



## Full wwPDB EM Validation Report ⓘ

Jul 21, 2024 – 08:23 PM EDT

PDB ID : 9BCT  
EMDB ID : EMD-44438  
Title : Cryo-EM structure of Thermococcus kodakarensis FttA-dependent transcription pre-termination complex containing 44 nt RNA  
Authors : You, L.; Ebright, R.H.  
Deposited on : 2024-04-09  
Resolution : 2.50 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:

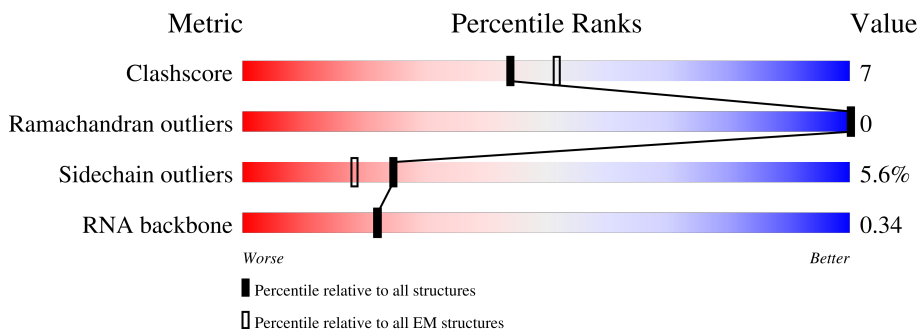
EMDB validation analysis : 0.0.1.dev92  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.37.1

# 1 Overall quality at a glance i

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

The reported resolution of this entry is 2.50 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.









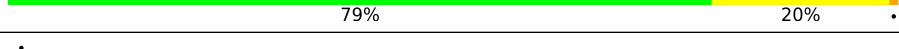
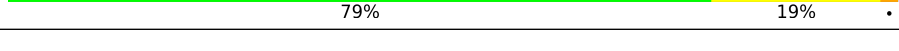
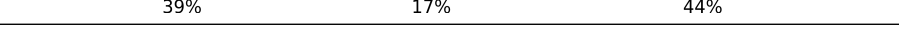
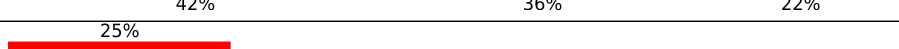
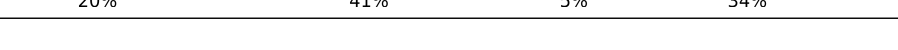
Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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	A	906	
2	B	1123	
3	C	391	
4	D	261	
5	E	190	
6	F	122	
7	H	82	

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Mol	Chain	Length	Quality of chain
8	K	60	 78% 12% 7%
9	L	94	 77% 21%
10	N	65	 78% 20%
11	P	49	 67% 27%
12	G	152	 75% 20%
13	I	73	 5% 71% 16% 11%
14	J	648	 79% 20%
14	M	648	 79% 19%
15	5	36	 39% 17% 44%
16	6	36	 42% 36% 22%
17	7	44	 25% 20% 41% 5% 34%

## 2 Entry composition [i](#)

There are 19 unique types of molecules in this entry. The entry contains 39772 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 DNA-directed RNA polymerase subunit A'.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	900	Total	C	N	O	S	0	0
			7182	4536	1277	1330	39		

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit B.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	1103	Total	C	N	O	S	0	0
			8804	5563	1574	1631	36		

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit A''.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	378	Total	C	N	O	S	0	0
			2958	1868	512	568	10		

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit D.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	256	Total	C	N	O	S	0	0
			2053	1322	339	388	4		

- Molecule 5 is a protein called DNA-directed RNA polymerase subunit E.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	E	183	Total	C	N	O	S	0	0
			1476	946	252	269	9		

- Molecule 6 is a protein called DNA-directed RNA polymerase subunit F.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	F	120	Total	C	N	O	S	0	0
			1000	641	167	189	3		

- Molecule 7 is a protein called DNA-directed RNA polymerase subunit H.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
7	H	77	636	412	106	118	0	0

- Molecule 8 is a protein called DNA-directed RNA polymerase subunit K.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	K	56	433	284	75	73	1	0	0

- Molecule 9 is a protein called DNA-directed RNA polymerase subunit L.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	L	94	776	493	134	147	2	0	0

- Molecule 10 is a protein called DNA-directed RNA polymerase subunit N.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	N	65	530	340	89	95	6	0	0

- Molecule 11 is a protein called DNA-directed RNA polymerase subunit P.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
11	P	48	373	233	74	62	4	0	0

- Molecule 12 is a protein called Transcription elongation factor Spt5.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
12	G	147	1142	739	190	213	0	0

- Molecule 13 is a protein called Transcription elongation factor Spt4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	I	65	515	321	95	95	4	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	-5	HIS	-	expression tag	UNP Q5JIY5
I	-4	HIS	-	expression tag	UNP Q5JIY5
I	-3	HIS	-	expression tag	UNP Q5JIY5
I	-2	HIS	-	expression tag	UNP Q5JIY5
I	-1	HIS	-	expression tag	UNP Q5JIY5
I	0	HIS	-	expression tag	UNP Q5JIY5

- Molecule 14 is a protein called Transcription termination factor FttA.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	J	645	Total	C	N	O	S	0	0
			5144	3284	903	941	16		
14	M	648	Total	C	N	O	S	0	0
			5161	3295	909	942	15		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	255	ALA	HIS	conflict	UNP Q5JH24
J	591	ALA	HIS	conflict	UNP Q5JH24
M	255	ALA	HIS	conflict	UNP Q5JH24
M	591	ALA	HIS	conflict	UNP Q5JH24

- Molecule 15 is a DNA chain called non-template strand DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	5	20	Total	C	N	O	P	0	0
			410	196	74	120	20		

- Molecule 16 is a DNA chain called template strand DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	6	28	Total	C	N	O	P	0	0
			575	273	102	172	28		

- Molecule 17 is a RNA chain called RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	7	29	Total	C	N	O	P	0	0
			595	267	83	216	29		

- Molecule 18 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
18	A	1	Total 1	Mg 1	0

- Molecule 19 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
19	A	2	Total 2	Zn 2	0
19	B	1	Total 1	Zn 1	0
19	N	1	Total 1	Zn 1	0
19	P	1	Total 1	Zn 1	0
19	I	1	Total 1	Zn 1	0
19	J	1	Total 1	Zn 1	0
19	M	1	Total 1	Zn 1	0

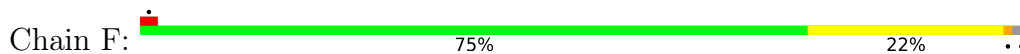








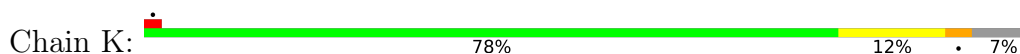
- Molecule 6: DNA-directed RNA polymerase subunit F



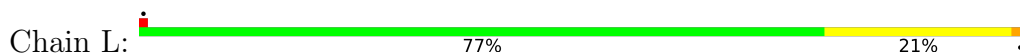
- Molecule 7: DNA-directed RNA polymerase subunit H



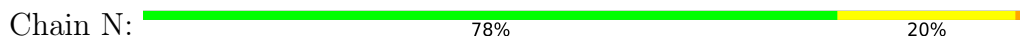
- Molecule 8: DNA-directed RNA polymerase subunit K



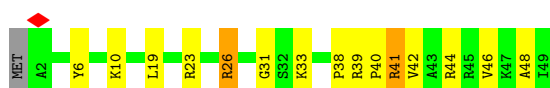
- Molecule 9: DNA-directed RNA polymerase subunit L



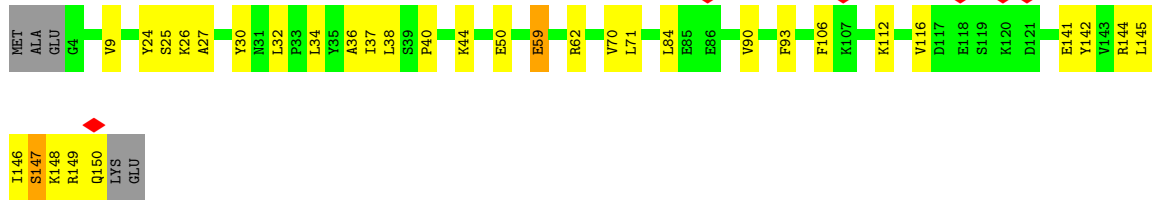
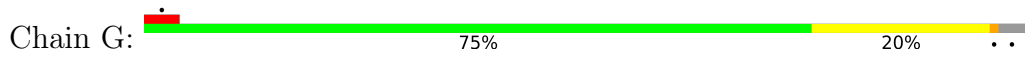
- Molecule 10: DNA-directed RNA polymerase subunit N



- Molecule 11: DNA-directed RNA polymerase subunit P



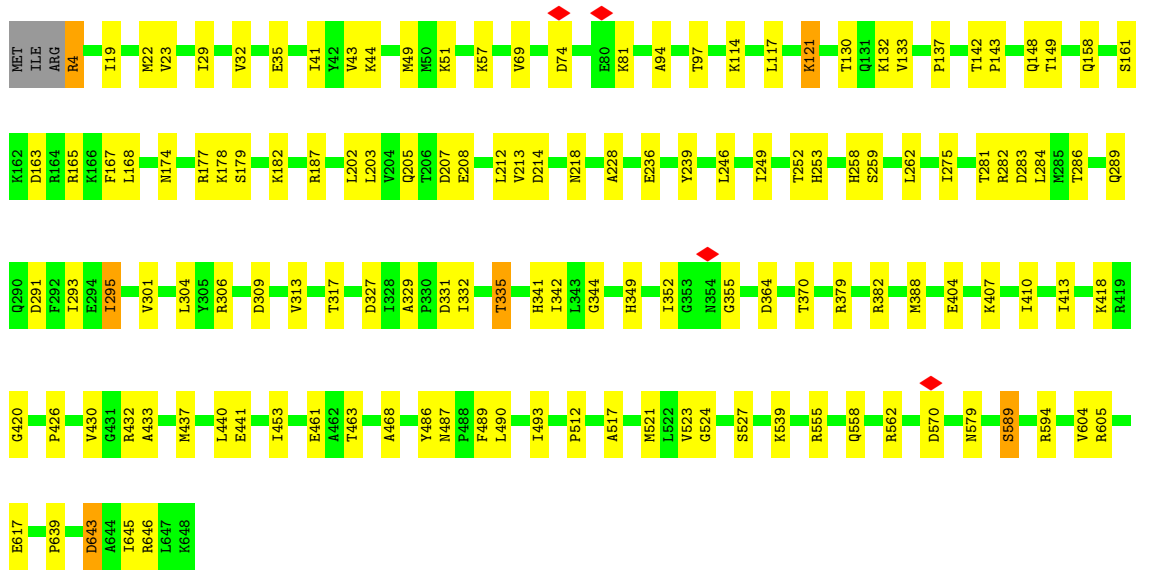
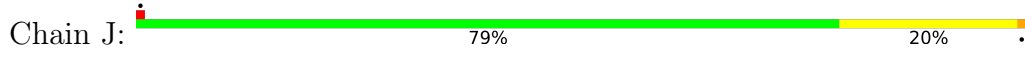
- Molecule 12: Transcription elongation factor Spt5



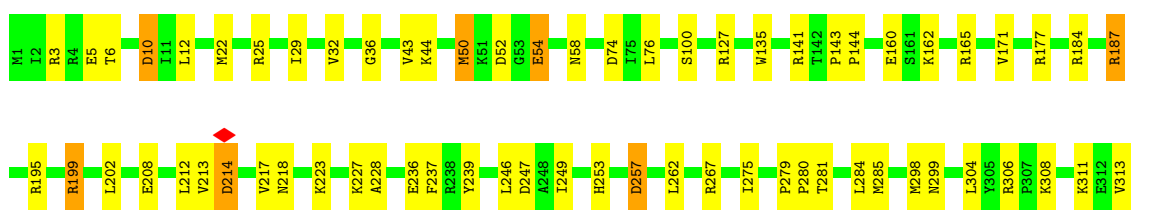
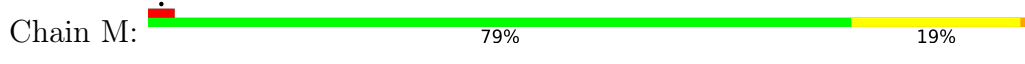
• Molecule 13: Transcription elongation factor Spt4

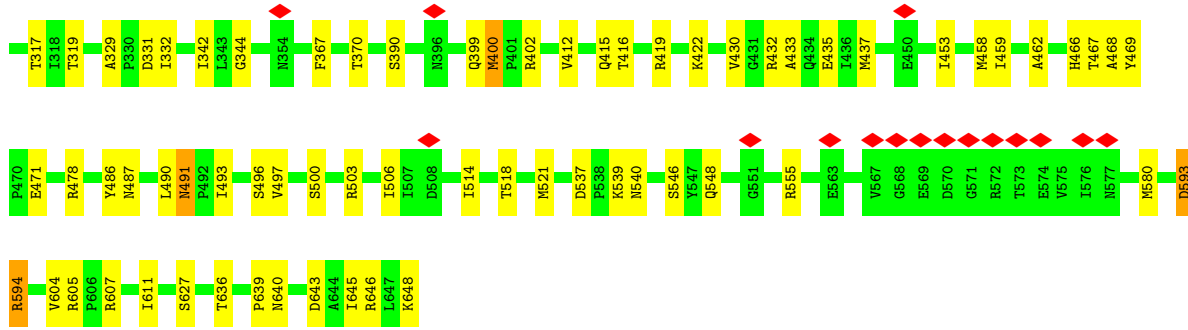


• Molecule 14: Transcription termination factor FttA

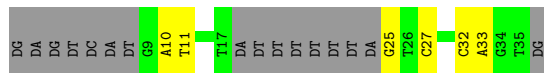


• Molecule 14: Transcription termination factor FttA

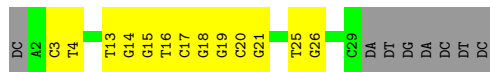




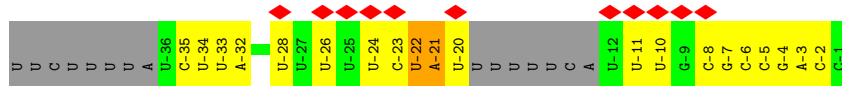
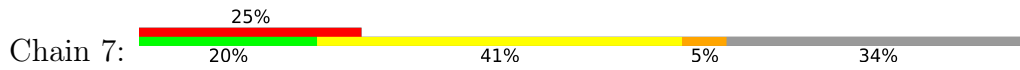
• Molecule 15: non-template strand DNA



• Molecule 16: template strand DNA



• Molecule 17: RNA



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	171269	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	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)	2200	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.657	Depositor
Minimum map value	-0.135	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.020	Depositor
Recommended contour level	0.088	Depositor
Map size ( $\text{\AA}$ )	330.0, 330.0, 330.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.825, 0.825, 0.825	Depositor

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ZN

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	0.25	0/7325	0.50	0/9887
2	B	0.25	0/8981	0.52	0/12135
3	C	0.24	0/2994	0.50	0/4042
4	D	0.24	0/2098	0.46	0/2841
5	E	0.26	0/1502	0.53	0/2023
6	F	0.24	0/1020	0.52	0/1373
7	H	0.27	0/650	0.49	0/877
8	K	0.24	0/441	0.49	0/598
9	L	0.24	0/791	0.50	0/1066
10	N	0.26	0/539	0.48	0/723
11	P	0.27	0/378	0.63	0/507
12	G	0.25	0/1164	0.52	0/1577
13	I	0.26	0/524	0.57	0/708
14	J	0.25	0/5253	0.51	1/7115 (0.0%)
14	M	0.25	0/5270	0.52	1/7139 (0.0%)
15	5	0.47	0/458	0.91	0/702
16	6	0.47	0/643	0.90	0/991
17	7	0.14	0/657	0.74	0/1014
All	All	0.26	0/40688	0.53	2/55318 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	M	214	ASP	CB-CG-OD1	7.10	124.69	118.30
14	J	214	ASP	CB-CG-OD1	5.30	123.07	118.30

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	A	7182	0	7247	117	0
2	B	8804	0	8858	148	0
3	C	2958	0	3077	52	0
4	D	2053	0	2063	27	0
5	E	1476	0	1512	31	0
6	F	1000	0	1000	14	0
7	H	636	0	646	11	0
8	K	433	0	466	6	0
9	L	776	0	770	14	0
10	N	530	0	540	8	0
11	P	373	0	392	11	0
12	G	1142	0	1192	20	0
13	I	515	0	506	8	0
14	J	5144	0	5246	64	0
14	M	5161	0	5264	70	0
15	5	410	0	228	4	0
16	6	575	0	317	16	0
17	7	595	0	304	13	0
18	A	1	0	0	0	0
19	A	2	0	0	0	0
19	B	1	0	0	0	0
19	I	1	0	0	0	0
19	J	1	0	0	0	0
19	M	1	0	0	0	0
19	N	1	0	0	0	0
19	P	1	0	0	0	0
All	All	39772	0	39628	543	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 7.

All (543) 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)
5:E:112:GLN:HE22	5:E:164:THR:HA	1.40	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:196:SER:HG	2:B:305:HIS:HD1	1.38	0.70
14:J:329:ALA:HB3	14:J:332:ILE:HB	1.74	0.70
2:B:820:ARG:HD3	2:B:820:ARG:H	1.56	0.70
1:A:432:ARG:HA	2:B:1039:VAL:HB	1.75	0.68
16:6:17:DC:H42	17:7:-4:G:H1	1.37	0.68
14:J:335:THR:HG23	14:J:349:HIS:HB3	1.76	0.67
2:B:450:GLN:OE1	15:5:25:DG:N2	2.27	0.67
12:G:26:LYS:O	12:G:30:TYR:HB2	1.95	0.67
1:A:568:THR:HA	1:A:611:ILE:O	1.94	0.67
1:A:319:LEU:HB3	3:C:366:LEU:HD22	1.75	0.66
14:J:487:ASN:HB3	14:J:490:LEU:HB2	1.78	0.65
1:A:54:ILE:HD11	1:A:224:THR:HA	1.78	0.65
5:E:127:PHE:HB2	5:E:136:LEU:HB3	1.78	0.65
14:J:202:LEU:HB3	14:J:213:VAL:HB	1.78	0.64
6:F:87:GLU:HB3	6:F:119:ARG:HH12	1.62	0.64
7:H:65:ARG:HD2	7:H:73:TYR:HD2	1.63	0.64
1:A:198:LEU:HD12	1:A:205:PRO:HA	1.80	0.64
2:B:339:ASP:HB3	2:B:342:HIS:HB2	1.79	0.64
9:L:9:GLU:HB2	9:L:12:LEU:HB3	1.80	0.64
2:B:700:PRO:HB2	2:B:719:PRO:HG2	1.78	0.64
14:M:422:LYS:H	14:M:540:ASN:HB3	1.63	0.64
2:B:758:ARG:NH1	16:6:19:DG:OP1	2.30	0.63
2:B:206:THR:HB	2:B:218:THR:HB	1.80	0.63
2:B:564:GLU:N	2:B:564:GLU:OE2	2.32	0.62
1:A:463:ASP:HA	2:B:890:GLY:HA2	1.82	0.62
2:B:738:ASP:OD1	2:B:922:ARG:NH2	2.33	0.62
14:J:332:ILE:HG12	14:J:352:ILE:HG12	1.81	0.62
2:B:452:HIS:HB3	2:B:455:ALA:HB3	1.82	0.62
12:G:145:LEU:HD21	12:G:148:LYS:HG2	1.80	0.61
14:M:486:TYR:OH	14:M:491:ASN:ND2	2.33	0.61
1:A:716:GLU:N	1:A:716:GLU:OE2	2.33	0.61
1:A:426:ARG:HD2	1:A:460:ALA:HB2	1.82	0.61
1:A:12:GLU:HB2	2:B:1116:SER:HB2	1.82	0.61
14:J:413:ILE:HD13	14:J:440:LEU:HD21	1.82	0.60
1:A:877:VAL:HG21	1:A:887:THR:HG23	1.84	0.60
2:B:614:VAL:HG12	2:B:615:ILE:HG13	1.82	0.60
2:B:1089:ASP:OD1	2:B:1089:ASP:N	2.35	0.59
3:C:125:ARG:HD3	3:C:240:GLU:HG2	1.83	0.59
14:M:54:GLU:OE2	14:M:58:ASN:ND2	2.33	0.59
3:C:227:SER:O	3:C:249:ASN:ND2	2.35	0.59
5:E:7:VAL:HG11	6:F:2:ILE:HG22	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:J:44:LYS:NZ	14:J:74:ASP:OD2	2.35	0.59
14:M:212:LEU:HB3	14:M:249:ILE:HG12	1.84	0.59
10:N:16:ASP:OD1	10:N:16:ASP:N	2.36	0.59
2:B:1051:GLU:HA	2:B:1055:GLU:HB2	1.86	0.58
14:M:202:LEU:HB3	14:M:213:VAL:HB	1.85	0.58
3:C:175:MET:O	3:C:178:VAL:HB	2.03	0.58
14:J:420:GLY:HA2	14:J:512:PRO:HG3	1.85	0.58
1:A:54:ILE:HG12	1:A:223:VAL:HG12	1.85	0.58
1:A:316:ARG:NH2	16:6:14:DG:OP1	2.37	0.58
1:A:160:GLU:HB3	1:A:164:ILE:HB	1.86	0.58
2:B:310:PRO:O	2:B:313:ARG:HB2	2.03	0.58
1:A:240:ASP:HB3	1:A:301:LEU:HB3	1.85	0.57
2:B:465:ARG:NH2	2:B:623:GLU:OE1	2.34	0.57
1:A:335:ASP:HB2	1:A:449:ARG:HG2	1.85	0.57
1:A:789:THR:HG21	2:B:623:GLU:HG2	1.85	0.57
14:J:212:LEU:HB3	14:J:249:ILE:HG12	1.86	0.57
2:B:448:ARG:O	2:B:456:ARG:NH2	2.35	0.57
14:M:471:GLU:HA	14:M:478:ARG:HD3	1.87	0.57
1:A:204:HIS:H	1:A:208:SER:HB2	1.69	0.57
3:C:161:VAL:HB	3:C:198:LEU:HB3	1.86	0.57
6:F:31:GLU:HG3	6:F:33:PRO:HD2	1.84	0.57
2:B:880:ASP:OD2	2:B:996:ARG:NH2	2.38	0.56
4:D:64:ALA:HB1	4:D:145:LYS:HD3	1.88	0.56
17:7:-5:C:H2'	17:7:-4:G:C8	2.40	0.56
5:E:82:GLN:HE22	5:E:147:ILE:HG22	1.71	0.56
1:A:20:GLU:OE1	2:B:1114:ARG:NH1	2.37	0.56
2:B:253:ILE:HG23	2:B:328:LEU:HD12	1.87	0.56
14:M:422:LYS:NZ	14:M:506:ILE:O	2.38	0.56
14:J:121:LYS:HE2	14:J:121:LYS:H	1.71	0.56
14:M:453:ILE:HD13	14:M:493:ILE:HD13	1.88	0.55
3:C:323:ARG:HD3	3:C:340:GLU:HG2	1.88	0.55
6:F:109:GLU:C	6:F:113:LYS:HZ2	2.09	0.55
9:L:33:HIS:NE2	9:L:41:ALA:O	2.39	0.55
4:D:91:LEU:HB2	4:D:104:LEU:HD23	1.89	0.55
1:A:879:PRO:HG3	3:C:303:ARG:HD2	1.87	0.55
3:C:160:VAL:HG13	3:C:199:VAL:HG22	1.89	0.55
7:H:67:SER:HB3	7:H:70:ALA:H	1.72	0.55
16:6:21:DG:H1	17:7:-8:C:H42	1.53	0.55
2:B:853:LYS:HG2	2:B:869:THR:HB	1.88	0.55
5:E:91:GLU:HB3	5:E:98:PHE:HB2	1.89	0.55
2:B:1006:HIS:NE2	2:B:1028:GLY:O	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:J:239:TYR:HB2	14:M:239:TYR:HB2	1.87	0.55
14:M:29:ILE:HA	14:M:43:VAL:HG12	1.88	0.55
3:C:360:GLY:O	3:C:364:ASN:ND2	2.40	0.55
1:A:51:LEU:HB2	1:A:223:VAL:HG23	1.89	0.54
1:A:85:HIS:HD2	1:A:87:GLY:H	1.55	0.54
1:A:771:GLN:NE2	1:A:772:GLN:O	2.40	0.54
5:E:115:ASP:OD1	5:E:115:ASP:N	2.40	0.54
1:A:781:ARG:NH1	2:B:454:GLU:O	2.40	0.54
2:B:119:VAL:HG22	2:B:391:GLN:HG3	1.90	0.54
4:D:96:PRO:HA	4:D:123:LEU:O	2.08	0.54
1:A:713:THR:OG1	1:A:714:LEU:N	2.40	0.54
3:C:62:GLU:OE2	8:K:6:ARG:NH1	2.40	0.54
5:E:150:SER:O	5:E:159:ASN:ND2	2.41	0.54
12:G:36:ALA:HB3	12:G:50:GLU:HB3	1.89	0.54
2:B:1064:VAL:HB	2:B:1119:LEU:HD11	1.90	0.54
16:6:25:DT:H2''	16:6:26:DG:N7	2.23	0.54
1:A:386:ASP:OD1	1:A:390:ARG:N	2.34	0.54
2:B:956:LYS:O	2:B:960:GLU:HG2	2.08	0.54
2:B:981:ARG:HG3	4:D:23:VAL:HG11	1.90	0.54
13:I:5:ARG:HB3	13:I:27:LEU:HG	1.90	0.53
14:M:177:ARG:HH21	14:M:208:GLU:HG3	1.73	0.53
1:A:818:PHE:O	1:A:822:MET:HG3	2.08	0.53
3:C:118:VAL:HA	3:C:264:THR:HG23	1.89	0.53
2:B:539:VAL:HG21	2:B:559:LEU:HB2	1.89	0.53
14:M:50:MET:SD	14:M:50:MET:N	2.81	0.53
5:E:20:MET:HE1	5:E:28:ILE:HD12	1.91	0.53
14:M:257:ASP:OD1	14:M:257:ASP:N	2.40	0.53
1:A:862:ASP:OD2	1:A:866:ILE:N	2.42	0.53
1:A:114:GLU:O	1:A:118:GLN:HG2	2.09	0.53
2:B:21:MET:HG3	2:B:645:MET:HB2	1.89	0.53
11:P:40:PRO:HB2	11:P:42:VAL:HG12	1.91	0.53
9:L:21:ASP:OD1	9:L:21:ASP:N	2.38	0.53
14:J:23:VAL:HG11	14:J:29:ILE:HD11	1.90	0.53
17:7:-8:C:H2'	17:7:-7:G:C8	2.44	0.53
1:A:156:PRO:HG2	1:A:168:LEU:HB2	1.91	0.52
2:B:680:GLY:HA3	2:B:720:ALA:HA	1.91	0.52
4:D:27:ASN:OD1	4:D:30:ARG:NH1	2.41	0.52
14:M:143:PRO:HB3	14:M:468:ALA:HB2	1.90	0.52
12:G:26:LYS:NZ	12:G:62:ARG:O	2.41	0.52
1:A:808:TYR:OH	1:A:820:HIS:NE2	2.40	0.52
2:B:186:ARG:NH1	15:5:27:DC:OP2	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:M:459:ILE:HD11	14:M:518:THR:HB	1.92	0.52
2:B:905:GLU:OE2	2:B:982:ARG:NH2	2.43	0.52
12:G:150:GLN:NE2	14:M:6:THR:O	2.43	0.52
13:I:18:ARG:HH12	13:I:23:GLY:HA2	1.75	0.52
14:J:291:ASP:O	14:J:295:ILE:HG23	2.09	0.52
1:A:266:LEU:HG	12:G:71:LEU:HD23	1.92	0.52
2:B:93:ARG:NH1	2:B:165:GLU:OE2	2.42	0.52
14:M:399:GLN:HG3	14:M:593:ASP:HB3	1.91	0.52
7:H:24:GLU:OE2	7:H:24:GLU:N	2.35	0.52
1:A:844:ARG:NH1	3:C:106:ASP:OD1	2.43	0.52
14:M:611:ILE:HB	14:M:636:THR:HG22	1.92	0.52
1:A:774:ILE:HG13	1:A:826:GLU:HG3	1.92	0.52
1:A:875:ASP:HB3	7:H:70:ALA:HB2	1.91	0.52
2:B:940:THR:HG23	2:B:964:LEU:HD11	1.92	0.52
2:B:606:TRP:HE1	2:B:617:TYR:HH	1.56	0.51
3:C:279:GLU:OE1	7:H:65:ARG:NH2	2.34	0.51
4:D:50:ASN:ND2	4:D:56:ASP:OD1	2.36	0.51
13:I:54:GLU:N	13:I:54:GLU:OE1	2.41	0.51
2:B:234:ALA:HA	2:B:270:GLN:HG3	1.93	0.51
5:E:46:LEU:HD21	5:E:103:PRO:HB3	1.92	0.51
2:B:852:ASP:HB2	2:B:871:ARG:HG2	1.93	0.51
10:N:36:ASP:OD2	10:N:46:ARG:NH2	2.44	0.51
14:J:426:PRO:HA	14:J:517:ALA:HB3	1.92	0.51
14:M:262:LEU:HD21	14:M:275:ILE:HG21	1.91	0.51
17:7:-4:G:H2'	17:7:-3:A:C8	2.45	0.51
14:J:94:ALA:HB2	14:J:117:LEU:HB2	1.92	0.51
1:A:147:CYS:HB3	1:A:150:CYS:SG	2.50	0.51
4:D:96:PRO:HG3	4:D:125:GLU:HG2	1.92	0.51
4:D:227:ASN:HD22	10:N:6:ARG:HH21	1.58	0.51
5:E:155:VAL:HB	5:E:158:GLU:HG3	1.92	0.51
14:J:562:ARG:HH21	14:J:579:ASN:HD22	1.59	0.51
14:M:212:LEU:HB2	14:M:246:LEU:HD11	1.92	0.51
12:G:37:ILE:HB	13:I:36:ILE:HB	1.93	0.51
4:D:14:ILE:HG12	4:D:16:PHE:HD1	1.76	0.51
4:D:56:ASP:HB3	11:P:46:VAL:HG21	1.93	0.51
2:B:802:PRO:HD2	11:P:40:PRO:HD3	1.93	0.50
16:6:18:DG:H2'	16:6:19:DG:H8	1.76	0.50
17:7:-5:C:H2'	17:7:-4:G:H8	1.76	0.50
14:J:236:GLU:OE2	14:M:187:ARG:NH2	2.35	0.50
14:J:313:VAL:O	14:J:317:THR:OG1	2.27	0.50
1:A:423:LEU:HB2	1:A:470:HIS:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:571:ASN:ND2	2:B:621:GLU:OE1	2.43	0.50
2:B:1074:GLU:HB2	2:B:1081:VAL:HG22	1.94	0.50
14:J:143:PRO:HB3	14:J:468:ALA:HB2	1.94	0.50
14:M:331:ASP:OD1	14:M:331:ASP:N	2.42	0.50
1:A:664:LYS:NZ	2:B:984:GLU:OE2	2.45	0.50
2:B:412:ALA:HB1	16:6:21:DG:H5"	1.92	0.50
2:B:1035:MET:O	2:B:1039:VAL:HG13	2.12	0.50
1:A:827:GLY:O	1:A:831:THR:HG23	2.11	0.50
12:G:93:PHE:O	14:M:3:ARG:NH2	2.45	0.50
12:G:150:GLN:NE2	14:M:5:GLU:O	2.45	0.50
1:A:476:GLU:OE2	8:K:11:ARG:NH2	2.45	0.50
14:M:267:ARG:NH1	14:M:304:LEU:O	2.45	0.50
1:A:195:ASP:OD1	1:A:209:ARG:NH1	2.45	0.50
2:B:844:ARG:NH2	16:6:19:DG:OP1	2.44	0.50
5:E:112:GLN:NE2	5:E:164:THR:HA	2.20	0.50
2:B:1077:ARG:HH22	2:B:1078:ARG:HH21	1.59	0.49
14:J:433:ALA:O	14:J:437:MET:HG3	2.12	0.49
7:H:50:VAL:HG13	7:H:55:ALA:HB3	1.93	0.49
2:B:881:LYS:HD3	2:B:997:LEU:HD12	1.94	0.49
14:J:282:ARG:O	14:J:286:THR:HG23	2.12	0.49
14:M:284:LEU:HD22	14:M:430:VAL:HG23	1.94	0.49
1:A:609:VAL:HA	1:A:619:GLY:HA3	1.94	0.49
11:P:41:ARG:HA	11:P:41:ARG:HH11	1.77	0.49
14:J:331:ASP:OD1	14:J:331:ASP:N	2.44	0.49
1:A:94:ARG:NH1	1:A:201:LEU:O	2.45	0.49
1:A:190:LEU:HD22	1:A:210:PRO:HB2	1.94	0.49
2:B:493:ILE:HD11	2:B:572:SER:HB2	1.94	0.49
14:J:306:ARG:NH2	14:J:309:ASP:OD1	2.44	0.49
14:M:50:MET:HG3	14:M:127:ARG:HB3	1.94	0.49
1:A:65:ALA:HB1	1:A:69:GLU:HG3	1.94	0.49
2:B:249:ASP:OD1	2:B:249:ASP:N	2.45	0.49
4:D:151:VAL:HG23	4:D:224:VAL:HG22	1.94	0.49
14:J:130:THR:HB	14:J:137:PRO:HD3	1.95	0.49
14:J:281:THR:OG1	14:J:344:GLY:O	2.31	0.49
1:A:247:ASP:OD1	1:A:250:ARG:NH2	2.45	0.49
2:B:237:LEU:HD23	2:B:314:ARG:HH21	1.77	0.49
8:K:31:PRO:HD2	8:K:34:ILE:HG12	1.94	0.49
10:N:47:ARG:O	10:N:51:SER:OG	2.22	0.49
12:G:59:GLU:HG2	12:G:62:ARG:HH21	1.77	0.49
14:M:400:MET:H	14:M:400:MET:CE	2.26	0.49
14:M:593:ASP:OD1	14:M:593:ASP:N	2.37	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:445:TYR:HB2	1:A:449:ARG:HH12	1.77	0.48
17:7:-7:G:H2'	17:7:-6:C:H6	1.78	0.48
1:A:323:ARG:NH2	2:B:1029:GLY:O	2.37	0.48
1:A:423:LEU:HB3	1:A:435:ILE:HD11	1.94	0.48
1:A:514:LEU:HB2	1:A:660:VAL:HG21	1.95	0.48
1:A:712:LYS:HE3	1:A:716:GLU:HG2	1.95	0.48
2:B:579:ARG:NH1	2:B:643:GLU:OE2	2.41	0.48
9:L:8:ARG:NH1	9:L:68:GLU:OE2	2.46	0.48
14:M:453:ILE:HG13	14:M:514:ILE:HB	1.95	0.48
1:A:527:GLU:OE1	9:L:33:HIS:NE2	2.46	0.48
1:A:808:TYR:O	2:B:662:HIS:ND1	2.44	0.48
6:F:18:LYS:HB2	6:F:48:ALA:HB1	1.96	0.48
1:A:323:ARG:HA	2:B:1031:ARG:HA	1.94	0.48
1:A:809:LYS:HG2	2:B:925:VAL:HG11	1.95	0.48
3:C:226:LEU:HD22	3:C:252:GLN:HB3	1.95	0.48
14:J:441:GLU:OE2	14:J:486:TYR:OH	2.32	0.48
16:6:14:DG:H2'	16:6:15:DG:C8	2.49	0.48
14:J:49:MET:HG2	14:J:69:VAL:HG11	1.94	0.48
2:B:740:VAL:HG12	2:B:914:VAL:HG12	1.94	0.48
14:M:160:GLU:OE2	14:M:311:LYS:NZ	2.43	0.48
1:A:599:GLU:HA	1:A:602:LYS:HE3	1.94	0.48
3:C:336:ARG:NH2	3:C:354:GLU:OE1	2.46	0.48
1:A:80:ALA:HB2	1:A:256:LYS:HG3	1.95	0.48
8:K:13:ILE:HD11	8:K:44:PHE:HB2	1.96	0.48
14:J:262:LEU:HD21	14:J:275:ILE:HG21	1.95	0.48
14:J:453:ILE:HD12	14:J:493:ILE:HD12	1.96	0.48
1:A:853:LYS:HB3	1:A:853:LYS:HE3	1.70	0.48
2:B:772:LYS:NZ	2:B:774:ASN:OD1	2.44	0.48
2:B:1108:LEU:HD21	3:C:346:LEU:HD13	1.96	0.48
5:E:127:PHE:HD2	5:E:136:LEU:HD23	1.79	0.47
14:J:259:SER:O	14:J:289:GLN:NE2	2.41	0.47
14:J:295:ILE:HD11	17:7:-22:U:C4	2.49	0.47
1:A:341:ASN:O	1:A:439:ARG:N	2.45	0.47
2:B:1005:MET:O	2:B:1024:ARG:NH1	2.41	0.47
14:J:187:ARG:NH2	14:M:236:GLU:OE1	2.42	0.47
14:J:555:ARG:NH1	14:J:558:GLN:OE1	2.45	0.47
14:M:10:ASP:N	14:M:10:ASP:OD1	2.47	0.47
2:B:581:LEU:HD12	2:B:618:LEU:HD12	1.96	0.47
2:B:659:TYR:HB3	2:B:662:HIS:HD2	1.79	0.47
14:J:252:THR:HG1	14:J:258:HIS:HD1	1.61	0.47
5:E:53:GLU:HB2	5:E:71:ILE:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:ARG:NH1	16:6:16:DT:OP1	2.43	0.47
2:B:257:LEU:O	2:B:261:LEU:HG	2.14	0.47
2:B:800:ILE:HD13	2:B:812:LEU:HD23	1.97	0.47
3:C:76:SER:HB3	3:C:98:LEU:HD13	1.96	0.47
5:E:54:ILE:HG13	5:E:70:ALA:HB2	1.97	0.47
1:A:195:ASP:HB3	1:A:205:PRO:HB3	1.97	0.47
3:C:103:GLU:HG2	3:C:108:ARG:HD2	1.96	0.47
3:C:384:LEU:HD11	5:E:13:ILE:HG21	1.96	0.47
6:F:90:LEU:O	6:F:94:VAL:HG13	2.14	0.47
12:G:144:ARG:NH1	12:G:146:ILE:HG22	2.29	0.47
14:M:12:LEU:HD22	14:M:32:VAL:HG12	1.96	0.47
14:M:648:LYS:HB3	14:M:648:LYS:HE3	1.76	0.47
3:C:116:MET:HA	3:C:266:THR:HA	1.97	0.47
9:L:83:ARG:HG2	9:L:83:ARG:HH11	1.80	0.47
2:B:726:VAL:HG22	2:B:912:LEU:HB3	1.97	0.47
2:B:962:GLU:OE2	4:D:201:ARG:N	2.48	0.47
14:J:97:THR:HG21	14:J:142:THR:HG21	1.96	0.47
14:J:212:LEU:HB2	14:J:246:LEU:HD11	1.97	0.47
1:A:432:ARG:O	2:B:1043:HIS:NE2	2.48	0.46
2:B:96:ASN:HA	2:B:163:GLY:HA3	1.98	0.46
2:B:172:GLU:OE2	2:B:481:LYS:NZ	2.41	0.46
14:J:355:GLY:O	14:J:382:ARG:NH2	2.49	0.46
4:D:64:ALA:HB2	11:P:48:ALA:HB1	1.97	0.46
5:E:35:GLU:HB3	6:F:38:PHE:HE2	1.80	0.46
3:C:133:GLU:OE2	3:C:137:ARG:NH2	2.48	0.46
13:I:39:ASP:OD1	13:I:42:SER:N	2.48	0.46
14:J:207:ASP:OD1	14:J:207:ASP:N	2.44	0.46
1:A:187:ARG:HG3	1:A:214:VAL:HB	1.98	0.46
2:B:85:ARG:HH22	11:P:33:LYS:HG3	1.81	0.46
12:G:38:LEU:HG	12:G:40:PRO:HD3	1.98	0.46
14:M:500:SER:HA	14:M:503:ARG:HG2	1.96	0.46
12:G:9:VAL:HG22	12:G:70:VAL:HG22	1.95	0.46
1:A:423:LEU:HD21	2:B:1048:LEU:HD21	1.98	0.46
2:B:239:THR:HG23	2:B:242:GLU:H	1.81	0.46
3:C:147:ALA:HA	3:C:163:ILE:HG22	1.97	0.46
3:C:182:LEU:HB3	3:C:186:PHE:HD2	1.81	0.46
3:C:250:PHE:HA	3:C:253:VAL:HG12	1.97	0.46
5:E:48:ILE:HD13	5:E:74:VAL:HG13	1.98	0.46
12:G:34:LEU:HD12	12:G:34:LEU:HA	1.80	0.46
2:B:342:HIS:NE2	2:B:624:GLU:OE2	2.48	0.46
3:C:352:ARG:HB3	3:C:354:GLU:HG3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:M:412:VAL:HA	14:M:415:GLN:HE21	1.79	0.46
14:M:487:ASN:HB3	14:M:490:LEU:HB2	1.98	0.46
1:A:341:ASN:HB3	1:A:439:ARG:HD3	1.97	0.46
2:B:490:THR:HB	2:B:554:VAL:HG13	1.98	0.46
1:A:709:LEU:HG	1:A:717:THR:HG23	1.97	0.46
1:A:844:ARG:HG3	3:C:339:PHE:HZ	1.81	0.46
2:B:703:PRO:HA	2:B:723:ASN:HD21	1.81	0.46
3:C:121:ASP:OD1	3:C:121:ASP:N	2.47	0.46
4:D:154:LYS:NZ	4:D:155:TYR:O	2.46	0.46
14:M:546:SER:O	14:M:548:GLN:NE2	2.42	0.46
17:7:-3:A:H2'	17:7:-2:C:C6	2.50	0.46
1:A:264:PRO:HD2	1:A:267:ILE:HG13	1.98	0.46
14:M:177:ARG:NH1	14:M:247:ASP:OD1	2.40	0.46
1:A:759:ILE:HD13	1:A:759:ILE:HA	1.85	0.45
1:A:769:LEU:HD11	1:A:808:TYR:CZ	2.51	0.45
14:M:555:ARG:HA	14:M:555:ARG:HD2	1.79	0.45
1:A:189:ARG:HE	1:A:189:ARG:HB3	1.53	0.45
1:A:671:ASP:HB3	2:B:969:SER:HB3	1.98	0.45
1:A:861:ARG:HG2	1:A:867:ILE:HA	1.98	0.45
1:A:771:GLN:HA	1:A:803:PHE:HA	1.99	0.45
2:B:536:ARG:NH1	2:B:536:ARG:HB2	2.31	0.45
2:B:743:ASN:HB3	2:B:746:SER:HB2	1.98	0.45
4:D:174:LEU:HD11	4:D:178:ARG:HE	1.80	0.45
5:E:181:LYS:HB3	5:E:181:LYS:HE3	1.71	0.45
14:M:44:LYS:NZ	14:M:74:ASP:OD2	2.49	0.45
1:A:344:GLY:HA3	1:A:449:ARG:HB2	1.98	0.45
2:B:190:VAL:HG13	2:B:330:LEU:HD23	1.98	0.45
3:C:181:LYS:HA	3:C:181:LYS:HD2	1.78	0.45
5:E:88:GLU:OE1	6:F:50:ARG:NH1	2.47	0.45
1:A:667:THR:H	1:A:746:HIS:HD1	1.63	0.45
4:D:62:ARG:NH1	10:N:3:VAL:O	2.40	0.45
14:J:29:ILE:HD13	14:J:43:VAL:HG12	1.99	0.45
1:A:284:ASN:OD1	1:A:306:GLN:NE2	2.50	0.45
1:A:764:GLN:HA	1:A:768:MET:O	2.17	0.45
14:M:416:THR:HG23	14:M:419:ARG:HH21	1.82	0.45
14:M:537:ASP:OD1	14:M:539:LYS:NZ	2.49	0.45
1:A:902:ARG:HA	1:A:902:ARG:HD3	1.77	0.45
3:C:220:LYS:HA	3:C:220:LYS:HD3	1.78	0.45
3:C:276:LEU:HA	7:H:9:ILE:HD11	1.98	0.45
14:J:32:VAL:HG22	14:J:41:ILE:HG12	1.99	0.45
1:A:844:ARG:HB3	3:C:321:ILE:HD12	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:919:ILE:HG21	2:B:928:LEU:HD21	1.98	0.45
5:E:5:LEU:HD22	6:F:6:LYS:HD2	1.98	0.45
17:7:-6:C:H2'	17:7:-5:C:C6	2.52	0.45
2:B:93:ARG:HG3	2:B:160:ILE:HD13	1.99	0.45
2:B:856:VAL:HG11	11:P:19:LEU:HD11	1.99	0.45
12:G:141:GLU:HG2	12:G:142:TYR:HD1	1.81	0.45
13:I:26:ASP:OD1	13:I:26:ASP:N	2.45	0.45
14:J:57:LYS:HA	14:J:57:LYS:HD2	1.82	0.45
2:B:485:LEU:HD13	2:B:709:ILE:HG21	1.99	0.44
2:B:754:SER:OG	2:B:996:ARG:NH1	2.50	0.44
7:H:20:ILE:HD11	7:H:55:ALA:HB2	2.00	0.44
17:7:-7:G:H2'	17:7:-6:C:C6	2.53	0.44
7:H:65:ARG:HG3	7:H:73:TYR:HB3	1.99	0.44
1:A:671:ASP:OD1	2:B:971:ARG:NH1	2.48	0.44
1:A:869:GLN:NE2	3:C:310:ASP:OD2	2.51	0.44
2:B:24:TYR:HB2	2:B:606:TRP:CE3	2.53	0.44
14:J:19:ILE:O	14:J:23:VAL:N	2.41	0.44
2:B:817:SER:O	2:B:837:ARG:N	2.48	0.44
2:B:1118:ARG:HA	2:B:1118:ARG:HD3	1.81	0.44
12:G:27:ALA:HA	12:G:32:LEU:HD12	2.00	0.44
14:J:218:ASN:HB3	14:J:228:ALA:HA	1.99	0.44
14:M:433:ALA:O	14:M:437:MET:HG2	2.17	0.44
14:M:36:GLY:HA2	14:M:171:VAL:HG12	1.98	0.44
3:C:235:ARG:NH1	3:C:237:GLU:OE2	2.50	0.44
1:A:845:LEU:HG	3:C:321:ILE:HD13	2.00	0.44
5:E:84:VAL:HG22	5:E:145:ARG:HB3	2.00	0.44
14:M:144:PRO:HD2	14:M:467:THR:HB	2.00	0.44
14:M:462:ALA:O	14:M:466:HIS:ND1	2.38	0.44
14:M:639:PRO:HG3	14:M:645:ILE:HD11	1.99	0.44
1:A:457:PRO:HG2	1:A:501:ILE:HG12	2.00	0.44
14:J:132:LYS:HA	14:J:132:LYS:HD2	1.81	0.44
14:J:463:THR:HG23	14:J:489:PHE:HD2	1.82	0.43
2:B:426:LEU:HD13	16:6:20:DC:H5''	1.99	0.43
3:C:17:PRO:HD2	3:C:20:ILE:HD12	1.99	0.43
5:E:111:SER:O	5:E:166:ARG:NH1	2.51	0.43
6:F:62:LEU:HD21	6:F:115:ILE:HD13	1.99	0.43
6:F:13:THR:OG1	6:F:16:GLU:OE1	2.31	0.43
14:J:174:ASN:O	14:J:177:ARG:NH1	2.51	0.43
14:M:399:GLN:HE21	14:M:593:ASP:HB3	1.82	0.43
1:A:388:GLU:OE2	1:A:390:ARG:NH1	2.51	0.43
2:B:16:GLU:HG2	2:B:596:VAL:HG11	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:144:GLU:HB3	3:C:219:LYS:HE2	2.00	0.43
4:D:87:VAL:HG21	4:D:111:VAL:HG11	2.01	0.43
2:B:885:ARG:NH2	2:B:930:GLU:OE2	2.42	0.43
3:C:96:LEU:H	3:C:96:LEU:HG	1.57	0.43
1:A:266:LEU:HD12	1:A:266:LEU:HA	1.86	0.43
2:B:248:SER:HB2	2:B:325:LEU:HD11	2.00	0.43
2:B:689:ARG:NE	2:B:691:ASP:OD2	2.40	0.43
2:B:857:THR:OG1	2:B:858:GLU:N	2.52	0.43
4:D:254:LEU:HD23	4:D:254:LEU:HA	1.85	0.43
5:E:180:LYS:HE3	5:E:180:LYS:HB3	1.89	0.43
15:5:10:DA:H2''	15:5:11:DT:H5''	2.01	0.43
1:A:107:LYS:HD3	1:A:154:GLN:HG3	1.99	0.43
1:A:477:ALA:HB1	2:B:1048:LEU:HD13	2.01	0.43
1:A:822:MET:HG2	2:B:458:LEU:HD23	1.99	0.43
2:B:701:GLN:HB3	10:N:51:SER:HB2	2.01	0.43
4:D:28:ALA:HA	9:L:23:THR:HG23	2.01	0.43
5:E:51:VAL:HG22	5:E:72:PHE:HB3	2.01	0.43
11:P:26:ARG:HH21	11:P:31:GLY:HA2	1.84	0.43
12:G:34:LEU:HB3	13:I:38:LEU:HD12	2.01	0.43
16:6:16:DT:H2'	16:6:17:DC:C6	2.54	0.43
1:A:804:VAL:HG11	1:A:812:LEU:HD22	2.01	0.43
2:B:58:PRO:HA	2:B:371:GLN:HG3	2.00	0.43
2:B:703:PRO:HB3	2:B:993:TYR:CG	2.54	0.43
14:J:94:ALA:HA	14:J:114:LYS:HB2	2.00	0.43
1:A:221:PRO:HB3	1:A:225:MET:SD	2.59	0.43
1:A:480:GLU:OE1	2:B:1048:LEU:N	2.42	0.43
1:A:873:GLY:HA3	3:C:303:ARG:HG2	2.00	0.43
2:B:462:HIS:HB3	2:B:466:ILE:HB	2.01	0.43
2:B:941:GLY:HA2	10:N:50:LEU:HD11	2.00	0.43
3:C:92:ILE:HD11	3:C:231:LYS:HD3	2.00	0.43
2:B:171:ILE:HD11	2:B:423:VAL:HG22	2.00	0.43
2:B:817:SER:HB3	2:B:837:ARG:HB2	2.01	0.43
4:D:105:LYS:HD3	4:D:105:LYS:HA	1.80	0.43
5:E:45:VAL:HA	5:E:76:VAL:HG12	1.99	0.43
10:N:52:HIS:NE2	10:N:54:GLU:OE2	2.45	0.43
12:G:112:LYS:NZ	14:M:5:GLU:OE2	2.43	0.43
2:B:962:GLU:HG3	2:B:968:HIS:HB3	2.00	0.42
14:J:604:VAL:O	14:J:605:ARG:NE	2.53	0.42
14:M:370:THR:HB	14:M:402:ARG:HH12	1.84	0.42
1:A:295:HIS:CD2	1:A:296:LYS:H	2.37	0.42
5:E:3:LYS:HD3	5:E:5:LEU:HD21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:M:218:ASN:HB3	14:M:228:ALA:HA	2.01	0.42
2:B:844:ARG:HB2	2:B:847:GLU:HG3	2.00	0.42
7:H:43:ILE:HB	7:H:77:ARG:HD3	2.01	0.42
14:J:4:ARG:NH2	14:J:208:GLU:O	2.53	0.42
14:J:35:GLU:HG2	14:J:168:LEU:HB3	2.01	0.42
14:M:141:ARG:HH22	17:7:-35:C:H41	1.67	0.42
14:M:521:MET:HG2	14:M:548:GLN:HA	2.02	0.42
15:5:32:DC:H2''	15:5:33:DA:C8	2.55	0.42
1:A:852:LEU:HB2	3:C:65:GLY:HA3	2.01	0.42
2:B:64:LYS:HG2	2:B:110:VAL:HB	2.02	0.42
2:B:590:LYS:HB3	2:B:614:VAL:HG22	2.02	0.42
2:B:1031:ARG:HG2	16:6:15:DG:H5''	2.00	0.42
6:F:116:ASP:HA	6:F:119:ARG:HE	1.84	0.42
14:M:44:LYS:HA	14:M:44:LYS:HD2	1.74	0.42
14:M:253:HIS:HB2	14:M:342:ILE:HD11	2.01	0.42
14:M:281:THR:OG1	14:M:344:GLY:O	2.38	0.42
14:M:604:VAL:O	14:M:605:ARG:NE	2.52	0.42
2:B:708:ARG:NH2	2:B:945:ASP:OD2	2.49	0.42
2:B:967:LYS:NZ	4:D:199:LEU:O	2.47	0.42
2:B:1120:LYS:HB3	2:B:1120:LYS:HE3	1.86	0.42
12:G:146:ILE:HD12	12:G:147:SER:HB3	2.01	0.42
14:M:280:PRO:HG3	14:M:469:TYR:CG	2.55	0.42
4:D:161:VAL:HG22	4:D:211:ILE:HG12	2.00	0.42
9:L:27:LEU:O	9:L:31:THR:OG1	2.31	0.42
14:J:432:ARG:HH21	14:J:589:SER:HB2	1.85	0.42
1:A:23:LYS:HB3	2:B:1085:VAL:HA	2.01	0.42
1:A:857:ASP:OD2	1:A:861:ARG:NH2	2.52	0.42
2:B:659:TYR:HB3	2:B:662:HIS:CD2	2.54	0.42
2:B:851:VAL:HG22	2:B:870:VAL:HG22	2.01	0.42
16:6:13:DT:H2'	16:6:14:DG:C8	2.55	0.42
2:B:161:ILE:HG13	2:B:411:LEU:HB3	2.01	0.42
2:B:546:ARG:NH2	2:B:616:GLU:OE1	2.47	0.42
5:E:41:ASP:OD1	5:E:41:ASP:N	2.33	0.42
11:P:6:TYR:HA	11:P:38:PRO:HD3	2.01	0.42
14:J:341:HIS:HD2	14:J:342:ILE:HG12	1.84	0.42
1:A:493:SER:HB2	1:A:500:LEU:HB2	2.00	0.42
12:G:24:TYR:HB2	13:I:36:ILE:HD13	2.00	0.42
1:A:457:PRO:HG3	1:A:500:LEU:HG	2.01	0.41
1:A:890:VAL:HG11	3:C:24:LEU:HA	2.02	0.41
14:J:284:LEU:HD22	14:J:430:VAL:HG11	2.02	0.41
14:J:304:LEU:HD12	14:J:304:LEU:HA	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:M:329:ALA:HB3	14:M:332:ILE:HB	2.01	0.41
1:A:362:PHE:HB3	8:K:30:VAL:HG21	2.02	0.41
2:B:968:HIS:NE2	4:D:204:GLU:OE2	2.45	0.41
3:C:32:ASN:HA	3:C:37:LEU:HG	2.01	0.41
3:C:321:ILE:HG22	3:C:326:ILE:HD12	2.02	0.41
6:F:55:LYS:HD3	6:F:55:LYS:HA	1.73	0.41
14:J:643:ASP:OD1	14:M:646:ARG:N	2.49	0.41
14:M:594:ARG:NH2	14:M:627:SER:OG	2.53	0.41
2:B:237:LEU:HD23	2:B:314:ARG:NH2	2.35	0.41
3:C:386:LEU:HD11	8:K:50:PRO:HG3	2.02	0.41
14:M:279:PRO:HG3	14:M:319:THR:HB	2.02	0.41
14:M:313:VAL:O	14:M:317:THR:OG1	2.30	0.41
1:A:302:LYS:NZ	1:A:302:LYS:HB3	2.35	0.41
2:B:38:TYR:OH	2:B:127:PRO:O	2.32	0.41
2:B:381:LYS:HB3	2:B:385:TYR:HB3	2.02	0.41
2:B:803:GLU:OE1	2:B:871:ARG:NH2	2.54	0.41
2:B:803:GLU:OE2	11:P:39:ARG:NH2	2.54	0.41
14:J:639:PRO:HG3	14:J:645:ILE:HD11	2.01	0.41
2:B:725:VAL:O	2:B:911:ASP:N	2.54	0.41
9:L:33:HIS:CD2	9:L:41:ALA:HB3	2.55	0.41
14:M:52:ASP:N	14:M:52:ASP:OD1	2.53	0.41
14:M:165:ARG:HD3	14:M:165:ARG:HA	1.83	0.41
1:A:207:LYS:O	2:B:1114:ARG:NH2	2.54	0.41
2:B:181:VAL:HG21	2:B:326:LYS:HB3	2.02	0.41
2:B:731:TYR:HB3	2:B:741:ILE:HD13	2.01	0.41
2:B:1065:CYS:HB2	2:B:1083:CYS:HB2	2.03	0.41
14:J:253:HIS:CE1	14:J:364:ASP:HB2	2.56	0.41
1:A:468:ASN:OD1	1:A:468:ASN:N	2.54	0.41
1:A:846:ILE:HD11	3:C:70:GLN:HG3	2.03	0.41
2:B:208:GLU:HG3	2:B:216:TYR:HB2	2.02	0.41
2:B:427:LEU:HD11	2:B:435:THR:HG23	2.03	0.41
2:B:904:THR:HG22	2:B:974:MET:HG2	2.02	0.41
4:D:24:PRO:HB3	9:L:30:GLU:HG3	2.02	0.41
7:H:37:ILE:HD12	7:H:37:ILE:HA	1.84	0.41
14:J:158:GLN:O	14:J:161:SER:OG	2.39	0.41
14:J:404:GLU:HA	14:J:407:LYS:HE3	2.03	0.41
14:J:524:GLY:HA3	17:7:-21:A:N3	2.36	0.41
2:B:905:GLU:HG2	2:B:975:TYR:HE1	1.85	0.41
4:D:28:ALA:O	4:D:32:THR:OG1	2.31	0.41
4:D:209:LYS:HD2	4:D:209:LYS:H	1.86	0.41
5:E:147:ILE:HD13	5:E:164:THR:HB	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:107:GLU:O	6:F:110:GLU:HG2	2.21	0.41
16:6:18:DG:H2'	16:6:19:DG:C8	2.56	0.41
1:A:325:ASN:ND2	2:B:1056:GLU:O	2.54	0.41
1:A:525:GLU:OE2	1:A:664:LYS:NZ	2.45	0.41
2:B:170:SER:HB2	2:B:442:VAL:HG22	2.01	0.41
2:B:178:LYS:NZ	2:B:345:ASN:OD1	2.54	0.41
2:B:279:ARG:HA	2:B:290:ARG:HD2	2.01	0.41
2:B:390:ILE:HD12	2:B:390:ILE:HA	1.88	0.41
2:B:886:HIS:NE2	2:B:930:GLU:OE1	2.40	0.41
3:C:170:LYS:HA	3:C:170:LYS:HD2	1.81	0.41
3:C:361:VAL:HG13	3:C:371:VAL:HB	2.02	0.41
14:J:148:GLN:N	14:J:461:GLU:OE2	2.44	0.41
14:M:199:ARG:HB3	14:M:218:ASN:HB2	2.03	0.41
1:A:35:TYR:CZ	1:A:47:MET:HG3	2.55	0.41
3:C:382:MET:HB2	5:E:59:ILE:HD11	2.02	0.41
14:J:410:ILE:HD13	14:J:410:ILE:HA	1.87	0.41
14:M:76:LEU:HD22	14:M:135:TRP:CE2	2.56	0.41
1:A:95:VAL:HG13	1:A:201:LEU:HD13	2.02	0.40
1:A:147:CYS:CB	1:A:150:CYS:SG	3.08	0.40
1:A:563:LEU:HB3	1:A:567:LEU:HD23	2.03	0.40
1:A:570:TRP:HB3	1:A:610:TYR:HD1	1.86	0.40
2:B:17:LEU:HD11	2:B:632:PRO:HG3	2.02	0.40
2:B:235:LEU:HD12	2:B:235:LEU:HA	1.87	0.40
2:B:237:LEU:HD13	2:B:242:GLU:HB3	2.03	0.40
2:B:247:VAL:HB	2:B:257:LEU:HD11	2.03	0.40
2:B:680:GLY:HA2	2:B:697:MET:HB2	2.02	0.40
3:C:72:ILE:HD12	3:C:308:VAL:HG21	2.03	0.40
9:L:14:GLU:HA	9:L:57:PHE:O	2.21	0.40
9:L:16:TYR:CE2	9:L:56:ARG:HB2	2.56	0.40
14:J:203:LEU:HD23	14:J:205:GLN:HE21	1.85	0.40
2:B:1037:ARG:HD2	3:C:365:VAL:HG11	2.02	0.40
5:E:145:ARG:HD2	5:E:167:GLN:HE22	1.85	0.40
11:P:19:LEU:HD23	11:P:19:LEU:HA	1.94	0.40
14:M:412:VAL:HA	14:M:415:GLN:NE2	2.37	0.40
1:A:110:ASP:N	1:A:110:ASP:OD1	2.54	0.40
1:A:131:ASP:OD2	1:A:131:ASP:N	2.55	0.40
2:B:971:ARG:HB3	2:B:986:ASP:HB3	2.04	0.40
4:D:2:VAL:HG11	9:L:85:VAL:HG22	2.04	0.40
14:J:539:LYS:HB2	14:J:539:LYS:HE3	1.82	0.40
1:A:134:ILE:HD13	1:A:134:ILE:HA	1.91	0.40
1:A:450:LEU:HD11	1:A:454:VAL:HB	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:9:GLU:HA	2:B:10:PRO:HD3	1.94	0.40
5:E:178:LYS:HB2	5:E:178:LYS:HE2	1.86	0.40
16:6:3:DC:H2''	16:6:4:DT:H5'	2.03	0.40
1:A:658:ILE:O	1:A:662:THR:OG1	2.34	0.40
1:A:765:MET:HE2	1:A:765:MET:HB2	1.79	0.40
2:B:747:ILE:HG23	2:B:876:PRO:HD2	2.04	0.40
3:C:124:HIS:CD2	3:C:130:LYS:HB3	2.57	0.40
3:C:214:ILE:HD12	3:C:217:LYS:HD2	2.03	0.40
9:L:77:LYS:HE3	9:L:77:LYS:HB2	1.90	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	898/906 (99%)	881 (98%)	17 (2%)	0	100	100
2	B	1099/1123 (98%)	1076 (98%)	23 (2%)	0	100	100
3	C	374/391 (96%)	363 (97%)	11 (3%)	0	100	100
4	D	254/261 (97%)	251 (99%)	3 (1%)	0	100	100
5	E	181/190 (95%)	179 (99%)	2 (1%)	0	100	100
6	F	118/122 (97%)	114 (97%)	4 (3%)	0	100	100
7	H	75/82 (92%)	75 (100%)	0	0	100	100
8	K	54/60 (90%)	50 (93%)	4 (7%)	0	100	100
9	L	92/94 (98%)	90 (98%)	2 (2%)	0	100	100
10	N	63/65 (97%)	63 (100%)	0	0	100	100
11	P	46/49 (94%)	44 (96%)	2 (4%)	0	100	100
12	G	145/152 (95%)	143 (99%)	2 (1%)	0	100	100
13	I	63/73 (86%)	62 (98%)	1 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
14	J	643/648 (99%)	628 (98%)	15 (2%)	0	100	100
14	M	646/648 (100%)	625 (97%)	21 (3%)	0	100	100
All	All	4751/4864 (98%)	4644 (98%)	107 (2%)	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	A	774/779 (99%)	743 (96%)	31 (4%)	31	56
2	B	949/969 (98%)	908 (96%)	41 (4%)	29	53
3	C	324/334 (97%)	301 (93%)	23 (7%)	14	28
4	D	226/230 (98%)	214 (95%)	12 (5%)	22	43
5	E	160/167 (96%)	147 (92%)	13 (8%)	11	23
6	F	104/107 (97%)	97 (93%)	7 (7%)	16	31
7	H	69/72 (96%)	64 (93%)	5 (7%)	14	28
8	K	45/49 (92%)	42 (93%)	3 (7%)	16	31
9	L	81/81 (100%)	75 (93%)	6 (7%)	13	27
10	N	59/59 (100%)	54 (92%)	5 (8%)	10	21
11	P	38/40 (95%)	33 (87%)	5 (13%)	4	7
12	G	127/131 (97%)	118 (93%)	9 (7%)	14	28
13	I	56/65 (86%)	52 (93%)	4 (7%)	14	28
14	J	562/565 (100%)	530 (94%)	32 (6%)	20	39
14	M	562/565 (100%)	525 (93%)	37 (7%)	16	32
All	All	4136/4213 (98%)	3903 (94%)	233 (6%)	25	40

All (233) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	ARG
1	A	100	CYS
1	A	110	ASP
1	A	123	MET
1	A	196	LYS
1	A	200	LEU
1	A	220	VAL
1	A	233	SER
1	A	281	THR
1	A	288	SER
1	A	302	LYS
1	A	388	GLU
1	A	424	PHE
1	A	431	HIS
1	A	434	SER
1	A	465	ASP
1	A	468	ASN
1	A	469	LEU
1	A	486	GLU
1	A	506	ASP
1	A	581	ARG
1	A	658	ILE
1	A	684	HIS
1	A	688	ARG
1	A	696	ARG
1	A	759	ILE
1	A	768	MET
1	A	850	GLN
1	A	853	LYS
1	A	855	ASP
1	A	871	LYS
2	B	28	LYS
2	B	56	VAL
2	B	66	LYS
2	B	69	LYS
2	B	78	GLN
2	B	121	VAL
2	B	153	LYS
2	B	183	ARG
2	B	186	ARG
2	B	196	SER
2	B	198	ARG
2	B	245	GLU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
2	B	261	LEU
2	B	280	LEU
2	B	296	HIS
2	B	308	VAL
2	B	335	ARG
2	B	338	ASP
2	B	343	TYR
2	B	355	LEU
2	B	368	LYS
2	B	384	ARG
2	B	403	LEU
2	B	416	TRP
2	B	504	LYS
2	B	566	LYS
2	B	585	GLU
2	B	612	MET
2	B	644	LEU
2	B	693	ARG
2	B	711	LYS
2	B	742	ILE
2	B	801	PHE
2	B	820	ARG
2	B	920	PRO
2	B	943	ARG
2	B	959	LYS
2	B	982	ARG
2	B	1031	ARG
2	B	1083	CYS
2	B	1089	ASP
3	C	5	LYS
3	C	33	GLU
3	C	94	VAL
3	C	96	LEU
3	C	108	ARG
3	C	128	ARG
3	C	136	ARG
3	C	144	GLU
3	C	145	ASN
3	C	157	MET
3	C	159	TYR
3	C	180	ARG
3	C	187	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
3	C	207	LYS
3	C	210	ASP
3	C	224	LYS
3	C	245	THR
3	C	250	PHE
3	C	340	GLU
3	C	344	GLN
3	C	345	HIS
3	C	352	ARG
3	C	384	LEU
4	D	76	SER
4	D	78	ASP
4	D	84	ASP
4	D	112	LYS
4	D	115	ASN
4	D	128	ARG
4	D	177	ARG
4	D	178	ARG
4	D	183	GLU
4	D	207	MET
4	D	209	LYS
4	D	214	GLU
5	E	16	ARG
5	E	32	GLU
5	E	39	ASP
5	E	41	ASP
5	E	69	GLU
5	E	105	ASP
5	E	115	ASP
5	E	118	VAL
5	E	124	ASN
5	E	153	SER
5	E	166	ARG
5	E	172	LYS
5	E	182	LYS
6	F	37	MET
6	F	43	VAL
6	F	44	SER
6	F	63	LYS
6	F	89	TYR
6	F	98	LYS
6	F	113	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
7	H	6	GLU
7	H	23	GLU
7	H	56	LYS
7	H	66	LYS
7	H	67	SER
8	K	6	ARG
8	K	11	ARG
8	K	46	LYS
9	L	6	ILE
9	L	8	ARG
9	L	21	ASP
9	L	64	LYS
9	L	66	THR
9	L	87	GLU
10	N	1	MET
10	N	22	LYS
10	N	33	LYS
10	N	37	ASP
10	N	47	ARG
11	P	10	LYS
11	P	23	ARG
11	P	26	ARG
11	P	41	ARG
11	P	44	ARG
12	G	25	SER
12	G	44	LYS
12	G	59	GLU
12	G	84	LEU
12	G	90	VAL
12	G	106	PHE
12	G	116	VAL
12	G	147	SER
12	G	149	ARG
13	I	39	ASP
13	I	41	GLU
13	I	47	LYS
13	I	65	ARG
14	J	4	ARG
14	J	22	MET
14	J	51	LYS
14	J	81	LYS
14	J	121	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
14	J	133	VAL
14	J	149	THR
14	J	163	ASP
14	J	165	ARG
14	J	167	PHE
14	J	178	LYS
14	J	179	SER
14	J	182	LYS
14	J	283	ASP
14	J	293	ILE
14	J	295	ILE
14	J	301	VAL
14	J	327	ASP
14	J	335	THR
14	J	370	THR
14	J	379	ARG
14	J	388	MET
14	J	418	LYS
14	J	521	MET
14	J	523	VAL
14	J	527	SER
14	J	570	ASP
14	J	589	SER
14	J	594	ARG
14	J	617	GLU
14	J	643	ASP
14	J	646	ARG
14	M	10	ASP
14	M	22	MET
14	M	25	ARG
14	M	50	MET
14	M	54	GLU
14	M	100	SER
14	M	162	LYS
14	M	184	ARG
14	M	187	ARG
14	M	195	ARG
14	M	199	ARG
14	M	214	ASP
14	M	217	VAL
14	M	223	LYS
14	M	227	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
14	M	237	PHE
14	M	257	ASP
14	M	285	MET
14	M	298	MET
14	M	299	ASN
14	M	306	ARG
14	M	308	LYS
14	M	367	PHE
14	M	390	SER
14	M	400	MET
14	M	432	ARG
14	M	435	GLU
14	M	458	MET
14	M	491	ASN
14	M	496	SER
14	M	497	VAL
14	M	580	MET
14	M	593	ASP
14	M	594	ARG
14	M	607	ARG
14	M	640	ASN
14	M	643	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	341	ASN
2	B	62	ASN
2	B	112	ASN
2	B	221	ASN
2	B	301	ASN
3	C	267	ASN
4	D	115	ASN
4	D	117	ASN
5	E	73	ASN
5	E	82	GLN
5	E	112	GLN
5	E	167	GLN
14	J	88	GLN
14	J	253	HIS
14	J	341	HIS
14	J	483	HIS

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Mol	Chain	Res	Type
14	J	579	ASN
14	M	399	GLN
14	M	483	HIS

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
17	7	27/44 (61%)	12 (44%)	0

All (12) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
17	7	-34	U
17	7	-33	U
17	7	-32	A
17	7	-28	U
17	7	-26	U
17	7	-24	U
17	7	-23	C
17	7	-22	U
17	7	-21	A
17	7	-20	U
17	7	-11	U
17	7	-10	U

There are no RNA pucker outliers to report.

### 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](#)

Of 9 ligands modelled in this entry, 9 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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.

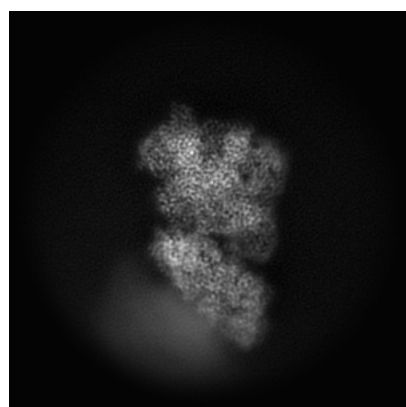
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-44438. These allow visual inspection of the internal detail of the map and identification of artifacts.

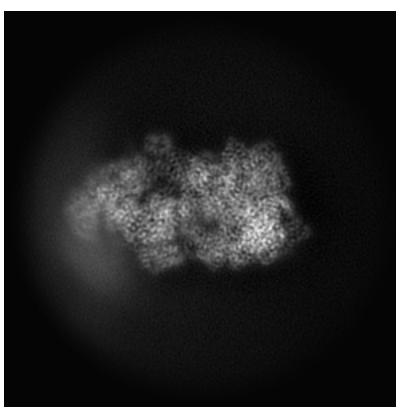
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

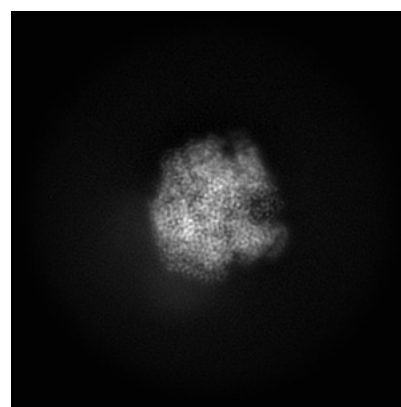
#### 6.1.1 Primary map



X



Y

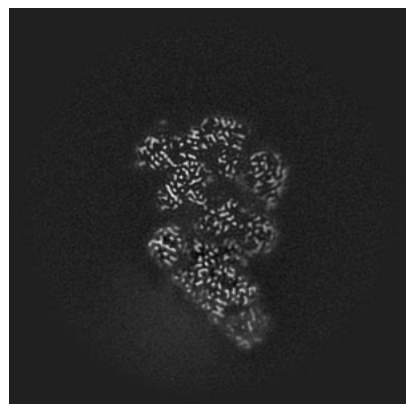


Z

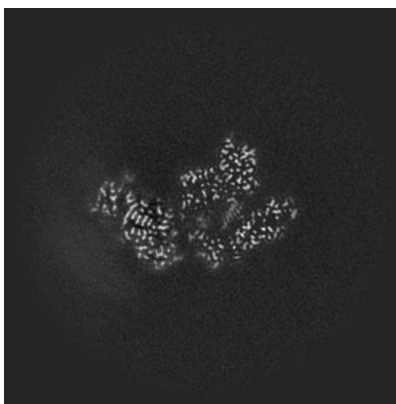
The images above show the map projected in three orthogonal directions.

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

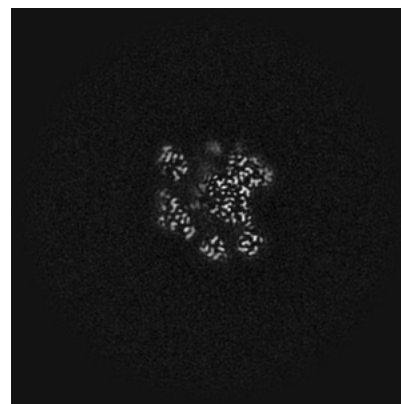
#### 6.2.1 Primary map



X Index: 200



Y Index: 200

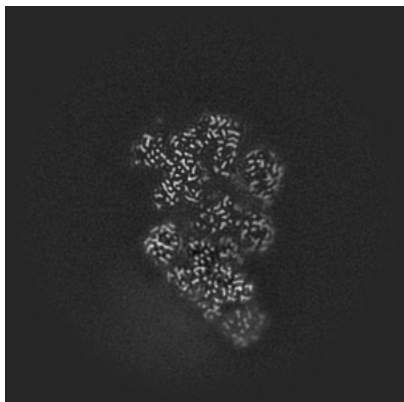


Z Index: 200

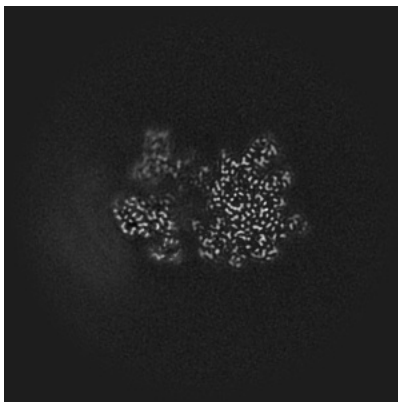
The images above show central slices of the map in three orthogonal directions.

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

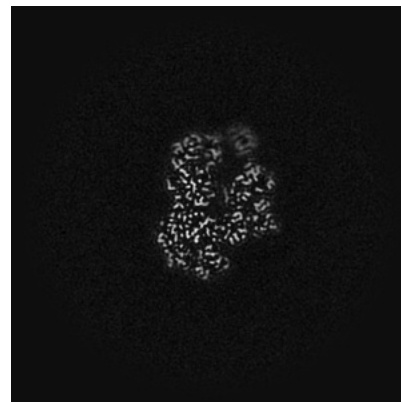
### 6.3.1 Primary map



X Index: 199



Y Index: 183

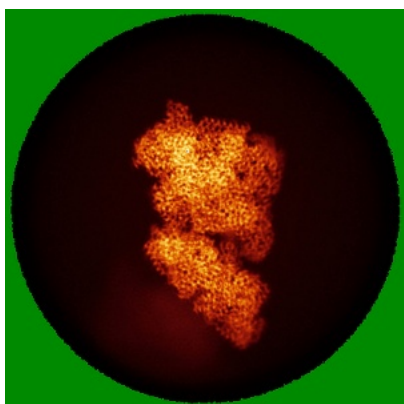


Z Index: 249

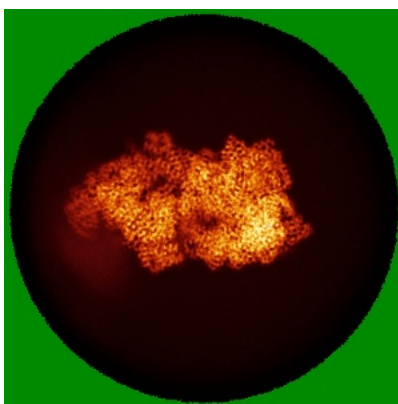
The images above show the largest variance slices of the map in three orthogonal directions.

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

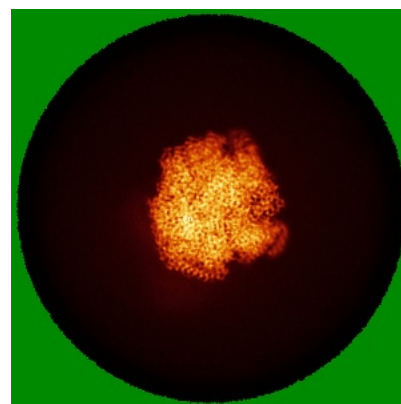
### 6.4.1 Primary map



X



Y



Z

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

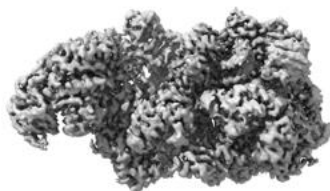


## 6.5 Orthogonal surface views [i](#)

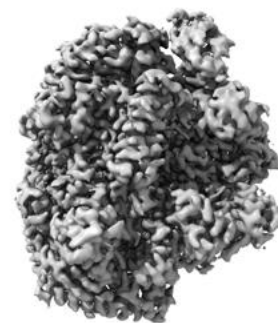
### 6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.088. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

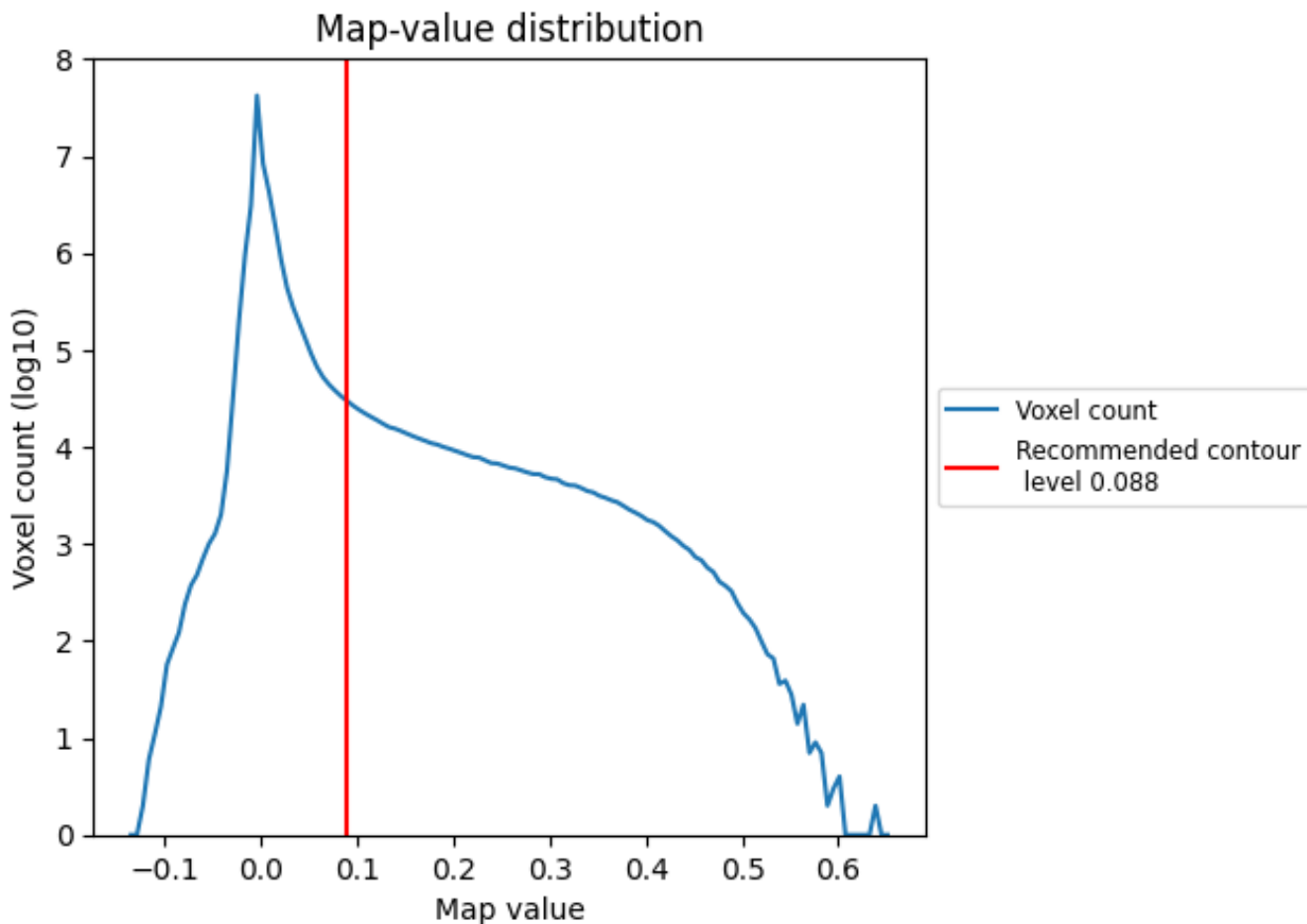
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

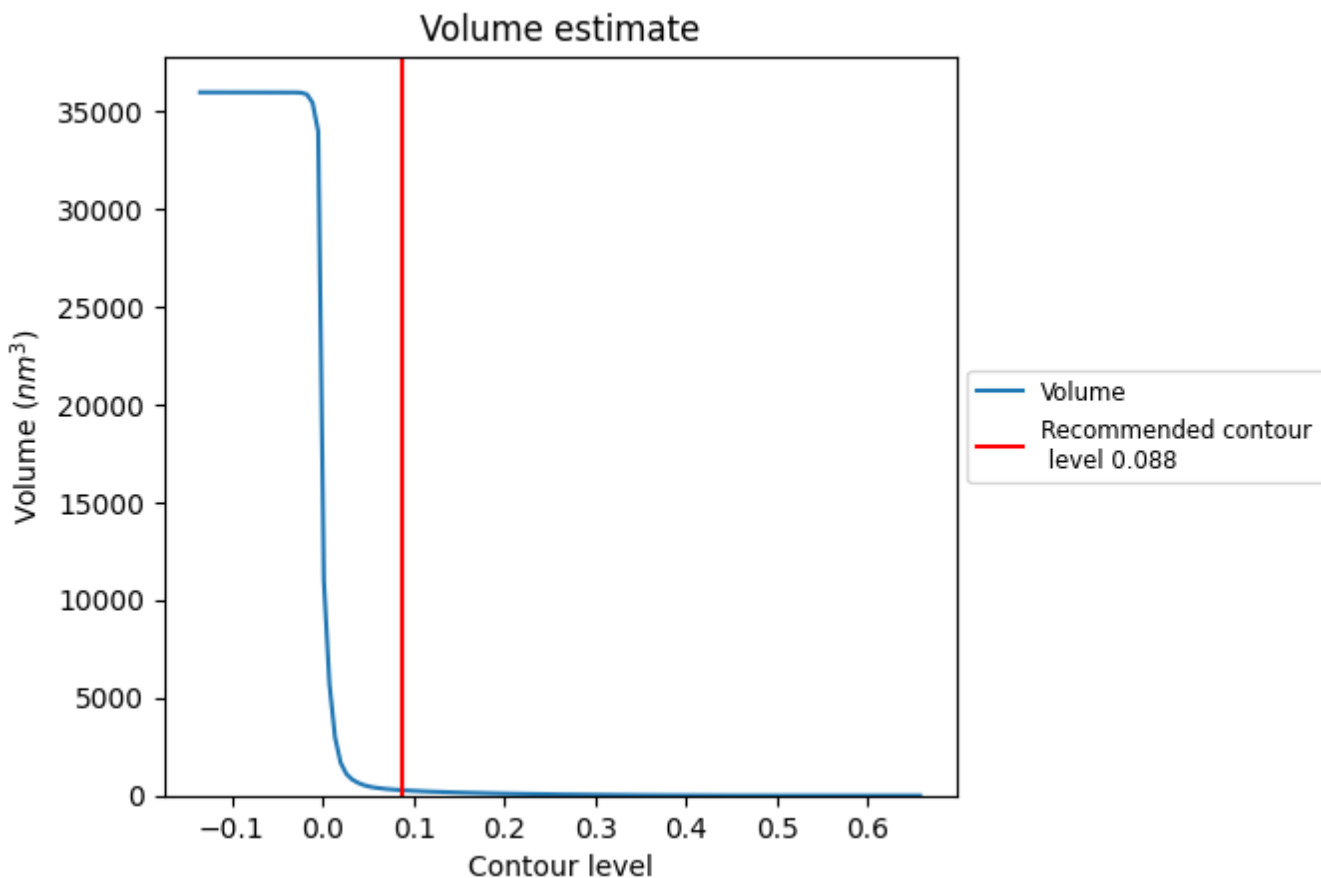
This section contains the results of statistical analysis of the map.

### 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.

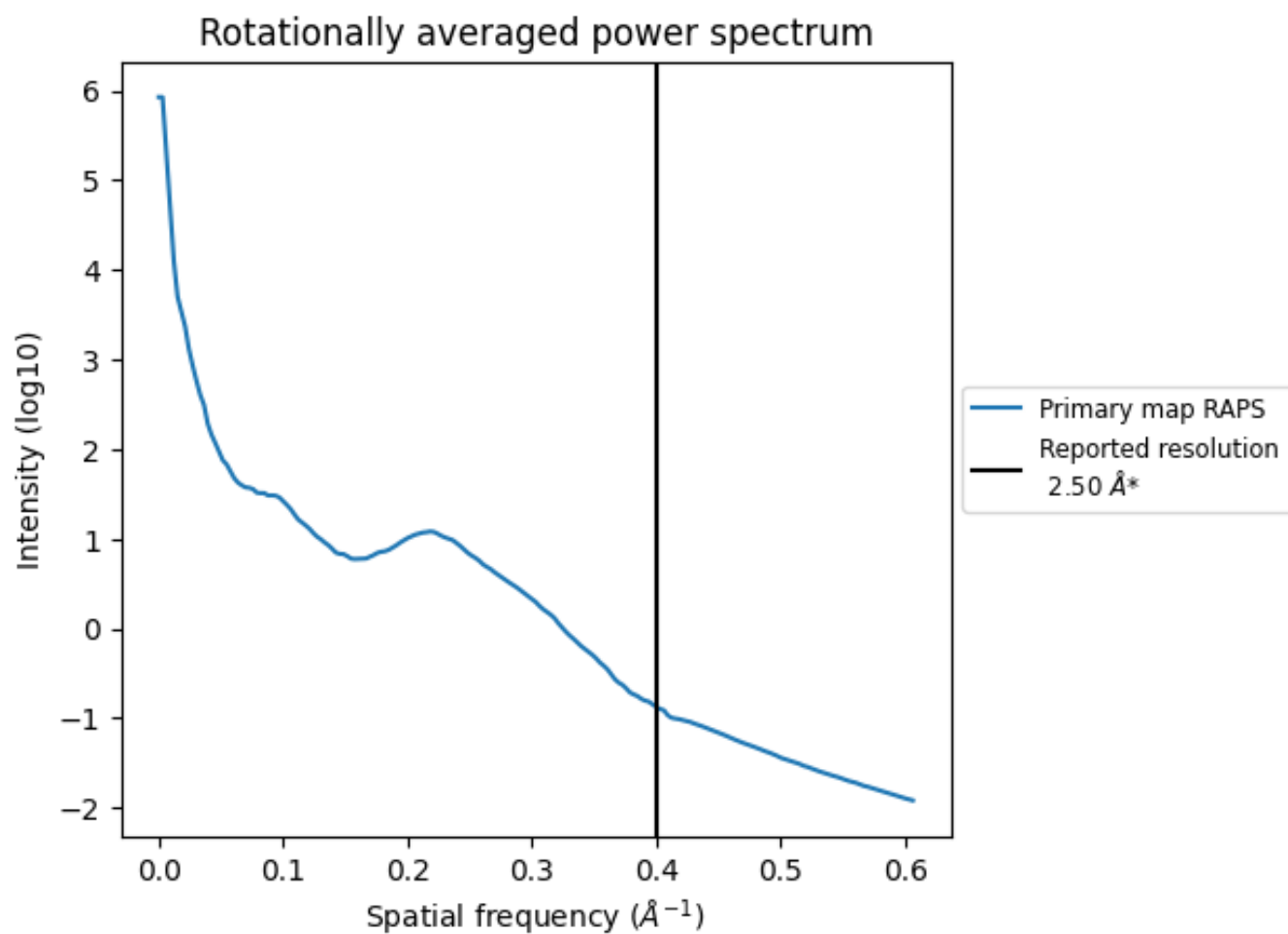
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 270 nm<sup>3</sup>; this corresponds to an approximate mass of 244 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](#)

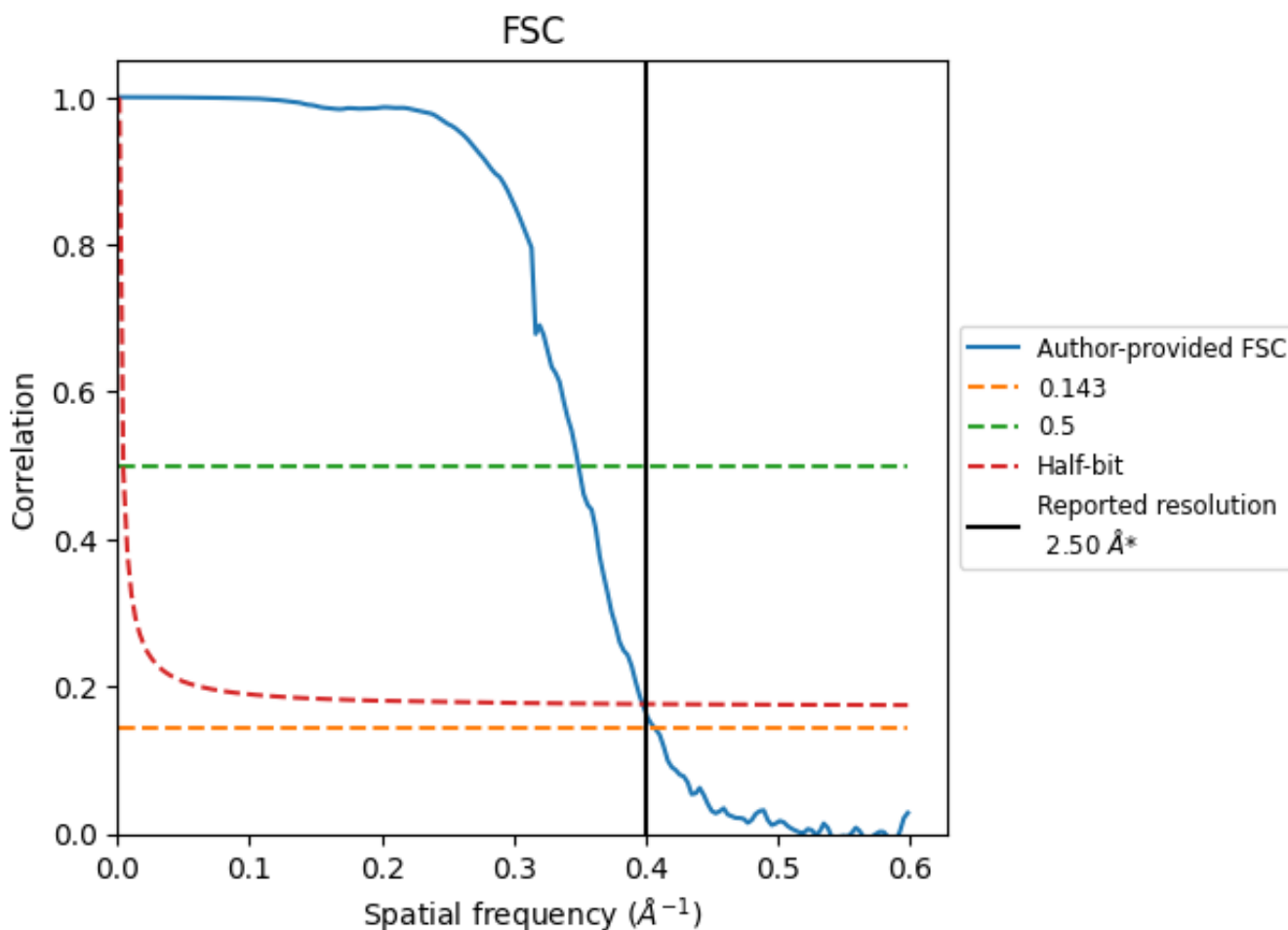


\*Reported resolution corresponds to spatial frequency of 0.400 Å<sup>-1</sup>

## 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 0.400 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

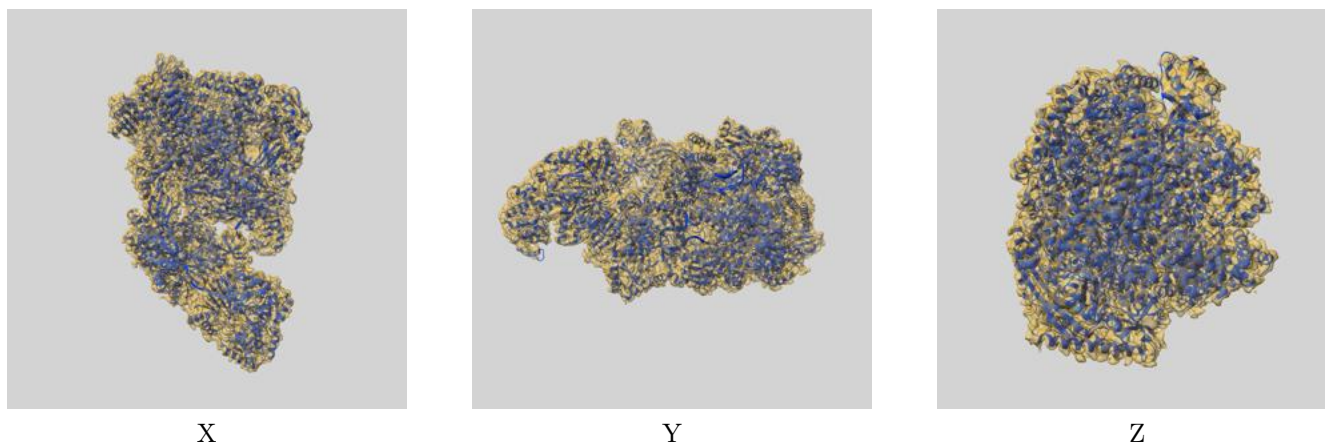
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.50	-	-
Author-provided FSC curve	2.46	2.87	2.51
Unmasked-calculated*	-	-	-

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

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

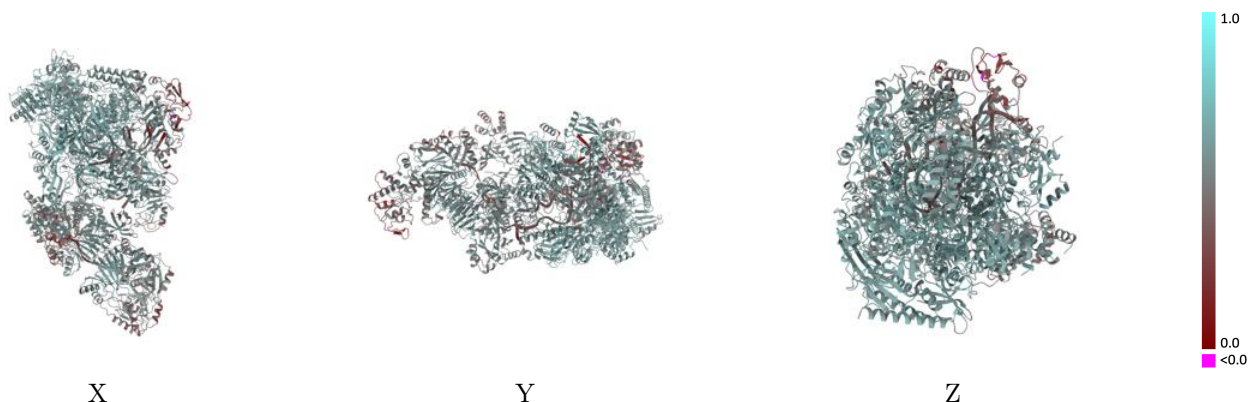
This section contains information regarding the fit between EMDB map EMD-44438 and PDB model 9BCT. Per-residue inclusion information can be found in section 3 on page 8.

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



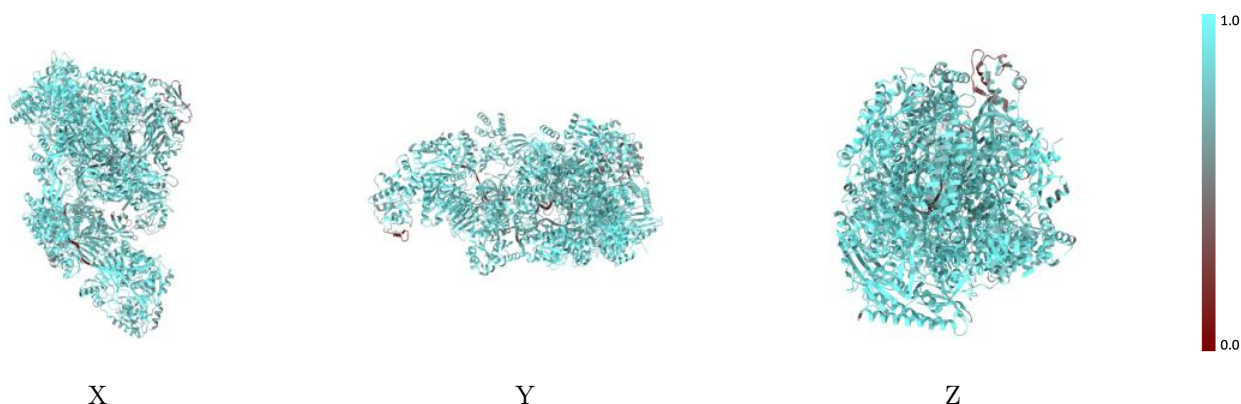
The images above show the 3D surface view of the map at the recommended contour level 0.088 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

## 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.

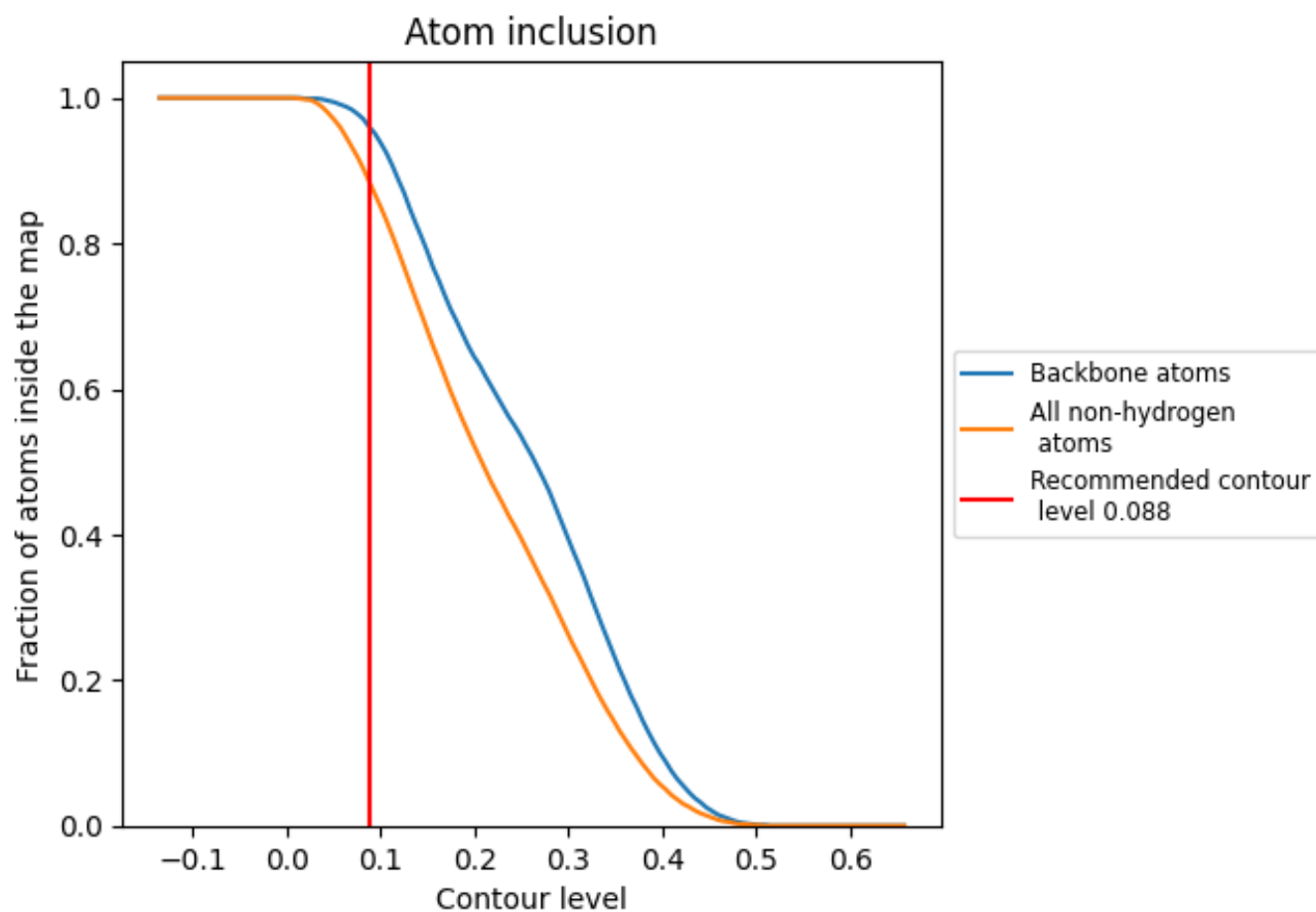
## 9.3 Atom inclusion mapped to coordinate model [i](#)



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 (0.088).









































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 88% 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 (0.088) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8850	 0.5540
5	 0.7540	 0.3700
6	 0.7860	 0.3940
7	 0.5630	 0.4440
A	 0.9210	 0.5990
B	 0.9150	 0.5920
C	 0.8340	 0.5180
D	 0.9310	 0.6070
E	 0.8580	 0.5120
F	 0.8220	 0.4320
G	 0.7920	 0.5200
H	 0.9260	 0.5960
I	 0.7500	 0.4140
J	 0.9110	 0.5610
K	 0.9360	 0.5970
L	 0.9270	 0.6020
M	 0.8590	 0.5020
N	 0.9280	 0.6110
P	 0.9220	 0.5660

