



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 04:38 PM UTC

PDB ID : 8K28 / pdb_00008k28
EMDB ID : EMD-36831
Title : ICP1 Csy-dsDNA complex (form 1)
Authors : Zhang, L.X.; Feng, Y.
Deposited on : 2023-07-12
Resolution : 3.54 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

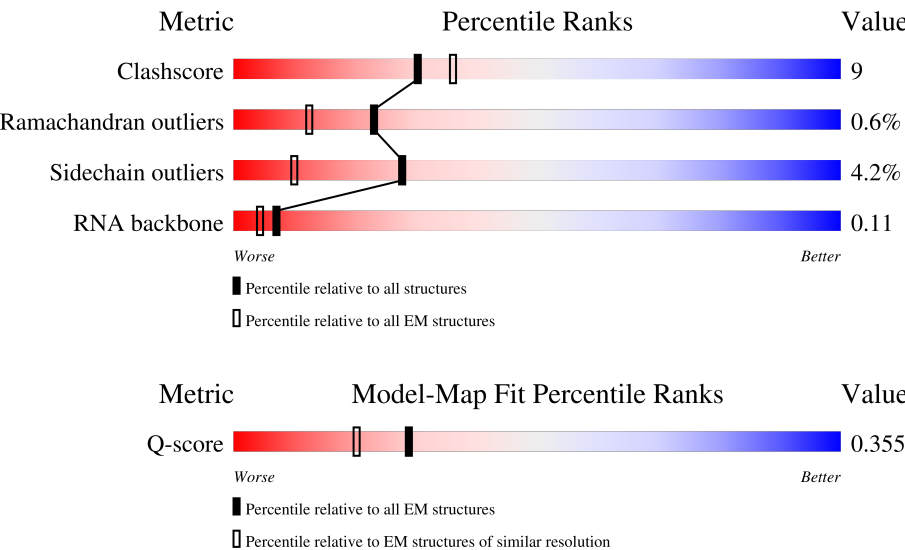
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.54 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.



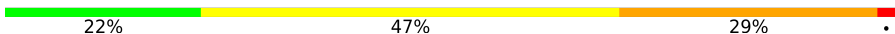
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	12891 (3.04 - 4.04)

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

Mol	Chain	Length	Quality of chain
1	A	179	
2	B	248	
3	C	306	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	D	306	 73%23%..
3	E	306	 71%26%..
3	F	306	 66%29%... .
3	G	306	 75%22%..
3	H	306	 5%78%19%..
4	I	168	 5%77%20%..
5	Q	33	 12%91%9%
6	R	27	 11%89%11%
7	P	59	 22%47%29%. .

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 20740 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 Csy1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	173	Total	C	N	O	S	0	0
			1345	851	226	259	9		

- Molecule 2 is a protein called Csy2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	246	Total	C	N	O	S	0	0
			1898	1211	319	355	13		

- Molecule 3 is a protein called Csy3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	302	Total	C	N	O	S	0	0
			2291	1441	389	454	7		
3	D	302	Total	C	N	O	S	0	0
			2289	1442	388	452	7		
3	E	302	Total	C	N	O	S	0	0
			2297	1448	390	452	7		
3	F	302	Total	C	N	O	S	0	0
			2297	1448	390	452	7		
3	G	302	Total	C	N	O	S	0	0
			2285	1438	386	454	7		
3	H	302	Total	C	N	O	S	0	0
			2277	1435	384	452	6		

- Molecule 4 is a protein called Csy4.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	I	166	Total	C	N	O	S	0	0
			1291	817	221	246	7		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	52	ILE	VAL	conflict	UNP F1D5V5

- Molecule 5 is a DNA chain called DNA (33-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
5	Q	33	Total	C	N	O	P	0	0
			677	322	125	197	33		

- Molecule 6 is a DNA chain called DNA (27-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	R	27	Total	C	N	O	P	0	0
			556	266	100	163	27		

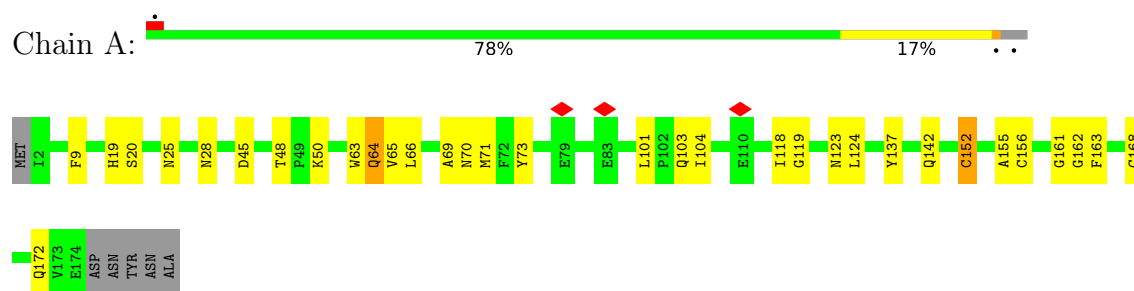
- Molecule 7 is a RNA chain called RNA (59-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
7	P	59	Total	C	N	O	P	0	0
			1237	555	205	418	59		

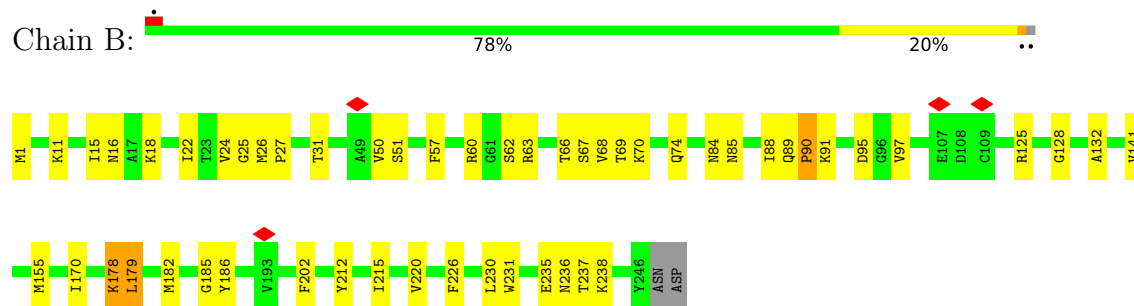
3 Residue-property plots

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.

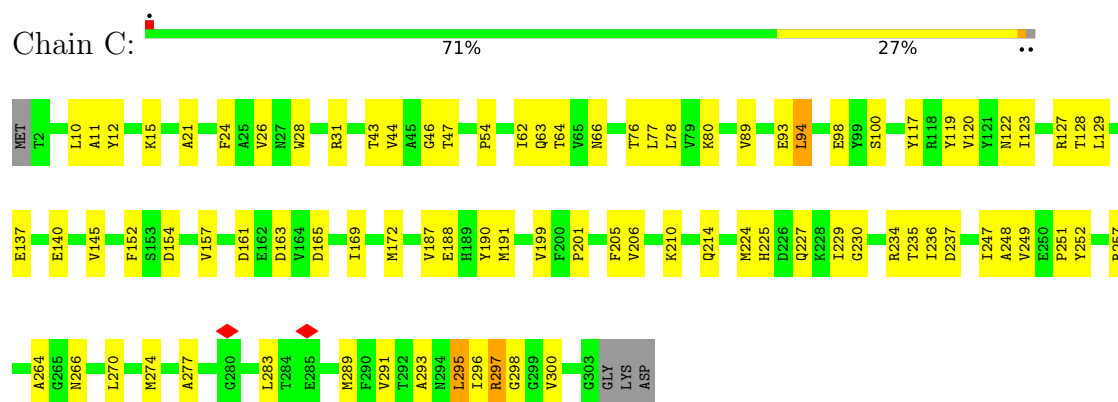
• Molecule 1: Csy1



• Molecule 2: Csy2



• Molecule 3: Csy3



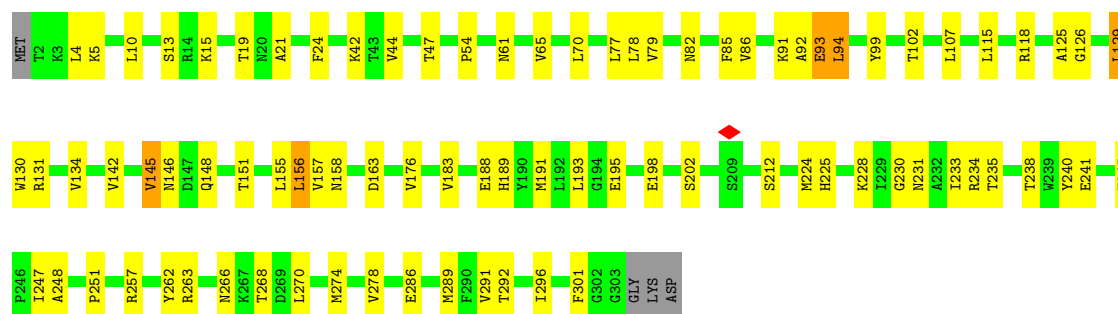
• Molecule 3: Csy3

Chain D:  73% 23% ..



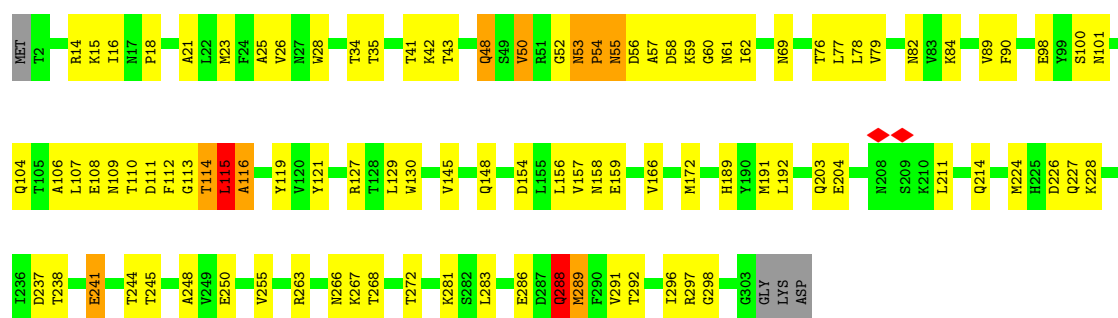
• Molecule 3: Csy3

Chain E:  71% 26% ..



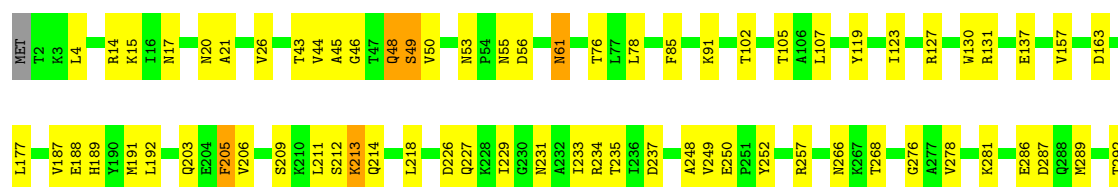
• Molecule 3: Csy3

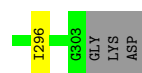
Chain F:  66% 29% ..



• Molecule 3: Csy3

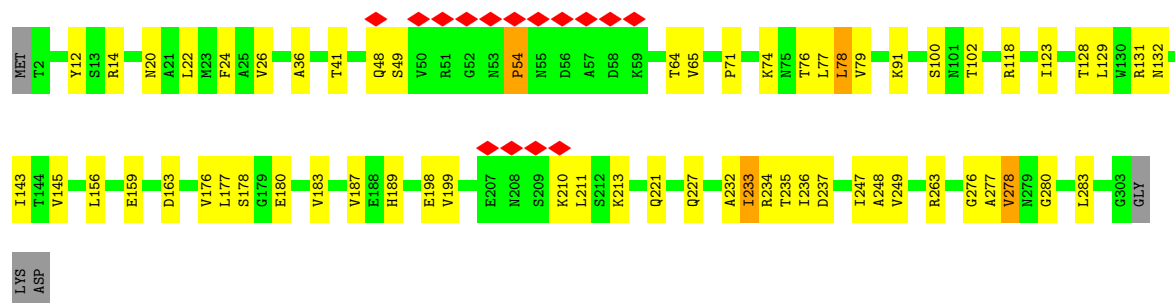
Chain G:  75% 22% ..





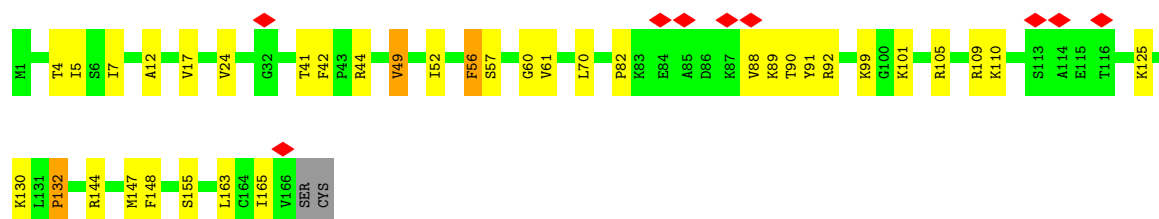
- Molecule 3: Csy3

Chain H: 5% 78% 19% ..



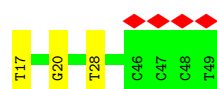
- Molecule 4: Csy4

Chain I: 5% 77% 20% ..



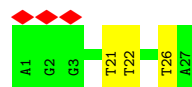
- Molecule 5: DNA (33-MER)

Chain Q: 12% 91% 9%



- Molecule 6: DNA (27-MER)

Chain R: 11% 89% 11%



- Molecule 7: RNA (59-MER)

Chain P: 22% 47% 29%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	90717	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.751	Depositor
Minimum map value	-1.678	Depositor
Average map value	0.007	Depositor
Map value standard deviation	0.083	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	264.0, 264.0, 264.0	wwPDB
Map dimensions	240, 240, 240	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	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	A	0.35	0/1374	0.69	3/1858 (0.2%)
2	B	0.25	0/1934	0.66	2/2615 (0.1%)
3	C	0.40	0/2330	0.66	4/3167 (0.1%)
3	D	0.56	3/2328 (0.1%)	0.83	8/3166 (0.3%)
3	E	0.50	2/2336 (0.1%)	0.68	4/3174 (0.1%)
3	F	0.67	7/2336 (0.3%)	0.93	19/3174 (0.6%)
3	G	0.55	1/2324 (0.0%)	0.75	5/3163 (0.2%)
3	H	0.42	0/2314	0.71	4/3145 (0.1%)
4	I	0.25	0/1309	0.69	0/1759
5	Q	0.22	0/759	0.45	0/1169
6	R	0.33	0/623	0.58	0/960
7	P	0.28	0/1378	0.53	1/2140 (0.0%)
All	All	0.45	13/21345 (0.1%)	0.72	50/29490 (0.2%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	125	ALA	CA-C	-7.45	1.42	1.52
3	D	93	GLU	CA-C	-7.01	1.43	1.52
3	D	6	ALA	CA-C	-6.95	1.45	1.53
3	E	92	ALA	CA-C	-6.66	1.44	1.53
3	D	99	TYR	CA-C	-6.65	1.43	1.52
3	G	205	PHE	CA-C	-6.47	1.44	1.52
3	F	116	ALA	CA-C	-6.19	1.44	1.52
3	F	115	LEU	CA-C	-5.82	1.45	1.52
3	F	48	GLN	CA-C	-5.64	1.45	1.52
3	F	61	ASN	CA-C	-5.53	1.46	1.52
3	F	61	ASN	N-CA	-5.53	1.39	1.46
3	F	113	GLY	CA-C	-5.41	1.46	1.52
3	F	62	ILE	CA-CB	-5.05	1.47	1.54

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	57	ALA	N-CA-C	-14.01	96.77	114.56
3	F	112	PHE	N-CA-C	-13.71	95.87	112.59
3	F	58	ASP	N-CA-C	-11.40	98.83	111.14
3	D	100	SER	N-CA-C	-10.23	101.36	114.04
3	D	6	ALA	CA-C-N	-8.89	110.87	119.76
3	D	6	ALA	C-N-CA	-8.89	110.87	119.76
3	C	94	LEU	N-CA-C	-8.10	96.35	109.96
3	F	108	GLU	N-CA-C	-7.62	104.00	113.38
3	F	61	ASN	N-CA-C	-7.54	96.98	108.96
3	D	37	VAL	N-CA-C	-7.38	99.15	108.89
3	F	54	PRO	N-CA-CB	7.36	110.97	103.25
3	C	54	PRO	N-CA-CB	7.18	110.14	103.46
3	F	288	GLN	N-CA-C	-7.18	103.53	111.36
3	D	54	PRO	N-CA-CB	6.95	110.18	103.51
3	H	54	PRO	N-CA-CB	6.91	110.50	103.25
3	D	99	TYR	N-CA-C	-6.76	96.41	110.80
3	F	113	GLY	N-CA-C	-6.68	104.22	112.77
3	G	209	SER	N-CA-C	6.62	118.61	109.54
1	A	155	ALA	N-CA-C	-6.58	104.04	111.14
3	E	54	PRO	N-CA-CB	6.50	109.52	103.41
3	F	50	VAL	N-CA-C	-6.30	96.24	109.34
3	F	238	THR	N-CA-C	-6.16	103.18	111.87
3	G	46	GLY	N-CA-C	6.11	119.73	111.72
3	F	245	THR	CA-C-N	-6.11	113.49	119.90
3	F	245	THR	C-N-CA	-6.11	113.49	119.90
3	G	211	LEU	N-CA-C	-6.09	100.78	108.45
3	F	53	ASN	N-CA-C	5.95	122.96	109.81
3	G	49	SER	N-CA-C	5.91	118.33	110.53
3	H	276	GLY	N-CA-C	-5.77	105.38	112.77
3	G	48	GLN	N-CA-C	-5.77	106.05	112.57
3	E	125	ALA	N-CA-C	-5.75	98.55	110.80
7	P	34	U	C2'-C3'-O3'	5.68	122.22	113.70
3	F	55	ASN	N-CA-C	-5.64	104.98	112.94
3	D	36	ALA	N-CA-C	5.47	118.28	110.24
2	B	90	PRO	CA-C-N	5.45	127.07	122.28
2	B	90	PRO	C-N-CA	5.45	127.07	122.28
3	F	106	ALA	N-CA-C	-5.29	106.34	112.89
1	A	65	VAL	N-CA-C	-5.27	105.18	110.30
3	F	60	GLY	CA-C-N	-5.21	115.27	122.30
3	F	60	GLY	C-N-CA	-5.21	115.27	122.30
3	F	113	GLY	CA-C-N	-5.17	112.84	122.09
3	F	113	GLY	C-N-CA	-5.17	112.84	122.09
3	D	7	PRO	CA-C-O	-5.14	115.14	121.31

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	64	GLN	N-CA-C	-5.13	106.28	112.54
3	E	94	LEU	N-CA-C	-5.13	102.47	110.42
3	H	145	VAL	CA-C-N	5.09	129.63	122.46
3	H	145	VAL	C-N-CA	5.09	129.63	122.46
3	C	145	VAL	CA-C-N	5.08	129.62	122.46
3	C	145	VAL	C-N-CA	5.08	129.62	122.46
3	E	158	ASN	N-CA-C	5.03	117.16	111.02

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1345	0	1274	19	0
2	B	1898	0	1912	39	0
3	C	2291	0	2199	57	0
3	D	2289	0	2202	47	0
3	E	2297	0	2224	53	0
3	F	2297	0	2224	62	0
3	G	2285	0	2184	36	0
3	H	2277	0	2198	35	0
4	I	1291	0	1325	23	0
5	Q	677	0	372	3	0
6	R	556	0	307	4	0
7	P	1237	0	630	38	0
All	All	20740	0	19051	353	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 9.

All (353) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:GLN:O	1:A:162:GLY:HA3	1.79	0.81

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:26:VAL:O	3:G:76:THR:HB	1.88	0.74
3:F:78:LEU:HA	3:F:189:HIS:O	1.88	0.71
3:C:11:ALA:HB3	3:C:94:LEU:H	1.57	0.70
3:D:94:LEU:HD13	7:P:2:U:H1'	1.73	0.68
3:H:26:VAL:O	3:H:76:THR:HB	1.92	0.68
4:I:89:LYS:HZ2	4:I:90:THR:H	1.40	0.67
3:F:129:LEU:HD13	3:F:191:MET:HE2	1.79	0.64
3:F:203:GLN:NE2	7:P:21:C:OP2	2.31	0.64
3:C:12:TYR:HE2	3:C:295:LEU:HD22	1.65	0.62
4:I:4:THR:HA	4:I:56:PHE:HA	1.81	0.62
2:B:15:ILE:HD11	2:B:27:PRO:HG3	1.82	0.62
3:G:78:LEU:HA	3:G:189:HIS:O	2.00	0.62
3:C:140:GLU:HG2	3:C:190:TYR:HB2	1.82	0.61
1:A:168:CYS:HB2	2:B:25:GLY:HA3	1.82	0.61
3:F:90:PHE:HB3	3:F:104:GLN:HG2	1.82	0.61
5:Q:20:DG:H1	7:P:14:U:H3	1.48	0.61
3:C:26:VAL:O	3:C:76:THR:HB	2.01	0.61
3:F:235:THR:HA	3:F:248:ALA:HA	1.83	0.60
3:E:19:THR:HG1	3:E:82:ASN:H	1.49	0.60
1:A:123:ASN:HB3	2:B:22:ILE:HG13	1.84	0.60
3:G:227:GLN:HB2	7:P:23:C:H5'	1.82	0.60
3:E:15:LYS:HB2	3:E:296:ILE:HG23	1.84	0.60
4:I:41:THR:HG21	4:I:88:VAL:HB	1.84	0.59
3:C:257:ARG:NH2	7:P:2:U:OP1	2.35	0.59
4:I:109:ARG:HG3	4:I:110:LYS:HD2	1.84	0.59
3:C:270:LEU:HG	3:C:274:MET:HE3	1.85	0.59
3:F:76:THR:HA	3:F:191:MET:O	2.03	0.59
3:E:24:PHE:HB2	3:E:78:LEU:HB3	1.84	0.58
3:F:69:ASN:ND2	3:G:20:ASN:OD1	2.36	0.58
3:C:122:ASN:ND2	3:C:235:THR:O	2.37	0.58
3:F:111:ASP:HB3	3:F:289:MET:HE3	1.85	0.58
3:E:79:VAL:HB	3:E:189:HIS:HB2	1.86	0.58
3:G:286:GLU:HA	3:G:289:MET:HE2	1.85	0.57
3:E:70:LEU:HD21	3:E:77:LEU:HB2	1.86	0.57
2:B:74:GLN:HE22	7:P:-2:A:H1'	1.68	0.57
3:D:3:LYS:HA	3:D:279:ASN:OD1	2.04	0.57
3:F:228:LYS:NZ	7:P:18:C:O2	2.38	0.57
3:G:14:ARG:NH1	3:G:17:ASN:OD1	2.38	0.57
4:I:144:ARG:HD3	3:H:198:GLU:HG3	1.85	0.57
7:P:29:A:H5''	3:H:227:GLN:HB2	1.86	0.57
2:B:11:LYS:HB3	2:B:132:ALA:HB3	1.87	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:250:GLU:HG3	3:G:252:TYR:H	1.68	0.57
3:D:96:GLY:HA2	3:D:100:SER:HB3	1.85	0.57
3:E:234:ARG:HE	3:E:248:ALA:HB1	1.69	0.57
3:D:15:LYS:HD2	3:D:89:VAL:HA	1.87	0.56
2:B:16:ASN:ND2	7:P:-4:A:OP2	2.38	0.56
3:F:26:VAL:O	3:F:76:THR:HB	2.05	0.56
3:C:47:THR:HG21	3:D:250:GLU:HG2	1.88	0.56
4:I:44:ARG:NH1	4:I:49:VAL:O	2.37	0.56
3:E:118:ARG:NH2	3:E:238:THR:OG1	2.39	0.56
3:F:42:LYS:NZ	3:G:226:ASP:OD2	2.38	0.56
3:H:65:VAL:HG12	3:H:213:LYS:H	1.71	0.56
3:H:12:TYR:HH	3:H:100:SER:HG	1.54	0.56
3:E:47:THR:HG21	3:F:250:GLU:HG2	1.86	0.55
3:E:241:GLU:OE2	3:E:266:ASN:ND2	2.39	0.55
3:F:79:VAL:HB	3:F:189:HIS:HB2	1.89	0.55
2:B:63:ARG:NH1	6:R:26:DT:OP2	2.39	0.55
1:A:28:ASN:HD22	1:A:101:LEU:HD21	1.70	0.55
3:D:78:LEU:HD21	3:D:80:LYS:HE3	1.89	0.55
3:C:251:PRO:HB3	3:C:270:LEU:HD13	1.89	0.55
3:E:191:MET:SD	3:E:191:MET:N	2.80	0.54
7:P:30:A:OP2	3:H:234:ARG:NH2	2.39	0.54
3:C:163:ASP:N	3:C:163:ASP:OD1	2.40	0.54
3:C:257:ARG:HD3	3:D:94:LEU:HD11	1.89	0.54
3:D:226:ASP:OD1	3:D:226:ASP:N	2.39	0.54
3:E:44:VAL:HA	3:F:227:GLN:HE22	1.73	0.54
2:B:185:GLY:HA3	2:B:238:LYS:HG2	1.90	0.54
3:F:263:ARG:NH1	3:F:268:THR:O	2.36	0.54
3:E:198:GLU:OE1	3:F:14:ARG:NH2	2.41	0.54
1:A:172:GLN:NE2	2:B:57:PHE:O	2.41	0.54
3:G:234:ARG:NH1	3:G:249:VAL:O	2.41	0.54
3:D:251:PRO:HB3	3:D:270:LEU:HD12	1.90	0.54
3:F:107:LEU:HD21	3:F:291:VAL:HG12	1.90	0.54
3:F:204:GLU:OE1	3:F:214:GLN:NE2	2.41	0.54
3:E:42:LYS:NZ	3:F:226:ASP:OD1	2.41	0.53
3:E:21:ALA:HB3	3:E:224:MET:HB2	1.89	0.53
1:A:137:TYR:O	6:R:26:DT:N3	2.41	0.53
3:E:257:ARG:NH2	7:P:12:G:O2'	2.39	0.53
4:I:4:THR:HG22	4:I:56:PHE:HB3	1.89	0.53
3:C:127:ARG:NH2	3:C:157:VAL:O	2.42	0.53
3:H:36:ALA:HB1	3:H:221:GLN:HG3	1.90	0.53
3:C:93:GLU:HB3	7:P:-4:A:H2	1.74	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:CYS:SG	1:A:156:CYS:CB	2.97	0.53
3:D:195:GLU:O	3:E:82:ASN:ND2	2.42	0.53
3:E:176:VAL:HG21	3:E:183:VAL:HB	1.91	0.53
4:I:57:SER:HB3	4:I:61:VAL:HG22	1.91	0.53
3:F:50:VAL:HG13	3:F:52:GLY:H	1.74	0.52
3:H:79:VAL:HB	3:H:189:HIS:HB2	1.91	0.52
3:G:43:THR:O	3:H:227:GLN:NE2	2.43	0.52
3:G:127:ARG:NH2	3:G:157:VAL:O	2.42	0.52
3:H:237:ASP:OD1	3:H:263:ARG:NH2	2.42	0.52
2:B:60:ARG:NH2	3:C:227:GLN:OE1	2.37	0.52
3:F:90:PHE:CB	3:F:104:GLN:HG2	2.40	0.52
3:F:286:GLU:HA	3:F:289:MET:HG3	1.91	0.52
1:A:124:LEU:HA	2:B:22:ILE:HD11	1.91	0.51
3:C:21:ALA:HB3	3:C:224:MET:HB2	1.92	0.51
3:E:247:ILE:HD13	3:E:262:TYR:HD2	1.75	0.51
3:G:231:ASN:HB2	7:P:24:U:H5'	1.92	0.51
3:H:77:LEU:HD22	3:H:199:VAL:HG12	1.92	0.51
3:C:129:LEU:HA	3:C:224:MET:HE1	1.92	0.51
3:E:4:LEU:CD1	3:E:102:THR:HG21	2.41	0.51
3:F:100:SER:O	3:F:104:GLN:HG3	2.10	0.51
3:D:142:VAL:HA	3:D:150:PHE:O	2.10	0.51
3:H:118:ARG:HG2	3:H:236:ILE:HG23	1.92	0.51
3:F:23:MET:HB3	3:F:77:LEU:HD11	1.92	0.51
4:I:99:LYS:HG3	4:I:125:LYS:HD2	1.93	0.51
3:C:230:GLY:O	3:C:234:ARG:HB3	2.10	0.51
3:E:13:SER:HB3	3:E:93:GLU:HG3	1.92	0.51
3:C:66:ASN:HB2	3:C:214:GLN:HA	1.93	0.51
2:B:170:ILE:HD12	2:B:212:TYR:HB2	1.93	0.51
3:F:16:ILE:HG22	3:F:18:PRO:HD3	1.93	0.51
3:H:143:ILE:HG12	3:H:187:VAL:HG22	1.93	0.51
3:C:43:THR:HG22	3:C:62:ILE:HG22	1.93	0.50
3:H:24:PHE:HB2	3:H:78:LEU:HB3	1.93	0.50
3:E:99:TYR:CE1	3:E:278:VAL:HG21	2.47	0.50
3:C:46:GLY:HA3	7:P:5:A:H1'	1.94	0.49
3:D:16:ILE:HG22	3:D:18:PRO:HD3	1.94	0.49
3:E:195:GLU:O	3:F:82:ASN:ND2	2.45	0.49
1:A:19:HIS:HB3	1:A:118:ILE:HD11	1.94	0.49
3:D:11:ALA:HB3	3:D:94:LEU:HB2	1.93	0.49
3:E:78:LEU:HA	3:E:189:HIS:O	2.13	0.49
3:F:283:LEU:HB2	3:F:288:GLN:HE21	1.78	0.49
2:B:62:SER:HB3	3:C:206:VAL:HG21	1.94	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:155:MET:HB2	2:B:220:VAL:HG21	1.95	0.49
3:D:46:GLY:HA3	7:P:11:U:H1'	1.93	0.49
2:B:84:ASN:OD1	2:B:85:ASN:N	2.43	0.49
3:C:298:GLY:HA3	7:P:-2:A:H5''	1.95	0.49
3:F:14:ARG:N	7:P:16:A:OP1	2.44	0.49
3:G:45:ALA:HA	3:G:61:ASN:O	2.13	0.49
3:H:163:ASP:N	3:H:163:ASP:OD1	2.46	0.49
2:B:62:SER:OG	2:B:63:ARG:N	2.45	0.49
3:D:237:ASP:OD1	3:D:263:ARG:NH2	2.46	0.48
3:D:26:VAL:HG23	3:D:76:THR:HB	1.94	0.48
2:B:68:VAL:HB	2:B:91:LYS:H	1.78	0.48
3:D:24:PHE:HB2	3:D:78:LEU:HD23	1.95	0.48
3:F:119:TYR:HE2	3:F:296:ILE:HG21	1.78	0.48
3:D:78:LEU:HA	3:D:189:HIS:O	2.13	0.48
3:F:26:VAL:HG23	3:F:76:THR:HB	1.96	0.48
3:C:119:TYR:HE1	3:C:236:ILE:HD13	1.78	0.48
3:C:123:ILE:HD11	3:C:187:VAL:HG21	1.96	0.48
3:D:118:ARG:NH2	3:D:238:THR:OG1	2.46	0.48
3:F:43:THR:HG21	3:G:205:PHE:HB3	1.96	0.48
6:R:21:DT:H2'	6:R:22:DT:H4'	1.96	0.48
3:D:4:LEU:HB3	3:D:278:VAL:CG1	2.44	0.48
3:D:55:ASN:OD1	3:D:55:ASN:N	2.44	0.48
3:E:129:LEU:H	3:E:129:LEU:HG	1.39	0.48
1:A:104:ILE:HD13	7:P:-5:U:H5''	1.96	0.48
3:F:121:TYR:HB2	3:F:166:VAL:HG21	1.95	0.48
3:F:127:ARG:HH22	3:F:158:ASN:HA	1.78	0.48
3:F:14:ARG:HH11	7:P:16:A:H5'	1.79	0.48
2:B:11:LYS:HD3	2:B:97:VAL:HG22	1.95	0.47
3:C:64:THR:OG1	3:C:210:LYS:O	2.32	0.47
3:C:225:HIS:ND1	7:P:1:G:OP1	2.47	0.47
3:C:274:MET:HE1	3:C:291:VAL:HG22	1.94	0.47
3:E:99:TYR:HE2	3:E:274:MET:HE3	1.79	0.47
3:E:65:VAL:HG12	3:E:212:SER:HB3	1.96	0.47
3:G:130:TRP:HH2	3:G:231:ASN:HB3	1.80	0.47
3:E:193:LEU:O	3:F:84:LYS:NZ	2.38	0.47
3:D:281:LYS:NZ	3:D:282:SER:O	2.43	0.47
3:E:10:LEU:HD23	3:E:301:PHE:HB2	1.96	0.47
3:F:255:VAL:HG13	7:P:18:C:H41	1.79	0.47
4:I:5:ILE:HG22	4:I:82:PRO:HA	1.96	0.47
4:I:105:ARG:HD3	7:P:41:C:H6	1.79	0.47
3:H:20:ASN:OD1	3:H:20:ASN:N	2.47	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:203:GLN:HG2	3:G:213:LYS:HG2	1.96	0.47
3:D:7:PRO:O	3:D:8:ALA:C	2.56	0.47
3:F:266:ASN:N	3:F:266:ASN:OD1	2.47	0.47
1:A:142:GLN:HE21	1:A:163:PHE:HB2	1.80	0.47
3:H:176:VAL:C	3:H:178:SER:H	2.23	0.47
3:G:234:ARG:HB2	3:G:249:VAL:HG22	1.96	0.47
1:A:142:GLN:O	1:A:162:GLY:CA	2.59	0.46
3:F:48:GLN:HG3	3:F:59:LYS:HA	1.98	0.46
3:C:15:LYS:HZ3	3:C:89:VAL:HG12	1.80	0.46
3:C:21:ALA:HB2	3:C:229:ILE:HD11	1.96	0.46
3:G:123:ILE:HD12	3:G:187:VAL:HG21	1.96	0.46
4:I:155:SER:HB3	4:I:163:LEU:HD22	1.97	0.46
2:B:67:SER:OG	3:C:205:PHE:O	2.30	0.46
3:C:277:ALA:HB2	3:C:283:LEU:HD11	1.97	0.46
3:H:102:THR:HG21	3:H:278:VAL:HG13	1.97	0.46
2:B:18:LYS:HA	2:B:24:VAL:HG23	1.97	0.46
3:D:152:PHE:HB3	3:D:155:LEU:HD11	1.98	0.46
3:E:85:PHE:HB2	3:E:183:VAL:HG12	1.97	0.46
3:F:121:TYR:OH	3:F:156:LEU:O	2.33	0.46
3:G:287:ASP:OD1	3:G:287:ASP:N	2.49	0.46
2:B:186:TYR:O	2:B:237:THR:OG1	2.34	0.46
3:C:199:VAL:HG23	3:C:201:PRO:HD3	1.97	0.46
4:I:12:ALA:HB3	4:I:17:VAL:HG22	1.97	0.46
3:D:7:PRO:HB2	3:D:9:VAL:O	2.16	0.46
3:F:127:ARG:NH2	3:F:157:VAL:O	2.49	0.46
4:I:101:LYS:NZ	7:P:39:A:OP2	2.39	0.46
2:B:50:VAL:O	2:B:231:TRP:N	2.43	0.45
2:B:68:VAL:HG13	2:B:88:ILE:HD11	1.97	0.45
2:B:235:GLU:HA	2:B:236:ASN:HA	1.76	0.45
3:D:153:SER:OG	3:D:154:ASP:N	2.49	0.45
3:E:266:ASN:OD1	3:E:266:ASN:N	2.50	0.45
3:H:118:ARG:HB3	3:H:236:ILE:HD12	1.98	0.45
1:A:50:LYS:HE3	1:A:50:LYS:HB2	1.75	0.45
3:D:235:THR:HA	3:D:248:ALA:HA	1.97	0.45
3:F:267:LYS:O	3:F:272:THR:OG1	2.32	0.45
4:I:57:SER:OG	4:I:60:GLY:O	2.33	0.45
4:I:92:ARG:NE	7:P:49:G:OP2	2.48	0.45
3:C:31:ARG:NH1	3:C:140:GLU:OE2	2.47	0.45
3:D:103:LEU:O	3:D:107:LEU:HB2	2.16	0.45
3:D:18:PRO:HB3	3:D:83:VAL:HG22	1.98	0.45
3:H:128:THR:HA	3:H:232:ALA:HB3	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:210:LYS:HG3	3:H:211:LEU:HG	1.98	0.45
2:B:31:THR:OG1	7:P:-5:U:OP1	2.34	0.45
3:E:202:SER:HG	3:E:225:HIS:CE1	2.34	0.45
3:E:230:GLY:HA2	3:E:233:ILE:HD12	1.97	0.45
3:G:102:THR:HA	3:G:105:THR:HG22	1.98	0.45
5:Q:28:DT:H6	5:Q:28:DT:H2'	1.57	0.45
2:B:11:LYS:HE2	2:B:11:LYS:HB2	1.78	0.45
2:B:60:ARG:O	2:B:66:THR:OG1	2.32	0.45
3:E:107:LEU:HD21	3:E:291:VAL:HG13	1.98	0.45
2:B:70:LYS:H	2:B:70:LYS:HG3	1.53	0.45
3:C:117:TYR:OH	3:C:161:ASP:O	2.34	0.45
3:D:94:LEU:CD1	7:P:2:U:H1'	2.45	0.45
3:E:156:LEU:H	3:E:156:LEU:HG	1.55	0.45
3:H:48:GLN:HB3	3:H:49:SER:H	1.61	0.45
1:A:20:SER:O	1:A:119:GLY:N	2.47	0.45
3:C:26:VAL:O	3:C:76:THR:CB	2.64	0.45
3:G:163:ASP:OD1	3:G:163:ASP:N	2.50	0.45
3:H:235:THR:HA	3:H:248:ALA:HA	1.98	0.45
3:E:231:ASN:O	3:E:235:THR:OG1	2.34	0.44
3:F:15:LYS:HE3	3:F:89:VAL:HA	1.98	0.44
3:G:235:THR:HA	3:G:248:ALA:HA	1.99	0.44
3:C:44:VAL:O	3:C:62:ILE:HA	2.18	0.44
3:C:152:PHE:HE1	3:C:165:ASP:HB3	1.82	0.44
3:C:154:ASP:OD1	3:C:154:ASP:N	2.50	0.44
3:E:61:ASN:ND2	5:Q:17:DT:O4'	2.49	0.44
2:B:125:ARG:HD3	2:B:128:GLY:HA2	1.99	0.44
3:F:25:ALA:HB3	3:F:35:THR:HG22	1.99	0.44
4:I:44:ARG:NH2	4:I:52:ILE:O	2.49	0.44
3:F:289:MET:HB3	3:F:289:MET:HE2	1.76	0.44
7:P:32:U:H3'	7:P:33:G:H8	1.82	0.44
3:H:237:ASP:HB3	3:H:247:ILE:HB	2.00	0.44
3:D:270:LEU:HG	3:D:294:ASN:HD22	1.83	0.44
1:A:25:ASN:OD1	2:B:89:GLN:NE2	2.51	0.44
3:H:176:VAL:HG11	3:H:183:VAL:HG12	1.99	0.44
2:B:95:ASP:OD1	2:B:95:ASP:N	2.50	0.44
3:F:15:LYS:HG3	3:F:89:VAL:HG23	2.00	0.44
3:G:213:LYS:HB3	3:G:213:LYS:HE3	1.71	0.44
3:C:249:VAL:HG21	3:C:293:ALA:HB1	2.00	0.43
4:I:24:VAL:HG22	4:I:70:LEU:HD11	2.00	0.43
3:H:78:LEU:HA	3:H:189:HIS:O	2.18	0.43
3:F:116:ALA:HA	3:F:119:TYR:HB2	2.00	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:I:7:ILE:HD12	4:I:7:ILE:HA	1.85	0.43
3:F:26:VAL:O	3:F:76:THR:CB	2.66	0.43
3:F:28:TRP:HB2	3:F:192:LEU:HD21	1.99	0.43
3:C:12:TYR:CD2	3:C:295:LEU:HB3	2.53	0.43
3:E:286:GLU:HA	3:E:289:MET:HE3	1.99	0.43
3:C:235:THR:HA	3:C:248:ALA:HA	2.01	0.43
3:F:121:TYR:OH	3:F:159:GLU:O	2.30	0.43
1:A:45:ASP:HA	1:A:63:TRP:HE1	1.83	0.43
3:C:252:TYR:HB3	3:C:264:ALA:HB2	2.00	0.43
3:D:78:LEU:HD11	3:D:188:GLU:HB3	2.01	0.43
3:F:114:THR:O	3:F:115:LEU:C	2.58	0.43
3:C:289:MET:SD	3:C:289:MET:N	2.87	0.43
3:E:231:ASN:HB2	7:P:12:G:H5'	2.00	0.43
3:F:145:VAL:O	3:F:148:GLN:HB3	2.18	0.43
3:F:241:GLU:H	3:F:241:GLU:HG2	1.40	0.43
3:D:153:SER:OG	3:D:154:ASP:OD1	2.36	0.43
3:H:22:LEU:HD23	3:H:22:LEU:HA	1.85	0.43
3:C:10:LEU:O	3:C:300:VAL:HA	2.19	0.42
3:C:26:VAL:HG23	3:C:76:THR:HB	2.00	0.42
3:D:7:PRO:HD2	3:D:10:LEU:HB2	1.99	0.42
3:E:142:VAL:HG23	3:E:188:GLU:HB2	2.01	0.42
3:D:254:SER:HA	3:D:261:ALA:HA	2.00	0.42
2:B:26:MET:HG3	2:B:27:PRO:HD2	2.01	0.42
3:C:11:ALA:HB1	7:P:-3:A:H1'	2.00	0.42
3:E:78:LEU:HD11	3:E:188:GLU:HB3	2.01	0.42
3:G:21:ALA:HB2	3:G:229:ILE:HD11	2.01	0.42
7:P:33:G:C5	7:P:34:U:H4'	2.54	0.42
2:B:68:VAL:HG11	2:B:90:PRO:HA	2.01	0.42
3:E:44:VAL:HG22	3:F:227:GLN:HE22	1.84	0.42
3:H:41:THR:HB	3:H:64:THR:HB	2.01	0.42
1:A:9:PHE:HE1	1:A:48:THR:HB	1.83	0.42
3:G:20:ASN:HD22	3:G:218:LEU:HD11	1.84	0.42
3:H:91:LYS:HE2	3:H:91:LYS:HB2	1.90	0.42
3:G:102:THR:HG23	3:G:278:VAL:HG22	2.02	0.42
3:C:24:PHE:HB2	3:C:78:LEU:HB3	2.02	0.42
3:D:134:VAL:HB	3:E:86:VAL:HG21	2.02	0.42
3:D:241:GLU:H	3:D:241:GLU:HG3	1.55	0.42
3:F:211:LEU:HD23	3:F:214:GLN:HB2	2.00	0.42
3:G:214:GLN:HE21	3:G:214:GLN:HB2	1.47	0.42
3:C:137:GLU:HA	3:D:182:PHE:HZ	1.84	0.42
3:D:263:ARG:HD2	3:D:268:THR:HB	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:163:ASP:OD1	3:E:163:ASP:N	2.52	0.42
3:F:130:TRP:CD1	3:F:228:LYS:HZ2	2.38	0.42
3:F:154:ASP:OD1	3:F:154:ASP:N	2.50	0.42
1:A:161:GLY:H	6:R:21:DT:H3	1.67	0.42
3:E:145:VAL:O	3:E:146:ASN:C	2.63	0.42
4:I:147:MET:H	4:I:147:MET:HG2	1.69	0.42
3:H:234:ARG:HB2	3:H:249:VAL:HG12	2.01	0.42
2:B:178:LYS:HB2	2:B:178:LYS:HE2	1.62	0.41
3:D:148:GLN:HG3	3:D:150:PHE:CE1	2.55	0.41
3:D:270:LEU:HD11	3:D:301:PHE:CZ	2.55	0.41
3:G:85:PHE:HB3	3:G:177:LEU:HD21	2.01	0.41
3:H:277:ALA:HB2	3:H:283:LEU:HD11	2.01	0.41
3:F:172:MET:HE2	3:F:172:MET:HB3	1.80	0.41
3:C:76:THR:HA	3:C:191:MET:O	2.20	0.41
3:D:120:VAL:HG11	3:D:169:ILE:HB	2.03	0.41
3:E:263:ARG:NH1	3:E:268:THR:O	2.47	0.41
3:G:91:LYS:HE2	3:G:91:LYS:HB2	1.94	0.41
3:G:131:ARG:NH1	7:P:27:U:OP2	2.53	0.41
3:C:234:ARG:HH11	3:C:297:ARG:HG3	1.85	0.41
3:D:111:ASP:HB3	3:D:289:MET:HE3	2.02	0.41
3:E:228:LYS:NZ	7:P:14:U:OP1	2.52	0.41
3:G:137:GLU:HB2	3:G:192:LEU:HB2	2.02	0.41
2:B:1:MET:HE2	2:B:141:VAL:HG13	2.02	0.41
2:B:51:SER:HB3	2:B:230:LEU:HD22	2.02	0.41
2:B:89:GLN:OE1	2:B:89:GLN:N	2.53	0.41
3:C:80:LYS:HG3	3:C:188:GLU:HG2	2.02	0.41
3:G:276:GLY:HA3	3:G:281:LYS:HE3	2.01	0.41
4:I:42:PHE:H	4:I:91:TYR:HE2	1.68	0.41
3:C:28:TRP:O	3:C:31:ARG:NE	2.52	0.41
3:C:266:ASN:OD1	3:C:266:ASN:N	2.54	0.41
3:D:272:THR:C	3:D:274:MET:H	2.29	0.41
3:F:297:ARG:NH1	3:F:298:GLY:O	2.53	0.41
3:E:115:LEU:HD12	3:E:115:LEU:HA	1.88	0.41
3:E:240:TYR:OH	3:E:245:THR:O	2.34	0.41
3:E:289:MET:HE3	3:E:289:MET:HB3	1.88	0.41
3:F:98:GLU:N	3:F:101:ASN:OD1	2.54	0.41
4:I:165:ILE:HG12	7:P:49:G:H4'	2.03	0.41
3:C:128:THR:HG23	3:C:129:LEU:HG	2.03	0.41
3:C:237:ASP:HB3	3:C:247:ILE:HG12	2.03	0.41
3:D:94:LEU:HD22	7:P:2:U:O2	2.21	0.41
3:E:131:ARG:O	3:E:134:VAL:HG12	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:P:2:U:H2'	7:P:3:C:C6	2.56	0.41
3:H:156:LEU:HD23	3:H:159:GLU:HB3	2.02	0.41
1:A:103:GLN:N	7:P:-4:A:N7	2.58	0.41
3:C:169:ILE:HA	3:C:172:MET:HG2	2.01	0.41
3:G:257:ARG:HA	3:G:257:ARG:HD3	1.91	0.41
4:I:130:LYS:HG3	4:I:132:PRO:HD3	2.02	0.41
2:B:179:LEU:HD13	2:B:179:LEU:HA	1.85	0.40
3:D:66:ASN:HB2	3:D:214:GLN:HA	2.04	0.40
3:D:4:LEU:HB3	3:D:278:VAL:HG12	2.02	0.40
3:F:21:ALA:HB3	3:F:224:MET:HB2	2.02	0.40
7:P:28:U:OP1	3:H:14:ARG:N	2.54	0.40
2:B:182:MET:SD	2:B:215:ILE:HD13	2.61	0.40
3:E:107:LEU:HD23	3:E:292:THR:HB	2.03	0.40
3:E:251:PRO:HA	3:E:270:LEU:HB2	2.03	0.40
3:E:292:THR:O	3:E:296:ILE:HG13	2.21	0.40
7:P:4:A:H2'	7:P:5:A:C8	2.55	0.40
3:H:71:PRO:HG2	3:H:74:LYS:HG2	2.03	0.40
3:F:115:LEU:HD22	3:F:292:THR:HG23	2.03	0.40
3:F:281:LYS:HE3	3:F:281:LYS:HB2	1.90	0.40
3:G:15:LYS:HB2	3:G:296:ILE:HG13	2.03	0.40
3:G:233:ILE:HD13	3:G:233:ILE:HA	1.84	0.40
3:C:120:VAL:HA	3:C:123:ILE:HG22	2.03	0.40
3:G:107:LEU:HD23	3:G:292:THR:HB	2.04	0.40
3:H:123:ILE:HD11	3:H:233:ILE:HD11	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	A	171/179 (96%)	150 (88%)	20 (12%)	1 (1%)	21 55

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	244/248 (98%)	227 (93%)	17 (7%)	0	100	100
3	C	300/306 (98%)	283 (94%)	16 (5%)	1 (0%)	36	66
3	D	300/306 (98%)	280 (93%)	19 (6%)	1 (0%)	36	66
3	E	300/306 (98%)	282 (94%)	16 (5%)	2 (1%)	18	52
3	F	300/306 (98%)	271 (90%)	27 (9%)	2 (1%)	18	52
3	G	300/306 (98%)	275 (92%)	23 (8%)	2 (1%)	18	52
3	H	300/306 (98%)	276 (92%)	19 (6%)	5 (2%)	7	35
4	I	164/168 (98%)	142 (87%)	21 (13%)	1 (1%)	21	55
All	All	2379/2431 (98%)	2186 (92%)	178 (8%)	15 (1%)	23	55

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	F	53	ASN
3	F	54	PRO
1	A	69	ALA
3	H	54	PRO
3	H	132	ASN
3	D	8	ALA
3	E	126	GLY
3	E	130	TRP
3	G	53	ASN
3	H	280	GLY
3	C	98	GLU
3	G	212	SER
3	H	131	ARG
3	H	177	LEU
4	I	132	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	A	143/157 (91%)	137 (96%)	6 (4%)	26	52
2	B	207/211 (98%)	202 (98%)	5 (2%)	43	64
3	C	238/251 (95%)	232 (98%)	6 (2%)	42	63
3	D	238/251 (95%)	222 (93%)	16 (7%)	15	41
3	E	240/251 (96%)	229 (95%)	11 (5%)	24	50
3	F	240/251 (96%)	227 (95%)	13 (5%)	20	47
3	G	237/251 (94%)	221 (93%)	16 (7%)	14	41
3	H	237/251 (94%)	232 (98%)	5 (2%)	47	66
4	I	147/151 (97%)	144 (98%)	3 (2%)	48	67
All	All	1927/2025 (95%)	1846 (96%)	81 (4%)	28	52

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	GLN
1	A	66	LEU
1	A	70	ASN
1	A	71	MET
1	A	73	TYR
1	A	152	CYS
2	B	69	THR
2	B	178	LYS
2	B	179	LEU
2	B	202	PHE
2	B	226	PHE
3	C	63	GLN
3	C	77	LEU
3	C	100	SER
3	C	295	LEU
3	C	296	ILE
3	C	297	ARG
3	D	4	LEU
3	D	35	THR
3	D	38	MET
3	D	39	VAL
3	D	86	VAL
3	D	91	LYS
3	D	94	LEU
3	D	100	SER
3	D	241	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	D	242	ASP
3	D	244	THR
3	D	245	THR
3	D	260	VAL
3	D	268	THR
3	D	270	LEU
3	D	274	MET
3	E	5	LYS
3	E	91	LYS
3	E	93	GLU
3	E	94	LEU
3	E	129	LEU
3	E	145	VAL
3	E	148	GLN
3	E	151	THR
3	E	155	LEU
3	E	156	LEU
3	E	157	VAL
3	F	34	THR
3	F	41	THR
3	F	55	ASN
3	F	56	ASP
3	F	109	ASN
3	F	110	THR
3	F	114	THR
3	F	115	LEU
3	F	237	ASP
3	F	241	GLU
3	F	244	THR
3	F	288	GLN
3	F	289	MET
3	G	4	LEU
3	G	44	VAL
3	G	48	GLN
3	G	49	SER
3	G	50	VAL
3	G	55	ASN
3	G	56	ASP
3	G	61	ASN
3	G	119	TYR
3	G	188	GLU
3	G	191	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	G	206	VAL
3	G	213	LYS
3	G	237	ASP
3	G	266	ASN
3	G	268	THR
4	I	49	VAL
4	I	56	PHE
4	I	148	PHE
3	H	78	LEU
3	H	129	LEU
3	H	180	GLU
3	H	233	ILE
3	H	278	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	A	70	ASN
1	A	103	GLN
1	A	122	HIS
1	A	123	ASN
1	A	142	GLN
1	A	154	HIS
1	A	157	ASN
2	B	9	ASN
2	B	122	ASN
2	B	135	ASN
2	B	180	GLN
3	C	146	ASN
3	D	61	ASN
3	D	104	GLN
3	D	294	ASN
3	E	20	ASN
3	E	109	ASN
3	E	146	ASN
3	F	17	ASN
3	F	69	ASN
3	F	72	HIS
3	F	214	GLN
3	F	227	GLN
3	F	288	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	G	20	ASN
3	G	214	GLN
3	G	266	ASN
3	G	288	GLN
3	G	294	ASN
4	I	35	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
7	P	58/59 (98%)	36 (62%)	4 (6%)

All (36) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
7	P	-5	U
7	P	-4	A
7	P	-2	A
7	P	1	G
7	P	2	U
7	P	4	A
7	P	7	C
7	P	8	C
7	P	9	U
7	P	13	C
7	P	18	C
7	P	19	U
7	P	20	U
7	P	23	C
7	P	24	U
7	P	25	A
7	P	26	U
7	P	27	U
7	P	28	U
7	P	29	A
7	P	30	A
7	P	31	A
7	P	32	U
7	P	34	U
7	P	35	U
7	P	37	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
7	P	38	C
7	P	39	A
7	P	40	G
7	P	41	C
7	P	42	C
7	P	43	G
7	P	44	C
7	P	45	A
7	P	46	U
7	P	47	A

All (4) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
7	P	24	U
7	P	29	A
7	P	33	G
7	P	34	U

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.

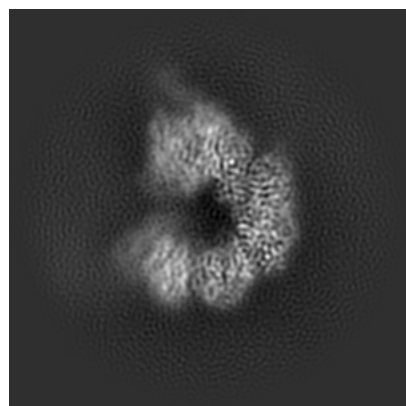
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-36831. These allow visual inspection of the internal detail of the map and identification of artifacts.

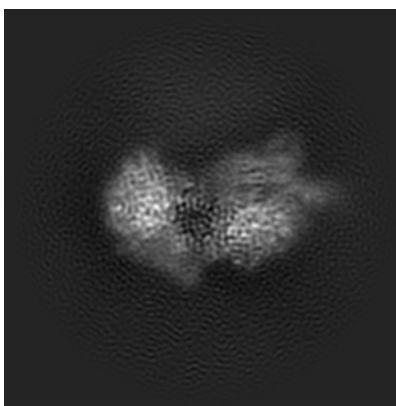
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

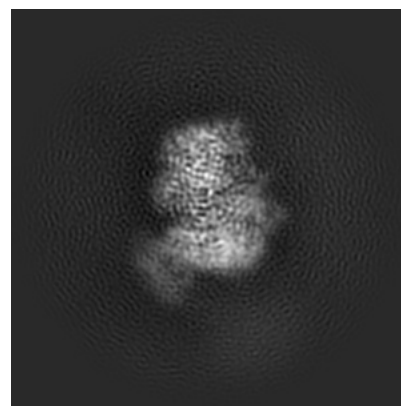
6.1.1 Primary map



X

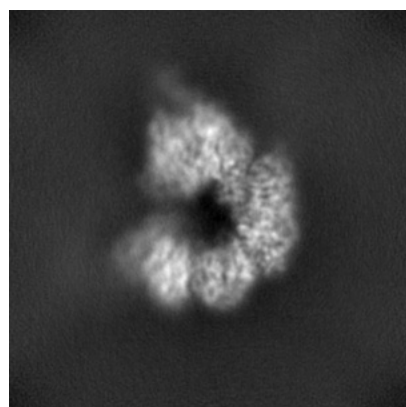


Y

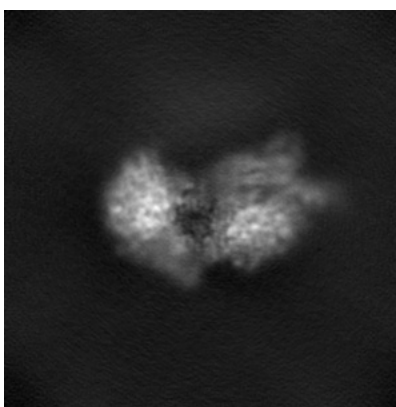


Z

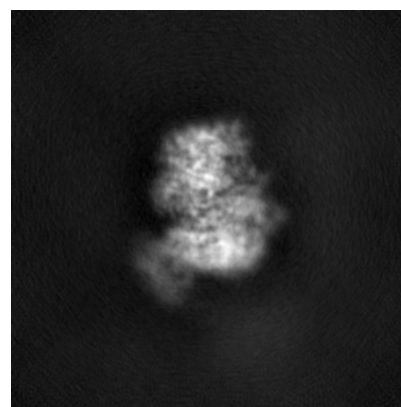
6.1.2 Raw map



X



Y

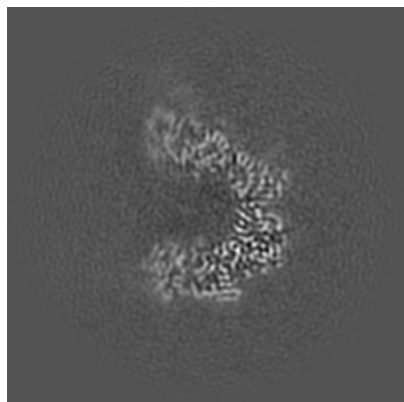


Z

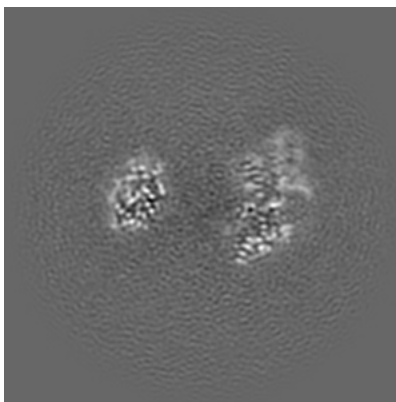
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

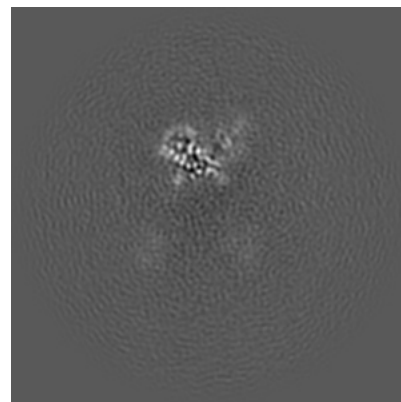
6.2.1 Primary map



X Index: 120

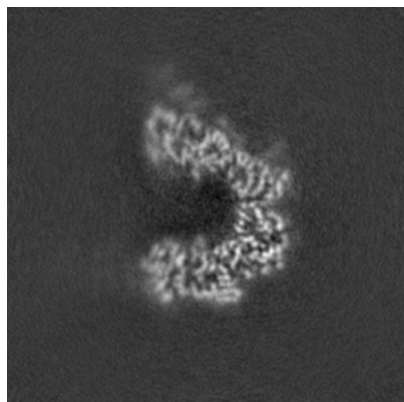


Y Index: 120

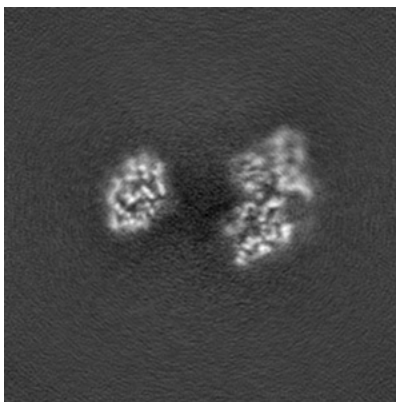


Z Index: 120

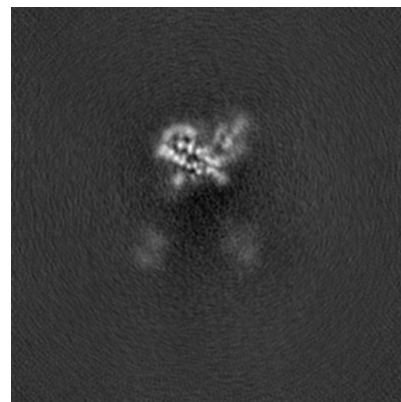
6.2.2 Raw map



X Index: 120



Y Index: 120

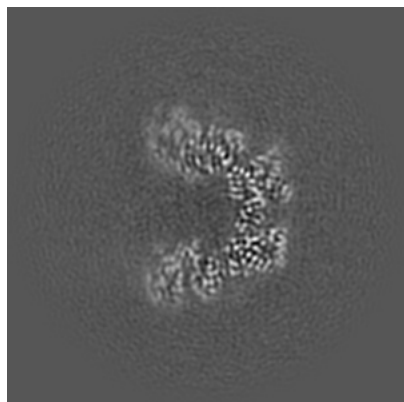


Z Index: 120

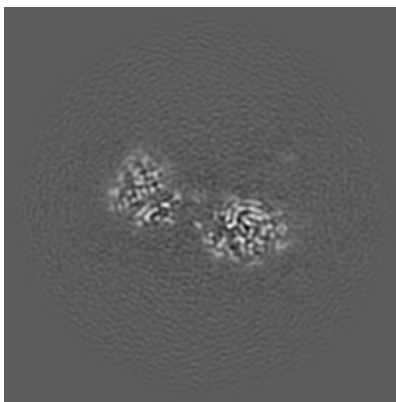
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

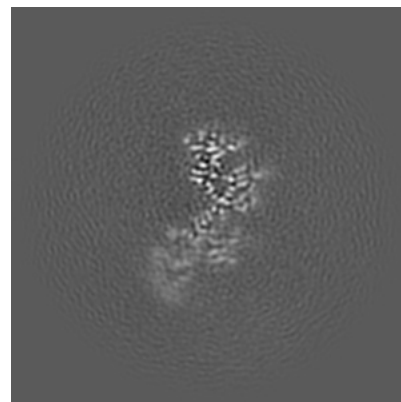
6.3.1 Primary map



X Index: 115

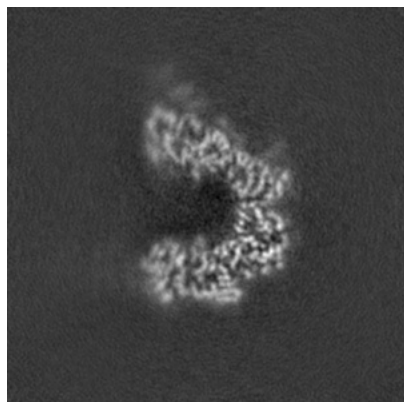


Y Index: 134

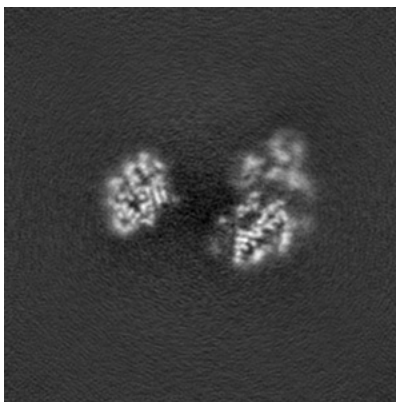


Z Index: 91

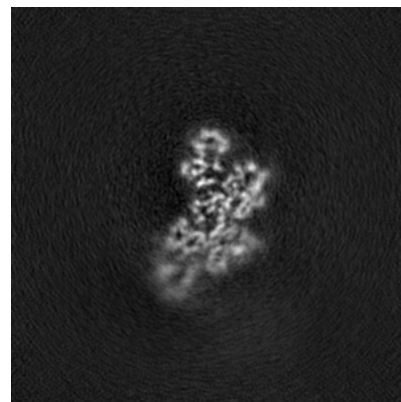
6.3.2 Raw map



X Index: 120



Y Index: 123

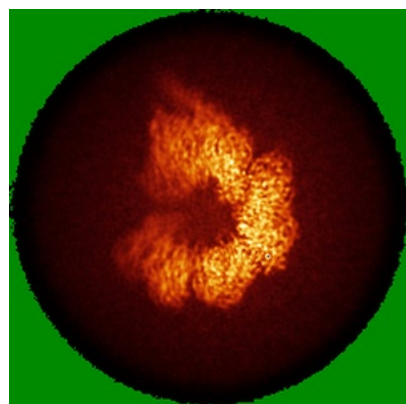


Z Index: 85

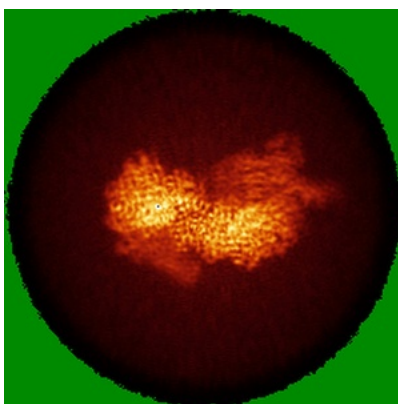
The images above show the largest variance slices of the map in three orthogonal directions.

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

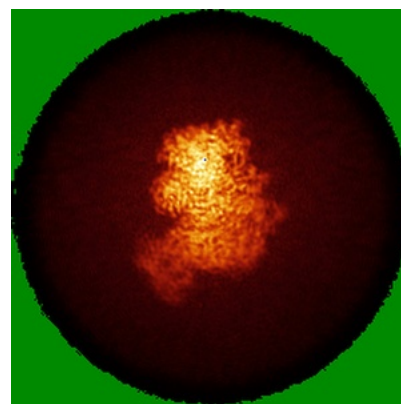
6.4.1 Primary map



X

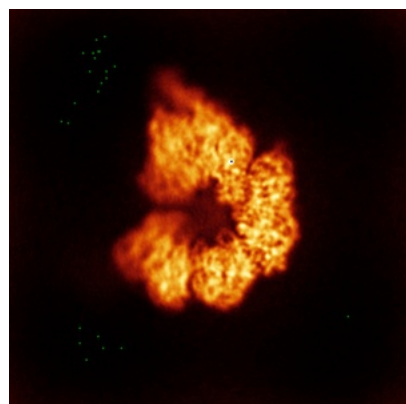


Y

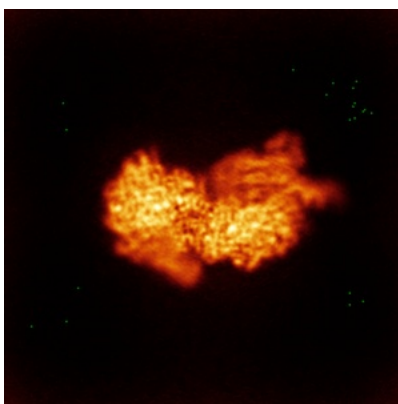


Z

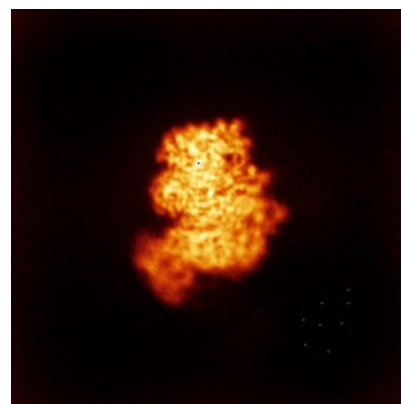
6.4.2 Raw map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.2. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

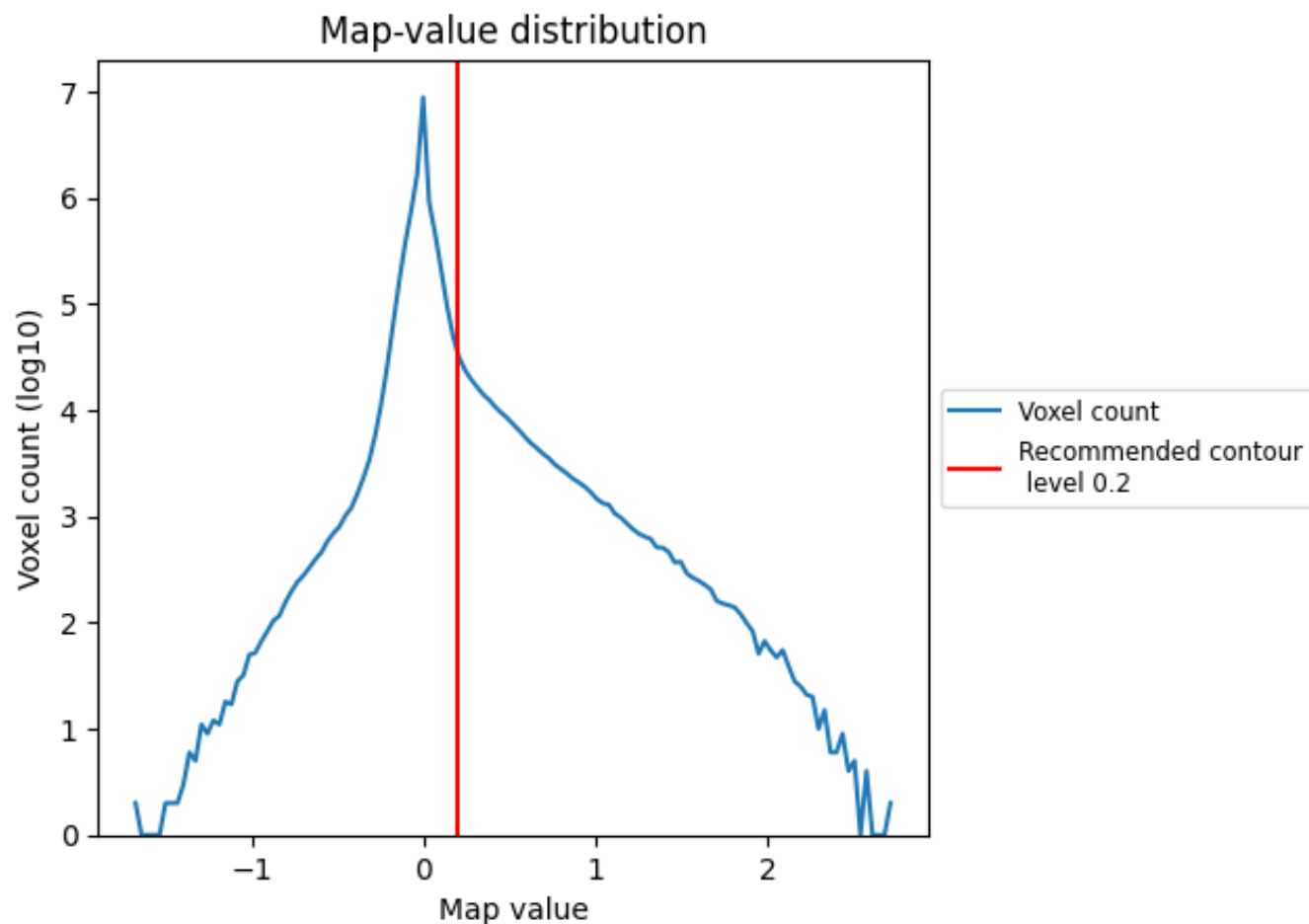
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

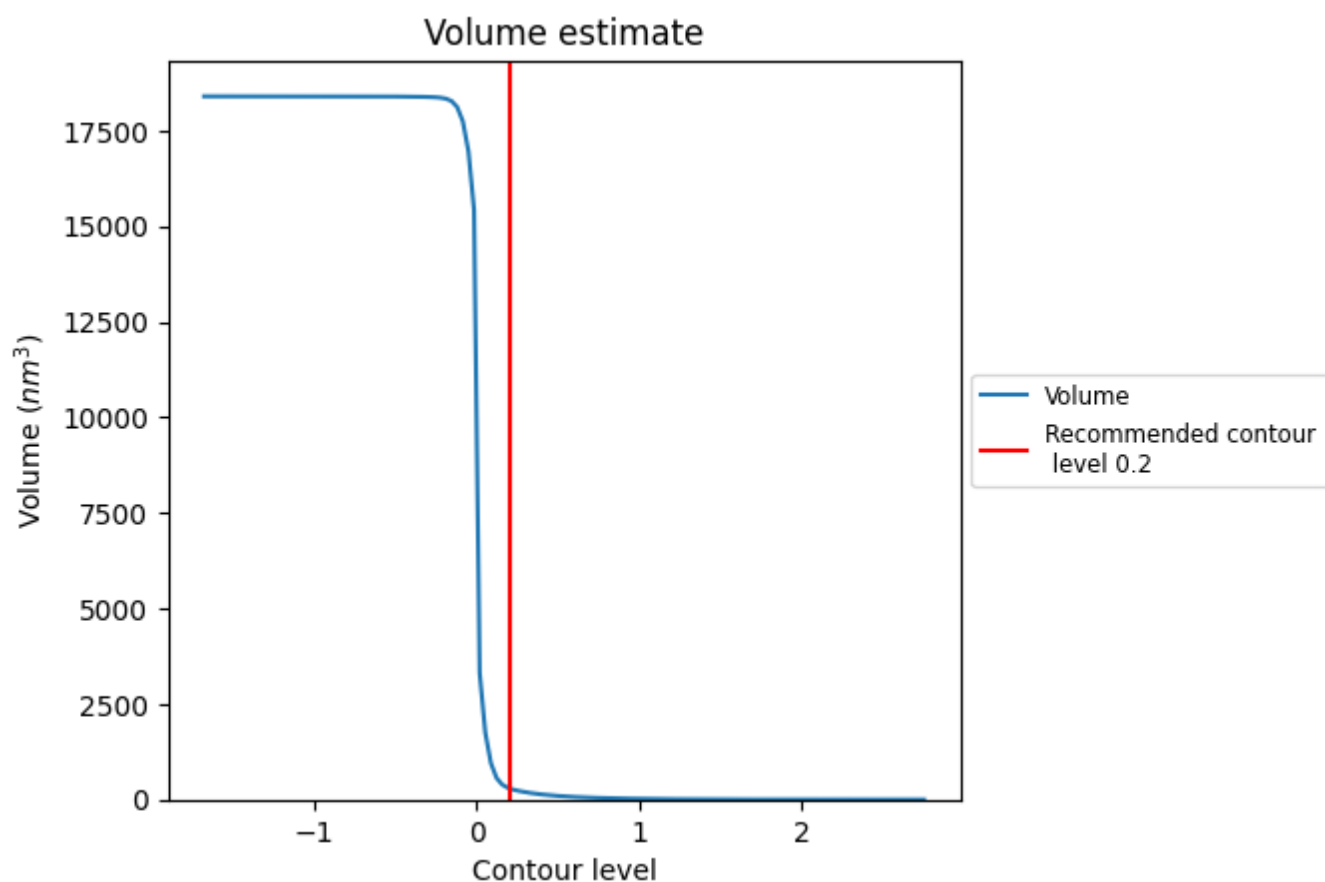
This section contains the results of statistical analysis of the map.

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.

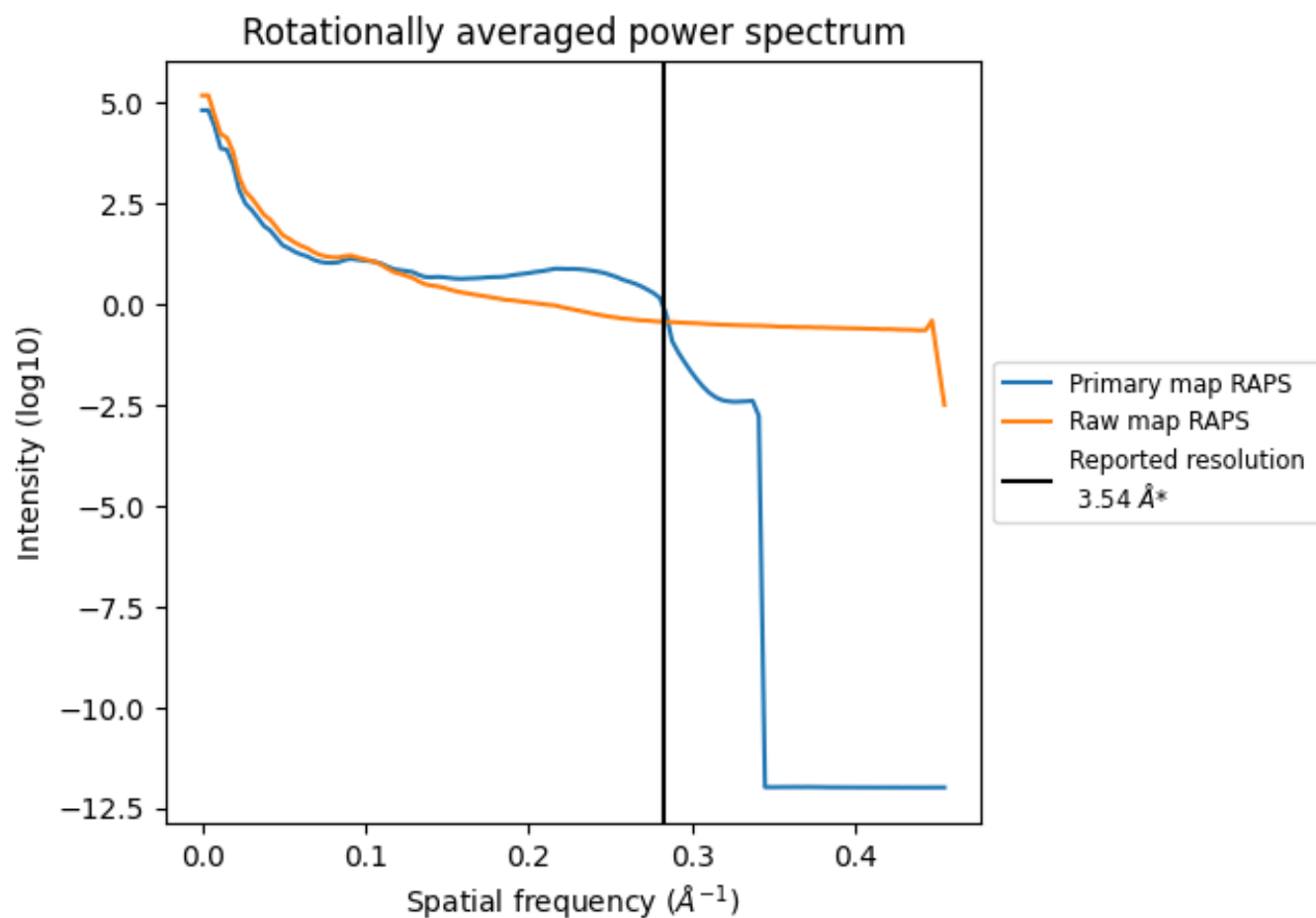
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 295 nm³; this corresponds to an approximate mass of 266 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 ⓘ

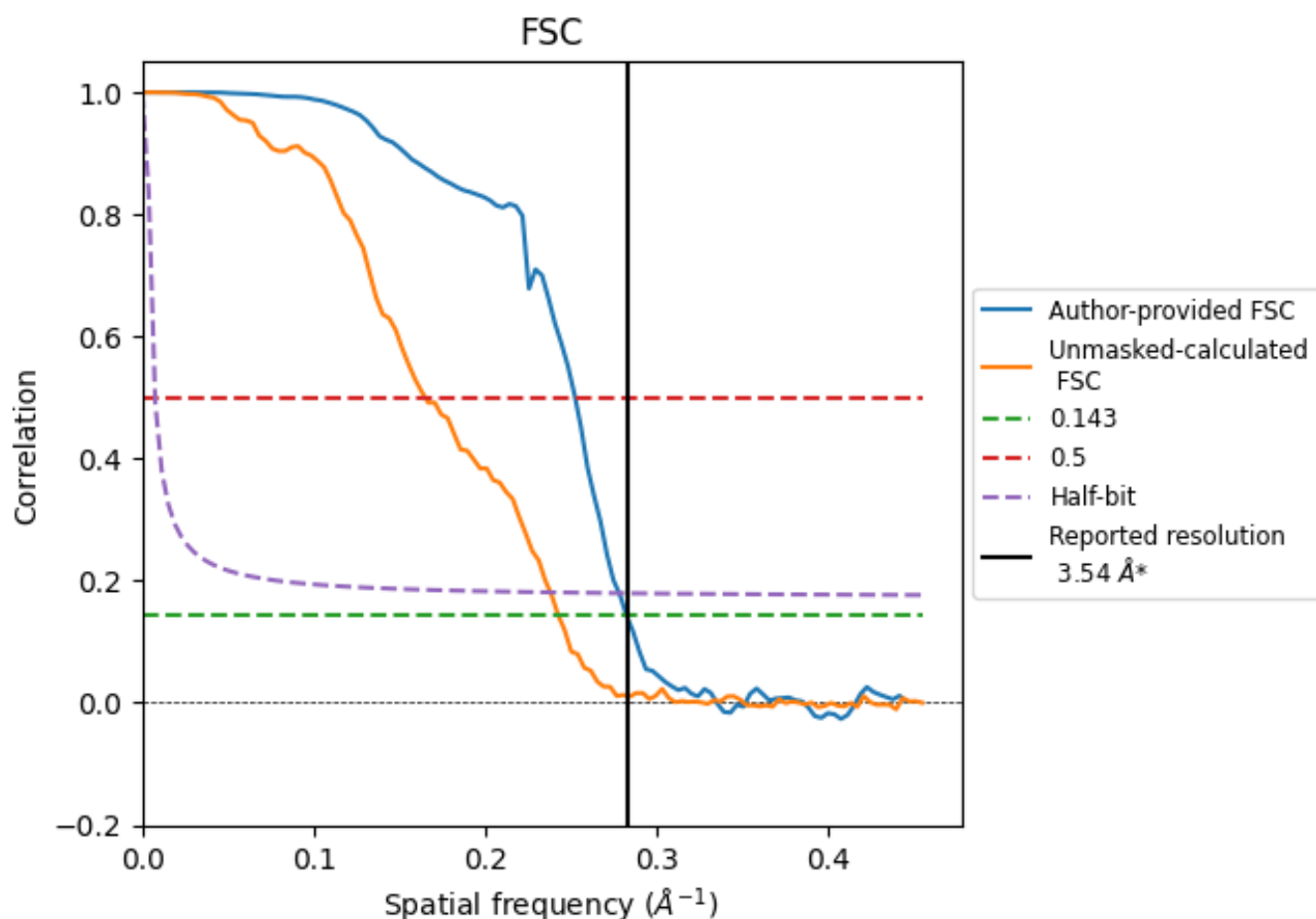


*Reported resolution corresponds to spatial frequency of $0.282 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.282 Å⁻¹

8.2 Resolution estimates [i](#)

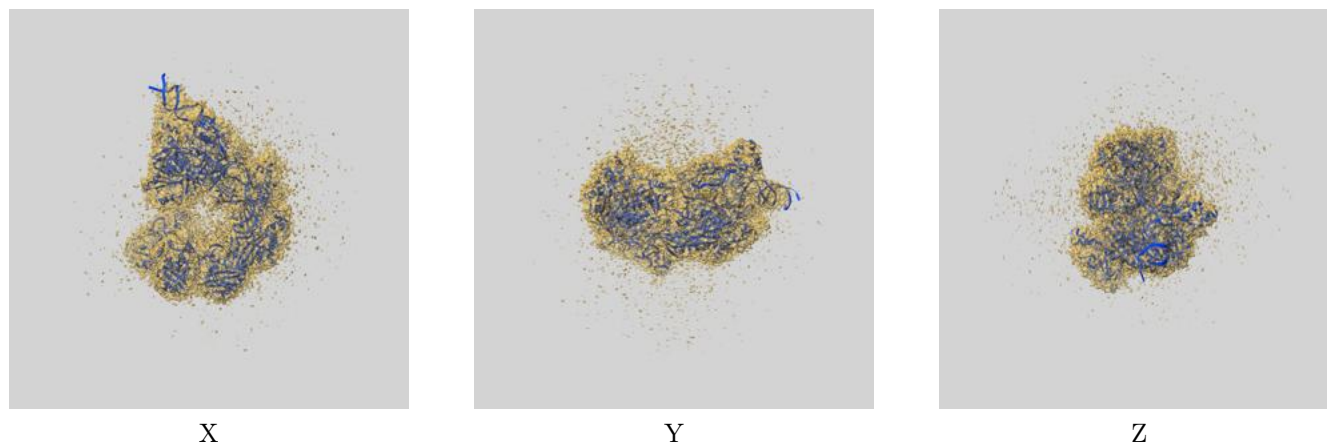
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.54	-	-
Author-provided FSC curve	3.54	3.96	3.60
Unmasked-calculated*	4.12	6.06	4.20

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.12 differs from the reported value 3.54 by more than 10 %

9 Map-model fit [i](#)

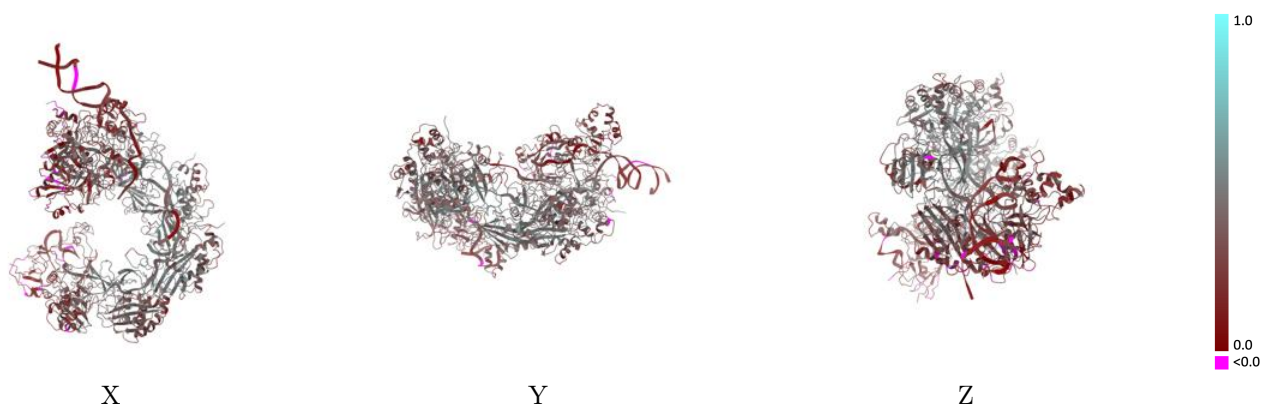
This section contains information regarding the fit between EMDB map EMD-36831 and PDB model 8K28. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.2 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](#)



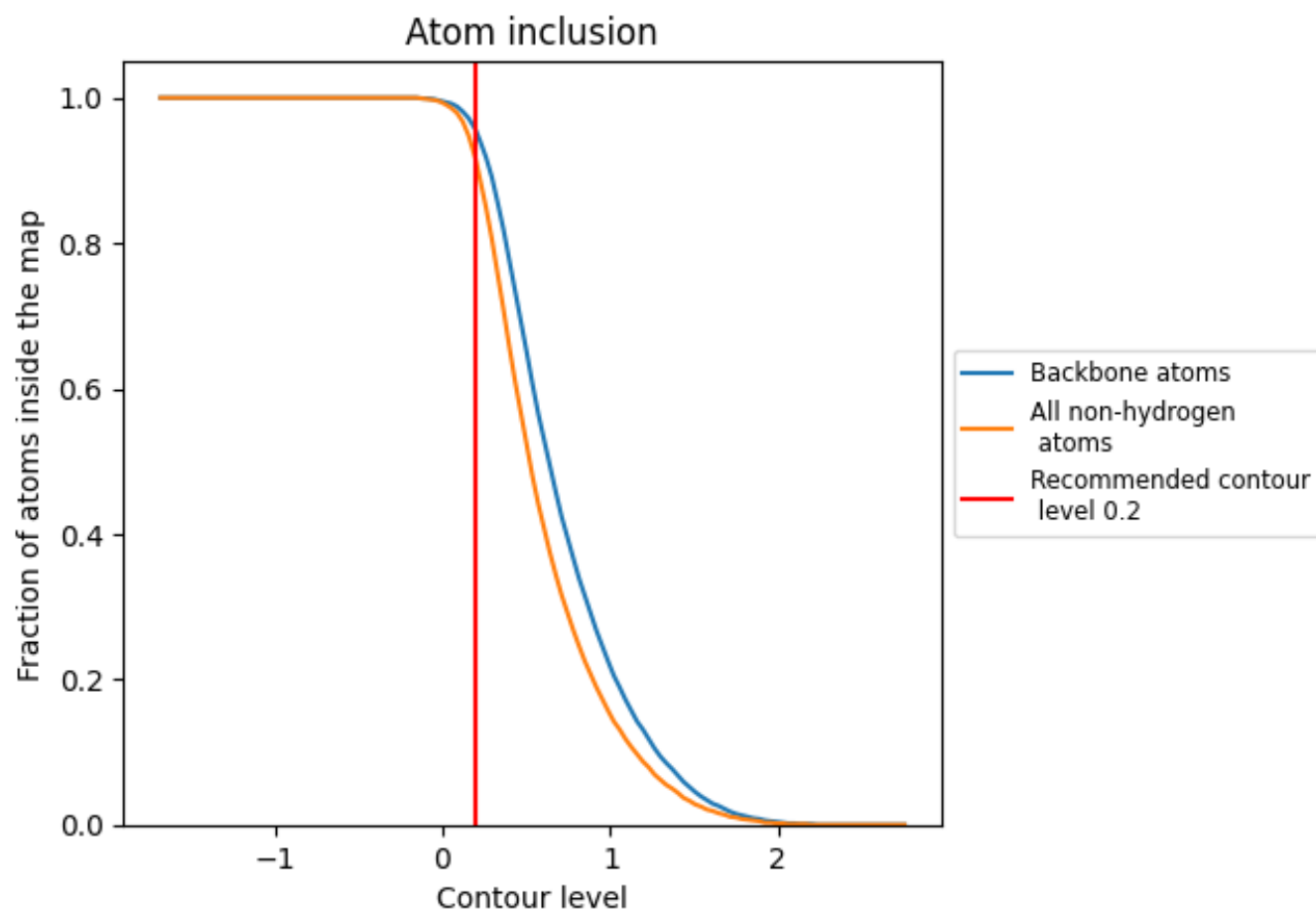
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 91% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9140	0.3550
A	0.8740	0.2840
B	0.8890	0.2860
C	0.9140	0.3480
D	0.9470	0.4270
E	0.9580	0.4460
F	0.9560	0.4460
G	0.9520	0.4110
H	0.8990	0.3260
I	0.8540	0.2030
P	0.9530	0.3340
Q	0.8030	0.2660
R	0.6850	0.1600

1.0

0.0

<0.0