



wwPDB EM Validation Summary Report ⓘ

Mar 24, 2026 – 04:41 PM UTC

PDB ID : 7EGB / pdb_00007egb
EMDB ID : EMD-31111
Title : TFIID-based holo PIC on SCP promoter
Authors : Chen, X.; Wu, Z.; Hou, H.; Qi, Y.; Wang, X.; Li, J.; Xu, Y.
Deposited on : 2021-03-24
Resolution : 3.30 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:

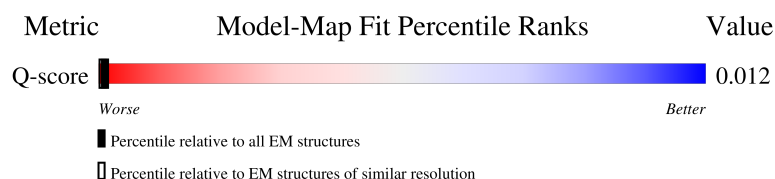
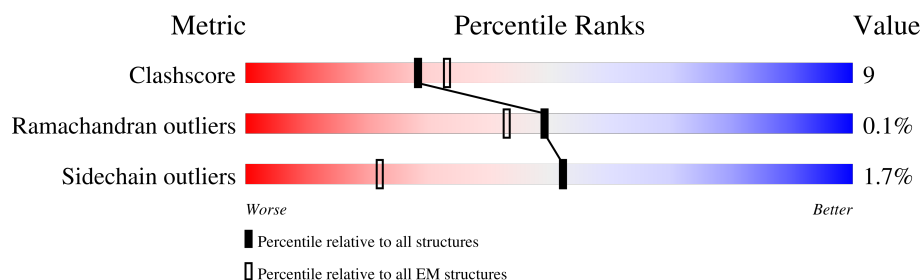
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 3.30 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	15087 (2.80 - 3.80)

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	
2	1	548	
3	2	395	
4	3	308	

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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	94	
30	Y	69	
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 112409 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	236	Total	C	N	O	S	0	0
			1932	1211	336	373	12		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	265	Total	C	N	O	S	0	0
			2167	1382	378	395	12		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	2	329	Total	C	N	O	S	0	0
			2567	1621	440	479	27		

- 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
			2066	1323	344	380	19		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
5	4	394	Total	C	N	O	S	0	0
			3189	2068	552	557	12		

- 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	605	Total	C	N	O	S	0	0
			4890	3127	848	885	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	714	Total	C	N	O	S	0	0
			5751	3683	999	1040	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
			2376	1537	404	423	12		

- Molecule 10 is a protein called Cyclin-H.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	9	281	Total	C	N	O	S	0	0
			2293	1465	394	416	18		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
11	A	559	Total	C	N	O	S	0	0
			4580	2924	795	834	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	160	Total	C	N	O	S	0	0
			1337	834	249	251	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	547	Total	C	N	O	S	0	0
			4376	2774	759	822	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	412	Total	C	N	O	S	0	0
			3143	1994	548	583	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	144	Total	C	N	O	S	0	0
			1171	742	215	210	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	76	Total	C	N	O	S	0	0
			622	388	109	122	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	99	Total	C	N	O	S	0	0
			806	510	142	151	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
22	P	179	Total	C	N	O	S	0	0
			1422	923	251	241	7		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
23	Q	113	Total	C	N	O	S	0	0
			930	585	152	189	4		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
24	R	252	Total	C	N	O	S	0	0
			1953	1224	346	366	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	184	Total	C	N	O	S	0	0
			1520	957	272	280	11		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
28	V	165	Total	C	N	O	S	0	0
			1357	865	235	253	4		

- Molecule 29 is a DNA chain called DNA (69-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
29	X	69	Total	C	N	O	P	0	0
			1429	672	279	409	69		

- Molecule 30 is a DNA chain called DNA (69-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
30	Y	69	Total	C	N	O	P	0	0
			1400	664	248	419	69		

- 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	1423	Total	C	N	O	S	0	0
			11274	7092	2016	2094	72		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
35	p	1135	Total	C	N	O	S	0	0
			9068	5735	1596	1673	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
			1050	656	178	212	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
			1721	1089	300	324	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
			636	406	108	117	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
			928	571	166	180	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	115	Total	C	N	O	S	0	0
			920	593	152	173	2		

- Molecule 45 is a protein called RPB12.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	z	44	Total	C	N	O	S	0	0
			373	231	72	64	6		

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

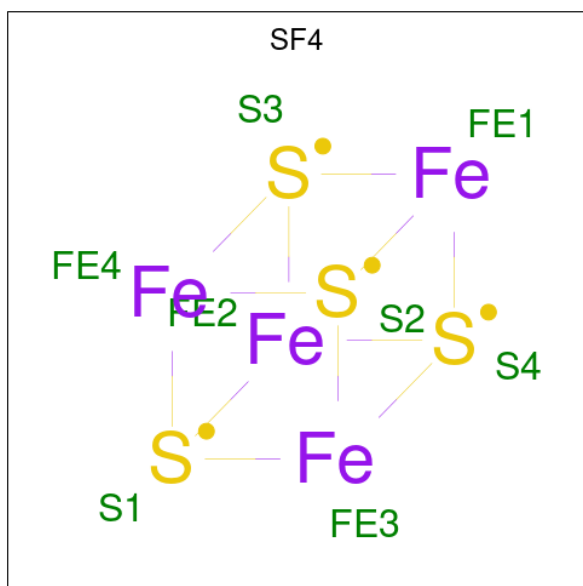
Mol	Chain	Residues	Atoms		AltConf
46	0	2	Total	Zn	0
			2	2	
46	2	3	Total	Zn	0
			3	3	
46	3	2	Total	Zn	0
			2	2	

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

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



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

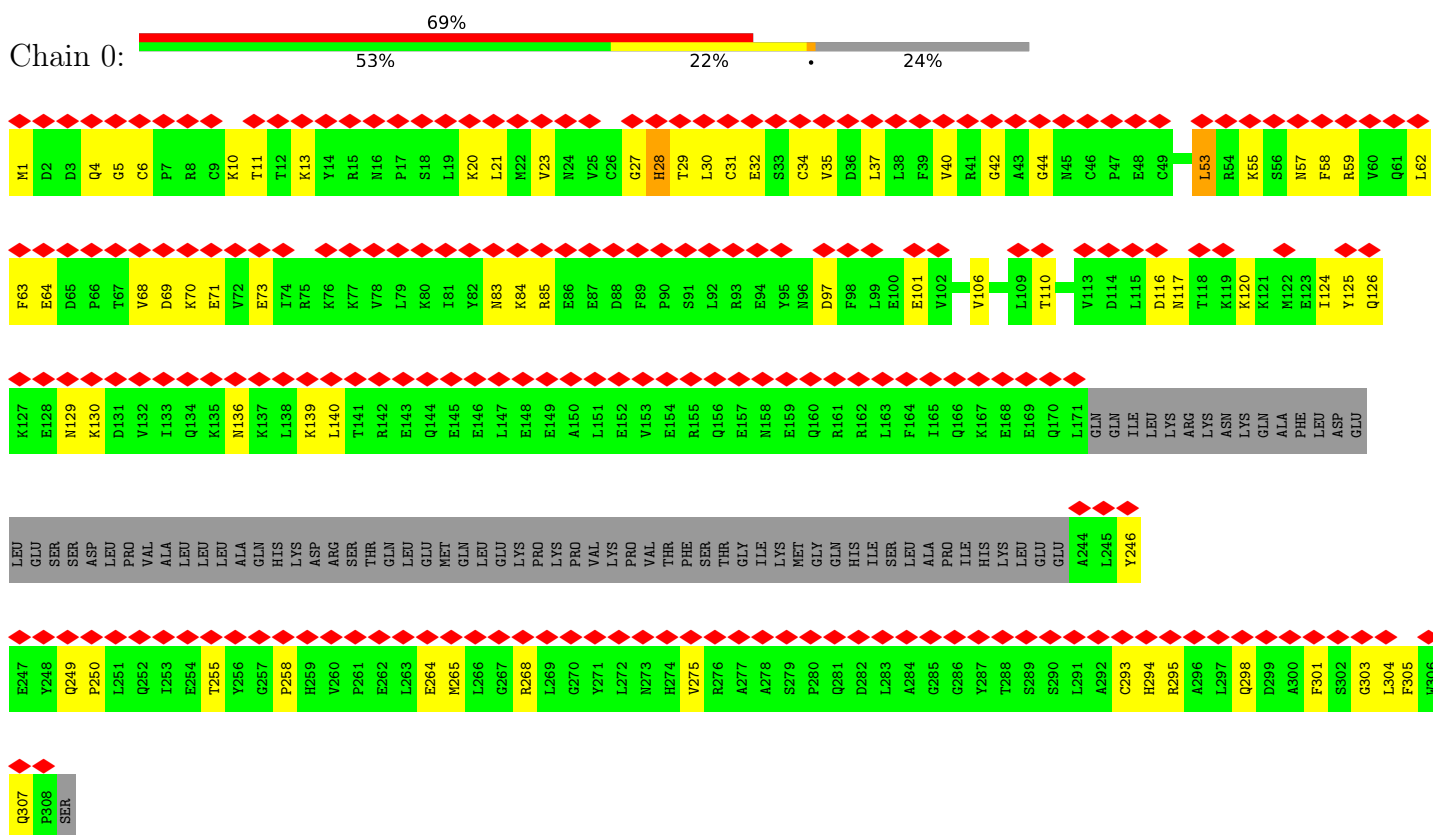
- Molecule 48 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

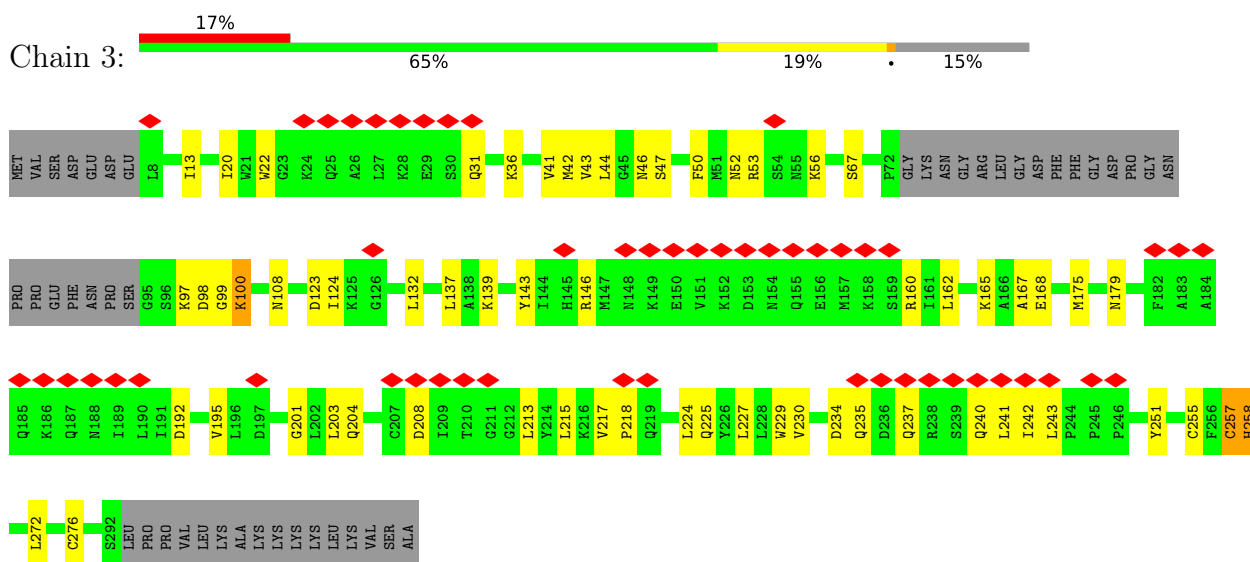
3 Residue-property plots

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.

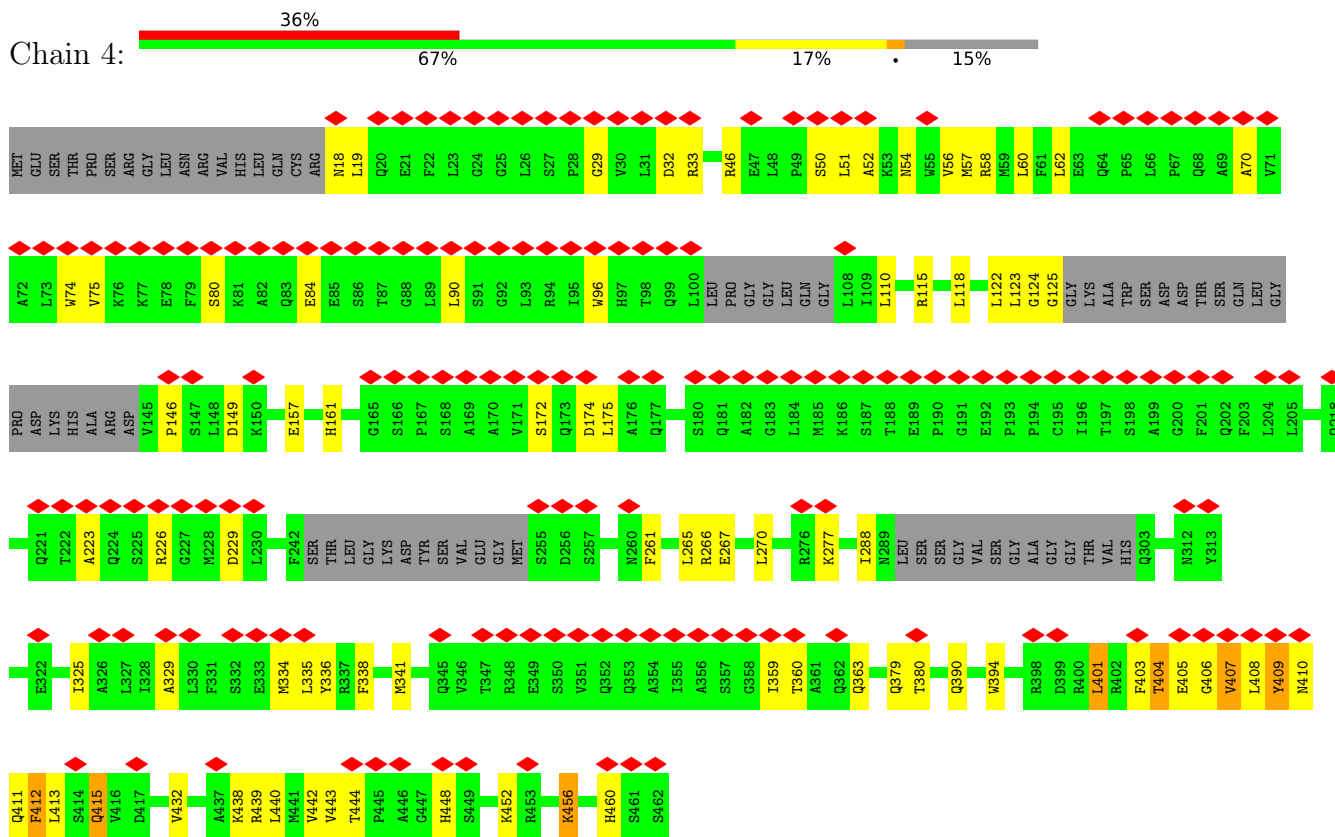
- Molecule 1: CDK-activating kinase assembly factor MAT1



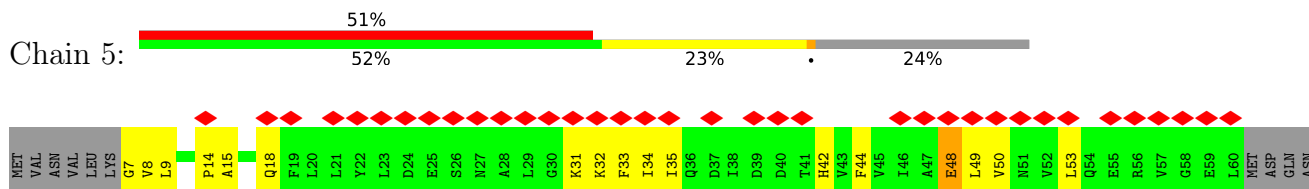


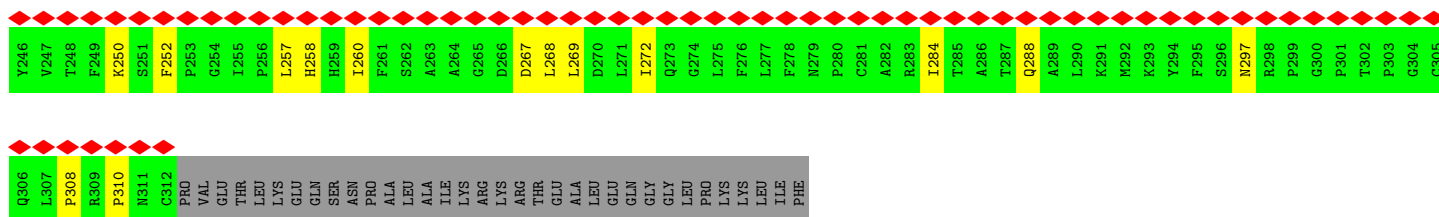


- Molecule 5: General transcription factor IIH subunit 4

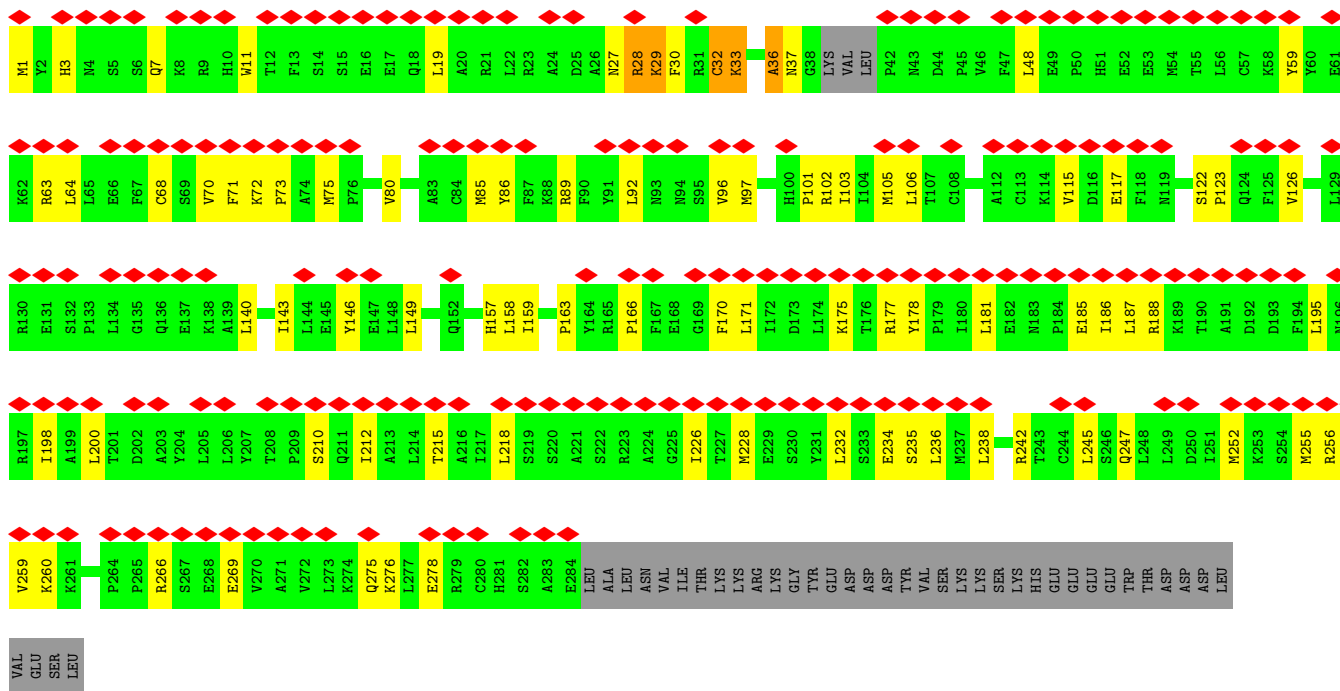


- Molecule 6: General transcription factor IIH subunit 5





• Molecule 10: Cyclin-H



• Molecule 11: Transcription initiation factor TFIID subunit 1





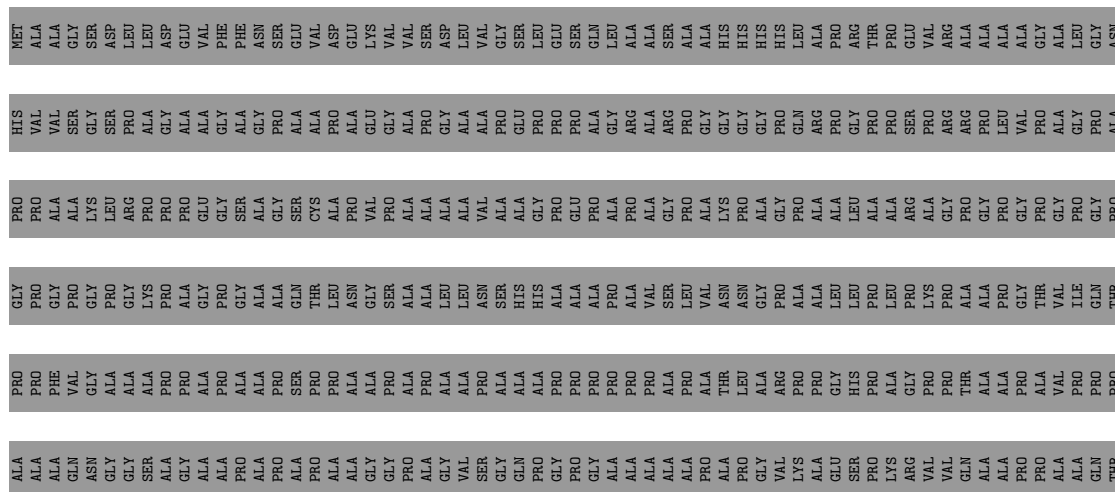
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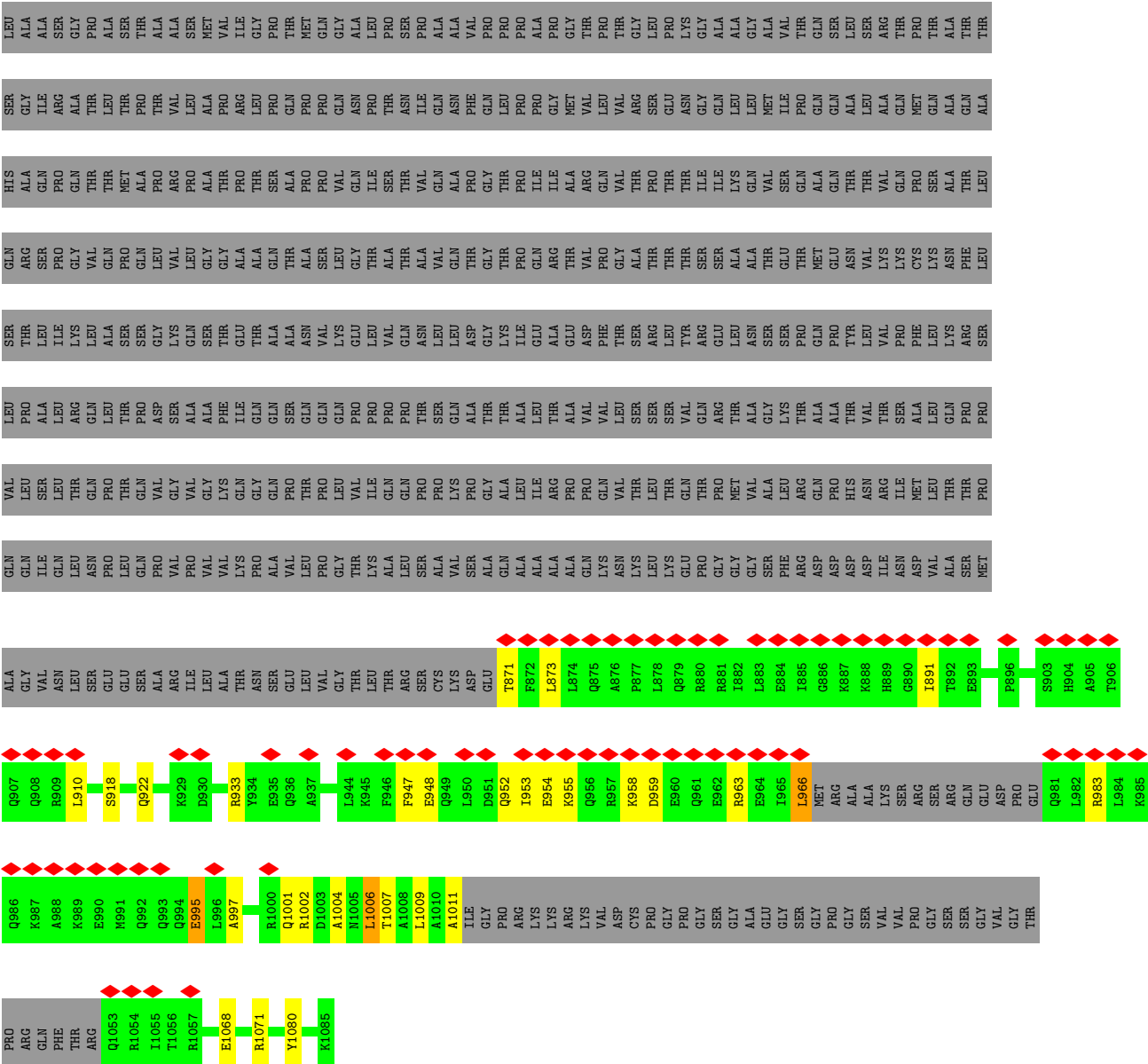
• Molecule 12: Transcription initiation factor TFIID subunit 2



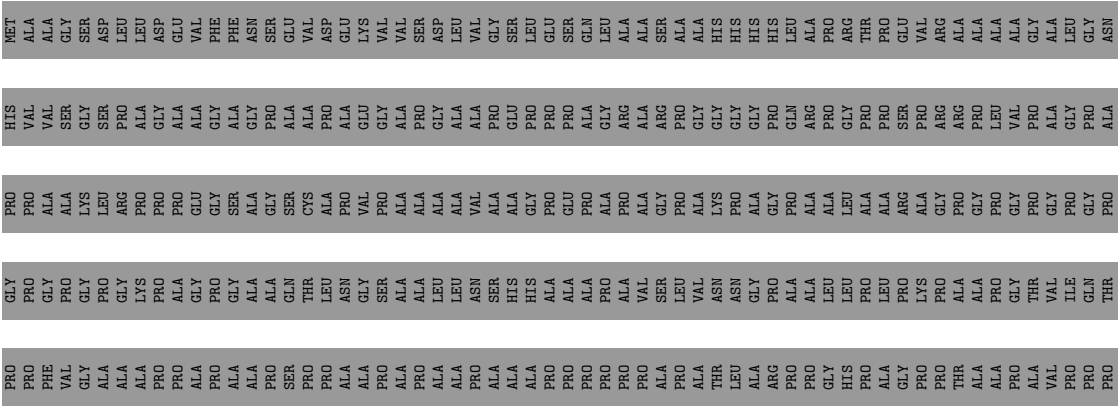
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V75	R76	I77	N78	D79	L80	E81	A82	A83	F84	I85	Y86	N87	T90	H95	S96	E97	S98	K99	Q100	R101	N102	Y105	N108	A112	P119	D120	A121	G122	N123	C127	I128	K129	V130	P131	S132	E133	L134	W135	K136	H137	V138	D139	E140	L141	K142	V143	L144	K145	I146	H147	I148	N149		
L152	P155	E167	H176	G181	N184	R187	W202	K203	S216	N217	G218	D219	L220	V221	E222	T226	H227	D228	M229	R230	H235	Y236	M237	L238	P241	S248	V258	D259	P260	Y261	H263	E264	V265	H266	H267	L274	L277	K278	Y283	L284	H285													
E286	V287	F288	E289	E292	E293	I294	L295	T296	C297	R298	S302	C303	F304	D310	Y319	A320	S321	M322	N328	L329	D337	T342	R343	R344	C345	Q348	Q353	R361	D366	E367	M382	K383	K384	N389	E397	I402	L414	I417	F418	GLY	GLY	GLY												

- Molecule 13: Transcription initiation factor TFIID subunit 4





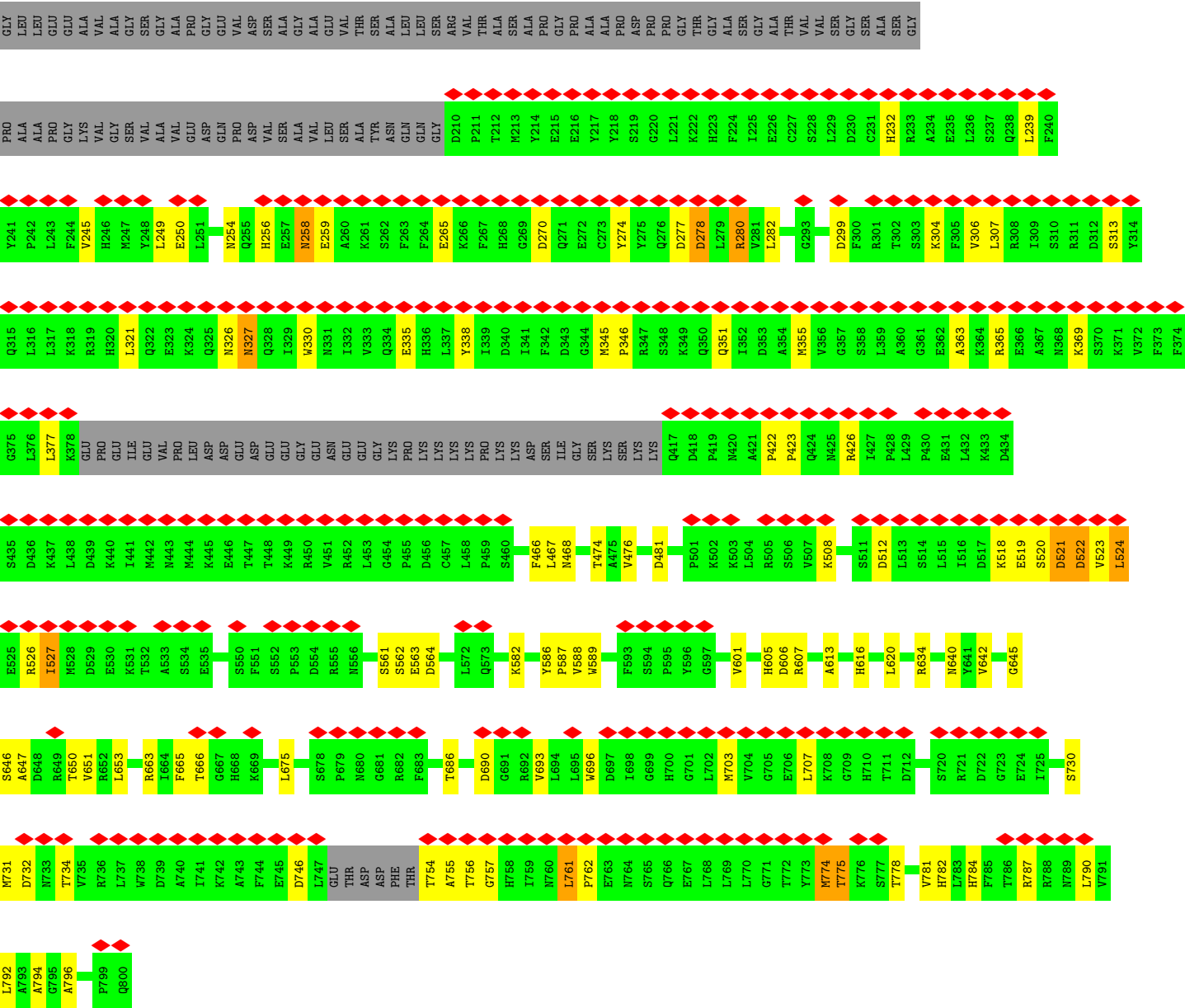
● Molecule 13: Transcription initiation factor TFIID subunit 4



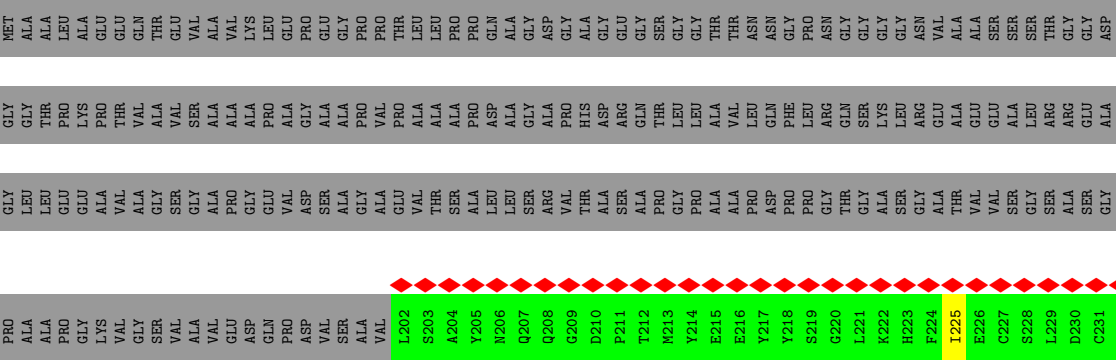
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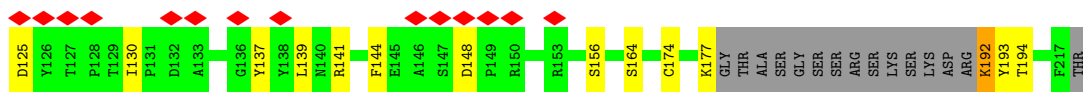
Chain E:

[illegible]

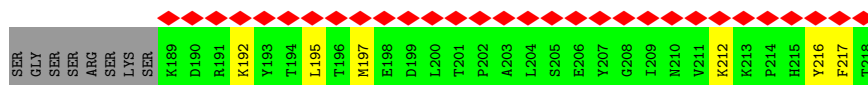
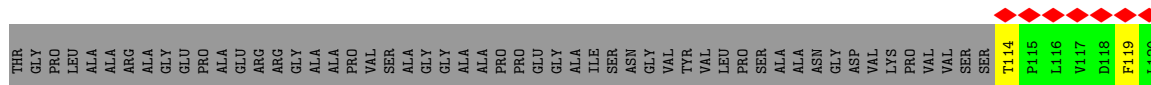
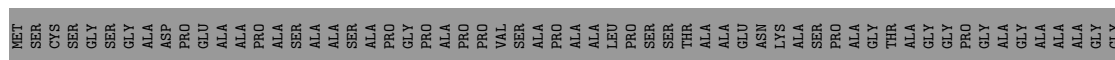
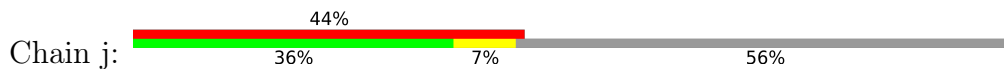


• Molecule 14: Transcription initiation factor TFIID subunit 5

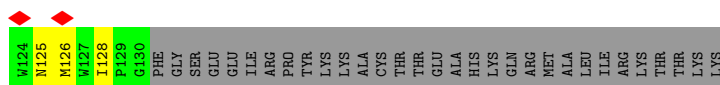
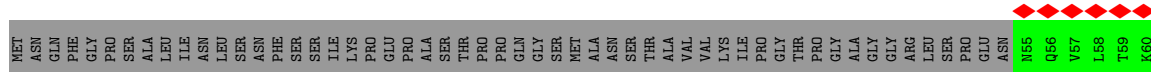
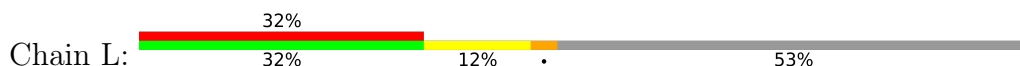




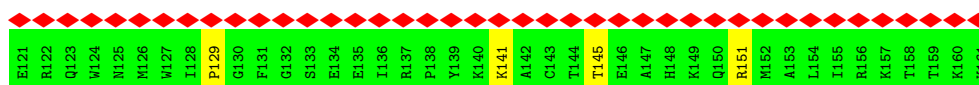
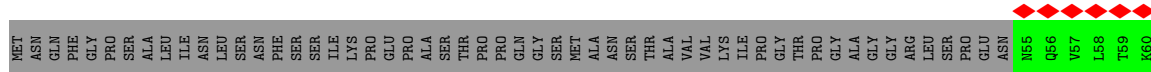
• Molecule 19: Transcription initiation factor TFIID subunit 10



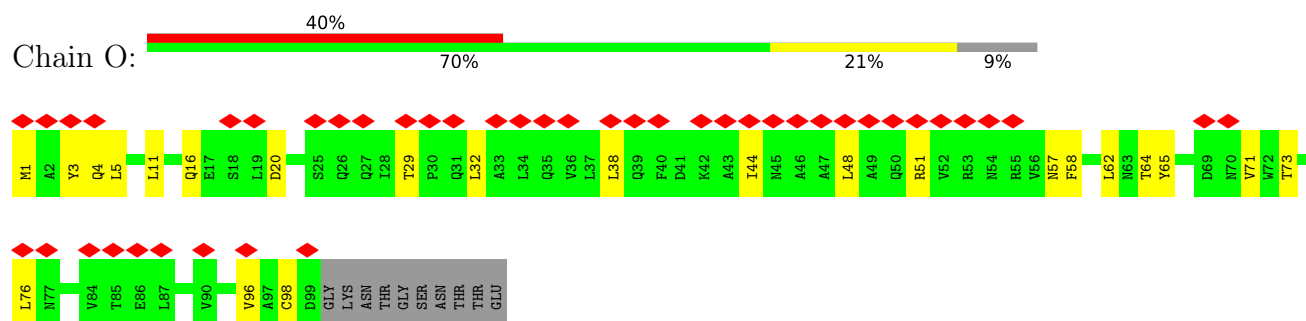
• Molecule 20: Transcription initiation factor TFIID subunit 12



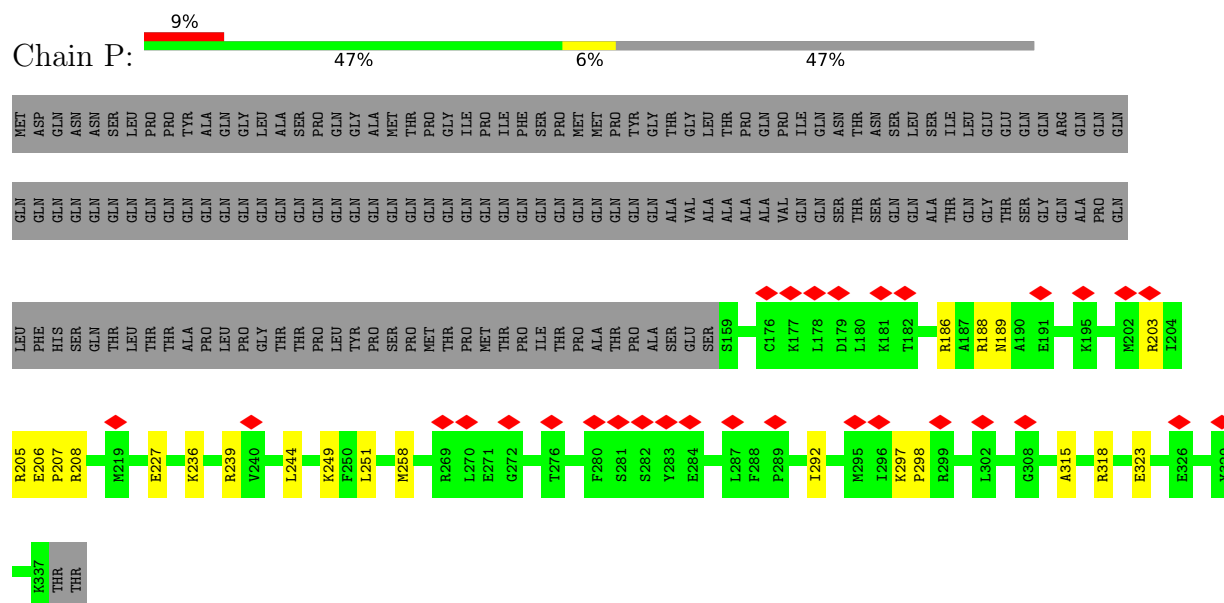
• Molecule 20: Transcription initiation factor TFIID subunit 12



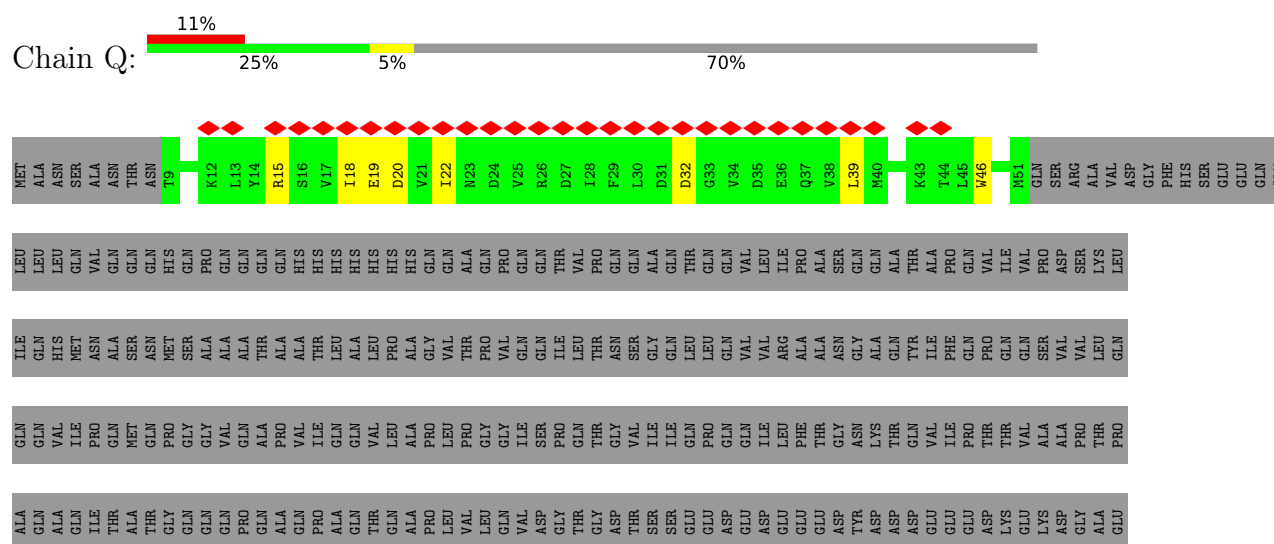
- Molecule 21: Transcription initiation factor IIA subunit 2



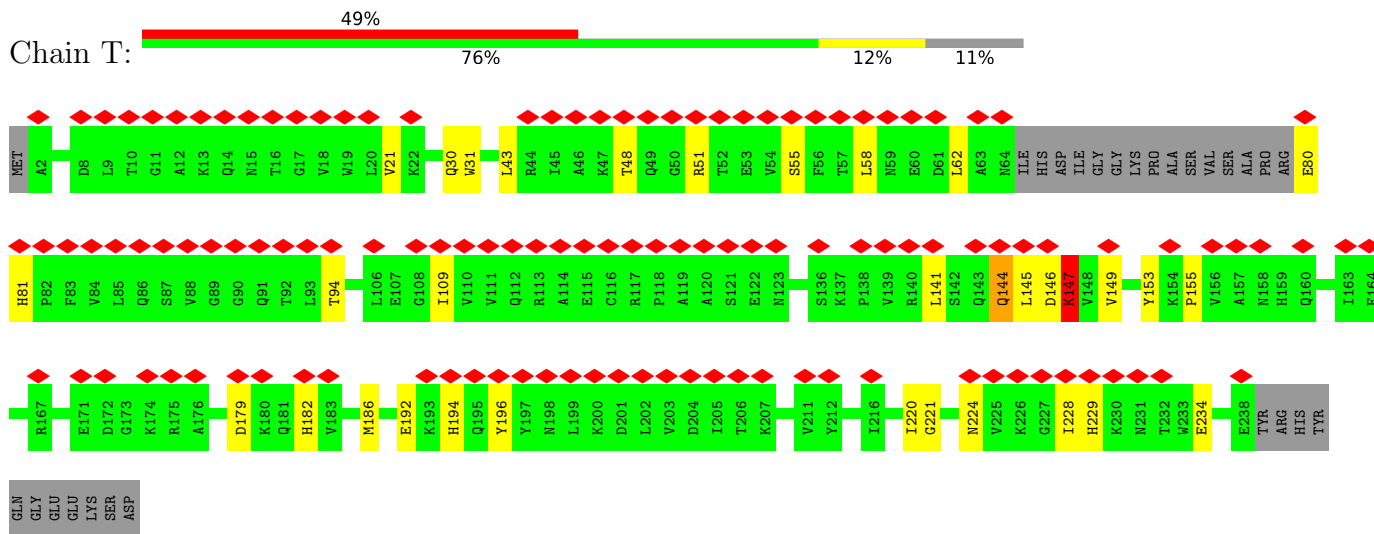
- Molecule 22: TATA-box-binding protein



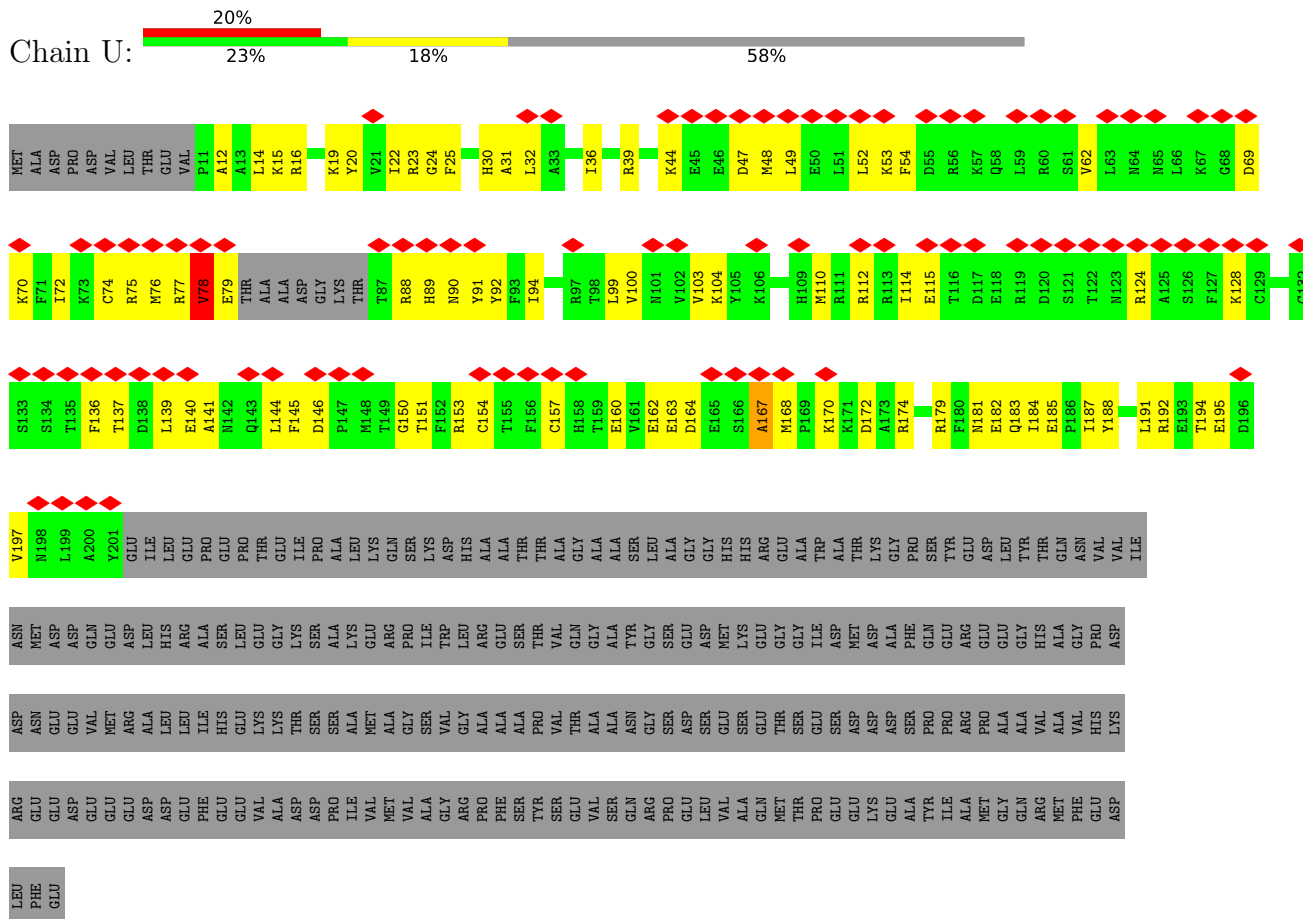
- Molecule 23: Transcription initiation factor IIA subunit 1



Chain T:



Chain U:

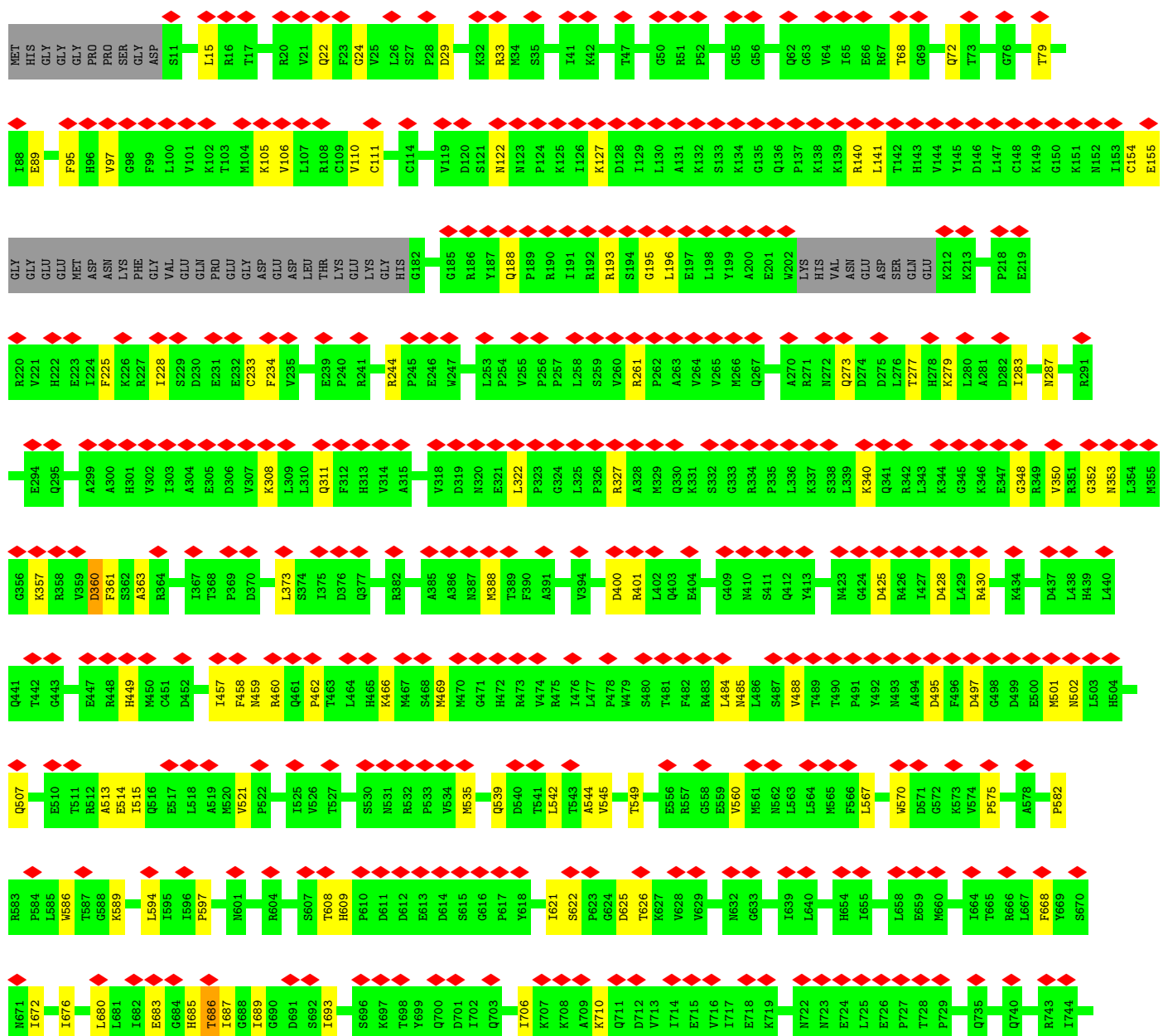
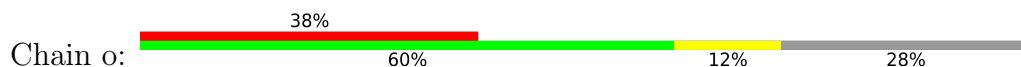


Chain V:





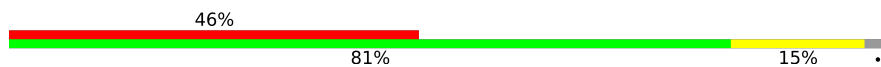
• Molecule 34: DNA-directed RNA polymerase II subunit RPB1

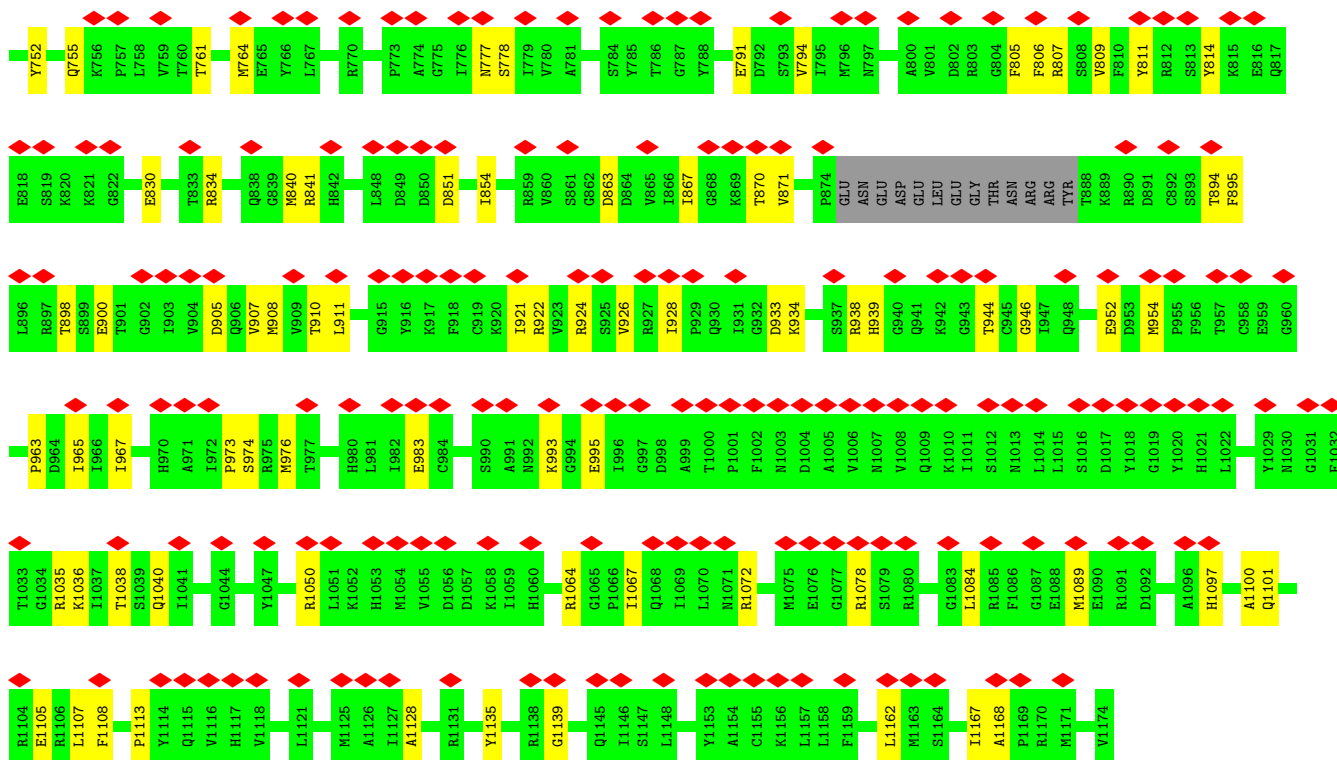




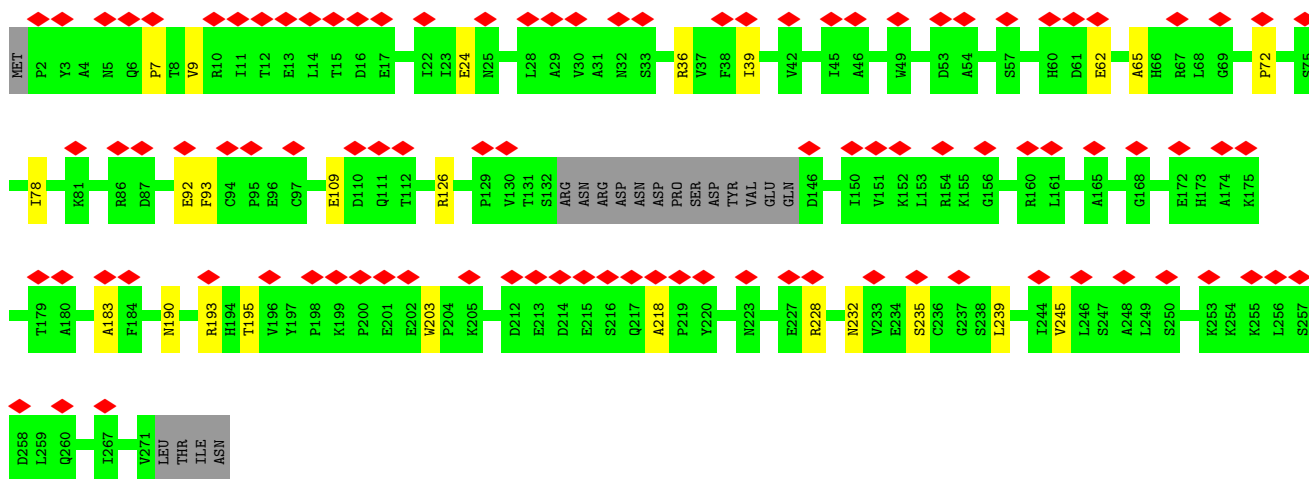
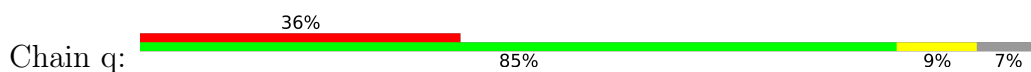
PRO	THR	TYR	SER	PRO	THR	SER	LYS	GLY	SER	THR	TYR	SER	PRO	THR	SER	GLY	TYR	SER	PRO	THR	SER	PRO	ALA	ILE	SER	ASP	ASP	ASP	ASP	GLU	VAL
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

Chain p:



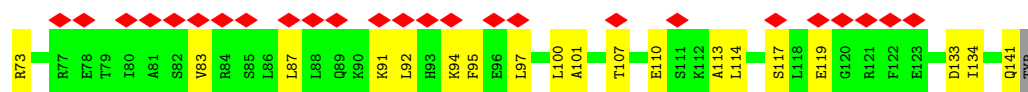


• Molecule 36: DNA-directed RNA polymerase II subunit RPB3

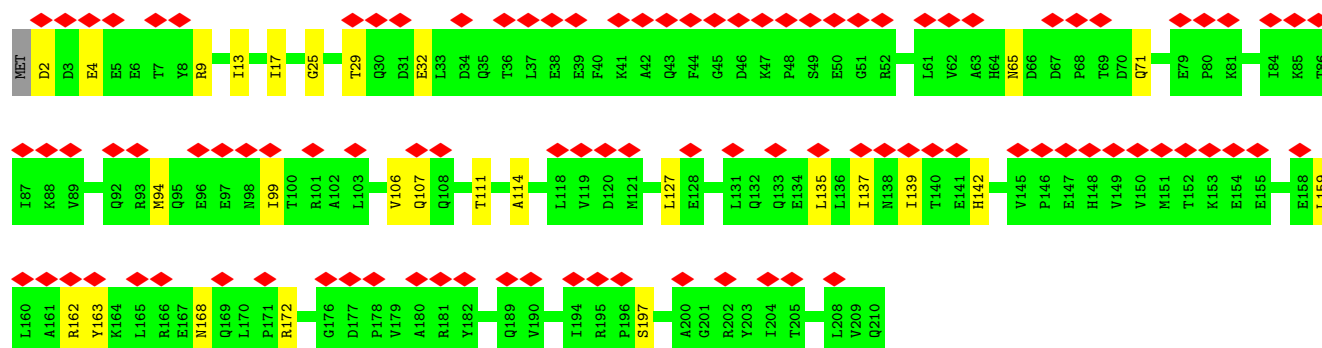
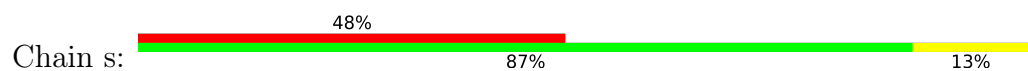


• Molecule 37: DNA-directed RNA polymerase II subunit RPB4

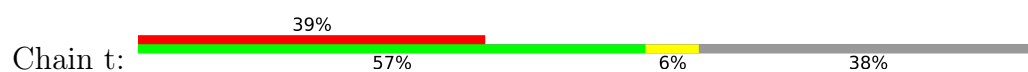




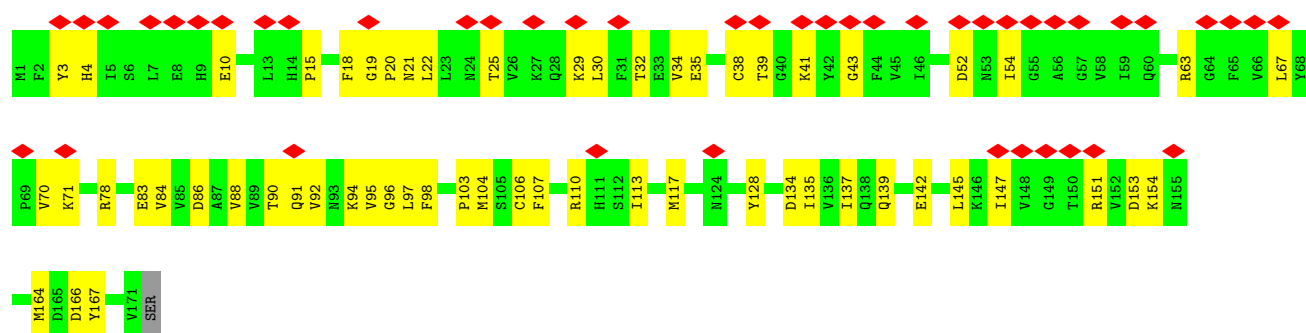
- Molecule 38: DNA-directed RNA polymerase II subunit E



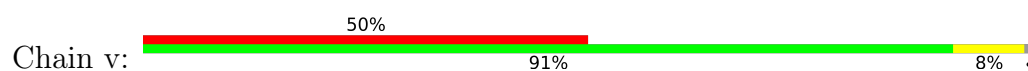
- Molecule 39: DNA-directed RNA polymerase II subunit F

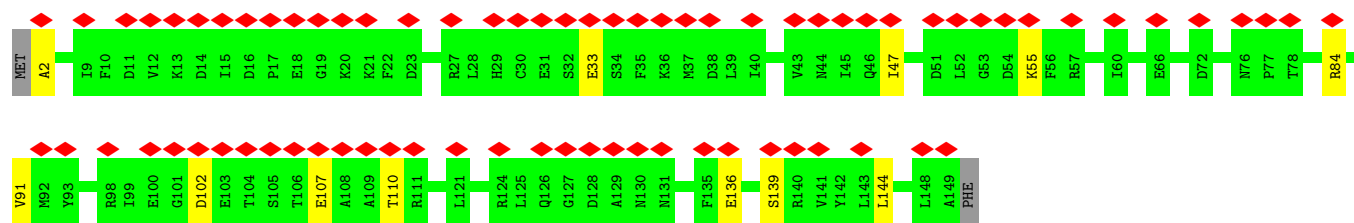


- Molecule 40: DNA-directed RNA polymerase II subunit RPB7

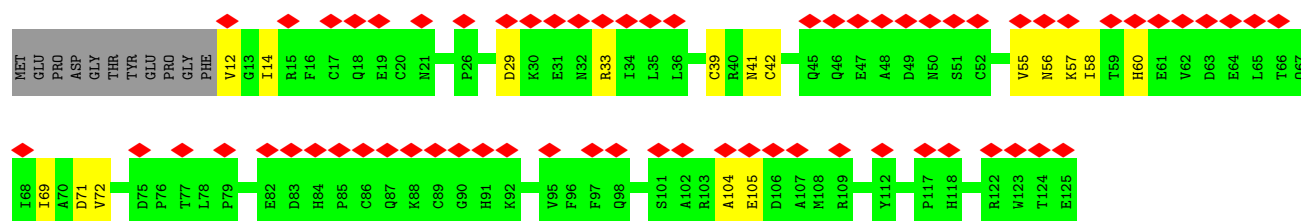
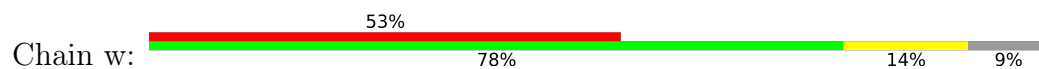


- Molecule 41: DNA-directed RNA polymerases I, II, and III subunit RPABC3

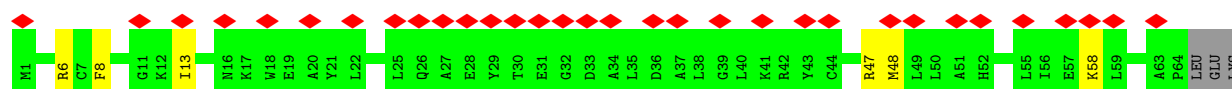
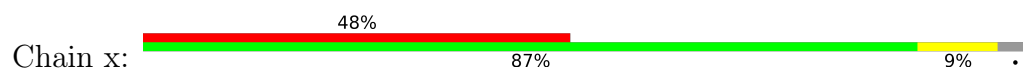




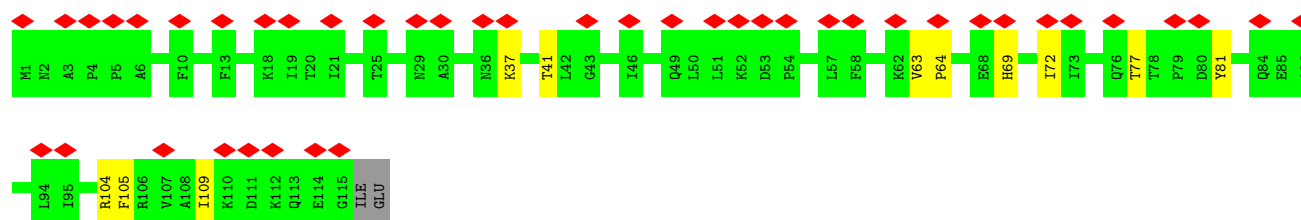
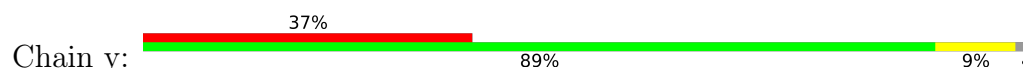
• Molecule 42: DNA-directed RNA polymerase II subunit RPB9



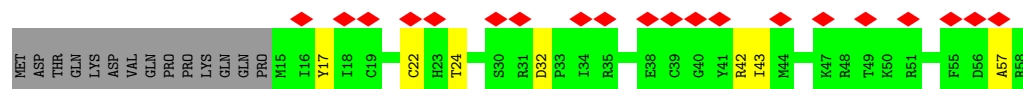
• Molecule 43: RPB10



• 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	92806	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)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	6.811	Depositor
Minimum map value	-4.382	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.084	Depositor
Recommended contour level	0.3	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: ZN, MG, SF4

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	0.39	0/1965	0.62	0/2646
2	1	0.23	0/2210	0.39	0/2975
3	2	0.16	0/2624	0.28	0/3555
4	3	0.30	0/2103	0.47	1/2846 (0.0%)
5	4	0.31	0/3262	0.51	2/4418 (0.0%)
6	5	0.34	0/433	0.60	0/585
7	6	0.29	0/4994	0.48	2/6745 (0.0%)
8	7	0.21	0/5875	0.37	1/7955 (0.0%)
9	8	0.35	0/2434	0.55	1/3300 (0.0%)
10	9	0.46	0/2342	0.72	6/3159 (0.2%)
11	A	0.63	0/4698	0.95	5/6345 (0.1%)
12	B	0.46	1/7993 (0.0%)	0.69	6/10836 (0.1%)
13	D	0.54	0/1350	0.81	2/1805 (0.1%)
13	d	0.48	0/1321	0.71	2/1772 (0.1%)
14	E	0.45	1/4482 (0.0%)	0.72	3/6069 (0.0%)
14	e	0.55	1/4433 (0.0%)	0.77	0/6004
15	F	0.65	1/3201 (0.0%)	1.07	13/4347 (0.3%)
15	f	0.47	1/3140 (0.0%)	0.80	6/4268 (0.1%)
16	G	0.68	0/1190	0.94	1/1601 (0.1%)
17	H	0.63	3/1673 (0.2%)	1.00	14/2285 (0.6%)
18	I	0.21	0/981	0.41	0/1332
18	i	0.21	0/989	0.40	0/1343
19	J	0.83	0/736	1.15	2/998 (0.2%)
19	j	0.73	0/775	0.96	0/1049
20	L	0.68	0/630	1.13	1/852 (0.1%)
20	l	0.59	0/888	0.88	0/1194
21	O	0.25	0/816	0.43	0/1105
22	P	0.16	0/1448	0.28	0/1948
23	Q	0.18	0/945	0.35	0/1274
24	R	0.23	0/1983	0.38	0/2679
25	S	0.12	0/896	0.25	0/1213
26	T	0.20	0/1817	0.39	1/2445 (0.0%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
27	U	0.41	0/1545	0.68	1/2075 (0.0%)
28	V	0.24	0/1380	0.58	0/1854
29	X	0.28	0/1607	0.58	0/2481
30	Y	0.29	0/1565	0.59	0/2410
31	c	0.48	0/1035	0.67	0/1406
32	k	0.29	0/799	0.53	1/1070 (0.1%)
33	m	0.85	0/733	1.18	1/977 (0.1%)
34	o	0.37	8/11479 (0.1%)	0.55	21/15496 (0.1%)
35	p	0.40	6/9249 (0.1%)	0.39	7/12482 (0.1%)
36	q	0.15	0/2102	0.29	0/2857
37	r	0.14	0/1064	0.22	0/1428
38	s	0.29	1/1752 (0.1%)	0.53	3/2366 (0.1%)
39	t	0.16	0/646	0.27	0/871
40	u	0.20	0/1382	0.28	0/1874
41	v	0.14	0/1207	0.28	0/1628
42	w	0.23	0/949	0.41	0/1284
43	x	0.16	0/516	0.25	0/696
44	y	0.22	0/939	0.37	0/1271
45	z	0.17	0/378	0.28	0/500
All	All	0.41	23/114954 (0.0%)	0.62	103/155974 (0.1%)

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	0	0	1
26	T	0	1
27	U	0	1
All	All	0	3

The worst 5 of 23 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
35	p	704	LEU	C-N	20.97	1.56	1.33
35	p	806	PHE	C-N	16.04	1.56	1.33
35	p	928	ILE	C-N	13.37	1.50	1.33
34	o	388	MET	C-N	12.26	1.50	1.33
34	o	1466	ALA	C-N	12.03	1.51	1.33

The worst 5 of 103 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	o	1082	HIS	CA-C-N	16.32	136.59	119.78
34	o	1082	HIS	C-N-CA	16.32	136.59	119.78
38	s	139	ILE	O-C-N	-15.11	106.61	121.87
34	o	1483	GLY	N-CA-C	14.43	129.23	112.50
17	H	156	PRO	N-CA-C	-13.59	90.10	111.03

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	0	126	GLN	Mainchain
26	T	144	GLN	Mainchain
27	U	78	VAL	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	0	1932	0	1905	86	0
2	1	2167	0	2175	35	0
3	2	2567	0	2539	49	0
4	3	2066	0	2097	89	0
5	4	3189	0	3243	131	0
6	5	428	0	434	45	0
7	6	4890	0	4949	214	0
8	7	5751	0	5794	115	0
9	8	2376	0	2402	73	0
10	9	2293	0	2315	76	0
11	A	4580	0	4502	97	0
12	B	7796	0	7751	113	0
13	D	1337	0	1358	44	0
13	d	1307	0	1318	16	0
14	E	4376	0	4256	78	0
14	e	4327	0	4197	64	0
15	F	3143	0	3093	103	0
15	f	3081	0	3012	70	0
16	G	1171	0	1222	17	0
17	H	1633	0	1621	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	I	959	0	969	13	0
18	i	967	0	977	10	0
19	J	720	0	719	12	0
19	j	759	0	759	14	0
20	L	622	0	620	21	0
20	l	876	0	893	10	0
21	O	806	0	818	22	0
22	P	1422	0	1514	29	0
23	Q	930	0	888	31	0
24	R	1953	0	1987	30	0
25	S	872	0	848	9	0
26	T	1788	0	1819	40	0
27	U	1520	0	1520	88	0
28	V	1357	0	1381	78	0
29	X	1429	0	770	13	0
30	Y	1400	0	775	9	0
31	c	1011	0	975	17	0
32	k	785	0	818	7	0
33	m	724	0	744	26	0
34	o	11274	0	11406	175	0
35	p	9068	0	9112	157	0
36	q	2059	0	2007	18	0
37	r	1050	0	1033	30	0
38	s	1721	0	1737	18	0
39	t	636	0	665	18	0
40	u	1351	0	1358	65	0
41	v	1186	0	1147	9	0
42	w	928	0	861	16	0
43	x	507	0	523	7	0
44	y	920	0	942	11	0
45	z	373	0	378	6	0
46	0	2	0	0	0	0
46	2	3	0	0	0	0
46	3	2	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
47	7	8	0	0	0	0
48	o	1	0	0	0	0
All	All	112409	0	111146	2117	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 9.

The worst 5 of 2117 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)
7:6:308:ILE:CG2	7:6:358:ARG:HA	1.25	1.62
6:5:33:PHE:CE1	6:5:49:LEU:HD12	1.44	1.51
6:5:31:LYS:HB3	12:B:540:LYS:NZ	1.31	1.39
4:3:50:PHE:CZ	5:4:118:LEU:HD11	1.57	1.37
6:5:31:LYS:CB	12:B:540:LYS:HZ3	1.37	1.37

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	232/309 (75%)	215 (93%)	17 (7%)	0	100	100
2	1	253/548 (46%)	245 (97%)	8 (3%)	0	100	100
3	2	325/395 (82%)	313 (96%)	12 (4%)	0	100	100
4	3	259/308 (84%)	251 (97%)	8 (3%)	0	100	100
5	4	384/462 (83%)	367 (96%)	17 (4%)	0	100	100
6	5	52/71 (73%)	47 (90%)	4 (8%)	1 (2%)	6	28
7	6	601/782 (77%)	576 (96%)	23 (4%)	2 (0%)	36	65
8	7	710/760 (93%)	683 (96%)	26 (4%)	1 (0%)	48	75

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	8	294/346 (85%)	277 (94%)	17 (6%)	0	100	100
10	9	277/323 (86%)	268 (97%)	9 (3%)	0	100	100
11	A	543/1872 (29%)	519 (96%)	23 (4%)	1 (0%)	43	71
12	B	959/1199 (80%)	911 (95%)	48 (5%)	0	100	100
13	D	154/1085 (14%)	148 (96%)	6 (4%)	0	100	100
13	d	154/1085 (14%)	149 (97%)	5 (3%)	0	100	100
14	E	541/800 (68%)	517 (96%)	23 (4%)	1 (0%)	43	71
14	e	531/800 (66%)	482 (91%)	48 (9%)	1 (0%)	43	71
15	F	408/677 (60%)	389 (95%)	18 (4%)	1 (0%)	43	71
15	f	399/677 (59%)	380 (95%)	19 (5%)	0	100	100
16	G	138/349 (40%)	135 (98%)	3 (2%)	0	100	100
17	H	207/310 (67%)	189 (91%)	14 (7%)	4 (2%)	6	28
18	I	118/264 (45%)	115 (98%)	3 (2%)	0	100	100
18	i	119/264 (45%)	115 (97%)	4 (3%)	0	100	100
19	J	86/218 (39%)	82 (95%)	4 (5%)	0	100	100
19	j	91/218 (42%)	85 (93%)	4 (4%)	2 (2%)	5	26
20	L	74/161 (46%)	72 (97%)	2 (3%)	0	100	100
20	l	105/161 (65%)	100 (95%)	5 (5%)	0	100	100
21	O	97/109 (89%)	97 (100%)	0	0	100	100
22	P	177/339 (52%)	175 (99%)	2 (1%)	0	100	100
23	Q	109/376 (29%)	102 (94%)	7 (6%)	0	100	100
24	R	248/316 (78%)	246 (99%)	2 (1%)	0	100	100
25	S	104/517 (20%)	104 (100%)	0	0	100	100
26	T	218/249 (88%)	211 (97%)	6 (3%)	1 (0%)	24	55
27	U	180/439 (41%)	152 (84%)	26 (14%)	2 (1%)	11	39
28	V	163/291 (56%)	139 (85%)	24 (15%)	0	100	100
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%)	82 (96%)	3 (4%)	0	100	100
34	o	1413/1970 (72%)	1369 (97%)	44 (3%)	0	100	100
35	p	1129/1174 (96%)	1089 (96%)	40 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
36	q	253/275 (92%)	241 (95%)	12 (5%)	0	100	100
37	r	126/142 (89%)	122 (97%)	4 (3%)	0	100	100
38	s	207/210 (99%)	203 (98%)	4 (2%)	0	100	100
39	t	77/127 (61%)	76 (99%)	1 (1%)	0	100	100
40	u	169/172 (98%)	166 (98%)	3 (2%)	0	100	100
41	v	146/150 (97%)	143 (98%)	3 (2%)	0	100	100
42	w	112/125 (90%)	103 (92%)	9 (8%)	0	100	100
43	x	62/67 (92%)	61 (98%)	1 (2%)	0	100	100
44	y	113/117 (97%)	112 (99%)	1 (1%)	0	100	100
45	z	42/58 (72%)	41 (98%)	1 (2%)	0	100	100
All	All	13465/22931 (59%)	12871 (96%)	577 (4%)	17 (0%)	49	75

5 of 17 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	5	48	GLU
14	E	523	VAL
17	H	128	PRO
14	e	522	ASP
19	j	124	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	216/283 (76%)	213 (99%)	3 (1%)	59	73
2	1	241/484 (50%)	241 (100%)	0	100	100
3	2	295/352 (84%)	295 (100%)	0	100	100
4	3	234/272 (86%)	229 (98%)	5 (2%)	47	67
5	4	346/399 (87%)	337 (97%)	9 (3%)	40	64
6	5	48/64 (75%)	48 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	6	536/688 (78%)	535 (100%)	1 (0%)	87	88
8	7	624/664 (94%)	623 (100%)	1 (0%)	87	88
9	8	259/299 (87%)	258 (100%)	1 (0%)	84	84
10	9	256/296 (86%)	254 (99%)	2 (1%)	73	79
11	A	496/1665 (30%)	452 (91%)	44 (9%)	9	32
12	B	876/1083 (81%)	862 (98%)	14 (2%)	55	72
13	D	145/815 (18%)	139 (96%)	6 (4%)	27	55
13	d	146/815 (18%)	144 (99%)	2 (1%)	59	73
14	E	480/657 (73%)	467 (97%)	13 (3%)	39	63
14	e	475/657 (72%)	462 (97%)	13 (3%)	39	63
15	F	328/574 (57%)	303 (92%)	25 (8%)	12	37
15	f	322/574 (56%)	312 (97%)	10 (3%)	35	60
16	G	132/322 (41%)	124 (94%)	8 (6%)	17	45
17	H	181/270 (67%)	173 (96%)	8 (4%)	25	54
18	I	106/235 (45%)	106 (100%)	0	100	100
18	i	107/235 (46%)	107 (100%)	0	100	100
19	J	79/154 (51%)	72 (91%)	7 (9%)	9	32
19	j	83/154 (54%)	82 (99%)	1 (1%)	63	75
20	L	71/141 (50%)	66 (93%)	5 (7%)	14	41
20	l	98/141 (70%)	97 (99%)	1 (1%)	68	76
21	O	90/98 (92%)	90 (100%)	0	100	100
22	P	154/293 (53%)	154 (100%)	0	100	100
23	Q	105/324 (32%)	105 (100%)	0	100	100
24	R	215/268 (80%)	214 (100%)	1 (0%)	81	83
25	S	93/448 (21%)	93 (100%)	0	100	100
26	T	196/218 (90%)	193 (98%)	3 (2%)	57	72
27	U	167/373 (45%)	163 (98%)	4 (2%)	43	65
28	V	150/261 (58%)	149 (99%)	1 (1%)	76	80
31	c	113/833 (14%)	111 (98%)	2 (2%)	51	70
32	k	87/182 (48%)	87 (100%)	0	100	100
33	m	80/106 (76%)	71 (89%)	9 (11%)	5	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
34	o	1254/1748 (72%)	1247 (99%)	7 (1%)	78	81
35	p	993/1027 (97%)	991 (100%)	2 (0%)	87	88
36	q	234/252 (93%)	234 (100%)	0	100	100
37	r	118/126 (94%)	118 (100%)	0	100	100
38	s	191/192 (100%)	191 (100%)	0	100	100
39	t	69/111 (62%)	69 (100%)	0	100	100
40	u	152/153 (99%)	151 (99%)	1 (1%)	76	80
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	104/106 (98%)	104 (100%)	0	100	100
45	z	41/55 (74%)	41 (100%)	0	100	100
All	All	12071/19766 (61%)	11862 (98%)	209 (2%)	52	71

5 of 209 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
15	F	326	HIS
19	J	164	SER
34	o	1291	ASN
15	F	368	THR
16	G	182	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 191 such sidechains are listed below:

Mol	Chain	Res	Type
22	P	229	GLN
15	f	134	HIS
25	S	32	ASN
31	c	50	HIS
32	k	202	ASN

5.3.3 RNA ⓘ

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 19 ligands modelled in this entry, 18 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.

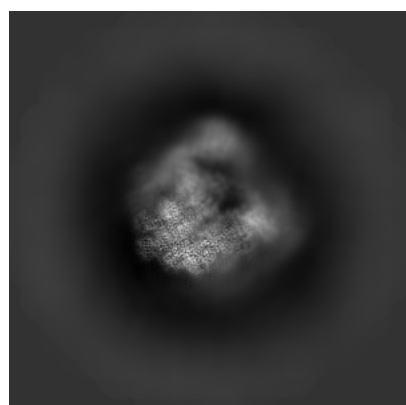
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31111. These allow visual inspection of the internal detail of the map and identification of artifacts.

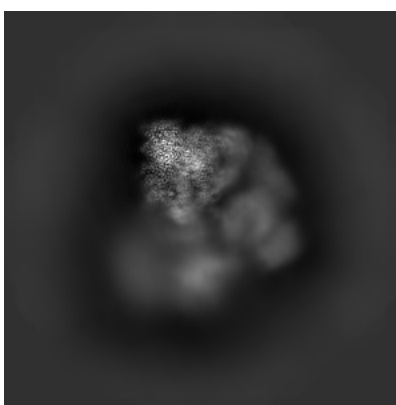
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

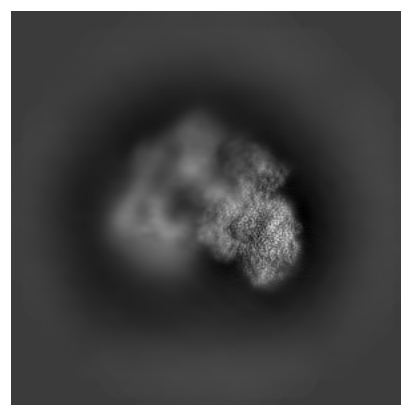
6.1.1 Primary map



X



Y

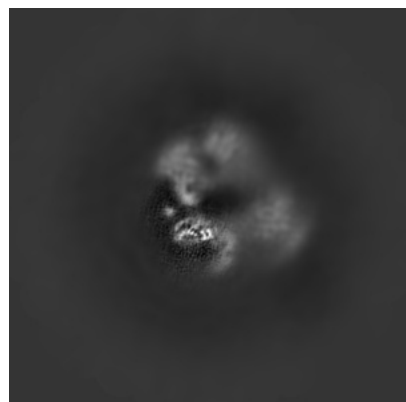


Z

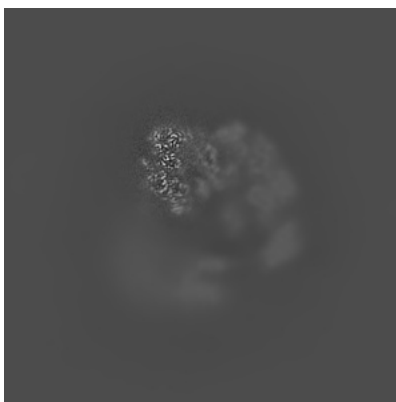
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

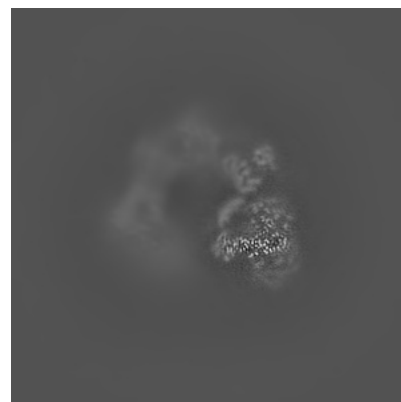
6.2.1 Primary map



X Index: 256



Y Index: 256

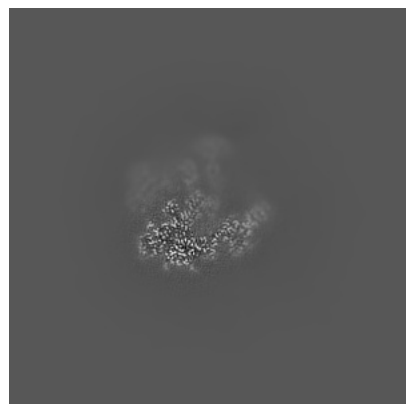


Z Index: 256

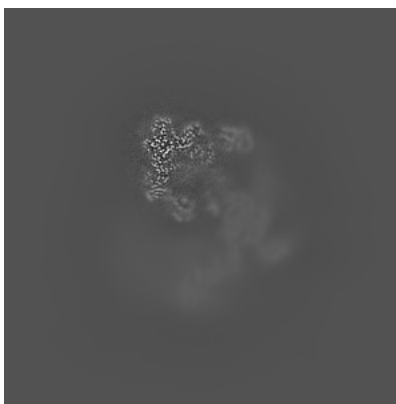
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

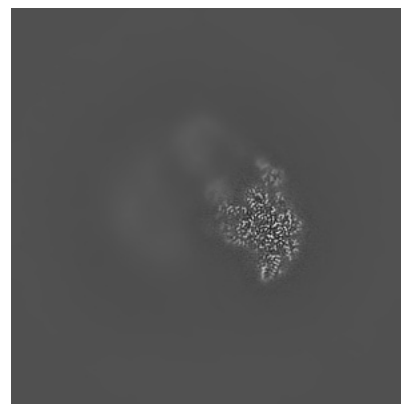
6.3.1 Primary map



X Index: 331



Y Index: 229

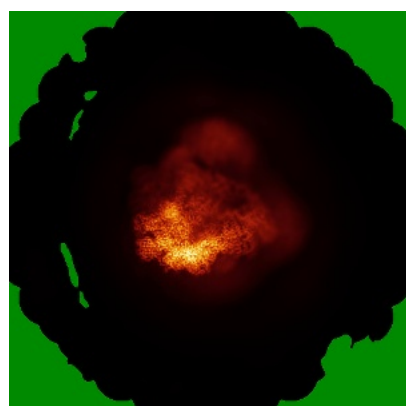


Z Index: 206

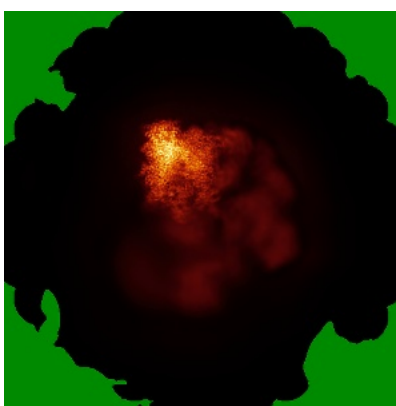
The images above show the largest variance slices of the map in three orthogonal directions.

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

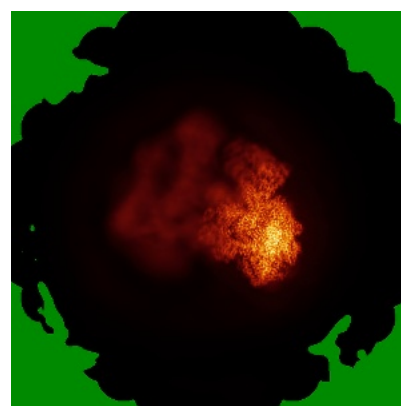
6.4.1 Primary map



X



Y

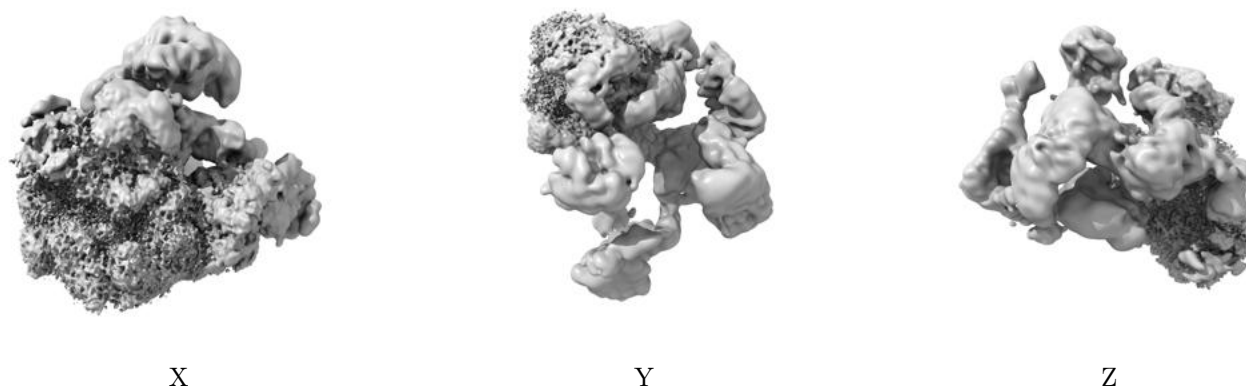


Z

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

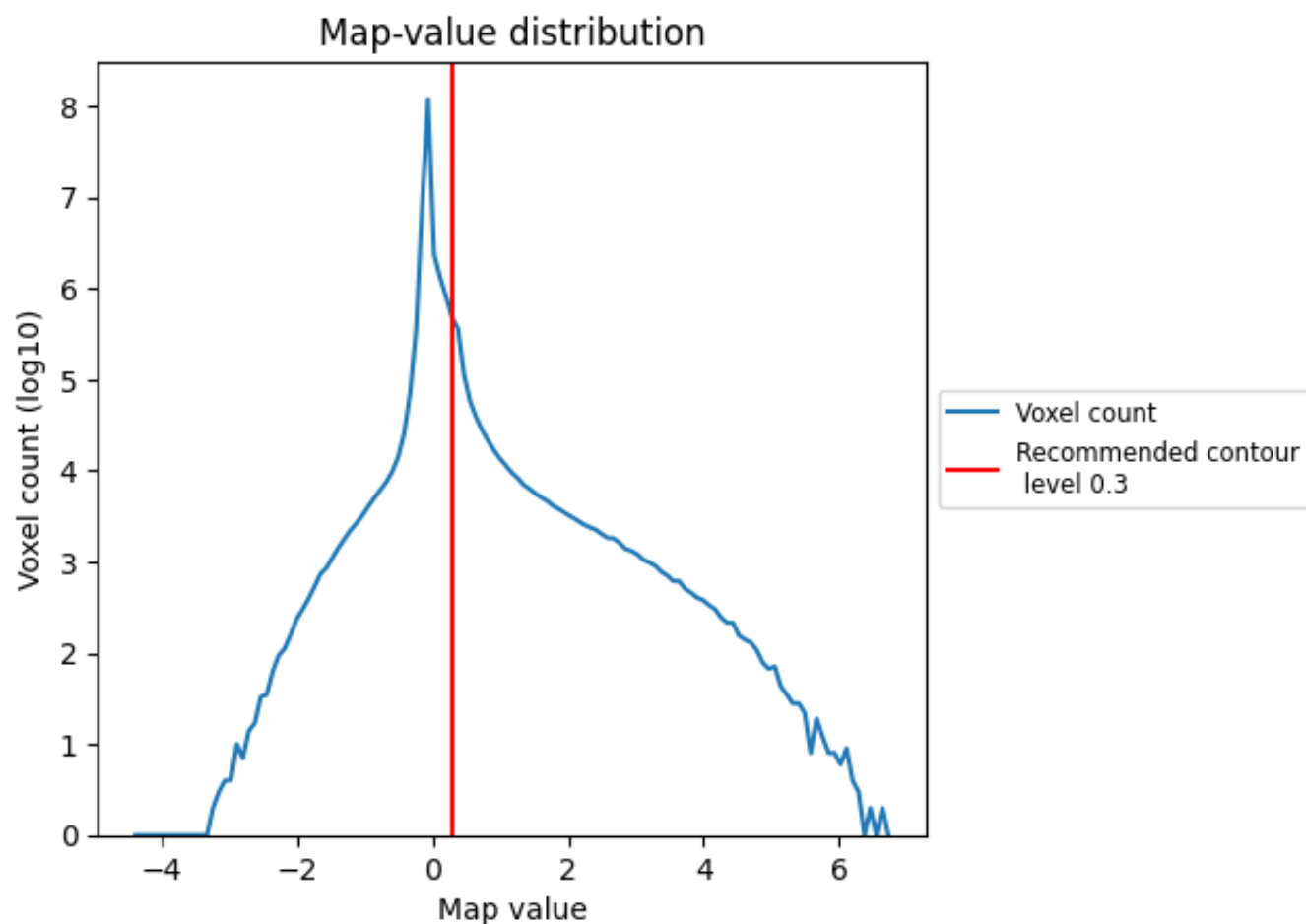
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

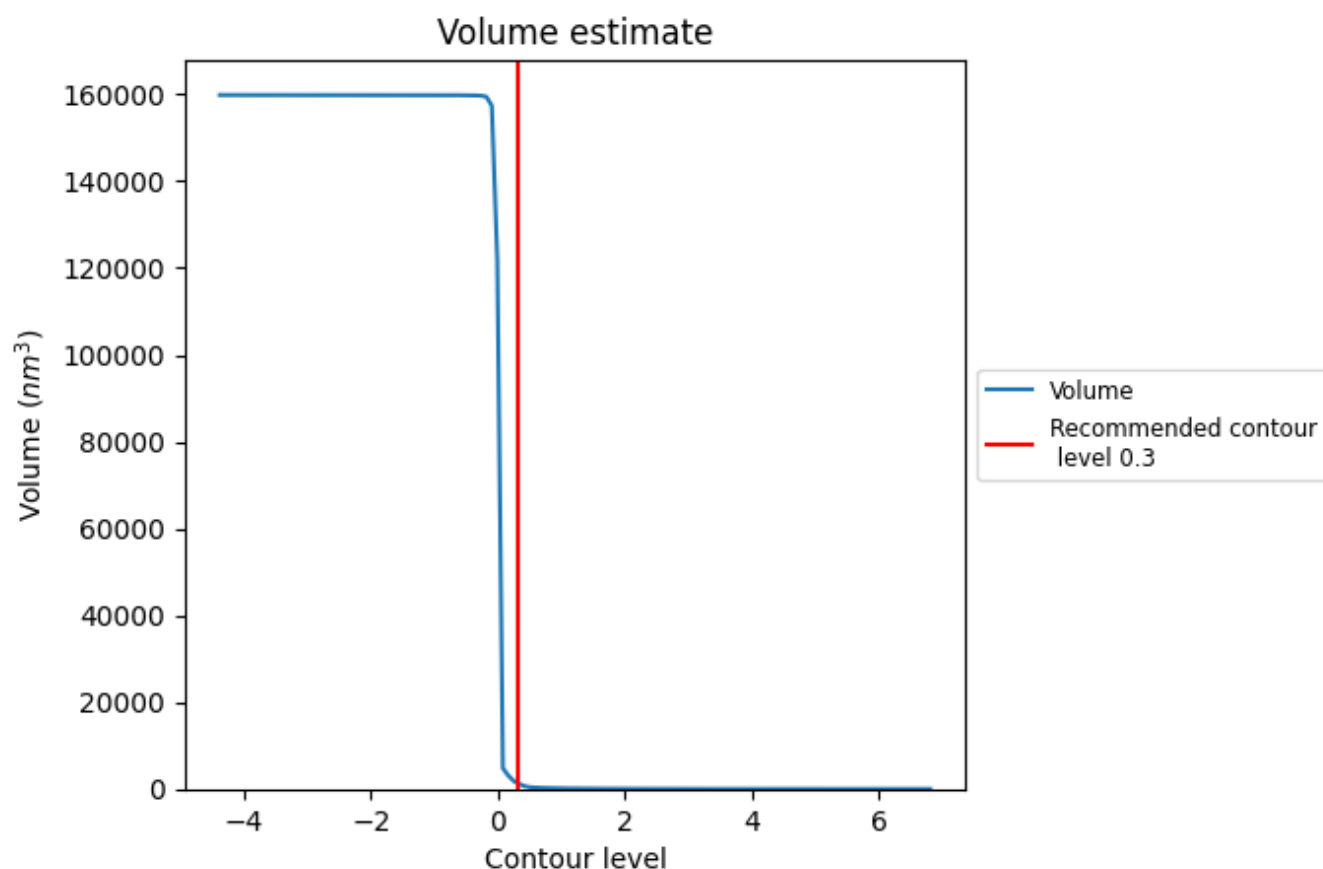
This section contains the results of statistical analysis of the map.

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.

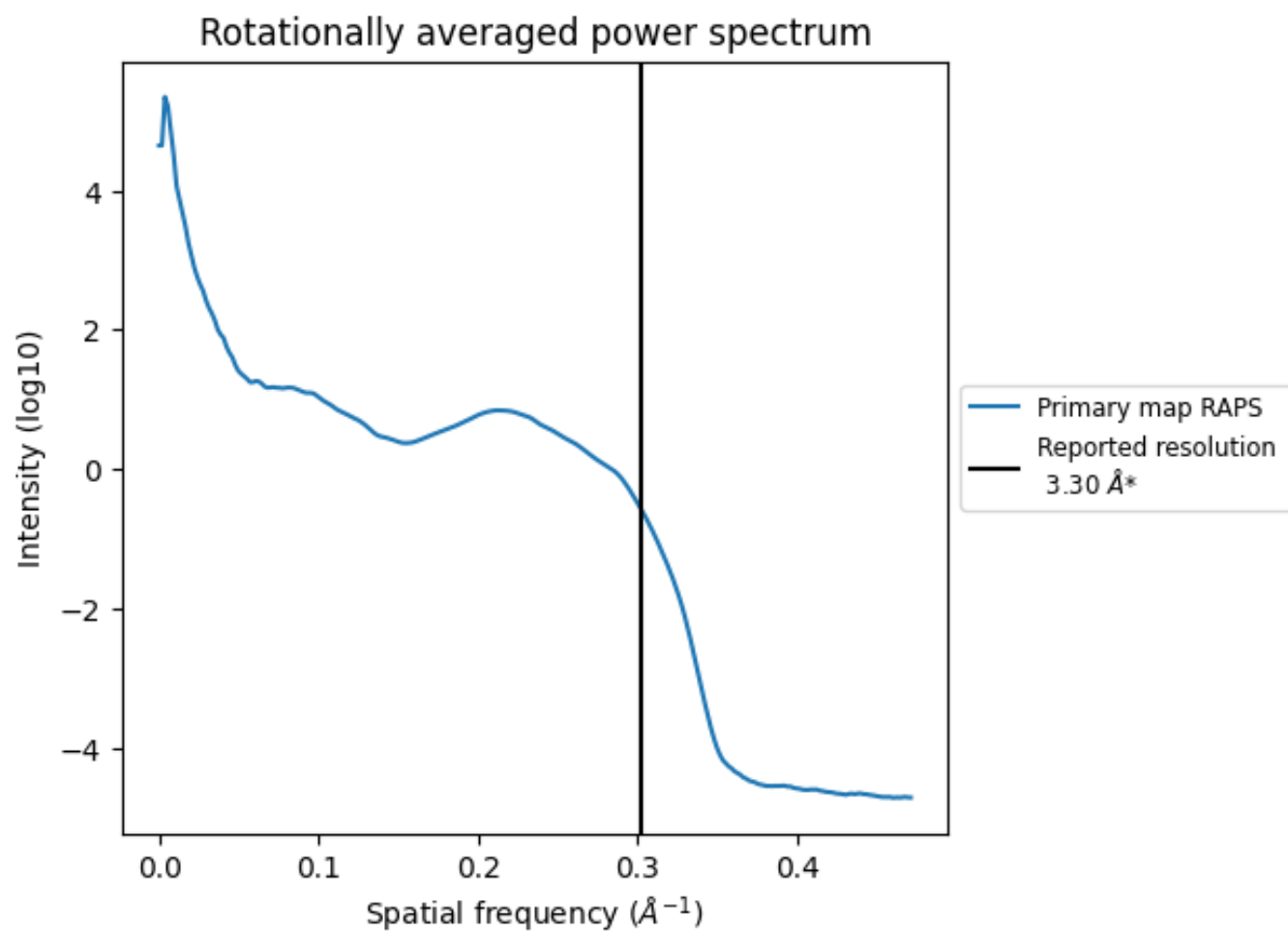
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1427 nm³; this corresponds to an approximate mass of 1289 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 0.303 Å⁻¹

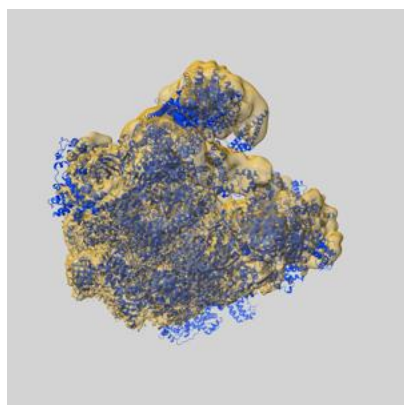
8 Fourier-Shell correlation

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

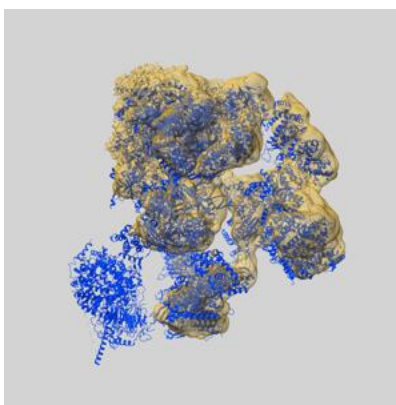
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-31111 and PDB model 7EGB. Per-residue inclusion information can be found in section [3](#) on page [14](#).

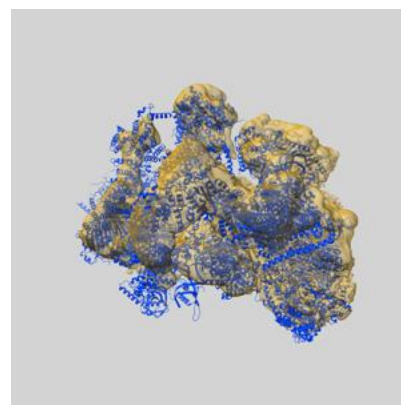
9.1 Map-model overlay [i](#)



X



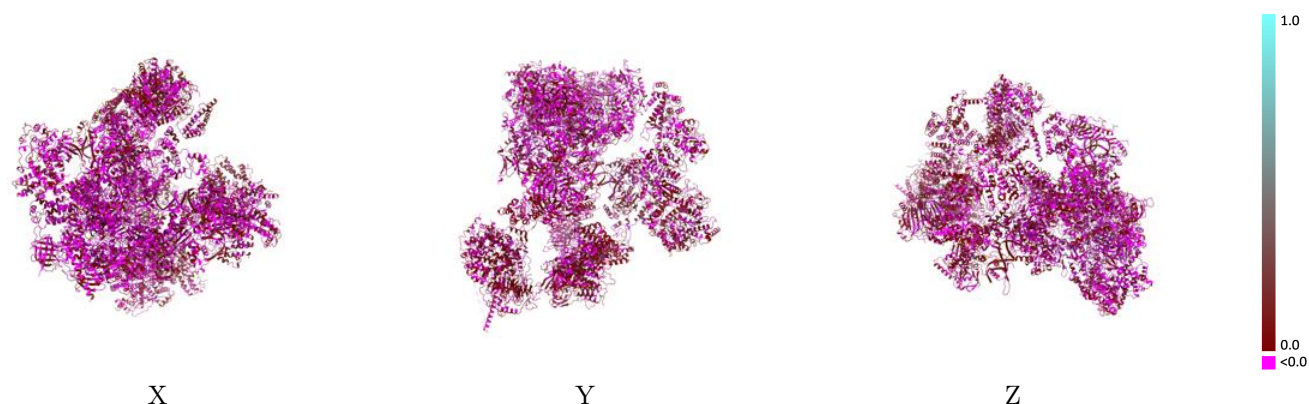
Y



Z

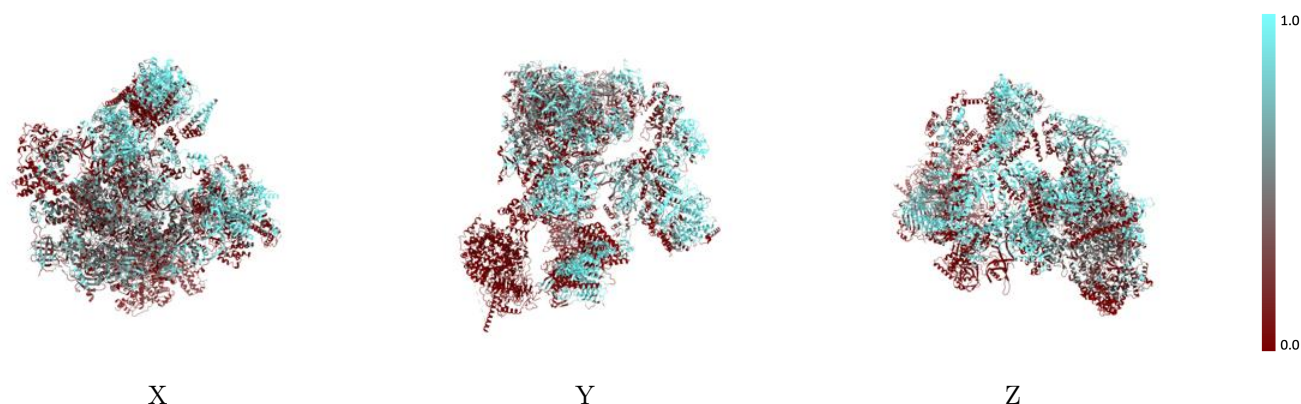
The images above show the 3D surface view of the map at the recommended contour level 0.3 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



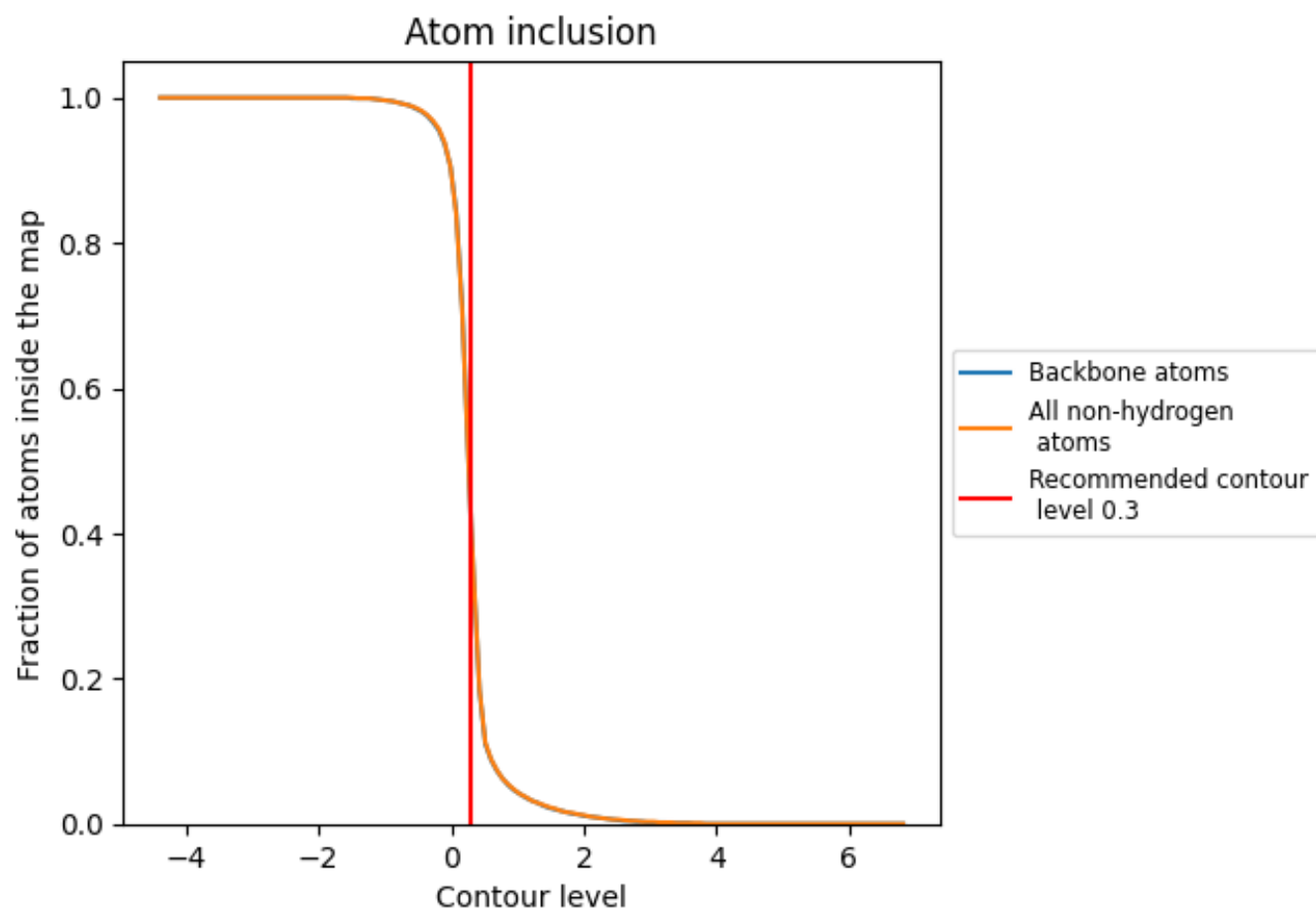
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.3).

9.4 Atom inclusion [i](#)



At the recommended contour level, 41% of all backbone atoms, 41% 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 (0.3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.4140	 0.0120
0	 0.0890	 -0.0020
1	 0.4320	 0.0430
2	 0.5810	 0.0260
3	 0.7830	 0.0500
4	 0.5720	 0.0510
5	 0.3110	 -0.0010
6	 0.6730	 0.0420
7	 0.5860	 0.0310
8	 0.1610	 0.0120
9	 0.2780	 0.0070
A	 0.1510	 0.0250
B	 0.6230	 0.0280
D	 0.5050	 0.0280
E	 0.4020	 0.0310
F	 0.3590	 0.0460
G	 0.0300	 0.0170
H	 0.6440	 0.0390
I	 0.4490	 -0.0180
J	 0.8020	 0.0260
L	 0.3170	 -0.0230
O	 0.5370	 -0.0150
P	 0.7170	 -0.0160
Q	 0.5570	 0.0160
R	 0.4940	 -0.0010
S	 0.5220	 0.0290
T	 0.4100	 0.0270
U	 0.4650	 -0.0380
V	 0.6180	 0.0190
X	 0.4470	 0.0500
Y	 0.4070	 0.0480
c	 0.0000	 0.0340
d	 0.0000	 0.0100
e	 0.0000	 0.0340
f	 0.2680	 0.0380



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Chain	Atom inclusion	Q-score
i	 0.0000	 0.0300
j	 0.0000	 0.0070
k	 0.0000	 0.0130
l	 0.0000	 0.0230
m	 0.0000	 0.0150
o	 0.4030	 -0.0240
p	 0.4300	 -0.0300
q	 0.4700	 0.0000
r	 0.5270	 -0.0370
s	 0.4330	 -0.0260
t	 0.3740	 -0.0310
u	 0.6640	 -0.0100
v	 0.4060	 -0.0110
w	 0.3780	 -0.0350
x	 0.4220	 0.0140
y	 0.4850	 -0.0190
z	 0.4710	 0.0040