



Full wwPDB EM Validation Report ⓘ

Sep 24, 2026 – 04:55 PM EDT

PDB ID : 9P7M / pdb_00009p7m
EMDB ID : EMD-71347
Title : LPS transport bridge LptBFGCADE from Escherichia coli
Authors : Taylor, R.J.; Pahil, K.S.; Walsh Jr., R.M.; Kahne, D.E.
Deposited on : 2025-06-20
Resolution : 4.00 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:

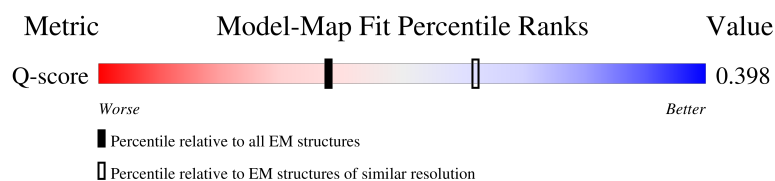
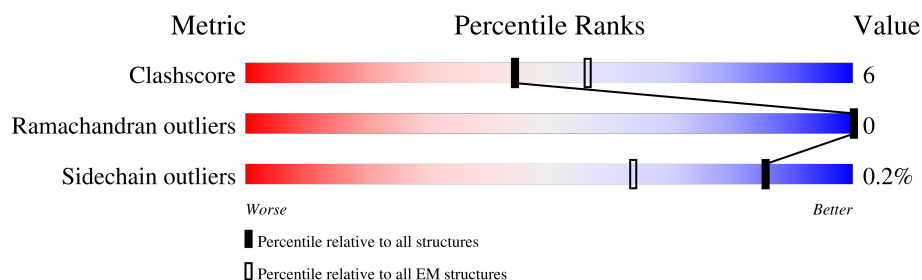
EMDB validation analysis : 0.0.1.dev133
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

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 4.00 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	7587 (3.50 - 4.50)

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	B	241	
1	H	241	
2	C	383	
3	D	784	

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Mol	Chain	Length	Quality of chain
4	E	193	
5	F	366	
6	G	360	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 33703 atoms, of which 16685 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 Lipopolysaccharide export system ATP-binding protein LptB.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	B	238	Total	C	H	N	O	S	0	0
			3733	1162	1876	338	352	5		
1	H	239	Total	C	H	N	O	S	0	0
			3740	1165	1878	339	353	5		

- Molecule 2 is a protein called Lipopolysaccharide export system protein LptC,Lipopolysaccharide export system protein LptA.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	C	305	Total	C	H	N	O	S	0	0
			3987	1286	1888	376	430	7		

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	194	LEU	-	linker	UNP P0ADV9
C	195	VAL	-	linker	UNP P0ADV9
C	196	PRO	-	linker	UNP P0ADV9
C	197	GLY	-	linker	UNP P0ADV9
C	198	GLY	-	linker	UNP P0ADV9
C	199	SER	-	linker	UNP P0ADV9
C	200	GLY	-	linker	UNP P0ADV9
C	201	GLY	-	linker	UNP P0ADV9
C	202	GLY	-	linker	UNP P0ADV9
C	203	GLY	-	linker	UNP P0ADV9
C	204	SER	-	linker	UNP P0ADV9
C	205	GLY	-	linker	UNP P0ADV9
C	206	GLY	-	linker	UNP P0ADV9
C	207	GLY	-	linker	UNP P0ADV9
C	208	GLY	-	linker	UNP P0ADV9
C	209	SER	-	linker	UNP P0ADV9
C	210	GLY	-	linker	UNP P0ADV9
C	211	GLY	-	linker	UNP P0ADV9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	212	GLY	-	linker	UNP P0ADV9
C	213	GLY	-	linker	UNP P0ADV9
C	214	SER	-	linker	UNP P0ADV9
C	373	LEU	-	expression tag	UNP P0ADV1
C	374	VAL	-	expression tag	UNP P0ADV1
C	375	PRO	-	expression tag	UNP P0ADV1
C	376	ARG	-	expression tag	UNP P0ADV1
C	377	GLY	-	expression tag	UNP P0ADV1
C	378	SER	-	expression tag	UNP P0ADV1
C	379	HIS	-	expression tag	UNP P0ADV1
C	380	HIS	-	expression tag	UNP P0ADV1
C	381	HIS	-	expression tag	UNP P0ADV1
C	382	HIS	-	expression tag	UNP P0ADV1
C	383	HIS	-	expression tag	UNP P0ADV1
C	384	HIS	-	expression tag	UNP P0ADV1
C	385	HIS	-	expression tag	UNP P0ADV1

- Molecule 3 is a protein called LPS-assembly protein LptD.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	D	759	Total	C	H	N	O	S	0	0
			11886	3860	5740	1054	1212	20		

- Molecule 4 is a protein called LPS-assembly lipoprotein LptE.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	E	151	Total	C	H	N	O	S	0	0
			2405	740	1218	216	223	8		

- Molecule 5 is a protein called Lipopolysaccharide export system permease protein LptF.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	F	247	Total	C	H	N	O	S	0	0
			4015	1278	2075	320	326	16		

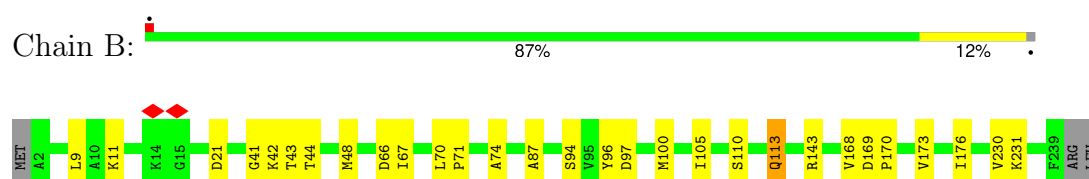
- Molecule 6 is a protein called Lipopolysaccharide export system permease protein LptG.

Mol	Chain	Residues	Atoms						AltConf	Trace
6	G	250	Total	C	H	N	O	S	0	0
			3937	1277	2010	307	325	18		

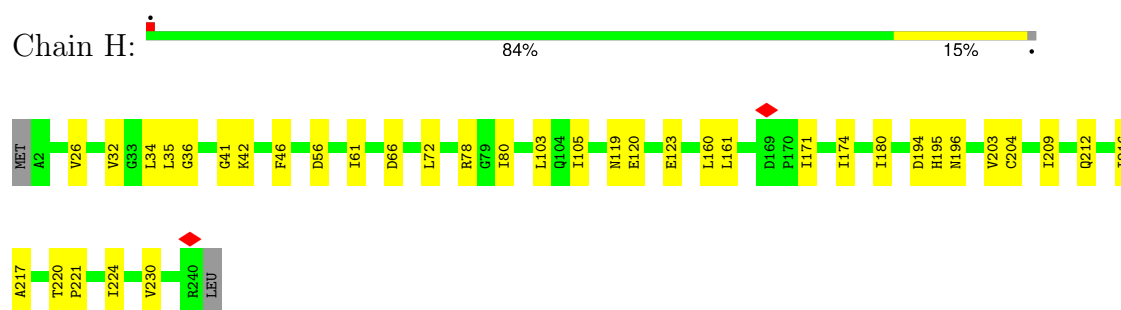
3 Residue-property plots [i](#)

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.

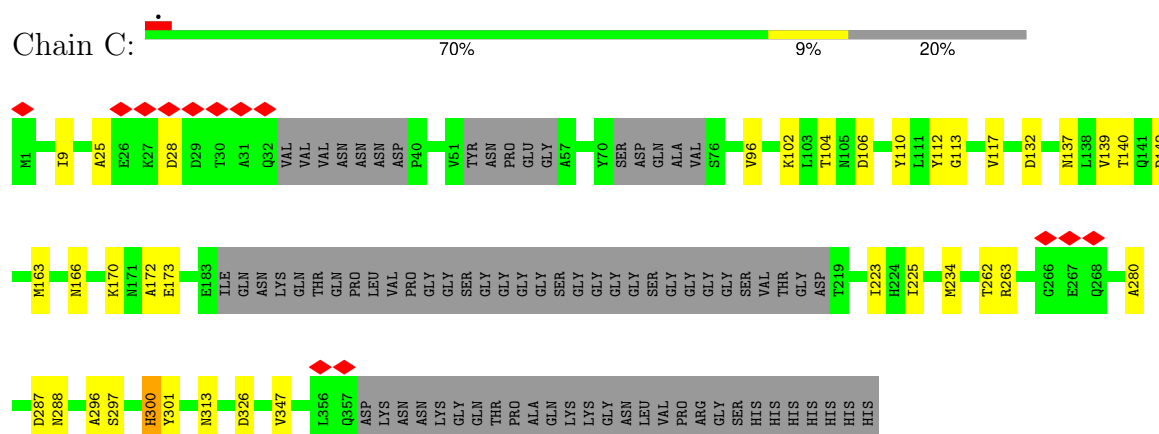
- Molecule 1: Lipopolysaccharide export system ATP-binding protein LptB



- Molecule 1: Lipopolysaccharide export system ATP-binding protein LptB

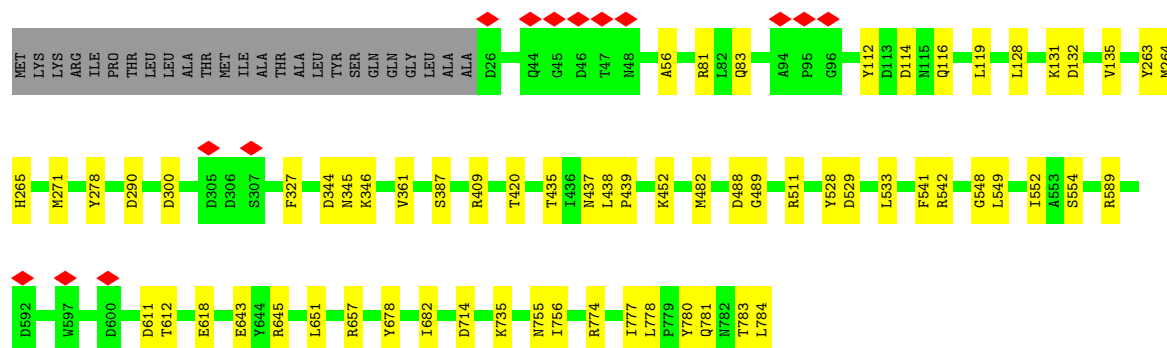


- Molecule 2: Lipopolysaccharide export system protein LptC, Lipopolysaccharide export system protein LptA



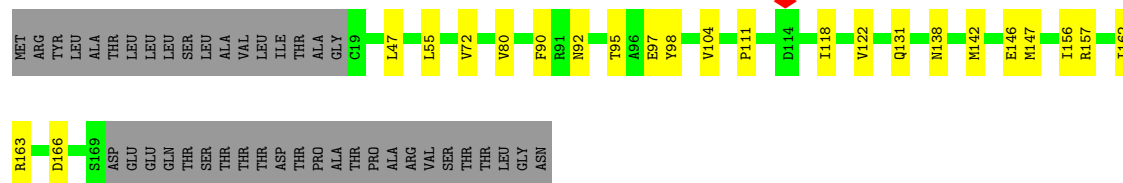
- Molecule 3: LPS-assembly protein LptD

Chain D: 



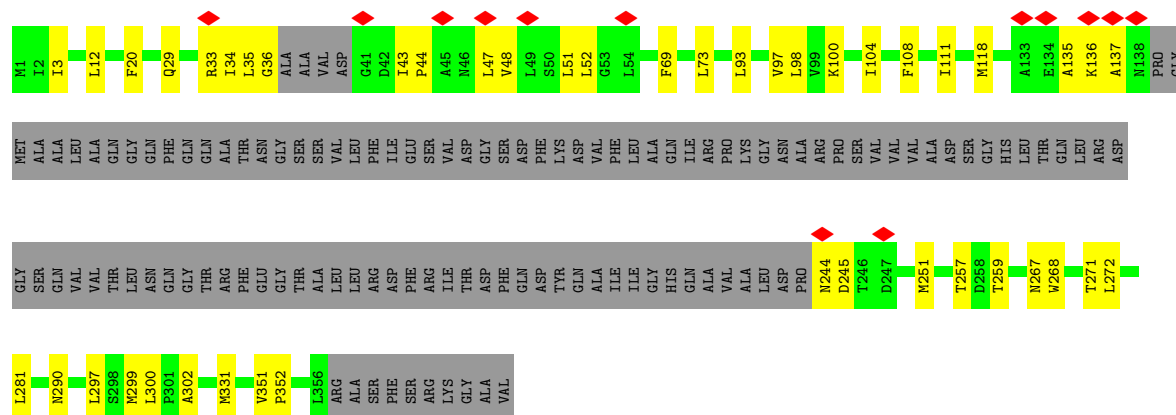
• Molecule 4: LPS-assembly lipoprotein LptE

Chain E: 



• Molecule 5: Lipopolysaccharide export system permease protein LptF

Chain F: 



• Molecule 6: Lipopolysaccharide export system permease protein LptG

Chain G: 





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	165070	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	63.918	Depositor
Minimum defocus (nm)	1100	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOCONTINUUM (6k x 4k)	Depositor
Maximum map value	0.323	Depositor
Minimum map value	-0.016	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.0342	Depositor
Map size (Å)	495.0, 495.0, 495.0	wwPDB
Map dimensions	600, 600, 600	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.825, 0.825, 0.825	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	B	0.26	0/1884	0.46	0/2548
1	H	0.26	0/1889	0.46	0/2555
2	C	0.20	0/2122	0.42	0/2865
3	D	0.20	0/6307	0.39	0/8587
4	E	0.22	0/1204	0.40	0/1625
5	F	0.24	0/1975	0.44	0/2685
6	G	0.23	0/1968	0.53	0/2661
All	All	0.22	0/17349	0.43	0/23526

There are no bond length outliers.

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	B	1857	1876	1875	21	0
1	H	1862	1878	1877	28	0
2	C	2099	1888	1841	25	0
3	D	6146	5740	5743	47	0
4	E	1187	1218	1217	17	0
5	F	1940	2075	2079	35	0
6	G	1927	2010	2030	55	0
All	All	17018	16685	16662	215	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 6.

All (215) 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:D:657:ARG:NH1	4:E:97:GLU:OE1	2.05	0.88
2:C:297:SER:OG	2:C:313:ASN:OD1	2.01	0.77
3:D:116:GLN:OE1	3:D:116:GLN:N	2.17	0.77
3:D:529:ASP:OD1	4:E:138:ASN:ND2	2.20	0.74
3:D:345:ASN:OD1	3:D:346:LYS:N	2.24	0.71
5:F:44:PRO:HD2	5:F:47:LEU:HD12	1.74	0.68
3:D:81:ARG:NH1	3:D:83:GLN:OE1	2.27	0.67
4:E:163:ARG:NH1	4:E:166:ASP:OD2	2.29	0.66
3:D:755:ASN:OD1	3:D:756:ILE:N	2.30	0.65
3:D:435:THR:HG22	3:D:452:LYS:HG2	1.80	0.64
6:G:62:LYS:HE2	6:G:62:LYS:HA	1.80	0.63
6:G:275:MET:HE2	6:G:275:MET:C	2.23	0.63
2:C:96:VAL:HG22	2:C:117:VAL:HG22	1.80	0.63
6:G:50:LEU:HD23	6:G:54:MET:CE	2.28	0.63
4:E:80:VAL:HG12	4:E:104:VAL:HG23	1.80	0.63
1:H:61:ILE:HG13	1:H:61:ILE:O	1.97	0.62
6:G:125:ALA:O	6:G:129:GLU:HG3	2.00	0.61
6:G:58:LEU:HD21	6:G:131:MET:CG	2.30	0.61
3:D:618:GLU:OE1	3:D:618:GLU:N	2.26	0.61
3:D:542:ARG:NH1	4:E:146:GLU:OE2	2.34	0.61
5:F:36:GLY:O	5:F:43:ILE:HG22	2.01	0.60
6:G:244:LEU:HB3	6:G:257:TYR:HE2	1.66	0.60
6:G:321:VAL:O	6:G:325:ILE:HG23	2.03	0.59
6:G:275:MET:HE1	6:G:276:TRP:CE3	2.40	0.57
1:H:103:LEU:HD12	1:H:103:LEU:O	2.05	0.57
1:H:209:ILE:HG13	1:H:224:ILE:HD11	1.87	0.56
4:E:122:VAL:HG21	4:E:147:MET:HG2	1.88	0.55
5:F:136:LYS:O	5:F:137:ALA:HB3	2.07	0.55
3:D:128:LEU:HD23	3:D:128:LEU:O	2.07	0.55
5:F:51:LEU:HD12	5:F:52:LEU:N	2.21	0.55
5:F:47:LEU:HD11	5:F:135:ALA:CB	2.37	0.54
3:D:387:SER:OG	3:D:420:THR:OG1	2.12	0.54
5:F:244:ASN:O	5:F:245:ASP:HB2	2.07	0.54
6:G:130:GLN:O	6:G:134:ASN:OD1	2.25	0.54
1:H:220:THR:HB	1:H:221:PRO:HD2	1.90	0.53
4:E:72:VAL:HG13	4:E:111:PRO:HD2	1.91	0.53
5:F:73:LEU:HD21	5:F:299:MET:HG3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:G:110:ALA:HB2	6:G:289:MET:CE	2.39	0.53
2:C:9:ILE:HD12	5:F:300:LEU:CD2	2.38	0.52
1:H:171:ILE:O	1:H:174:ILE:HG22	2.09	0.52
6:G:357:MET:HA	6:G:357:MET:HE2	1.92	0.52
6:G:20:MET:HA	6:G:20:MET:HE2	1.92	0.52
1:B:21:ASP:CG	1:B:21:ASP:O	2.52	0.51
6:G:57:LEU:O	6:G:60:VAL:HG23	2.10	0.51
6:G:110:ALA:HB2	6:G:289:MET:HE1	1.93	0.51
6:G:122:GLU:HG3	6:G:123:TRP:CD1	2.46	0.51
5:F:267:ASN:ND2	5:F:331:MET:HE2	2.26	0.51
5:F:351:VAL:HG22	5:F:352:PRO:HD3	1.93	0.51
1:B:71:PRO:HD2	5:F:351:VAL:HG12	1.93	0.50
2:C:300:HIS:CD2	2:C:301:TYR:N	2.80	0.50
3:D:781:GLN:N	3:D:781:GLN:OE1	2.45	0.50
1:H:161:LEU:CD1	1:H:180:ILE:HD11	2.42	0.50
3:D:777:ILE:HG13	3:D:778:LEU:HD12	1.93	0.50
6:G:58:LEU:HD21	6:G:131:MET:HG3	1.94	0.50
6:G:40:LYS:O	6:G:40:LYS:HG2	2.12	0.50
2:C:223:ILE:H	2:C:223:ILE:HD12	1.76	0.49
3:D:755:ASN:OD1	3:D:755:ASN:C	2.54	0.49
4:E:122:VAL:HG21	4:E:147:MET:CG	2.42	0.49
6:G:50:LEU:HD23	6:G:54:MET:HE1	1.93	0.49
6:G:97:PHE:HB2	1:H:105:ILE:HD11	1.94	0.49
1:B:11:LYS:HB2	1:B:48:MET:HE1	1.95	0.49
2:C:234:MET:HE2	2:C:234:MET:O	2.13	0.49
3:D:554:SER:OG	3:D:589:ARG:O	2.26	0.49
1:H:26:VAL:HG22	1:H:32:VAL:HG21	1.95	0.49
1:H:78:ARG:NE	1:H:78:ARG:HA	2.26	0.48
3:D:678:TYR:OH	4:E:92:ASN:ND2	2.46	0.48
1:H:194:ASP:OD1	1:H:195:HIS:N	2.46	0.48
2:C:262:THR:HG22	2:C:263:ARG:N	2.28	0.48
2:C:173:GLU:OE1	2:C:173:GLU:N	2.46	0.47
3:D:488:ASP:OD1	3:D:489:GLY:N	2.47	0.47
5:F:47:LEU:HD11	5:F:135:ALA:HB2	1.95	0.47
3:D:300:ASP:OD1	3:D:300:ASP:O	2.32	0.47
4:E:47:LEU:HD22	4:E:80:VAL:HG11	1.96	0.47
6:G:274:ASN:OD1	6:G:274:ASN:C	2.58	0.47
6:G:348:PHE:HA	6:G:351:ILE:HG22	1.95	0.47
6:G:122:GLU:OE2	6:G:123:TRP:HD1	1.97	0.47
1:B:169:ASP:HB2	1:B:170:PRO:HD2	1.96	0.47
6:G:16:PHE:O	6:G:19:ILE:HG22	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:G:107:MET:HE1	6:G:290:LEU:HD12	1.95	0.47
1:B:168:VAL:HG11	1:B:176:ILE:HD12	1.97	0.47
4:E:55:LEU:HD21	4:E:156:ILE:HD11	1.97	0.47
3:D:300:ASP:OD1	3:D:300:ASP:C	2.58	0.47
1:B:110:SER:OG	1:B:113:GLN:CD	2.58	0.47
3:D:511:ARG:HD3	3:D:541:PHE:CE2	2.50	0.47
3:D:409:ARG:HG2	3:D:435:THR:OG1	2.15	0.46
6:G:50:LEU:CD2	6:G:54:MET:HE1	2.45	0.46
1:H:80:ILE:N	1:H:80:ILE:HD12	2.30	0.46
6:G:261:LEU:HD11	6:G:271:TYR:CD1	2.50	0.46
1:B:87:ALA:HB2	1:B:143:ARG:NH1	2.30	0.46
3:D:344:ASP:OD2	3:D:735:LYS:NZ	2.40	0.46
1:B:67:ILE:HG22	1:B:67:ILE:O	2.15	0.46
3:D:651:LEU:C	3:D:651:LEU:HD23	2.41	0.46
5:F:118:MET:HG2	5:F:251:MET:HE3	1.97	0.46
1:H:72:LEU:O	1:H:72:LEU:HD12	2.16	0.46
1:B:230:VAL:HG13	1:B:231:LYS:N	2.30	0.46
3:D:774:ARG:NH1	3:D:780:TYR:O	2.41	0.46
5:F:34:ILE:C	5:F:35:LEU:HD23	2.40	0.46
5:F:43:ILE:HD11	5:F:51:LEU:HD21	1.97	0.46
1:H:209:ILE:CG1	1:H:224:ILE:HD11	2.45	0.46
6:G:254:LEU:HD13	6:G:275:MET:HB3	1.98	0.46
3:D:643:GLU:OE1	3:D:783:THR:OG1	2.25	0.45
5:F:48:VAL:HA	5:F:51:LEU:HG	1.99	0.45
1:B:94:SER:HB3	1:B:97:ASP:OD2	2.16	0.45
6:G:25:MET:O	6:G:29:LEU:HD13	2.16	0.45
3:D:131:LYS:HG3	3:D:132:ASP:OD2	2.17	0.45
3:D:271:MET:HE3	3:D:290:ASP:HB3	1.98	0.45
5:F:97:VAL:HG13	5:F:98:LEU:HD12	1.99	0.45
1:H:56:ASP:OD1	1:H:56:ASP:C	2.60	0.45
5:F:290:ASN:OD1	5:F:290:ASN:C	2.60	0.45
6:G:133:ARG:HG2	6:G:133:ARG:HH11	1.82	0.45
6:G:281:GLN:HG3	6:G:282:PRO:HD3	1.98	0.45
1:H:203:VAL:HG23	1:H:204:CYS:N	2.32	0.45
3:D:263:TYR:HE1	3:D:265:HIS:HA	1.82	0.45
1:H:212:GLN:O	1:H:212:GLN:HG3	2.17	0.45
1:B:9:LEU:HD12	1:B:48:MET:SD	2.57	0.45
3:D:645:ARG:HG3	3:D:645:ARG:O	2.16	0.45
3:D:112:TYR:OH	3:D:114:ASP:OD2	2.35	0.44
5:F:12:LEU:C	5:F:12:LEU:HD13	2.43	0.44
6:G:328:PRO:O	6:G:331:LEU:HG	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:272:LEU:HD23	5:F:272:LEU:O	2.17	0.44
6:G:275:MET:SD	6:G:276:TRP:N	2.90	0.44
5:F:272:LEU:HD23	5:F:272:LEU:C	2.43	0.44
6:G:19:ILE:HD11	6:G:72:ALA:HA	1.99	0.44
6:G:281:GLN:N	6:G:282:PRO:HD2	2.33	0.44
6:G:55:TYR:CD1	6:G:132:ALA:HB1	2.52	0.44
6:G:251:ILE:HG13	6:G:252:SER:N	2.32	0.44
3:D:611:ASP:OD1	3:D:612:THR:N	2.50	0.44
2:C:113:GLY:N	2:C:132:ASP:O	2.47	0.44
6:G:106:VAL:O	6:G:109:THR:HG22	2.17	0.44
4:E:118:ILE:HG21	4:E:162:ILE:HD11	1.99	0.44
1:H:34:LEU:C	1:H:35:LEU:HD12	2.43	0.44
3:D:783:THR:O	3:D:784:LEU:C	2.62	0.43
2:C:25:ALA:O	2:C:28:ASP:OD1	2.36	0.43
1:H:119:ASN:OD1	1:H:119:ASN:C	2.60	0.43
1:H:221:PRO:HA	1:H:224:ILE:HG22	2.00	0.43
3:D:533:LEU:HD11	4:E:142:MET:HE1	2.00	0.43
5:F:20:PHE:C	5:F:20:PHE:CD2	2.96	0.43
2:C:9:ILE:HD12	5:F:300:LEU:HD22	1.99	0.43
3:D:548:GLY:C	3:D:549:LEU:HD22	2.44	0.43
6:G:38:GLN:NE2	6:G:55:TYR:CE2	2.86	0.43
1:B:41:GLY:O	1:B:42:LYS:C	2.61	0.43
1:B:70:LEU:HD22	1:B:74:ALA:CB	2.49	0.43
2:C:142:ASP:HA	2:C:166:ASN:HA	2.01	0.43
5:F:297:LEU:HD13	5:F:297:LEU:O	2.17	0.43
6:G:37:ASP:O	6:G:37:ASP:OD1	2.36	0.43
2:C:137:ASN:OD1	2:C:139:VAL:HG22	2.19	0.43
3:D:278:TYR:CD1	3:D:278:TYR:C	2.97	0.43
3:D:528:TYR:OH	4:E:131:GLN:O	2.34	0.43
3:D:682:ILE:HD11	4:E:90:PHE:CE1	2.54	0.43
5:F:108:PHE:O	5:F:111:ILE:HG22	2.18	0.43
1:H:46:PHE:CE2	1:H:160:LEU:HD12	2.54	0.43
6:G:77:LEU:HD12	6:G:308:ARG:HG2	2.01	0.42
2:C:287:ASP:OD1	2:C:288:ASN:N	2.52	0.42
5:F:29:GLN:O	5:F:33:ARG:HG3	2.19	0.42
5:F:281:LEU:HD13	5:F:281:LEU:C	2.45	0.42
1:H:41:GLY:O	1:H:42:LYS:C	2.61	0.42
1:H:194:ASP:OD2	1:H:196:ASN:O	2.38	0.42
2:C:110:TYR:HB3	2:C:112:TYR:CE1	2.54	0.42
4:E:95:THR:HG21	4:E:98:TYR:CZ	2.54	0.42
1:B:105:ILE:HD12	5:F:93:LEU:CD2	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:G:80:LEU:HB2	6:G:296:ILE:HG13	2.00	0.42
6:G:122:GLU:OE2	6:G:123:TRP:CD1	2.72	0.42
6:G:276:TRP:HA	6:G:279:ILE:HG22	2.02	0.42
1:B:70:LEU:HD13	1:B:74:ALA:HB1	2.00	0.42
2:C:326:ASP:OD1	2:C:326:ASP:C	2.62	0.42
3:D:482:MET:HE1	3:D:552:ILE:HB	2.02	0.42
5:F:257:THR:HG22	5:F:259:THR:HG22	2.00	0.42
6:G:244:LEU:N	6:G:244:LEU:HD12	2.35	0.42
1:B:87:ALA:CB	1:B:143:ARG:NH1	2.83	0.42
3:D:119:LEU:HB2	3:D:135:VAL:HG21	2.01	0.42
6:G:116:LEU:C	6:G:116:LEU:HD13	2.44	0.42
6:G:275:MET:HE1	6:G:276:TRP:CD2	2.55	0.42
2:C:163:MET:SD	2:C:163:MET:C	3.03	0.42
5:F:100:LYS:O	5:F:104:ILE:HG12	2.20	0.42
6:G:135:TYR:O	6:G:138:GLN:HG2	2.21	0.41
6:G:326:PHE:O	6:G:327:GLY:C	2.63	0.41
1:B:170:PRO:O	1:B:173:VAL:HG12	2.20	0.41
2:C:102:LYS:O	2:C:110:TYR:HD1	2.03	0.41
1:B:96:TYR:CE1	1:B:100:MET:HE3	2.55	0.41
3:D:651:LEU:HD11	3:D:780:TYR:CZ	2.56	0.41
1:H:212:GLN:O	1:H:212:GLN:CG	2.68	0.41
3:D:131:LYS:HG3	3:D:132:ASP:CG	2.45	0.41
3:D:437:ASN:CG	3:D:439:PRO:HD3	2.45	0.41
6:G:291:MET:HE1	6:G:315:PHE:HD2	1.85	0.41
2:C:170:LYS:HE2	2:C:234:MET:SD	2.61	0.41
6:G:140:MET:O	6:G:141:TYR:HB2	2.20	0.41
1:B:43:THR:CG2	1:B:44:THR:N	2.83	0.41
2:C:172:ALA:C	2:C:173:GLU:OE1	2.64	0.41
4:E:157:ARG:HG3	4:E:157:ARG:HH11	1.86	0.41
3:D:327:PHE:CE1	3:D:361:VAL:HG23	2.56	0.41
5:F:3:ILE:HD13	5:F:93:LEU:HD13	2.03	0.41
2:C:347:VAL:N	3:D:56:ALA:O	2.54	0.41
6:G:59:SER:O	6:G:60:VAL:C	2.63	0.41
1:H:120:GLU:O	1:H:123:GLU:HG3	2.21	0.41
3:D:263:TYR:CD1	3:D:264:MET:N	2.89	0.41
1:H:36:GLY:N	1:H:42:LYS:HD3	2.36	0.41
2:C:104:THR:HG22	2:C:106:ASP:H	1.85	0.40
3:D:437:ASN:OD1	3:D:438:LEU:N	2.54	0.40
6:G:63:ASP:HA	6:G:66:ILE:HG22	2.03	0.40
6:G:272:GLN:O	6:G:275:MET:SD	2.79	0.40
6:G:315:PHE:HA	6:G:318:VAL:HG22	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:G:343:LEU:HG	6:G:344:PRO:HD3	2.02	0.40
1:B:66:ASP:OD1	1:B:66:ASP:C	2.64	0.40
3:D:714:ASP:OD1	3:D:714:ASP:C	2.64	0.40
1:H:216:ILE:HG21	1:H:230:VAL:HG12	2.03	0.40
2:C:140:THR:OG1	2:C:142:ASP:OD2	2.29	0.40
2:C:280:ALA:N	2:C:296:ALA:O	2.54	0.40
5:F:69:PHE:HZ	5:F:302:ALA:HB1	1.85	0.40
1:H:66:ASP:OD1	1:H:66:ASP:C	2.64	0.40
2:C:225:ILE:N	2:C:225:ILE:HD12	2.36	0.40
1:H:216:ILE:HG22	1:H:217:ALA:N	2.36	0.40
1:B:170:PRO:HA	1:B:173:VAL:HG12	2.04	0.40
5:F:268:TRP:O	5:F:271:THR:HG22	2.20	0.40
5:F:297:LEU:HD13	5:F:297:LEU:C	2.47	0.40
5:F:351:VAL:N	5:F:352:PRO:CD	2.84	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	B	236/241 (98%)	225 (95%)	11 (5%)	0	100	100
1	H	237/241 (98%)	225 (95%)	12 (5%)	0	100	100
2	C	295/383 (77%)	285 (97%)	10 (3%)	0	100	100
3	D	757/784 (97%)	722 (95%)	35 (5%)	0	100	100
4	E	149/193 (77%)	146 (98%)	3 (2%)	0	100	100
5	F	241/366 (66%)	230 (95%)	11 (5%)	0	100	100
6	G	244/360 (68%)	233 (96%)	11 (4%)	0	100	100
All	All	2159/2568 (84%)	2066 (96%)	93 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	B	199/202 (98%)	198 (100%)	1 (0%)	81	82
1	H	199/202 (98%)	199 (100%)	0	100	100
2	C	195/327 (60%)	194 (100%)	1 (0%)	81	82
3	D	662/681 (97%)	662 (100%)	0	100	100
4	E	132/167 (79%)	132 (100%)	0	100	100
5	F	214/307 (70%)	214 (100%)	0	100	100
6	G	207/300 (69%)	206 (100%)	1 (0%)	81	82
All	All	1808/2186 (83%)	1805 (100%)	3 (0%)	85	87

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	113	GLN
2	C	300	HIS
6	G	27	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	118	ASN
2	C	255	ASN
2	C	300	HIS
3	D	55	ASN
3	D	479	ASN
5	F	346	ASN
6	G	38	GLN
6	G	130	GLN
1	H	129	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

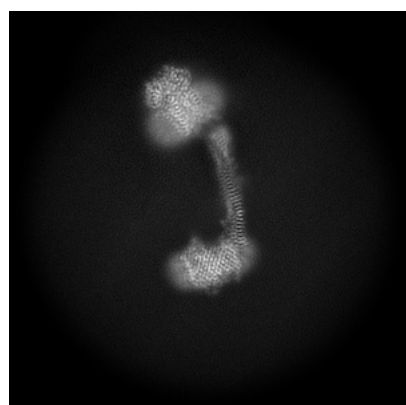
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-71347. These allow visual inspection of the internal detail of the map and identification of artifacts.

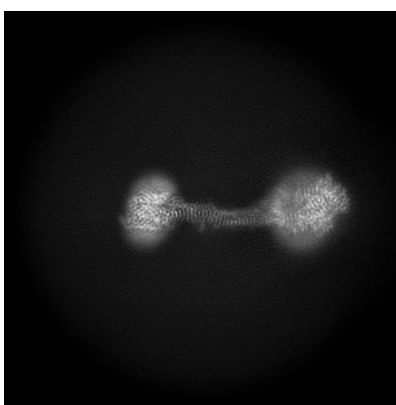
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

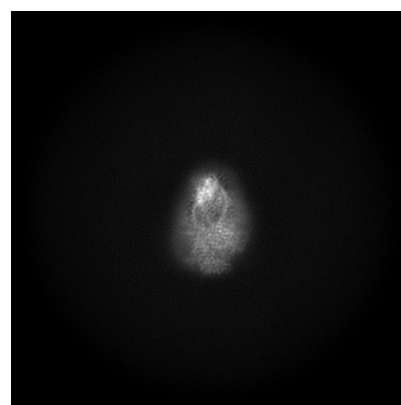
6.1.1 Primary map



X



Y

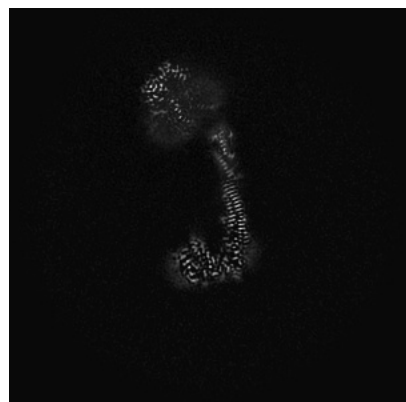


Z

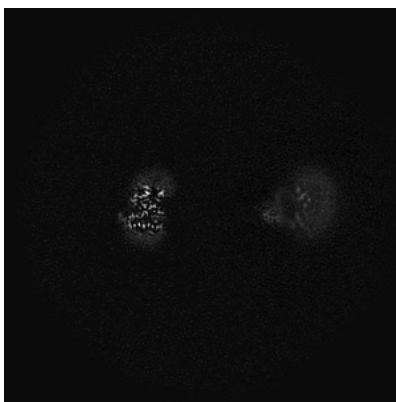
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 300



Y Index: 300

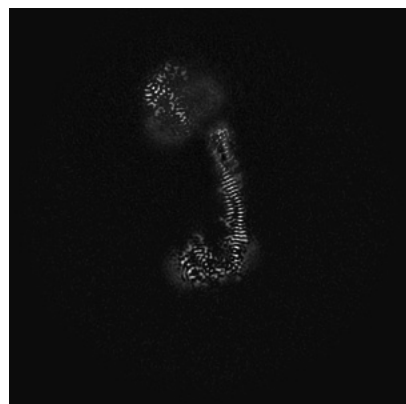


Z Index: 300

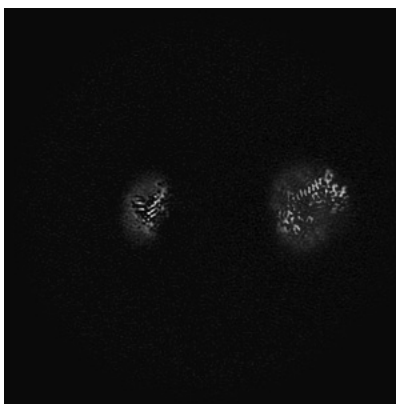
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

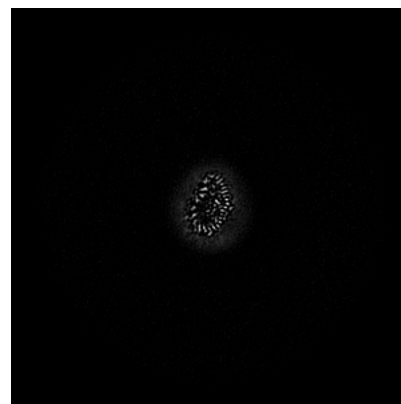
6.3.1 Primary map



X Index: 296



Y Index: 265

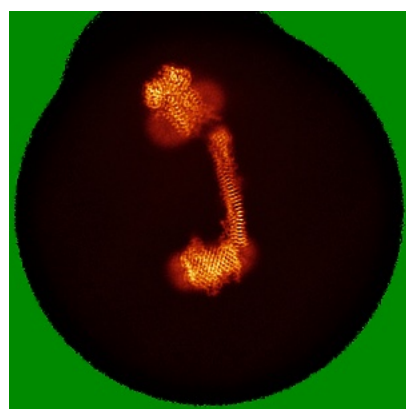


Z Index: 216

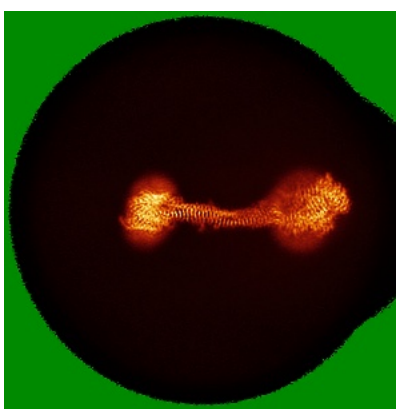
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



X



Y

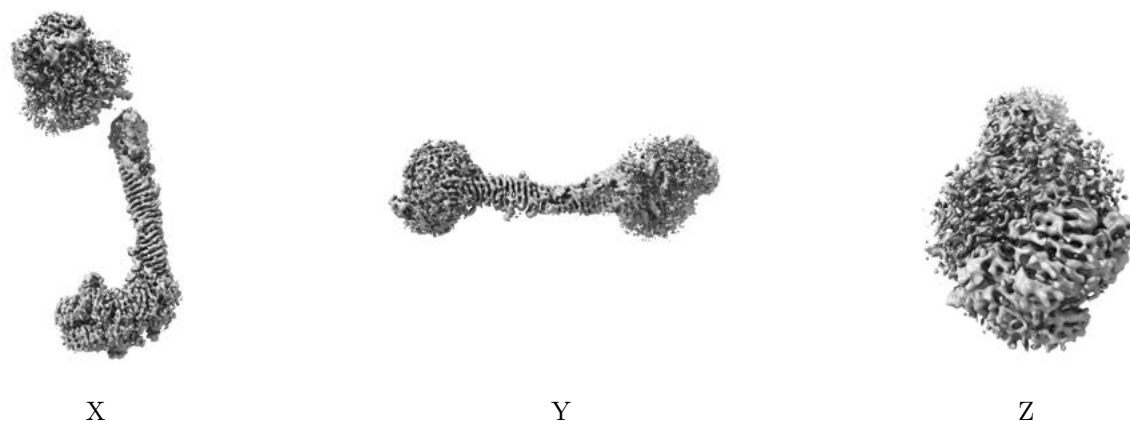


Z

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 0.0342. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

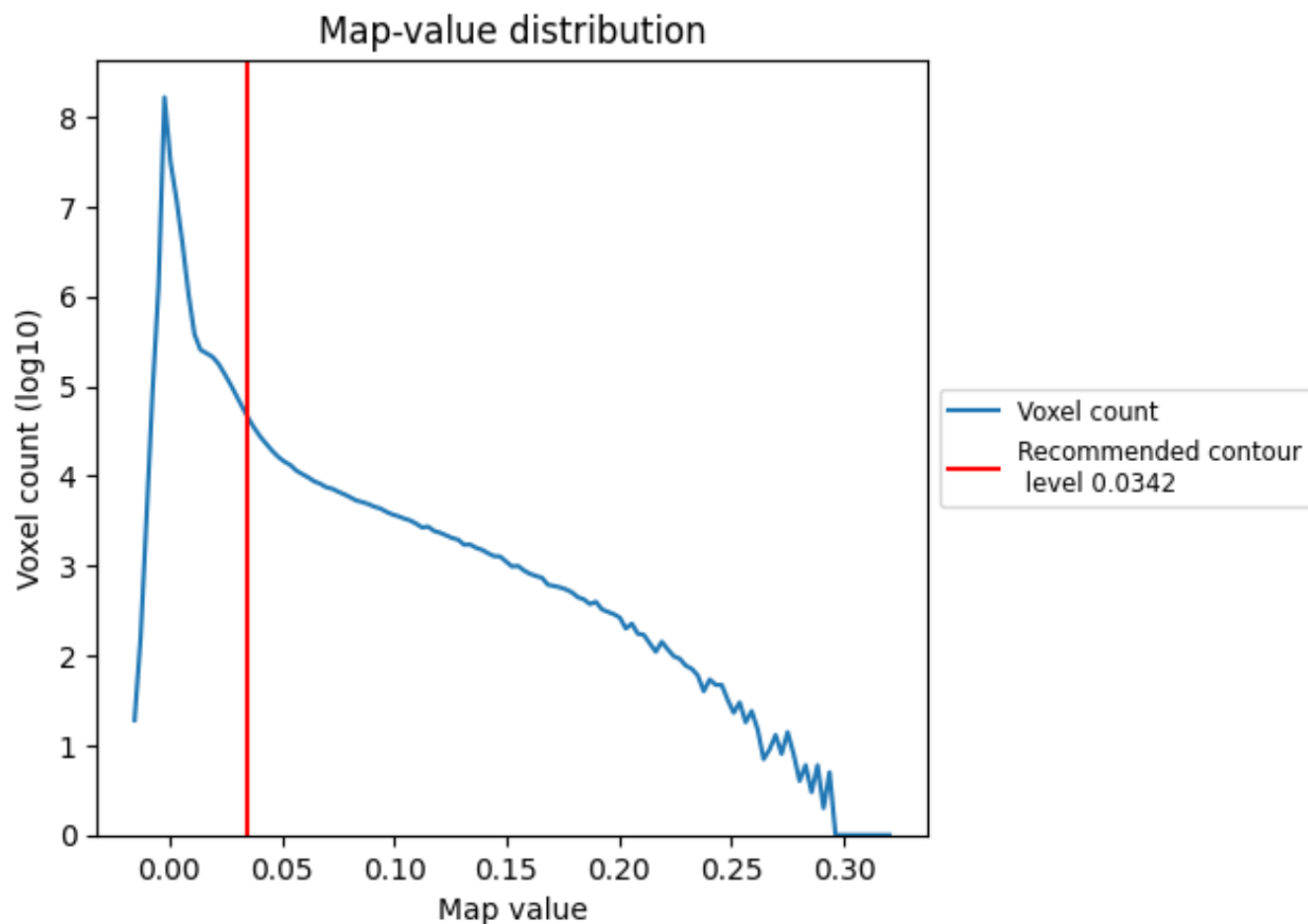
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

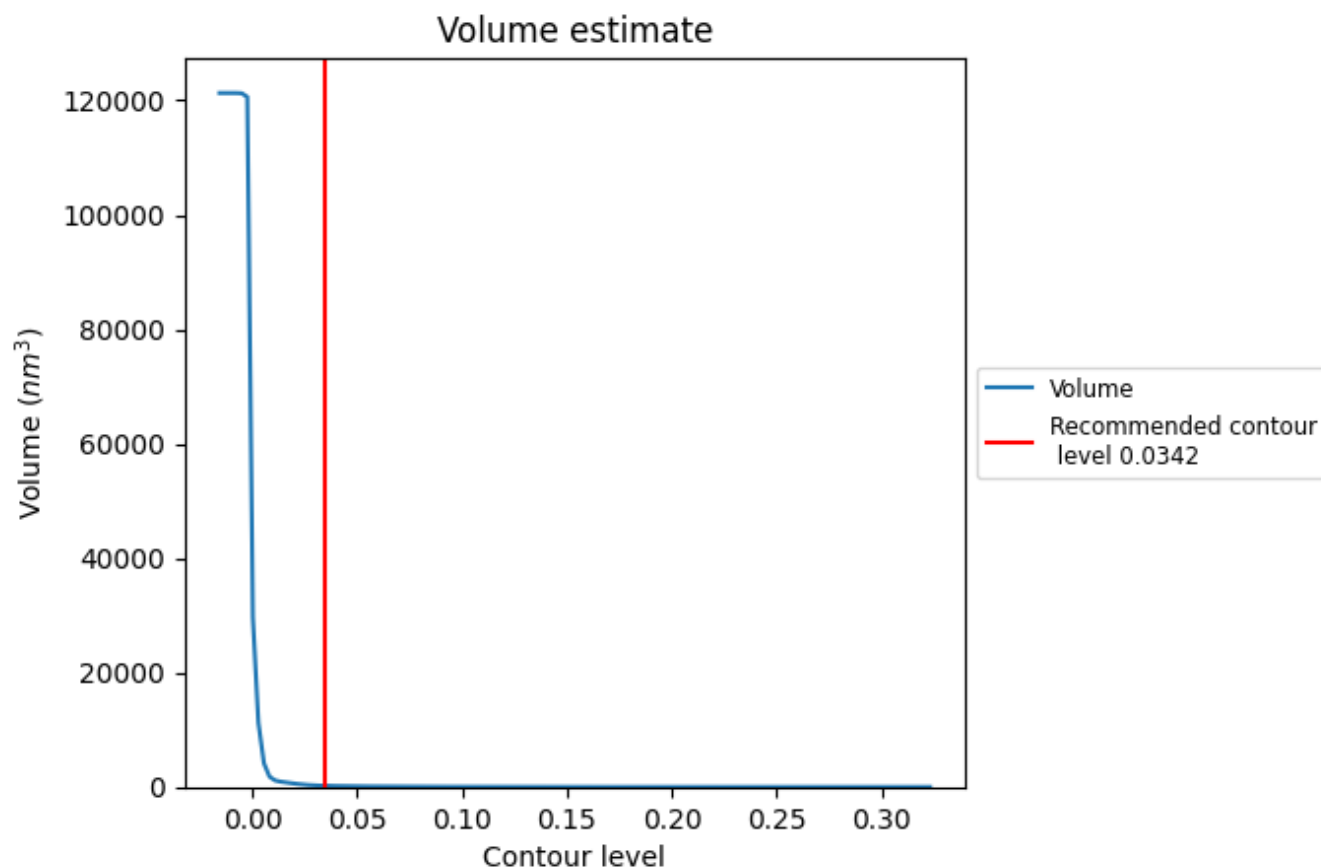
This section contains the results of statistical analysis of the map.

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.

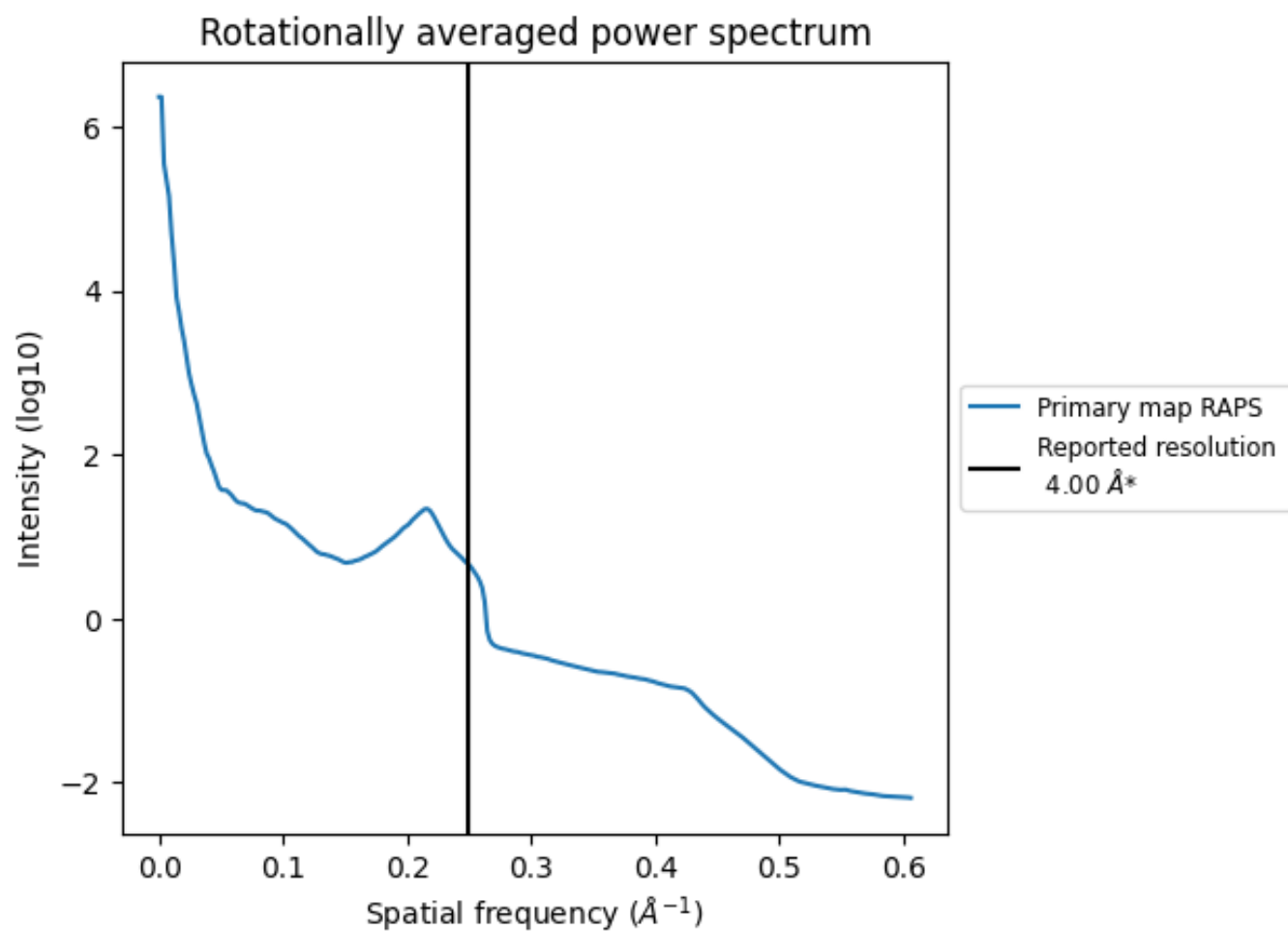
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 211 nm^3 ; this corresponds to an approximate mass of 191 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 0.250 Å⁻¹

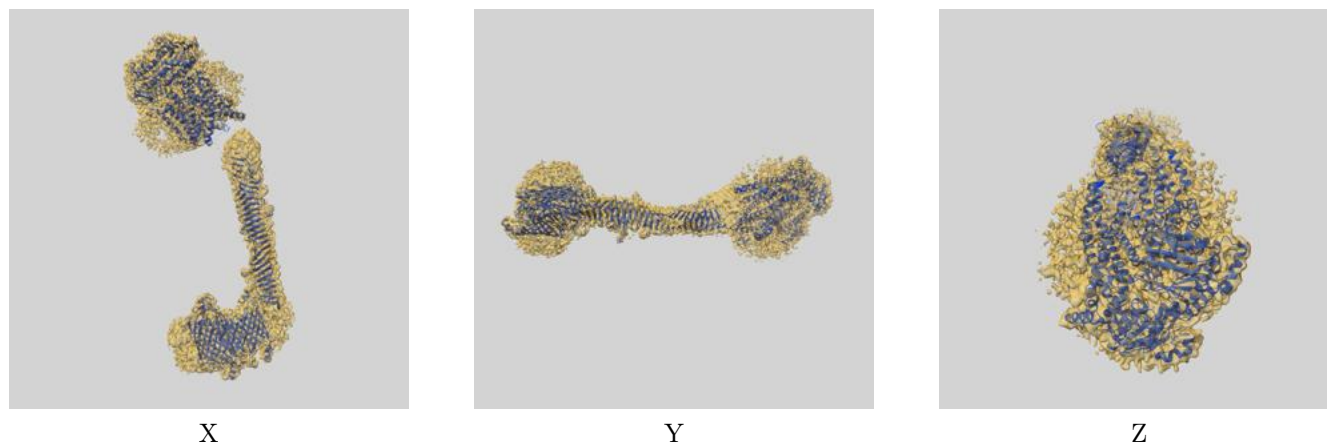
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

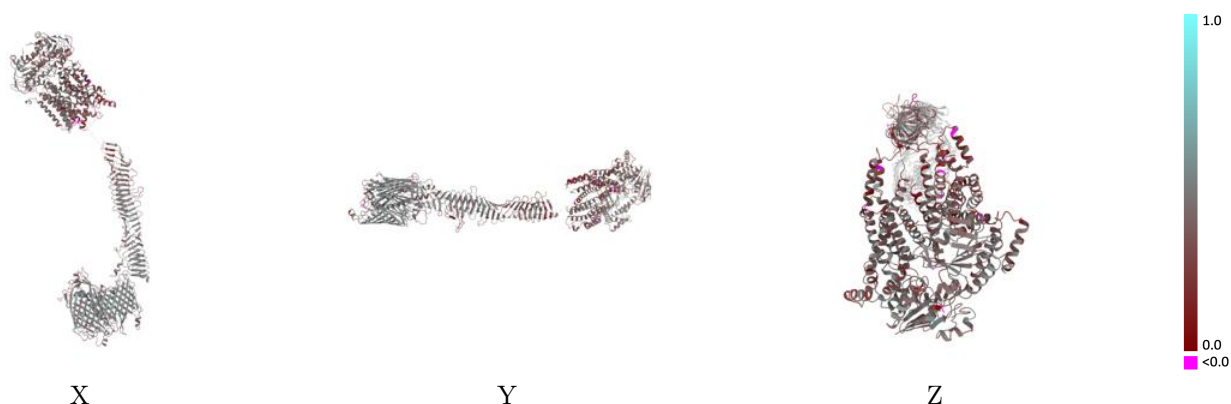
This section contains information regarding the fit between EMDB map EMD-71347 and PDB model 9P7M. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



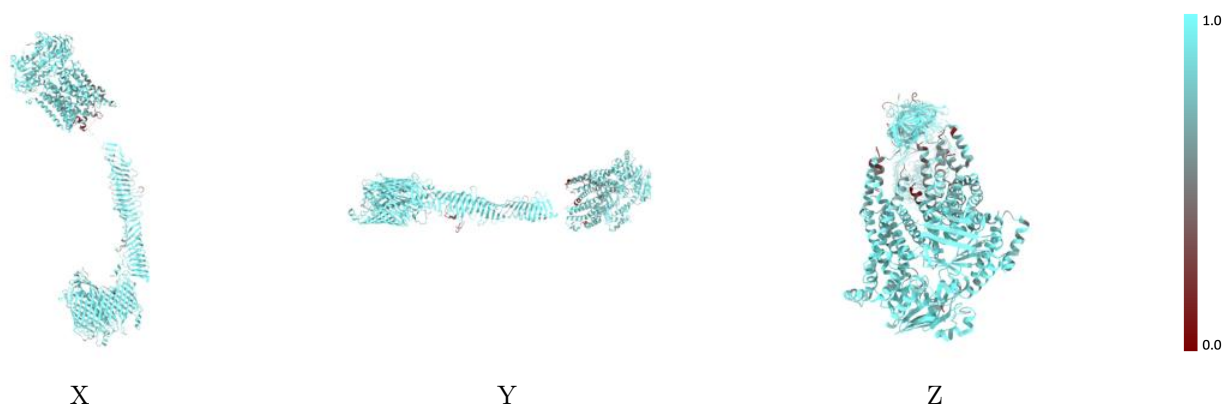
The images above show the 3D surface view of the map at the recommended contour level 0.0342 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



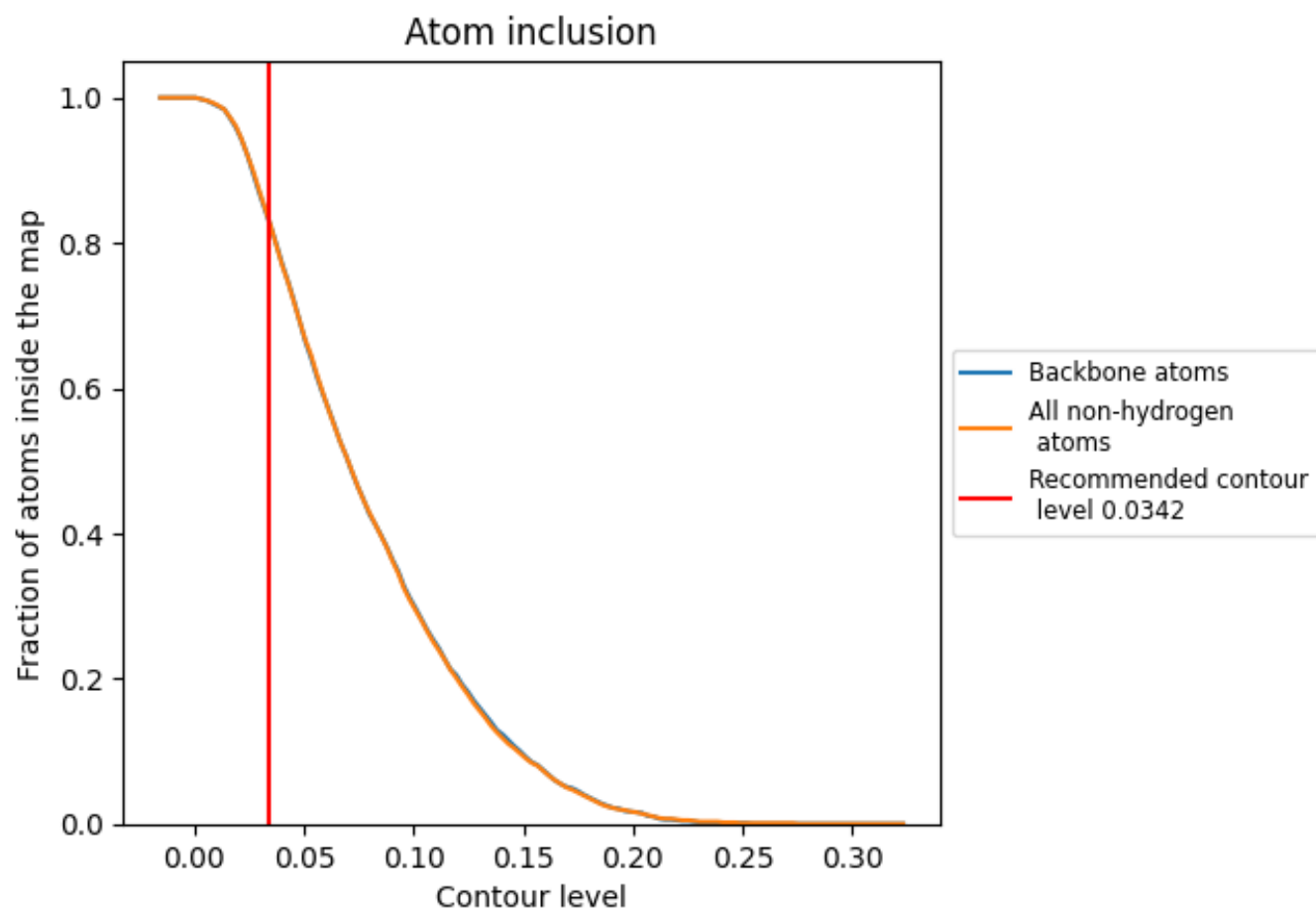
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0342).

9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 83% 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 (0.0342) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8280	0.3980
B	0.8790	0.3970
C	0.7980	0.3520
D	0.8490	0.4460
E	0.8330	0.4300
F	0.8370	0.3660
G	0.7210	0.3160
H	0.8740	0.3860

1.0

0.0

<0.0