Full wwPDB EM Validation Report ⓘ

Nov 22, 2022 – 11:24 AM JST

PDB ID	:	7ECV
EMDB ID	:	EMD-31058
Title	:	The Csy-AcrIF14 complex
Authors	:	Zhang, L.X.; Feng, Y.
Deposited on	:	2021-03-13
Resolution	:	3.43 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:

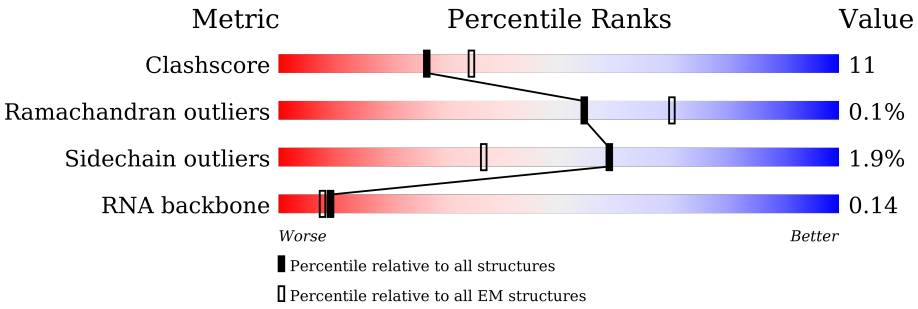
EMDB validation analysis	:	0.0.1.dev43
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.43 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 >=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 <=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	L	187	69% 78% 20%
2	A	434	15% 94%
3	B	327	73% 20% 7%
4	C	342	5% 70% 15% 14%
4	D	342	81% 15%
4	E	342	75% 20%
4	F	342	79% 19%

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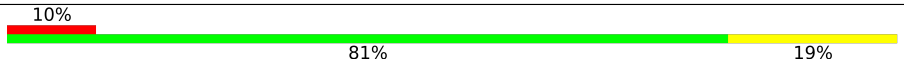
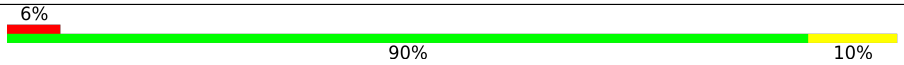
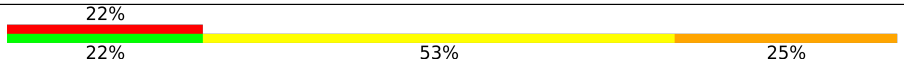
Mol	Chain	Length	Quality of chain
4	G	342	
4	H	342	
5	I	124	
5	J	124	
6	M	60	

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 24153 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 Type I-F CRISPR-associated endoribonuclease Cas6/Csy4.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	L	185	Total	C	N	O	S	1	0
			1423	906	257	255	5		

- Molecule 2 is a protein called Type I-F CRISPR-associated protein Csy1.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	A	424	Total	C	N	O	0	0
			2065	1138	466	461		

- Molecule 3 is a protein called CRISPR type I-F/YPEST-associated protein Csy2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	305	Total	C	N	O	S	0	0
			2368	1500	439	424	5		

- Molecule 4 is a protein called CRISPR-associated protein Csy3.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	C	293	Total	C	N	O	S	0	0
			2265	1425	408	430	2		
4	D	333	Total	C	N	O	S	0	0
			2548	1601	463	482	2		
4	E	334	Total	C	N	O	S	0	0
			2560	1611	466	481	2		
4	F	335	Total	C	N	O	S	0	0
			2574	1617	469	486	2		
4	G	333	Total	C	N	O	S	0	0
			2560	1610	466	482	2		
4	H	331	Total	C	N	O	S	0	0
			2547	1601	466	478	2		

- Molecule 5 is a protein called AcrIF14.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	I	124	Total 992	C 628	N 169	O 189	S 6	0	0
5	J	124	Total 979	C 620	N 166	O 188	S 5	0	0

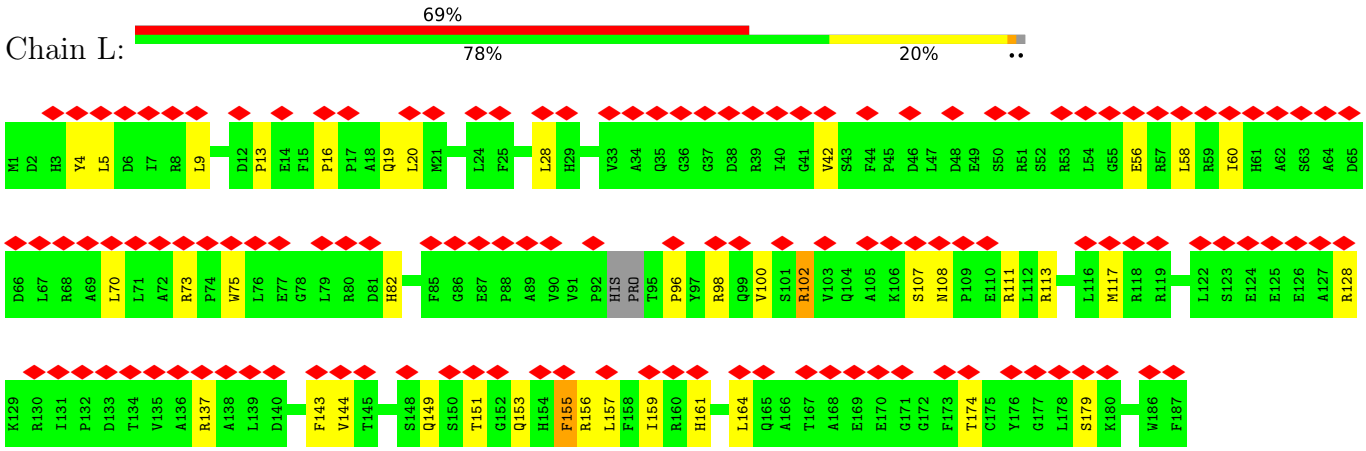
- Molecule 6 is a RNA chain called RNA (60-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	M	60	Total 1272	C 569	N 223	O 421	P 59	0	0

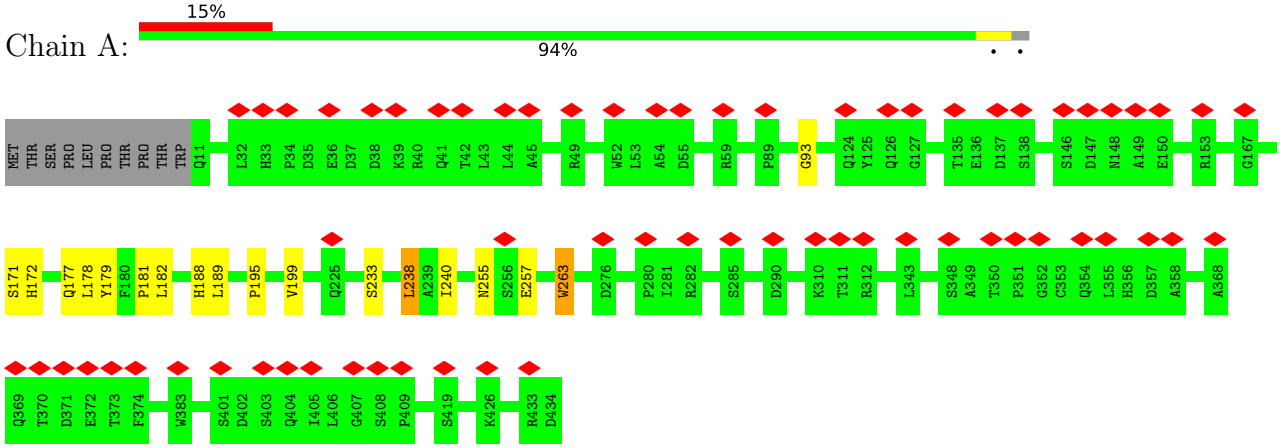
3 Residue-property plots (i)

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.

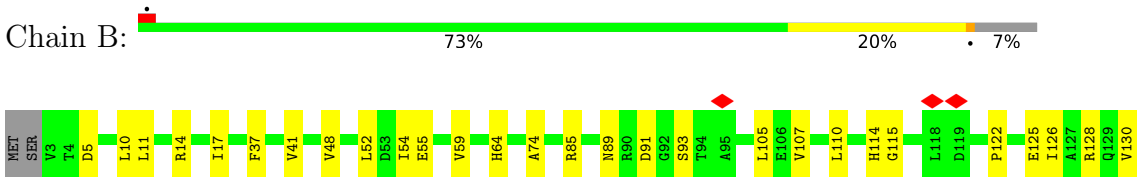
• Molecule 1: Type I-F CRISPR-associated endoribonuclease Cas6/Csy4

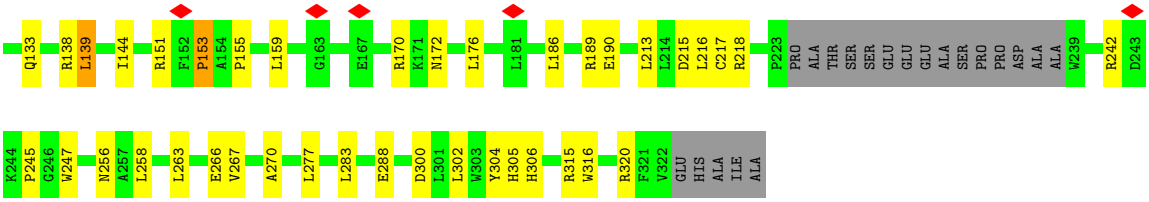


• Molecule 2: Type I-F CRISPR-associated protein Csy1

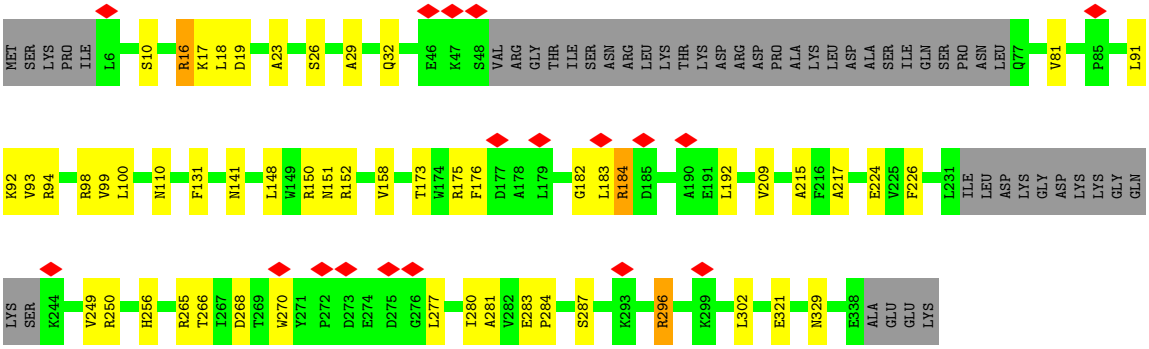


• Molecule 3: CRISPR type I-F/YPEST-associated protein Csy2

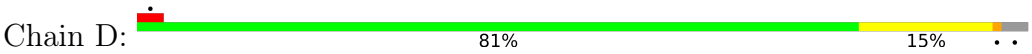


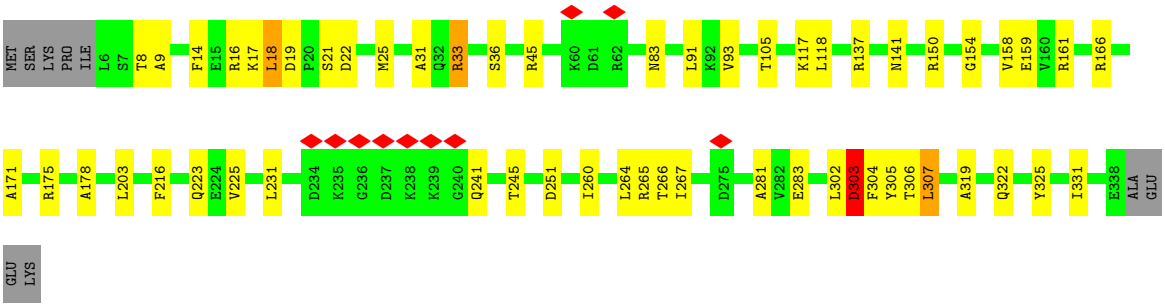


• Molecule 4: CRISPR-associated protein Csy3

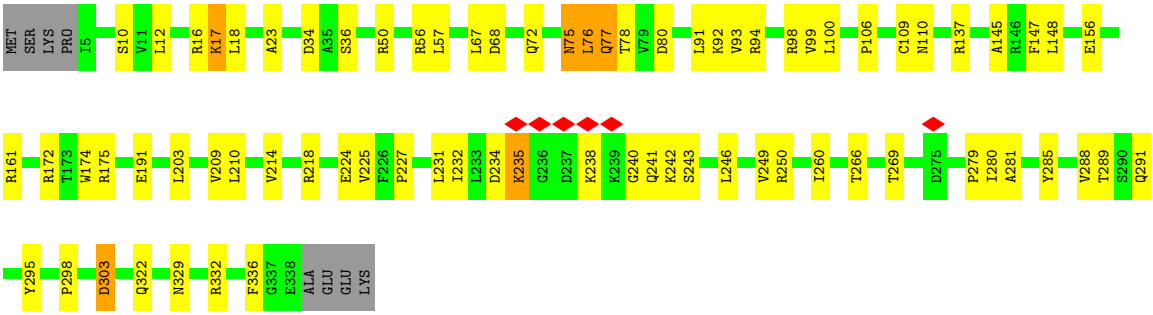


• Molecule 4: CRISPR-associated protein Csy3



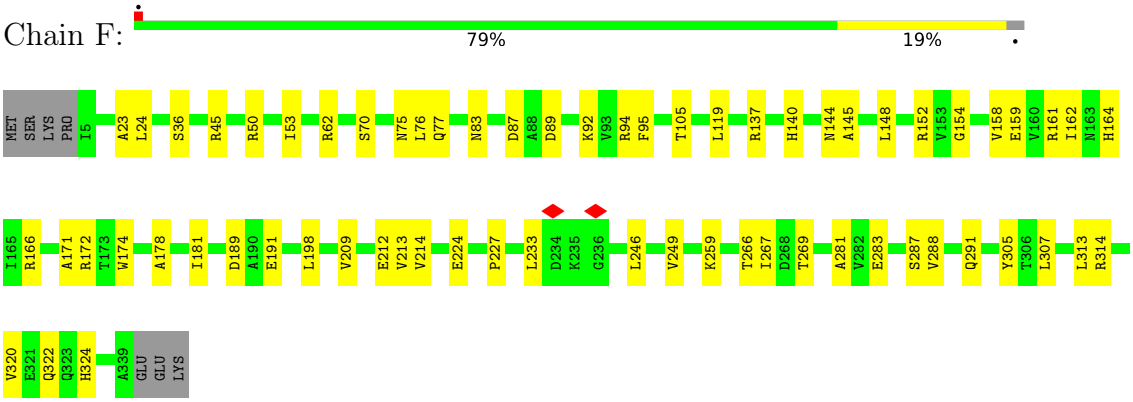


• Molecule 4: CRISPR-associated protein Csy3



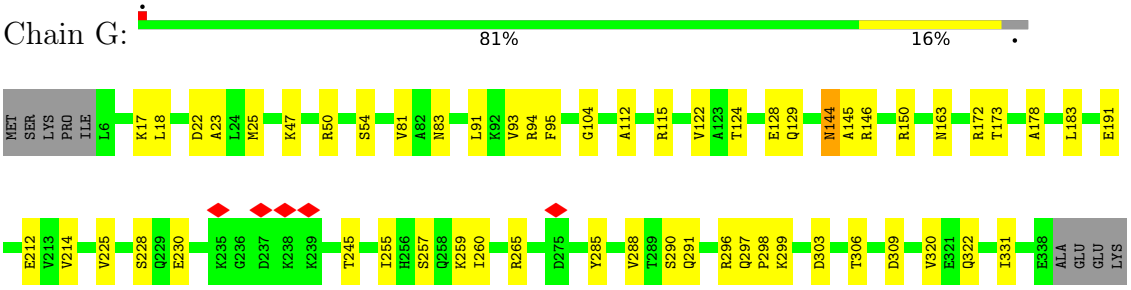
• Molecule 4: CRISPR-associated protein Csy3

Chain F:



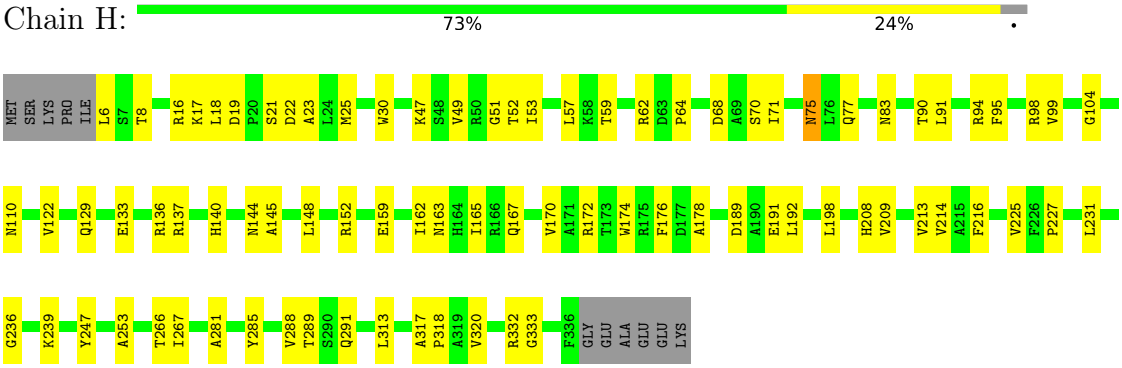
• Molecule 4: CRISPR-associated protein Csy3

Chain G:



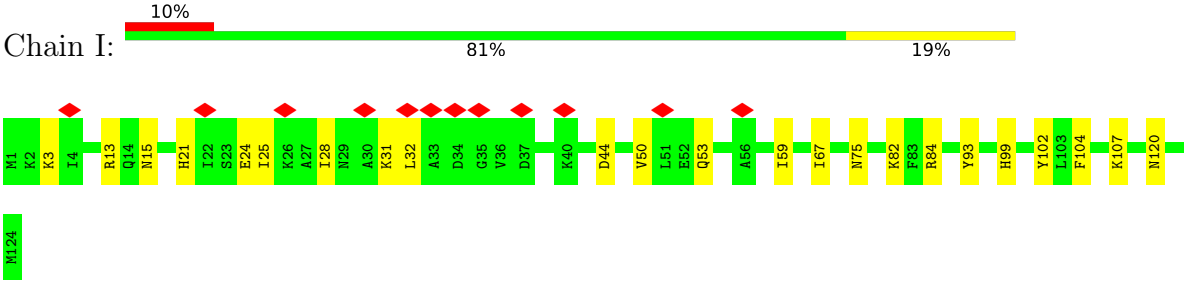
• Molecule 4: CRISPR-associated protein Csy3

Chain H:



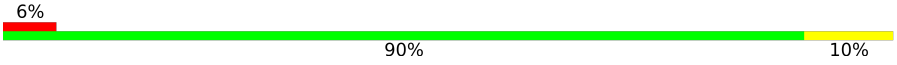
• Molecule 5: AcrIF14

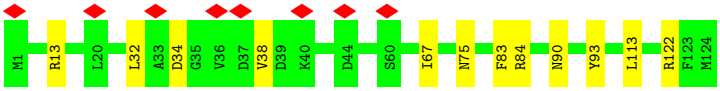
Chain I:



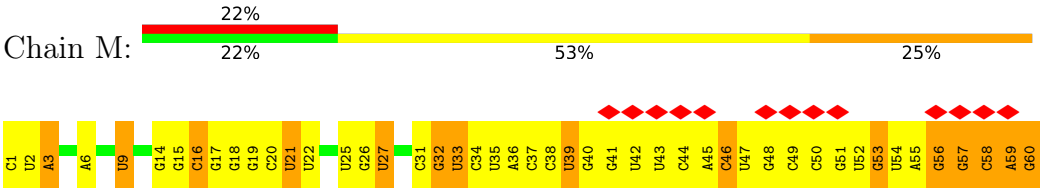
• Molecule 5: AcrIF14

Chain J:





• Molecule 6: RNA (60-MER)



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	126583	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	50	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	0.389	Depositor
Minimum map value	-0.241	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.013	Depositor
Recommended contour level	0.028	Depositor
Map size (Å)	220.0, 220.0, 220.0	wwPDB
Map dimensions	200, 200, 200	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	Depositor

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|>5$	RMSZ	#$\|Z\|>5$
1	L	0.29	0/1462	0.58	0/1984
2	A	0.26	0/2090	0.49	0/2698
3	B	0.33	0/2425	0.57	0/3302
4	C	0.28	0/2308	0.52	0/3135
4	D	0.32	0/2595	0.55	0/3525
4	E	0.35	0/2607	0.54	0/3539
4	F	0.36	0/2621	0.54	0/3558
4	G	0.37	0/2607	0.54	0/3537
4	H	0.35	0/2594	0.55	0/3521
5	I	0.31	0/1007	0.50	0/1346
5	J	0.31	0/995	0.47	0/1333
6	M	0.54	0/1420	0.91	0/2212
All	All	0.34	0/24731	0.57	0/33690

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

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	L	1423	0	1364	59	0
2	A	2065	0	1073	37	0
3	B	2368	0	2332	66	0
4	C	2265	0	2214	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	2548	0	2513	65	0
4	E	2560	0	2542	83	0
4	F	2574	0	2553	47	0
4	G	2560	0	2546	48	0
4	H	2547	0	2533	67	0
5	I	992	0	988	17	0
5	J	979	0	946	6	0
6	M	1272	0	644	69	0
All	All	24153	0	22248	513	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 11.

All (513) 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:A:179:TYR:HE1	2:A:189:LEU:CD1	1.03	1.62
2:A:179:TYR:CE1	2:A:189:LEU:CD1	1.82	1.56
1:L:102:ARG:NH2	1:L:155:PHE:HZ	1.05	1.47
1:L:102:ARG:NH2	1:L:155:PHE:CZ	1.82	1.45
2:A:179:TYR:CD1	2:A:189:LEU:HD12	1.64	1.31
4:E:76:LEU:O	4:E:240:GLY:HA3	1.23	1.30
4:D:303:ASP:O	4:D:307:LEU:HD21	1.10	1.28
4:D:303:ASP:O	4:D:307:LEU:CD2	1.83	1.23
4:H:159:GLU:HG3	4:H:176:PHE:O	1.40	1.22
1:L:100:VAL:HG13	1:L:159:ILE:HD11	1.24	1.14
2:A:179:TYR:CE1	2:A:189:LEU:HD12	1.59	1.13
4:E:227:PRO:HG2	4:E:246:LEU:HD13	1.11	1.09
2:A:179:TYR:HD2	6:M:1:C:C2	1.73	1.07
4:H:94:ARG:HG3	4:H:214:VAL:HG22	1.32	1.06
2:A:179:TYR:HE1	2:A:189:LEU:HD13	1.16	1.06
3:B:17:ILE:HG21	3:B:139:LEU:HD23	1.10	1.05
2:A:179:TYR:CE1	2:A:189:LEU:HD11	1.68	1.02
4:D:303:ASP:OD1	4:D:305:TYR:N	1.91	1.02
4:E:76:LEU:O	4:E:240:GLY:CA	2.08	1.01
2:A:179:TYR:HE1	2:A:189:LEU:HD11	1.01	1.01
1:L:155:PHE:CZ	1:L:156:ARG:O	2.13	1.01
1:L:100:VAL:CG1	1:L:159:ILE:HD11	1.91	1.00
4:G:17:LYS:O	4:G:18:LEU:HD12	1.61	1.00
2:A:178:LEU:HD22	6:M:2:U:H5"	1.45	0.99
4:E:227:PRO:HG2	4:E:246:LEU:CD1	1.93	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:76:LEU:HD11	6:M:33:U:O2	1.63	0.98
4:D:18:LEU:HD21	4:D:331:ILE:CG2	1.95	0.97
1:L:13:PRO:HG3	4:C:277:LEU:CD1	1.95	0.97
4:D:158:VAL:HG13	4:D:178:ALA:HB3	1.44	0.96
2:A:179:TYR:CD2	6:M:1:C:N3	2.33	0.95
4:D:307:LEU:H	4:D:307:LEU:HD23	1.26	0.94
1:L:42:VAL:HG21	1:L:58:LEU:HD11	1.49	0.94
4:G:297:GLN:NE2	4:G:298:PRO:HD2	1.82	0.93
1:L:155:PHE:CD2	6:M:46:C:C4	2.56	0.92
4:E:227:PRO:CG	4:E:246:LEU:HD13	1.98	0.92
2:A:179:TYR:CD2	6:M:1:C:C2	2.58	0.91
4:D:303:ASP:OD1	4:D:304:PHE:N	2.03	0.90
3:B:17:ILE:HG21	3:B:139:LEU:CD2	2.02	0.90
1:L:102:ARG:NH2	1:L:155:PHE:CE2	2.39	0.90
4:F:162:ILE:HG13	4:F:213:VAL:HG23	1.53	0.90
4:D:18:LEU:HD21	4:D:331:ILE:HG21	1.54	0.89
2:A:179:TYR:HD2	6:M:1:C:O2	1.56	0.88
1:L:155:PHE:CD2	6:M:46:C:N3	2.42	0.88
2:A:181:PRO:HG2	3:B:305:HIS:NE2	1.90	0.87
2:A:179:TYR:HD1	2:A:189:LEU:HD12	1.36	0.87
2:A:179:TYR:CE2	6:M:1:C:N3	2.42	0.86
1:L:155:PHE:CE2	1:L:156:ARG:O	2.28	0.86
3:B:17:ILE:HD11	3:B:105:LEU:HG	1.59	0.85
1:L:42:VAL:CG2	1:L:58:LEU:HD11	2.08	0.84
4:E:80:ASP:OD1	4:E:242:LYS:CE	2.25	0.84
6:M:56:G:O2'	6:M:57:G:C8	2.30	0.83
4:G:18:LEU:HD13	4:G:331:ILE:CG2	2.07	0.83
6:M:58:C:C1'	6:M:59:A:OP2	2.27	0.82
4:E:77:GLN:CG	4:E:243:SER:HB3	2.09	0.81
4:E:77:GLN:HG3	4:E:243:SER:HB3	1.61	0.80
3:B:110:LEU:CD1	3:B:302:LEU:HD21	2.13	0.78
3:B:110:LEU:HD12	3:B:110:LEU:O	1.84	0.78
4:C:23:ALA:HB1	4:C:93:VAL:CG1	2.13	0.78
2:A:178:LEU:HD11	6:M:3:A:H5"	1.66	0.78
4:D:307:LEU:CD2	4:D:307:LEU:H	1.97	0.77
4:D:303:ASP:OD2	4:D:306:THR:CB	2.33	0.77
4:D:303:ASP:C	4:D:307:LEU:HD21	2.02	0.76
4:E:249:VAL:HG12	4:E:250:ARG:H	1.48	0.76
1:L:155:PHE:CE2	6:M:46:C:N4	2.54	0.76
2:A:179:TYR:CE1	2:A:189:LEU:HD13	2.00	0.75
4:D:231:LEU:HD22	4:E:76:LEU:HB3	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:139:LEU:HD12	3:B:139:LEU:O	1.87	0.75
6:M:58:C:H1'	6:M:59:A:OP2	1.86	0.74
4:E:80:ASP:OD1	4:E:242:LYS:HE2	1.88	0.74
4:E:76:LEU:HD21	6:M:33:U:O2	1.87	0.74
4:E:77:GLN:HG2	4:E:243:SER:CB	2.17	0.74
4:G:17:LYS:C	4:G:18:LEU:HD12	2.06	0.74
6:M:58:C:O4'	6:M:59:A:OP2	2.05	0.74
3:B:17:ILE:CG2	3:B:139:LEU:HD23	2.05	0.73
4:E:99:VAL:CG2	4:E:203:LEU:HD23	2.17	0.73
1:L:155:PHE:HD2	6:M:46:C:N3	1.85	0.72
4:E:75:ASN:OD1	4:E:75:ASN:N	2.22	0.72
1:L:149:GLN:CB	6:M:60:G:H2'	2.19	0.72
3:B:110:LEU:HD12	3:B:302:LEU:HD21	1.70	0.72
4:D:17:LYS:O	4:D:18:LEU:HG	1.89	0.71
1:L:13:PRO:HG3	4:C:277:LEU:HD13	1.73	0.71
4:E:99:VAL:HG12	4:E:209:VAL:CG2	2.21	0.71
6:M:56:G:H1'	6:M:57:G:C6	2.26	0.71
4:H:163:ASN:ND2	4:H:170:VAL:HG22	2.06	0.71
4:E:78:THR:HB	4:E:241:GLN:O	1.92	0.70
4:G:50:ARG:NH1	6:M:21:U:O2'	2.24	0.70
1:L:155:PHE:HD2	6:M:46:C:C4	2.06	0.70
4:E:12:LEU:HB3	4:E:336:PHE:O	1.92	0.70
6:M:58:C:H1'	6:M:59:A:P	2.32	0.69
1:L:42:VAL:CG2	1:L:58:LEU:CD1	2.71	0.69
4:G:47:LYS:HE2	4:G:81:VAL:HG21	1.75	0.69
4:D:303:ASP:OD2	4:D:306:THR:OG1	2.09	0.69
4:D:8:THR:HG22	4:D:9:ALA:N	2.07	0.69
4:D:307:LEU:HD23	4:D:307:LEU:N	2.06	0.68
2:A:178:LEU:HD11	6:M:3:A:C5'	2.24	0.68
4:D:303:ASP:O	4:D:307:LEU:HD22	1.93	0.67
4:F:181:ILE:HD11	4:F:189:ASP:OD2	1.95	0.67
3:B:213:LEU:HD23	3:B:216:LEU:HD21	1.77	0.66
4:H:57:LEU:HD11	4:H:62:ARG:HB3	1.77	0.66
4:E:99:VAL:CG2	4:E:203:LEU:CD2	2.73	0.66
1:L:155:PHE:CE2	6:M:46:C:C4	2.84	0.66
4:F:320:VAL:HG12	4:F:324:HIS:CE1	2.29	0.66
3:B:17:ILE:HD11	3:B:105:LEU:CG	2.25	0.66
4:E:76:LEU:HD21	6:M:33:U:C2	2.30	0.66
3:B:10:LEU:HD21	3:B:130:VAL:HG11	1.76	0.66
2:A:195:PRO:O	2:A:199:VAL:HG23	1.95	0.66
3:B:216:LEU:HD11	3:B:218:ARG:HB2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:260:ILE:O	4:D:264:LEU:HG	1.97	0.65
3:B:242:ARG:HH12	3:B:245:PRO:HD3	1.61	0.65
4:H:18:LEU:O	4:H:18:LEU:HD12	1.95	0.64
6:M:56:G:O2'	6:M:57:G:N7	2.24	0.64
3:B:17:ILE:HG22	3:B:144:ILE:HA	1.79	0.64
3:B:267:VAL:HG11	3:B:277:LEU:HD13	1.79	0.64
3:B:216:LEU:HD12	3:B:218:ARG:N	2.12	0.64
4:D:302:LEU:O	4:D:303:ASP:O	2.16	0.64
4:C:148:LEU:HD22	4:C:151:ASN:HB2	1.80	0.64
4:C:284:PRO:HG3	4:C:329:ASN:HD21	1.62	0.63
5:I:53:GLN:HE21	5:I:59:ILE:HD13	1.63	0.63
4:F:288:VAL:HG23	4:F:291:GLN:HB3	1.80	0.63
6:M:58:C:C1'	6:M:59:A:P	2.87	0.63
2:A:179:TYR:CD2	6:M:1:C:O2	2.45	0.63
1:L:73:ARG:HB3	1:L:75:TRP:HD1	1.63	0.63
1:L:108:ASN:HD22	6:M:46:C:H5'	1.64	0.63
6:M:57:G:O2'	6:M:58:C:OP1	2.14	0.63
4:H:68:ASP:HA	4:H:71:ILE:HG22	1.81	0.62
2:A:178:LEU:CD1	6:M:3:A:H5"	2.29	0.62
4:D:231:LEU:HD13	4:E:76:LEU:HD23	1.82	0.62
4:E:76:LEU:HD21	6:M:33:U:H3	1.64	0.62
4:G:23:ALA:HB2	4:G:260:ILE:HD13	1.81	0.62
6:M:53:G:H8	6:M:55:A:H61	1.46	0.62
4:G:18:LEU:HD13	4:G:331:ILE:HG22	1.81	0.61
2:A:238:LEU:HD12	2:A:238:LEU:O	2.00	0.61
3:B:216:LEU:CD1	3:B:218:ARG:N	2.64	0.61
4:E:288:VAL:CG2	4:E:291:GLN:HB2	2.30	0.61
4:H:17:LYS:O	4:H:18:LEU:HG	2.01	0.61
4:D:141:ASN:ND2	4:D:266:THR:OG1	2.34	0.61
4:E:76:LEU:CD1	6:M:33:U:O2	2.45	0.61
3:B:54:ILE:HD13	3:B:126:ILE:HD11	1.82	0.61
4:C:93:VAL:HG21	4:C:148:LEU:HD21	1.83	0.60
4:D:303:ASP:OD2	4:D:306:THR:HB	2.00	0.60
4:C:266:THR:HA	4:C:281:ALA:HA	1.83	0.60
4:H:163:ASN:HD22	4:H:170:VAL:HG22	1.64	0.60
3:B:89:ASN:HD22	3:B:93:SER:HB2	1.66	0.60
4:G:18:LEU:CD1	4:G:331:ILE:CG2	2.79	0.60
2:A:179:TYR:CD1	2:A:189:LEU:CD1	2.43	0.60
4:C:26:SER:O	4:C:91:LEU:HD12	2.01	0.60
4:C:283:GLU:HG2	4:C:287:SER:H	1.66	0.60
1:L:151:THR:OG1	1:L:153:GLN:NE2	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:99:VAL:HG21	4:E:203:LEU:CD2	2.31	0.59
4:E:17:LYS:HG3	4:E:100:LEU:HB2	1.83	0.59
3:B:17:ILE:CD1	3:B:105:LEU:HG	2.30	0.59
4:H:75:ASN:ND2	4:H:75:ASN:H	2.01	0.59
4:E:249:VAL:HG12	4:E:250:ARG:N	2.15	0.59
4:G:297:GLN:HE21	4:G:298:PRO:HD2	1.64	0.58
4:D:19:ASP:OD1	4:D:19:ASP:O	2.22	0.58
4:E:227:PRO:CG	4:E:246:LEU:CD1	2.71	0.58
4:G:288:VAL:HG23	4:G:291:GLN:HB2	1.86	0.58
4:F:288:VAL:CG2	4:F:291:GLN:HB3	2.34	0.58
4:E:77:GLN:CG	4:E:243:SER:CB	2.76	0.58
4:G:265:ARG:NH2	6:M:14:G:OP2	2.36	0.58
6:M:58:C:C4'	6:M:59:A:OP2	2.51	0.58
4:C:150:ARG:HD3	4:C:226:PHE:HD2	1.69	0.58
4:D:267:ILE:O	4:D:267:ILE:HG22	2.04	0.58
4:E:76:LEU:HD21	6:M:33:U:N3	2.19	0.57
4:E:298:PRO:HB2	4:F:62:ARG:HD3	1.86	0.57
1:L:100:VAL:CG1	1:L:159:ILE:CD1	2.75	0.57
5:I:50:VAL:HA	5:I:53:GLN:HG2	1.87	0.57
4:H:104:GLY:HA3	4:H:122:VAL:HG11	1.87	0.57
4:E:94:ARG:HG3	4:E:214:VAL:HG22	1.87	0.57
3:B:41:VAL:HG21	3:B:59:VAL:CG1	2.35	0.57
4:E:77:GLN:HG2	4:E:243:SER:OG	2.04	0.57
4:F:162:ILE:CG1	4:F:213:VAL:HG23	2.32	0.57
4:H:53:ILE:HG13	4:H:53:ILE:O	2.05	0.57
4:C:16:ARG:HB3	6:M:36:A:H5"	1.86	0.56
4:E:99:VAL:HG12	4:E:209:VAL:HG23	1.86	0.56
4:G:146:ARG:CZ	4:G:183:LEU:HD12	2.35	0.56
4:F:95:PHE:CZ	4:F:213:VAL:HG11	2.40	0.56
4:E:288:VAL:HG23	4:E:291:GLN:HB2	1.87	0.56
4:D:45:ARG:NH2	4:D:83:ASN:OD1	2.38	0.56
4:D:302:LEU:C	4:D:303:ASP:O	2.43	0.56
4:C:173:THR:OG1	4:C:175:ARG:NH1	2.39	0.56
4:G:144:ASN:HD21	4:G:183:LEU:HD12	1.71	0.56
2:A:255:ASN:ND2	2:A:257:GLU:OE2	2.38	0.56
4:C:23:ALA:HB1	4:C:93:VAL:HG12	1.84	0.56
4:D:8:THR:CG2	4:D:9:ALA:N	2.69	0.56
4:G:230:GLU:HB2	4:G:245:THR:HG23	1.87	0.56
4:H:163:ASN:HD22	4:H:170:VAL:CG2	2.19	0.56
3:B:256:ASN:HD22	3:B:283:LEU:HD21	1.71	0.56
4:E:161:ARG:HG2	4:E:175:ARG:HG2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:152:ARG:NH2	4:H:178:ALA:O	2.39	0.55
5:J:32:LEU:HD21	5:J:67:ILE:HG21	1.87	0.55
2:A:172:HIS:O	2:A:177:GLN:NE2	2.40	0.55
3:B:55:GLU:HB2	3:B:114:HIS:HB2	1.87	0.55
1:L:13:PRO:HG3	4:C:277:LEU:HD11	1.84	0.55
1:L:144:VAL:O	1:L:156:ARG:HA	2.06	0.55
4:E:18:LEU:HD23	4:E:99:VAL:HB	1.87	0.55
4:C:296:ARG:NH1	4:C:302:LEU:O	2.38	0.55
6:M:57:G:C2'	6:M:58:C:OP1	2.54	0.55
4:E:68:ASP:O	4:E:72:GLN:NE2	2.40	0.55
4:E:106:PRO:HG2	4:E:109:CYS:SG	2.47	0.55
4:E:16:ARG:NH2	4:F:224:GLU:OE2	2.38	0.54
6:M:56:G:O2'	6:M:57:G:C5	2.52	0.54
1:L:155:PHE:CE1	1:L:156:ARG:C	2.80	0.54
3:B:14:ARG:HG3	3:B:64:HIS:HD2	1.72	0.54
4:C:23:ALA:HB1	4:C:93:VAL:HG13	1.87	0.54
4:F:320:VAL:O	4:F:324:HIS:ND1	2.40	0.54
4:H:51:GLY:O	4:H:75:ASN:ND2	2.40	0.54
4:C:10:SER:OG	4:C:110:ASN:ND2	2.41	0.54
4:D:166:ARG:HB2	4:D:171:ALA:HB2	1.88	0.54
4:F:145:ALA:H	4:F:178:ALA:HB1	1.73	0.54
4:G:94:ARG:HG3	4:G:214:VAL:HG22	1.90	0.54
5:I:15:ASN:ND2	5:I:44:ASP:OD2	2.40	0.54
4:D:241:GLN:HB3	4:D:245:THR:HG22	1.90	0.54
4:H:70:SER:O	5:J:84:ARG:NH1	2.40	0.54
3:B:91:ASP:OD2	3:B:93:SER:OG	2.26	0.54
3:B:144:ILE:O	4:H:98:ARG:NH1	2.41	0.54
3:B:270:ALA:O	4:H:110:ASN:ND2	2.41	0.54
2:A:179:TYR:HE2	6:M:1:C:C4	2.26	0.54
4:D:21:SER:OG	4:D:22:ASP:N	2.40	0.54
4:F:95:PHE:CZ	4:F:213:VAL:CG1	2.91	0.54
4:H:59:THR:HA	4:H:62:ARG:HE	1.72	0.54
3:B:105:LEU:HD13	3:B:107:VAL:HB	1.90	0.53
2:A:178:LEU:HD22	6:M:2:U:C5'	2.27	0.53
4:D:18:LEU:CD2	4:D:203:LEU:HD21	2.38	0.53
1:L:100:VAL:H	1:L:179:SER:HB3	1.73	0.53
4:H:129:GLN:HE22	4:H:320:VAL:HG23	1.73	0.53
5:I:13:ARG:HD3	5:I:75:ASN:HB3	1.89	0.53
4:C:265:ARG:NH2	6:M:38:C:OP2	2.42	0.53
4:D:303:ASP:CG	4:D:306:THR:H	2.11	0.53
4:F:140:HIS:O	4:F:144:ASN:ND2	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:181:PRO:CG	3:B:305:HIS:NE2	2.69	0.52
4:E:289:THR:OG1	4:F:50:ARG:NH2	2.40	0.52
1:L:20:LEU:HD21	1:L:82:HIS:HB3	1.90	0.52
4:C:249:VAL:HG12	4:C:250:ARG:N	2.25	0.52
4:F:87:ASP:OD1	4:F:87:ASP:N	2.42	0.52
4:D:137:ARG:HB3	4:D:267:ILE:HD12	1.92	0.52
4:E:99:VAL:HG21	4:E:203:LEU:HG	1.92	0.52
1:L:70:LEU:O	1:L:75:TRP:NE1	2.36	0.52
4:G:128:GLU:HG2	4:G:129:GLN:HG2	1.91	0.52
4:G:146:ARG:NH1	4:G:183:LEU:CD1	2.72	0.52
1:L:155:PHE:CZ	1:L:157:LEU:HA	2.45	0.52
4:E:235:LYS:CA	4:E:235:LYS:HE2	2.39	0.52
4:D:158:VAL:CG1	4:D:178:ALA:HB3	2.29	0.51
4:D:303:ASP:CG	4:D:304:PHE:N	2.60	0.51
4:F:159:GLU:OE2	4:F:161:ARG:NH2	2.41	0.51
1:L:111:ARG:HE	6:M:47:U:H5	1.57	0.51
3:B:74:ALA:HB3	4:H:236:GLY:HA3	1.91	0.51
3:B:153:PRO:HB2	3:B:155:PRO:HD3	1.92	0.51
4:G:81:VAL:HG12	4:G:83:ASN:OD1	2.10	0.51
3:B:189:ARG:HG3	3:B:288:GLU:HG3	1.92	0.51
4:G:23:ALA:HA	4:G:95:PHE:HB3	1.93	0.51
2:A:179:TYR:CE2	6:M:1:C:C4	2.99	0.51
3:B:11:LEU:HD12	3:B:110:LEU:HB3	1.92	0.51
4:C:17:LYS:HG3	4:C:100:LEU:HB2	1.92	0.51
3:B:85:ARG:HD3	4:H:289:THR:HG21	1.92	0.51
4:E:10:SER:OG	4:E:110:ASN:OD1	2.29	0.51
4:F:174:TRP:HD1	4:F:191:GLU:HB3	1.76	0.51
4:H:49:VAL:HG23	4:H:77:GLN:HG3	1.92	0.51
4:H:137:ARG:NH1	4:H:267:ILE:O	2.44	0.51
4:C:81:VAL:HG23	4:C:224:GLU:HB3	1.93	0.51
4:C:19:ASP:HB3	4:D:223:GLN:HE22	1.76	0.51
4:D:159:GLU:OE1	4:D:175:ARG:NH1	2.44	0.51
1:L:113:ARG:HE	1:L:128:ARG:HA	1.76	0.50
3:B:305:HIS:HA	3:B:316:TRP:HA	1.93	0.50
4:H:172:ARG:NH2	4:H:191:GLU:OE2	2.44	0.50
3:B:128:ARG:HH11	4:H:167:GLN:HE22	1.57	0.50
4:D:8:THR:HG22	4:D:9:ALA:H	1.76	0.50
4:G:25:MET:HB2	4:G:91:LEU:HD11	1.92	0.50
4:G:144:ASN:HD21	4:G:183:LEU:CD1	2.24	0.50
4:D:303:ASP:OD1	4:D:304:PHE:C	2.50	0.50
4:E:99:VAL:CG1	4:E:209:VAL:CG2	2.90	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:45:ARG:NH2	4:F:83:ASN:OD1	2.44	0.50
4:F:137:ARG:NH1	4:F:267:ILE:O	2.44	0.50
2:A:171:SER:OG	2:A:177:GLN:NE2	2.43	0.50
4:C:256:HIS:ND1	6:M:39:U:OP1	2.44	0.50
4:E:280:ILE:HG23	4:E:288:VAL:CG1	2.42	0.50
4:G:91:LEU:HD13	4:G:225:VAL:HG11	1.94	0.50
4:C:99:VAL:HG22	4:C:209:VAL:HG22	1.93	0.50
4:F:305:TYR:OH	4:G:54:SER:O	2.24	0.50
4:H:227:PRO:HG3	4:H:253:ALA:HB1	1.94	0.50
2:A:179:TYR:HE2	6:M:1:C:N3	2.06	0.50
4:D:33:ARG:HA	4:D:36:SER:HB3	1.93	0.50
4:D:304:PHE:HA	4:D:307:LEU:HD21	1.94	0.50
4:C:92:LYS:HA	4:C:215:ALA:O	2.12	0.50
4:G:146:ARG:HG3	4:G:183:LEU:HD11	1.92	0.50
4:H:133:GLU:OE1	4:H:136:ARG:NH2	2.45	0.49
1:L:155:PHE:CE1	1:L:156:ARG:O	2.63	0.49
2:A:240:ILE:HD13	2:A:263:TRP:HB3	1.93	0.49
5:I:21:HIS:HB3	5:I:24:GLU:HG3	1.95	0.49
4:H:6:LEU:HD11	4:H:313:LEU:CD2	2.43	0.49
4:H:52:THR:HB	4:H:71:ILE:HD11	1.95	0.49
4:D:25:MET:HG2	4:D:93:VAL:HG22	1.94	0.49
4:C:16:ARG:NH2	6:M:37:C:OP1	2.45	0.49
4:E:147:PHE:HB2	4:E:260:ILE:HG13	1.93	0.49
4:H:57:LEU:CD1	4:H:62:ARG:HB3	2.42	0.49
4:C:158:VAL:HG13	4:C:217:ALA:HB2	1.95	0.49
4:D:158:VAL:HA	4:D:216:PHE:O	2.12	0.49
4:E:174:TRP:HD1	4:E:191:GLU:HG2	1.77	0.49
4:E:280:ILE:HG23	4:E:288:VAL:HG13	1.95	0.49
1:L:155:PHE:HB2	6:M:46:C:O2	2.13	0.49
2:A:233:SER:HB2	3:B:216:LEU:HD13	1.95	0.49
4:D:251:ASP:OD1	4:D:251:ASP:N	2.46	0.49
4:E:23:ALA:HB1	4:E:93:VAL:HG13	1.95	0.49
4:H:8:THR:HG22	4:H:313:LEU:HD11	1.94	0.49
4:E:76:LEU:CD2	6:M:33:U:H3	2.25	0.49
4:C:270:TRP:HB3	4:C:321:GLU:HG2	1.95	0.48
4:C:98:ARG:NH1	4:D:154:GLY:O	2.46	0.48
1:L:13:PRO:HB3	4:C:280:ILE:HG22	1.95	0.48
4:D:137:ARG:CB	4:D:267:ILE:HD12	2.43	0.48
4:F:95:PHE:CE1	4:F:213:VAL:HG13	2.48	0.48
3:B:110:LEU:HD11	3:B:302:LEU:HD21	1.93	0.48
3:B:304:TYR:OH	3:B:320:ARG:NH2	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:105:THR:O	4:D:105:THR:HG23	2.14	0.48
4:H:25:MET:HB3	4:H:91:LEU:HD11	1.96	0.48
4:H:30:TRP:HE3	4:H:90:THR:HG21	1.78	0.48
6:M:1:C:O2	6:M:1:C:H3'	2.13	0.48
4:C:18:LEU:CD2	4:C:99:VAL:HG12	2.43	0.48
4:G:129:GLN:HE22	4:G:320:VAL:HG13	1.78	0.48
4:G:299:LYS:HG3	4:H:64:PRO:HG3	1.95	0.48
4:D:303:ASP:C	4:D:307:LEU:CD2	2.70	0.48
1:L:4:TYR:HA	1:L:60:ILE:O	2.13	0.48
4:G:25:MET:HB3	4:G:93:VAL:HG22	1.95	0.48
4:G:94:ARG:NE	4:G:212:GLU:OE1	2.43	0.48
4:G:150:ARG:NH2	6:M:18:G:OP1	2.44	0.48
4:G:22:ASP:OD1	4:H:83:ASN:ND2	2.47	0.48
4:C:29:ALA:HB3	4:C:32:GLN:HG3	1.96	0.47
4:D:8:THR:CG2	4:D:9:ALA:H	2.27	0.47
4:H:77:GLN:NE2	5:J:93:TYR:OH	2.47	0.47
4:D:265:ARG:NH2	6:M:32:G:OP2	2.43	0.47
1:L:155:PHE:CD1	1:L:155:PHE:C	2.85	0.47
3:B:263:LEU:HB2	3:B:266:GLU:HB2	1.96	0.47
4:H:99:VAL:O	4:H:208:HIS:HA	2.15	0.47
3:B:105:LEU:HD12	3:B:105:LEU:O	2.14	0.47
5:J:38:VAL:HG23	5:J:38:VAL:O	2.14	0.47
3:B:17:ILE:HG22	3:B:144:ILE:HG12	1.97	0.47
4:E:34:ASP:HA	4:E:161:ARG:HH12	1.80	0.47
4:E:235:LYS:CE	4:E:235:LYS:HA	2.45	0.47
4:H:165:ILE:HG22	4:H:170:VAL:HA	1.96	0.47
6:M:56:G:H1'	6:M:57:G:C5	2.49	0.47
1:L:96:PRO:HB2	1:L:164:LEU:HD22	1.97	0.47
1:L:149:GLN:N	6:M:60:G:H5'	2.28	0.47
3:B:122:PRO:HG2	3:B:125:GLU:HB3	1.96	0.47
4:E:80:ASP:OD1	4:E:242:LYS:NZ	2.48	0.47
4:H:266:THR:HA	4:H:281:ALA:HA	1.97	0.47
3:B:215:ASP:OD1	3:B:215:ASP:N	2.48	0.46
4:F:77:GLN:NE2	5:I:93:TYR:OH	2.42	0.46
4:G:228:SER:HB2	4:G:255:ILE:HA	1.96	0.46
4:E:291:GLN:OE1	4:E:295:TYR:OH	2.33	0.46
4:H:137:ARG:HB3	4:H:267:ILE:HB	1.96	0.46
2:A:182:LEU:HD21	2:A:188:HIS:HD2	1.80	0.46
4:F:227:PRO:HG2	4:F:246:LEU:HD22	1.97	0.46
4:F:307:LEU:HD21	4:F:322:GLN:HB3	1.97	0.46
4:H:174:TRP:HH2	4:H:198:LEU:HD13	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:149:GLN:CB	6:M:60:G:C2'	2.92	0.46
4:E:36:SER:O	4:E:92:LYS:NZ	2.47	0.46
4:G:257:SER:OG	4:H:47:LYS:NZ	2.45	0.46
4:C:249:VAL:HG12	4:C:250:ARG:H	1.80	0.46
4:E:266:THR:HA	4:E:281:ALA:HA	1.97	0.46
5:I:32:LEU:HD21	5:I:67:ILE:HG12	1.96	0.46
4:F:36:SER:O	4:F:92:LYS:NZ	2.40	0.46
4:F:89:ASP:OD1	4:F:89:ASP:N	2.43	0.46
1:L:98:ARG:HH11	1:L:161:HIS:HB3	1.80	0.46
2:A:233:SER:CB	3:B:216:LEU:HD13	2.46	0.46
4:D:283:GLU:O	4:D:325:TYR:OH	2.29	0.46
4:E:50:ARG:O	6:M:31:C:O2'	2.33	0.46
4:F:75:ASN:OD1	5:I:84:ARG:NH2	2.47	0.46
4:G:172:ARG:NH2	4:G:191:GLU:OE1	2.48	0.46
4:G:259:LYS:NZ	6:M:16:C:OP2	2.47	0.46
4:E:76:LEU:CD2	6:M:33:U:O2	2.62	0.46
4:H:288:VAL:HG12	4:H:291:GLN:H	1.81	0.46
1:L:143:PHE:HA	1:L:157:LEU:O	2.16	0.46
4:C:152:ARG:NH2	4:C:182:GLY:O	2.48	0.46
1:L:155:PHE:CZ	1:L:156:ARG:C	2.88	0.46
4:D:303:ASP:OD2	4:D:306:THR:N	2.43	0.46
4:E:91:LEU:HD13	4:E:225:VAL:HG11	1.96	0.46
4:E:100:LEU:HD22	4:F:154:GLY:HA3	1.98	0.46
4:H:140:HIS:O	4:H:144:ASN:ND2	2.48	0.45
4:D:31:ALA:HA	4:D:33:ARG:HH11	1.82	0.45
4:F:166:ARG:HG2	4:F:171:ALA:HB2	1.98	0.45
4:E:156:GLU:HB2	4:E:218:ARG:HB3	1.98	0.45
4:H:68:ASP:HA	4:H:71:ILE:CG2	2.45	0.45
4:D:161:ARG:HE	4:D:175:ARG:HD2	1.81	0.45
4:E:231:LEU:HD13	4:F:76:LEU:HD13	1.98	0.45
4:F:53:ILE:HD13	5:I:84:ARG:HD3	1.98	0.45
4:H:91:LEU:HD13	4:H:225:VAL:HG11	1.97	0.45
2:A:93:GLY:HA3	3:B:190:GLU:HB3	1.98	0.45
3:B:300:ASP:HA	3:B:320:ARG:HG3	1.99	0.45
4:D:266:THR:HA	4:D:281:ALA:HA	1.99	0.45
4:H:94:ARG:HG3	4:H:214:VAL:CG2	2.24	0.45
1:L:16:PRO:HG2	1:L:19:GLN:HG3	1.99	0.44
1:L:155:PHE:HD2	6:M:46:C:C2	2.35	0.44
1:L:155:PHE:HB2	6:M:46:C:C2	2.52	0.44
4:H:6:LEU:CD1	4:H:313:LEU:CD2	2.95	0.44
4:H:189:ASP:HB3	4:H:192:LEU:HD13	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:17:ILE:HG12	3:B:139:LEU:HD21	1.99	0.44
1:L:13:PRO:CG	4:C:277:LEU:HD13	2.46	0.44
1:L:164:LEU:HD23	1:L:164:LEU:HA	1.90	0.44
3:B:5:ASP:OD1	3:B:5:ASP:N	2.51	0.44
3:B:85:ARG:NH2	6:M:9:U:O2	2.48	0.44
4:F:164:HIS:HE1	4:F:209:VAL:HG12	1.82	0.44
4:G:146:ARG:NH1	4:G:183:LEU:HD11	2.32	0.44
5:I:3:LYS:HE3	5:I:3:LYS:HB3	1.80	0.44
1:L:113:ARG:HB3	1:L:117:MET:HE3	1.99	0.44
3:B:41:VAL:HG21	3:B:59:VAL:HG12	1.99	0.44
4:F:266:THR:HA	4:F:281:ALA:HA	1.99	0.44
3:B:159:LEU:HD11	3:B:172:ASN:HB3	2.00	0.44
4:G:104:GLY:HA3	4:G:122:VAL:HG11	1.99	0.44
4:G:112:ALA:HA	4:G:115:ARG:HG2	1.99	0.44
4:G:288:VAL:CG2	4:G:291:GLN:HB2	2.47	0.44
4:H:99:VAL:HB	4:H:209:VAL:HG12	2.00	0.44
4:H:163:ASN:HB2	4:H:170:VAL:HG13	1.99	0.44
4:E:249:VAL:CG1	4:E:250:ARG:H	2.22	0.44
3:B:216:LEU:HD12	3:B:217:CYS:C	2.38	0.44
4:F:148:LEU:HB2	4:F:152:ARG:HB2	2.00	0.44
4:F:259:LYS:NZ	6:M:22:U:OP1	2.40	0.44
3:B:54:ILE:HA	3:B:115:GLY:HA3	1.98	0.44
2:A:233:SER:HB3	3:B:218:ARG:HA	1.99	0.44
1:L:155:PHE:CE1	1:L:157:LEU:N	2.86	0.43
4:E:145:ALA:HB1	4:E:148:LEU:HD12	2.00	0.43
4:D:91:LEU:HD13	4:D:225:VAL:HG11	2.00	0.43
4:G:145:ALA:H	4:G:178:ALA:HB1	1.82	0.43
1:L:9:LEU:HD12	1:L:56:GLU:HA	2.00	0.43
1:L:107:SER:HA	1:L:137:ARG:H	1.84	0.43
4:D:16:ARG:NH2	4:E:224:GLU:OE1	2.37	0.43
4:H:332:ARG:NH1	4:H:333:GLY:O	2.52	0.43
3:B:48:VAL:HB	3:B:133:GLN:HG2	2.00	0.43
4:E:235:LYS:HE2	4:E:235:LYS:HA	2.00	0.43
4:E:329:ASN:HA	4:E:332:ARG:HB3	2.01	0.43
5:J:83:PHE:HB3	5:J:113:LEU:HD11	2.01	0.43
3:B:52:LEU:HD13	3:B:126:ILE:HG13	1.99	0.43
4:D:33:ARG:NE	4:D:159:GLU:OE2	2.39	0.43
4:F:94:ARG:HB2	4:F:214:VAL:HG22	2.00	0.43
4:E:280:ILE:HD12	4:E:288:VAL:HG11	2.00	0.43
4:F:24:LEU:HD11	4:F:249:VAL:HG21	1.99	0.43
4:F:137:ARG:HH12	4:F:269:THR:HG23	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:163:ASN:HD22	4:G:173:THR:HG22	1.84	0.43
1:L:155:PHE:CG	1:L:156:ARG:N	2.84	0.43
4:C:18:LEU:HD23	4:C:99:VAL:HG12	2.00	0.43
4:E:57:LEU:HD11	4:E:67:LEU:HD21	2.01	0.43
4:F:50:ARG:HB2	6:M:27:U:H5"	2.00	0.43
4:G:303:ASP:OD1	4:G:306:THR:N	2.52	0.43
4:H:162:ILE:HG22	4:H:213:VAL:HA	2.00	0.43
5:I:104:PHE:HZ	5:I:107:LYS:HG2	1.82	0.43
4:E:234:ASP:OD1	4:E:234:ASP:N	2.47	0.43
4:G:298:PRO:HB2	4:H:62:ARG:HD2	2.01	0.43
4:E:99:VAL:HG21	4:E:203:LEU:CG	2.49	0.42
4:H:6:LEU:C	4:H:6:LEU:HD12	2.39	0.42
5:I:28:ILE:HA	5:I:31:LYS:HB3	2.01	0.42
4:C:92:LYS:HE2	4:C:94:ARG:HB2	2.01	0.42
3:B:138:ARG:HH22	4:H:16:ARG:HH21	1.67	0.42
4:C:268:ASP:HB3	4:C:280:ILE:HG13	2.00	0.42
4:F:313:LEU:HB2	4:F:314:ARG:HD2	2.02	0.42
1:L:28:LEU:HD11	1:L:60:ILE:HD13	2.01	0.42
4:C:141:ASN:OD1	4:C:266:THR:OG1	2.36	0.42
4:D:18:LEU:CD2	4:D:331:ILE:CG2	2.82	0.42
4:H:145:ALA:H	4:H:178:ALA:HB1	1.83	0.42
3:B:267:VAL:CG1	3:B:277:LEU:HD13	2.47	0.42
4:G:146:ARG:CZ	4:G:183:LEU:CD1	2.97	0.42
4:E:137:ARG:HH12	4:E:269:THR:HG23	1.84	0.42
4:G:296:ARG:HH12	4:G:322:GLN:HE22	1.68	0.42
4:F:23:ALA:HA	4:F:94:ARG:O	2.19	0.42
4:F:70:SER:O	5:I:84:ARG:NH1	2.53	0.42
4:H:214:VAL:HG11	4:H:216:PHE:CZ	2.54	0.42
4:H:19:ASP:OD1	4:H:19:ASP:N	2.39	0.42
4:H:23:ALA:HA	4:H:95:PHE:HB3	2.02	0.42
4:H:52:THR:HB	4:H:71:ILE:CD1	2.50	0.42
5:I:93:TYR:HB3	5:I:102:TYR:HB3	2.01	0.42
5:J:13:ARG:HE	5:J:75:ASN:HB3	1.84	0.42
4:F:50:ARG:O	6:M:25:U:O2'	2.35	0.42
4:F:94:ARG:HH21	4:F:212:GLU:HG3	1.84	0.42
3:B:306:HIS:CE1	3:B:315:ARG:HH21	2.38	0.41
4:G:290:SER:OG	4:G:291:GLN:N	2.52	0.41
3:B:105:LEU:HD12	3:B:105:LEU:C	2.41	0.41
4:H:162:ILE:HD13	4:H:162:ILE:HG21	1.86	0.41
3:B:216:LEU:CD1	3:B:217:CYS:C	2.88	0.41
4:D:17:LYS:C	4:D:18:LEU:HG	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:317:ALA:HA	4:H:318:PRO:HD3	1.87	0.41
5:I:82:LYS:NZ	5:I:120:ASN:O	2.54	0.41
3:B:37:PHE:CE1	3:B:139:LEU:HD11	2.55	0.41
4:F:283:GLU:OE2	4:F:287:SER:OG	2.39	0.41
3:B:186:LEU:HA	3:B:288:GLU:O	2.20	0.41
4:E:232:ILE:O	5:I:99:HIS:NE2	2.48	0.41
4:C:176:PHE:HE2	4:C:192:LEU:HD12	1.85	0.41
4:C:183:LEU:HD12	4:C:184:ARG:HB3	2.02	0.41
5:I:25:ILE:HA	5:I:28:ILE:HG12	2.02	0.41
1:L:155:PHE:CD2	6:M:46:C:N4	2.81	0.41
3:B:159:LEU:HD22	3:B:176:LEU:HD22	2.02	0.41
4:E:67:LEU:HD23	4:E:67:LEU:HA	1.90	0.41
4:E:98:ARG:HB3	4:E:210:LEU:HD12	2.03	0.41
4:E:174:TRP:CD1	4:E:191:GLU:HG2	2.56	0.41
4:H:6:LEU:HD12	4:H:313:LEU:HD22	2.02	0.41
4:H:239:LYS:HE2	4:H:239:LYS:HB2	1.87	0.41
6:M:56:G:H2'	6:M:57:G:OP2	2.20	0.41
1:L:5:LEU:HB3	1:L:60:ILE:HG22	2.02	0.41
4:D:305:TYR:HD1	4:E:56:ARG:HH21	1.68	0.41
4:G:18:LEU:HD13	4:G:331:ILE:HG23	1.95	0.41
4:D:14:PHE:HZ	4:D:118:LEU:HD21	1.87	0.40
4:D:117:LYS:HA	4:D:117:LYS:HD3	1.92	0.40
4:D:319:ALA:HB3	4:D:322:GLN:HG2	2.04	0.40
4:E:235:LYS:HA	4:E:235:LYS:HD3	1.91	0.40
4:F:158:VAL:HB	4:F:178:ALA:HB3	2.03	0.40
4:H:68:ASP:CA	4:H:71:ILE:HG22	2.50	0.40
4:D:18:LEU:HD21	4:D:331:ILE:HG23	1.93	0.40
4:E:303:ASP:OD1	4:E:303:ASP:N	2.52	0.40
4:F:105:THR:HA	4:F:119:LEU:HD11	2.03	0.40
4:E:269:THR:HG22	4:E:279:PRO:HB3	2.03	0.40
1:L:149:GLN:CB	6:M:60:G:H3'	2.51	0.40
1:L:155:PHE:CE1	1:L:157:LEU:CA	3.05	0.40
4:E:80:ASP:OD1	4:E:242:LYS:HE3	2.18	0.40
4:G:124:THR:O	4:G:128:GLU:HB3	2.21	0.40
4:H:22:ASP:OD2	4:H:247:TYR:OH	2.32	0.40
6:M:1:C:O2	6:M:1:C:C3'	2.70	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	L	182/187 (97%)	176 (97%)	6 (3%)	0	100	100
2	A	422/434 (97%)	369 (87%)	53 (13%)	0	100	100
3	B	301/327 (92%)	269 (89%)	31 (10%)	1 (0%)	41	75
4	C	287/342 (84%)	269 (94%)	18 (6%)	0	100	100
4	D	331/342 (97%)	311 (94%)	19 (6%)	1 (0%)	41	75
4	E	332/342 (97%)	316 (95%)	16 (5%)	0	100	100
4	F	333/342 (97%)	305 (92%)	28 (8%)	0	100	100
4	G	331/342 (97%)	308 (93%)	23 (7%)	0	100	100
4	H	329/342 (96%)	304 (92%)	25 (8%)	0	100	100
5	I	122/124 (98%)	114 (93%)	8 (7%)	0	100	100
5	J	122/124 (98%)	116 (95%)	6 (5%)	0	100	100
All	All	3092/3248 (95%)	2857 (92%)	233 (8%)	2 (0%)	54	83

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	303	ASP
3	B	153	PRO

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	L	142/159 (89%)	140 (99%)	2 (1%)	67	85
2	A	76/365 (21%)	74 (97%)	2 (3%)	46	74
3	B	243/270 (90%)	238 (98%)	5 (2%)	53	78
4	C	228/274 (83%)	224 (98%)	4 (2%)	59	81
4	D	257/274 (94%)	252 (98%)	5 (2%)	57	80
4	E	259/274 (94%)	249 (96%)	10 (4%)	32	64
4	F	261/274 (95%)	258 (99%)	3 (1%)	73	88
4	G	260/274 (95%)	257 (99%)	3 (1%)	71	87
4	H	259/274 (94%)	254 (98%)	5 (2%)	57	80
5	I	106/108 (98%)	106 (100%)	0	100	100
5	J	101/108 (94%)	98 (97%)	3 (3%)	41	71
All	All	2192/2654 (83%)	2150 (98%)	42 (2%)	59	80

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	102	ARG
1	L	155	PHE
2	A	238	LEU
2	A	263	TRP
3	B	139	LEU
3	B	151	ARG
3	B	170	ARG
3	B	247	TRP
3	B	258	LEU
4	C	16	ARG
4	C	131	PHE
4	C	184	ARG
4	C	296	ARG
4	D	18	LEU
4	D	33	ARG
4	D	150	ARG
4	D	303	ASP
4	D	307	LEU
4	E	17	LYS
4	E	75	ASN
4	E	76	LEU
4	E	77	GLN
4	E	172	ARG

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Mol	Chain	Res	Type
4	E	235	LYS
4	E	238	LYS
4	E	285	TYR
4	E	303	ASP
4	E	322	GLN
4	F	172	ARG
4	F	198	LEU
4	F	233	LEU
4	G	144	ASN
4	G	285	TYR
4	G	309	ASP
4	H	21	SER
4	H	75	ASN
4	H	148	LEU
4	H	231	LEU
4	H	285	TYR
5	J	34	ASP
5	J	90	ASN
5	J	122	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	108	ASN
1	L	153	GLN
1	L	165	GLN
2	A	177	GLN
2	A	188	HIS
2	A	225	GLN
3	B	18	GLN
3	B	64	HIS
3	B	82	ASN
3	B	89	ASN
3	B	104	HIS
3	B	172	ASN
3	B	173	GLN
3	B	256	ASN
4	C	300	GLN
4	C	329	ASN
4	D	141	ASN
4	D	163	ASN
4	D	164	HIS

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Mol	Chain	Res	Type
4	D	223	GLN
4	E	127	ASN
4	E	163	ASN
4	E	223	GLN
4	F	144	ASN
4	F	163	ASN
4	F	208	HIS
4	G	144	ASN
4	G	163	ASN
4	G	297	GLN
4	G	310	ASN
4	G	322	GLN
4	G	329	ASN
4	H	75	ASN
4	H	77	GLN
4	H	129	GLN
4	H	167	GLN
4	H	208	HIS
4	H	310	ASN
4	H	329	ASN
5	I	53	GLN
5	J	29	ASN
5	J	55	GLN

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
6	M	59/60 (98%)	35 (59%)	2 (3%)

All (35) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
6	M	3	A
6	M	6	A
6	M	9	U
6	M	15	G
6	M	16	C
6	M	17	G
6	M	19	G
6	M	20	C
6	M	21	U

Continued on next page...

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Mol	Chain	Res	Type
6	M	26	G
6	M	27	U
6	M	32	G
6	M	33	U
6	M	34	C
6	M	35	U
6	M	39	U
6	M	40	G
6	M	41	G
6	M	42	U
6	M	43	U
6	M	44	C
6	M	45	A
6	M	46	C
6	M	48	G
6	M	49	C
6	M	50	C
6	M	51	G
6	M	52	U
6	M	53	G
6	M	54	U
6	M	56	G
6	M	57	G
6	M	58	C
6	M	59	A
6	M	60	G

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
6	M	56	G
6	M	58	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 (i)

There are no chain breaks in this entry.

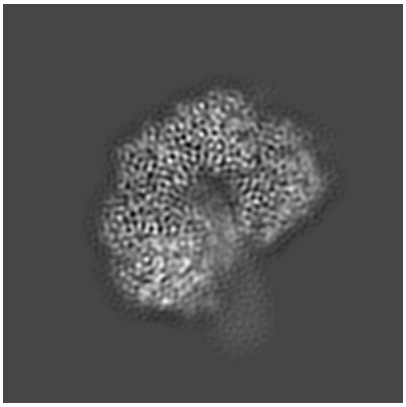
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-31058. These allow visual inspection of the internal detail of the map and identification of artifacts.

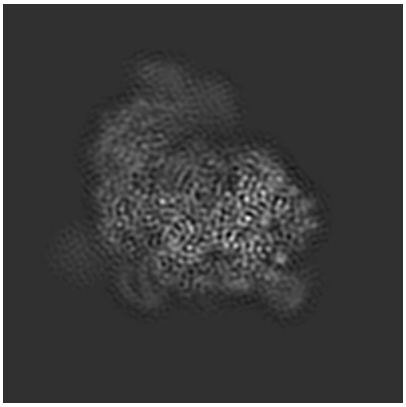
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

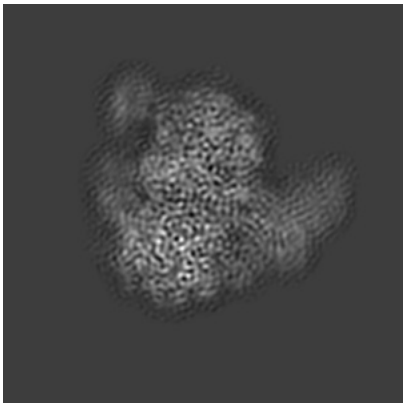
6.1.1 Primary map



X



Y

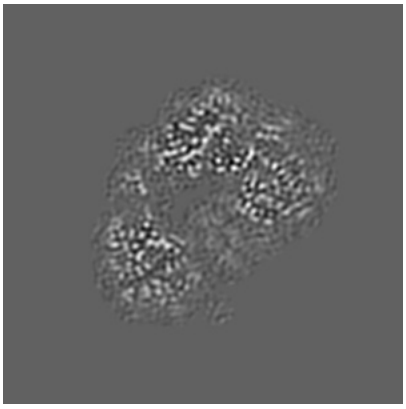


Z

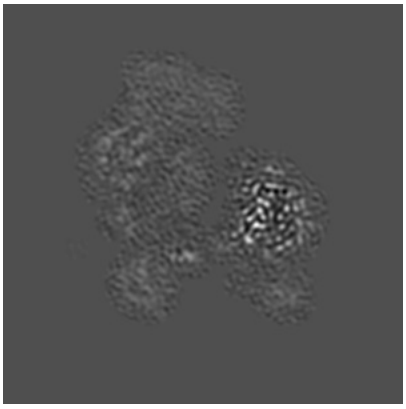
The images above show the map projected in three orthogonal directions.

6.2 Central slices

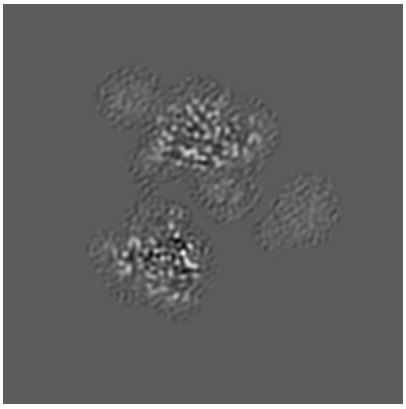
6.2.1 Primary map



X Index: 100



Y Index: 100

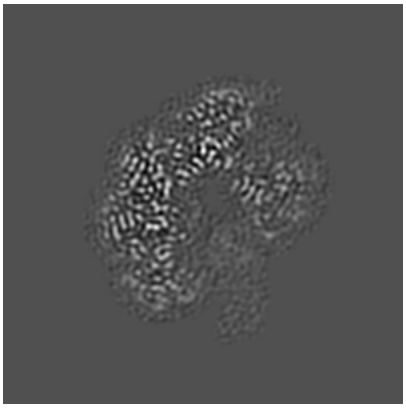


Z Index: 100

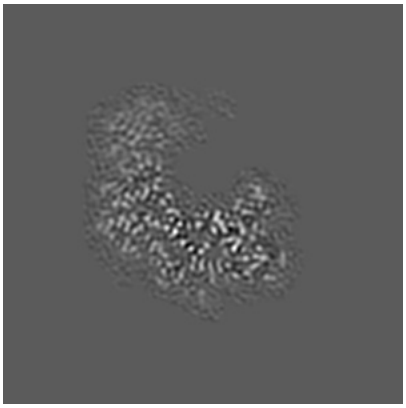
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

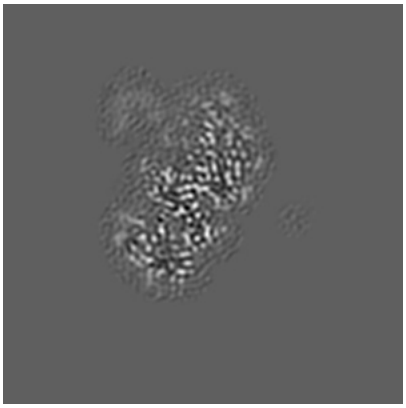
6.3.1 Primary map



X Index: 89



Y Index: 77



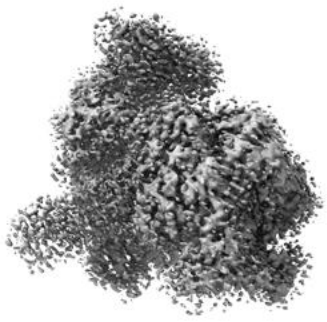
Z Index: 121

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

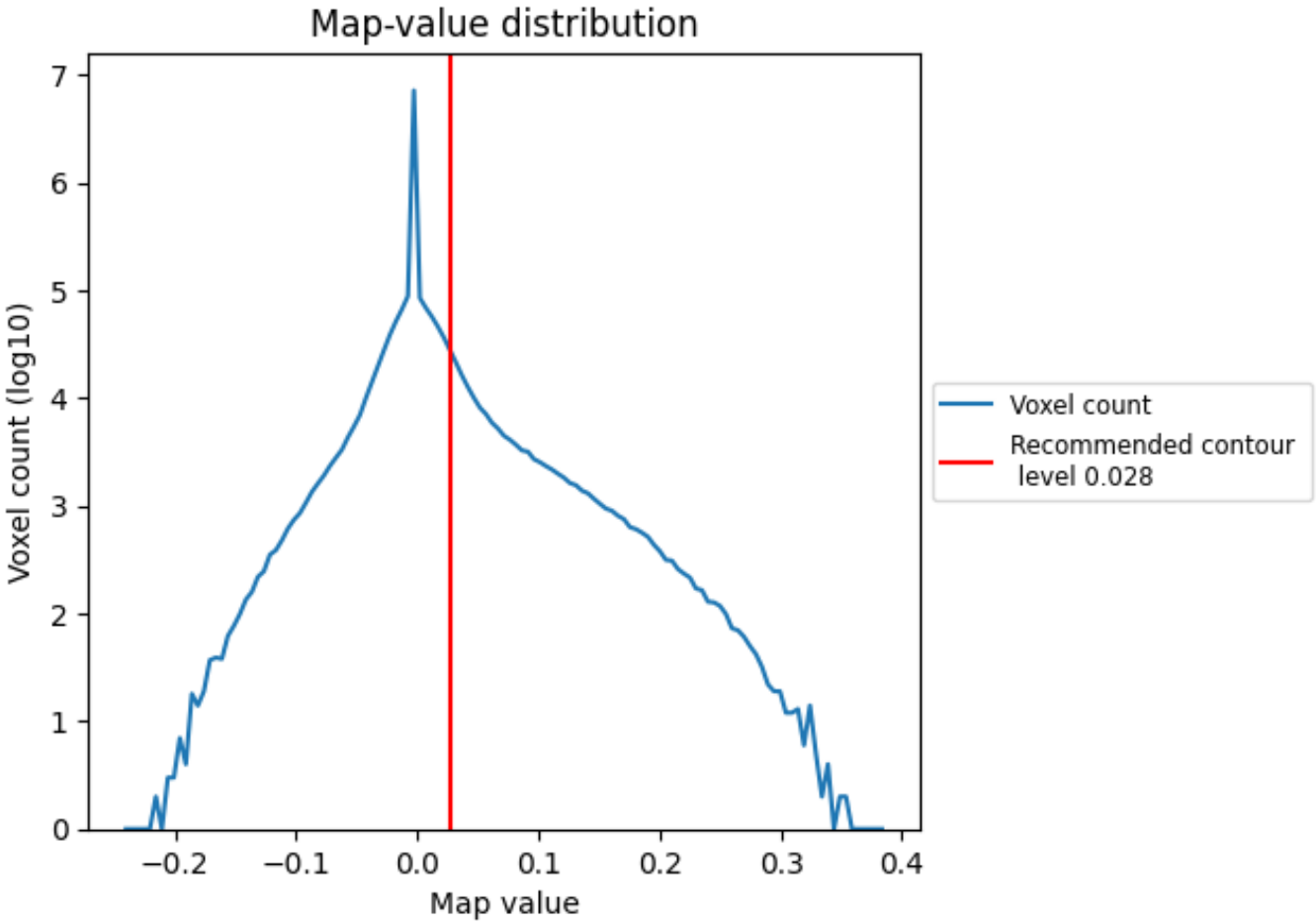
6.5 Mask visualisation (i)

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

7 Map analysis

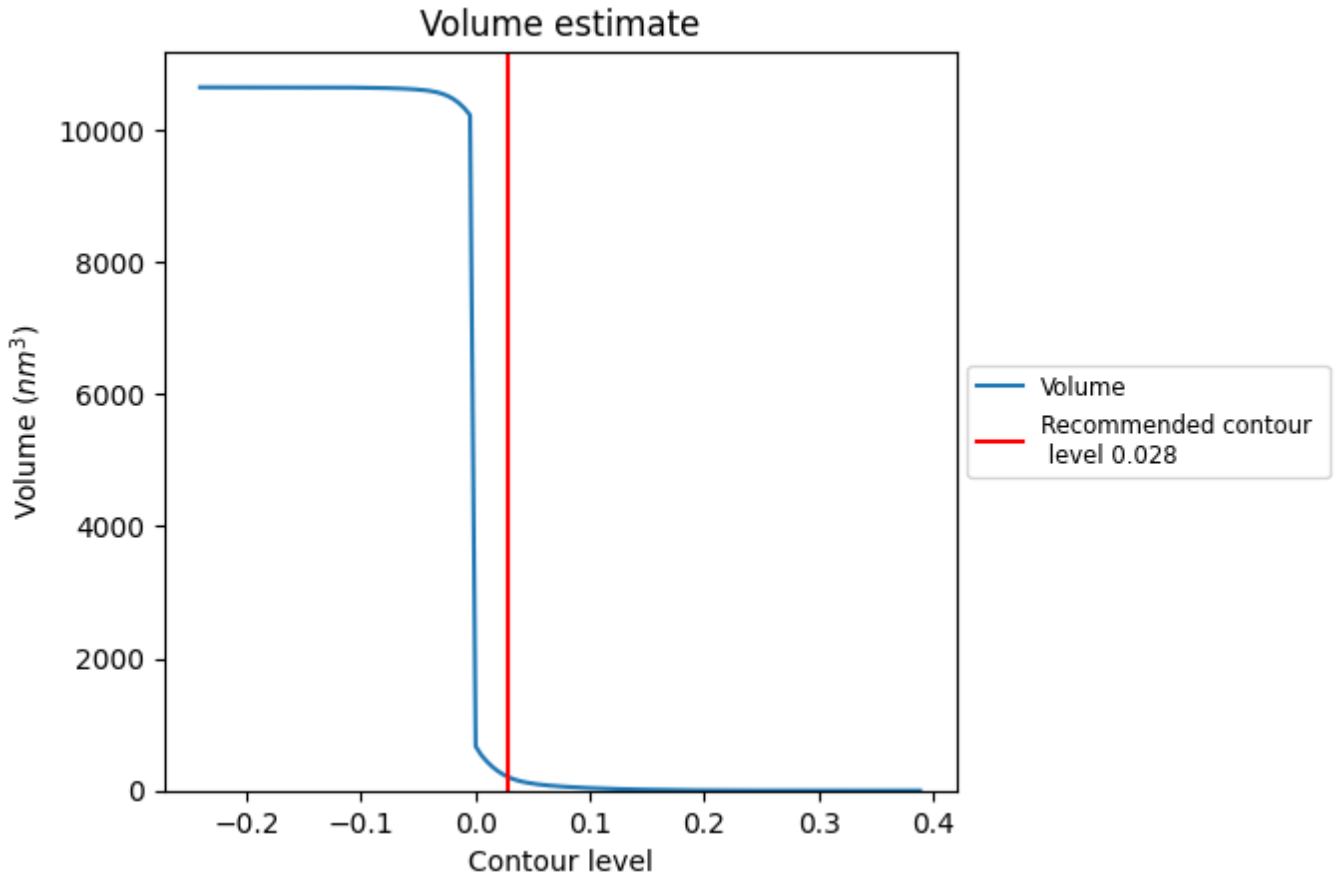
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

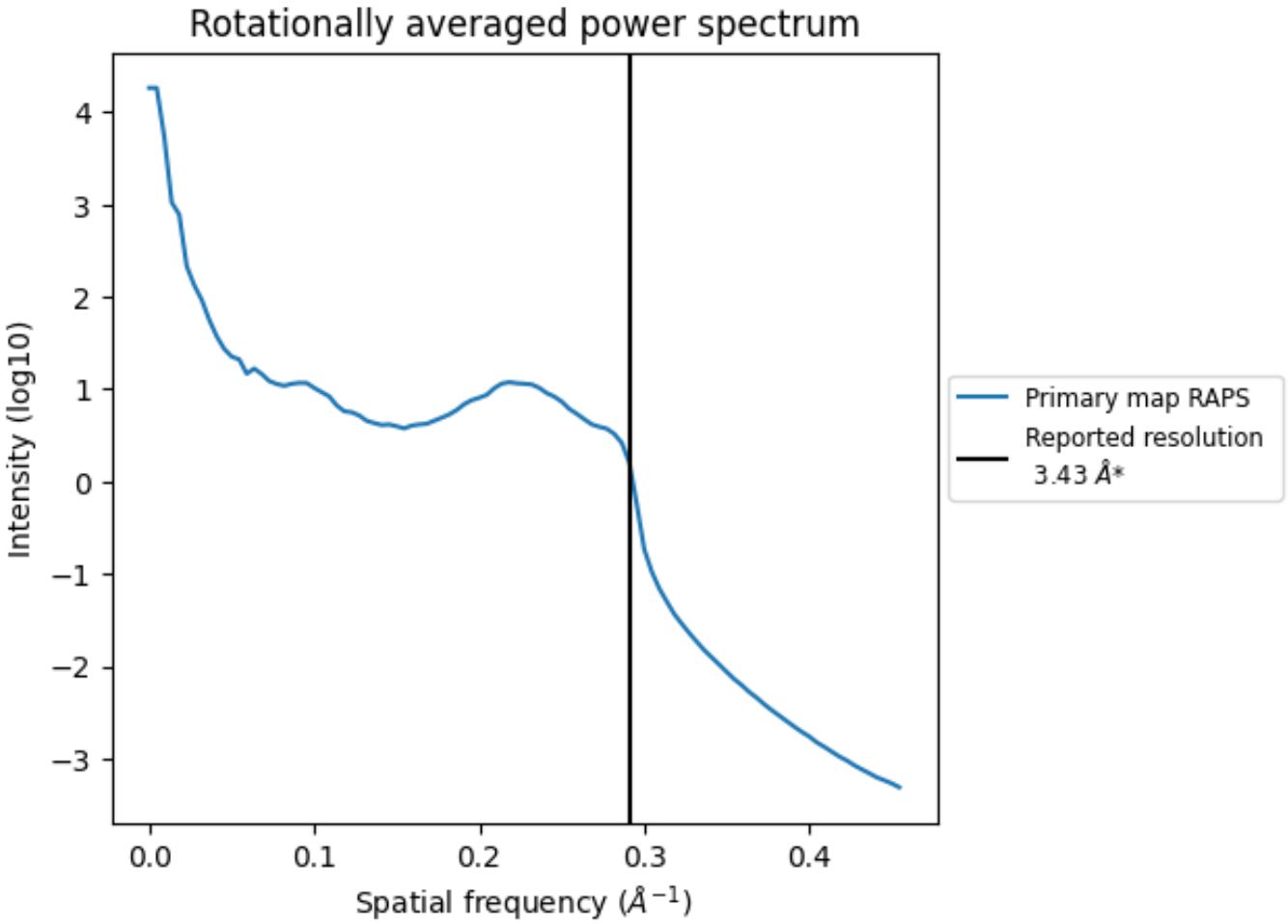
7.2 Volume estimate (i)



The volume at the recommended contour level is 212 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 (i)

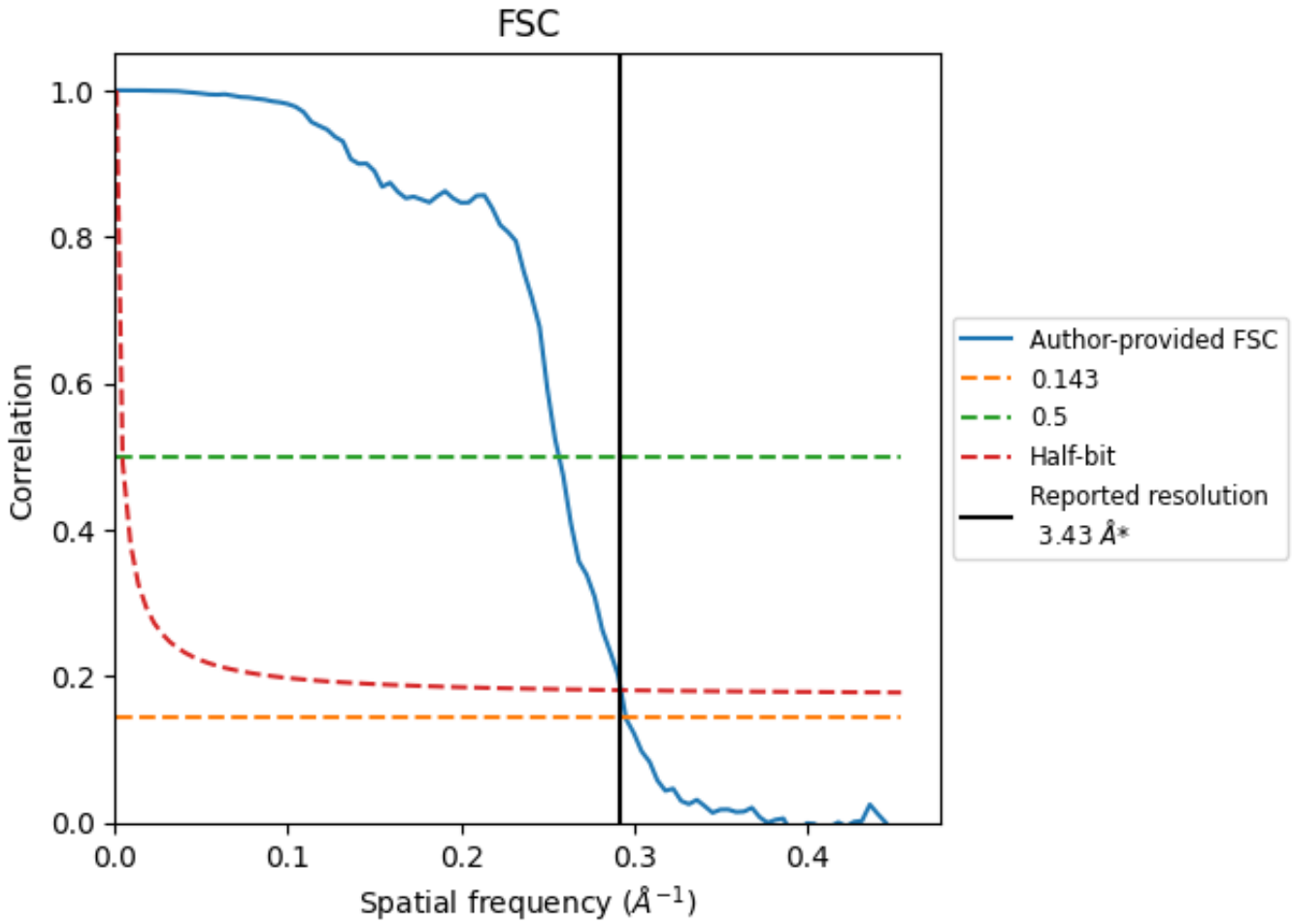


*Reported resolution corresponds to spatial frequency of 0.292 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.292 Å^{-1}

8.2 Resolution estimates (i)

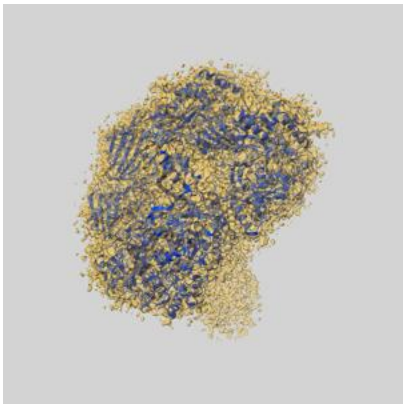
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.43	-	-
Author-provided FSC curve	3.39	3.90	3.42
Unmasked-calculated*	-	-	-

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

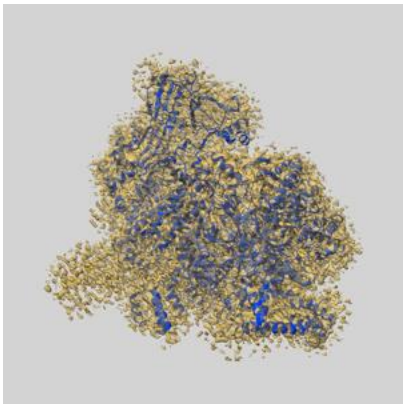
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-31058 and PDB model 7ECV. Per-residue inclusion information can be found in section 3 on page 6.

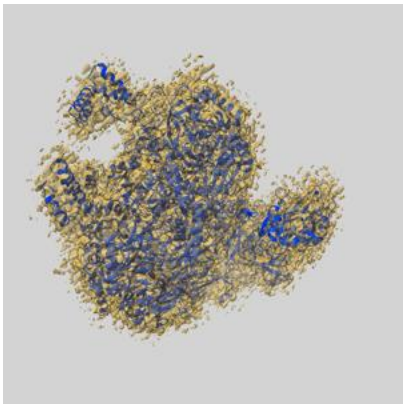
9.1 Map-model overlay



X



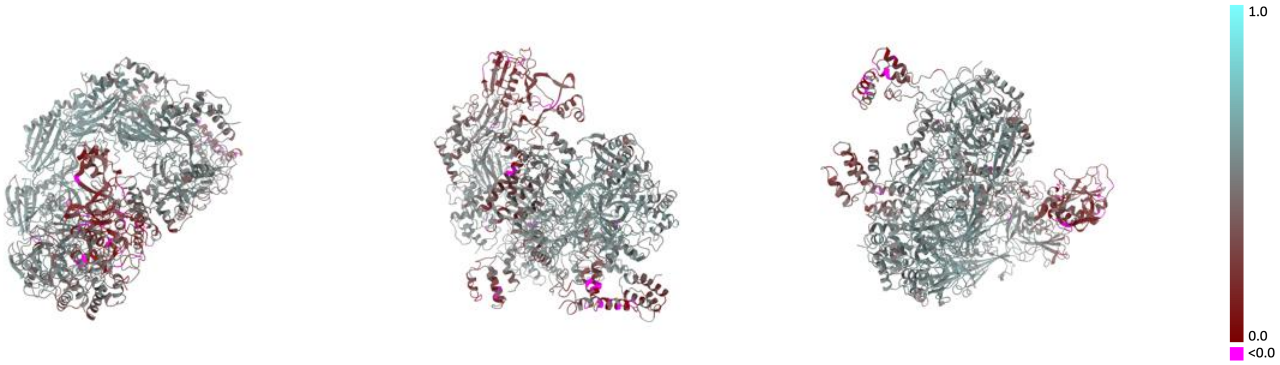
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.028 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.

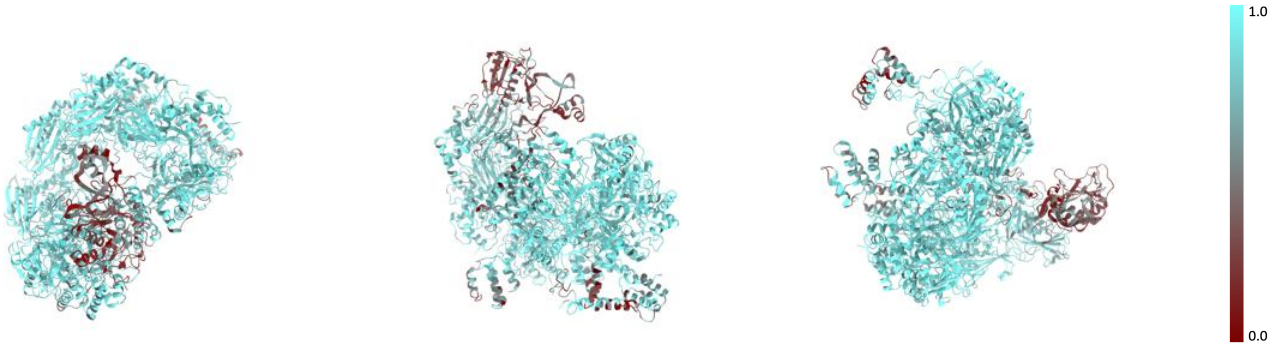
9.2 Q-score mapped to coordinate model (i)



X Y Z

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

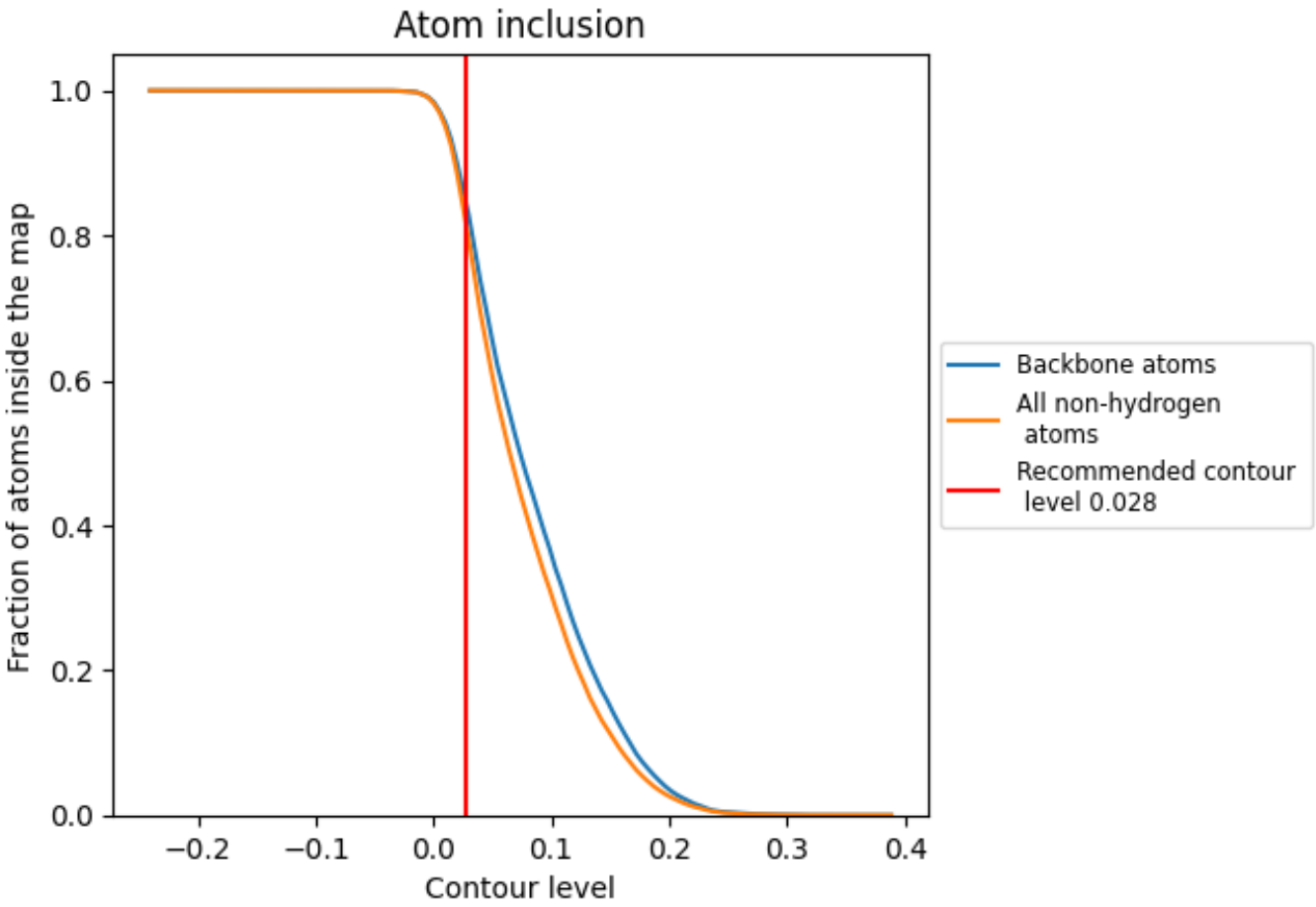
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.028).

9.4 Atom inclusion (i)



At the recommended contour level, 84% of all backbone atoms, 81% 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 (0.028) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8136	0.4710
A	0.7463	0.3890
B	0.8498	0.4860
C	0.7770	0.4290
D	0.8620	0.5050
E	0.8931	0.5290
F	0.9067	0.5400
G	0.9095	0.5410
H	0.9005	0.5290
I	0.7449	0.4110
J	0.7952	0.4390
L	0.3206	0.2100
M	0.7319	0.4280