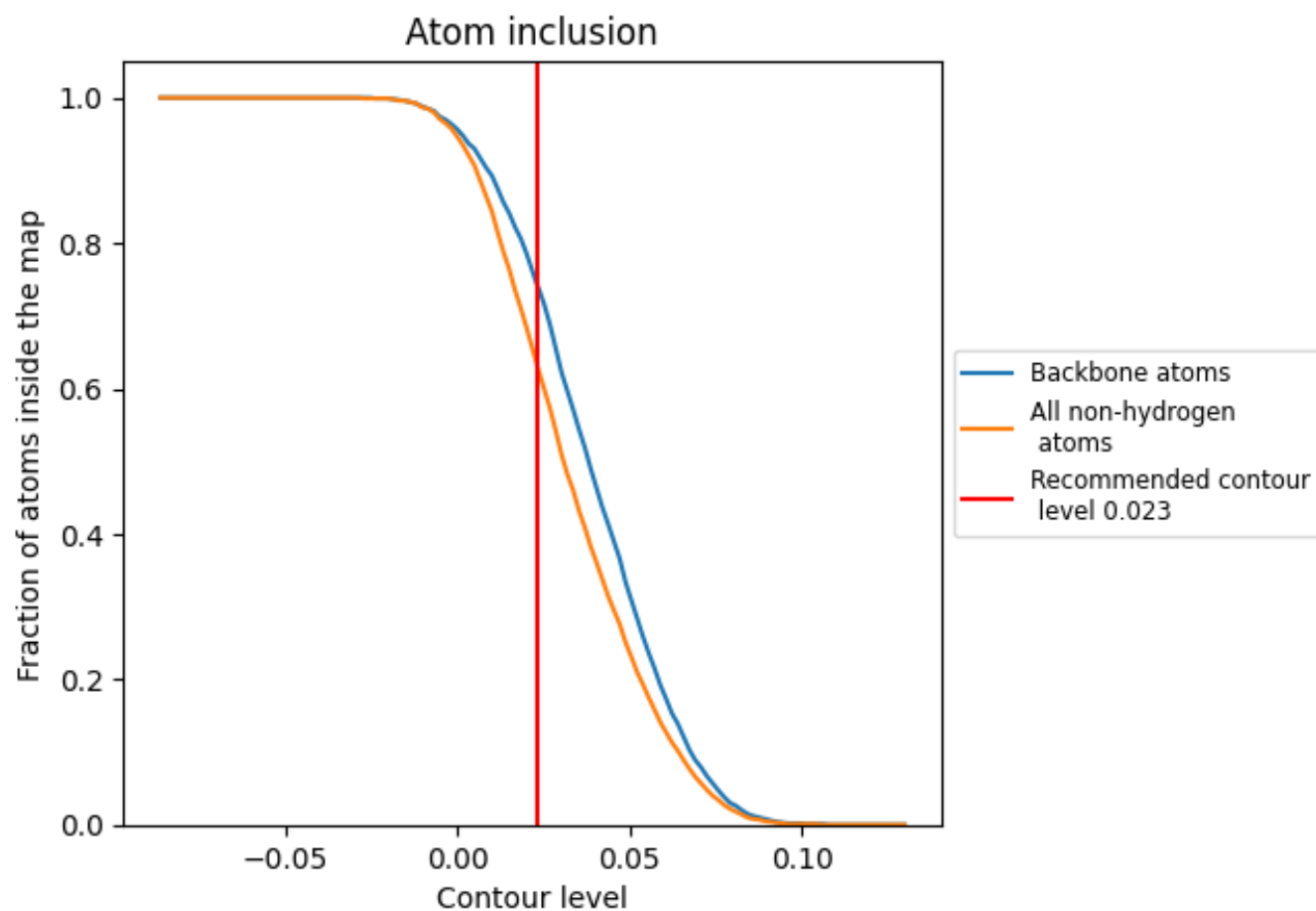
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.023).

9.4 Atom inclusion [i](#)



At the recommended contour level, 74% of all backbone atoms, 63% 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 (0.023) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6340	0.4040
1	0.6970	0.4060
2	0.7000	0.4170
3	0.7070	0.4060
4	0.5890	0.4120
H	0.4880	0.3970
L	0.4280	0.3780

1.0

0.0

<0.0