



## Full wwPDB EM Validation Report ⓘ

Nov 13, 2022 – 05:14 PM EST

PDB ID : 7JZX  
EMDB ID : EMD-22583  
Title : Cryo-EM structure of CRISPR-Cas surveillance complex with AcrIF7  
Authors : Chang, L.; Li, Z.; Gabel, C.  
Deposited on : 2020-09-02  
Resolution : 3.40 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

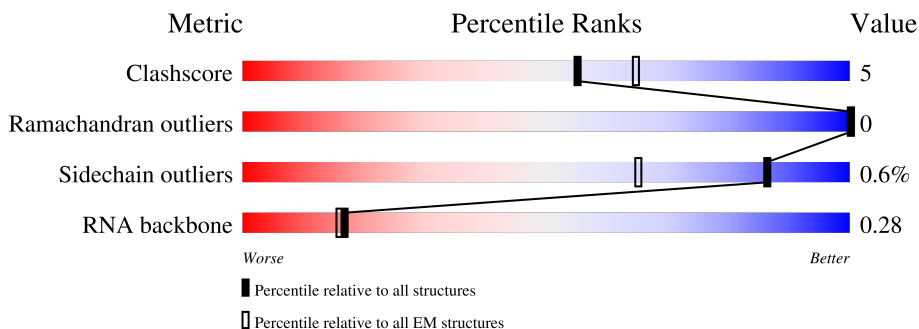
EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

# 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 3.40 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.






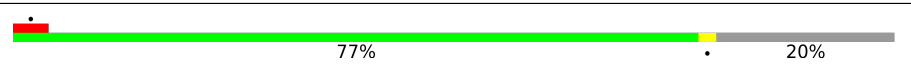
Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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  $\geq 3$ , 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	C	187	
2	B	327	
3	D	342	
3	E	342	
3	F	342	
3	G	342	
3	H	342	

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Mol	Chain	Length	Quality of chain
3	I	342	 <p>85% 13%</p>
4	A	434	 <p>14% 88% 10%</p>
5	M	61	 <p>43% 39% 18%</p>
6	J	83	 <p>77% 20%</p>

## 2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 23948 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 CRISPR-associated endonuclease Cas6/Csy4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	C	187	1448	922	264	257	5	1	0

- Molecule 2 is a protein called Type I-F CRISPR-associated protein Csy2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	306	2411	1529	445	432	5	0	0

- Molecule 3 is a protein called CRISPR type I-F/YPEST-associated protein Csy3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	E	335	2592	1626	471	493	2	0	0
3	D	298	2307	1453	415	437	2	0	0
3	F	335	2592	1626	471	493	2	0	0
3	G	335	2592	1626	471	493	2	0	0
3	H	330	2554	1604	464	484	2	0	0
3	I	335	2592	1626	471	493	2	0	0

- Molecule 4 is a protein called CRISPR-associated protein Csy1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	A	427	3079	1930	582	564	3	0	0

- Molecule 5 is a RNA chain called RNA (61-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
5	M	61	1273	569	223	421	60	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	41	A	G	conflict	GB 313291946
M	53	A	G	conflict	GB 313291946

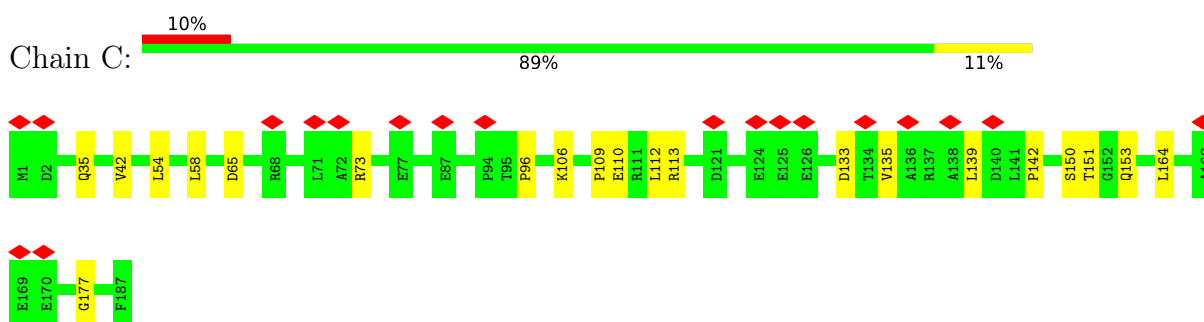
- Molecule 6 is a protein called AcrF7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	J	66	508	321	77	109	1	0	0

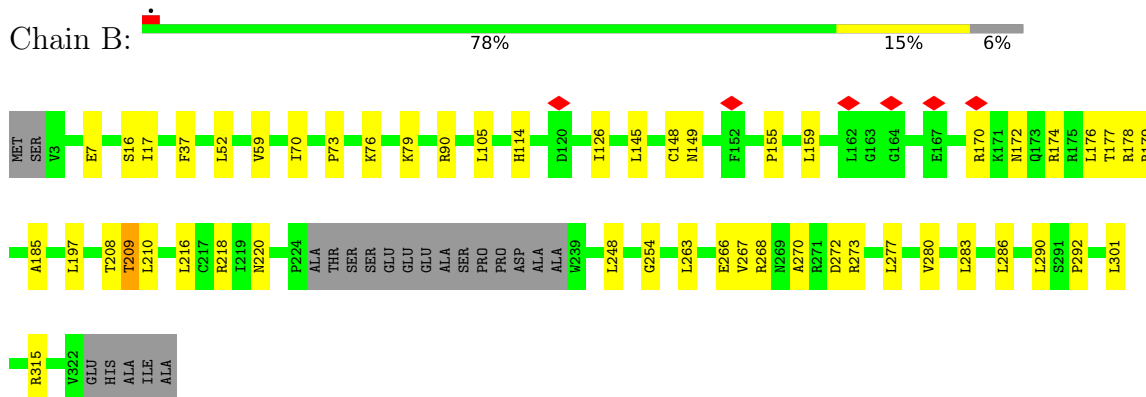
### 3 Residue-property plots [i](#)

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.

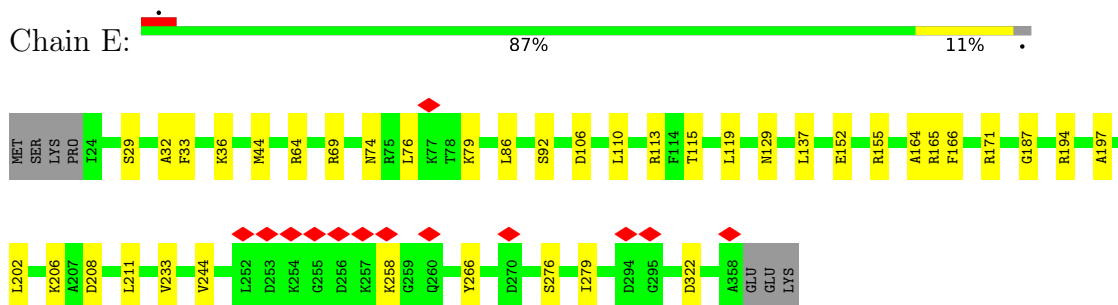
- Molecule 1: CRISPR-associated endonuclease Cas6/Csy4




- Molecule 2: Type I-F CRISPR-associated protein Csy2

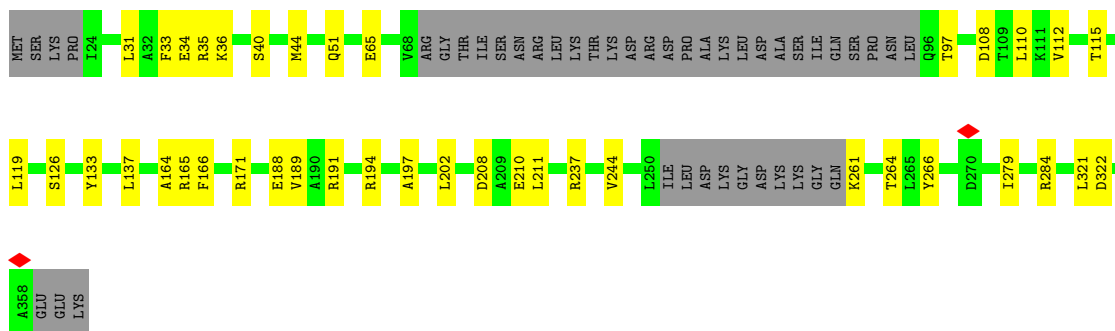


- Molecule 3: CRISPR type I-F/YPEST-associated protein Csy3




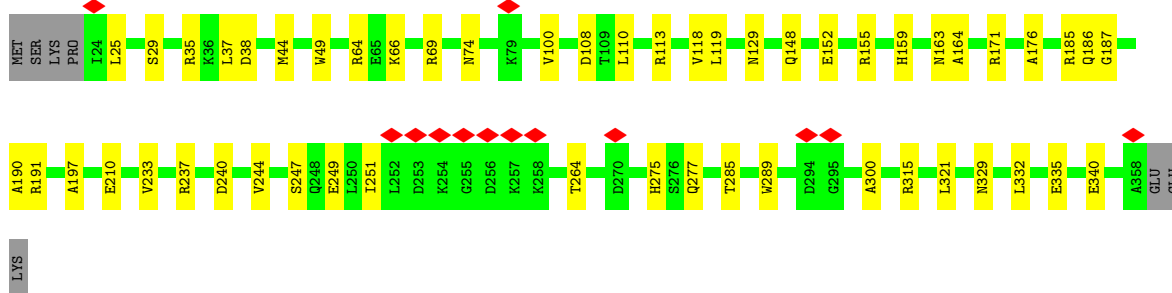
- Molecule 3: CRISPR type I-F/YPEST-associated protein Csy3

Chain D:  75% 12% 13%




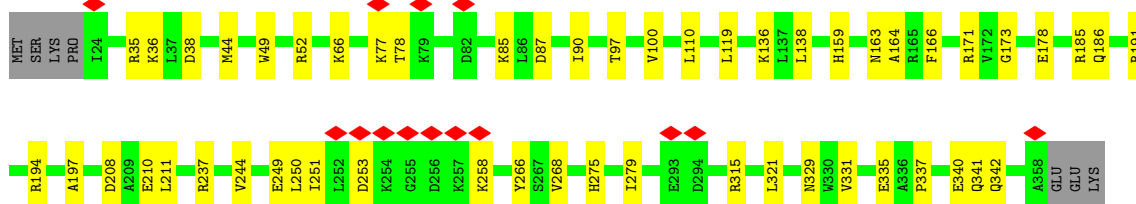
- Molecule 3: CRISPR type I-F/YPEST-associated protein Csy3

Chain F:  83% 15%




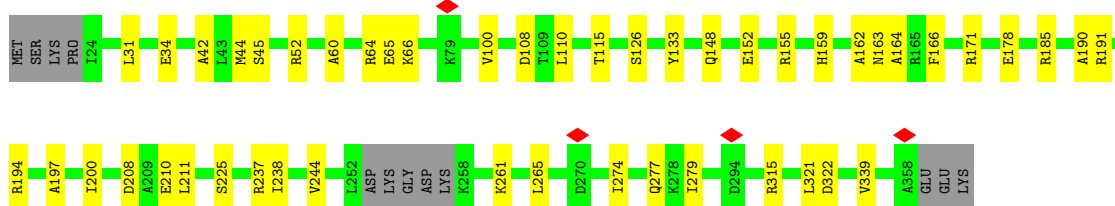
- Molecule 3: CRISPR type I-F/YPEST-associated protein Csy3

Chain G:  82% 15%

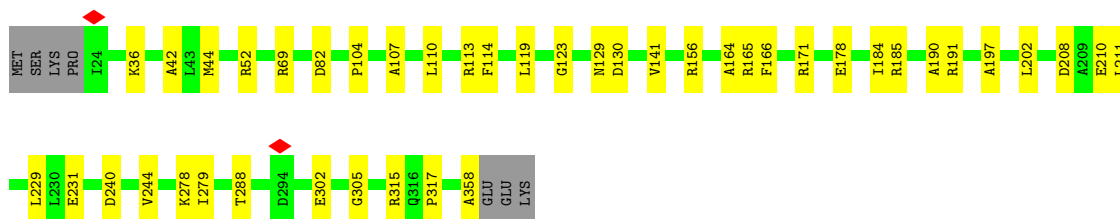
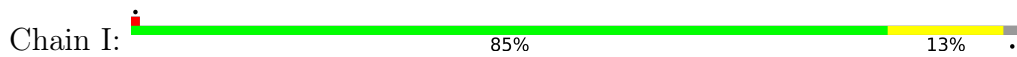


- Molecule 3: CRISPR type I-F/YPEST-associated protein Csy3

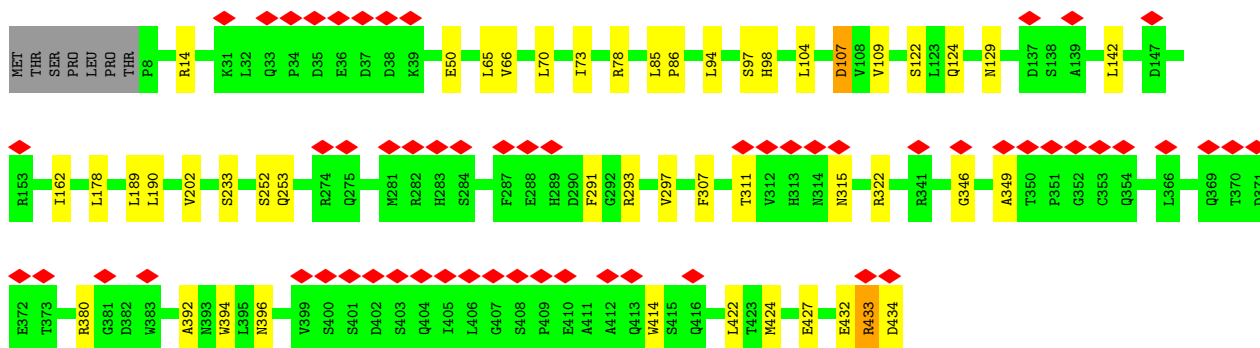
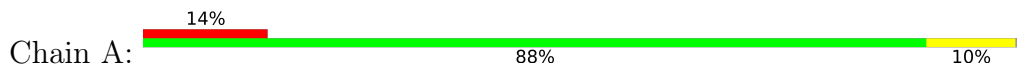
Chain H:  82% 14%



- Molecule 3: CRISPR type I-F/YPEST-associated protein Csy3



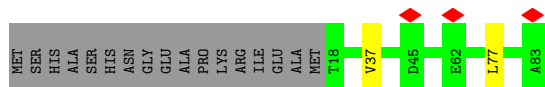
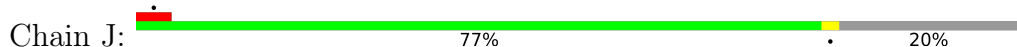
• Molecule 4: CRISPR-associated protein Csy1



• Molecule 5: RNA (61-MER)



• Molecule 6: AcrF7





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	502177	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	54	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.217	Depositor
Minimum map value	-0.027	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.0264	Depositor
Map size ( $\text{\AA}$ )	270.9, 270.9, 270.9	wwPDB
Map dimensions	258, 258, 258	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.05, 1.05, 1.05	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	C	0.36	0/1490	0.49	0/2024
2	B	0.45	0/2471	0.58	0/3365
3	D	0.44	0/2350	0.56	0/3190
3	E	0.44	0/2639	0.55	0/3579
3	F	0.43	0/2639	0.55	0/3579
3	G	0.41	0/2639	0.56	0/3579
3	H	0.47	0/2600	0.56	0/3527
3	I	0.48	0/2639	0.56	0/3579
4	A	0.37	0/3143	0.54	0/4286
5	M	0.72	1/1421 (0.1%)	0.92	0/2211
6	J	0.36	0/520	0.49	0/708
All	All	0.45	1/24551 (0.0%)	0.58	0/33627

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	M	36	A	O3'-P	5.44	1.67	1.61

There are no bond angle outliers.

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	C	1448	0	1394	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2411	0	2408	39	0
3	D	2307	0	2275	25	0
3	E	2592	0	2581	28	0
3	F	2592	0	2581	37	0
3	G	2592	0	2581	34	0
3	H	2554	0	2543	33	0
3	I	2592	0	2581	28	0
4	A	3079	0	2757	28	0
5	M	1273	0	646	19	0
6	J	508	0	454	1	0
All	All	23948	0	22801	247	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (247) 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)
2:B:185:ALA:HB1	2:B:301:LEU:HD11	1.20	1.16
2:B:185:ALA:CB	2:B:301:LEU:HD11	1.90	1.01
3:G:250:LEU:HA	5:M:21:U:O4	1.72	0.90
2:B:185:ALA:HB1	2:B:301:LEU:CD1	2.08	0.81
5:M:37:C:O2	5:M:37:C:H5 <sup>''</sup>	1.80	0.81
3:D:35:ARG:HD3	5:M:37:C:OP1	1.79	0.81
3:H:185:ARG:NH2	3:H:190:ALA:HB1	2.12	0.64
3:I:208:ASP:HB3	3:I:211:LEU:HB2	1.80	0.64
3:G:208:ASP:HB3	3:G:211:LEU:HB2	1.80	0.64
3:G:250:LEU:HA	5:M:21:U:C4	2.33	0.63
2:B:148:CYS:HB2	3:I:184:ILE:HG13	1.78	0.63
3:G:36:LYS:HG3	3:G:119:LEU:HB2	1.80	0.62
3:E:36:LYS:HG3	3:E:119:LEU:HB2	1.80	0.62
3:G:52:ARG:NH1	3:G:178:GLU:OE1	2.34	0.61
3:D:171:ARG:NH2	3:D:197:ALA:O	2.34	0.60
2:B:301:LEU:HD23	2:B:301:LEU:O	2.02	0.59
3:H:148:GLN:HE22	3:H:339:VAL:HG13	1.68	0.59
2:B:268:ARG:NH1	3:I:358:ALA:O	2.36	0.58
3:E:165:ARG:NH2	3:E:202:LEU:O	2.36	0.58
2:B:185:ALA:CB	2:B:301:LEU:CD1	2.76	0.58
3:H:65:GLU:OE2	3:H:261:LYS:NZ	2.35	0.58
2:B:254:GLY:HA3	2:B:315:ARG:HA	1.86	0.57
3:D:284:ARG:NH2	5:M:38:C:OP2	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:86:PRO:HG2	4:A:98:HIS:HB3	1.87	0.57
3:H:108:ASP:OD2	3:H:237:ARG:NH1	2.38	0.57
3:D:191:ARG:NH2	3:D:210:GLU:OE1	2.38	0.57
3:G:44:MET:HB2	3:G:110:LEU:HD11	1.85	0.57
2:B:70:ILE:HD12	2:B:79:LYS:HD3	1.86	0.57
3:E:208:ASP:HB2	3:E:211:LEU:HB2	1.86	0.57
3:H:171:ARG:NH2	3:H:197:ALA:O	2.38	0.57
3:G:340:GLU:HG3	3:G:341:GLN:HE21	1.71	0.56
2:B:155:PRO:O	2:B:179:ARG:NH2	2.37	0.56
2:B:149:ASN:ND2	3:I:231:GLU:OE1	2.39	0.56
2:B:273:ARG:NH1	3:I:130:ASP:OD1	2.39	0.55
3:I:110:LEU:HD13	3:I:244:VAL:HG11	1.89	0.55
3:E:266:TYR:OH	3:F:64:ARG:NH1	2.39	0.55
3:E:33:PHE:HZ	3:E:137:LEU:HD21	1.71	0.55
3:H:110:LEU:HD13	3:H:244:VAL:HG11	1.89	0.54
6:J:37:VAL:HB	6:J:77:LEU:HB3	1.88	0.54
2:B:267:VAL:HG21	2:B:277:LEU:HD22	1.89	0.54
3:I:52:ARG:NH1	3:I:178:GLU:OE1	2.40	0.54
2:B:7:GLU:H	2:B:114:HIS:HB3	1.73	0.54
3:F:191:ARG:NH2	3:F:210:GLU:OE1	2.39	0.54
3:G:66:LYS:HE2	3:G:100:VAL:HG11	1.90	0.54
3:H:185:ARG:HH21	3:H:190:ALA:HB1	1.72	0.54
3:I:165:ARG:NH2	3:I:202:LEU:O	2.37	0.54
2:B:177:THR:HG23	2:B:292:PRO:HB2	1.89	0.54
3:F:110:LEU:HD13	3:F:244:VAL:HG11	1.90	0.53
3:E:110:LEU:HD13	3:E:244:VAL:HG11	1.90	0.53
1:C:96:PRO:HB2	1:C:164:LEU:HD22	1.89	0.53
2:B:197:LEU:HD22	2:B:209:THR:HB	1.89	0.53
3:G:191:ARG:NH2	3:G:210:GLU:OE1	2.42	0.53
3:E:29:SER:OG	3:E:129:ASN:OD1	2.26	0.53
4:A:14:ARG:NH1	4:A:50:GLU:OE1	2.41	0.53
4:A:307:PHE:O	4:A:322:ARG:NH1	2.42	0.53
3:D:322:ASP:OD1	3:D:322:ASP:N	2.34	0.53
2:B:73:PRO:HG2	2:B:76:LYS:HB2	1.91	0.53
3:D:165:ARG:NH2	3:D:202:LEU:O	2.40	0.53
3:F:277:GLN:NE2	5:M:27:U:OP2	2.35	0.52
4:A:252:SER:OG	4:A:253:GLN:N	2.43	0.52
3:H:238:ILE:HD12	3:H:244:VAL:HG12	1.90	0.52
2:B:248:LEU:HD22	2:B:286:LEU:HD21	1.92	0.52
2:B:270:ALA:O	3:I:129:ASN:ND2	2.37	0.52
3:D:33:PHE:HZ	3:D:137:LEU:HD21	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:191:ARG:NH2	3:I:210:GLU:OE1	2.41	0.52
3:F:66:LYS:HE2	3:F:100:VAL:HG11	1.92	0.52
3:I:123:GLY:HA3	3:I:141:VAL:HG11	1.91	0.52
3:E:64:ARG:NH1	3:D:266:TYR:OH	2.42	0.51
3:F:113:ARG:HG3	3:F:233:VAL:HG22	1.92	0.51
3:E:44:MET:HB2	3:E:110:LEU:HD11	1.92	0.51
3:E:171:ARG:NH2	3:E:197:ALA:O	2.44	0.51
3:I:171:ARG:NH2	3:I:197:ALA:O	2.44	0.51
1:C:35:GLN:OE1	1:C:73:ARG:NH1	2.42	0.51
3:E:276:SER:OG	3:F:66:LYS:NZ	2.44	0.51
2:B:37:PHE:HB3	2:B:59:VAL:HG11	1.92	0.51
4:A:291:PHE:HB3	4:A:394:TRP:HE1	1.76	0.50
3:D:110:LEU:HD13	3:D:244:VAL:HG11	1.92	0.50
3:F:315:ARG:NH1	3:F:321:LEU:O	2.42	0.50
3:I:69:ARG:O	5:M:13:C:O2'	2.28	0.50
3:D:36:LYS:HG3	3:D:119:LEU:HB2	1.92	0.50
3:G:266:TYR:OH	3:H:64:ARG:NH1	2.44	0.50
4:A:311:THR:O	4:A:315:ASN:N	2.44	0.50
4:A:392:ALA:HB2	4:A:422:LEU:HD21	1.92	0.50
3:G:171:ARG:NH2	3:G:197:ALA:O	2.45	0.50
3:H:115:THR:HG21	3:I:240:ASP:HB2	1.94	0.50
3:H:208:ASP:HB2	3:H:211:LEU:HB2	1.94	0.50
5:M:21:U:H3'	5:M:21:U:O2	2.12	0.50
3:H:191:ARG:NH2	3:H:210:GLU:OE1	2.43	0.49
3:I:166:PHE:HB2	3:I:279:ILE:HD12	1.94	0.49
3:G:159:HIS:O	3:G:163:ASN:ND2	2.45	0.49
2:B:148:CYS:HB3	3:I:229:LEU:HD21	1.94	0.49
1:C:113:ARG:NH2	1:C:133:ASP:OD1	2.40	0.49
4:A:346:GLY:HA2	4:A:349:ALA:HB2	1.95	0.49
1:C:106:LYS:HB3	1:C:112:LEU:HD11	1.93	0.49
2:B:159:LEU:HD13	2:B:176:LEU:HD22	1.95	0.49
3:E:113:ARG:HG3	3:E:233:VAL:HG22	1.95	0.49
3:E:115:THR:HG21	3:F:240:ASP:HB2	1.95	0.48
3:I:44:MET:HB2	3:I:110:LEU:HD11	1.95	0.48
3:E:69:ARG:O	5:M:37:C:O2'	2.31	0.48
1:C:109:PRO:HA	1:C:112:LEU:HD12	1.95	0.48
3:F:35:ARG:NE	3:F:38:ASP:OD1	2.46	0.48
3:E:258:LYS:NZ	5:M:39:U:O4	2.46	0.48
3:G:110:LEU:HD13	3:G:244:VAL:HG11	1.95	0.48
4:A:178:LEU:N	4:A:189:LEU:O	2.47	0.48
2:B:210:LEU:HD21	4:A:202:VAL:HG22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:65:LEU:N	4:A:162:ILE:O	2.44	0.47
3:F:69:ARG:NH1	5:M:33:U:O2'	2.48	0.47
3:E:32:ALA:HB1	5:M:29:C:H1'	1.95	0.47
3:I:278:LYS:NZ	5:M:10:U:OP1	2.41	0.47
3:G:315:ARG:NH1	3:G:321:LEU:O	2.42	0.47
4:A:97:SER:O	4:A:97:SER:OG	2.32	0.47
3:D:65:GLU:OE2	3:D:97:THR:OG1	2.31	0.47
3:F:37:LEU:HD23	3:F:118:VAL:HG22	1.96	0.47
3:F:249:GLU:HB2	3:F:264:THR:HG23	1.97	0.47
3:G:337:PRO:HG2	3:G:342:GLN:HB3	1.97	0.47
3:I:42:ALA:HA	3:I:114:PHE:HB3	1.96	0.47
4:A:293:ARG:HD3	4:A:297:VAL:HG21	1.97	0.47
2:B:16:SER:HB2	2:B:145:LEU:HB2	1.96	0.46
3:F:187:GLY:HA2	3:G:237:ARG:HH22	1.80	0.46
4:A:380:ARG:O	4:A:433:ARG:NH1	2.48	0.46
3:F:186:GLN:O	3:G:237:ARG:NH2	2.46	0.46
2:B:148:CYS:O	3:I:113:ARG:NH2	2.46	0.46
3:H:277:GLN:NE2	5:M:15:G:OP1	2.49	0.46
3:H:322:ASP:N	3:H:322:ASP:OD1	2.45	0.46
2:B:208:THR:OG1	2:B:209:THR:N	2.49	0.46
3:D:31:LEU:HD11	3:D:133:TYR:HE2	1.81	0.46
3:H:66:LYS:HE2	3:H:100:VAL:HG11	1.97	0.46
3:F:159:HIS:O	3:F:163:ASN:ND2	2.49	0.46
3:F:171:ARG:NH2	3:F:197:ALA:O	2.48	0.46
2:B:178:ARG:NH2	4:A:434:ASP:OXT	2.42	0.46
3:H:31:LEU:HD11	3:H:133:TYR:HE2	1.80	0.46
3:D:34:GLU:HG3	3:D:126:SER:HB3	1.98	0.46
3:F:25:LEU:HD23	3:F:332:LEU:HA	1.96	0.46
3:F:185:ARG:HB2	3:F:190:ALA:HB2	1.98	0.46
3:I:104:PRO:HD2	3:I:107:ALA:HB3	1.98	0.46
1:C:150:SER:OG	1:C:151:THR:N	2.49	0.45
3:E:164:ALA:H	3:E:197:ALA:HB1	1.81	0.45
3:F:49:TRP:HB2	3:F:237:ARG:HD3	1.98	0.45
3:E:74:ASN:N	3:E:74:ASN:OD1	2.47	0.45
3:G:49:TRP:HB2	3:G:237:ARG:HD3	1.98	0.45
3:D:208:ASP:HB3	3:D:211:LEU:HB2	1.98	0.45
2:B:52:LEU:HD22	2:B:126:ILE:HG22	1.99	0.45
3:F:108:ASP:OD1	3:F:108:ASP:N	2.48	0.45
4:A:291:PHE:HB2	4:A:293:ARG:HG2	1.98	0.45
1:C:54:LEU:HG	1:C:142:PRO:HB2	1.98	0.45
3:G:85:LYS:HB2	3:G:85:LYS:HE2	1.75	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:42:ALA:HB3	3:H:274:ILE:HB	1.97	0.45
3:F:152:GLU:OE2	3:F:155:ARG:NH2	2.50	0.45
3:F:44:MET:HB2	3:F:110:LEU:HD11	1.99	0.44
2:B:17:ILE:HB	2:B:105:LEU:HB2	1.99	0.44
3:F:285:THR:HA	3:F:300:ALA:HA	1.99	0.44
3:G:35:ARG:NE	3:G:38:ASP:OD1	2.42	0.44
3:G:185:ARG:HG3	3:G:186:GLN:H	1.83	0.44
3:D:40:SER:HB3	3:D:115:THR:H	1.81	0.44
3:G:97:THR:O	3:G:97:THR:OG1	2.35	0.44
4:A:124:GLN:NE2	4:A:129:ASN:OD1	2.50	0.44
3:E:92:SER:O	3:E:92:SER:OG	2.31	0.44
2:B:218:ARG:HH11	4:A:233:SER:HB3	1.82	0.44
3:D:44:MET:HB3	3:D:112:VAL:HG22	1.99	0.44
3:F:251:ILE:O	3:G:258:LYS:NZ	2.51	0.44
3:H:159:HIS:O	3:H:163:ASN:ND2	2.51	0.44
3:I:305:GLY:H	3:I:315:ARG:HB2	1.83	0.44
3:G:136:LYS:NZ	3:G:331:VAL:O	2.43	0.44
3:H:315:ARG:NH1	3:H:321:LEU:O	2.47	0.44
3:I:302:GLU:HB2	3:I:305:GLY:HA2	2.00	0.44
4:A:85:LEU:HD21	4:A:104:LEU:HD11	1.99	0.44
2:B:280:VAL:HG11	4:A:190:LEU:HD22	1.99	0.43
3:E:187:GLY:HA2	3:F:237:ARG:HH22	1.82	0.43
4:A:70:LEU:HD12	4:A:73:ILE:HD12	2.00	0.43
3:D:188:GLU:HG2	3:D:189:VAL:H	1.84	0.43
3:H:185:ARG:HH21	3:H:190:ALA:CB	2.31	0.43
1:C:65:ASP:OD2	1:C:65:ASP:N	2.47	0.43
3:F:247:SER:HG	3:F:275:HIS:CG	2.37	0.43
3:G:87:ASP:HA	3:G:90:ILE:HG12	2.00	0.43
3:G:164:ALA:H	3:G:197:ALA:HB1	1.82	0.43
3:I:164:ALA:H	3:I:197:ALA:HB1	1.84	0.43
3:D:44:MET:HB2	3:D:110:LEU:HD11	2.00	0.43
3:G:194:ARG:HE	3:G:194:ARG:HB2	1.52	0.43
3:H:52:ARG:NH1	3:H:178:GLU:OE1	2.51	0.43
4:A:122:SER:O	4:A:122:SER:OG	2.34	0.43
2:B:263:LEU:HB2	2:B:266:GLU:HB2	2.00	0.43
3:E:194:ARG:HE	3:E:194:ARG:HB2	1.52	0.43
2:B:170:ARG:HH11	2:B:174:ARG:HH11	1.67	0.43
3:H:225:SER:O	3:H:225:SER:OG	2.30	0.43
3:I:36:LYS:HG3	3:I:119:LEU:HB2	2.00	0.43
3:E:322:ASP:OD1	3:E:322:ASP:N	2.44	0.43
3:G:138:LEU:HD23	3:G:138:LEU:HA	1.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:329:ASN:O	3:G:335:GLU:N	2.51	0.43
3:E:76:LEU:HD11	3:E:86:LEU:HD13	2.01	0.43
3:G:251:ILE:HG22	3:G:253:ASP:H	1.84	0.43
3:H:152:GLU:OE2	3:H:155:ARG:NH2	2.51	0.43
3:F:329:ASN:O	3:F:335:GLU:N	2.52	0.43
1:C:42:VAL:HG13	1:C:177:GLY:HA3	2.01	0.43
3:F:164:ALA:H	3:F:197:ALA:HB1	1.84	0.43
3:D:261:LYS:HE2	3:D:264:THR:HG22	2.00	0.42
3:F:176:ALA:H	3:F:237:ARG:HB2	1.83	0.42
1:C:139:LEU:N	5:M:45:A:N1	2.57	0.42
3:G:249:GLU:OE2	3:G:275:HIS:NE2	2.52	0.42
2:B:185:ALA:N	2:B:290:LEU:O	2.53	0.42
3:D:194:ARG:HE	3:D:194:ARG:HB2	1.61	0.42
1:C:110:GLU:HA	1:C:113:ARG:HD2	2.01	0.42
2:B:90:ARG:HD3	3:I:317:PRO:HG3	2.02	0.42
3:I:156:ARG:HH12	3:I:288:THR:HG22	1.85	0.42
2:B:216:LEU:HD22	2:B:248:LEU:HD13	2.02	0.42
3:E:166:PHE:HB2	3:E:279:ILE:HD12	2.01	0.42
3:D:164:ALA:H	3:D:197:ALA:HB1	1.84	0.42
3:F:289:TRP:CG	3:F:340:GLU:HB3	2.54	0.42
1:C:42:VAL:HB	1:C:58:LEU:HD11	2.01	0.42
2:B:283:LEU:HB3	4:A:94:LEU:HD12	2.01	0.42
3:H:45:SER:HA	3:H:60:ALA:HA	2.01	0.42
3:H:162:ALA:HB1	3:H:200:ILE:HD13	2.02	0.42
4:A:396:ASN:HD22	4:A:414:TRP:HH2	1.67	0.41
4:A:424:MET:HA	4:A:427:GLU:HG2	2.02	0.41
3:E:106:ASP:OD1	3:E:106:ASP:N	2.53	0.41
3:E:206:LYS:HE3	3:E:206:LYS:HB2	1.74	0.41
3:F:119:LEU:HD22	3:G:173:GLY:HA3	2.03	0.41
3:F:237:ARG:HA	3:F:237:ARG:HD2	1.79	0.41
3:H:321:LEU:HD23	3:H:321:LEU:HA	1.91	0.41
3:F:74:ASN:OD1	3:F:74:ASN:N	2.50	0.41
3:G:166:PHE:HB2	3:G:279:ILE:HD12	2.02	0.41
1:C:150:SER:N	5:M:61:C:OP1	2.53	0.41
3:F:185:ARG:HA	3:F:185:ARG:HD3	1.81	0.41
3:F:29:SER:HB3	3:F:129:ASN:HD22	1.86	0.41
3:H:44:MET:HB2	3:H:110:LEU:HD11	2.02	0.41
3:I:185:ARG:HB2	3:I:190:ALA:HB2	2.03	0.41
1:C:153:GLN:OE1	5:M:46:C:O2'	2.37	0.41
3:E:152:GLU:OE2	3:E:155:ARG:NH2	2.54	0.41
3:D:321:LEU:HD23	3:D:321:LEU:HA	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:321:LEU:HD23	3:F:321:LEU:HA	1.92	0.41
3:H:164:ALA:H	3:H:197:ALA:HB1	1.85	0.41
1:C:135:VAL:HG21	5:M:54:U:H4'	2.02	0.41
3:H:108:ASP:OD1	3:H:108:ASP:N	2.54	0.41
4:A:66:VAL:HG21	4:A:109:VAL:HG13	2.02	0.41
4:A:107:ASP:OD1	4:A:252:SER:OG	2.39	0.41
3:E:79:LYS:HB3	3:E:79:LYS:HE2	1.88	0.41
3:D:166:PHE:HB2	3:D:279:ILE:HD12	2.02	0.41
2:B:220:ASN:OD1	2:B:220:ASN:N	2.53	0.40
3:H:34:GLU:HG2	3:H:126:SER:HB3	2.02	0.40
3:H:194:ARG:HE	3:H:194:ARG:HB2	1.57	0.40
4:A:432:GLU:HG3	4:A:433:ARG:HD3	2.03	0.40
2:B:272:ASP:OD2	2:B:272:ASP:N	2.51	0.40
3:D:108:ASP:OD1	3:D:108:ASP:N	2.51	0.40
3:D:188:GLU:OE2	3:D:188:GLU:N	2.48	0.40
3:H:166:PHE:HB2	3:H:279:ILE:HD12	2.03	0.40
5:M:41:A:H5'	5:M:42:U:H5''	2.04	0.40
3:G:237:ARG:HD2	3:G:237:ARG:HA	1.91	0.40
3:H:265:LEU:HD23	3:H:265:LEU:HA	1.96	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	C	186/187 (100%)	181 (97%)	5 (3%)	0	100	100
2	B	302/327 (92%)	279 (92%)	23 (8%)	0	100	100
3	D	292/342 (85%)	271 (93%)	21 (7%)	0	100	100
3	E	333/342 (97%)	307 (92%)	26 (8%)	0	100	100
3	F	333/342 (97%)	307 (92%)	26 (8%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	G	333/342 (97%)	310 (93%)	23 (7%)	0	100	100
3	H	326/342 (95%)	308 (94%)	18 (6%)	0	100	100
3	I	333/342 (97%)	311 (93%)	22 (7%)	0	100	100
4	A	425/434 (98%)	385 (91%)	40 (9%)	0	100	100
6	J	64/83 (77%)	63 (98%)	1 (2%)	0	100	100
All	All	2927/3083 (95%)	2722 (93%)	205 (7%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	C	146/160 (91%)	146 (100%)	0	100	100
2	B	255/270 (94%)	253 (99%)	2 (1%)	81	91
3	D	234/273 (86%)	232 (99%)	2 (1%)	78	90
3	E	266/273 (97%)	266 (100%)	0	100	100
3	F	266/273 (97%)	265 (100%)	1 (0%)	91	95
3	G	266/273 (97%)	263 (99%)	3 (1%)	73	86
3	H	262/273 (96%)	262 (100%)	0	100	100
3	I	266/273 (97%)	265 (100%)	1 (0%)	91	95
4	A	262/365 (72%)	258 (98%)	4 (2%)	65	82
6	J	53/66 (80%)	53 (100%)	0	100	100
All	All	2276/2499 (91%)	2263 (99%)	13 (1%)	86	94

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	172	ASN
2	B	209	THR

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Mol	Chain	Res	Type
3	D	51	GLN
3	D	237	ARG
3	F	148	GLN
3	G	77	LYS
3	G	78	THR
3	G	268	VAL
3	I	82	ASP
4	A	78	ARG
4	A	107	ASP
4	A	142	LEU
4	A	433	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	29	HIS
2	B	156	ASN
3	E	51	GLN
3	G	163	ASN
3	H	148	GLN
4	A	124	GLN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
5	M	59/61 (96%)	29 (49%)	1 (1%)

All (29) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	M	3	A
5	M	6	A
5	M	9	U
5	M	10	U
5	M	11	C
5	M	14	G
5	M	15	G
5	M	16	C
5	M	17	G
5	M	20	C

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Mol	Chain	Res	Type
5	M	21	U
5	M	22	U
5	M	23	G
5	M	24	A
5	M	27	U
5	M	28	C
5	M	30	G
5	M	33	U
5	M	34	C
5	M	35	U
5	M	36	A
5	M	37	C
5	M	39	U
5	M	41	A
5	M	42	U
5	M	43	U
5	M	48	G
5	M	54	U
5	M	55	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
5	M	16	C

#### 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 monosaccharides 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.

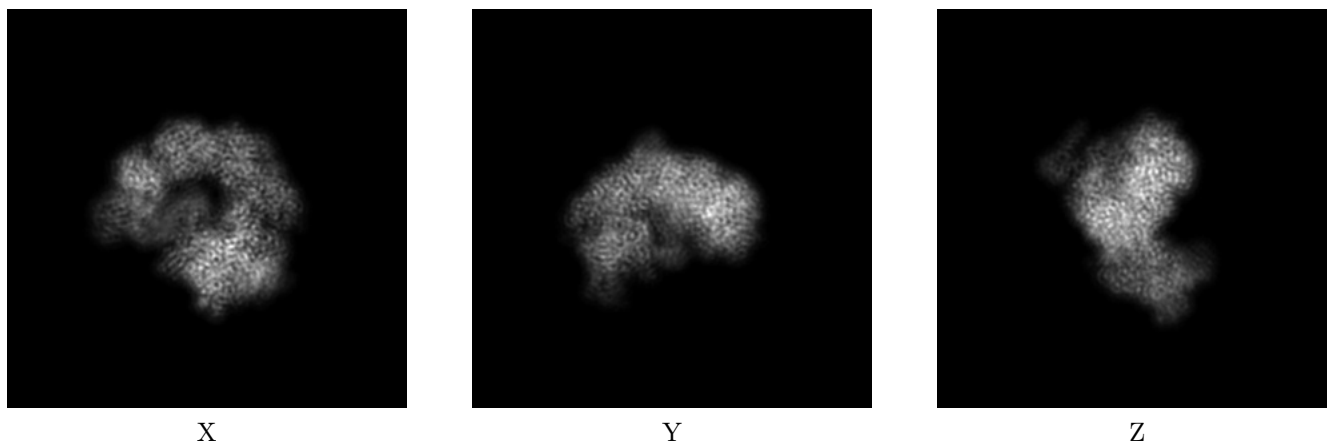
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-22583. 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](#)

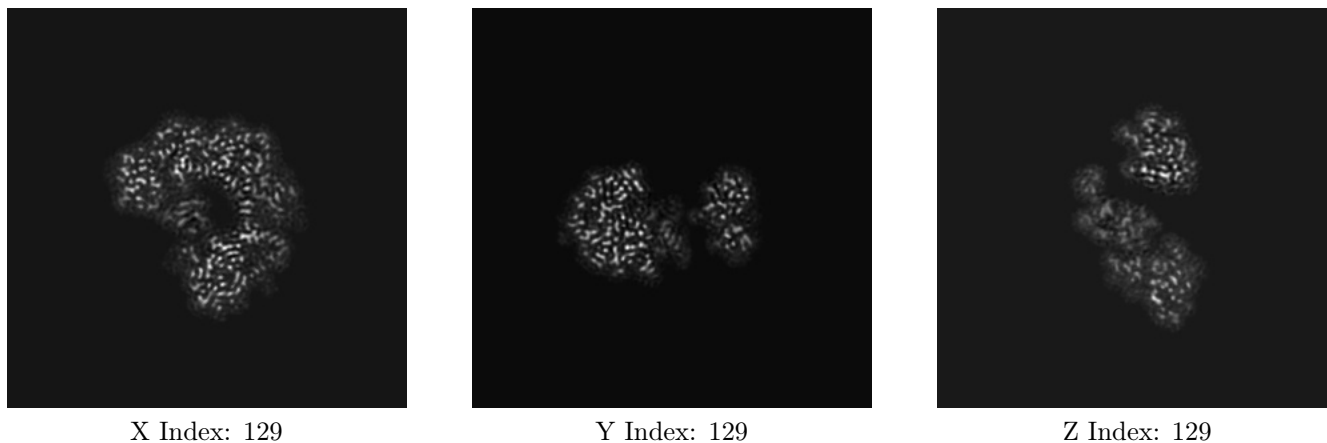
#### 6.1.1 Primary map



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

### 6.2 Central slices [i](#)

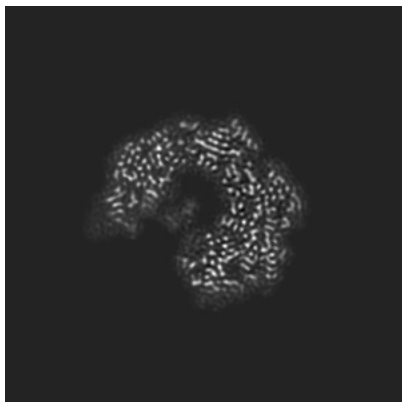
#### 6.2.1 Primary map



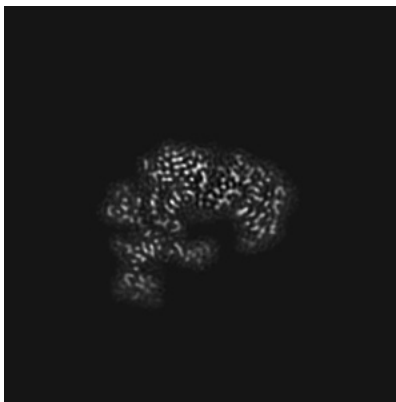
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

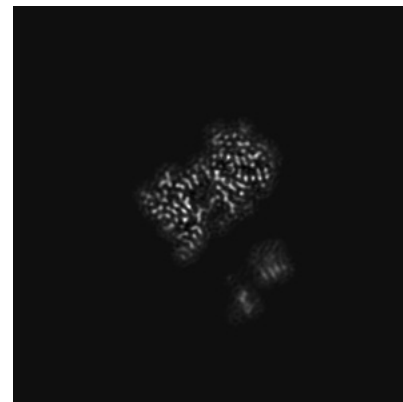
### 6.3.1 Primary map



X Index: 139



Y Index: 150



Z Index: 108

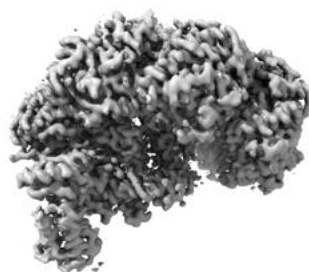
The images above show the largest variance slices of the map in three orthogonal directions.

## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0264. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

## 6.5 Mask visualisation

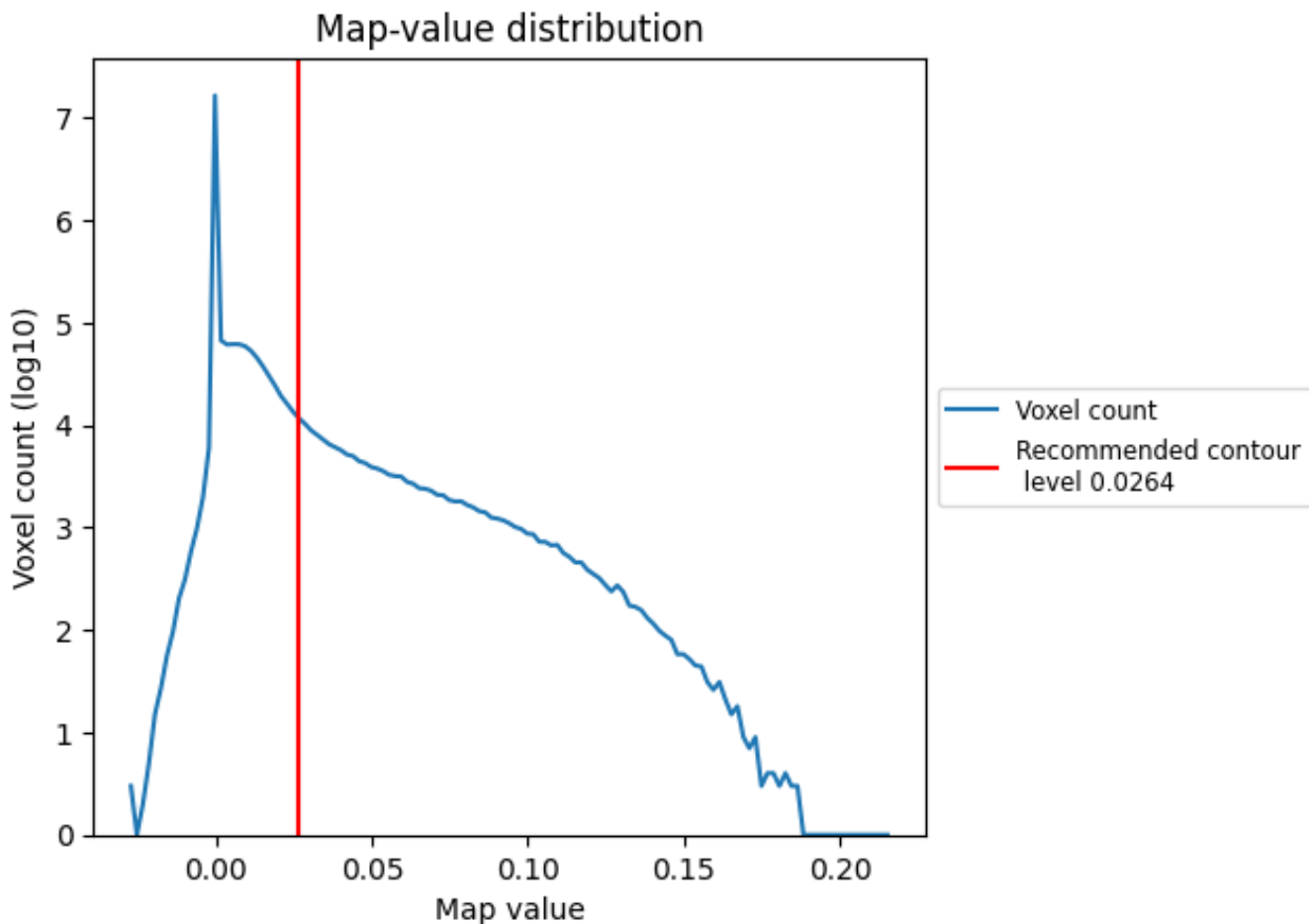
This section was not generated. No masks/segmentation were deposited.



## 7 Map analysis [i](#)

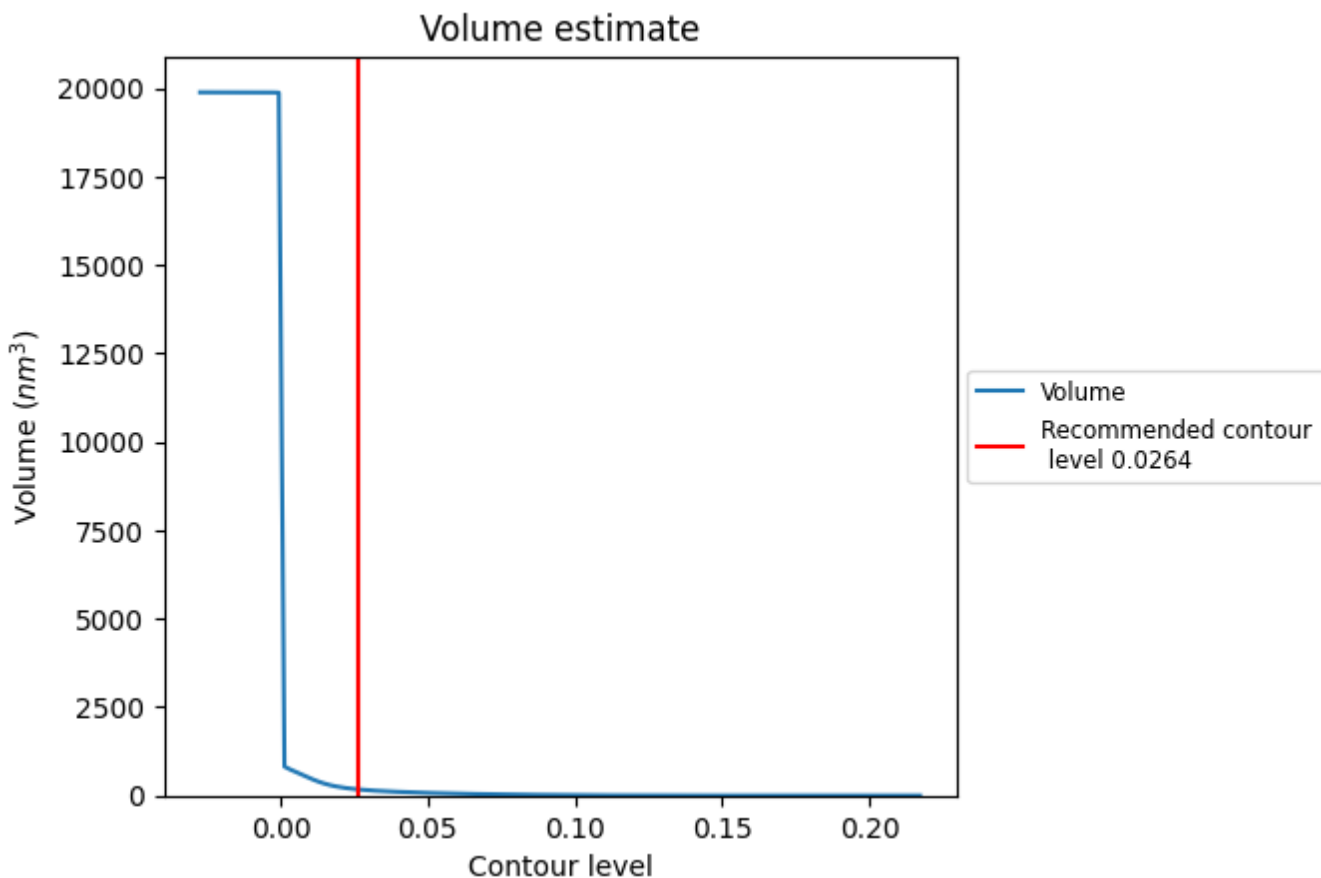
This section contains the results of statistical analysis of the map.

### 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.

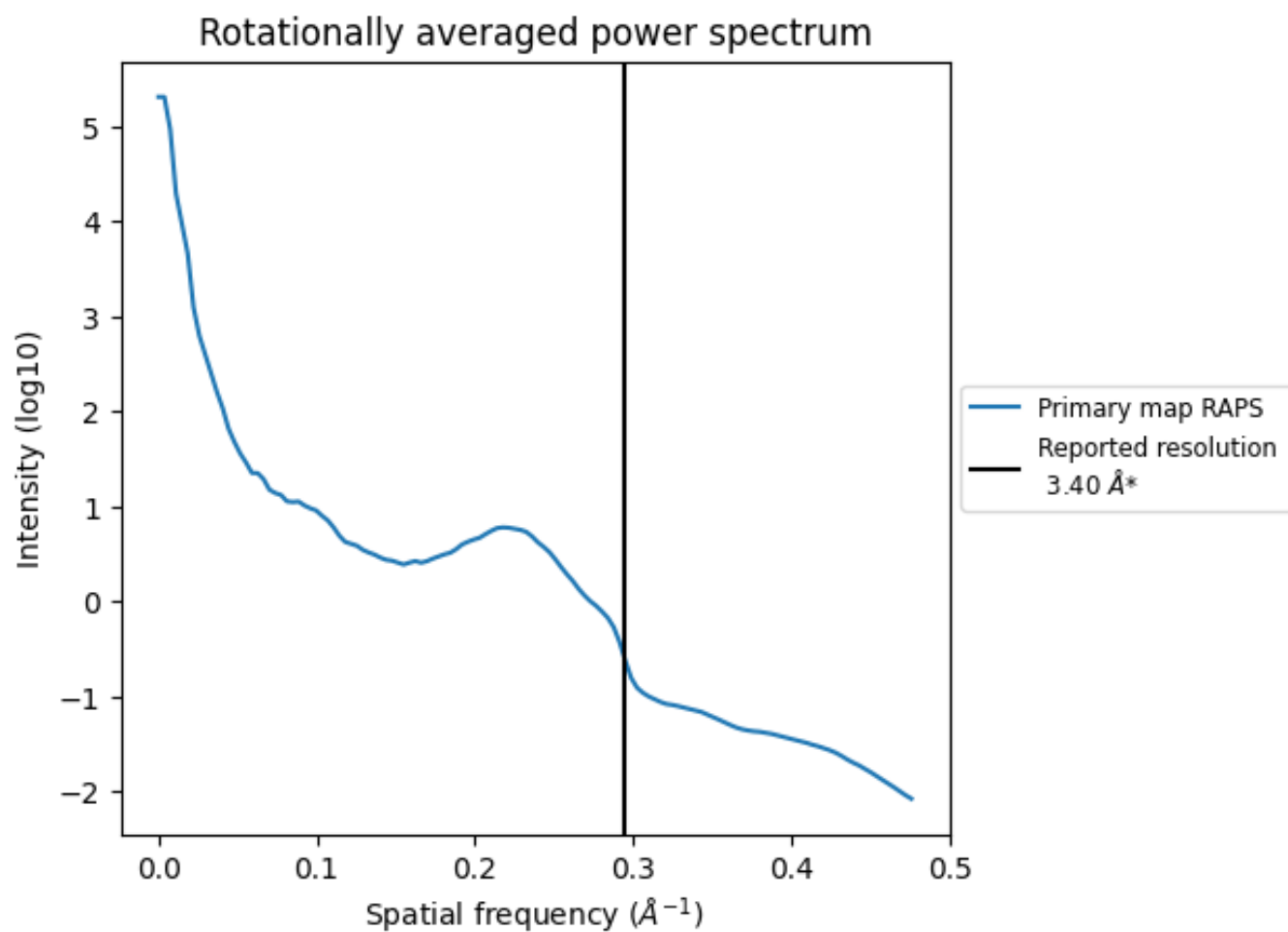
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 174  $\text{nm}^3$ ; this corresponds to an approximate mass of 157 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 [i](#)



\*Reported resolution corresponds to spatial frequency of 0.294 Å<sup>-1</sup>

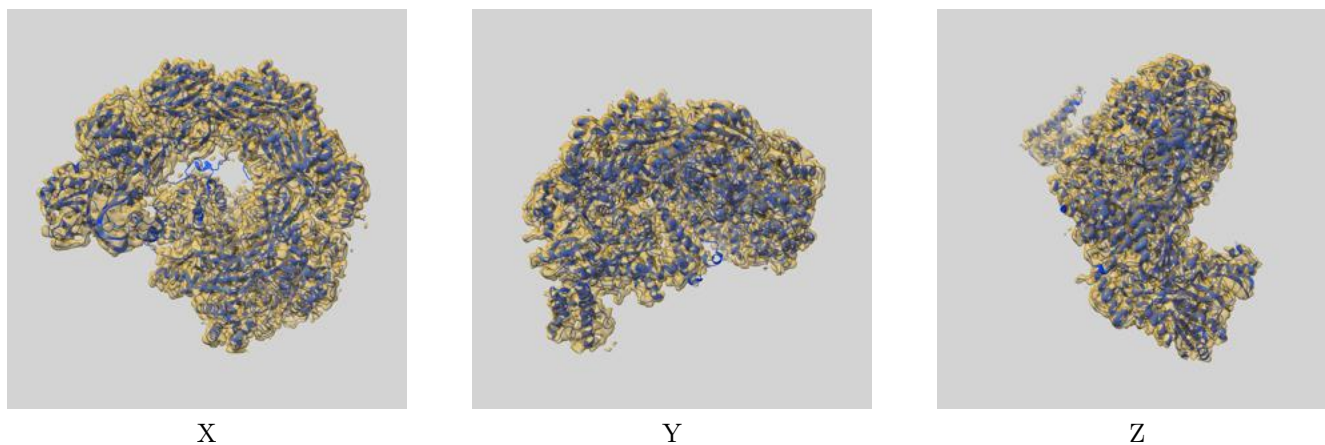
## 8 Fourier-Shell correlation

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

## 9 Map-model fit [i](#)

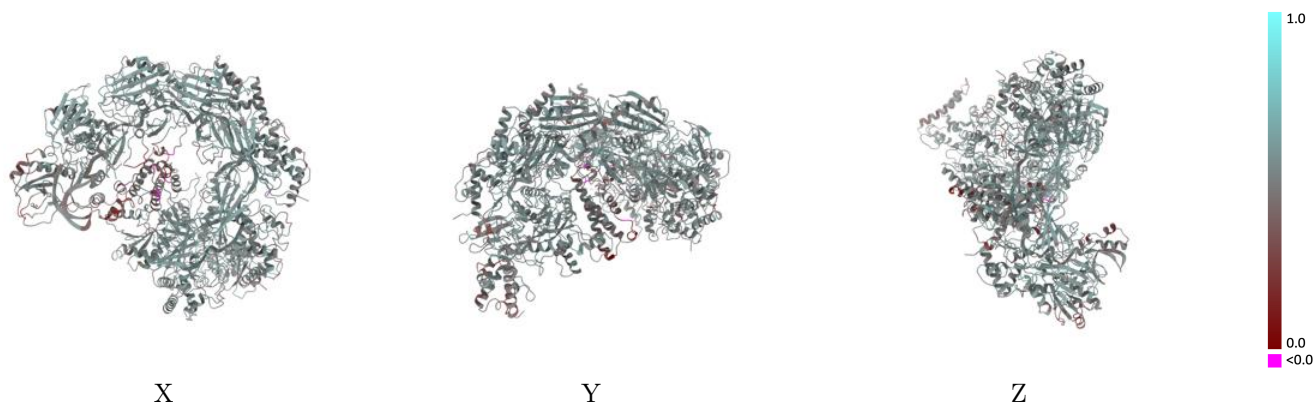
This section contains information regarding the fit between EMDB map EMD-22583 and PDB model 7JZX. Per-residue inclusion information can be found in section 3 on page 6.

### 9.1 Map-model overlay [i](#)



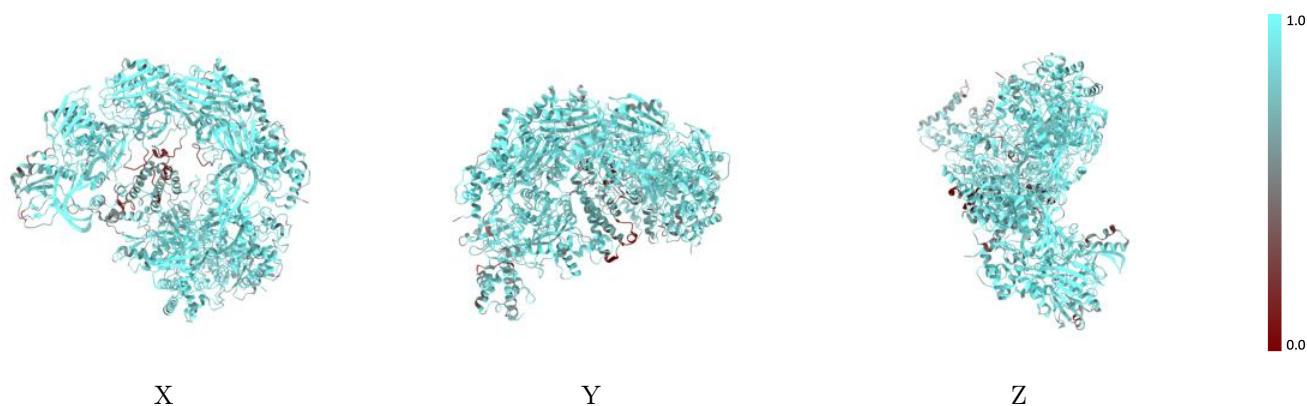
The images above show the 3D surface view of the map at the recommended contour level 0.0264 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



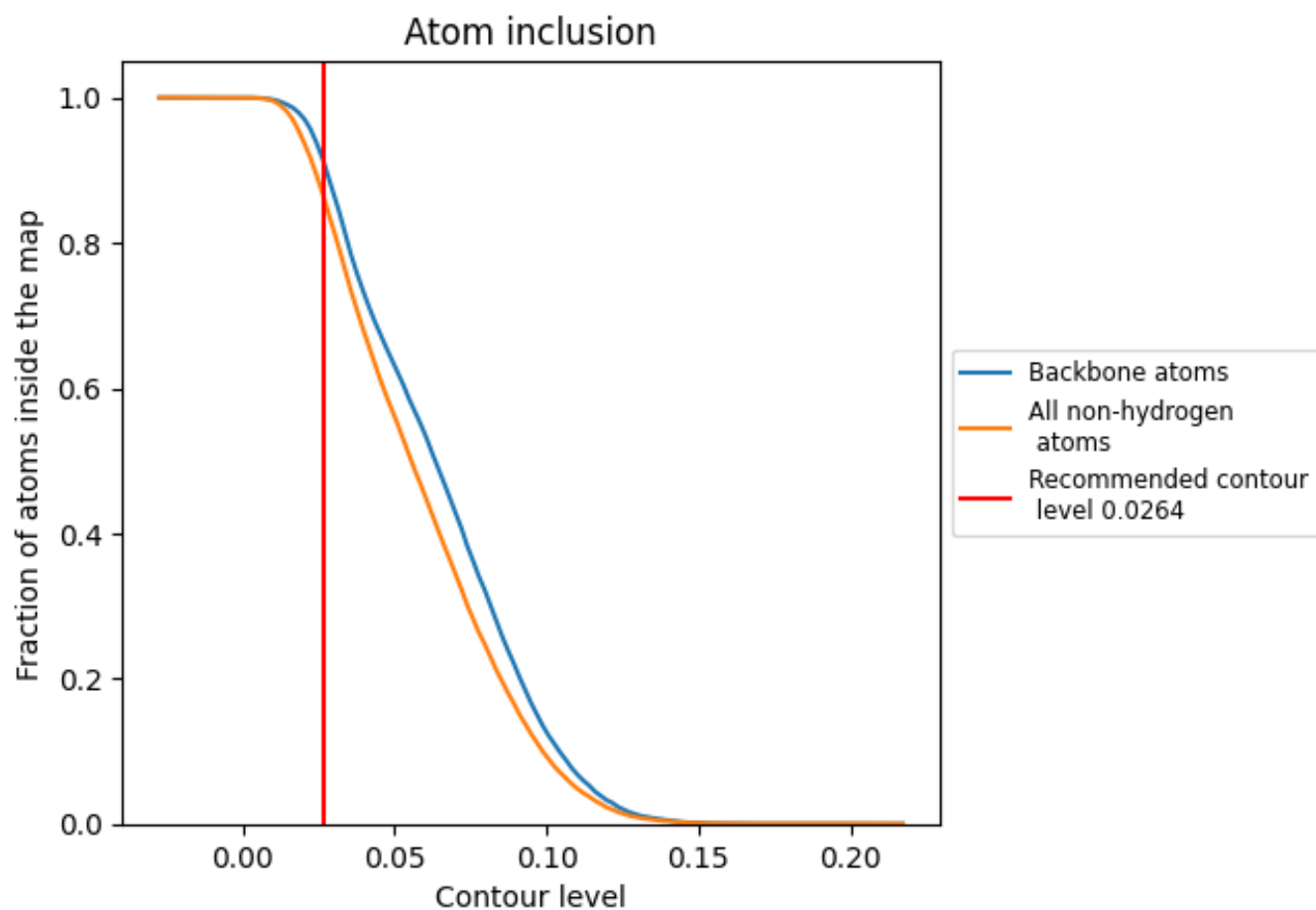
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0264).

























## 9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 87% of all non-hydrogen atoms, are inside the map.

## 9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.0264) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8673	 0.5150
A	 0.7619	 0.4600
B	 0.8997	 0.5260
C	 0.7749	 0.4600
D	 0.9115	 0.5330
E	 0.8805	 0.5280
F	 0.8714	 0.5230
G	 0.8710	 0.5270
H	 0.9040	 0.5340
I	 0.9106	 0.5410
J	 0.7560	 0.4950
M	 0.9262	 0.5100

