



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 12:44 AM UTC

PDB ID : 6TLJ / pdb_00006tlj
EMDB ID : EMD-10516
Title : Cryo-EM structure of the Anaphase-promoting complex/Cyclosome, in complex with the Mitotic checkpoint complex (APC/C-MCC) at 3.8 angstrom resolution
Authors : Alfieri, C.; Barford, D.
Deposited on : 2019-12-02
Resolution : 3.80 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

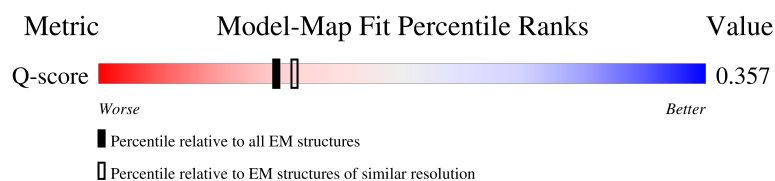
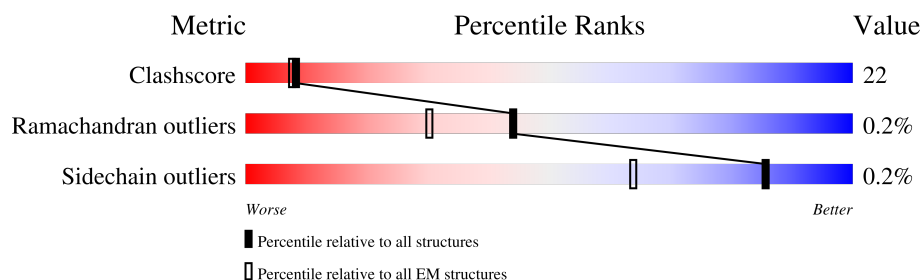
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 3.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10198 (3.30 - 4.30)

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

Mol	Chain	Length	Quality of chain
1	A	1944	19% 43% 30% 26%
2	B	84	92% 43% 49% 6%
3	C	597	21% 53% 35% 12%
3	P	597	17% 46% 36% 18%

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Mol	Chain	Length	Quality of chain
4	D	121	
5	E	110	
6	F	824	
6	H	824	
7	G	85	
7	W	85	
8	I	808	
9	J	620	
9	K	620	
10	L	185	
11	M	74	
12	N	822	
13	O	755	
14	Q	374	
15	R	499	
16	S	1050	
17	X	599	
17	Y	599	
18	Z	205	

2 Entry composition [i](#)

There are 18 unique types of molecules in this entry. The entry contains 72305 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 Anaphase-promoting complex subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1441	Total	C	N	O	S	0	0
			10949	7039	1853	1983	74		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	356	PHE	PRO	conflict	UNP Q9H1A4

- Molecule 2 is a protein called Anaphase-promoting complex subunit 11.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	79	Total	C	N	O	S	0	0
			643	411	116	100	16		

- Molecule 3 is a protein called Cell division cycle protein 23 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	524	Total	C	N	O	S	0	0
			4306	2774	727	781	24		
3	P	491	Total	C	N	O	S	0	0
			4043	2611	679	729	24		

- Molecule 4 is a protein called Anaphase-promoting complex subunit 15.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	D	18	Total	C	N	O	0	0
			153	104	23	26		

- Molecule 5 is a protein called Anaphase-promoting complex subunit 16.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	56	Total	C	N	O	S	0	0
			450	290	74	85	1		

- Molecule 6 is a protein called Cell division cycle protein 27 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	483	Total	C	N	O	S	0	0
			3849	2470	649	704	26		
6	H	483	Total	C	N	O	S	0	0
			3853	2473	650	704	26		

- Molecule 7 is a protein called Anaphase-promoting complex subunit CDC26.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	25	Total	C	N	O	S	0	0
			213	133	40	39	1		
7	W	25	Total	C	N	O	S	0	0
			213	133	40	39	1		

- Molecule 8 is a protein called Anaphase-promoting complex subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	I	733	Total	C	N	O	S	0	0
			5728	3675	953	1067	33		

- Molecule 9 is a protein called Cell division cycle protein 16 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	J	504	Total	C	N	O	S	0	0
			4047	2601	684	737	25		
9	K	493	Total	C	N	O	S	0	0
			3988	2563	672	729	24		

- Molecule 10 is a protein called Anaphase-promoting complex subunit 10.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	L	182	Total	C	N	O	S	0	0
			1435	898	263	268	6		

- Molecule 11 is a protein called Anaphase-promoting complex subunit 13.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	M	59	Total	C	N	O	S	0	0
			493	310	79	102	2		

- Molecule 12 is a protein called Anaphase-promoting complex subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	N	703	Total	C	N	O	S	0	0
			5403	3436	971	971	25		

- Molecule 13 is a protein called Anaphase-promoting complex subunit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	O	685	Total	C	N	O	S	0	0
			5402	3446	940	988	28		

- Molecule 14 is a protein called Cell division cycle protein 20 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	Q	365	Total	C	N	O	S	0	0
			2739	1720	503	505	11		

- Molecule 15 is a protein called Cell division cycle protein 20 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	R	383	Total	C	N	O	S	0	0
			2953	1855	538	548	12		

- Molecule 16 is a protein called Mitotic checkpoint serine/threonine-protein kinase BUB1 beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	S	293	Total	C	N	O	S	0	0
			2227	1397	405	420	5		

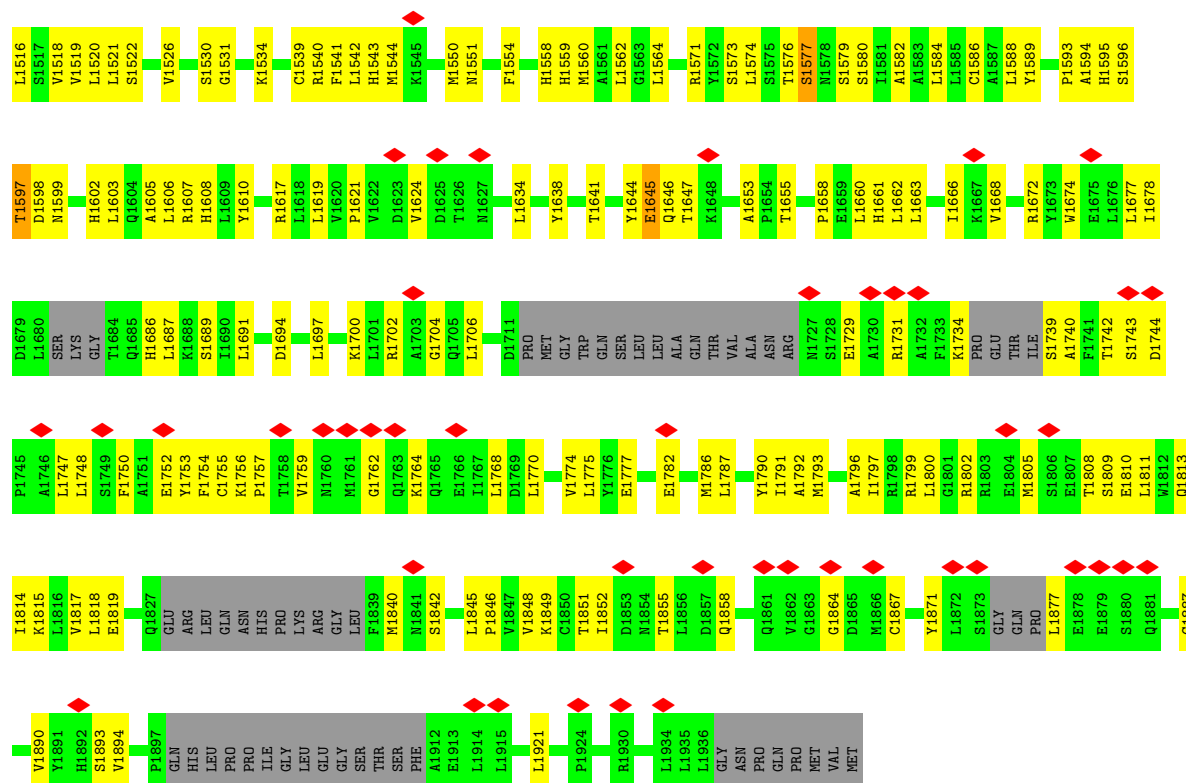
- Molecule 17 is a protein called Anaphase-promoting complex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	X	484	Total	C	N	O	S	0	0
			3773	2393	652	704	24		
17	Y	496	Total	C	N	O	S	0	0
			3868	2449	669	724	26		

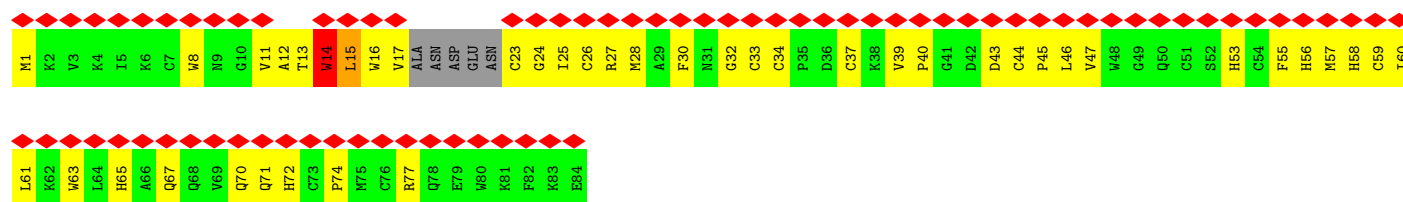
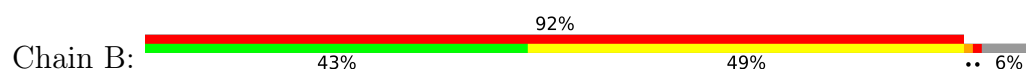
- Molecule 18 is a protein called Mitotic spindle assembly checkpoint protein MAD2A.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	Z	195	Total	C	N	O	S	0	0
			1577	1012	256	305	4		

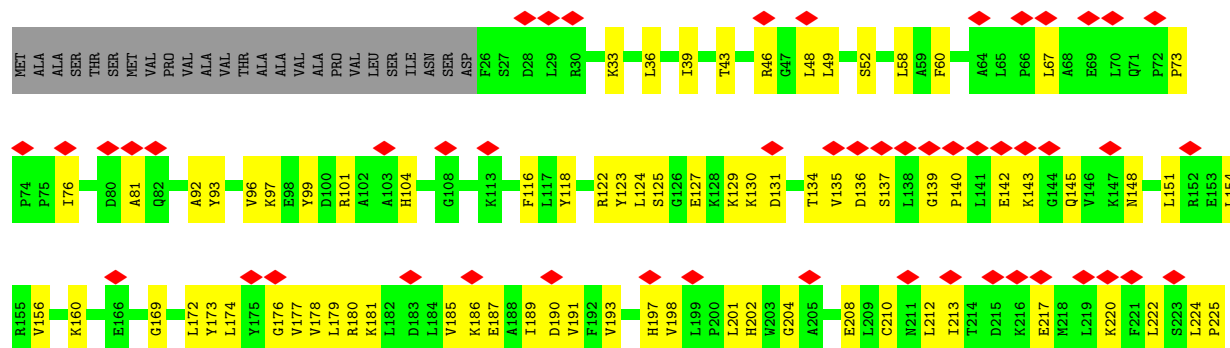


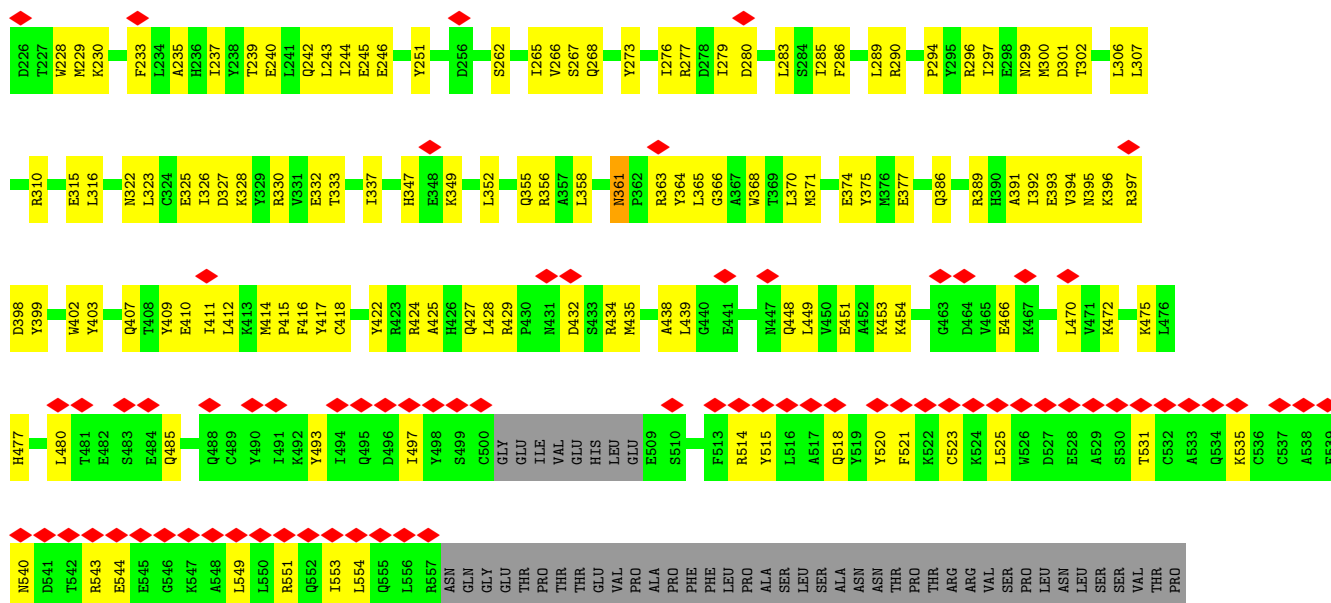


• Molecule 2: Anaphase-promoting complex subunit 11

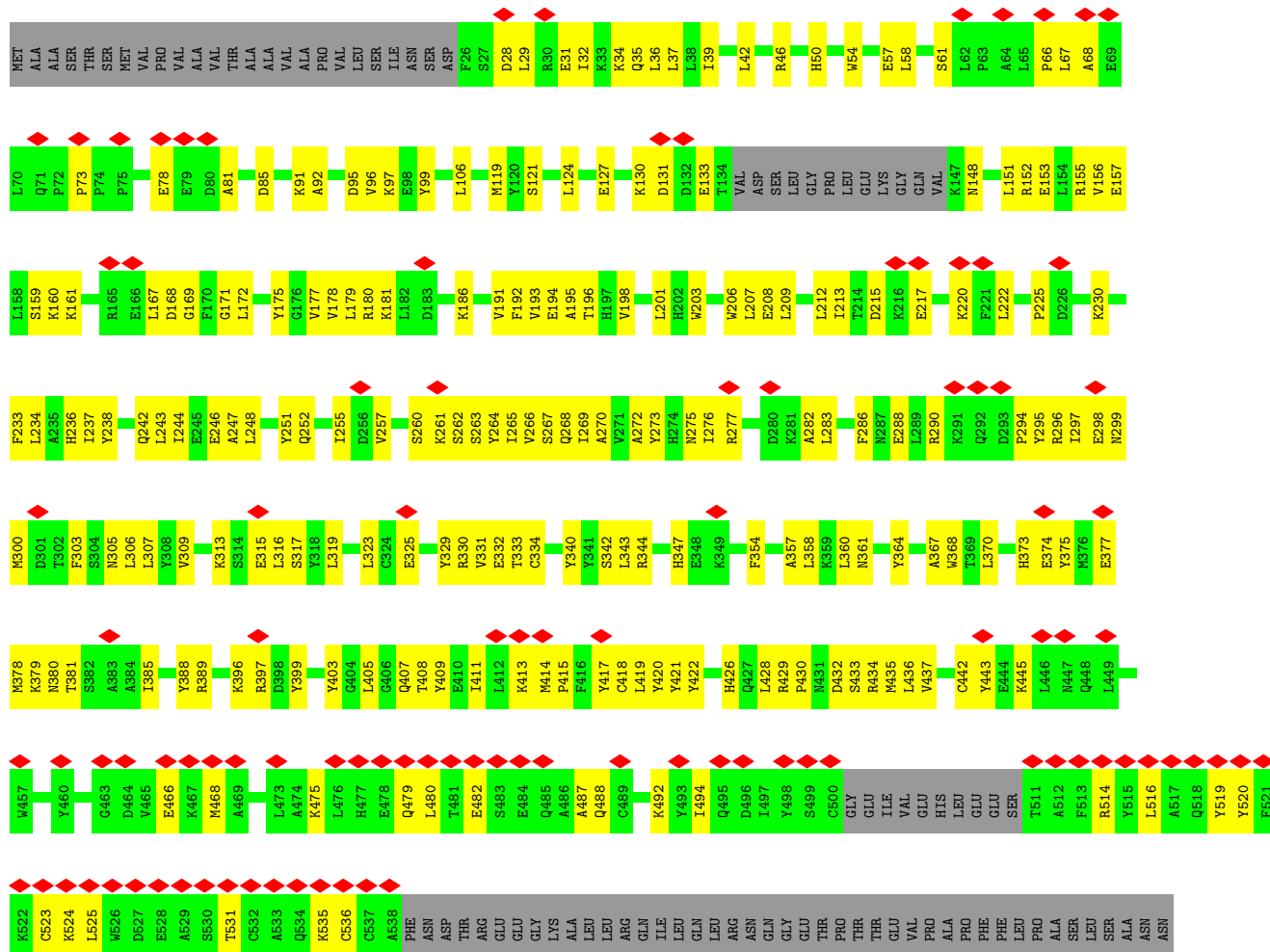


• Molecule 3: Cell division cycle protein 23 homolog

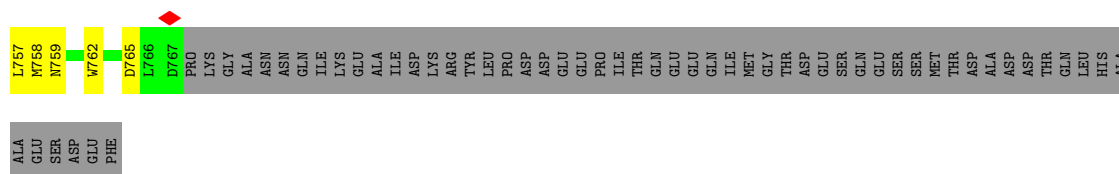




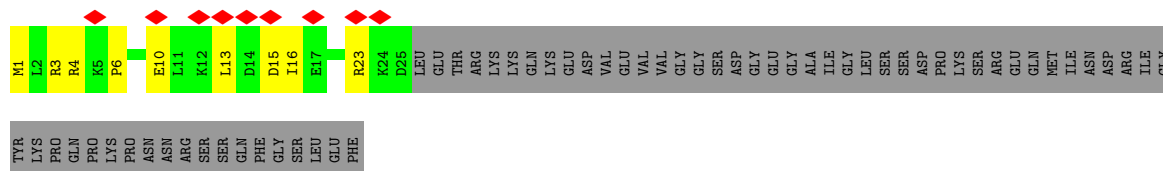
• Molecule 3: Cell division cycle protein 23 homolog



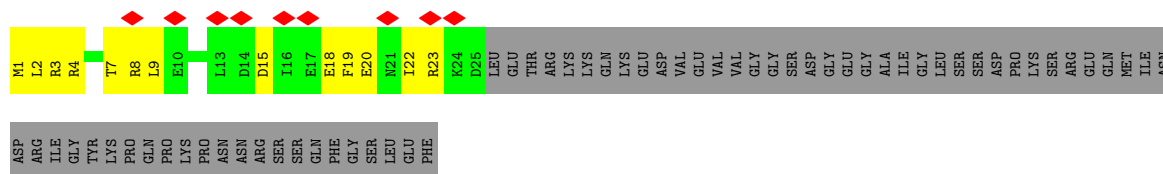




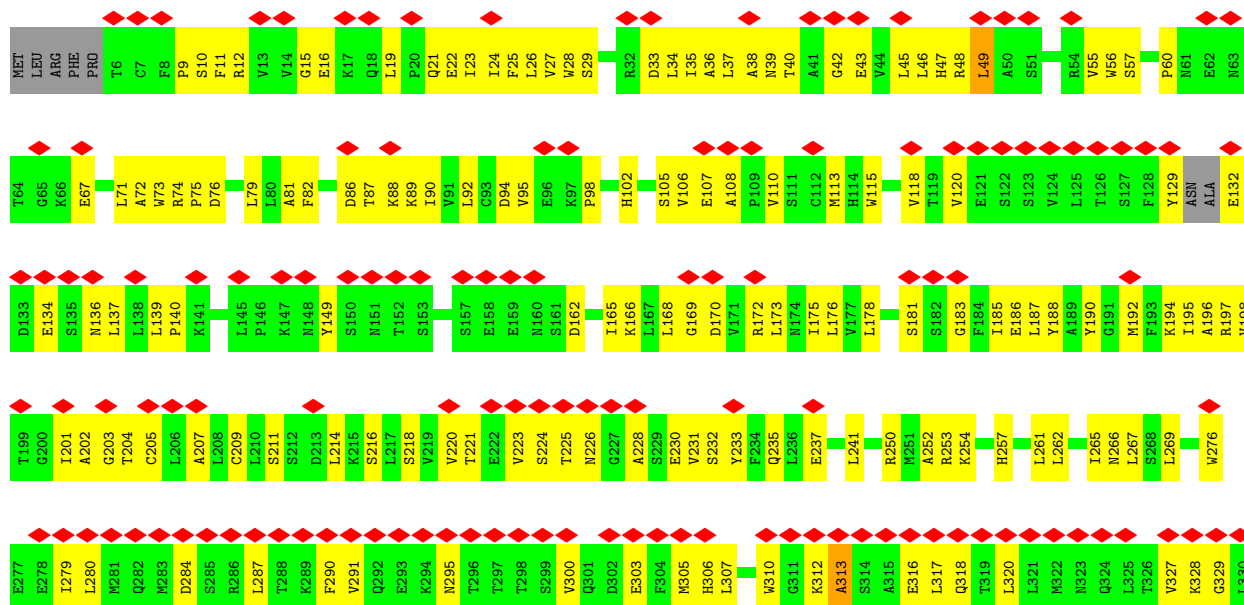
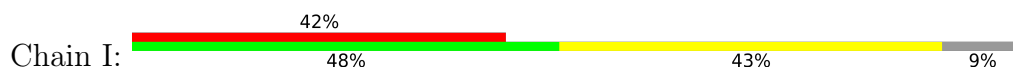
• Molecule 7: Anaphase-promoting complex subunit CDC26

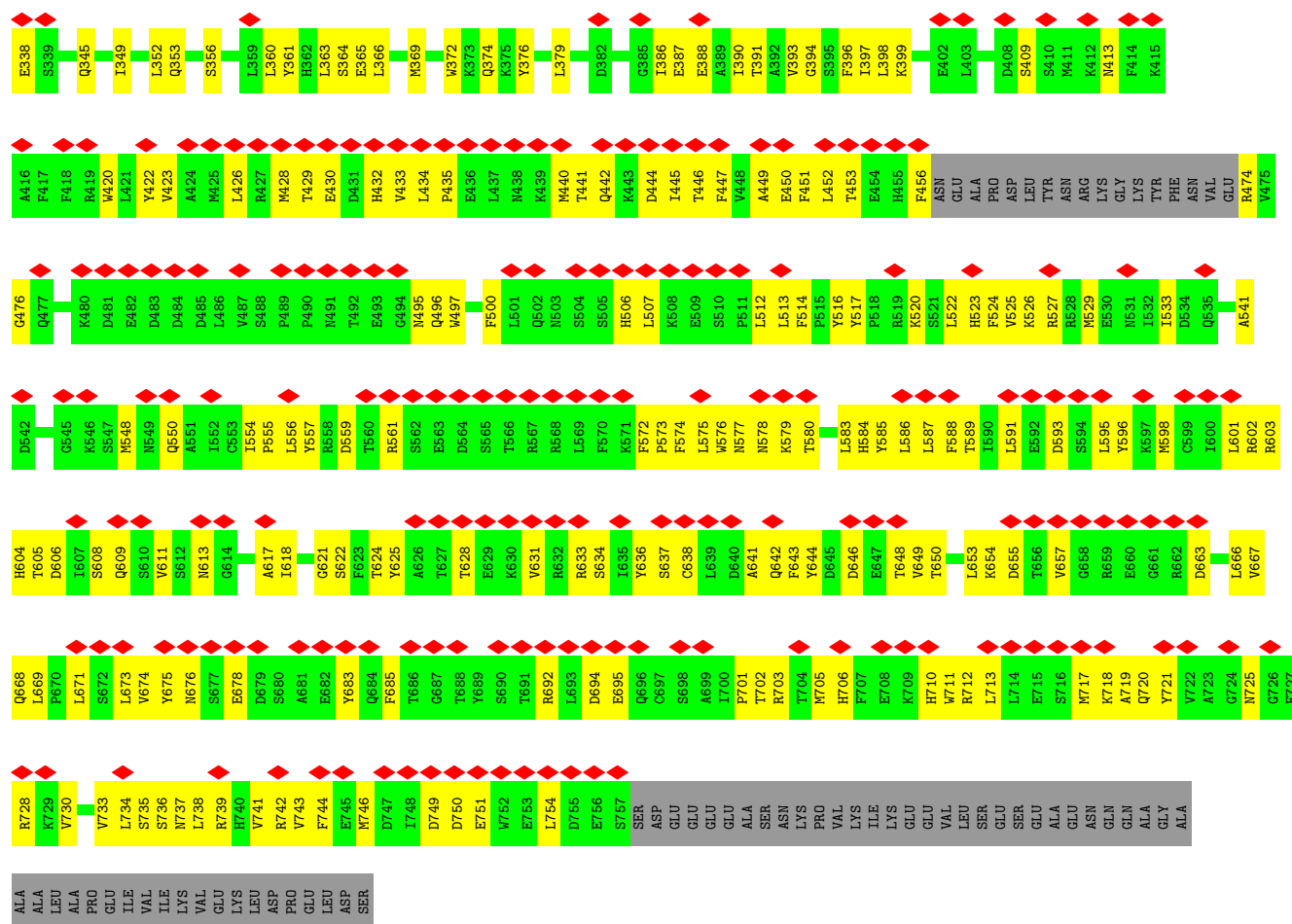


• Molecule 7: Anaphase-promoting complex subunit CDC26

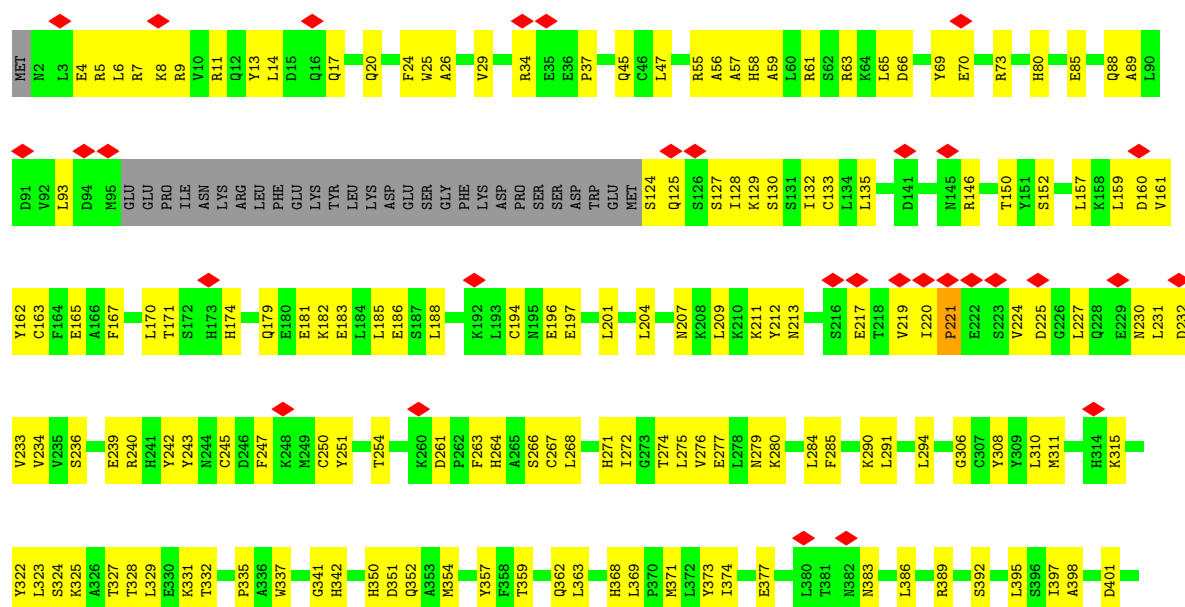


• Molecule 8: Anaphase-promoting complex subunit 4

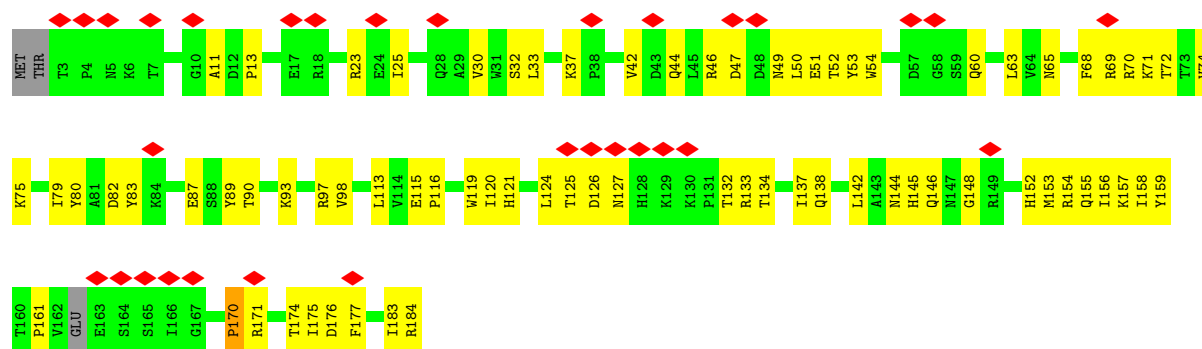




• Molecule 9: Cell division cycle protein 16 homolog







• Molecule 11: Anaphase-promoting complex subunit 13

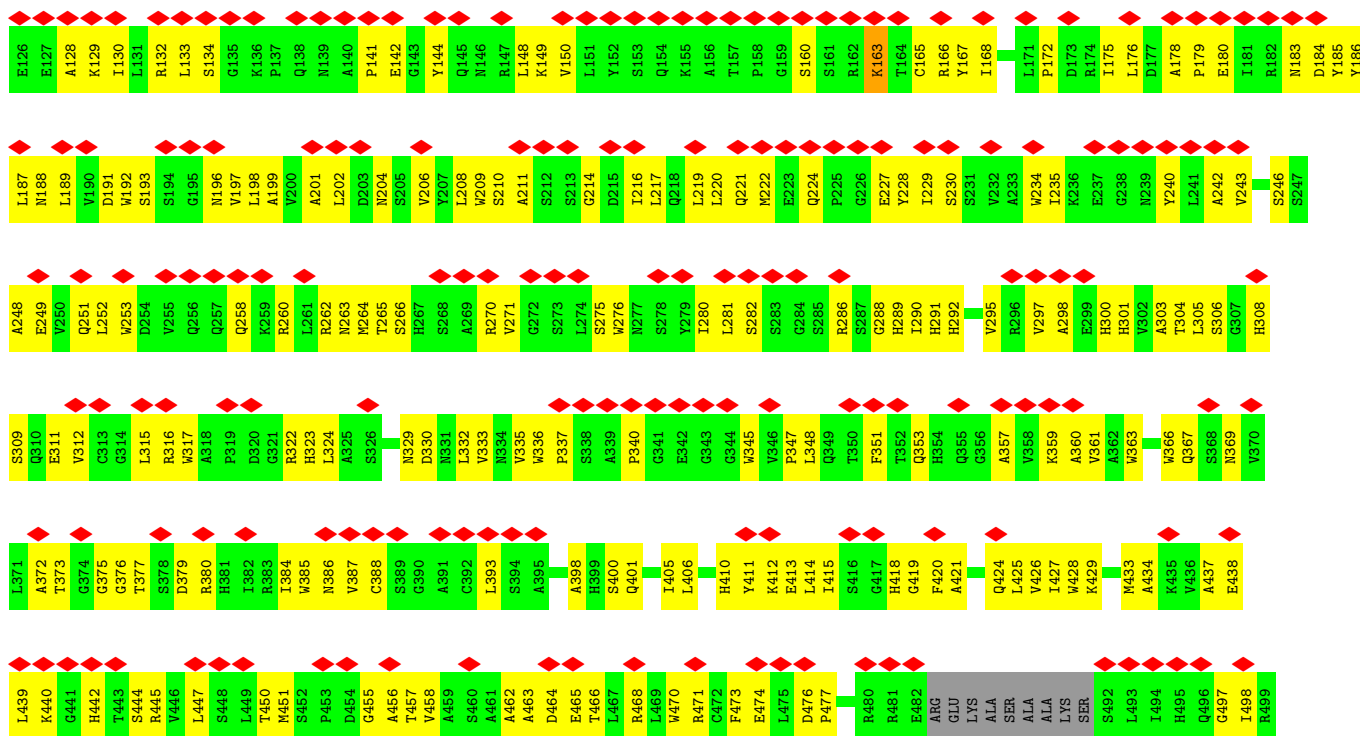


• Molecule 12: Anaphase-promoting complex subunit 2

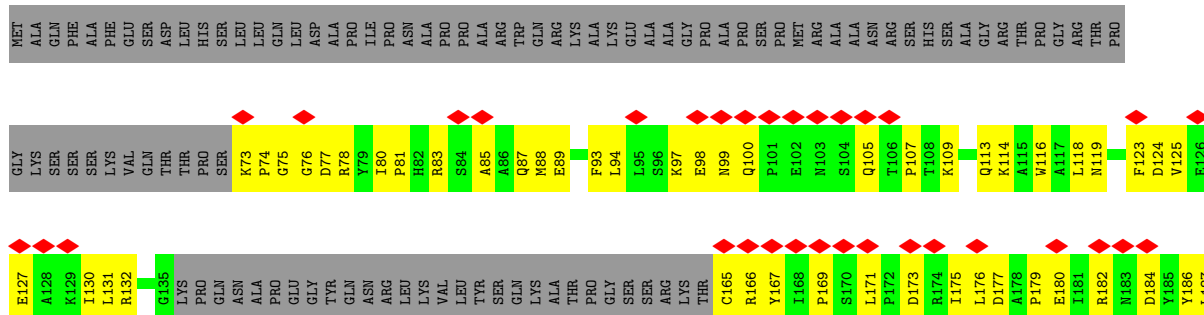


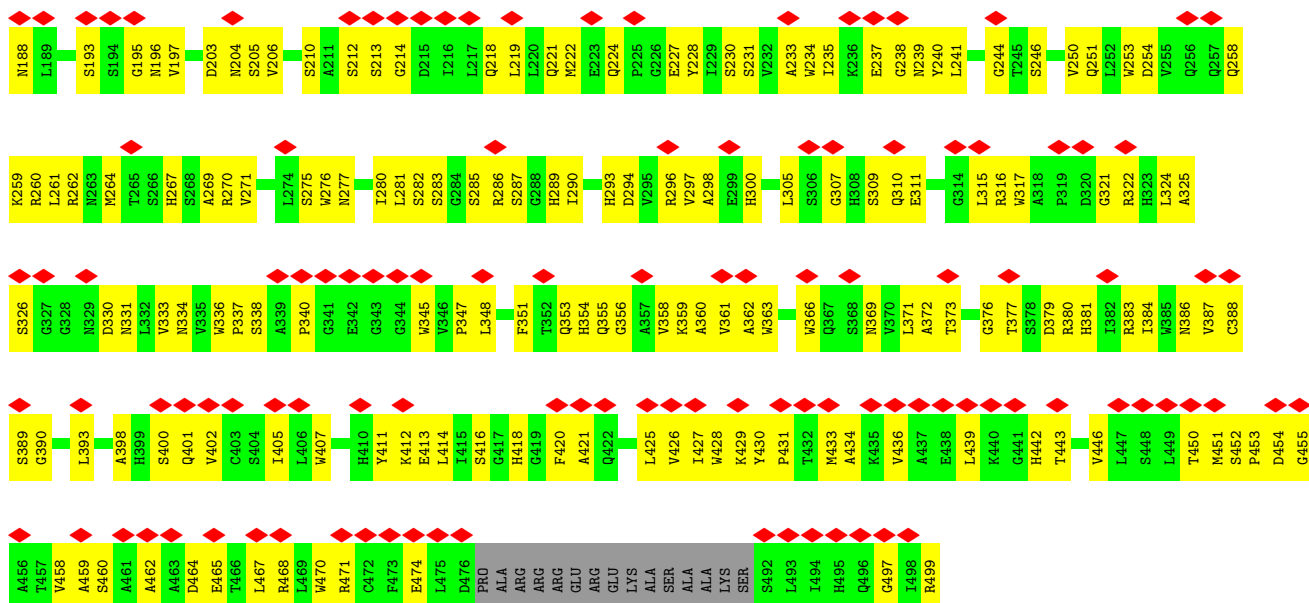


- Molecule 14: Cell division cycle protein 20 homolog

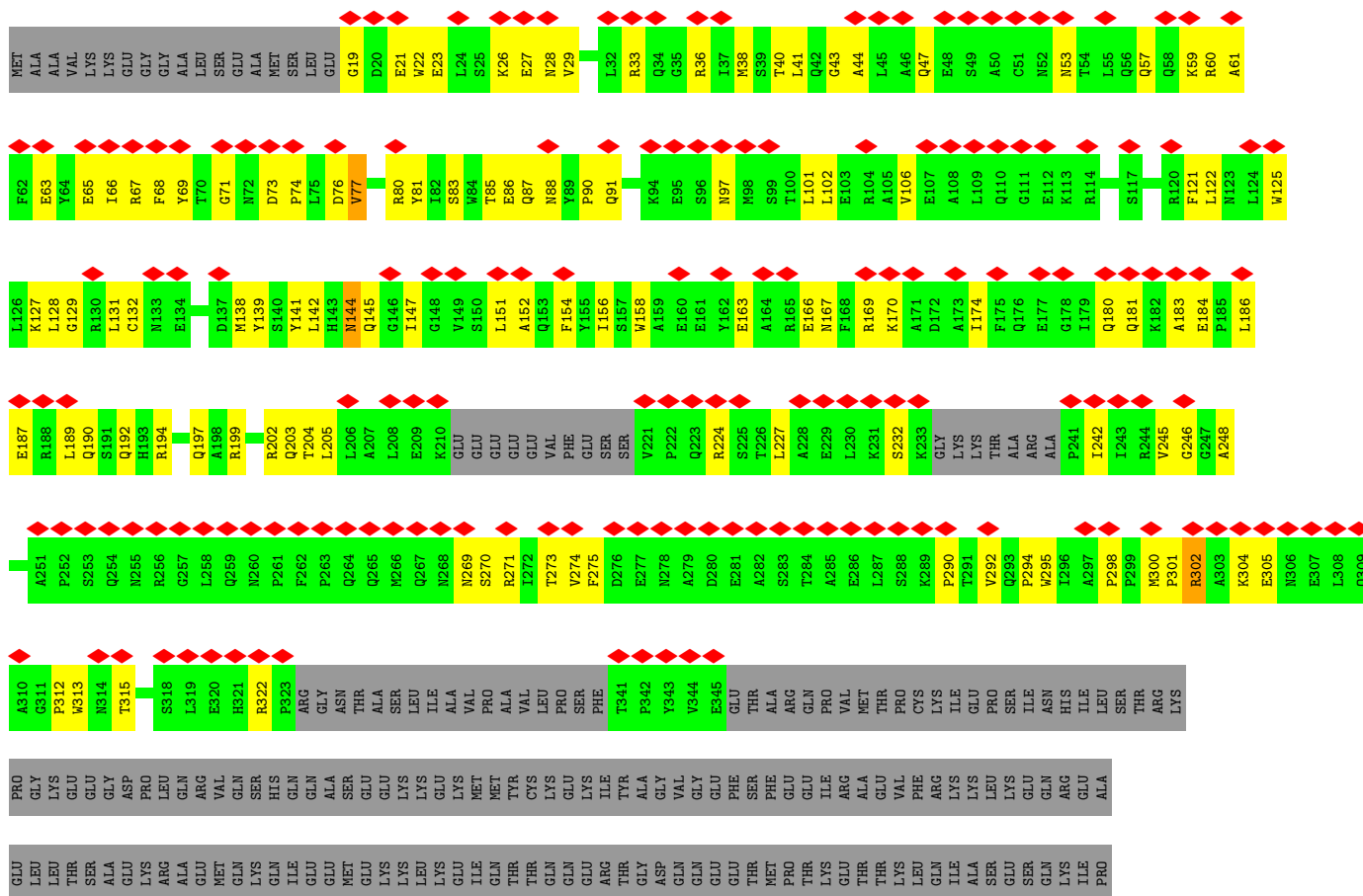


- Molecule 15: Cell division cycle protein 20 homolog



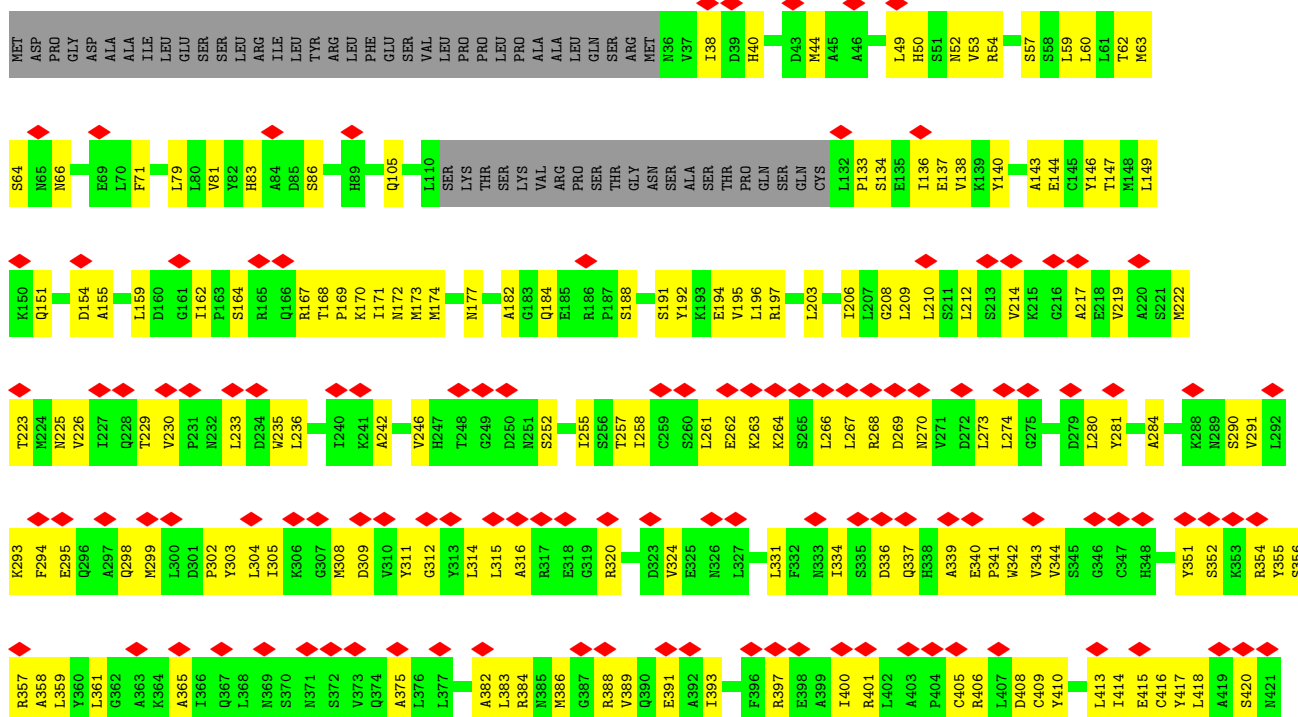


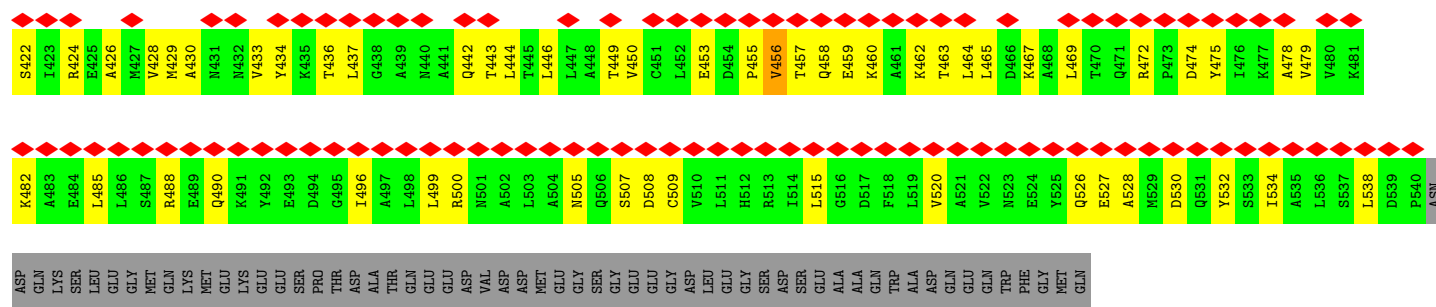
• Molecule 16: Mitotic checkpoint serine/threonine-protein kinase BUB1 beta



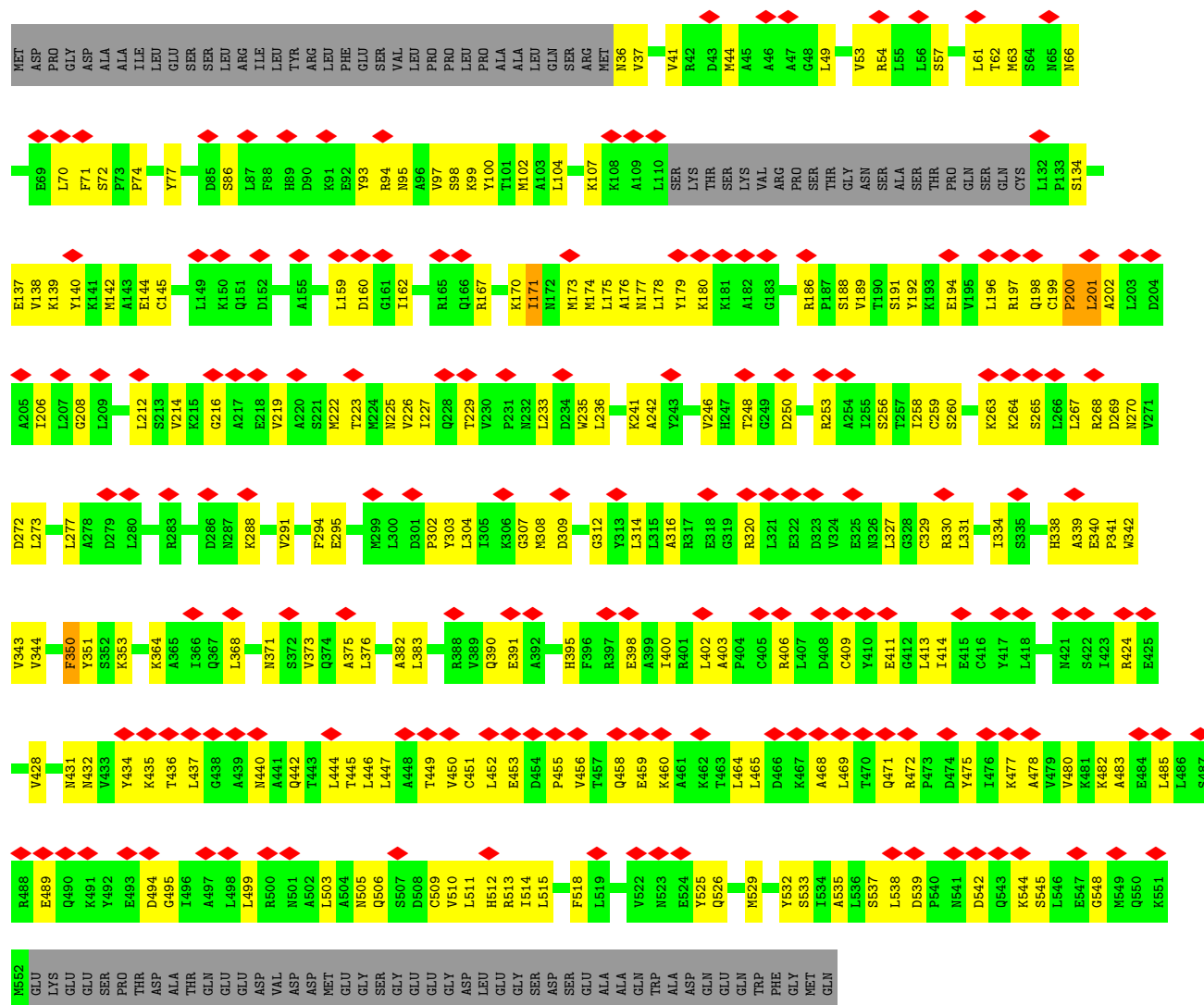
[illegible]

- Molecule 17: Anaphase-promoting complex subunit 7



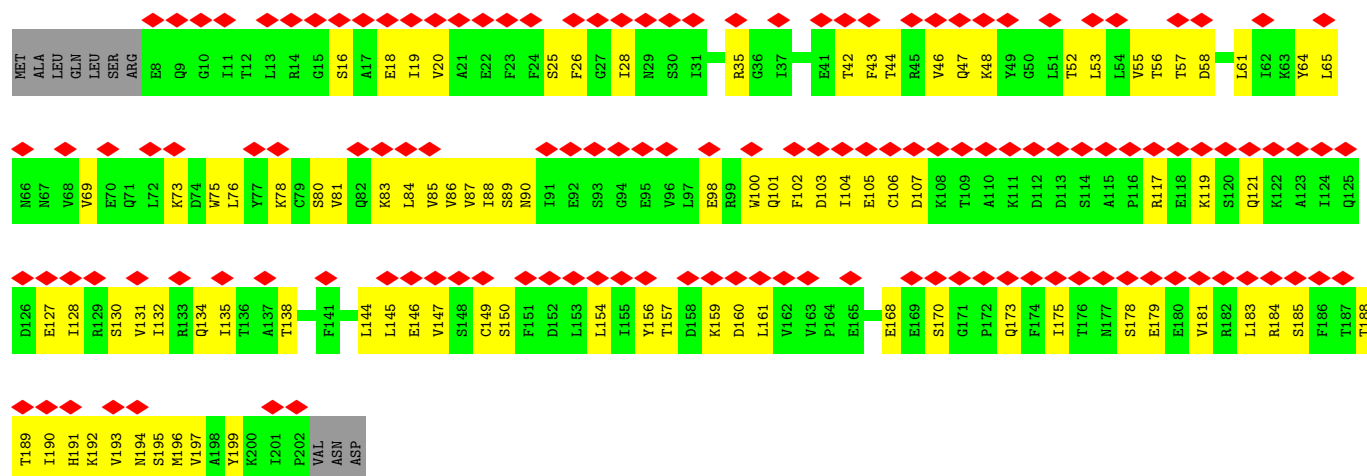


• Molecule 17: Anaphase-promoting complex subunit 7



• Molecule 18: Mitotic spindle assembly checkpoint protein MAD2A





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	343551	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	27	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.206	Depositor
Minimum map value	-0.097	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	359.04, 359.04, 359.04	wwPDB
Map dimensions	264, 264, 264	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.36, 1.36, 1.36	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	A	0.30	1/11191 (0.0%)	0.55	2/15239 (0.0%)
2	B	0.22	0/665	0.56	0/896
3	C	0.29	0/4404	0.50	2/5945 (0.0%)
3	P	0.27	0/4138	0.48	0/5587
4	D	0.26	0/159	0.50	0/218
5	E	0.28	0/459	0.45	0/619
6	F	0.29	0/3939	0.46	1/5325 (0.0%)
6	H	0.32	0/3943	0.47	0/5329
7	G	0.26	0/214	0.55	0/284
7	W	0.27	0/214	0.55	0/284
8	I	0.22	0/5849	0.49	0/7932
9	J	0.31	0/4146	0.51	0/5616
9	K	0.29	0/4086	0.49	0/5534
10	L	0.31	0/1468	0.54	0/1993
11	M	0.31	0/502	0.61	0/680
12	N	0.24	0/5496	0.67	9/7444 (0.1%)
13	O	0.26	0/5501	0.51	2/7432 (0.0%)
14	Q	0.22	0/2807	0.51	0/3830
15	R	0.25	0/3029	0.51	0/4124
16	S	0.23	0/2272	0.55	1/3088 (0.0%)
17	X	0.22	0/3833	0.47	0/5187
17	Y	0.22	0/3928	0.48	0/5311
18	Z	0.17	0/1605	0.42	0/2176
All	All	0.27	1/73848 (0.0%)	0.52	17/100073 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
2	B	0	2

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Mol	Chain	#Chirality outliers	#Planarity outliers
8	I	0	2
9	J	0	2
9	K	0	3
12	N	0	9
13	O	0	1
14	Q	0	1
16	S	0	1
All	All	0	25

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1278	GLY	C-N	-5.98	1.25	1.33

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	N	125	TYR	CA-C-N	10.42	140.45	121.70
12	N	125	TYR	C-N-CA	10.42	140.45	121.70
12	N	63	ALA	N-CA-C	8.32	128.51	110.80
12	N	508	ILE	CA-C-N	-6.68	108.78	121.54
12	N	508	ILE	C-N-CA	-6.68	108.78	121.54

There are no chirality outliers.

5 of 25 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1356	ASP	Peptide
1	A	274	VAL	Peptide
1	A	856	GLY	Peptide
1	A	859	PRO	Peptide
2	B	14	TRP	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	10949	0	10691	486	0
2	B	643	0	617	59	0
3	C	4306	0	4275	189	0
3	P	4043	0	4000	175	0
4	D	153	0	148	6	0
5	E	450	0	435	19	0
6	F	3849	0	3783	124	0
6	H	3853	0	3794	161	0
7	G	213	0	220	14	0
7	W	213	0	220	19	0
8	I	5728	0	5615	295	0
9	J	4047	0	3949	186	0
9	K	3988	0	3908	164	0
10	L	1435	0	1382	67	0
11	M	493	0	469	33	0
12	N	5403	0	5104	261	0
13	O	5402	0	5436	258	0
14	Q	2739	0	2582	186	0
15	R	2953	0	2839	192	0
16	S	2227	0	2028	114	0
17	X	3773	0	3831	181	0
17	Y	3868	0	3926	188	0
18	Z	1577	0	1593	74	0
All	All	72305	0	70845	3110	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 22.

The worst 5 of 3110 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)
16:S:290:PRO:HB2	16:S:292:VAL:HG23	1.37	1.01
14:Q:141:PRO:HB2	16:S:202:ARG:HE	1.26	0.95
9:J:264:HIS:HE2	9:J:266:SER:HG	1.11	0.93
1:A:1134:TRP:HD1	1:A:1597:THR:HA	1.36	0.90
2:B:11:VAL:HG22	12:N:642:GLY:HA2	1.54	0.89

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	1399/1944 (72%)	1246 (89%)	147 (10%)	6 (0%)	30	62
2	B	75/84 (89%)	64 (85%)	11 (15%)	0	100	100
3	C	520/597 (87%)	490 (94%)	30 (6%)	0	100	100
3	P	485/597 (81%)	455 (94%)	30 (6%)	0	100	100
4	D	16/121 (13%)	13 (81%)	3 (19%)	0	100	100
5	E	54/110 (49%)	52 (96%)	2 (4%)	0	100	100
6	F	479/824 (58%)	448 (94%)	31 (6%)	0	100	100
6	H	479/824 (58%)	443 (92%)	34 (7%)	2 (0%)	30	62
7	G	23/85 (27%)	23 (100%)	0	0	100	100
7	W	23/85 (27%)	23 (100%)	0	0	100	100
8	I	727/808 (90%)	674 (93%)	53 (7%)	0	100	100
9	J	500/620 (81%)	454 (91%)	45 (9%)	1 (0%)	43	73
9	K	489/620 (79%)	445 (91%)	44 (9%)	0	100	100
10	L	180/185 (97%)	159 (88%)	19 (11%)	2 (1%)	11	41
11	M	55/74 (74%)	43 (78%)	12 (22%)	0	100	100
12	N	681/822 (83%)	575 (84%)	104 (15%)	2 (0%)	36	67
13	O	677/755 (90%)	639 (94%)	37 (6%)	1 (0%)	48	79
14	Q	361/374 (96%)	327 (91%)	32 (9%)	2 (1%)	21	54
15	R	377/499 (76%)	338 (90%)	39 (10%)	0	100	100
16	S	285/1050 (27%)	240 (84%)	45 (16%)	0	100	100
17	X	480/599 (80%)	456 (95%)	23 (5%)	1 (0%)	43	73
17	Y	492/599 (82%)	464 (94%)	26 (5%)	2 (0%)	30	62
18	Z	193/205 (94%)	187 (97%)	6 (3%)	0	100	100
All	All	9050/12481 (72%)	8258 (91%)	773 (8%)	19 (0%)	44	73

5 of 19 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1099	PRO
6	H	147	PHE
10	L	175	ILE
13	O	130	SER
17	X	456	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	1151/1720 (67%)	1148 (100%)	3 (0%)	86	84
2	B	71/75 (95%)	70 (99%)	1 (1%)	59	70
3	C	452/520 (87%)	452 (100%)	0	100	100
3	P	422/520 (81%)	421 (100%)	1 (0%)	87	87
4	D	18/115 (16%)	18 (100%)	0	100	100
5	E	47/89 (53%)	47 (100%)	0	100	100
6	F	407/727 (56%)	407 (100%)	0	100	100
6	H	408/727 (56%)	408 (100%)	0	100	100
7	G	23/77 (30%)	23 (100%)	0	100	100
7	W	23/77 (30%)	23 (100%)	0	100	100
8	I	621/730 (85%)	620 (100%)	1 (0%)	87	87
9	J	424/548 (77%)	424 (100%)	0	100	100
9	K	423/548 (77%)	423 (100%)	0	100	100
10	L	155/170 (91%)	155 (100%)	0	100	100
11	M	55/67 (82%)	54 (98%)	1 (2%)	51	67
12	N	518/724 (72%)	516 (100%)	2 (0%)	84	83
13	O	577/650 (89%)	576 (100%)	1 (0%)	87	87
14	Q	275/310 (89%)	275 (100%)	0	100	100
15	R	311/411 (76%)	311 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
16	S	208/938 (22%)	206 (99%)	2 (1%)	68	74
17	X	407/513 (79%)	407 (100%)	0	100	100
17	Y	418/513 (82%)	416 (100%)	2 (0%)	81	81
18	Z	181/190 (95%)	181 (100%)	0	100	100
All	All	7595/10959 (69%)	7581 (100%)	14 (0%)	85	87

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
12	N	368	THR
13	O	420	ILE
17	Y	350	PHE
16	S	302	ARG
17	Y	171	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 198 such sidechains are listed below:

Mol	Chain	Res	Type
12	N	726	ASN
3	P	287	ASN
13	O	69	GLN
13	O	500	ASN
3	P	477	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
12	N	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	N	92:TRP	C	93:ASN	N	2.97

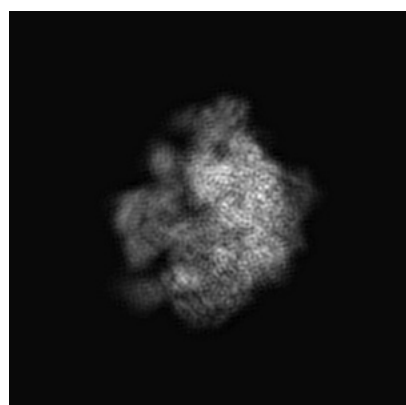
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-10516. These allow visual inspection of the internal detail of the map and identification of artifacts.

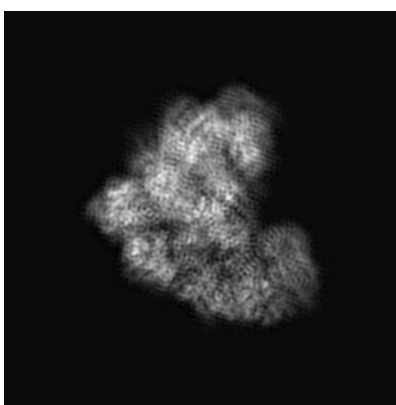
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

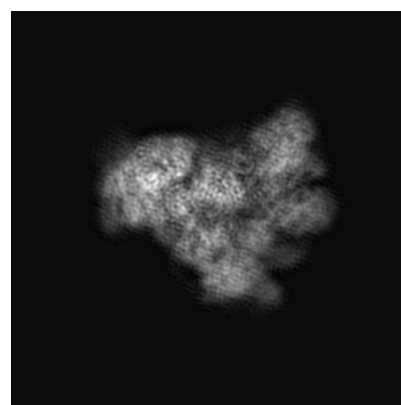
6.1.1 Primary map



X



Y

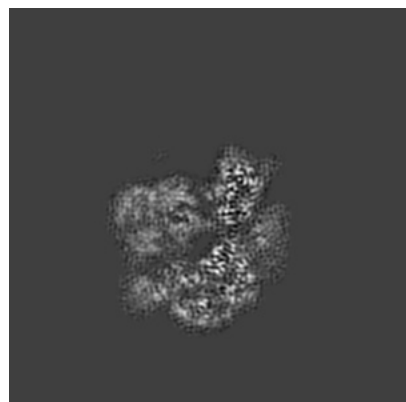


Z

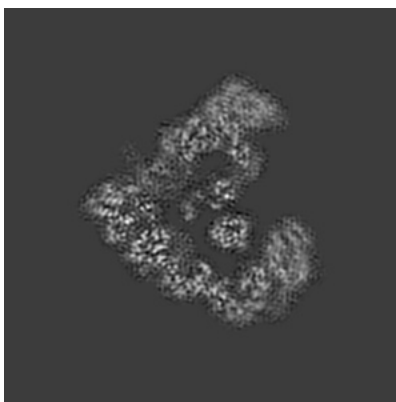
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

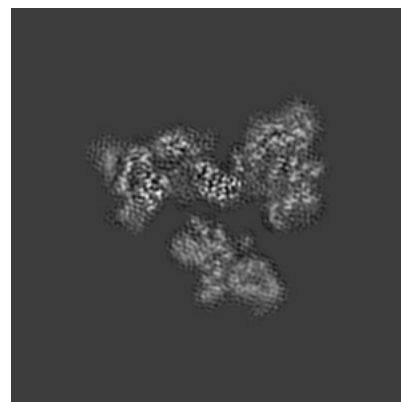
6.2.1 Primary map



X Index: 132



Y Index: 132

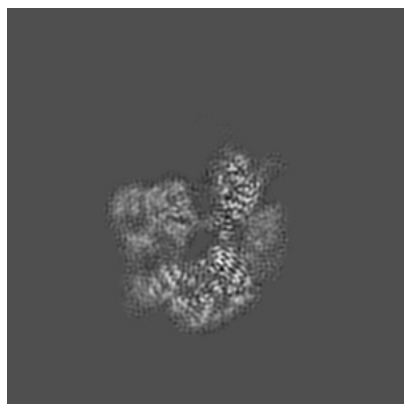


Z Index: 132

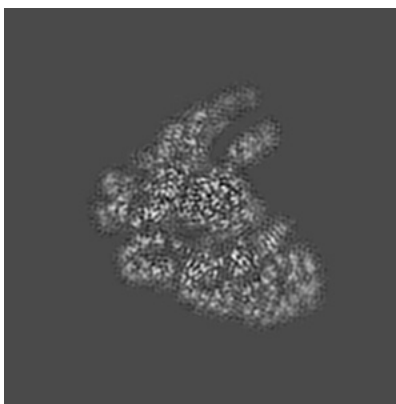
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

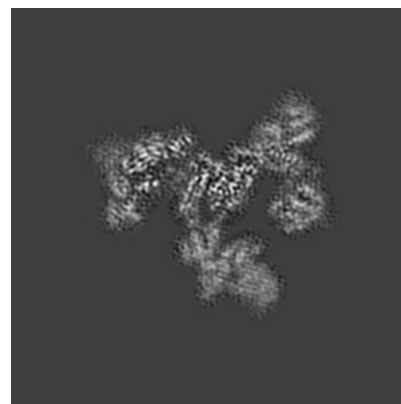
6.3.1 Primary map



X Index: 129



Y Index: 147

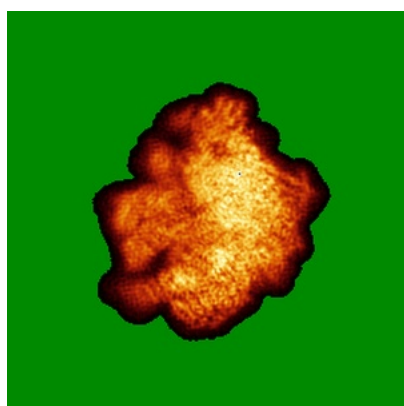


Z Index: 139

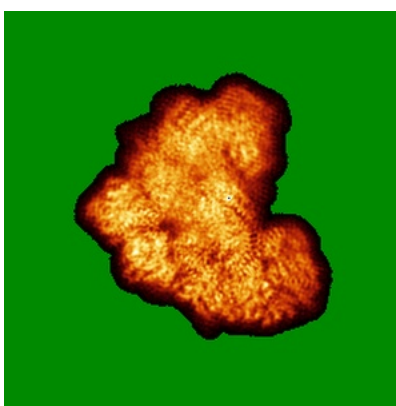
The images above show the largest variance slices of the map in three orthogonal directions.

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

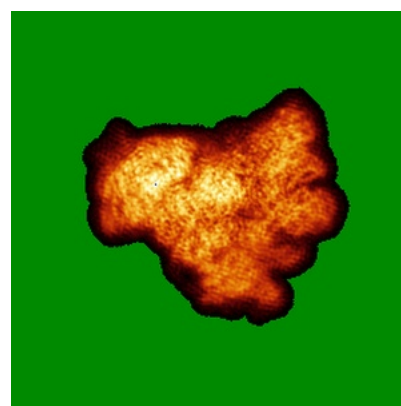
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.06. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

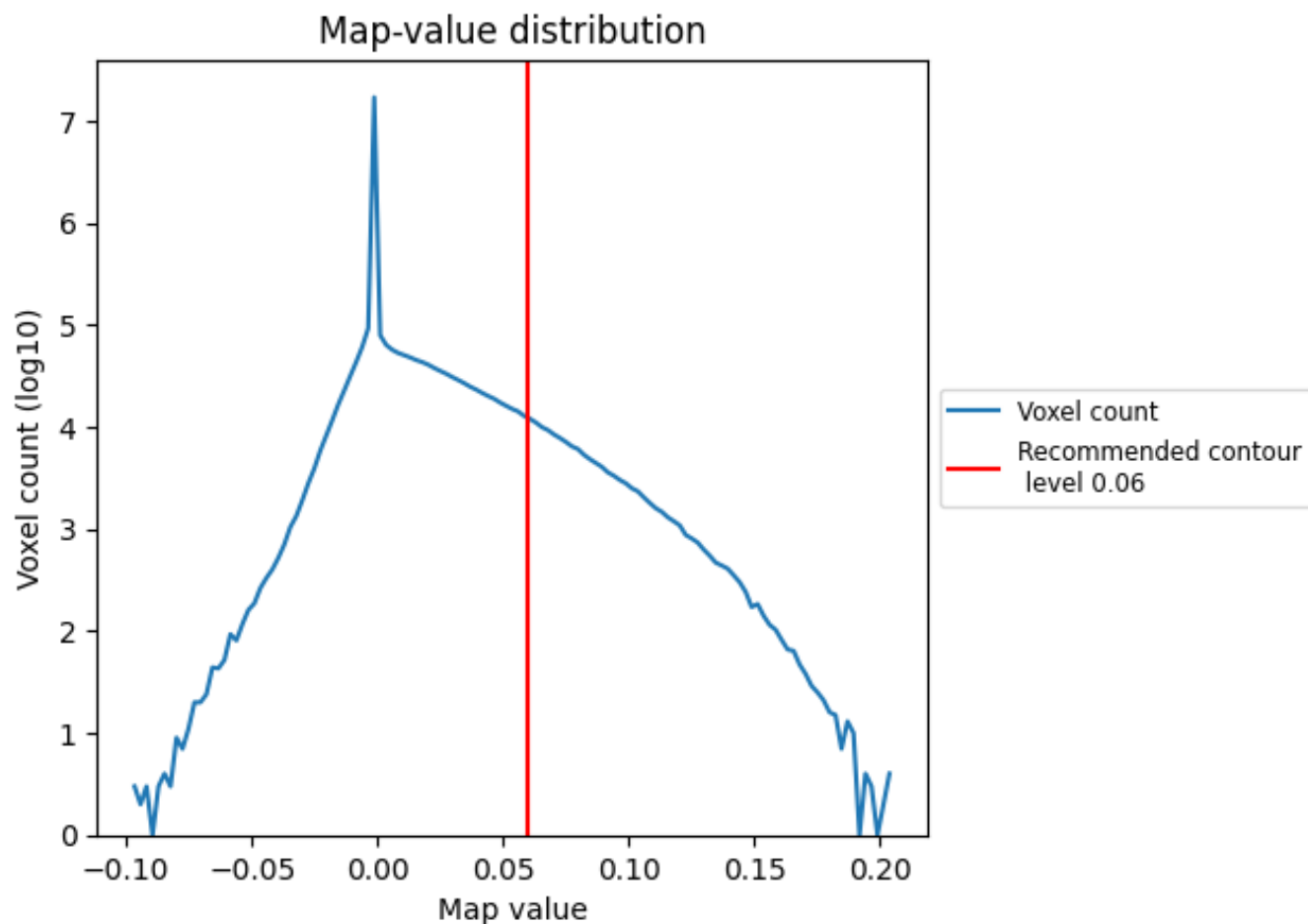
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

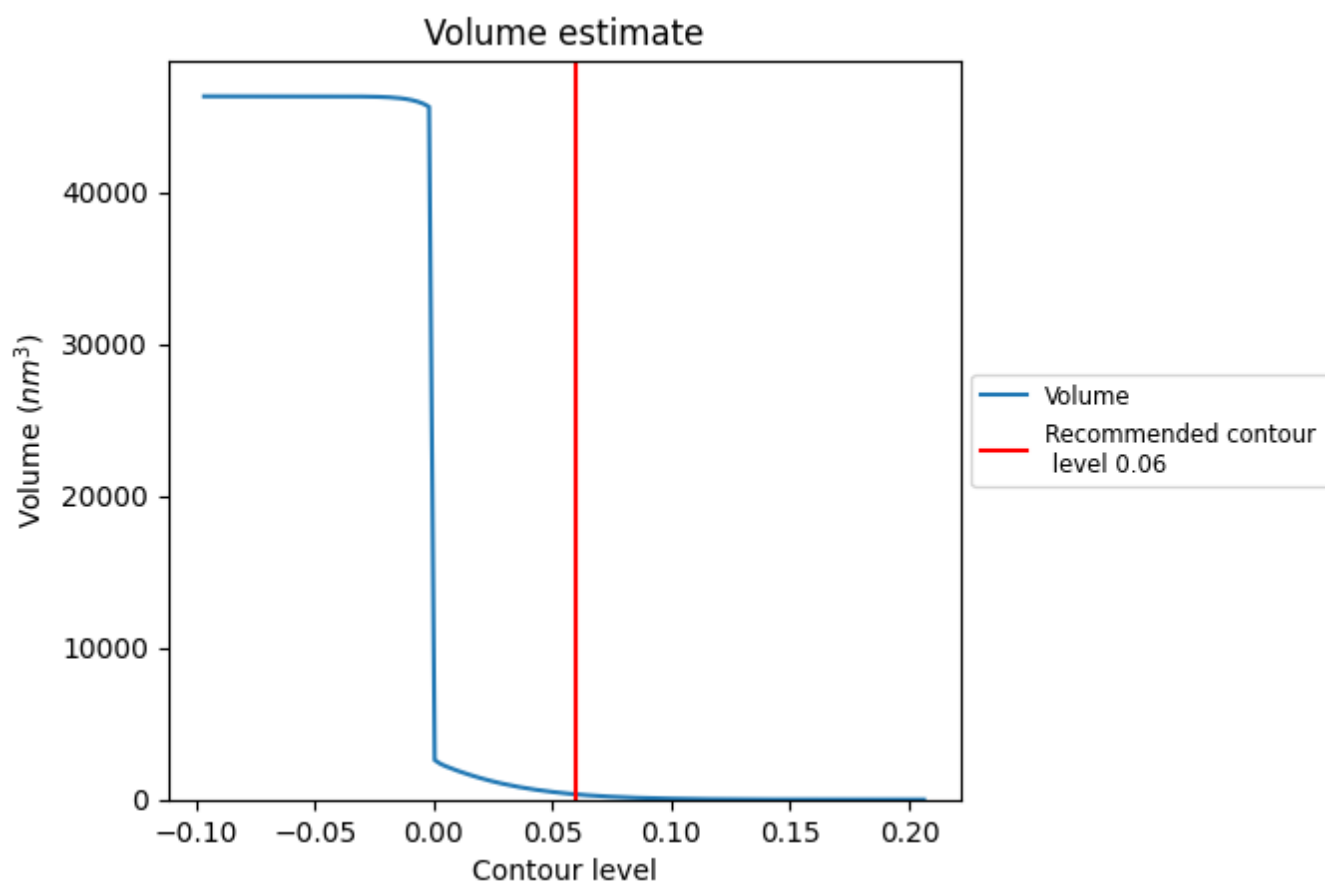
This section contains the results of statistical analysis of the map.

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.

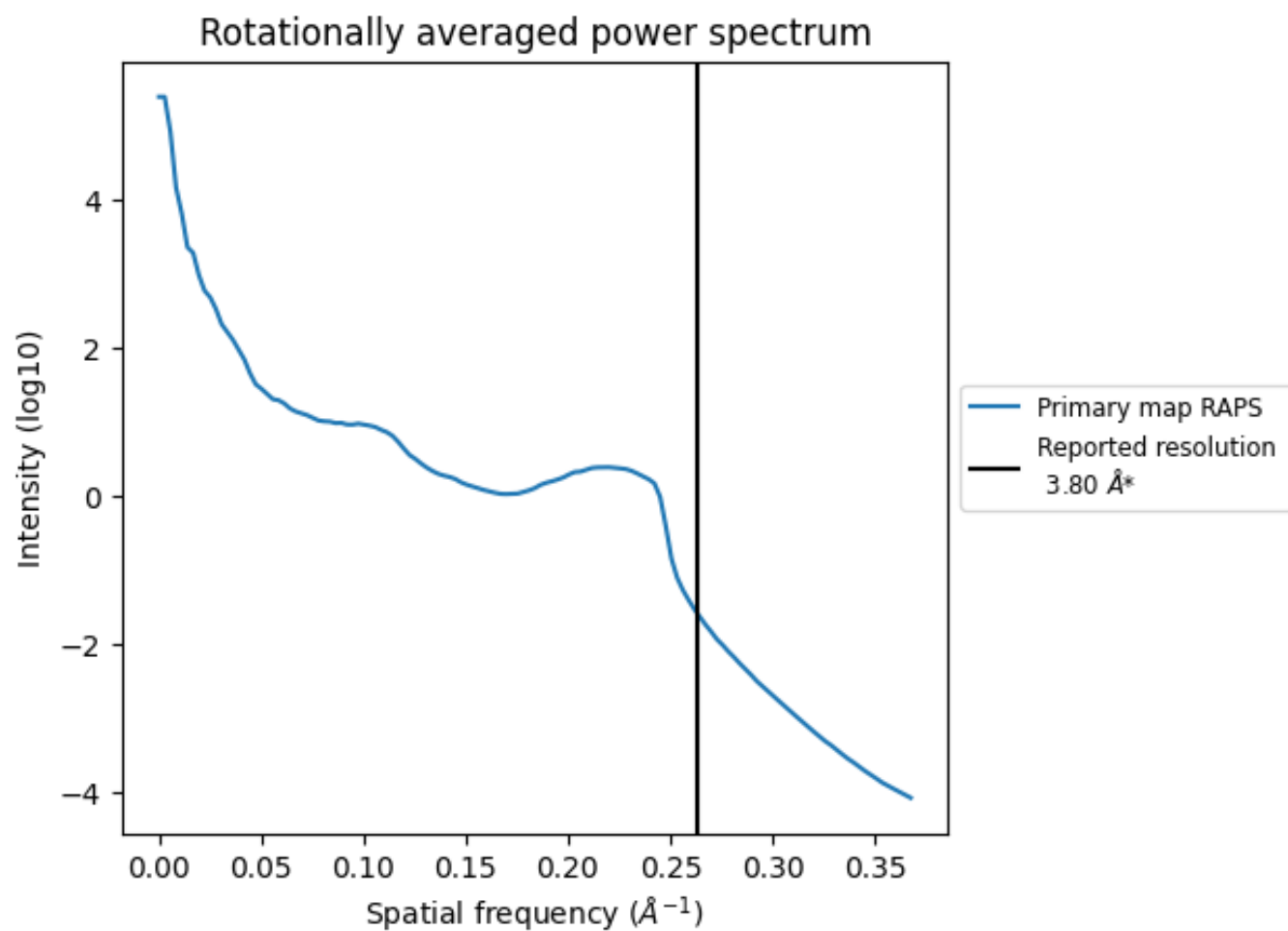
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 343 nm³; this corresponds to an approximate mass of 310 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.263 Å⁻¹

8 Fourier-Shell correlation ⓘ

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

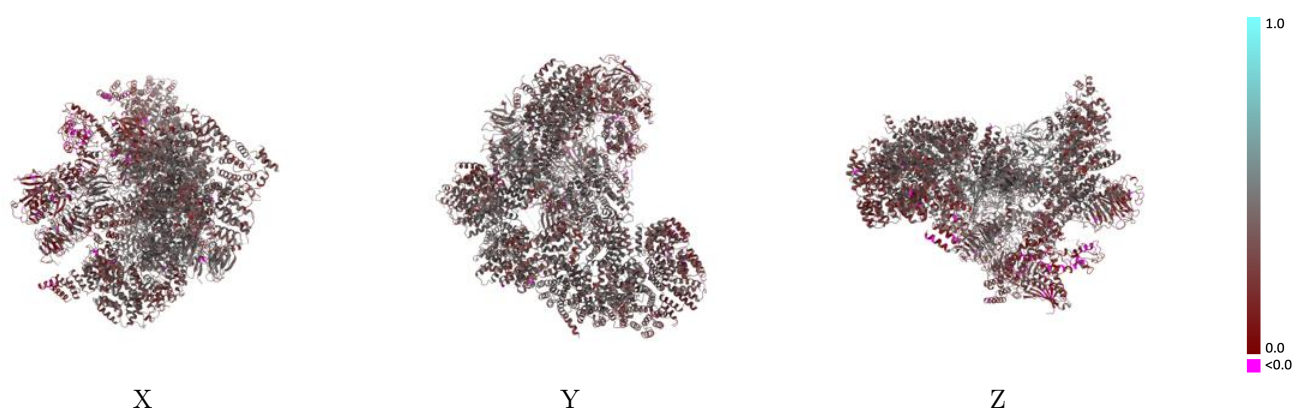
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-10516 and PDB model 6TLJ. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)

This section was not generated.

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

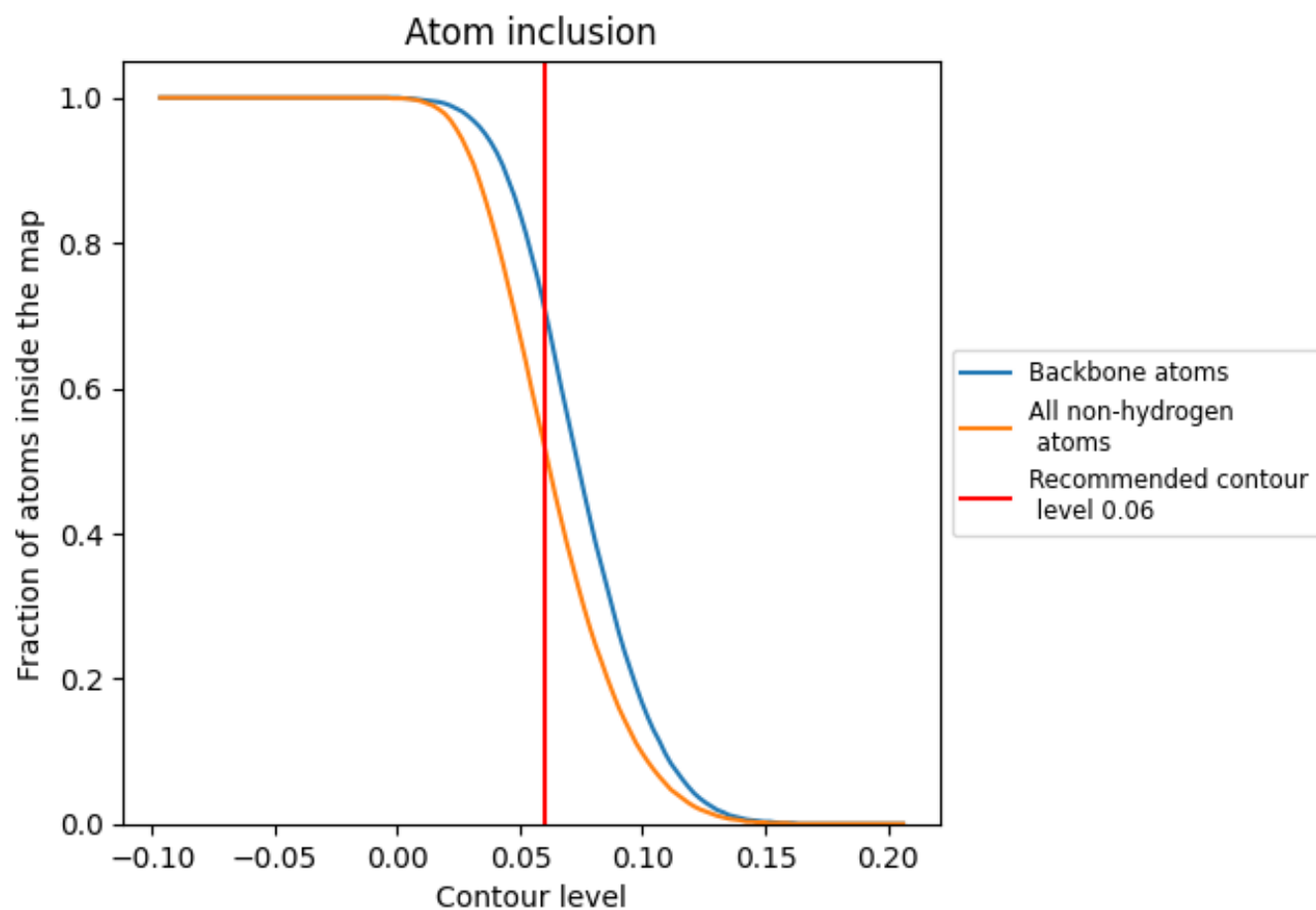


The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 52% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5200	 0.3570
A	 0.5550	 0.4080
B	 0.0250	 0.1810
C	 0.5580	 0.3750
D	 0.4600	 0.4830
E	 0.6330	 0.4370
F	 0.6350	 0.4030
G	 0.5270	 0.4060
H	 0.6880	 0.4100
I	 0.4270	 0.3240
J	 0.6720	 0.4070
K	 0.6300	 0.3790
L	 0.6040	 0.4320
M	 0.5030	 0.4120
N	 0.3890	 0.2730
O	 0.4950	 0.3820
P	 0.5830	 0.3540
Q	 0.4320	 0.2950
R	 0.4840	 0.3450
S	 0.3730	 0.2790
W	 0.5410	 0.4030
X	 0.4380	 0.3000
Y	 0.4940	 0.3270
Z	 0.2870	 0.2380

