



wwPDB EM Validation Summary Report ⓘ

Mar 5, 2026 – 05:30 PM UTC

PDB ID : 6WGI / pdb_00006wgi
EMDB ID : EMD-21666
Title : Atomic model of the mutant OCCM (ORC-Cdc6-Cdt1-Mcm2-7 with Mcm6 WHD truncation) loaded on DNA at 10.5 Å resolution
Authors : Yuan, Z.; Schneider, S.; Dodd, T.; Riera, A.; Bai, L.; Yan, C.; Magdalou, I.; Ivanov, I.; Stillman, B.; Li, H.; Speck, C.
Deposited on : 2020-04-05
Resolution : 10.00 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
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

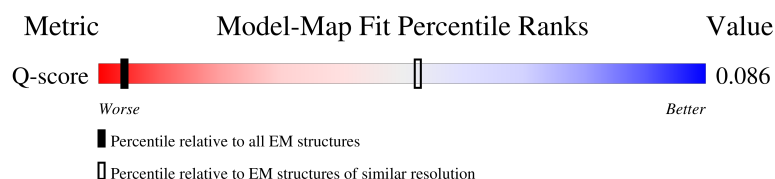
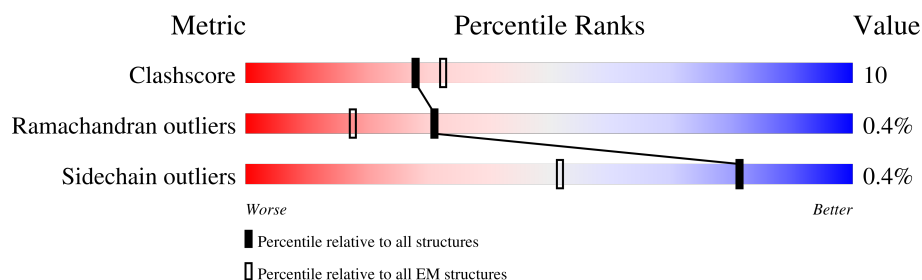
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 10.00 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	147 (9.50 - 10.50)

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

Mol	Chain	Length	Quality of chain
1	G	34	
2	H	34	
3	9	513	
4	A	913	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
5	B	620	
6	C	616	
7	E	479	
8	D	529	
9	F	435	
10	2	868	
11	3	971	
12	4	933	
13	5	775	
14	L	604	
15	6	1017	
16	7	845	

2 Entry composition [i](#)

There are 18 unique types of molecules in this entry. The entry contains 52948 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 DNA chain called DNA (34-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
1	G	34	Total	C	N	O	P	0	0
			694	340	101	220	33		

- Molecule 2 is a DNA chain called DNA (34-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	34	Total	C	N	O	P	0	0
			697	334	143	186	34		

- Molecule 3 is a protein called Cell division control protein 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	9	373	Total	C	N	O	S	0	0
			2962	1901	493	551	17		

- Molecule 4 is a protein called Origin recognition complex subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	A	420	Total	C	N	O	S	0	0
			3323	2125	556	624	18		

- Molecule 5 is a protein called Origin recognition complex subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	B	325	Total	C	N	O	S	0	0
			2640	1705	439	480	16		

- Molecule 6 is a protein called Origin recognition complex subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	C	544	Total	C	N	O	S	0	0
			4501	2906	742	838	15		

- Molecule 7 is a protein called Origin recognition complex subunit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	E	422	Total	C	N	O	S	0	0
			3418	2220	545	640	13		

- Molecule 8 is a protein called Origin recognition complex subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	D	436	Total	C	N	O	S	0	0
			3534	2262	601	658	13		

- Molecule 9 is a protein called Origin recognition complex subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	F	157	Total	C	N	O	S	0	0
			1315	846	222	235	12		

- Molecule 10 is a protein called DNA replication licensing factor MCM2.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	2	573	Total	C	N	O	S	0	0
			4274	2685	754	817	18		

- Molecule 11 is a protein called DNA replication licensing factor MCM3.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	3	637	Total	C	N	O	S	0	0
			4773	3025	819	917	12		

- Molecule 12 is a protein called DNA replication licensing factor MCM4.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	4	650	Total	C	N	O	S	0	0
			4950	3104	846	975	25		

- Molecule 13 is a protein called Minichromosome maintenance protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	5	529	Total	C	N	O	S	0	0
			3955	2484	680	772	19		

- Molecule 14 is a protein called Cell division cycle protein CDT1.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	L	367	Total	C	N	O	S	0	0
			2898	1860	487	538	13		

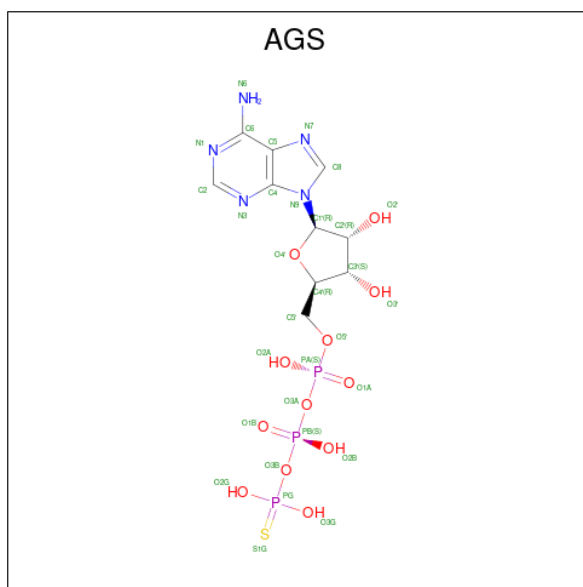
- Molecule 15 is a protein called DNA replication licensing factor MCM6.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	6	563	Total	C	N	O	S	0	0
			4187	2643	723	799	22		

- Molecule 16 is a protein called DNA replication licensing factor MCM7.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	7	625	Total	C	N	O	S	0	0
			4701	2966	798	908	29		

- Molecule 17 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (CCD ID: AGS) (formula: $C_{10}H_{16}N_5O_{12}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						AltConf
17	9	1	Total	C	N	O	P	S	0
			31	10	5	12	3	1	
17	A	1	Total	C	N	O	P	S	0
			31	10	5	12	3	1	
17	E	1	Total	C	N	O	P	S	0
			31	10	5	12	3	1	
17	D	1	Total	C	N	O	P	S	0
			31	10	5	12	3	1	

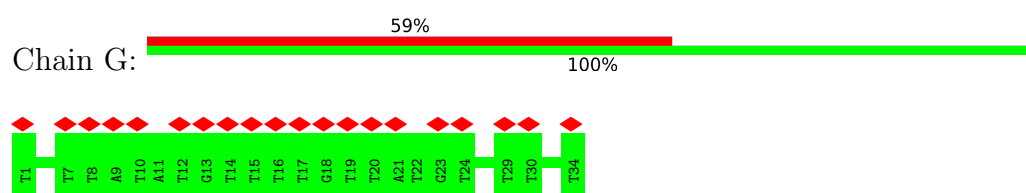
- Molecule 18 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

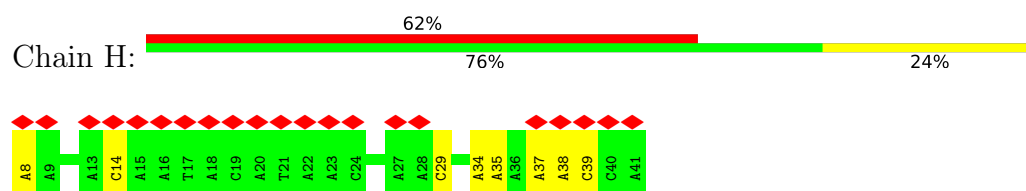
3 Residue-property plots

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.

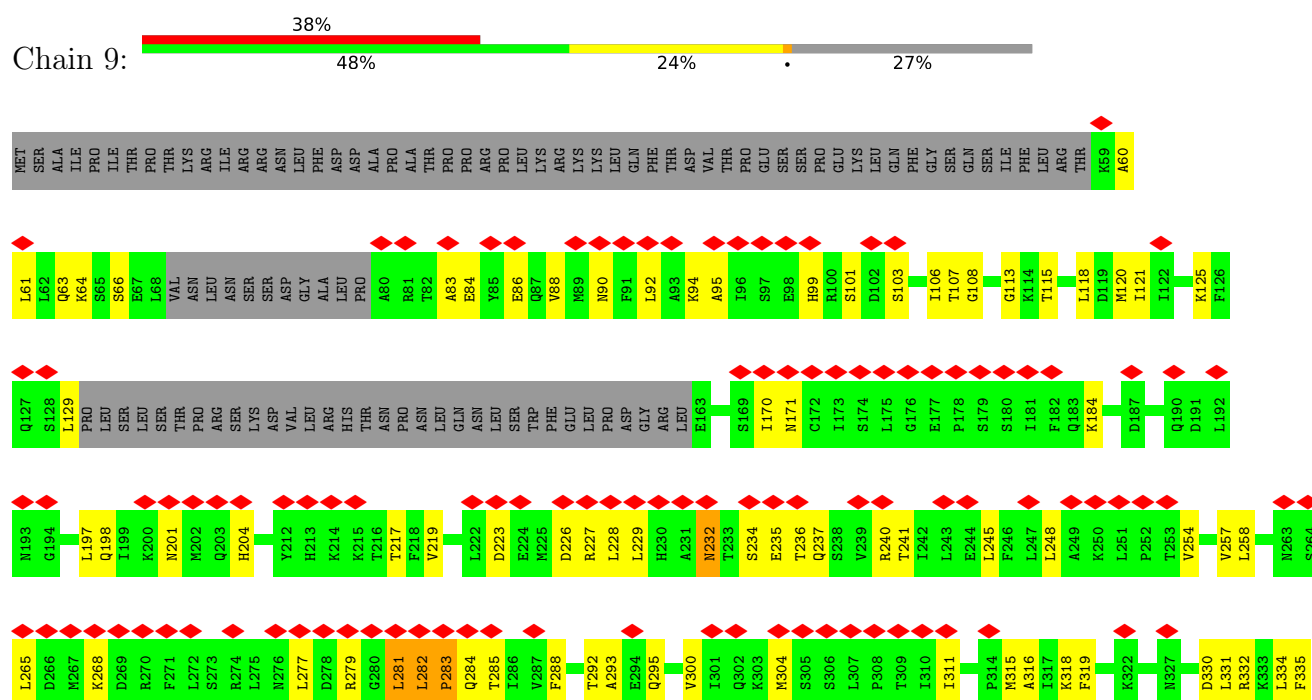
• Molecule 1: DNA (34-MER)

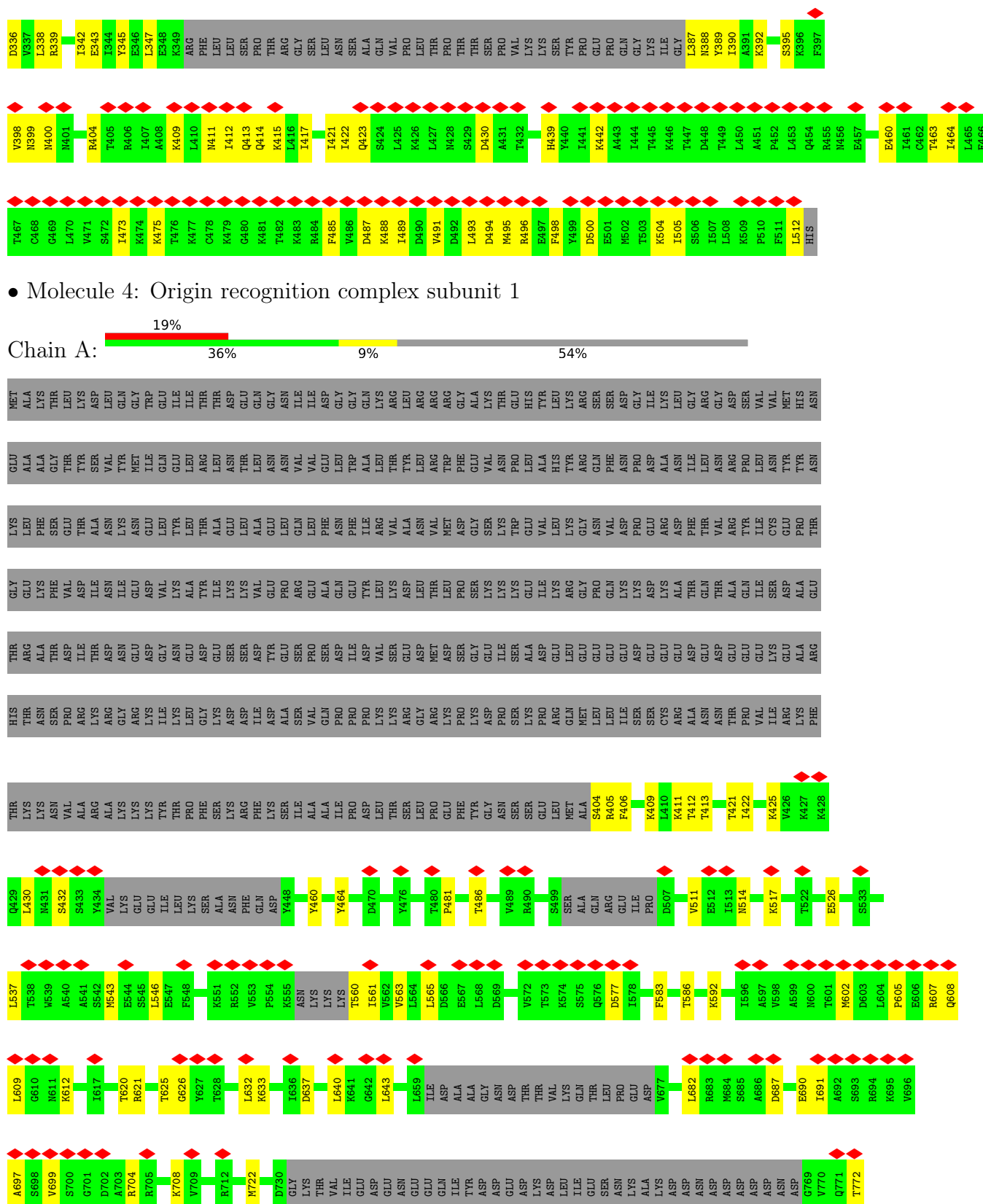


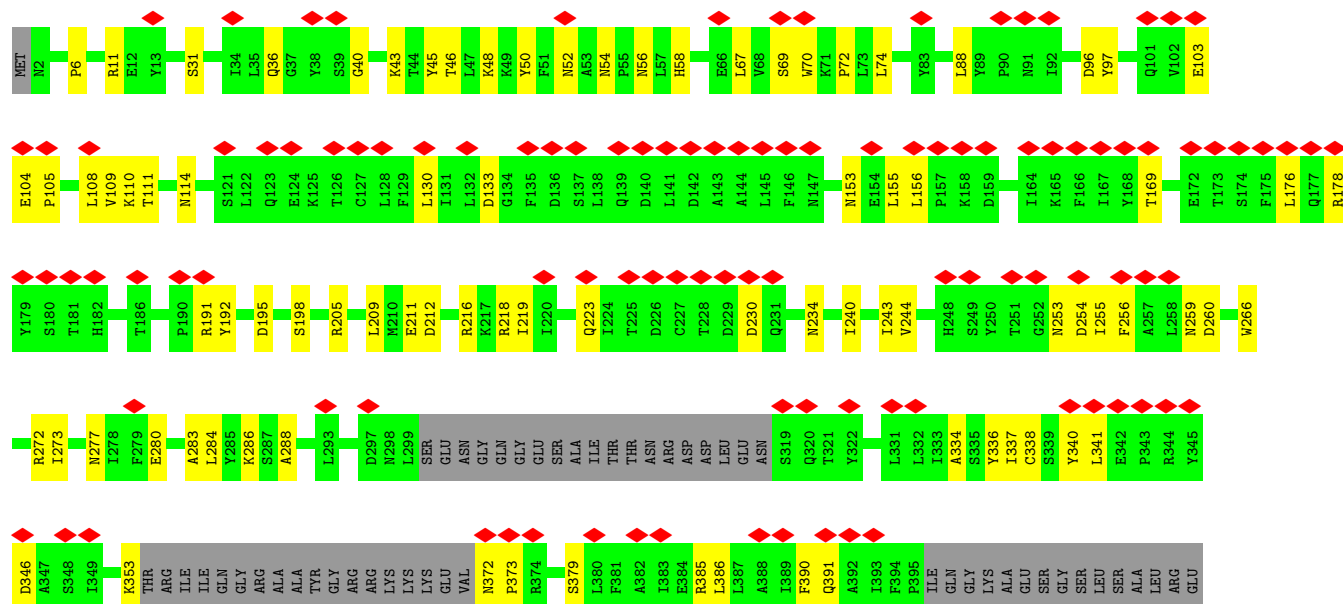
• Molecule 2: DNA (34-MER)



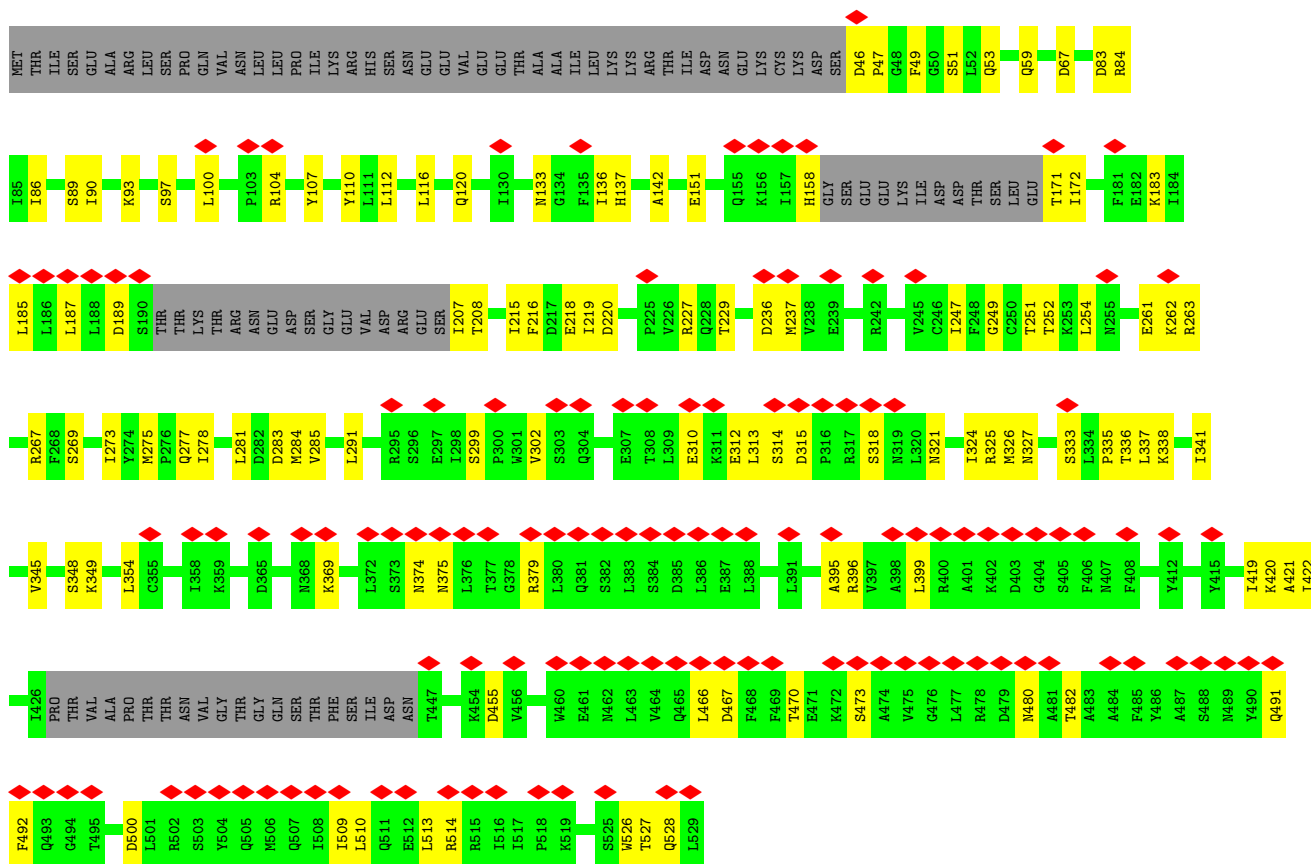
• Molecule 3: Cell division control protein 6



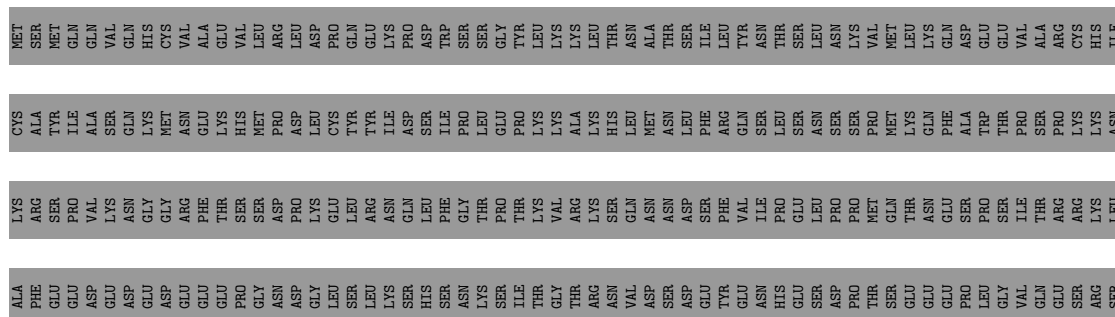




- Molecule 8: Origin recognition complex subunit 4

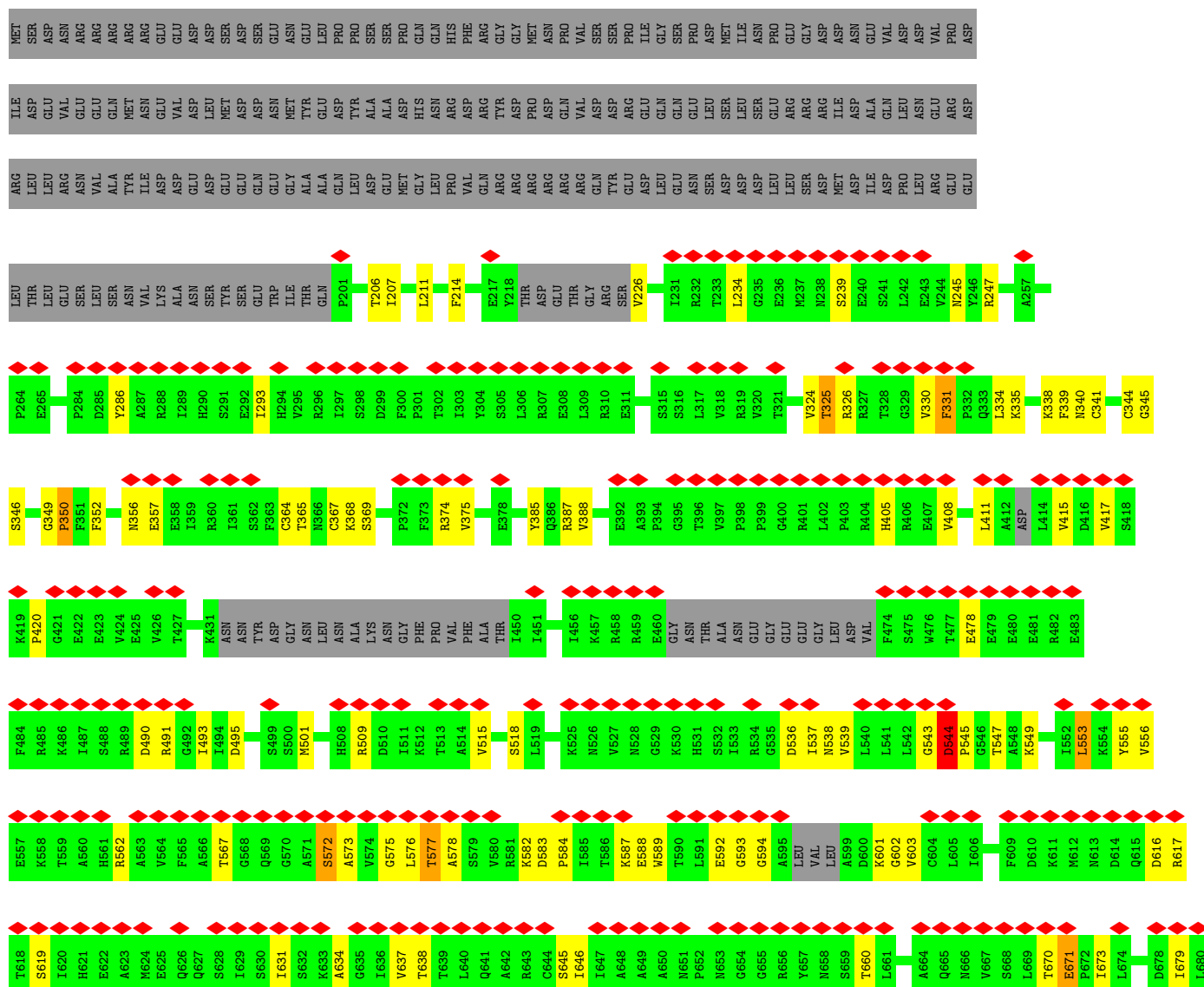


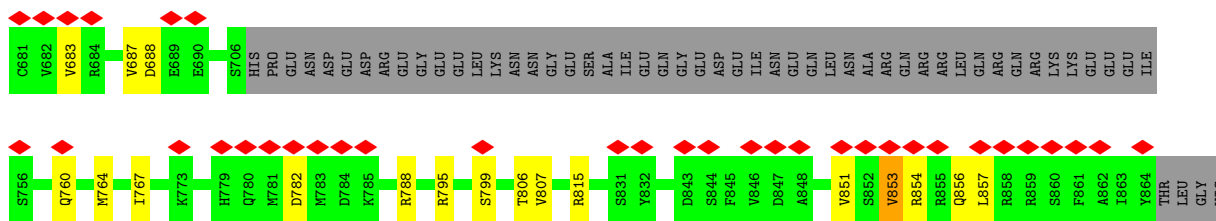
- Molecule 9: Origin recognition complex subunit 6



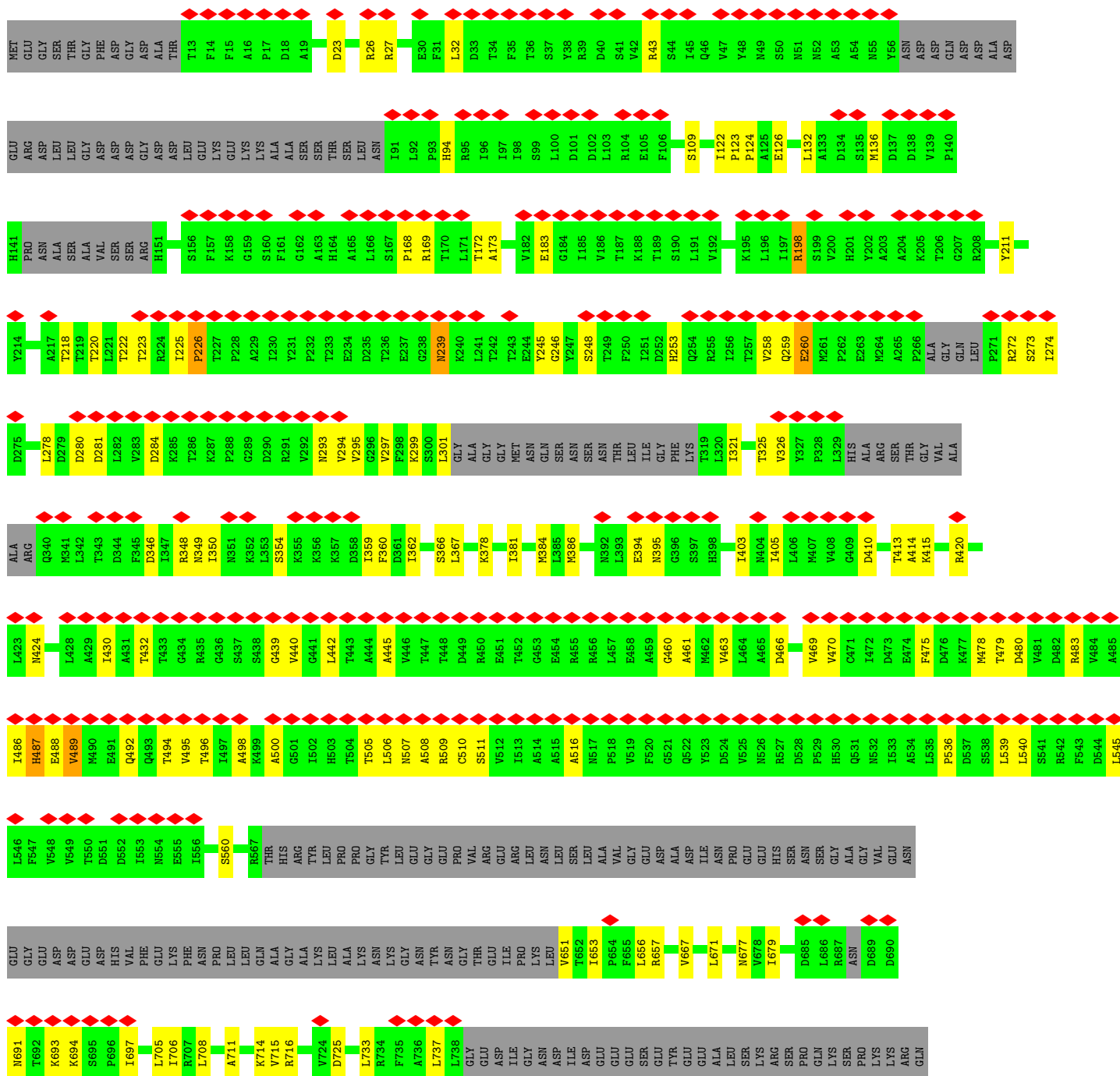


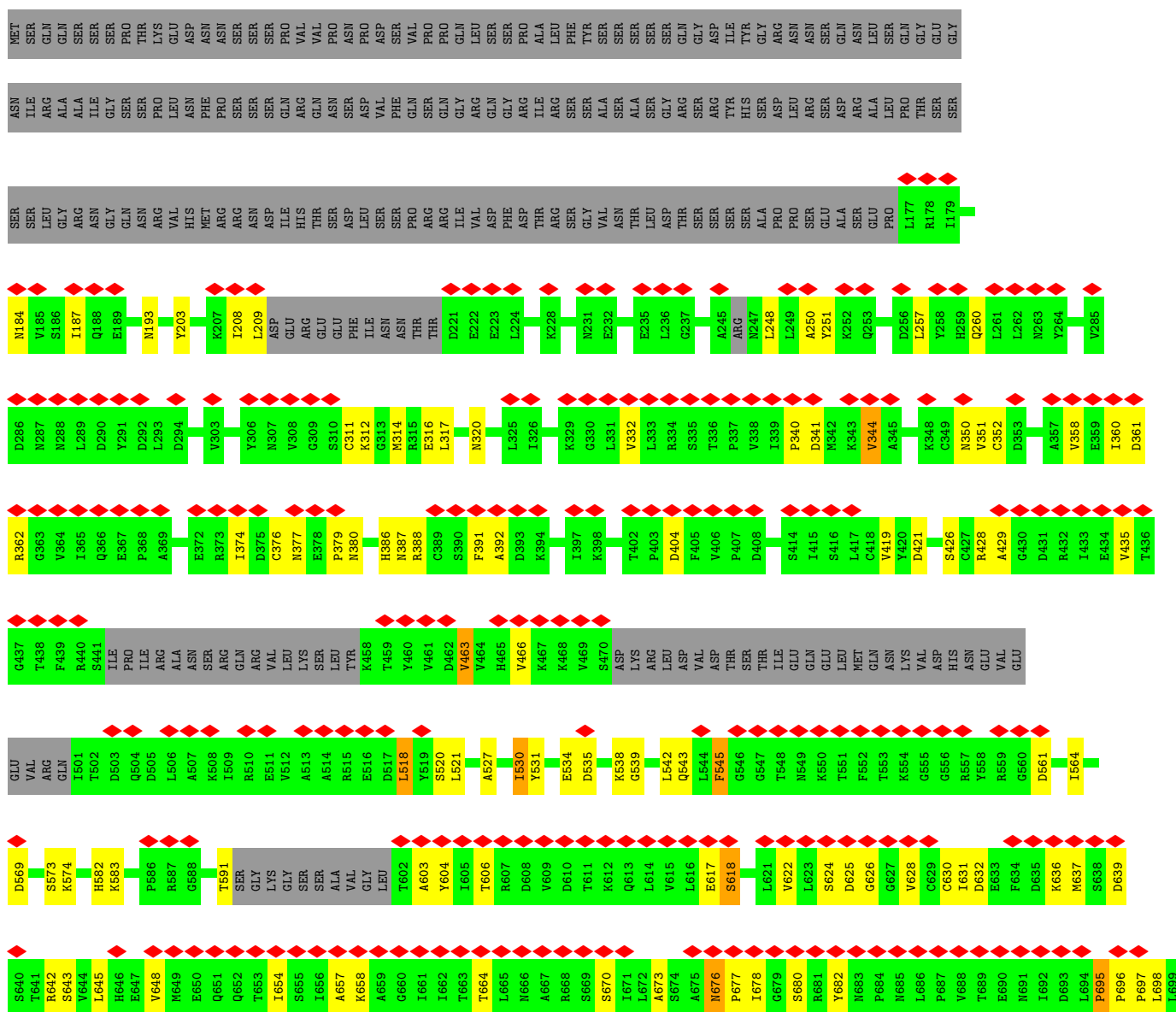
• Molecule 10: DNA replication licensing factor MCM2

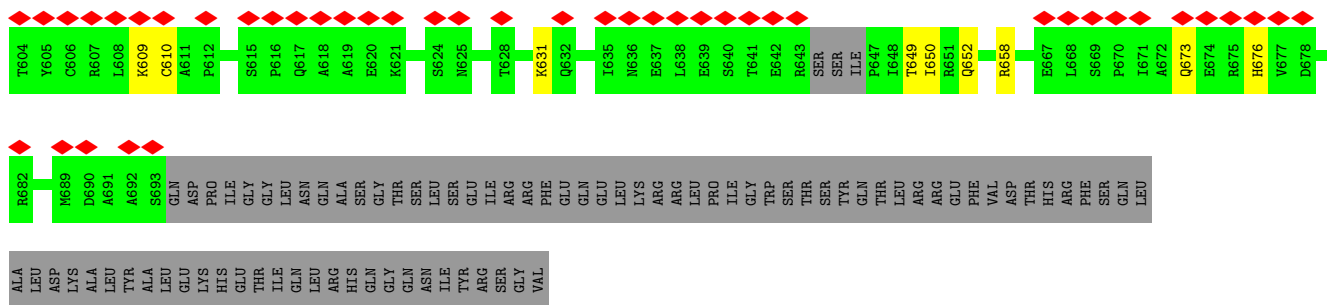




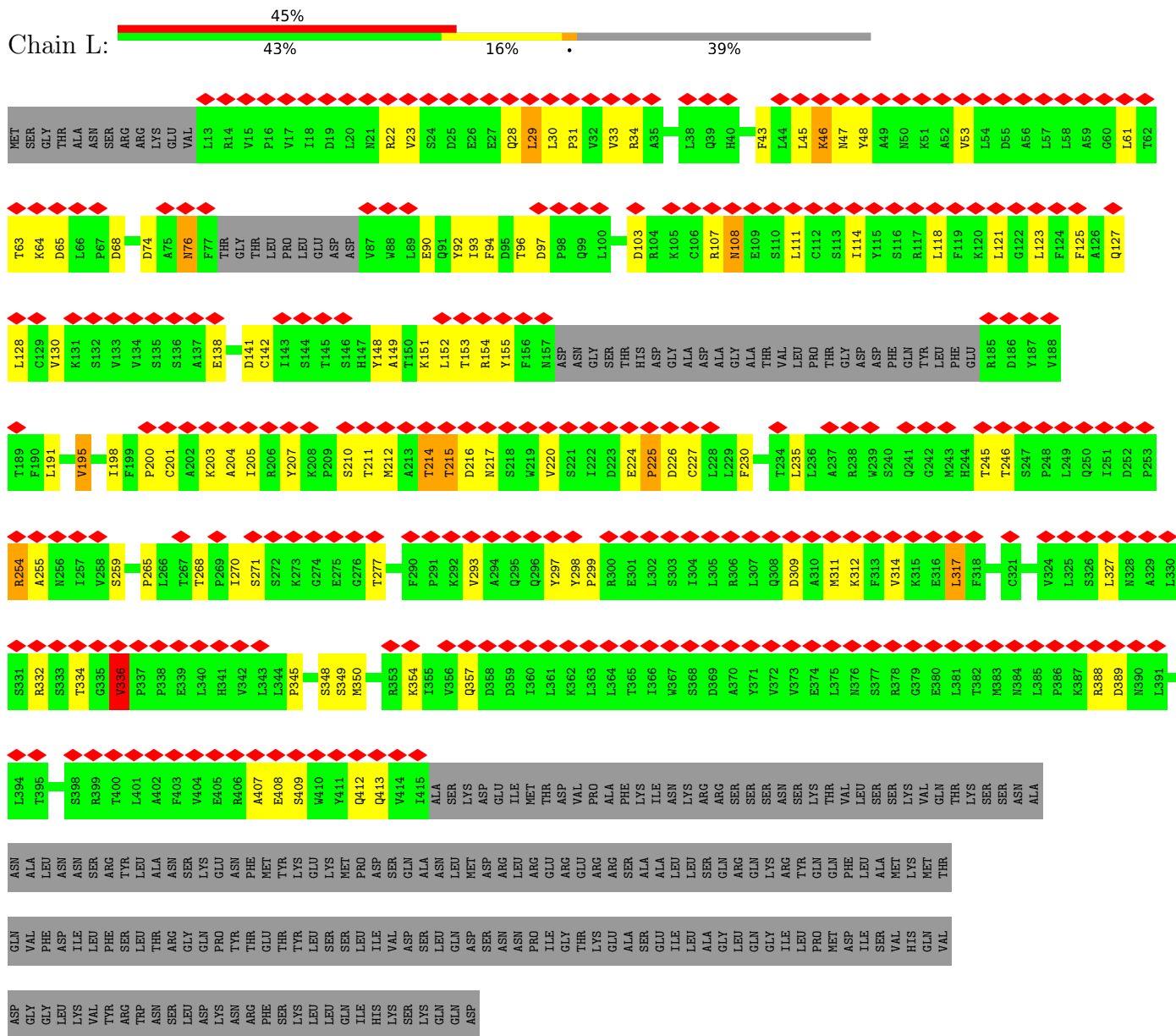
• Molecule 11: DNA replication licensing factor MCM3



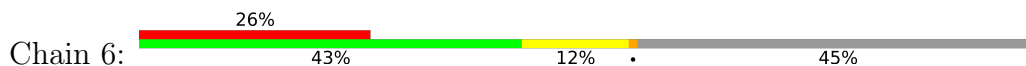




- Molecule 14: Cell division cycle protein CDT1



- Molecule 15: DNA replication licensing factor MCM6





Chain 7:



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	27830	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.050	Depositor
Minimum map value	-0.010	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0173	Depositor
Map size (Å)	335.36, 335.36, 335.36	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.31, 1.31, 1.31	Depositor

5 Model quality

5.1 Standard geometry

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

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	G	0.34	0/772	0.62	0/1193
2	H	0.33	0/787	0.48	0/1209
3	9	0.30	0/3004	0.68	4/4042 (0.1%)
4	A	0.29	0/3369	0.60	0/4536
5	B	0.29	0/2692	0.60	0/3629
6	C	0.28	0/4598	0.58	0/6208
7	E	0.33	0/3497	0.60	1/4756 (0.0%)
8	D	0.32	0/3595	0.60	2/4859 (0.0%)
9	F	0.24	0/1336	0.59	1/1798 (0.1%)
10	2	0.42	0/4340	1.03	25/5873 (0.4%)
11	3	0.44	1/4850 (0.0%)	1.05	28/6597 (0.4%)
12	4	0.40	0/5014	1.07	35/6791 (0.5%)
13	5	0.40	0/4001	0.98	14/5423 (0.3%)
14	L	0.50	0/2958	1.24	42/4028 (1.0%)
15	6	0.38	0/4256	1.02	15/5771 (0.3%)
16	7	0.39	0/4761	0.98	16/6448 (0.2%)
All	All	0.37	1/53830 (0.0%)	0.88	183/73161 (0.3%)

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
11	3	0	1
12	4	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	3	260	GLU	C-N	-10.90	1.10	1.33

The worst 5 of 183 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	6	743	GLU	CA-C-N	12.24	132.99	120.38
15	6	743	GLU	C-N-CA	12.24	132.99	120.38
12	4	618	SER	N-CA-C	10.33	121.51	108.19
14	L	336	VAL	CA-C-N	-9.48	110.62	120.38
14	L	336	VAL	C-N-CA	-9.48	110.62	120.38

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
11	3	260	GLU	Mainchain
12	4	617	GLU	Mainchain

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	G	694	0	399	0	0
2	H	697	0	380	14	0
3	9	2962	0	3063	98	0
4	A	3323	0	3366	70	0
5	B	2640	0	2646	53	0
6	C	4501	0	4445	75	0
7	E	3418	0	3394	75	0
8	D	3534	0	3571	85	0
9	F	1315	0	1353	33	0
10	2	4274	0	4110	100	0
11	3	4773	0	4608	118	0
12	4	4950	0	4771	107	0
13	5	3955	0	3878	54	0
14	L	2898	0	2912	56	0
15	6	4187	0	3940	100	0
16	7	4701	0	4581	86	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
17	9	31	0	12	4	0
17	A	31	0	12	7	0
17	D	31	0	12	3	0
17	E	31	0	12	2	0
18	A	1	0	0	0	0
18	E	1	0	0	0	0
All	All	52948	0	51465	1006	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 10.

The worst 5 of 1006 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:H:35:DA:OP1	3:9:232:ASN:CB	1.95	1.14
10:2:576:LEU:CB	10:2:619:SER:HB3	1.78	1.13
7:E:340:TYR:CB	7:E:373:PRO:HB2	1.77	1.12
14:L:30:LEU:CD1	14:L:128:LEU:O	2.04	1.06
15:6:608:LEU:O	15:6:628:LEU:CB	2.04	1.05

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	
3	9	363/513 (71%)	320 (88%)	41 (11%)	2 (1%)	21	59
4	A	406/913 (44%)	374 (92%)	31 (8%)	1 (0%)	43	78
5	B	313/620 (50%)	290 (93%)	23 (7%)	0	100	100
6	C	536/616 (87%)	506 (94%)	29 (5%)	1 (0%)	43	78
7	E	414/479 (86%)	386 (93%)	27 (6%)	1 (0%)	43	78

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
8	D	428/529 (81%)	402 (94%)	25 (6%)	1 (0%)	43	78
9	F	153/435 (35%)	146 (95%)	7 (5%)	0	100	100
10	2	559/868 (64%)	533 (95%)	23 (4%)	3 (0%)	24	63
11	3	619/971 (64%)	574 (93%)	43 (7%)	2 (0%)	36	72
12	4	634/933 (68%)	594 (94%)	37 (6%)	3 (0%)	24	63
13	5	511/775 (66%)	493 (96%)	15 (3%)	3 (1%)	21	59
14	L	361/604 (60%)	328 (91%)	32 (9%)	1 (0%)	36	72
15	6	551/1017 (54%)	519 (94%)	26 (5%)	6 (1%)	11	46
16	7	605/845 (72%)	578 (96%)	26 (4%)	1 (0%)	43	78
All	All	6453/10118 (64%)	6043 (94%)	385 (6%)	25 (0%)	31	67

5 of 25 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	9	284	GLN
6	C	605	TYR
7	E	445	SER
10	2	350	PRO
10	2	572	SER

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	
3	9	335/470 (71%)	334 (100%)	1 (0%)	86	86
4	A	362/812 (45%)	360 (99%)	2 (1%)	78	83
5	B	294/573 (51%)	294 (100%)	0	100	100
6	C	504/576 (88%)	504 (100%)	0	100	100
7	E	386/440 (88%)	386 (100%)	0	100	100
8	D	397/488 (81%)	397 (100%)	0	100	100
9	F	151/406 (37%)	151 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
10	2	438/770 (57%)	435 (99%)	3 (1%)	76	81
11	3	493/835 (59%)	490 (99%)	3 (1%)	78	83
12	4	528/848 (62%)	525 (99%)	3 (1%)	78	83
13	5	424/688 (62%)	424 (100%)	0	100	100
14	L	325/545 (60%)	319 (98%)	6 (2%)	51	68
15	6	416/886 (47%)	413 (99%)	3 (1%)	76	81
16	7	498/753 (66%)	497 (100%)	1 (0%)	87	87
All	All	5551/9090 (61%)	5529 (100%)	22 (0%)	81	84

5 of 22 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
14	L	130	VAL
14	L	336	VAL
14	L	277	THR
15	6	355	ASP
11	3	239	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 111 such sidechains are listed below:

Mol	Chain	Res	Type
8	D	511	GLN
16	7	585	ASN
11	3	517	ASN
16	7	568	ASN
14	L	412	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
17	AGS	D	2001	-	32,33,33	1.26	3 (9%)	45,52,52	1.08	1 (2%)
17	AGS	E	2001	18	32,33,33	0.96	2 (6%)	45,52,52	0.74	1 (2%)
17	AGS	9	2001	-	32,33,33	0.86	2 (6%)	45,52,52	0.89	1 (2%)
17	AGS	A	2001	-	32,33,33	1.05	3 (9%)	45,52,52	0.94	1 (2%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
17	AGS	D	2001	-	-	2/21/38/38	0/3/3/3
17	AGS	E	2001	18	-	7/21/38/38	0/3/3/3
17	AGS	9	2001	-	-	7/21/38/38	0/3/3/3
17	AGS	A	2001	-	-	7/21/38/38	0/3/3/3

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	D	2001	AGS	PB-O3B	-4.10	1.55	1.59
17	D	2001	AGS	PB-O3A	-3.92	1.55	1.59
17	E	2001	AGS	PB-O3A	-3.67	1.55	1.59
17	A	2001	AGS	PB-O3B	-3.63	1.55	1.59
17	D	2001	AGS	PA-O3A	-3.50	1.55	1.59

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
17	D	2001	AGS	PB-O3B-PG	-6.59	109.06	133.17
17	A	2001	AGS	PB-O3B-PG	-5.36	113.54	133.17
17	9	2001	AGS	PB-O3B-PG	-4.85	115.44	133.17
17	E	2001	AGS	PB-O3B-PG	-3.38	120.79	133.17

There are no chirality outliers.

5 of 23 torsion outliers are listed below:

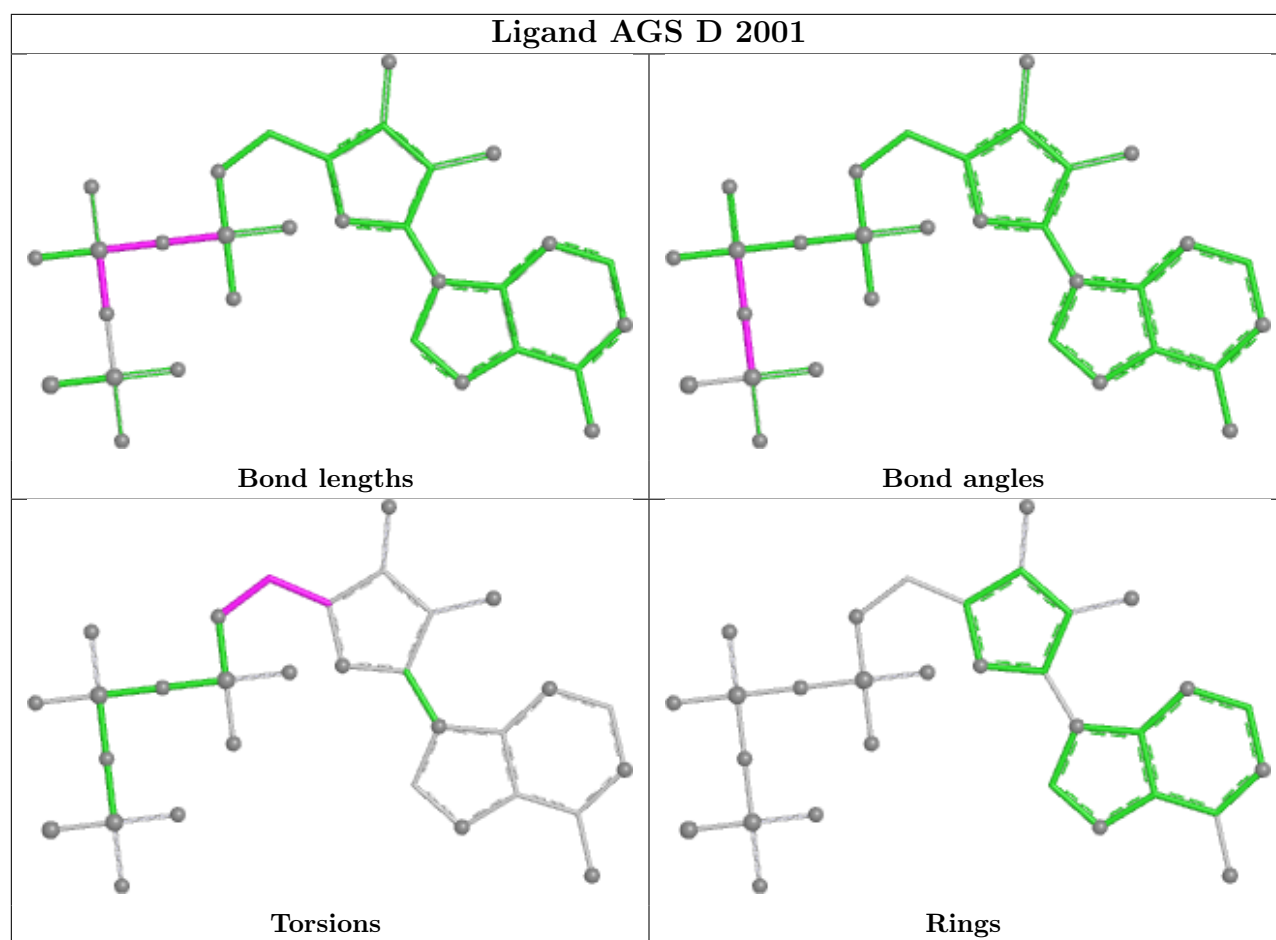
Mol	Chain	Res	Type	Atoms
17	9	2001	AGS	C5'-O5'-PA-O2A
17	9	2001	AGS	C5'-O5'-PA-O3A
17	A	2001	AGS	C5'-O5'-PA-O1A
17	A	2001	AGS	C5'-O5'-PA-O2A
17	A	2001	AGS	C5'-O5'-PA-O3A

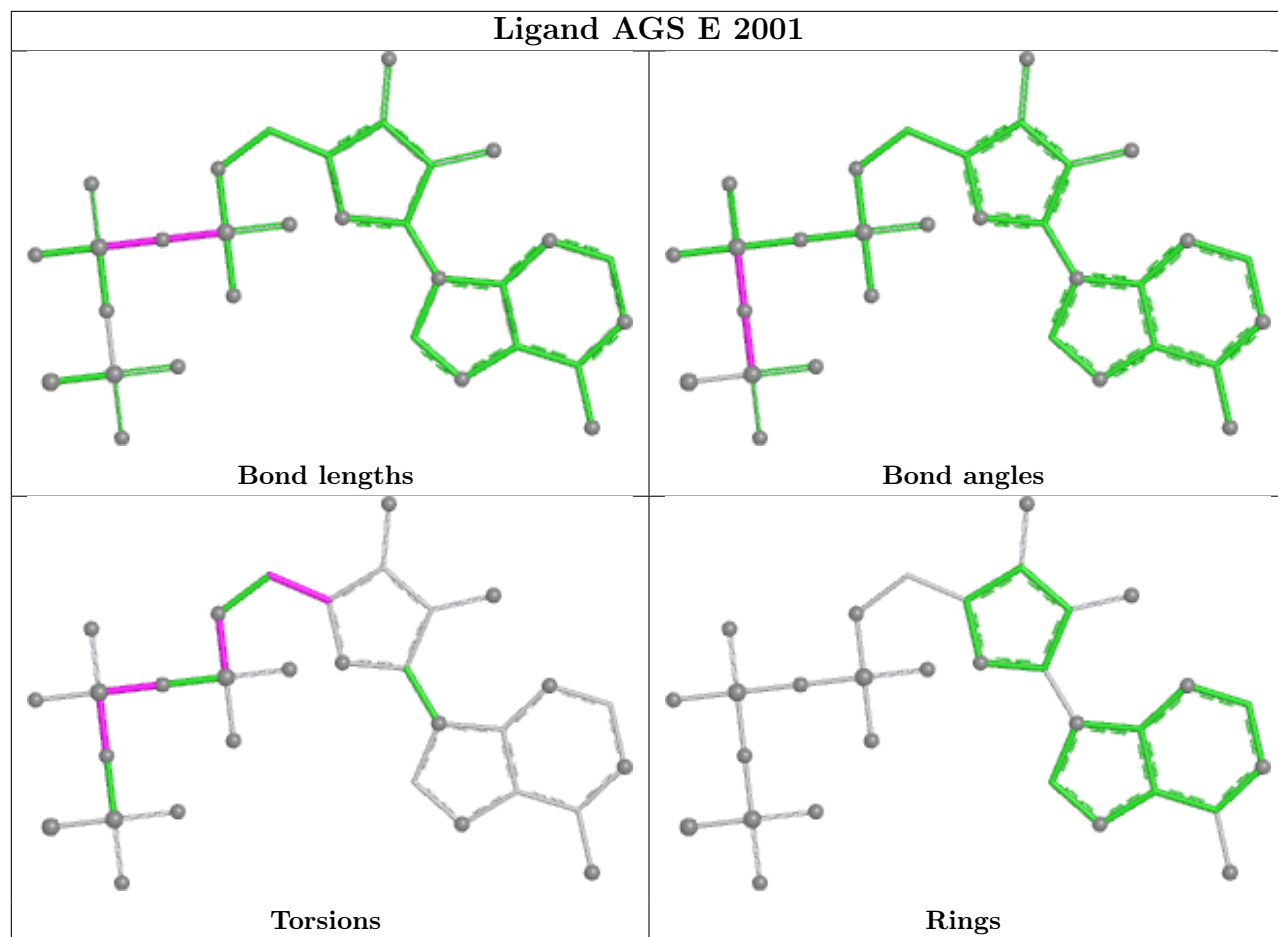
There are no ring outliers.

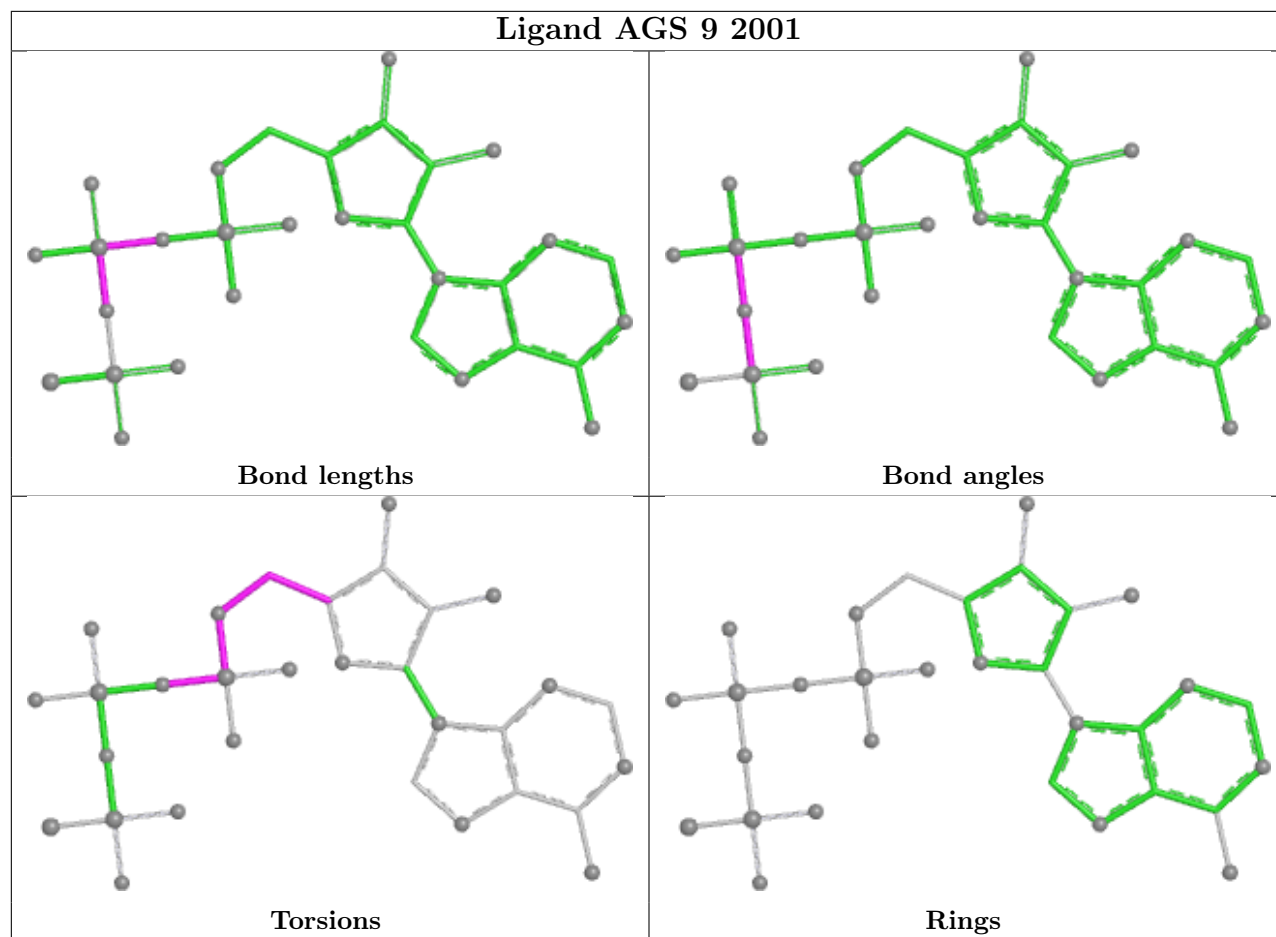
4 monomers are involved in 16 short contacts:

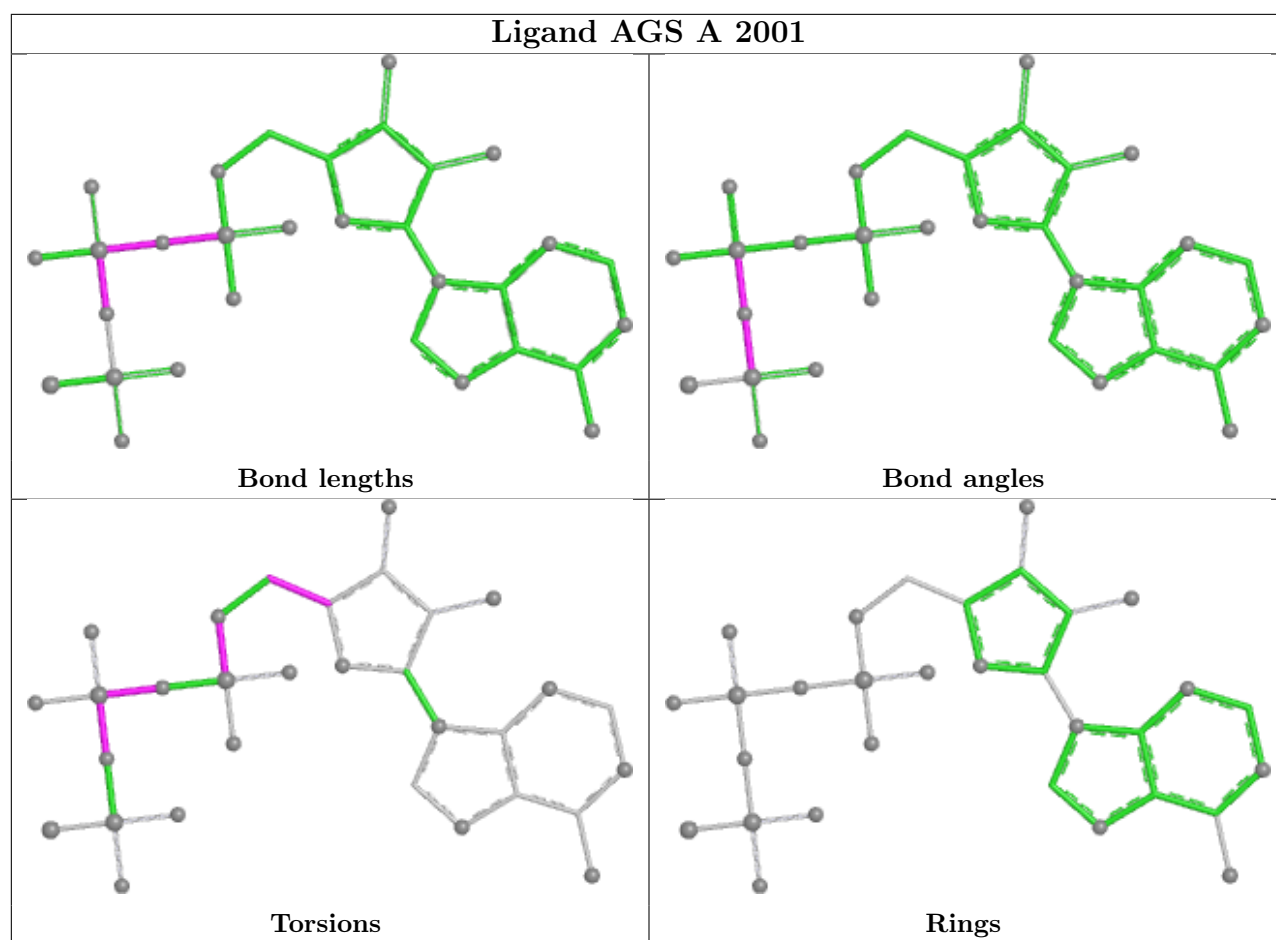
Mol	Chain	Res	Type	Clashes	Symm-Clashes
17	D	2001	AGS	3	0
17	E	2001	AGS	2	0
17	9	2001	AGS	4	0
17	A	2001	AGS	7	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









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
3	9	1
11	3	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	9	399:ASN	C	400:ASN	N	3.97
1	3	260:GLU	C	261:MET	N	1.11

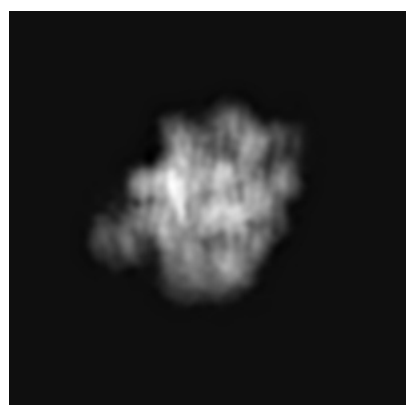
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21666. These allow visual inspection of the internal detail of the map and identification of artifacts.

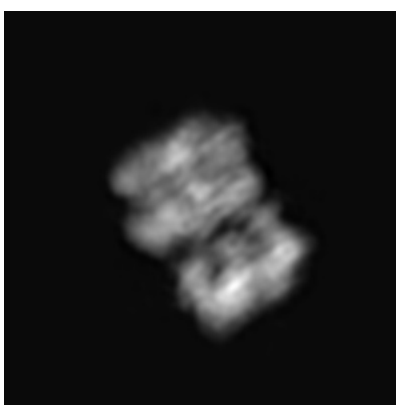
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

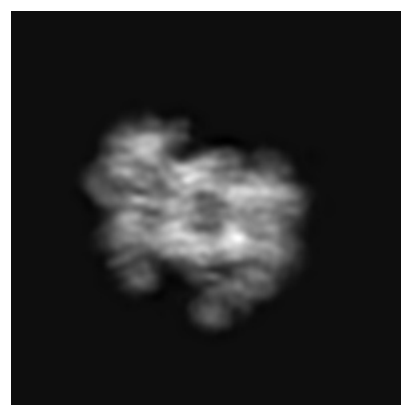
6.1.1 Primary map



X



Y

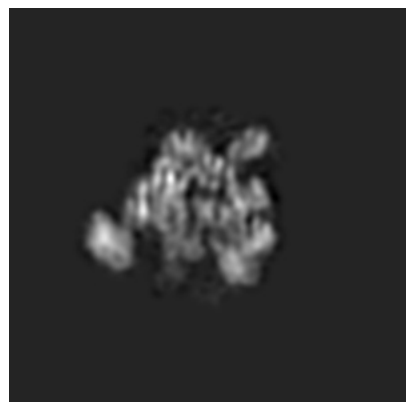


Z

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

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 128



Y Index: 128



Z Index: 128

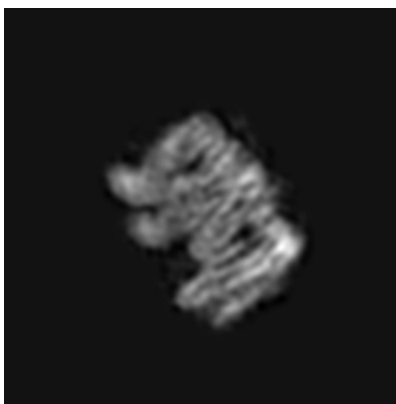
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 85



Y Index: 109

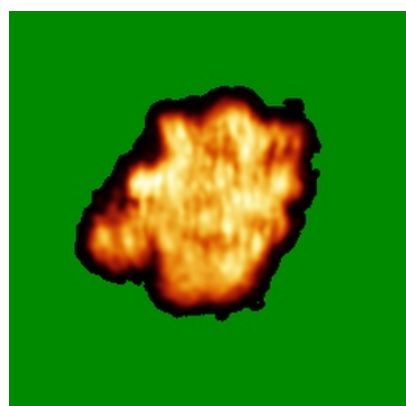


Z Index: 145

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

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

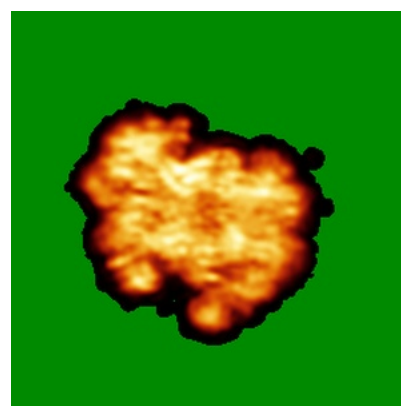
6.4.1 Primary map



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

This section was not generated.

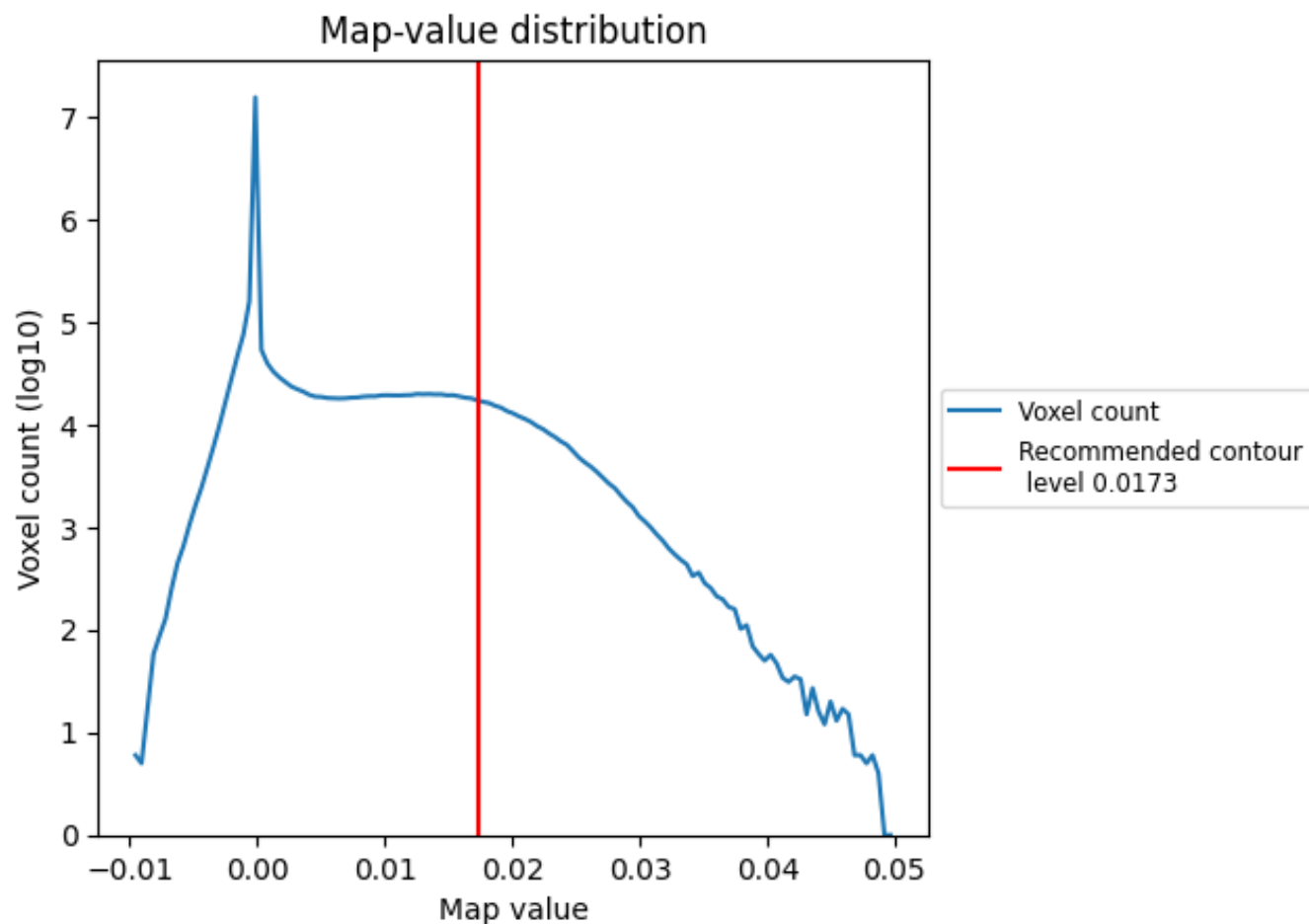
6.6 Mask visualisation

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

7 Map analysis [i](#)

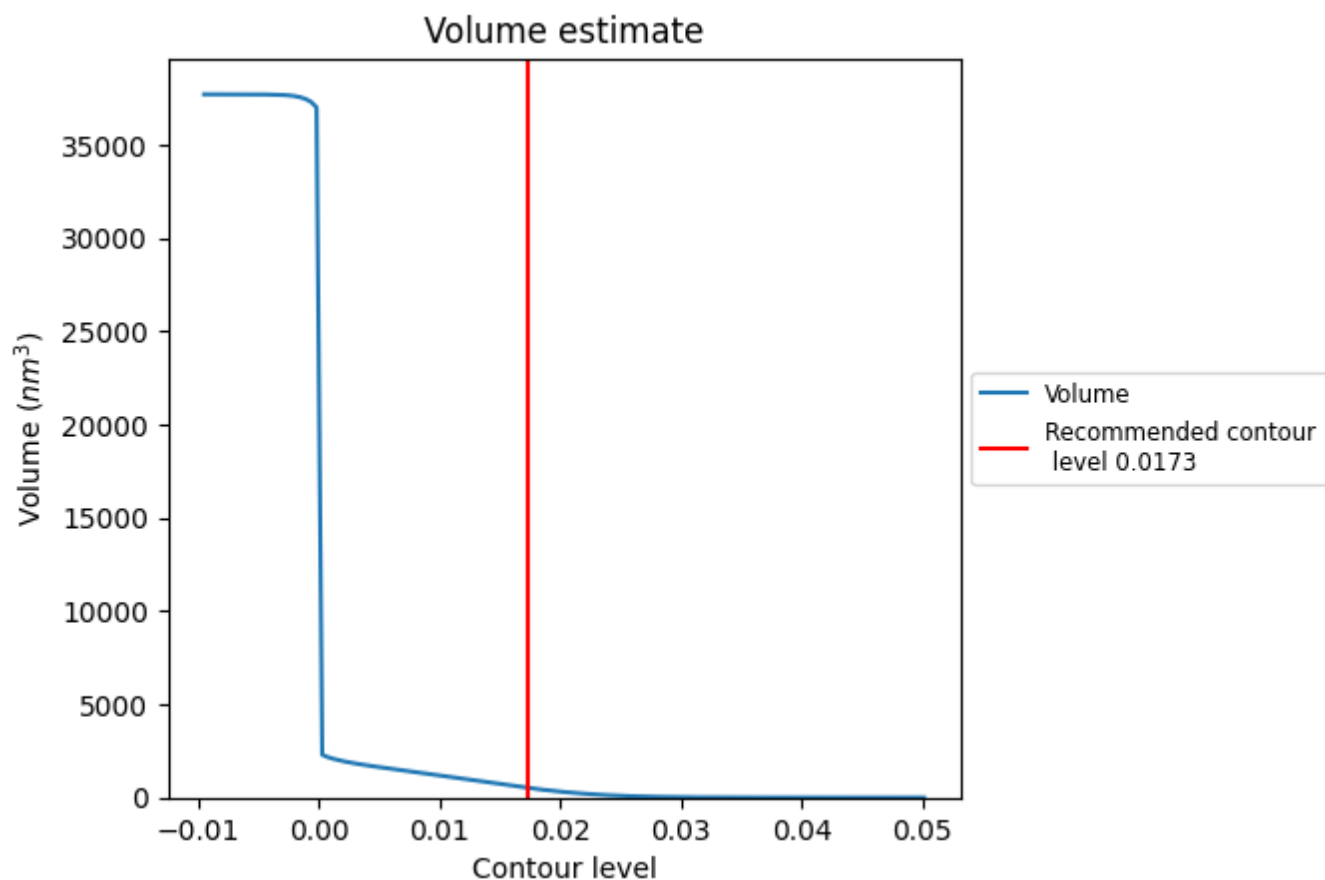
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

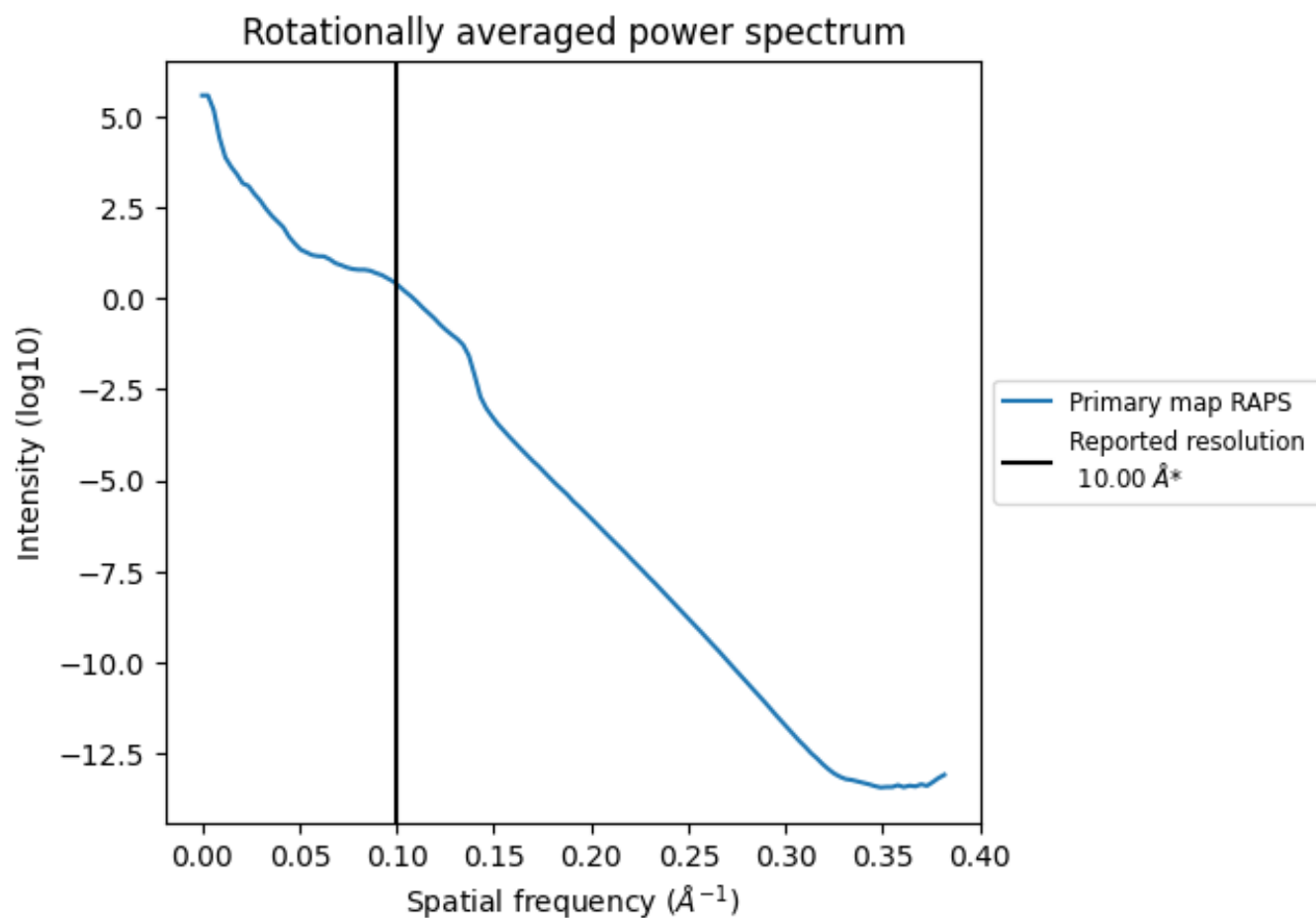
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 519 nm³; this corresponds to an approximate mass of 469 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 0.100 Å⁻¹

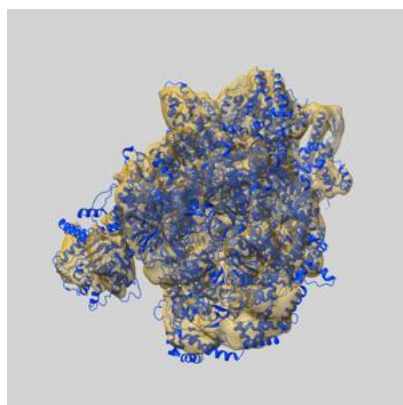
8 Fourier-Shell correlation

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

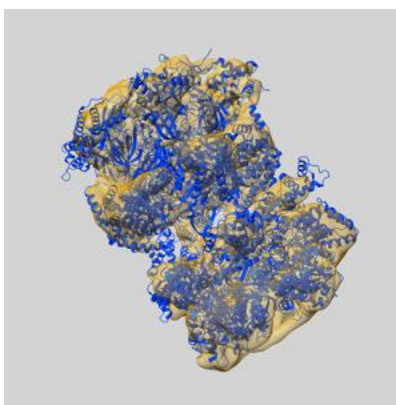
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-21666 and PDB model 6WGI. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

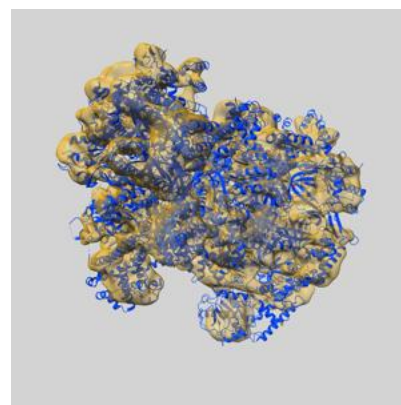
9.1 Map-model overlay [i](#)



X



Y



Z

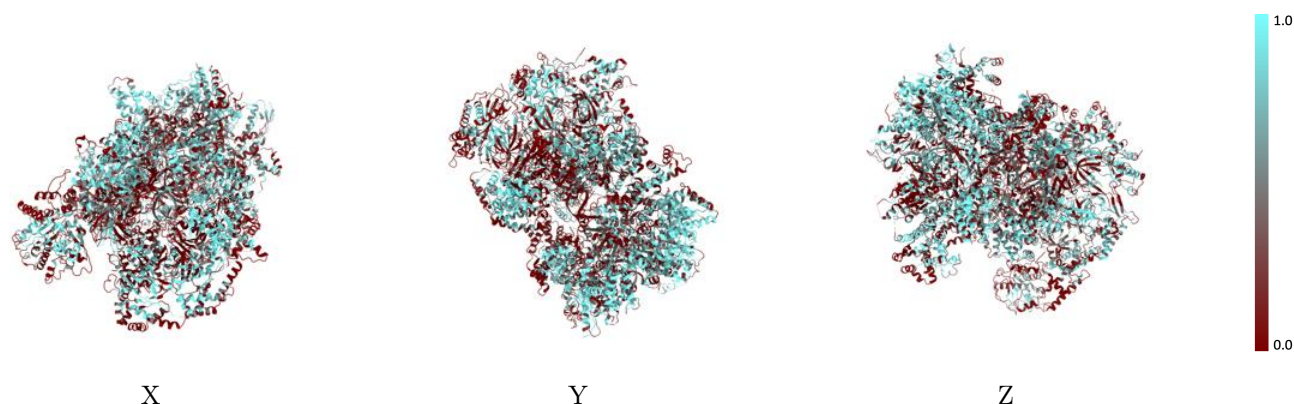
The images above show the 3D surface view of the map at the recommended contour level 0.0173 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



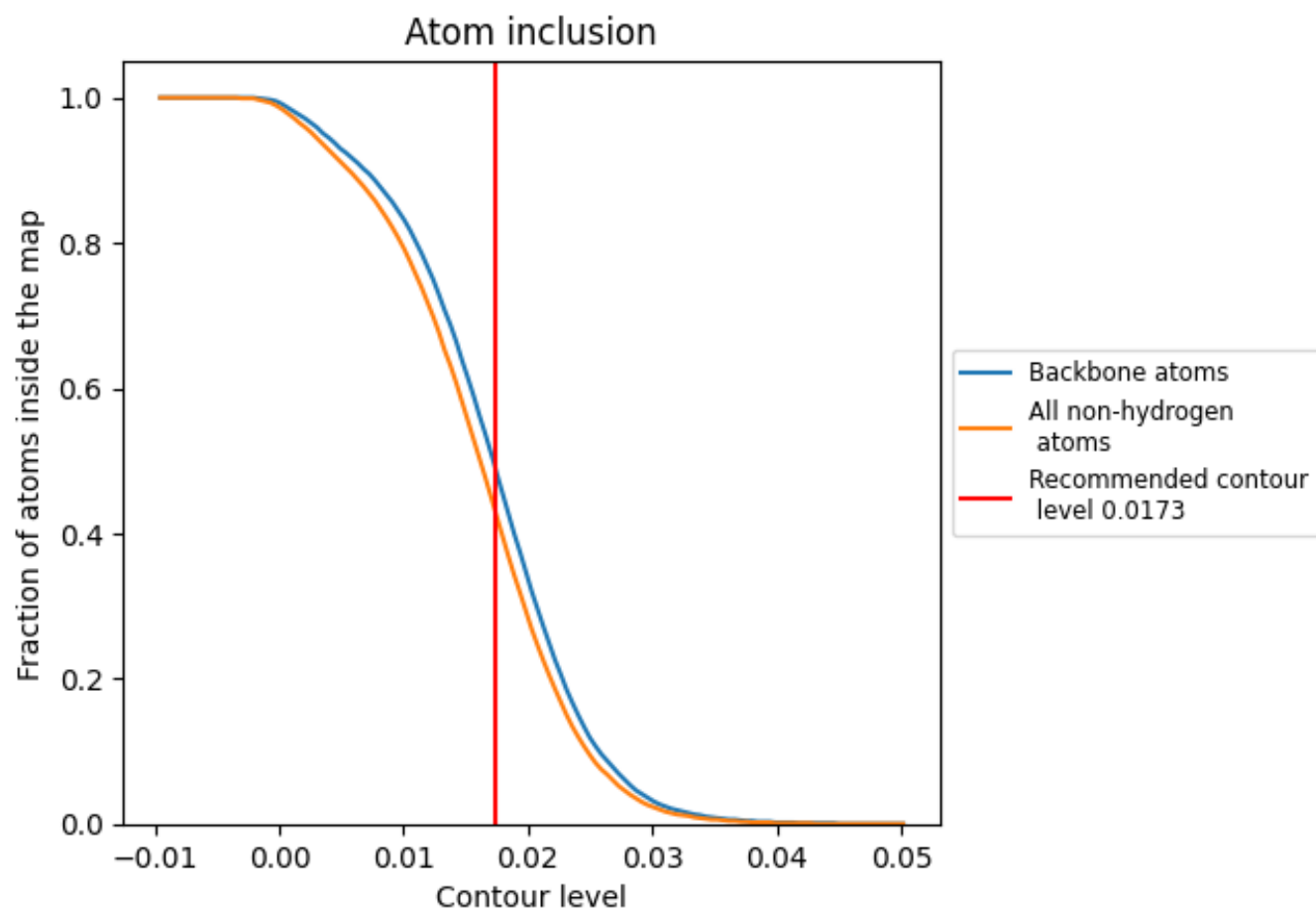
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0173).

9.4 Atom inclusion [i](#)



At the recommended contour level, 50% of all backbone atoms, 43% 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 (0.0173) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.4340	 0.0860
2	 0.4520	 0.0970
3	 0.4410	 0.0710
4	 0.4210	 0.0990
5	 0.2950	 0.0450
6	 0.4890	 0.0900
7	 0.4520	 0.0890
9	 0.4550	 0.0710
A	 0.5080	 0.0980
B	 0.3500	 0.0900
C	 0.4870	 0.0980
D	 0.6170	 0.1050
E	 0.5630	 0.1060
F	 0.0060	 0.0330
G	 0.3130	 0.0950
H	 0.3530	 0.1190
L	 0.2380	 0.0700

