



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 08:27 AM UTC

PDB ID : 7EGC / pdb_00007egc
EMDB ID : EMD-31112
Title : p53-bound TFIID-based holo PIC on HDM2 promoter
Authors : Chen, X.; Wu, Z.; Hou, H.; Qi, Y.; Wang, X.; Li, J.; Xu, Y.
Deposited on : 2021-03-24
Resolution : 3.90 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:

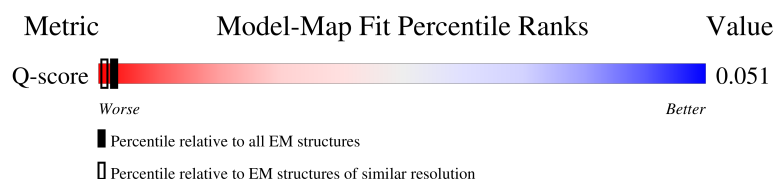
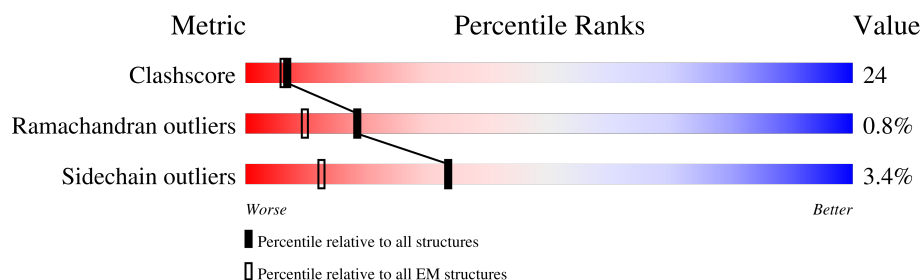
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 3.90 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	8855 (3.40 - 4.40)

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	0	309	20% 63% 12% 24%
2	1	548	16% 54% 17% 26%
3	2	395	5% 62% 20% 16%
4	3	308	6% 61% 23% 15%

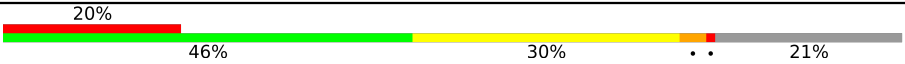
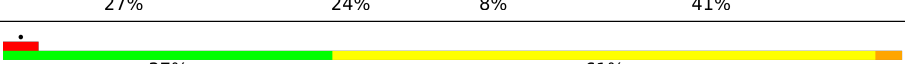
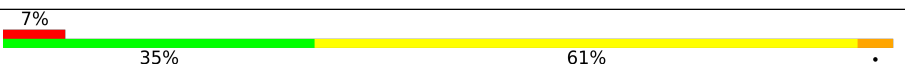
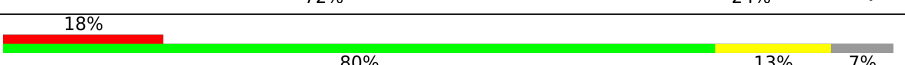
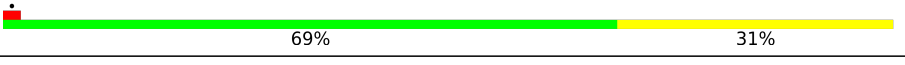
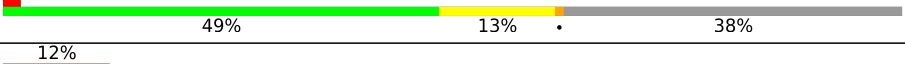
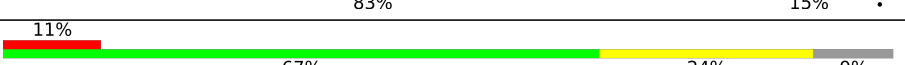
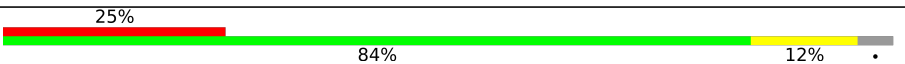
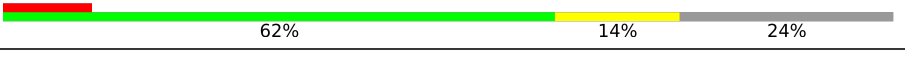
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Mol	Chain	Length	Quality of chain
5	4	462	
6	5	71	
7	6	782	
8	7	760	
9	8	346	
10	9	323	
11	A	1872	
12	B	1199	
13	D	1085	
13	d	1085	
14	E	800	
14	e	800	
15	F	677	
15	f	677	
16	G	349	
17	H	310	
18	I	264	
18	i	264	
19	J	218	
19	j	218	
20	L	161	
20	l	161	
21	O	109	
22	P	339	
23	Q	376	

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Mol	Chain	Length	Quality of chain
24	R	316	
25	S	517	
26	T	249	
27	U	439	
28	V	291	
29	X	71	
30	Y	71	
31	c	929	
32	k	211	
33	m	124	
34	o	1970	
35	p	1174	
36	q	275	
37	r	142	
38	s	210	
39	t	127	
40	u	172	
41	v	150	
42	w	125	
43	x	67	
44	y	117	
45	z	58	

2 Entry composition

There are 48 unique types of molecules in this entry. The entry contains 113129 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 CDK-activating kinase assembly factor MAT1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0	234	Total	C	N	O	S	0	0
			1774	1108	309	347	10		

- Molecule 2 is a protein called General transcription factor IIH subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	405	Total	C	N	O	S	0	0
			2634	1640	486	501	7		

- Molecule 3 is a protein called General transcription factor IIH subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	2	331	Total	C	N	O	S	0	0
			2534	1597	441	470	26		

- Molecule 4 is a protein called General transcription factor IIH subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	3	263	Total	C	N	O	S	0	0
			2065	1323	344	379	19		

- Molecule 5 is a protein called General transcription factor IIH subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	4	449	Total	C	N	O	S	0	0
			3579	2303	624	638	14		

- Molecule 6 is a protein called General transcription factor IIH subunit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	5	54	Total	C	N	O	S	0	0
			428	277	67	82	2		

- Molecule 7 is a protein called General transcription and DNA repair factor IIIH helicase subunit XPB.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	6	606	Total	C	N	O	S	0	0
			4880	3117	849	884	30		

- Molecule 8 is a protein called General transcription and DNA repair factor IIIH helicase subunit XPD.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	7	734	Total	C	N	O	S	0	0
			5833	3727	1022	1055	29		

- Molecule 9 is a protein called Cyclin-dependent kinase 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	8	298	Total	C	N	O	S	0	0
			2370	1531	404	424	11		

- Molecule 10 is a protein called Cyclin-H.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	9	287	Total	C	N	O	S	0	0
			2334	1493	402	422	17		

- Molecule 11 is a protein called Transcription initiation factor TFIID subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	A	550	Total	C	N	O	S	0	0
			4511	2882	782	820	27		

- Molecule 12 is a protein called Transcription initiation factor TFIID subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	B	963	Total	C	N	O	S	0	0
			7796	5011	1315	1412	58		

- Molecule 13 is a protein called Transcription initiation factor TFIID subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	D	158	Total	C	N	O	S	0	0
			1322	824	247	248	3		

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Mol	Chain	Residues	Atoms					AltConf	Trace
13	d	158	Total	C	N	O	S	0	0
			1307	814	238	252	3		

- Molecule 14 is a protein called Transcription initiation factor TFIID subunit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	E	546	Total	C	N	O	S	0	0
			4364	2766	757	820	21		
14	e	539	Total	C	N	O	S	0	0
			4327	2746	748	814	19		

- Molecule 15 is a protein called Transcription initiation factor TFIID subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	F	408	Total	C	N	O	S	0	0
			3109	1970	542	579	18		
15	f	403	Total	C	N	O	S	0	0
			3081	1954	533	576	18		

- Molecule 16 is a protein called Transcription initiation factor TFIID subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	G	145	Total	C	N	O	S	0	0
			1180	748	217	211	4		

- Molecule 17 is a protein called Transcription initiation factor TFIID subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	H	209	Total	C	N	O	S	0	0
			1633	1034	283	311	5		

- Molecule 18 is a protein called Transcription initiation factor TFIID subunit 9.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	I	120	Total	C	N	O	S	0	0
			959	610	166	177	6		
18	i	121	Total	C	N	O	S	0	0
			967	615	167	178	7		

- Molecule 19 is a protein called Transcription initiation factor TFIID subunit 10.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	J	90	Total	C	N	O	S	0	0
			720	466	115	135	4		
19	j	95	Total	C	N	O	S	0	0
			759	488	124	143	4		

- Molecule 20 is a protein called Transcription initiation factor TFIID subunit 12.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	L	75	Total	C	N	O	S	0	0
			614	384	107	120	3		
20	l	107	Total	C	N	O	S	0	0
			876	547	158	166	5		

- Molecule 21 is a protein called Transcription initiation factor IIA subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	O	97	Total	C	N	O	S	0	0
			771	491	133	145	2		

- Molecule 22 is a protein called TATA-box-binding protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	P	177	Total	C	N	O	S	0	0
			1412	918	249	238	7		

- Molecule 23 is a protein called Transcription initiation factor IIA subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	Q	122	Total	C	N	O	S	0	0
			996	623	162	207	4		

- Molecule 24 is a protein called Transcription initiation factor IIB.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	R	250	Total	C	N	O	S	0	0
			1929	1209	343	360	17		

- Molecule 25 is a protein called General transcription factor IIF subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	S	108	Total	C	N	O	S	0	0
			872	558	153	159	2		

- Molecule 26 is a protein called General transcription factor IIF subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	T	222	Total	C	N	O	S	0	0
			1788	1127	320	338	3		

- Molecule 27 is a protein called General transcription factor IIE subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	U	179	Total	C	N	O	S	0	0
			1476	932	261	272	11		

- Molecule 28 is a protein called Transcription initiation factor IIE subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	V	172	Total	C	N	O	S	0	0
			1404	893	243	264	4		

- Molecule 29 is a DNA chain called DNA (71-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
29	X	71	Total	C	N	O	P	0	0
			1464	695	265	433	71		

- Molecule 30 is a DNA chain called DNA (71-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
30	Y	71	Total	C	N	O	P	0	0
			1447	687	270	419	71		

- Molecule 31 is a protein called Transcription initiation factor TFIID subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	c	127	Total	C	N	O	S	0	0
			1011	638	174	193	6		

- Molecule 32 is a protein called Transcription initiation factor TFIID subunit 11.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	k	98	Total	C	N	O	S	0	0
			785	499	142	139	5		

- Molecule 33 is a protein called Transcription initiation factor TFIID subunit 13.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	m	87	Total	C	N	O	S	0	0
			724	456	131	131	6		

- Molecule 34 is a protein called DNA-directed RNA polymerase II subunit RPB1.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	o	1427	Total	C	N	O	S	0	0
			11308	7114	2023	2099	72		

- Molecule 35 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	p	1134	Total	C	N	O	S	0	0
			9062	5732	1595	1671	64		

- Molecule 36 is a protein called DNA-directed RNA polymerase II subunit RPB3.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	q	257	Total	C	N	O	S	0	0
			2059	1294	351	408	6		

- Molecule 37 is a protein called DNA-directed RNA polymerase II subunit RPB4.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	r	128	Total	C	N	O	S	0	0
			1005	632	172	197	4		

- Molecule 38 is a protein called DNA-directed RNA polymerase II subunit E.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	s	209	Total	C	N	O	S	0	0
			1720	1089	300	323	8		

- Molecule 39 is a protein called DNA-directed RNA polymerase II subunit F.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	t	79	Total	C	N	O	S	0	0
			635	406	108	116	5		

- Molecule 40 is a protein called DNA-directed RNA polymerase II subunit RPB7.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	u	171	Total	C	N	O	S	0	0
			1351	875	219	249	8		

- Molecule 41 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC3.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	v	148	Total	C	N	O	S	0	0
			1186	750	194	237	5		

- Molecule 42 is a protein called DNA-directed RNA polymerase II subunit RPB9.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	w	114	Total	C	N	O	S	0	0
			927	571	166	179	11		

- Molecule 43 is a protein called RPB10.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	x	64	Total	C	N	O	S	0	0
			507	328	86	87	6		

- Molecule 44 is a protein called RNA_pol_L_2 domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	y	117	Total	C	N	O	S	0	0
			937	604	154	177	2		

- Molecule 45 is a protein called RPB12.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	z	44	Total	C	N	O	S	0	0
			372	231	72	63	6		

- Molecule 46 is ZINC ION (CCD ID: ZN) (formula: Zn).

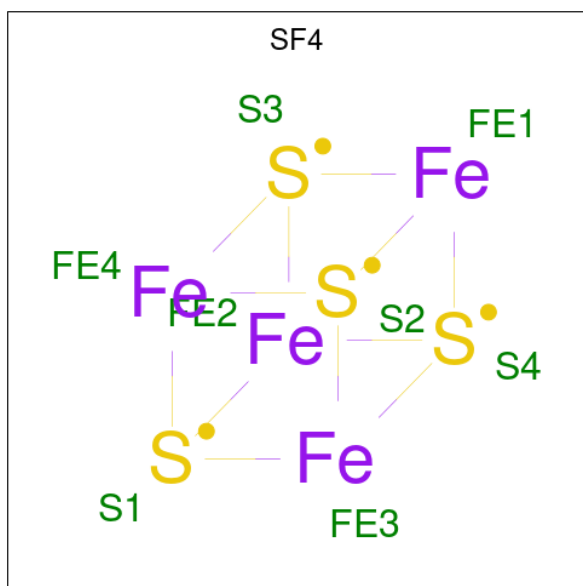
Mol	Chain	Residues	Atoms		AltConf
46	0	2	Total	Zn	0
			2	2	
46	2	3	Total	Zn	0
			3	3	
46	3	1	Total	Zn	0
			1	1	

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Mol	Chain	Residues	Atoms		AltConf
46	R	1	Total 1	Zn 1	0
46	U	1	Total 1	Zn 1	0
46	o	2	Total 2	Zn 2	0
46	p	1	Total 1	Zn 1	0
46	q	1	Total 1	Zn 1	0
46	w	2	Total 2	Zn 2	0
46	x	1	Total 1	Zn 1	0
46	z	1	Total 1	Zn 1	0

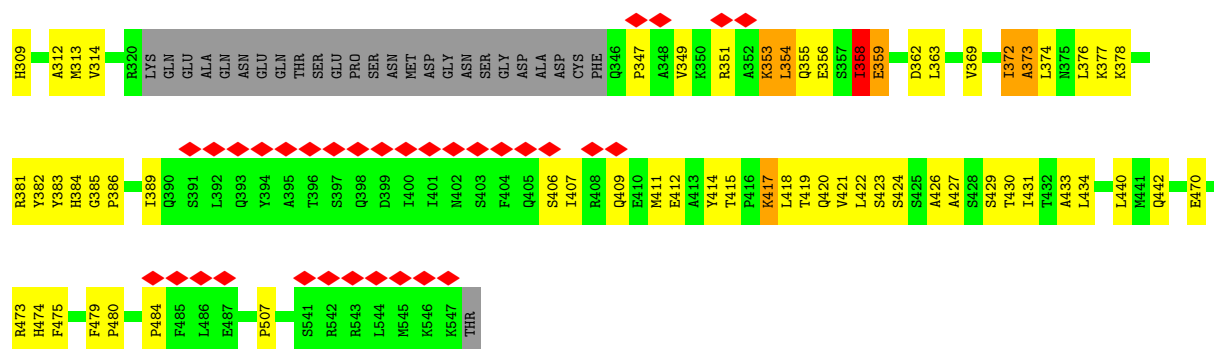
- Molecule 47 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



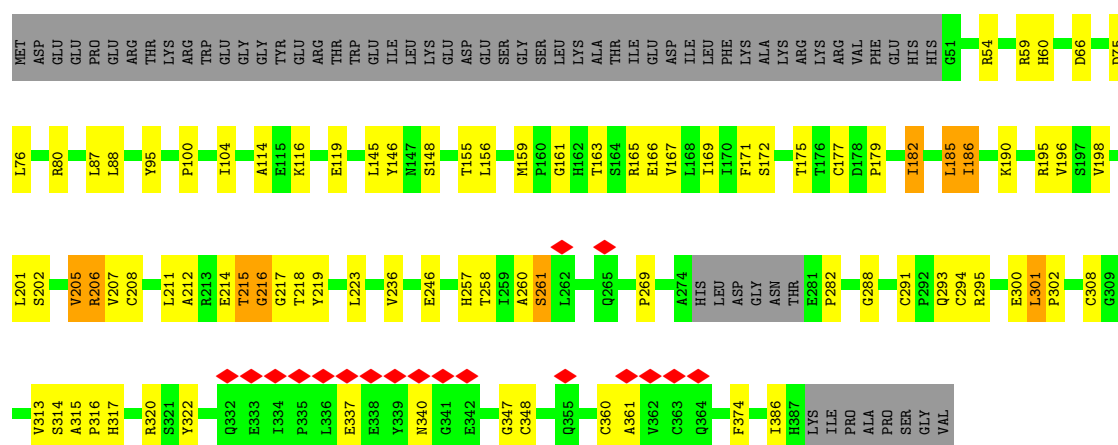
Mol	Chain	Residues	Atoms			AltConf
47	7	1	Total 8	Fe 4	S 4	0

- Molecule 48 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

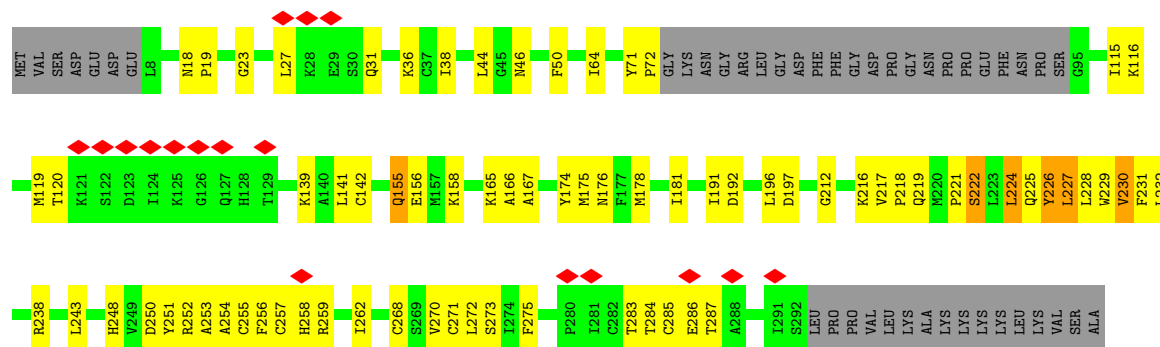
Mol	Chain	Residues	Atoms		AltConf
48	o	1	Total	Mg	0
			1	1	



• Molecule 3: General transcription factor IIH subunit 2

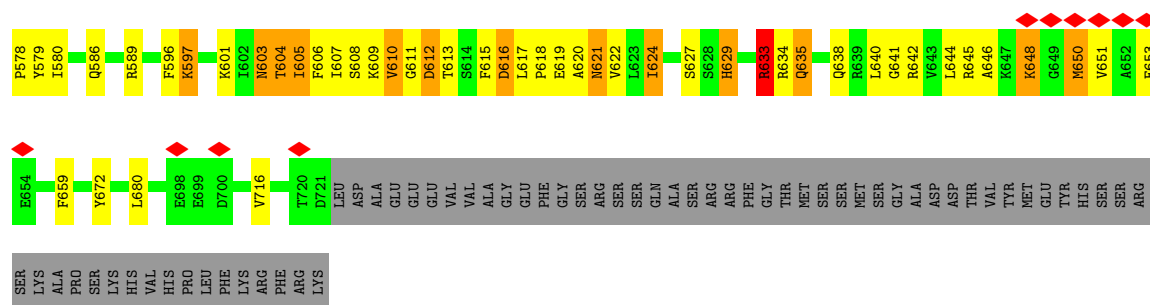


• Molecule 4: General transcription factor IIH subunit 3

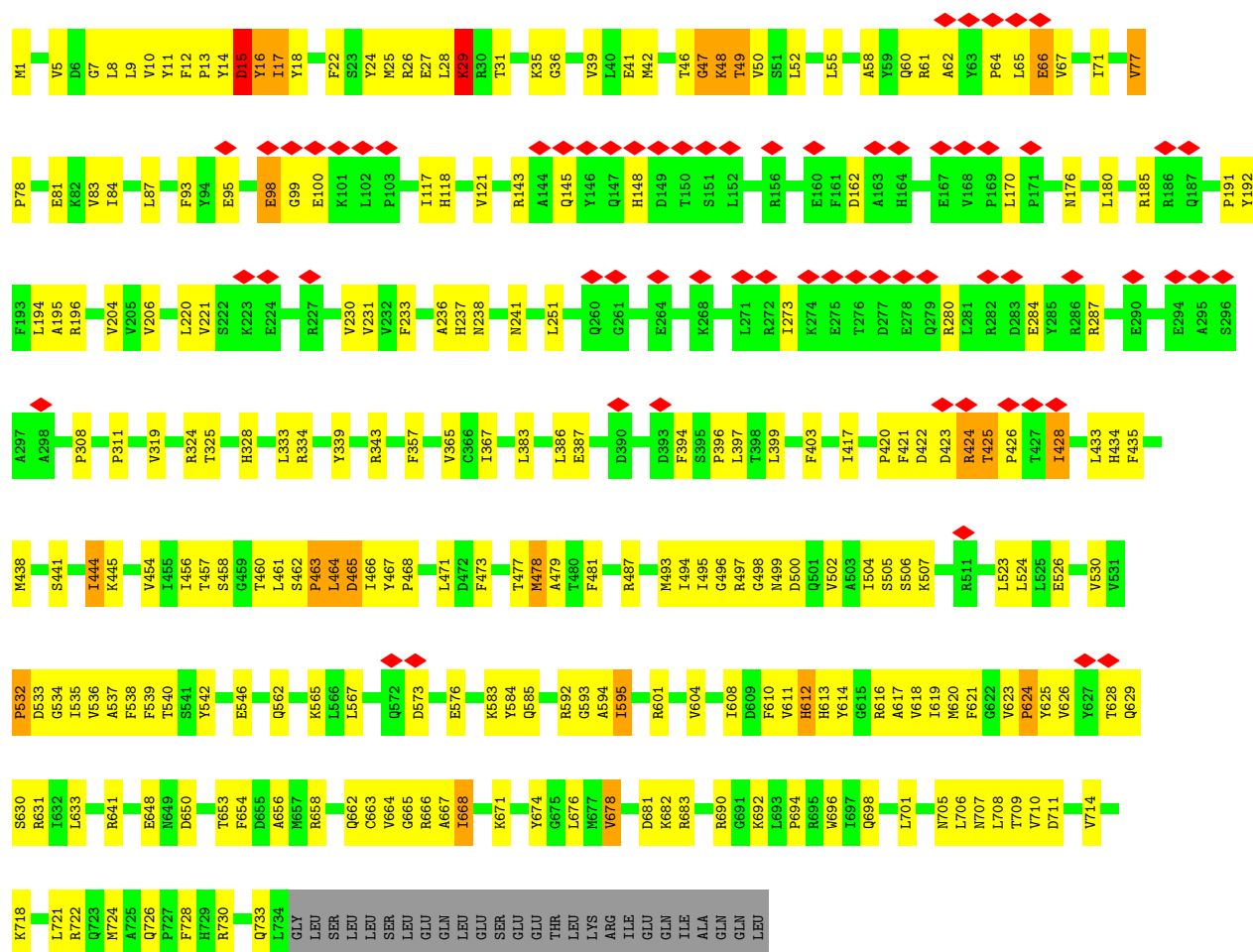


• Molecule 5: General transcription factor IIH subunit 4

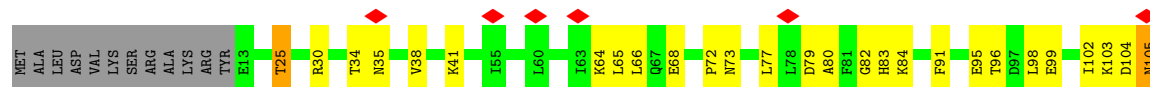


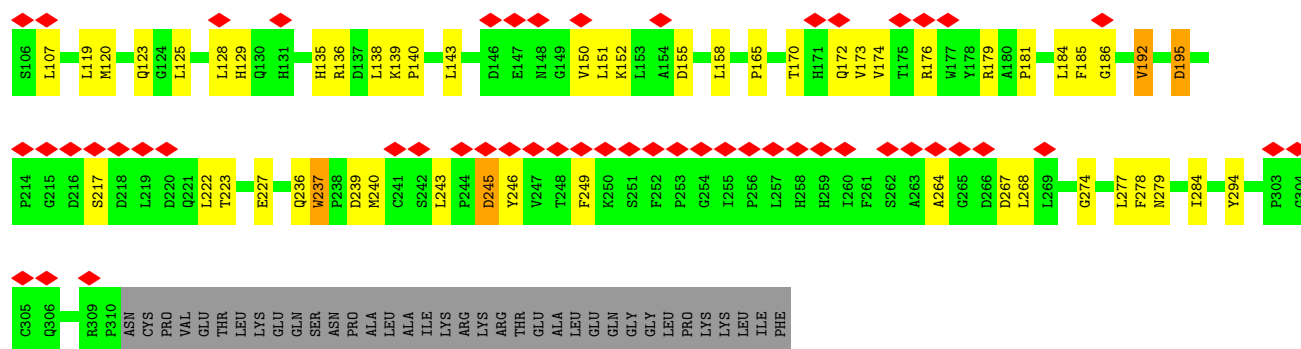


• Molecule 8: General transcription and DNA repair factor IIH helicase subunit XPD

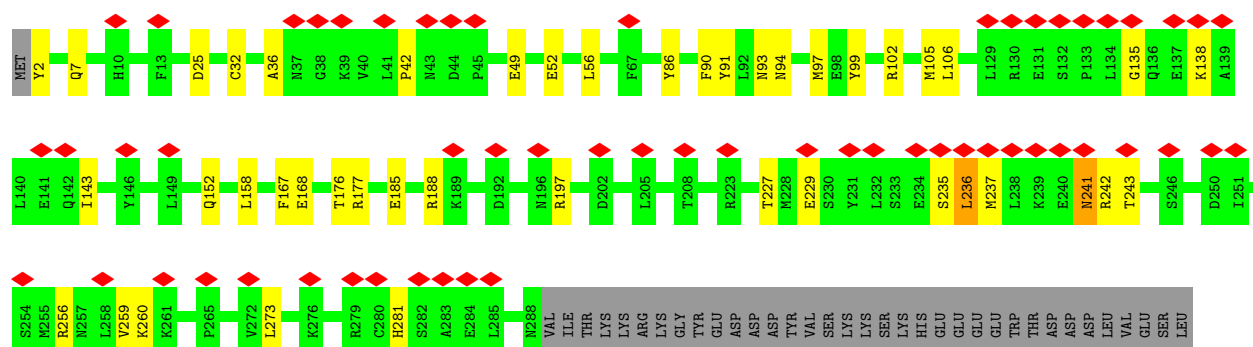
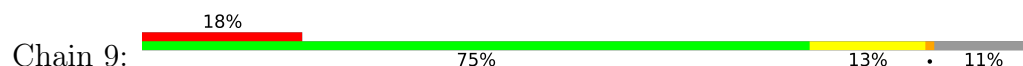


• Molecule 9: Cyclin-dependent kinase 7

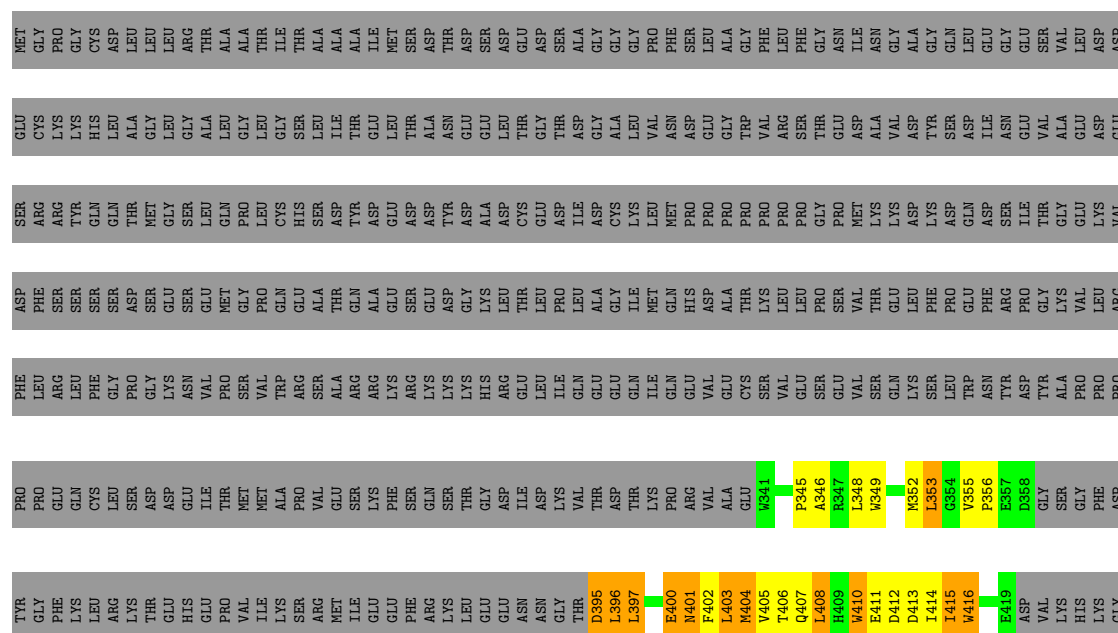




• Molecule 10: Cyclin-H



• Molecule 11: Transcription initiation factor TFIID subunit 1



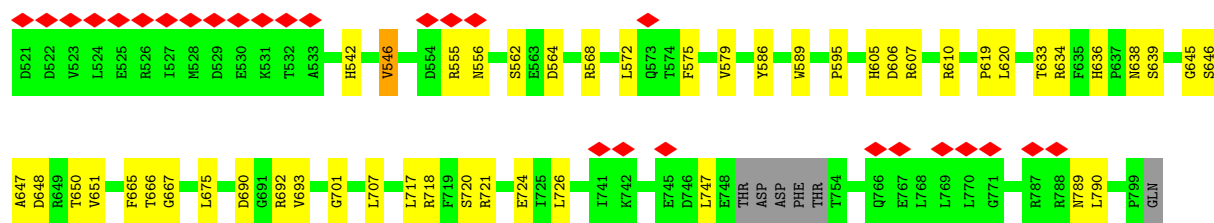


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Chain d: 7% 12% 85%

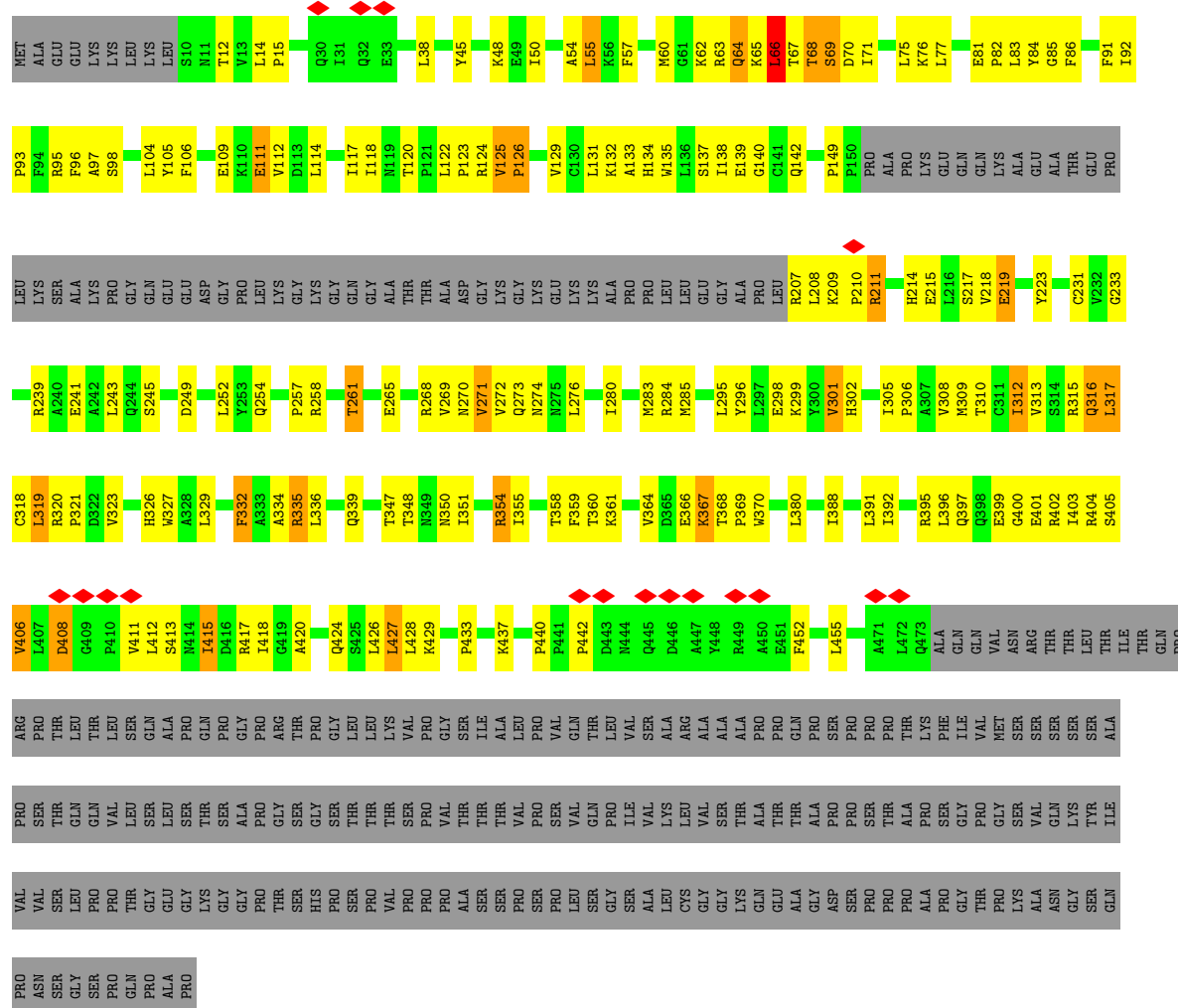
[illegible]

- Molecule 14: Transcription initiation factor TFIID subunit 5



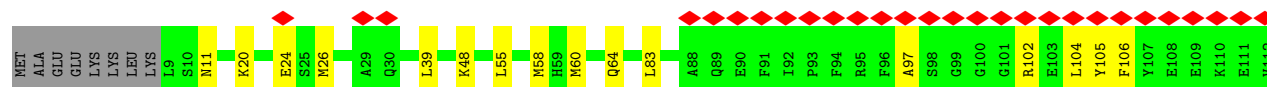
• Molecule 15: Transcription initiation factor TFIID subunit 6

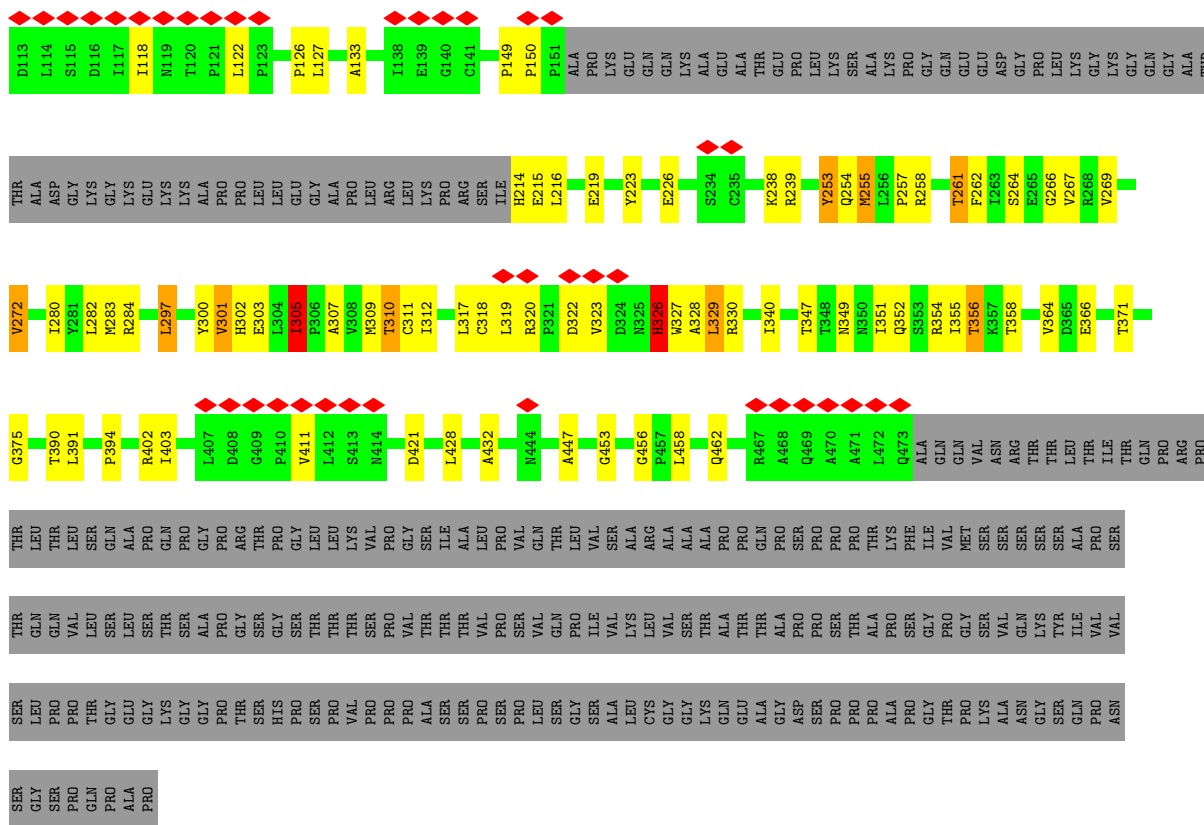
Chain F: 34% 23% 40%



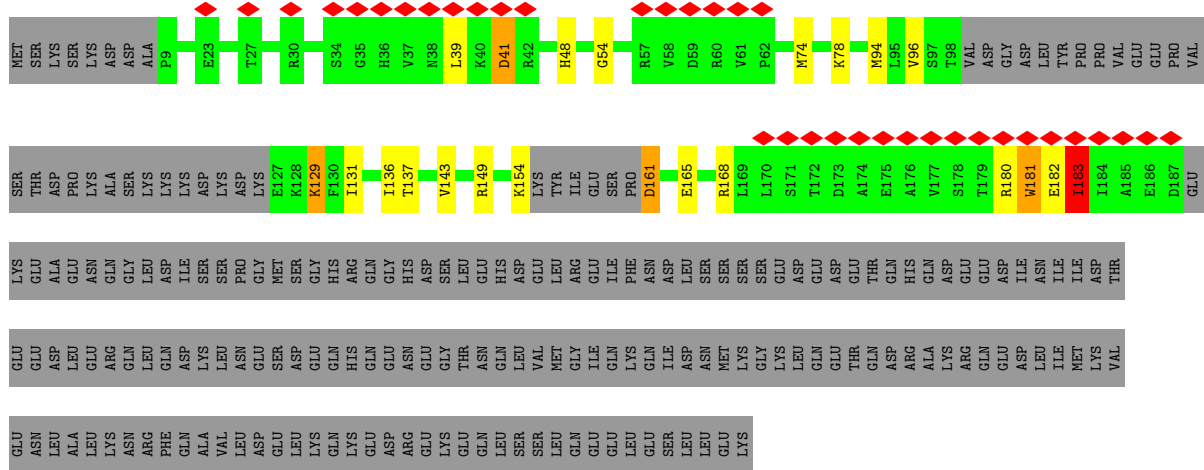
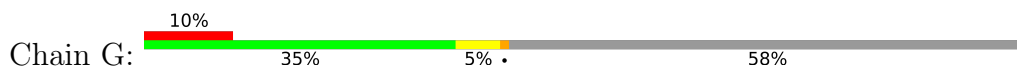
• Molecule 15: Transcription initiation factor TFIID subunit 6

Chain f: 10% 45% 13% 40%

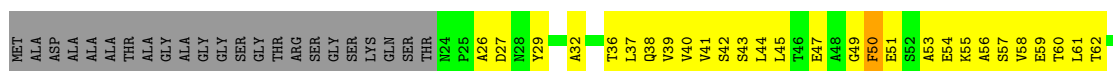


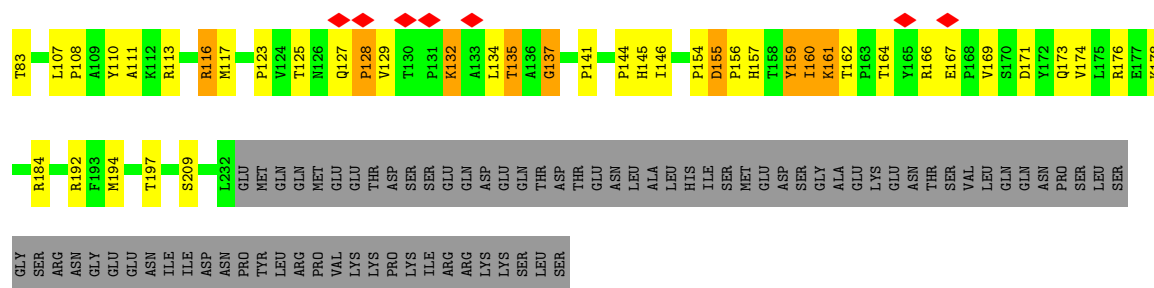


• Molecule 16: Transcription initiation factor TFIID subunit 7

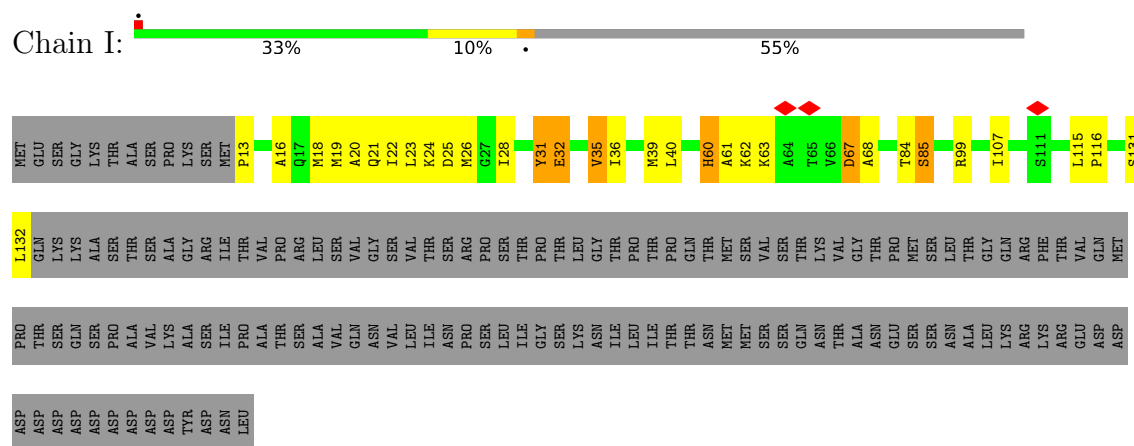


• Molecule 17: Transcription initiation factor TFIID subunit 8

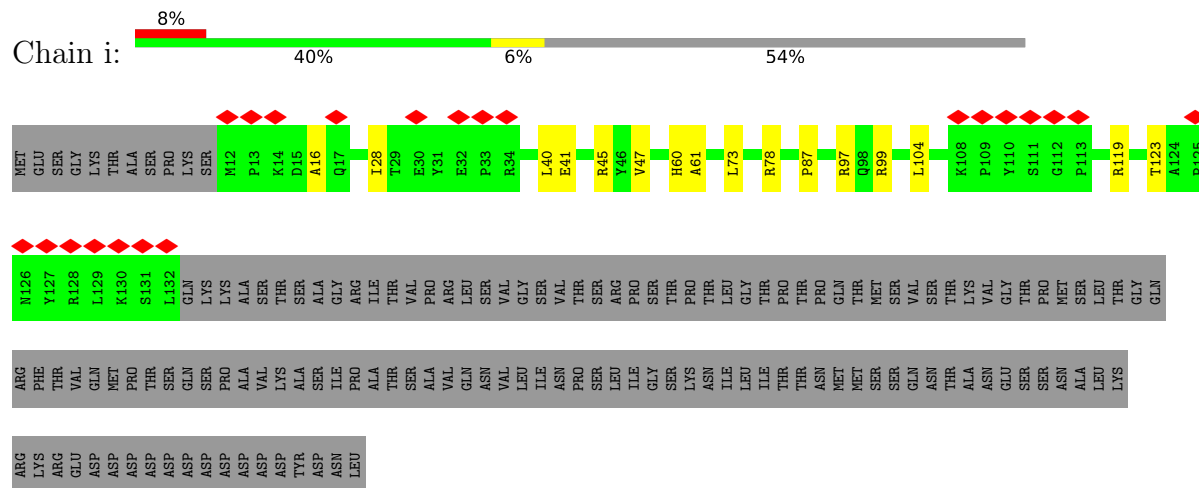




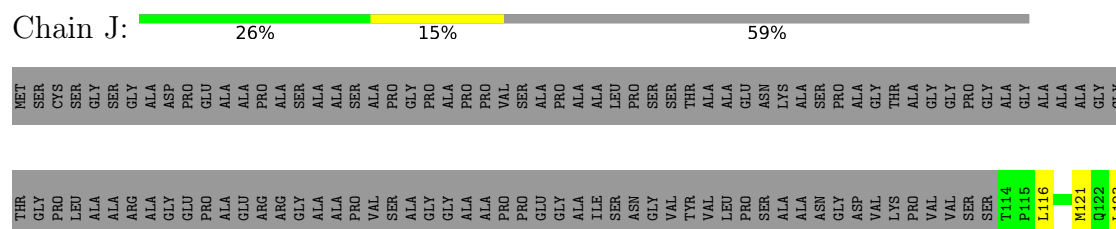
• Molecule 18: Transcription initiation factor TFIID subunit 9

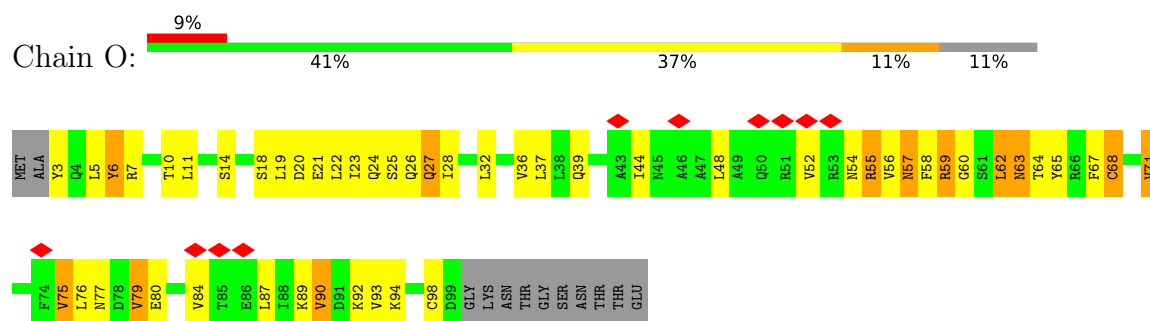


• Molecule 18: Transcription initiation factor TFIID subunit 9

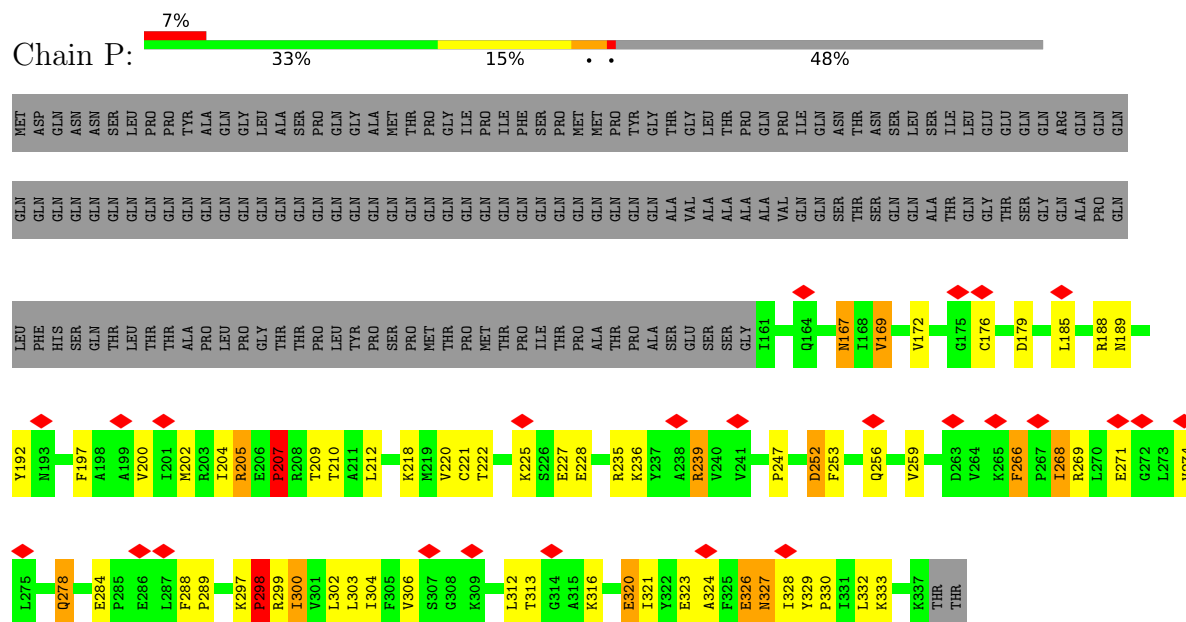


• Molecule 19: Transcription initiation factor TFIID subunit 10

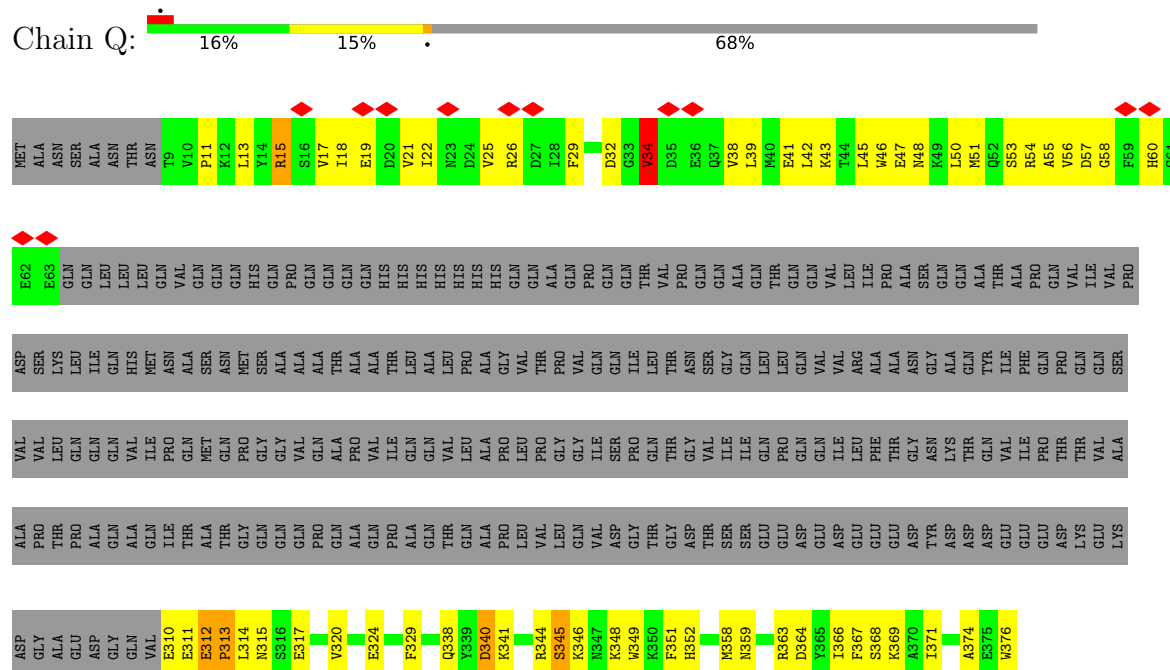




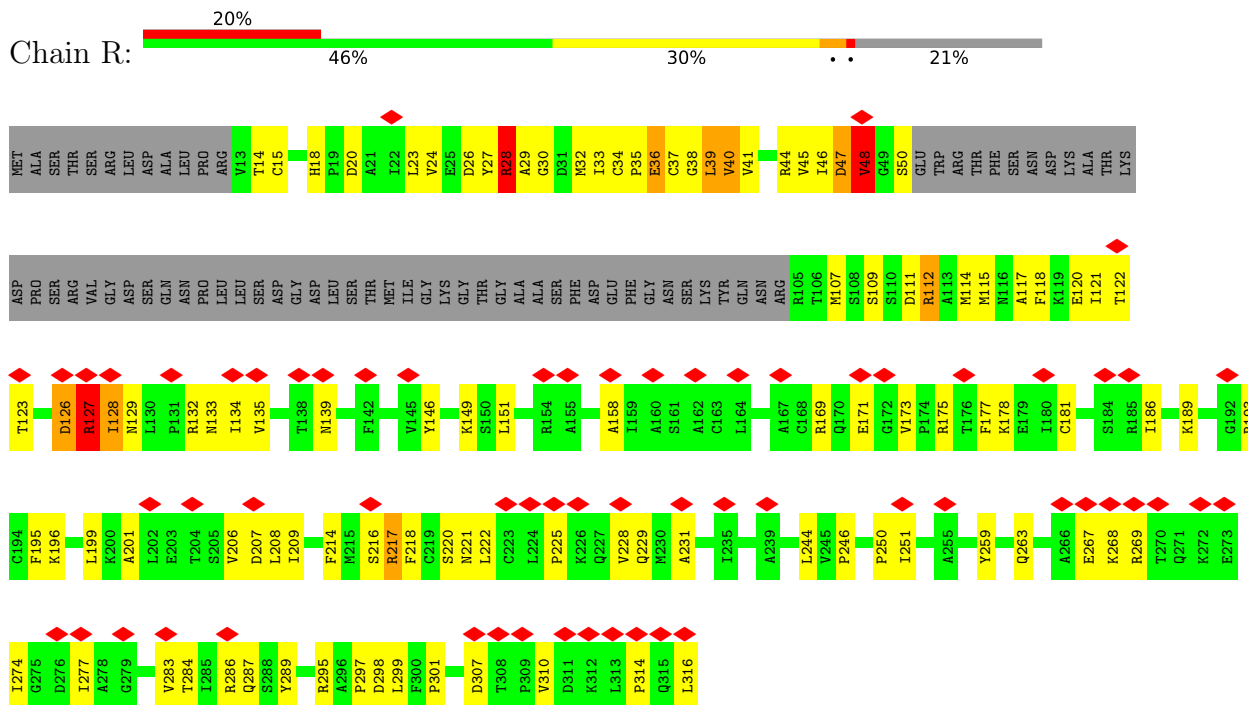
- Molecule 22: TATA-box-binding protein



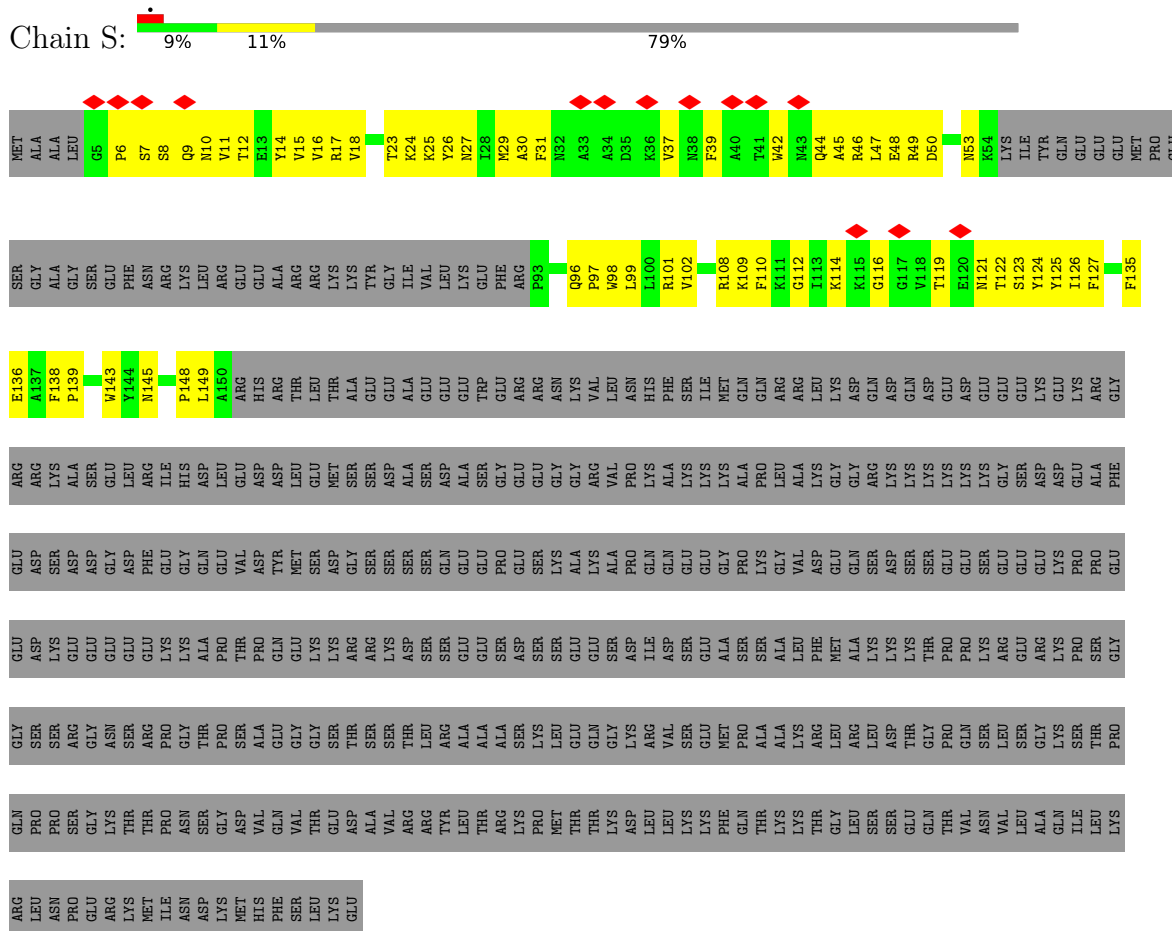
- Molecule 23: Transcription initiation factor IIA subunit 1



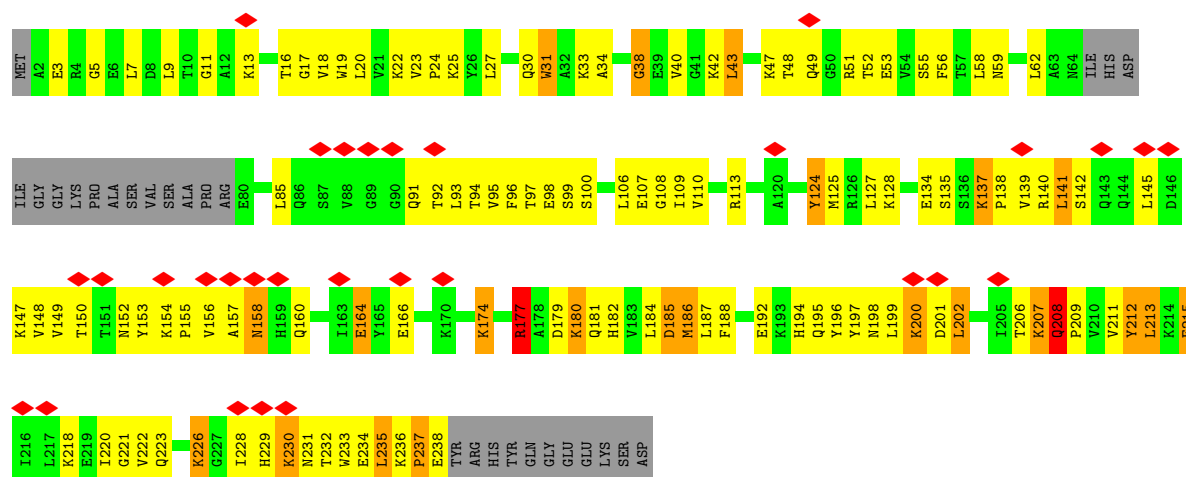
Chain R:



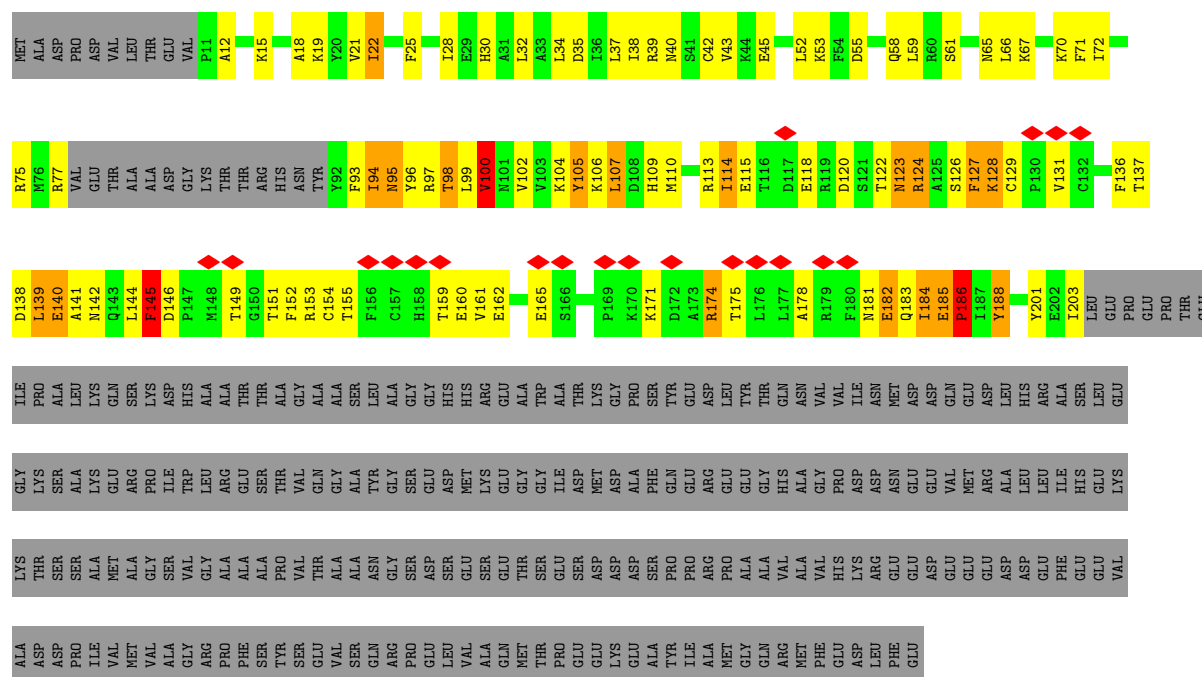
Chain S:



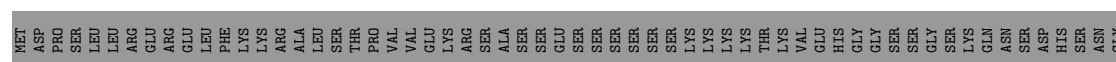
Chain T:

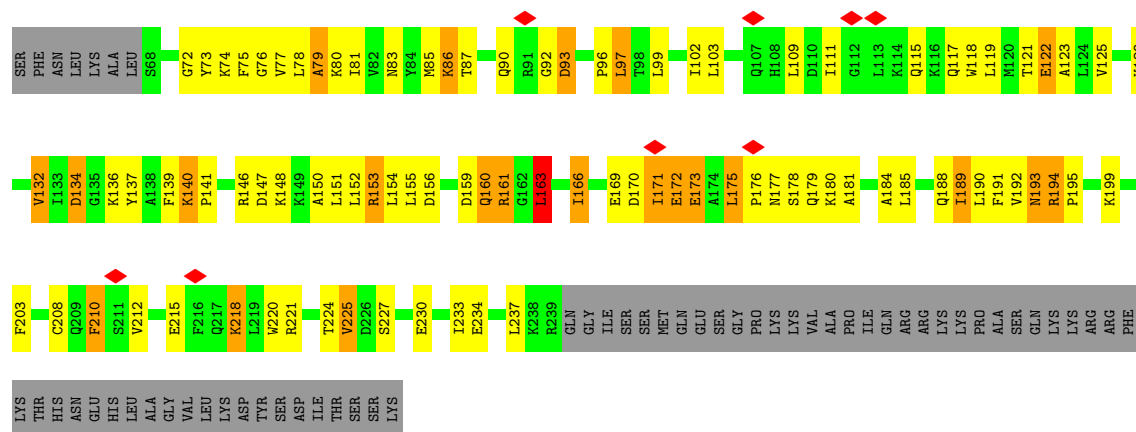


Chain U:

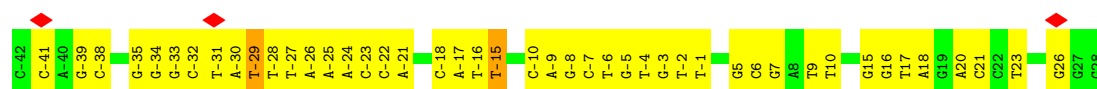
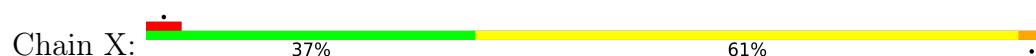


Chain V:





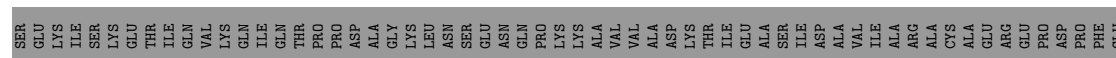
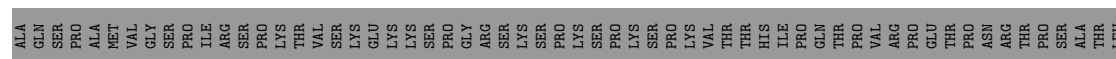
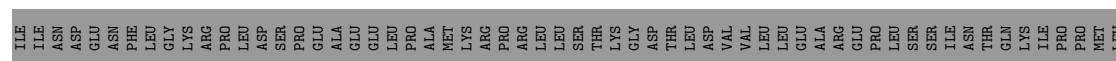
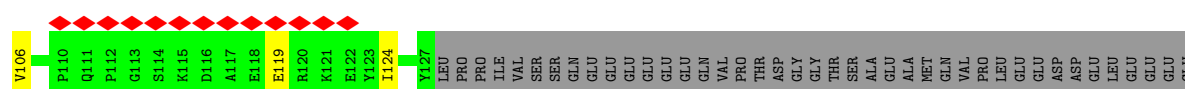
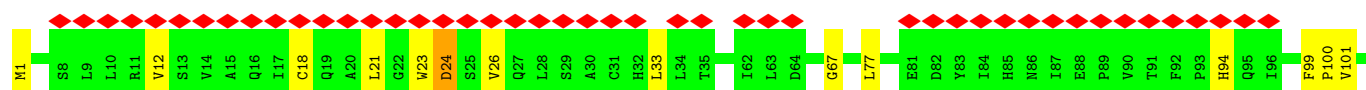
• Molecule 29: DNA (71-MER)



• Molecule 30: DNA (71-MER)



• Molecule 31: Transcription initiation factor TFIID subunit 3

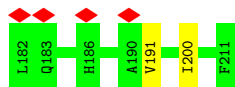


[illegible]

- Molecule 32: Transcription initiation factor TFIID subunit 11



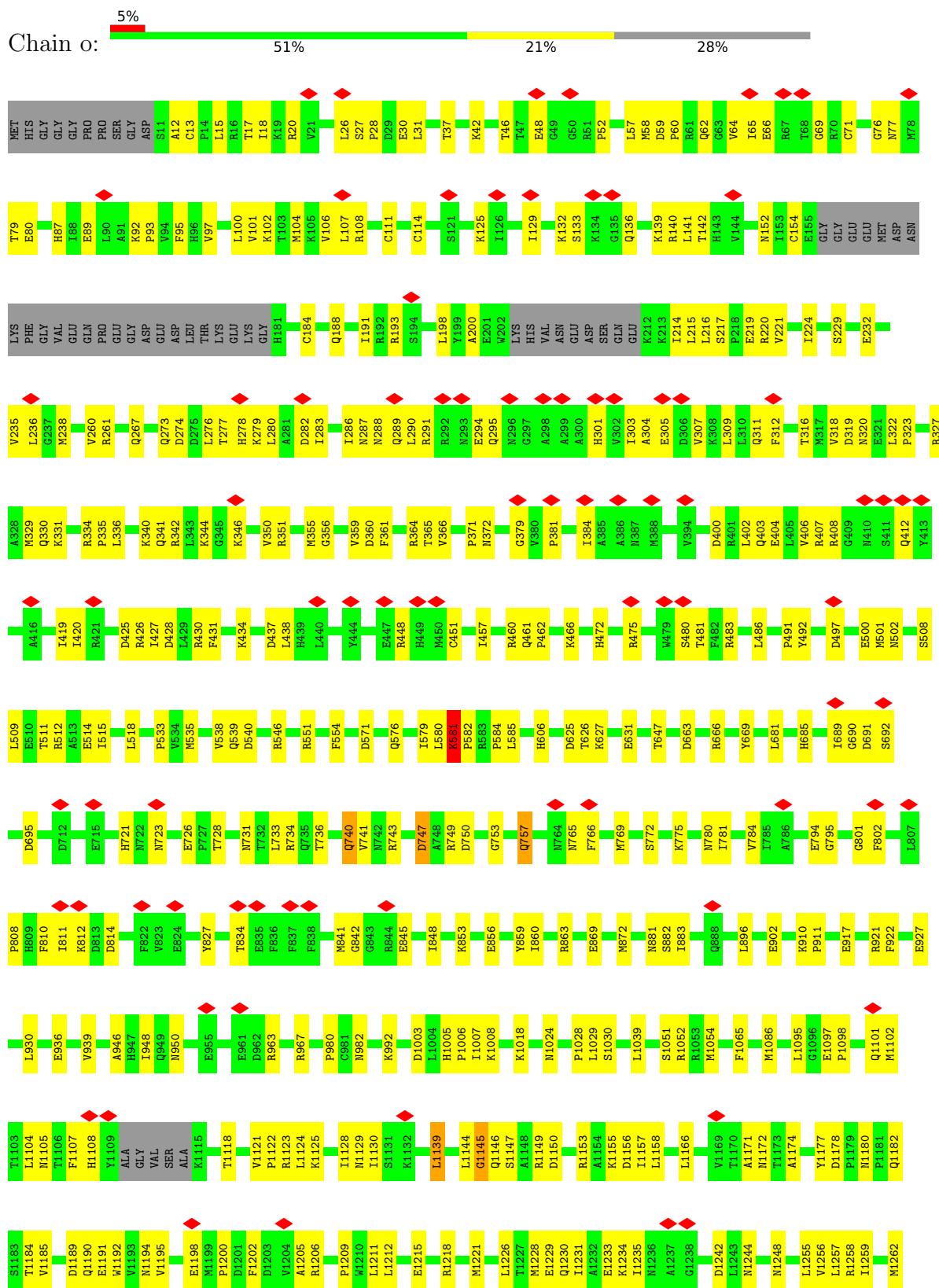
Y121	Met	GLN	ASP
E122	Val	ASP	ASP
M123	Ser	Val	ALA
Y124	Asp	ASP	HIS
R125	Leu	THR	GLU
R126	Thr	THR	SER
S127	Val	THR	PRO
A128	Glu	GLU	ASP
F129	Arg	ARG	LYS
P130	Asp	ASP	GLY
K131	Ser	SER	GLU
A132	Ser	SER	THR
A133	Leu	LEU	GLY
I134	Asn	ASN	SER
K135	Pro	PRO	ASP
R136	Ala	ALA	GLU
L137	Ala	ALA	THR
I138	Lys	LYS	ALA
Q139	Leu	LEU	VAL
T142	Lys	LYS	PRO
G143	Ile	ILE	GLY
T144	Thr	THR	ASP
S145	Lys	LYS	PRO
V146	Glu	GLU	GLY
Q148	Lys	LYS	ALA
N149	Lys	LYS	THR
V150	Gln	GLN	ASP
V151	Lys	LYS	GLY
I152	Val	VAL	ILE
A153	Asp	ASP	PRO
M154	Glu	GLU	GLU
S155	Ile	ILE	THR
G156	Gln	GLN	ASP
I157	Lys	LYS	GLY
S158	Met	MET	ASP
K159	Gln	GLN	ASP
V160	Ile	ILE	LEU
F161	Leu	LEU	LYS
V162	Val	VAL	GLU
G163	Ser	SER	ALA
E164	Ser	SER	ALA
V165	Phe	PHE	ALA
V166	S114	S114	GLU
E167	E115	E115	GLY
S168	E116	E116	GLY
	E117	E117	LEU
	L118	L118	SER

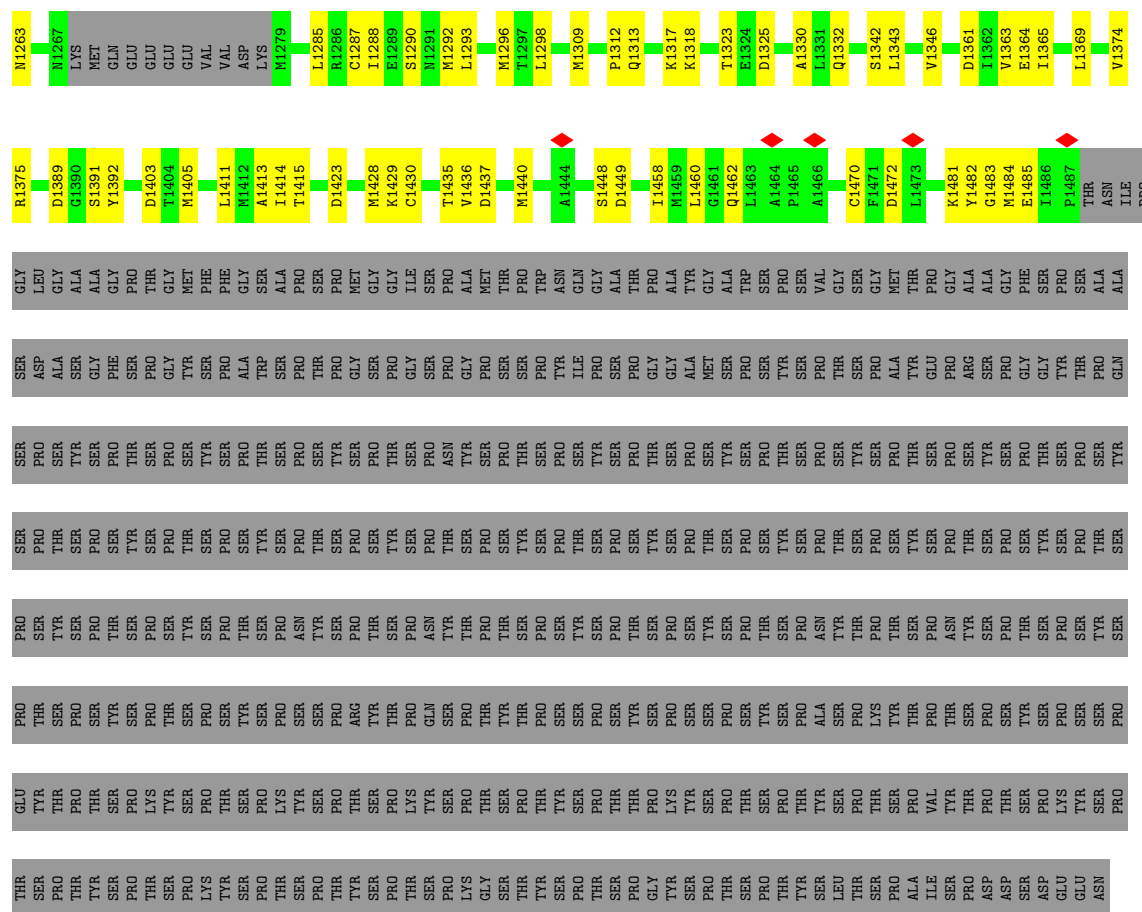


- Molecule 33: Transcription initiation factor TFIID subunit 13

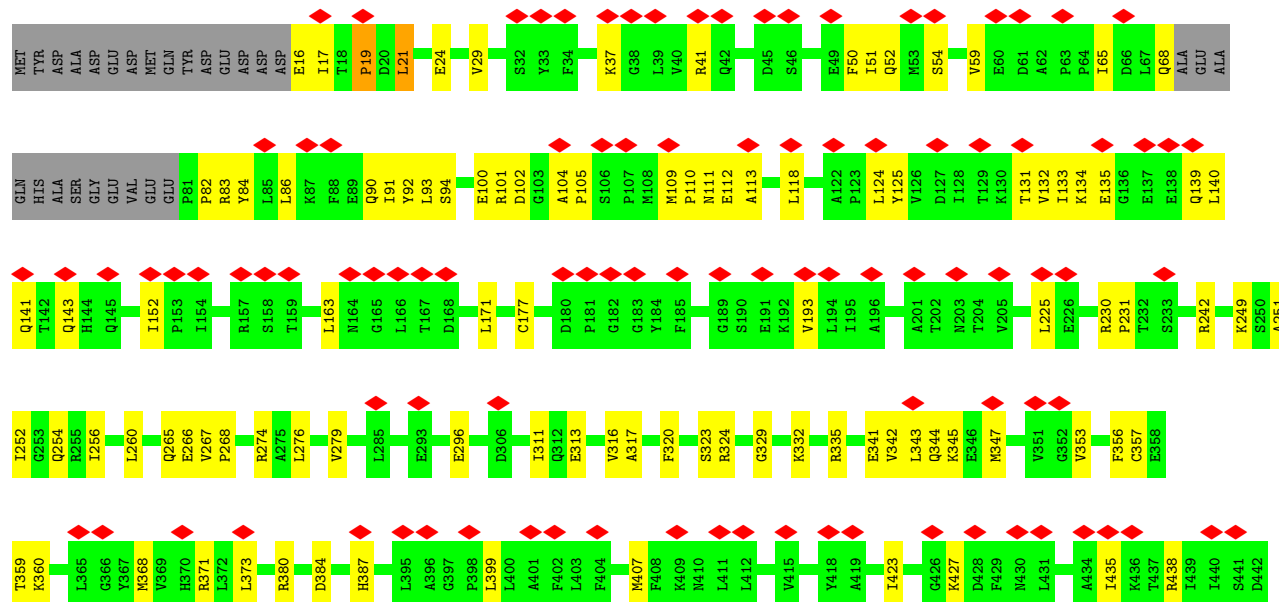
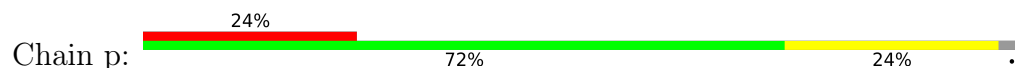
[illegible]

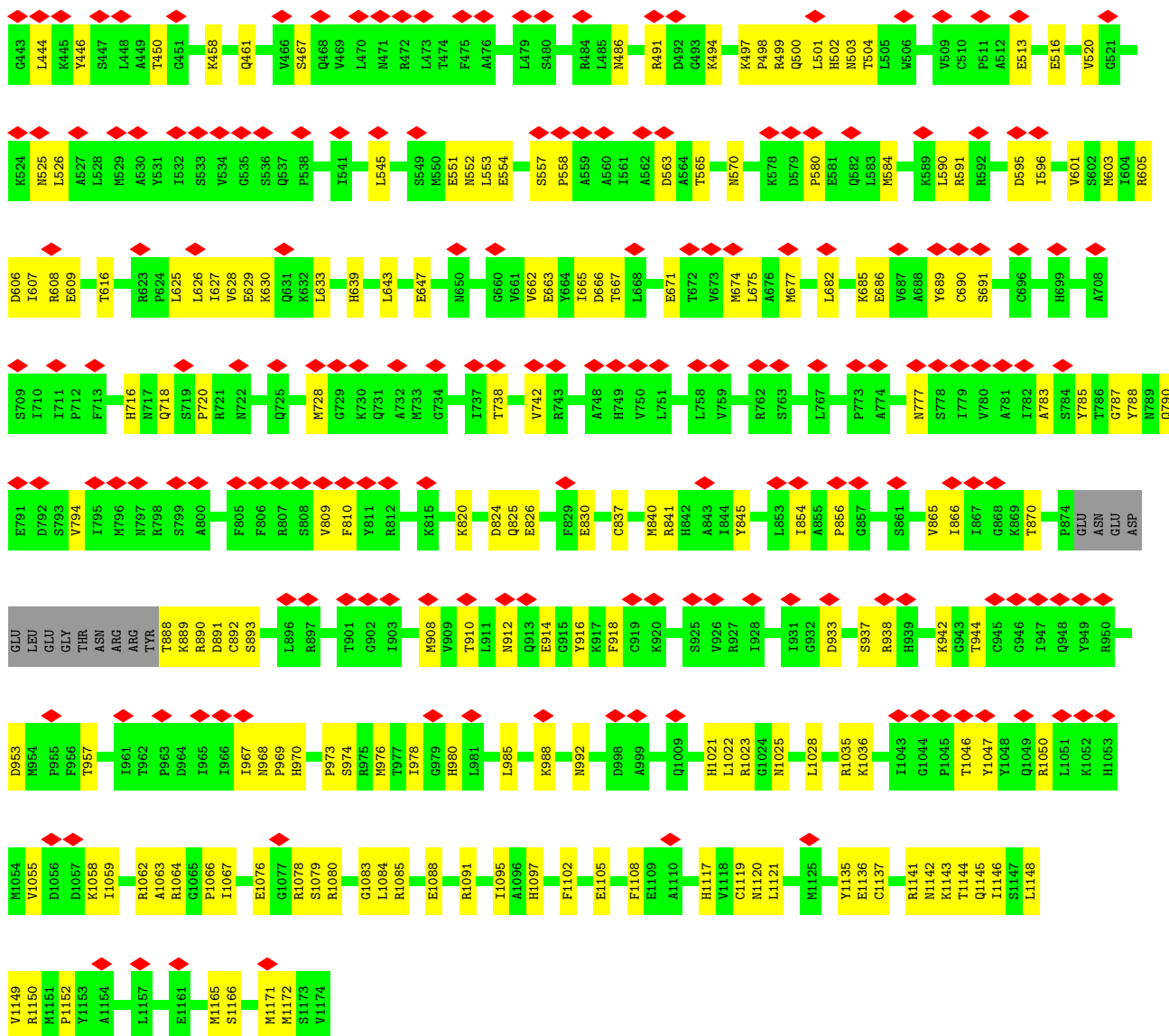
● Molecule 34: DNA-directed RNA polymerase II subunit RPB1



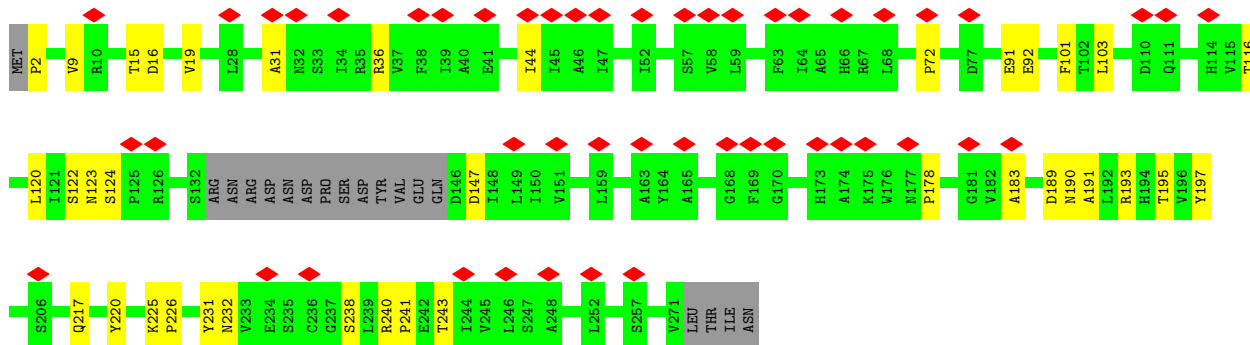
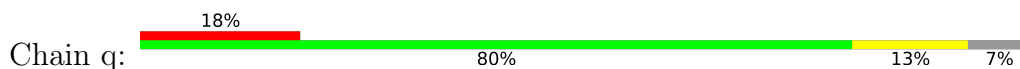


• Molecule 35: DNA-directed RNA polymerase subunit beta

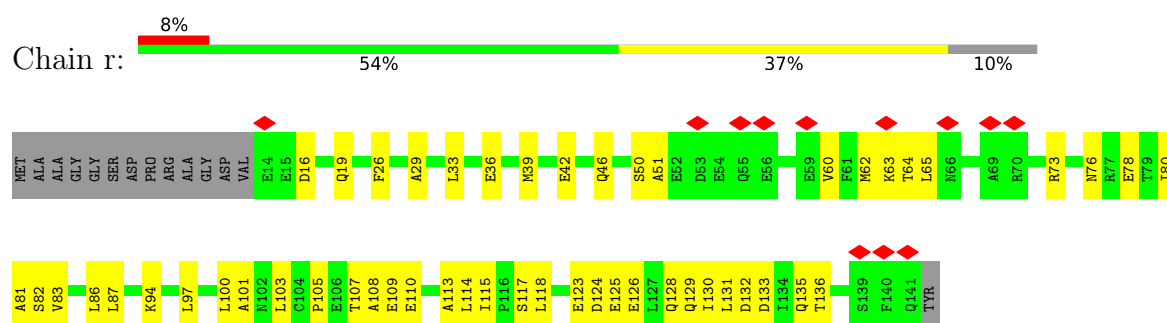




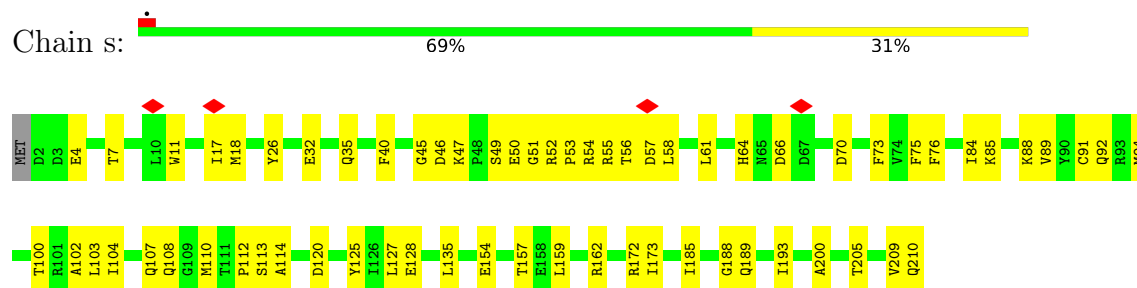
• Molecule 36: DNA-directed RNA polymerase II subunit RPB3



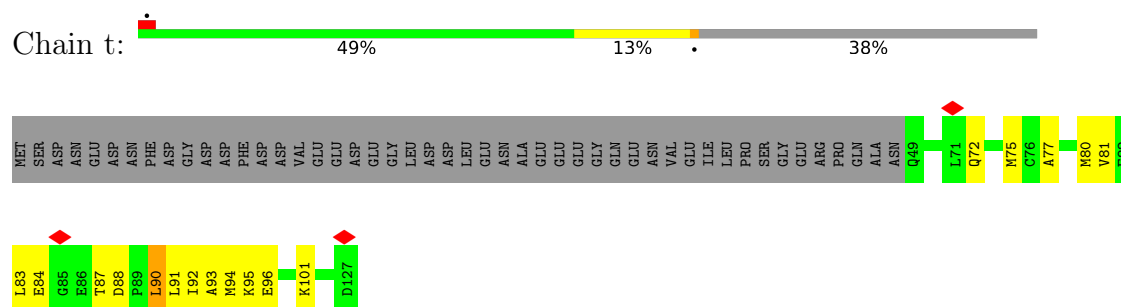
• Molecule 37: DNA-directed RNA polymerase II subunit RPB4



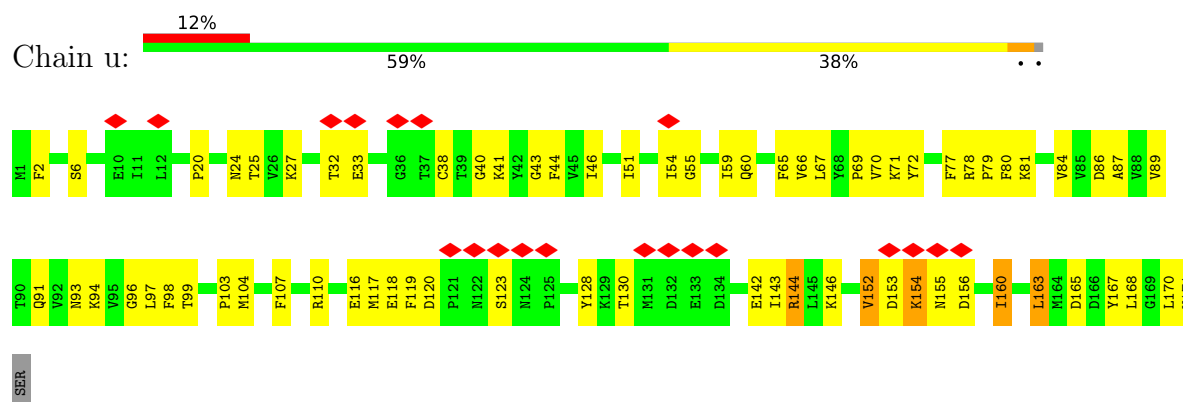
- Molecule 38: DNA-directed RNA polymerase II subunit E



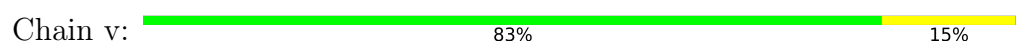
- Molecule 39: DNA-directed RNA polymerase II subunit F

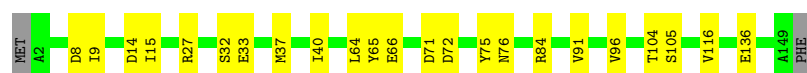


- Molecule 40: DNA-directed RNA polymerase II subunit RPB7

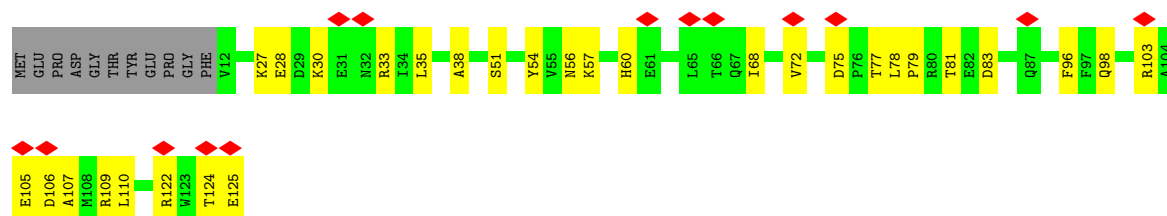


- Molecule 41: DNA-directed RNA polymerases I, II, and III subunit RPABC3

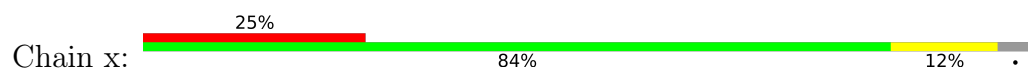




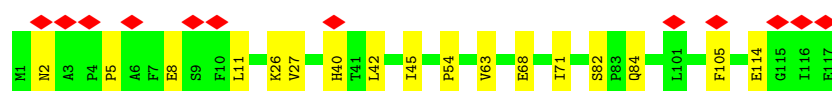
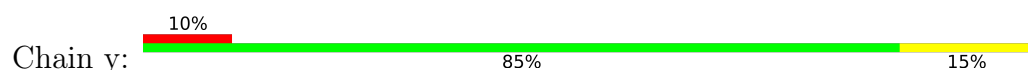
- Molecule 42: DNA-directed RNA polymerase II subunit RPB9



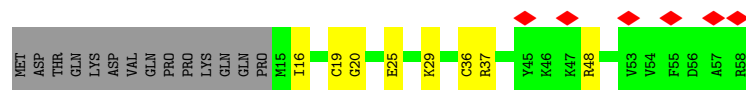
- Molecule 43: RPB10



- Molecule 44: RNA_pol_L_2 domain-containing protein



- Molecule 45: RPB12



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	42939	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50.0	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	0.212	Depositor
Minimum map value	-0.116	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0051	Depositor
Map size ($\text{\AA}$)	542.72, 542.72, 542.72	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.06, 1.06, 1.06	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, MG, ZN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	0	0.35	0/1804	0.68	0/2443
2	1	0.58	1/2674 (0.0%)	1.00	13/3660 (0.4%)
3	2	0.50	0/2588	0.78	6/3509 (0.2%)
4	3	0.46	0/2102	0.73	3/2844 (0.1%)
5	4	0.50	0/3663	0.81	5/4965 (0.1%)
6	5	0.45	0/433	0.73	0/585
7	6	0.58	1/4983 (0.0%)	0.88	15/6731 (0.2%)
8	7	0.49	0/5957	0.81	6/8071 (0.1%)
9	8	0.54	0/2429	0.78	0/3295
10	9	0.49	0/2384	0.69	2/3220 (0.1%)
11	A	0.61	0/4627	0.92	13/6248 (0.2%)
12	B	0.49	1/7993 (0.0%)	0.70	9/10836 (0.1%)
13	D	0.52	0/1335	0.82	2/1784 (0.1%)
13	d	0.21	0/1321	0.44	0/1772
14	E	0.38	0/4469	0.64	5/6050 (0.1%)
14	e	0.29	0/4433	0.52	0/6004
15	F	0.68	1/3167 (0.0%)	1.07	7/4303 (0.2%)
15	f	0.51	0/3140	0.86	8/4268 (0.2%)
16	G	0.70	0/1199	0.94	1/1612 (0.1%)
17	H	0.44	0/1673	0.68	1/2285 (0.0%)
18	I	0.74	0/981	0.98	0/1332
18	i	0.21	0/989	0.40	0/1343
19	J	0.23	0/736	0.44	0/998
19	j	0.22	0/775	0.45	0/1049
20	L	0.59	0/622	0.99	1/841 (0.1%)
20	l	0.44	0/888	0.71	0/1194
21	O	0.77	0/781	1.20	5/1061 (0.5%)
22	P	1.02	3/1438 (0.2%)	1.37	12/1935 (0.6%)
23	Q	0.59	0/1013	1.07	6/1366 (0.4%)
24	R	0.50	0/1957	1.11	10/2643 (0.4%)
25	S	0.35	0/896	0.78	0/1213
26	T	0.76	0/1817	1.26	11/2445 (0.4%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
27	U	0.81	2/1499 (0.1%)	1.39	16/2012 (0.8%)
28	V	0.69	0/1428	1.10	4/1917 (0.2%)
29	X	0.45	0/1641	0.75	4/2534 (0.2%)
30	Y	0.46	0/1623	0.72	4/2499 (0.2%)
31	c	0.48	0/1035	0.67	0/1406
32	k	0.29	0/799	0.53	1/1070 (0.1%)
33	m	0.30	0/733	0.55	0/977
34	o	0.35	0/11516	0.58	4/15548 (0.0%)
35	p	0.33	0/9243	0.51	2/12475 (0.0%)
36	q	0.32	0/2102	0.44	0/2857
37	r	0.18	0/1019	0.45	0/1374
38	s	0.25	0/1751	0.52	1/2366 (0.0%)
39	t	0.63	0/645	0.92	1/871 (0.1%)
40	u	0.42	0/1382	0.82	3/1874 (0.2%)
41	v	0.28	0/1207	0.47	0/1628
42	w	0.23	0/948	0.45	0/1284
43	x	0.36	0/516	0.44	0/696
44	y	0.30	0/956	0.47	0/1294
45	z	0.28	0/377	0.43	0/500
All	All	0.49	9/115687 (0.0%)	0.77	181/157087 (0.1%)

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
27	U	183	GLN	C-O	8.28	1.33	1.24
7	6	432	THR	C-N	7.30	1.43	1.33
22	P	326	GLU	C-O	6.92	1.32	1.24
27	U	124	ARG	C-N	6.86	1.42	1.33
22	P	323	GLU	C-O	6.68	1.32	1.24

The worst 5 of 181 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
27	U	186	PRO	CA-N-CD	-13.20	93.52	112.00
2	1	421	VAL	N-CA-C	-12.30	98.89	110.82
15	F	81	GLU	CA-C-N	10.40	130.60	119.05
15	F	81	GLU	C-N-CA	10.40	130.60	119.05
12	B	491	HIS	N-CA-C	-9.78	100.62	111.28

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	0	1774	0	1601	65	0
2	1	2634	0	2042	285	0
3	2	2534	0	2456	149	0
4	3	2065	0	2098	239	0
5	4	3579	0	3620	239	0
6	5	428	0	434	58	0
7	6	4880	0	4923	542	0
8	7	5833	0	5796	444	0
9	8	2370	0	2391	58	0
10	9	2334	0	2357	60	0
11	A	4511	0	4428	439	0
12	B	7796	0	7751	148	0
13	D	1322	0	1340	110	0
13	d	1307	0	1318	20	0
14	E	4364	0	4244	182	0
14	e	4327	0	4197	43	0
15	F	3109	0	3045	564	0
15	f	3081	0	3012	190	0
16	G	1180	0	1235	19	0
17	H	1633	0	1621	282	0
18	I	959	0	969	77	0
18	i	967	0	977	46	0
19	J	720	0	719	158	0
19	j	759	0	759	18	0
20	L	614	0	614	84	0
20	l	876	0	891	68	0
21	O	771	0	764	169	0
22	P	1412	0	1506	96	0
23	Q	996	0	926	163	0
24	R	1929	0	1970	185	0
25	S	872	0	848	159	0
26	T	1788	0	1819	261	0
27	U	1476	0	1486	160	0
28	V	1404	0	1423	186	0
29	X	1464	0	802	92	0
30	Y	1447	0	795	102	0
31	c	1011	0	975	24	0

Continued on next page...

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
32	k	785	0	818	7	0
33	m	724	0	744	12	0
34	o	11308	0	11428	408	0
35	p	9062	0	9107	314	0
36	q	2059	0	2008	27	0
37	r	1005	0	964	82	0
38	s	1720	0	1737	48	0
39	t	635	0	665	24	0
40	u	1351	0	1358	132	0
41	v	1186	0	1147	15	0
42	w	927	0	859	35	0
43	x	507	0	523	5	0
44	y	937	0	959	14	0
45	z	372	0	379	5	0
46	0	2	0	0	0	0
46	2	3	0	0	0	0
46	3	1	0	0	0	0
46	R	1	0	0	0	0
46	U	1	0	0	0	0
46	o	2	0	0	0	0
46	p	1	0	0	0	0
46	q	1	0	0	0	0
46	w	2	0	0	0	0
46	x	1	0	0	0	0
46	z	1	0	0	0	0
47	7	8	0	0	0	0
48	o	1	0	0	0	0
All	All	113129	0	110848	5455	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 24.

The worst 5 of 5455 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)
7:6:557:LYS:HB3	7:6:596:PHE:CZ	1.28	1.69
14:E:507:VAL:CG2	14:E:513:LEU:HD11	1.25	1.65
8:7:5:VAL:HG11	8:7:55:LEU:CD1	1.17	1.64
15:F:412:LEU:CD1	15:F:417:ARG:HD2	1.21	1.62
11:A:410:TRP:HB3	15:F:306:PRO:CD	1.27	1.62

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	0	230/309 (74%)	199 (86%)	29 (13%)	2 (1%)	14	47
2	1	399/548 (73%)	311 (78%)	70 (18%)	18 (4%)	2	19
3	2	327/395 (83%)	288 (88%)	37 (11%)	2 (1%)	21	55
4	3	259/308 (84%)	243 (94%)	16 (6%)	0	100	100
5	4	447/462 (97%)	399 (89%)	38 (8%)	10 (2%)	5	31
6	5	52/71 (73%)	47 (90%)	4 (8%)	1 (2%)	6	33
7	6	602/782 (77%)	512 (85%)	79 (13%)	11 (2%)	6	34
8	7	732/760 (96%)	649 (89%)	74 (10%)	9 (1%)	10	41
9	8	296/346 (86%)	261 (88%)	33 (11%)	2 (1%)	18	53
10	9	285/323 (88%)	270 (95%)	14 (5%)	1 (0%)	30	64
11	A	534/1872 (28%)	505 (95%)	26 (5%)	3 (1%)	21	55
12	B	959/1199 (80%)	912 (95%)	47 (5%)	0	100	100
13	D	152/1085 (14%)	146 (96%)	5 (3%)	1 (1%)	18	53
13	d	154/1085 (14%)	150 (97%)	4 (3%)	0	100	100
14	E	540/800 (68%)	506 (94%)	33 (6%)	1 (0%)	43	74
14	e	531/800 (66%)	484 (91%)	47 (9%)	0	100	100
15	F	404/677 (60%)	380 (94%)	17 (4%)	7 (2%)	7	35
15	f	399/677 (59%)	378 (95%)	20 (5%)	1 (0%)	36	69
16	G	139/349 (40%)	135 (97%)	4 (3%)	0	100	100
17	H	207/310 (67%)	190 (92%)	13 (6%)	4 (2%)	6	33
18	I	118/264 (45%)	111 (94%)	5 (4%)	2 (2%)	7	35
18	i	119/264 (45%)	115 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
19	J	86/218 (39%)	82 (95%)	4 (5%)	0	100	100
19	j	91/218 (42%)	89 (98%)	2 (2%)	0	100	100
20	L	73/161 (45%)	64 (88%)	4 (6%)	5 (7%)	1	14
20	l	105/161 (65%)	101 (96%)	4 (4%)	0	100	100
21	O	95/109 (87%)	82 (86%)	8 (8%)	5 (5%)	1	17
22	P	175/339 (52%)	168 (96%)	5 (3%)	2 (1%)	11	43
23	Q	118/376 (31%)	106 (90%)	9 (8%)	3 (2%)	4	29
24	R	246/316 (78%)	234 (95%)	10 (4%)	2 (1%)	16	50
25	S	104/517 (20%)	102 (98%)	2 (2%)	0	100	100
26	T	218/249 (88%)	204 (94%)	11 (5%)	3 (1%)	9	38
27	U	175/439 (40%)	165 (94%)	9 (5%)	1 (1%)	21	55
28	V	170/291 (58%)	144 (85%)	17 (10%)	9 (5%)	1	17
31	c	125/929 (14%)	116 (93%)	9 (7%)	0	100	100
32	k	96/211 (46%)	91 (95%)	5 (5%)	0	100	100
33	m	85/124 (68%)	79 (93%)	6 (7%)	0	100	100
34	o	1417/1970 (72%)	1310 (92%)	106 (8%)	1 (0%)	48	80
35	p	1128/1174 (96%)	1054 (93%)	73 (6%)	1 (0%)	48	80
36	q	253/275 (92%)	225 (89%)	28 (11%)	0	100	100
37	r	126/142 (89%)	119 (94%)	7 (6%)	0	100	100
38	s	207/210 (99%)	195 (94%)	12 (6%)	0	100	100
39	t	77/127 (61%)	73 (95%)	3 (4%)	1 (1%)	9	40
40	u	169/172 (98%)	156 (92%)	12 (7%)	1 (1%)	21	55
41	v	146/150 (97%)	132 (90%)	14 (10%)	0	100	100
42	w	112/125 (90%)	103 (92%)	9 (8%)	0	100	100
43	x	62/67 (92%)	59 (95%)	3 (5%)	0	100	100
44	y	115/117 (98%)	109 (95%)	6 (5%)	0	100	100
45	z	42/58 (72%)	38 (90%)	4 (10%)	0	100	100
All	All	13701/22931 (60%)	12591 (92%)	1001 (7%)	109 (1%)	18	50

5 of 109 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	1	106	LYS

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Mol	Chain	Res	Type
2	1	108	ASN
2	1	110	GLU
2	1	114	LYS
2	1	120	GLU

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	0	173/283 (61%)	171 (99%)	2 (1%)	63	72
2	1	176/484 (36%)	168 (96%)	8 (4%)	24	49
3	2	279/352 (79%)	272 (98%)	7 (2%)	42	62
4	3	234/272 (86%)	229 (98%)	5 (2%)	47	65
5	4	387/399 (97%)	378 (98%)	9 (2%)	44	63
6	5	48/64 (75%)	46 (96%)	2 (4%)	26	50
7	6	533/688 (78%)	479 (90%)	54 (10%)	7	27
8	7	616/664 (93%)	598 (97%)	18 (3%)	37	58
9	8	258/299 (86%)	251 (97%)	7 (3%)	39	60
10	9	259/296 (88%)	257 (99%)	2 (1%)	73	77
11	A	488/1665 (29%)	443 (91%)	45 (9%)	8	30
12	B	876/1083 (81%)	859 (98%)	17 (2%)	50	66
13	D	143/815 (18%)	135 (94%)	8 (6%)	19	45
13	d	146/815 (18%)	146 (100%)	0	100	100
14	E	478/657 (73%)	467 (98%)	11 (2%)	44	63
14	e	475/657 (72%)	473 (100%)	2 (0%)	84	83
15	F	324/574 (56%)	294 (91%)	30 (9%)	8	30
15	f	322/574 (56%)	309 (96%)	13 (4%)	28	51
16	G	133/322 (41%)	125 (94%)	8 (6%)	17	44
17	H	181/270 (67%)	170 (94%)	11 (6%)	17	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
18	I	106/235 (45%)	100 (94%)	6 (6%)	18	44
18	i	107/235 (46%)	107 (100%)	0	100	100
19	J	79/154 (51%)	77 (98%)	2 (2%)	42	62
19	j	83/154 (54%)	83 (100%)	0	100	100
20	L	70/141 (50%)	59 (84%)	11 (16%)	2	15
20	l	98/141 (70%)	97 (99%)	1 (1%)	68	75
21	O	84/98 (86%)	76 (90%)	8 (10%)	8	29
22	P	153/293 (52%)	140 (92%)	13 (8%)	10	33
23	Q	111/324 (34%)	106 (96%)	5 (4%)	24	49
24	R	212/268 (79%)	200 (94%)	12 (6%)	18	44
25	S	93/448 (21%)	92 (99%)	1 (1%)	65	74
26	T	196/218 (90%)	166 (85%)	30 (15%)	3	15
27	U	163/373 (44%)	144 (88%)	19 (12%)	5	22
28	V	155/261 (59%)	135 (87%)	20 (13%)	4	20
31	c	113/833 (14%)	111 (98%)	2 (2%)	51	67
32	k	87/182 (48%)	87 (100%)	0	100	100
33	m	80/106 (76%)	79 (99%)	1 (1%)	61	71
34	o	1257/1748 (72%)	1247 (99%)	10 (1%)	73	77
35	p	993/1027 (97%)	991 (100%)	2 (0%)	87	87
36	q	234/252 (93%)	234 (100%)	0	100	100
37	r	106/126 (84%)	106 (100%)	0	100	100
38	s	191/192 (100%)	191 (100%)	0	100	100
39	t	69/111 (62%)	68 (99%)	1 (1%)	59	71
40	u	152/153 (99%)	150 (99%)	2 (1%)	61	71
41	v	129/131 (98%)	129 (100%)	0	100	100
42	w	103/112 (92%)	103 (100%)	0	100	100
43	x	53/56 (95%)	53 (100%)	0	100	100
44	y	106/106 (100%)	106 (100%)	0	100	100
45	z	41/55 (74%)	41 (100%)	0	100	100
All	All	11953/19766 (60%)	11548 (97%)	405 (3%)	33	55

5 of 405 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
16	G	182	GLU
23	Q	320	VAL
34	o	1423	ASP
17	H	159	TYR
20	L	110	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 236 such sidechains are listed below:

Mol	Chain	Res	Type
15	F	316	GLN
35	p	1117	HIS
24	R	229	GLN
35	p	1021	HIS
34	o	1005	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 17 are monoatomic - leaving 1 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
47	SF4	7	1000	8	0,12,12	-	-	-		

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
47	SF4	7	1000	8	-	-	0/6/5/5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

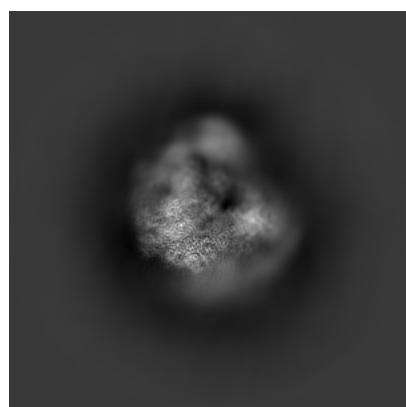
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31112. These allow visual inspection of the internal detail of the map and identification of artifacts.

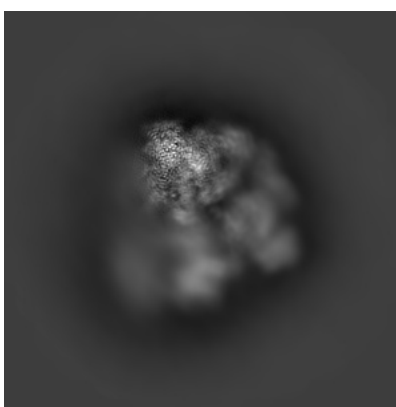
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

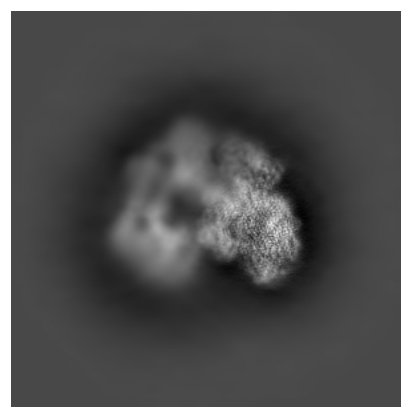
6.1.1 Primary map



X



Y

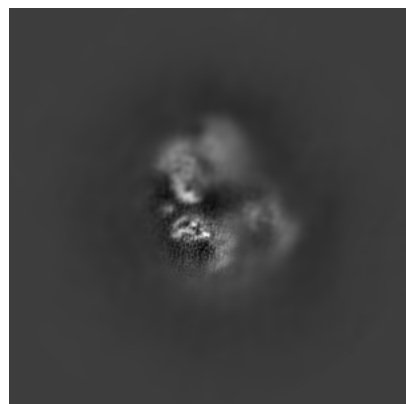


Z

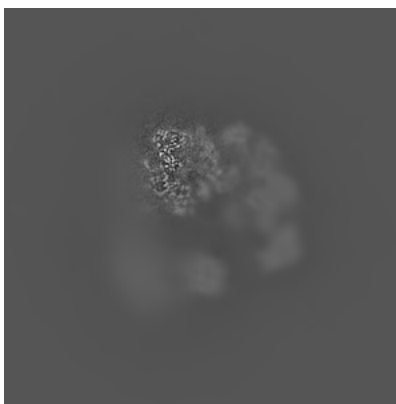
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

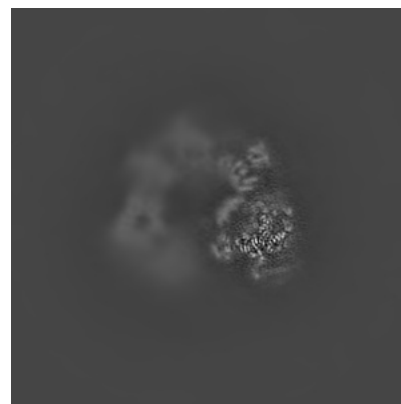
6.2.1 Primary map



X Index: 256



Y Index: 256

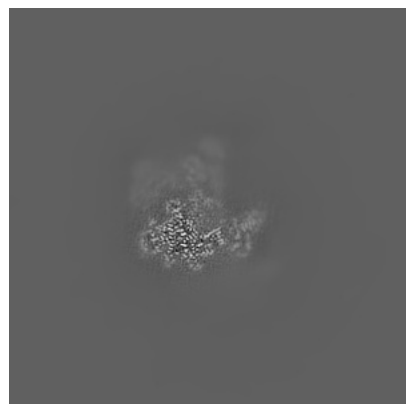


Z Index: 256

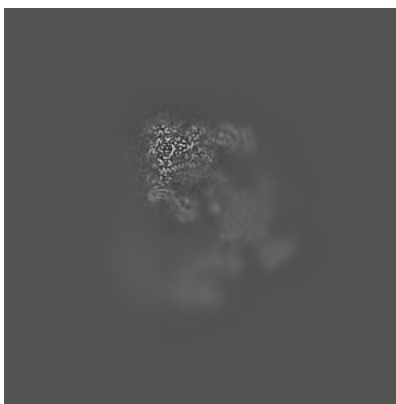
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

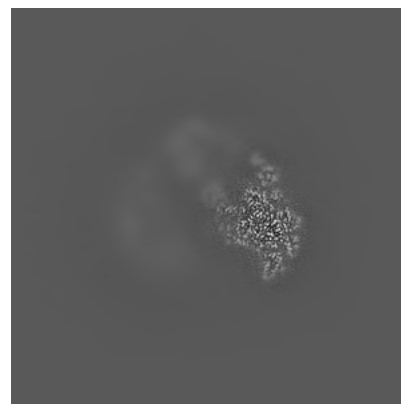
6.3.1 Primary map



X Index: 334



Y Index: 228

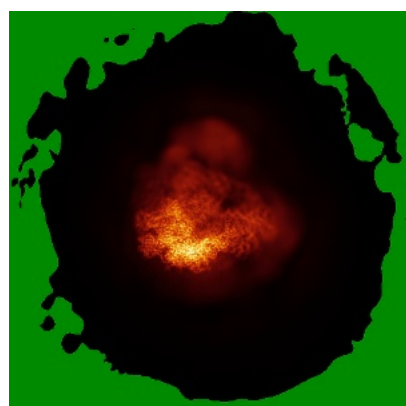


Z Index: 209

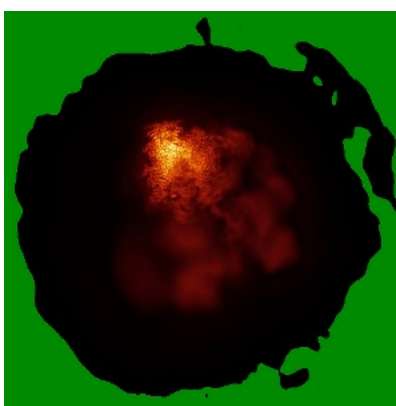
The images above show the largest variance slices of the map in three orthogonal directions.

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

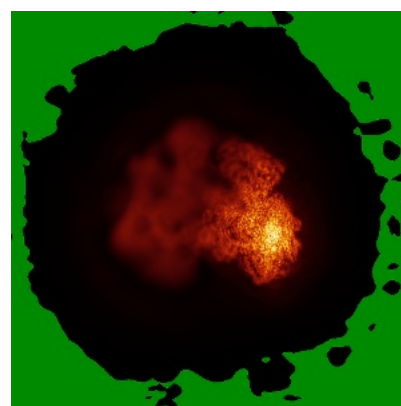
6.4.1 Primary map



X



Y

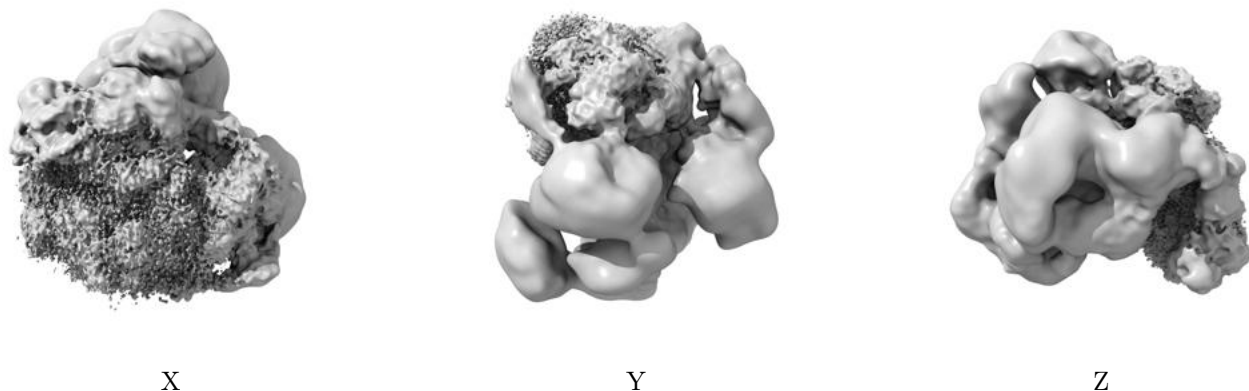


Z

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 0.0051. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

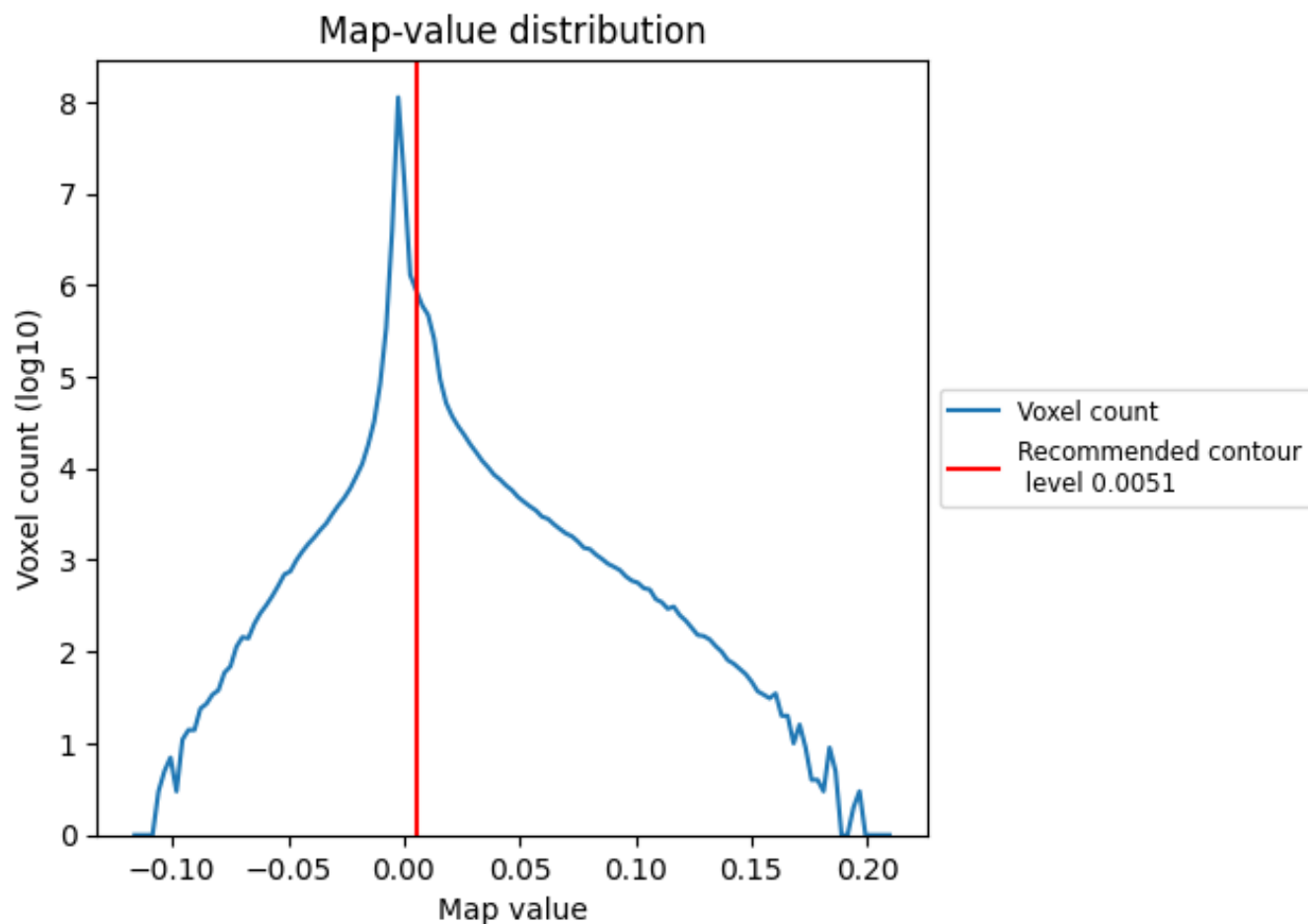
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

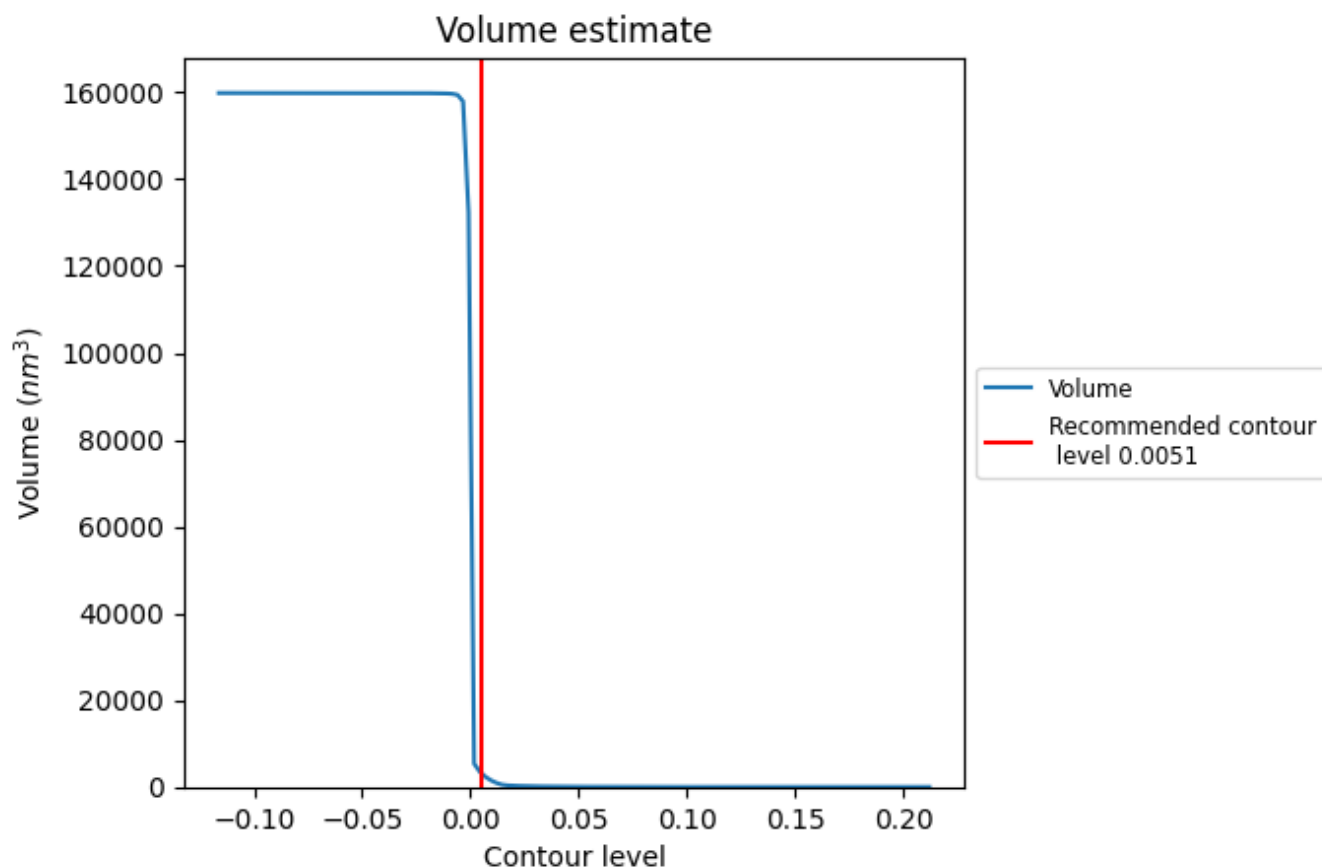
This section contains the results of statistical analysis of the map.

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.

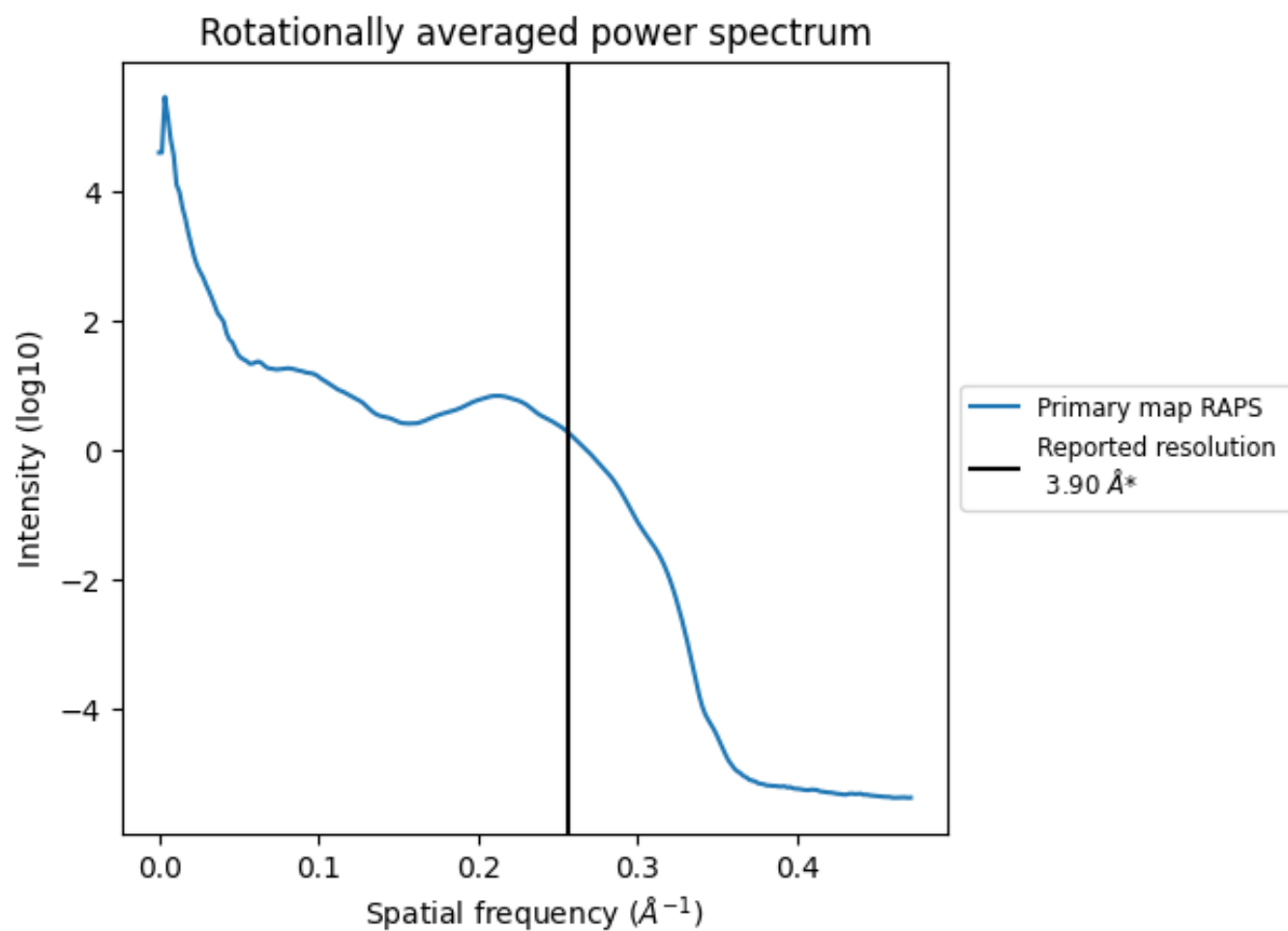
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 3205 nm^3 ; this corresponds to an approximate mass of 2895 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 0.256 Å⁻¹

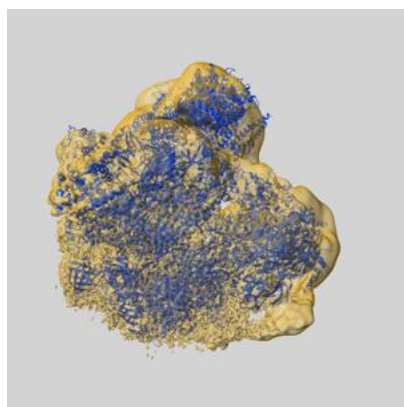
8 Fourier-Shell correlation

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

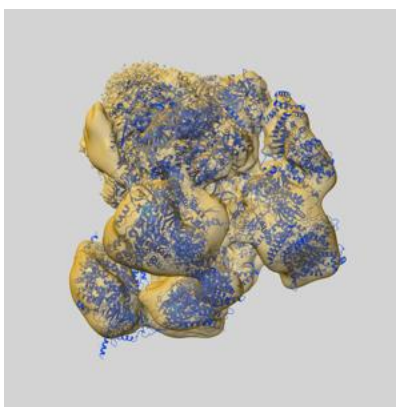
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-31112 and PDB model 7EGC. Per-residue inclusion information can be found in section [3](#) on page [14](#).

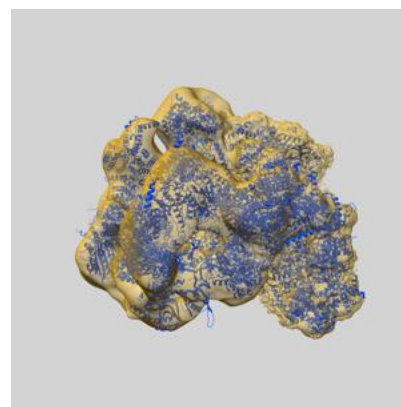
9.1 Map-model overlay [i](#)



X



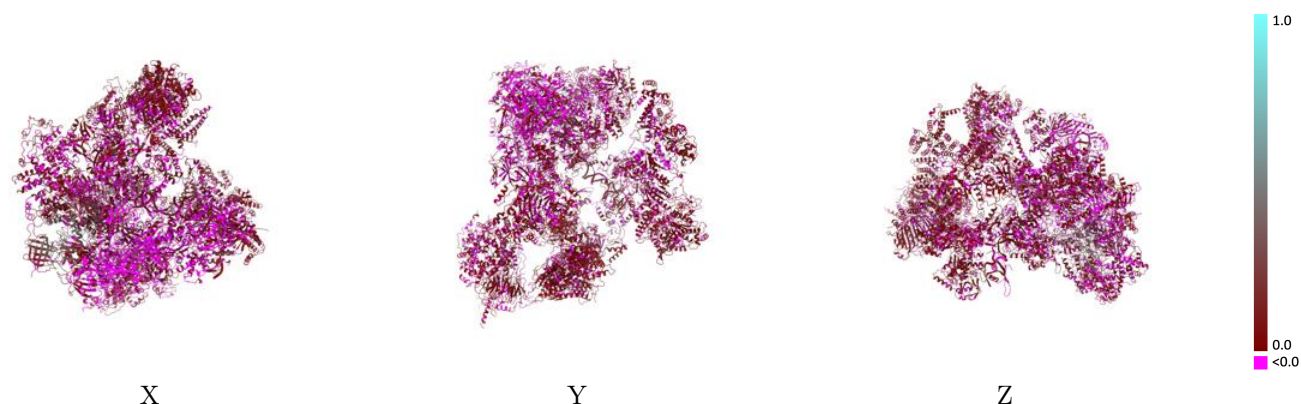
Y



Z

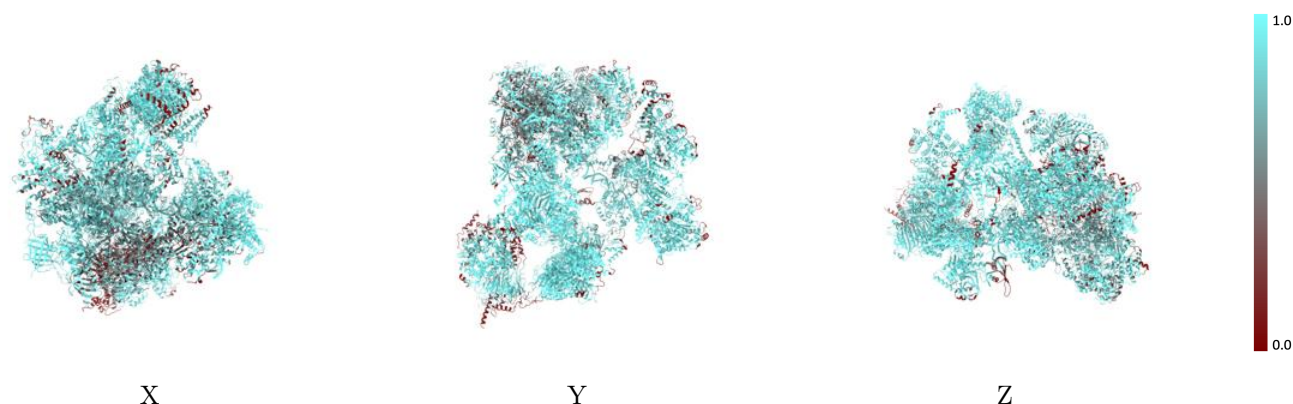
The images above show the 3D surface view of the map at the recommended contour level 0.0051 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



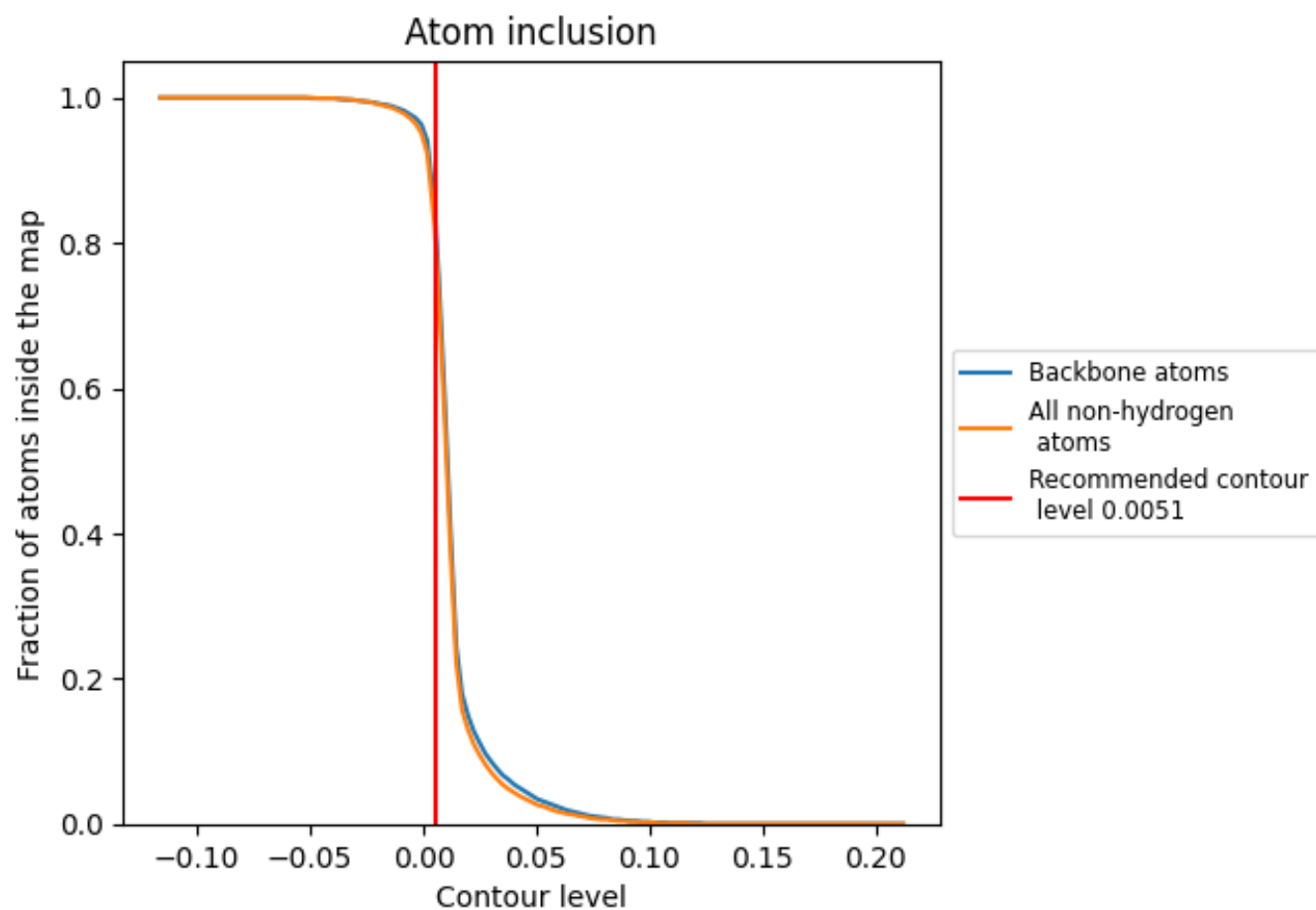
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0051).

9.4 Atom inclusion ⓘ



At the recommended contour level, 85% of all backbone atoms, 82% 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 (0.0051) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8210	 0.0510
0	 0.6780	 0.0570
1	 0.7750	 0.0270
2	 0.9400	 0.0590
3	 0.9280	 0.0540
4	 0.8880	 0.0660
5	 0.9290	 0.0540
6	 0.9350	 0.0560
7	 0.8980	 0.0570
8	 0.7650	 0.0310
9	 0.7140	 0.0580
A	 0.9050	 0.0510
B	 0.9610	 0.0590
D	 0.9570	 0.0370
E	 0.9470	 0.0620
F	 0.9620	 0.0700
G	 0.7490	 0.0490
H	 0.9540	 0.0570
I	 0.9500	 0.0490
J	 1.0000	 0.0740
L	 0.9600	 0.0400
O	 0.8670	 0.0250
P	 0.7910	 -0.0350
Q	 0.8340	 0.0430
R	 0.6570	 -0.0250
S	 0.8680	 0.0190
T	 0.7950	 0.0260
U	 0.8710	 0.0620
V	 0.9120	 0.0630
X	 0.8320	 0.0580
Y	 0.8250	 0.0620
c	 0.5150	 0.0220
d	 0.5320	 0.0350
e	 0.8140	 0.0440
f	 0.8270	 0.0520



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Chain	Atom inclusion	Q-score
i	 0.7930	 0.0320
j	 0.4240	 0.0270
k	 0.2970	 0.0090
l	 0.7300	 0.0210
m	 0.3460	 0.0180
o	 0.8150	 0.1210
p	 0.6110	 -0.0170
q	 0.6790	 -0.0350
r	 0.8940	 0.0770
s	 0.8650	 0.1030
t	 0.7920	 0.1140
u	 0.8390	 0.0330
v	 0.9310	 0.2230
w	 0.7750	 0.0030
x	 0.5760	 -0.0350
y	 0.7790	 0.1080
z	 0.6400	 -0.0420