



wwPDB EM Validation Summary Report ⓘ

Mar 9, 2026 – 06:48 AM UTC

PDB ID : 7QYB / pdb_00007qyb
EMDB ID : EMD-14211
Title : Proteasome-ZFAND5 Complex Z-C state
Authors : Zhu, Y.; Lu, Y.
Deposited on : 2022-01-27
Resolution : 4.10 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

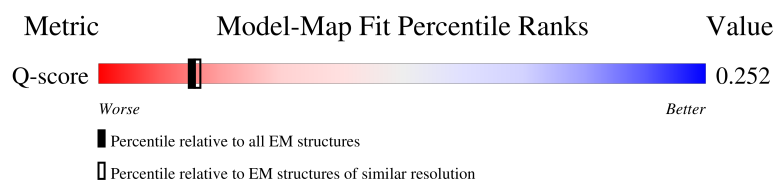
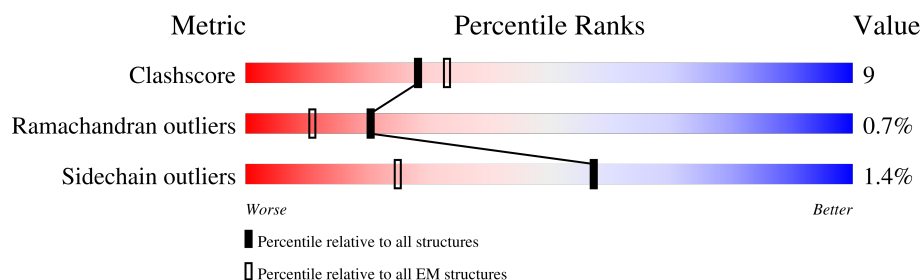
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.10 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	6458 (3.60 - 4.60)

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	U	953	34% 72% 17% 8%
2	V	533	48% 67% 22% 10%
3	W	456	41% 75% 22% •
4	X	422	25% 77% 12% 10%

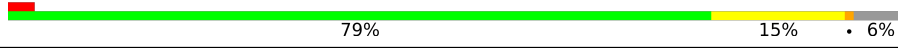
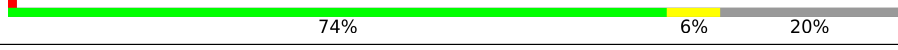
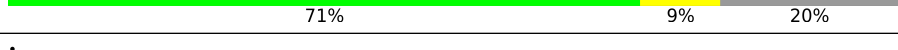
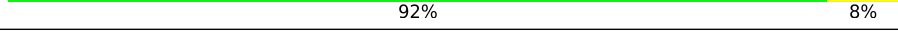
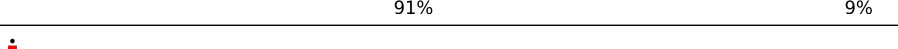
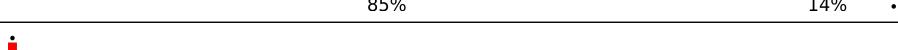
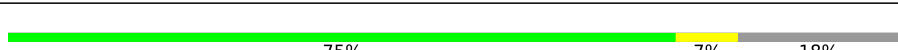
Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
5	Y	389	
6	Z	324	
7	a	376	
8	b	377	
9	c	309	
10	d	349	
11	e	70	
12	f	908	
13	A	433	
14	B	440	
15	C	398	
16	D	418	
17	E	403	
18	F	439	
19	G	245	
19	g	245	
20	H	233	
20	h	233	
21	I	260	
21	i	260	
22	J	247	
22	j	247	
23	K	240	
23	k	240	
24	L	268	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
24	l	268	
25	M	254	
25	m	254	
26	N	238	
26	n	238	
27	O	276	
27	o	276	
28	P	204	
28	p	204	
29	Q	201	
29	q	201	
30	R	262	
30	r	262	
31	S	240	
31	s	240	
32	T	263	
32	t	263	

2 Entry composition

There are 36 unique types of molecules in this entry. The entry contains 105037 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 26S proteasome non-ATPase regulatory subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	U	872	Total	C	N	O	S	0	0
			6828	4328	1157	1298	45		

- Molecule 2 is a protein called 26S proteasome non-ATPase regulatory subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	V	480	Total	C	N	O	S	0	0
			3852	2444	684	710	14		

- Molecule 3 is a protein called 26S proteasome non-ATPase regulatory subunit 12.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	W	456	Total	C	N	O	S	0	0
			3703	2339	635	704	25		

- Molecule 4 is a protein called 26S proteasome non-ATPase regulatory subunit 11.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	X	380	Total	C	N	O	S	0	0
			3009	1918	509	570	12		

- Molecule 5 is a protein called 26S proteasome non-ATPase regulatory subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	Y	378	Total	C	N	O	S	0	0
			3115	1987	533	578	17		

- Molecule 6 is a protein called 26S proteasome non-ATPase regulatory subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	Z	286	Total	C	N	O	S	0	0
			2276	1454	390	427	5		

- Molecule 7 is a protein called 26S proteasome non-ATPase regulatory subunit 13.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	a	373	Total	C	N	O	S	0	0
			2995	1911	510	559	15		

- Molecule 8 is a protein called 26S proteasome non-ATPase regulatory subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	b	191	Total	C	N	O	S	0	0
			1458	910	261	279	8		

- Molecule 9 is a protein called 26S proteasome non-ATPase regulatory subunit 14.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	c	287	Total	C	N	O	S	0	0
			2260	1430	389	422	19		

- Molecule 10 is a protein called 26S proteasome non-ATPase regulatory subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	d	257	Total	C	N	O	S	0	0
			2116	1371	346	390	9		

- Molecule 11 is a protein called 26S proteasome complex subunit SEM1.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	e	40	Total	C	N	O	S	0	0
			334	200	55	77	2		

- Molecule 12 is a protein called 26S proteasome non-ATPase regulatory subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	f	889	Total	C	N	O	S	0	0
			6866	4315	1174	1331	46		

- Molecule 13 is a protein called 26S protease regulatory subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	A	413	Total	C	N	O	S	0	0
			3229	2034	566	611	18		

- Molecule 14 is a protein called 26S protease regulatory subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	B	411	Total	C	N	O	S	0	0
			3207	2022	548	622	15		

- Molecule 15 is a protein called Isoform 2 of 26S proteasome regulatory subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	C	396	Total	C	N	O	S	0	0
			3105	1954	558	576	17		

- Molecule 16 is a protein called 26S protease regulatory subunit 6B.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	D	380	Total	C	N	O	S	0	0
			3040	1923	524	580	13		

- Molecule 17 is a protein called 26S proteasome regulatory subunit 10B.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	E	389	Total	C	N	O	S	0	0
			3097	1947	552	581	17		

- Molecule 18 is a protein called 26S protease regulatory subunit 6A.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	F	395	Total	C	N	O	S	0	0
			3098	1951	533	596	18		

- Molecule 19 is a protein called Proteasome subunit alpha type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	G	239	Total	C	N	O	S	0	0
			1820	1157	304	346	13		
19	g	240	Total	C	N	O	S	0	0
			1826	1160	305	348	13		

- Molecule 20 is a protein called Proteasome subunit alpha type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	H	230	Total	C	N	O	S	0	0
			1692	1073	285	329	5		
20	h	232	Total	C	N	O	S	0	0
			1708	1081	289	333	5		

- Molecule 21 is a protein called Proteasome subunit alpha type-4.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	I	248	Total	C	N	O	S	0	0
			1895	1195	324	368	8		
21	i	250	Total	C	N	O	S	0	0
			1912	1204	329	371	8		

- Molecule 22 is a protein called Proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	J	239	Total	C	N	O	S	0	0
			1704	1056	308	335	5		
22	j	239	Total	C	N	O	S	0	0
			1704	1056	308	335	5		

- Molecule 23 is a protein called Proteasome subunit alpha type-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	K	228	Total	C	N	O	S	0	0
			1729	1086	284	349	10		
23	k	228	Total	C	N	O	S	0	0
			1722	1080	284	348	10		

- Molecule 24 is a protein called Isoform Long of Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	L	238	Total	C	N	O	S	0	0
			1850	1159	334	346	11		
24	l	238	Total	C	N	O	S	0	0
			1850	1159	334	346	11		

- Molecule 25 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	M	240	Total	C	N	O	S	0	0
			1856	1178	314	353	11		
25	m	240	Total	C	N	O	S	0	0
			1856	1178	314	353	11		

- Molecule 26 is a protein called Proteasome subunit beta type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	N	191	Total	C	N	O	S	0	0
			1430	893	245	280	12		
26	n	191	Total	C	N	O	S	0	0
			1430	893	245	280	12		

- Molecule 27 is a protein called Proteasome subunit beta type-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	O	220	Total	C	N	O	S	0	0
			1643	1033	280	318	12		
27	o	220	Total	C	N	O	S	0	0
			1643	1033	280	318	12		

- Molecule 28 is a protein called Proteasome subunit beta type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	P	204	Total	C	N	O	S	0	0
			1585	1010	262	294	19		
28	p	204	Total	C	N	O	S	0	0
			1585	1010	262	294	19		

- Molecule 29 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	Q	199	Total	C	N	O	S	0	0
			1570	1006	265	290	9		
29	q	199	Total	C	N	O	S	0	0
			1570	1006	265	290	9		

- Molecule 30 is a protein called Proteasome subunit beta type-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	R	201	Total	C	N	O	S	0	0
			1548	974	273	292	9		
30	r	201	Total	C	N	O	S	0	0
			1548	974	273	292	9		

- Molecule 31 is a protein called Proteasome subunit beta type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	S	213	Total	C	N	O	S	0	0
			1641	1036	282	313	10		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					AltConf	Trace
31	s	213	Total	C	N	O	S	0	0
			1641	1036	282	313	10		

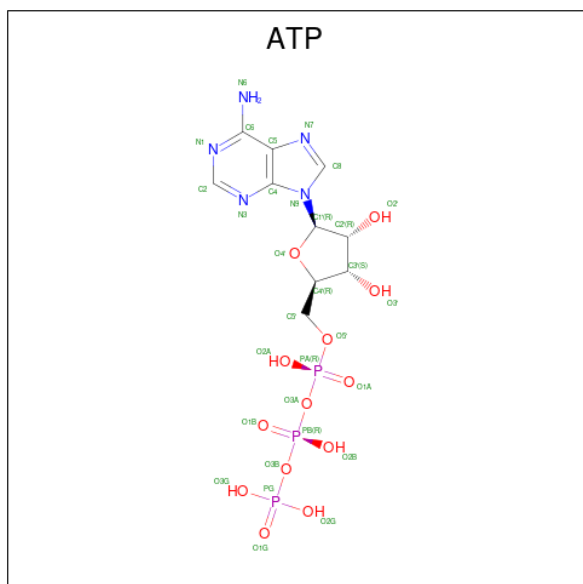
- Molecule 32 is a protein called Proteasome subunit beta type-4.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	T	215	Total	C	N	O	S	0	0
			1667	1052	285	318	12		
32	t	215	Total	C	N	O	S	0	0
			1667	1052	285	318	12		

- Molecule 33 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
33	c	1	Total	Zn	0
			1	1	

- Molecule 34 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
34	A	1	Total	C	N	O	P	0
			31	10	5	13	3	
34	B	1	Total	C	N	O	P	0
			31	10	5	13	3	

Continued on next page...

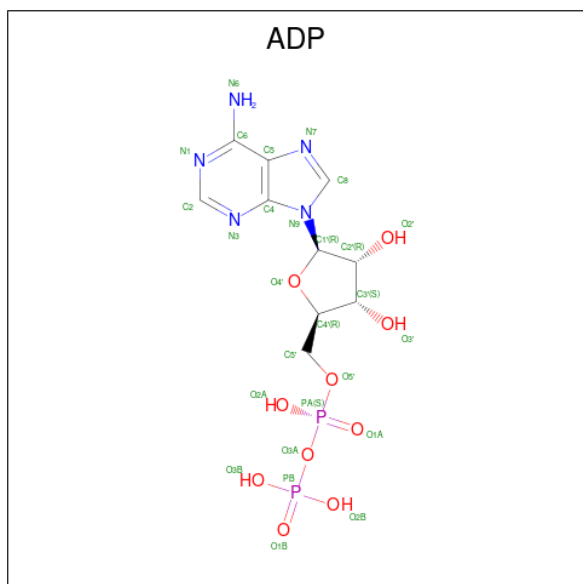
Continued from previous page...

Mol	Chain	Residues	Atoms					AltConf
34	C	1	Total	C	N	O	P	0
			31	10	5	13	3	
34	F	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 35 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
35	A	1	Total	Mg	0
			1	1	
35	B	1	Total	Mg	0
			1	1	
35	C	1	Total	Mg	0
			1	1	
35	D	1	Total	Mg	0
			1	1	
35	F	1	Total	Mg	0
			1	1	

- Molecule 36 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂) (labeled as "Ligand of Interest" by depositor).

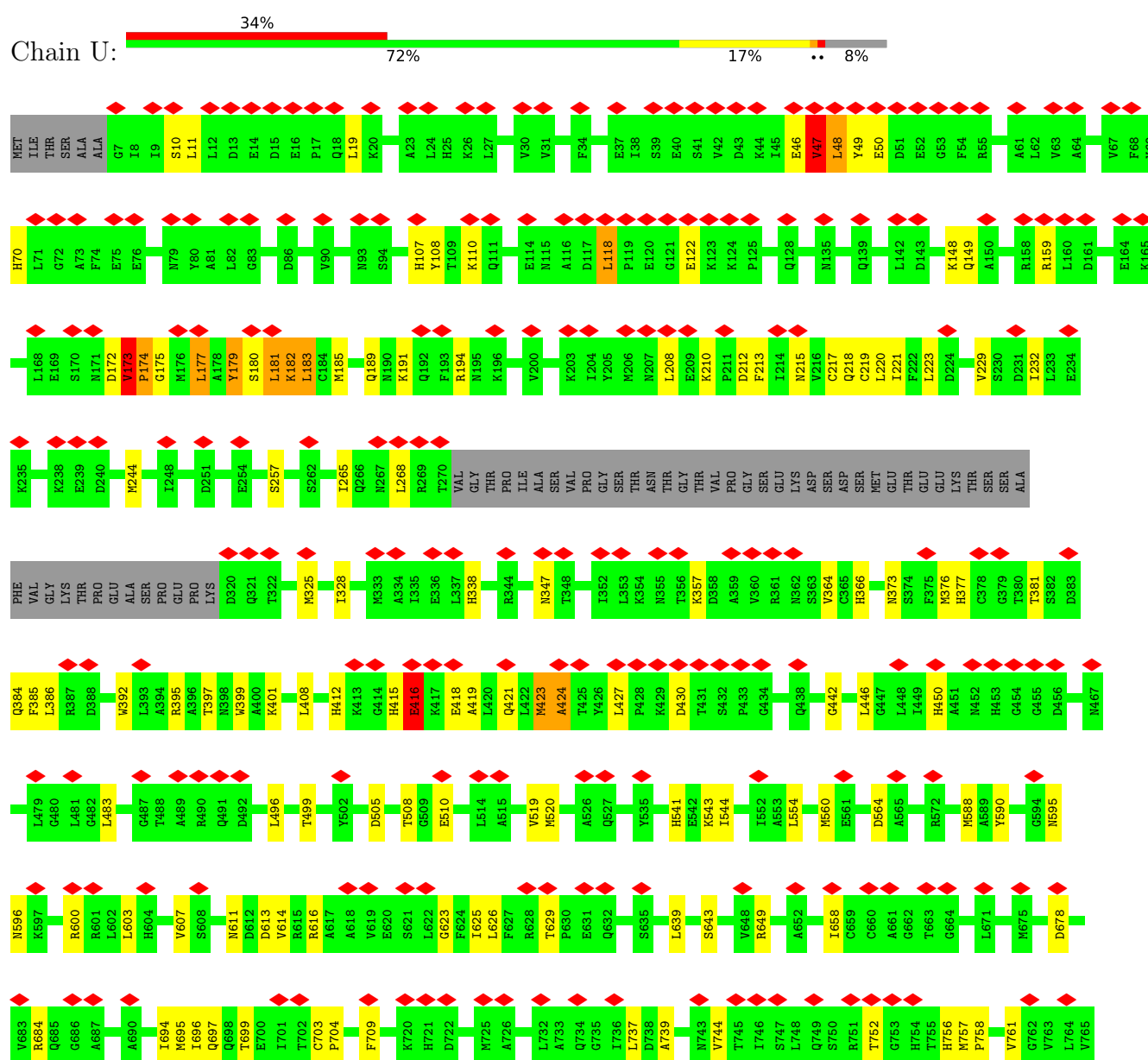


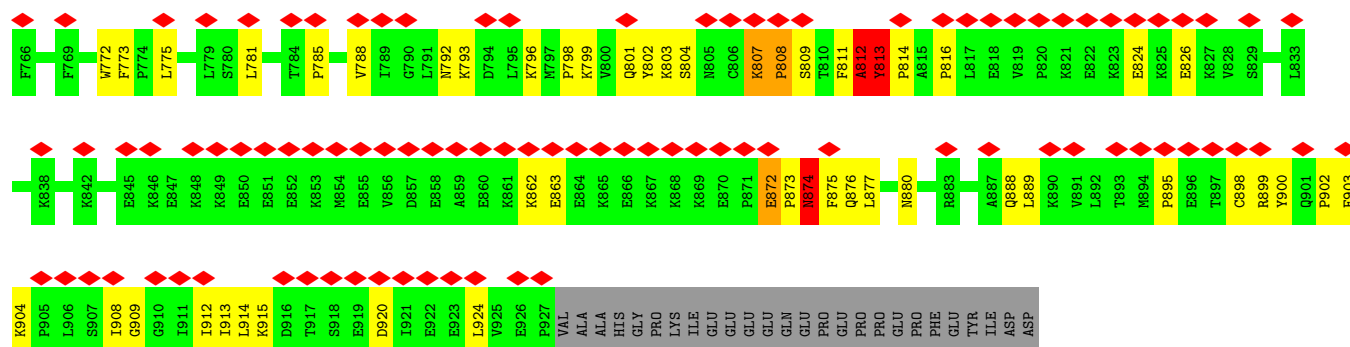
Mol	Chain	Residues	Atoms					AltConf
36	D	1	Total	C	N	O	P	0
			27	10	5	10	2	

3 Residue-property plots [i](#)

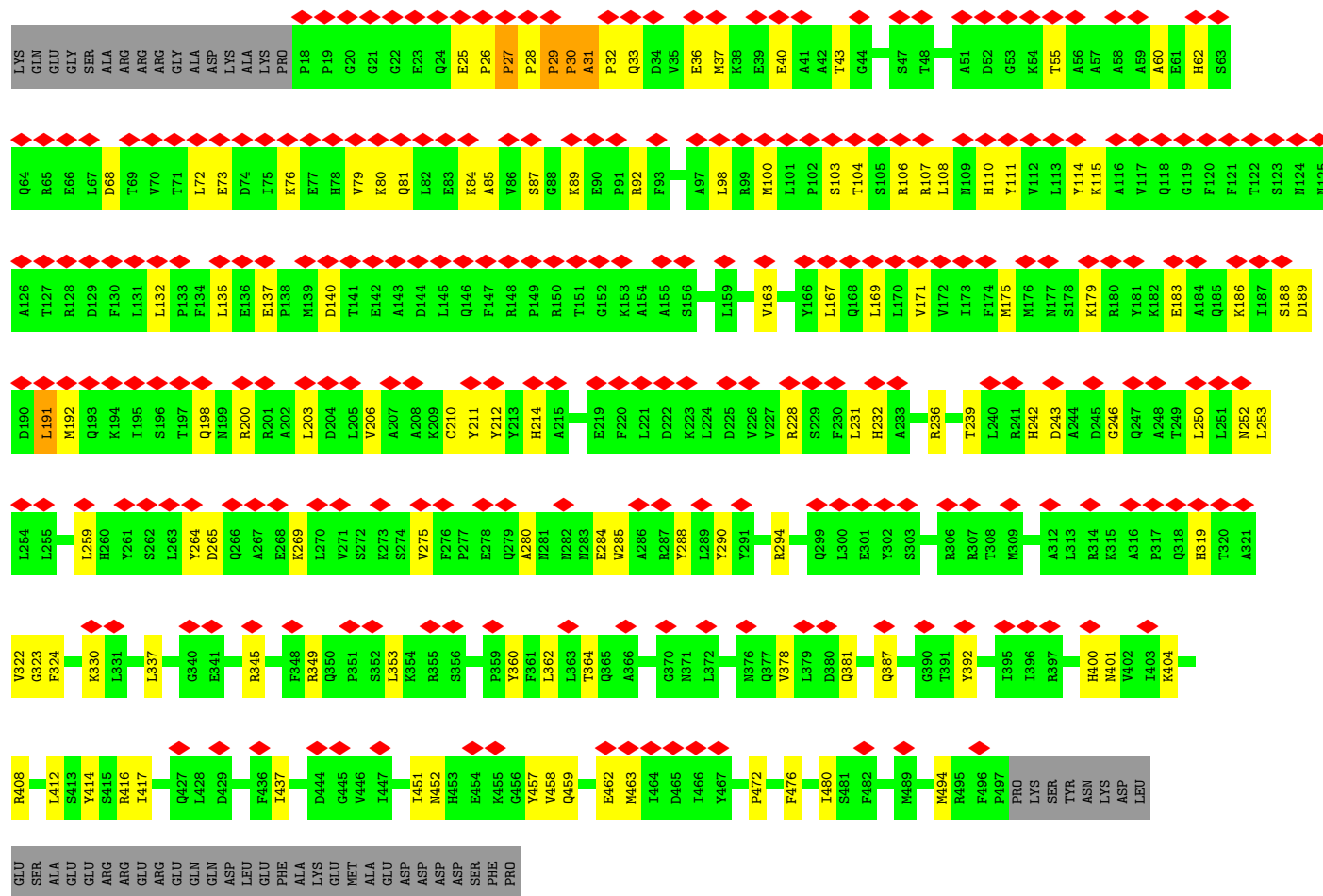
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 26S proteasome non-ATPase regulatory subunit 1

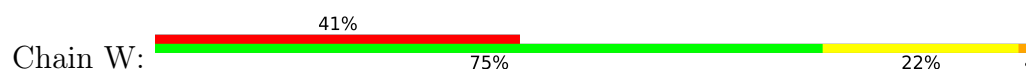


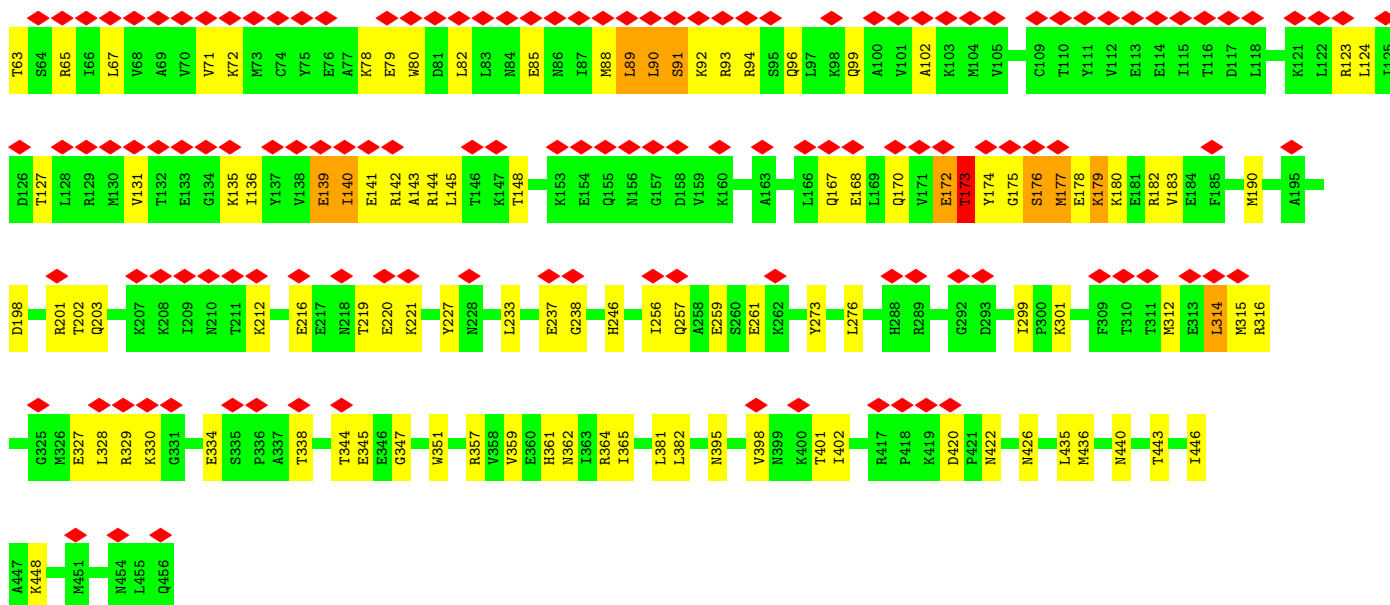


• Molecule 2: 26S proteasome non-ATPase regulatory subunit 3

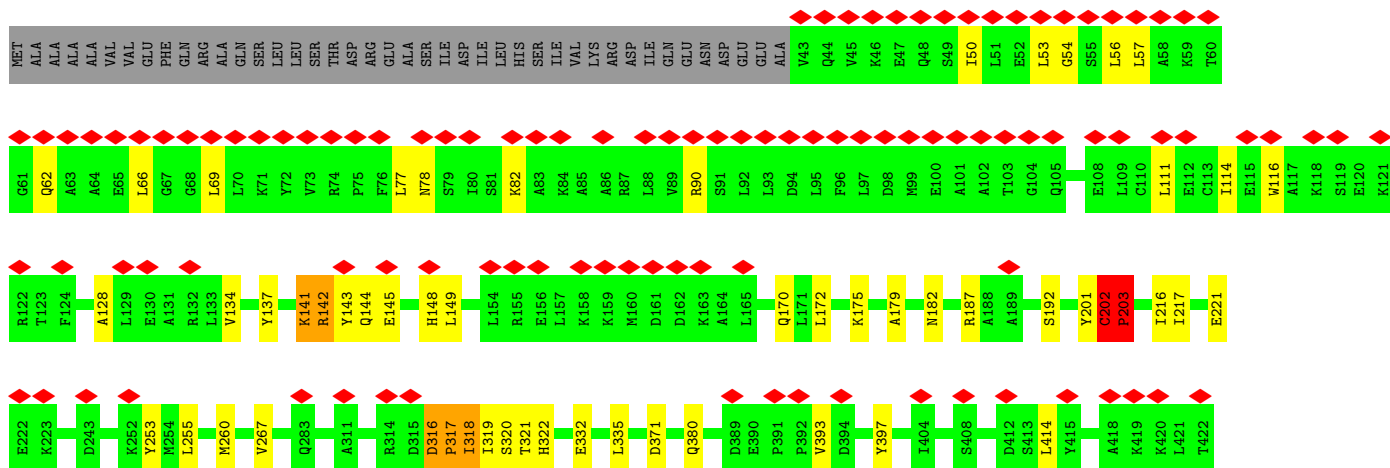
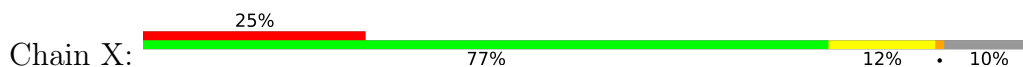


• Molecule 3: 26S proteasome non-ATPase regulatory subunit 12

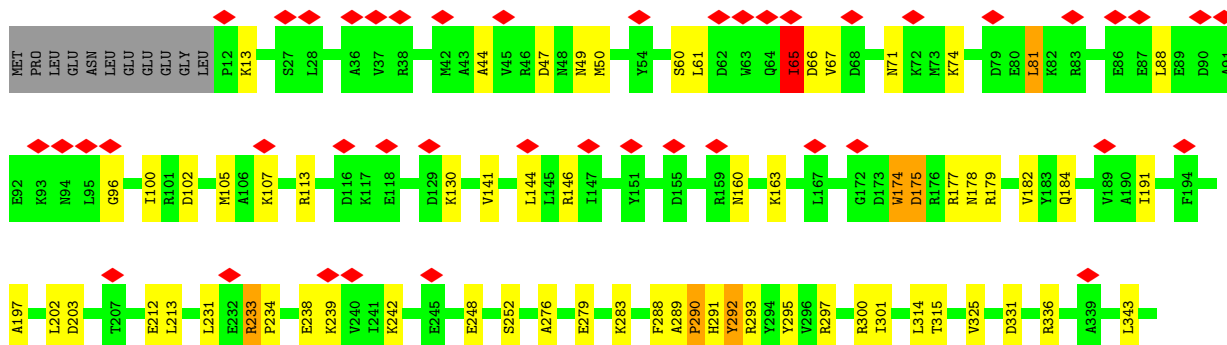
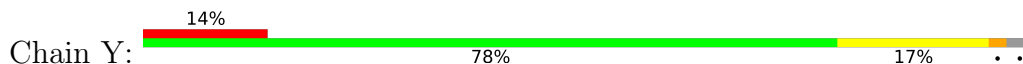


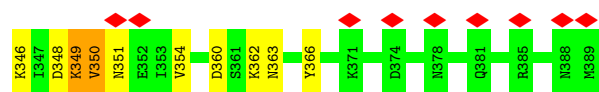


• Molecule 4: 26S proteasome non-ATPase regulatory subunit 11

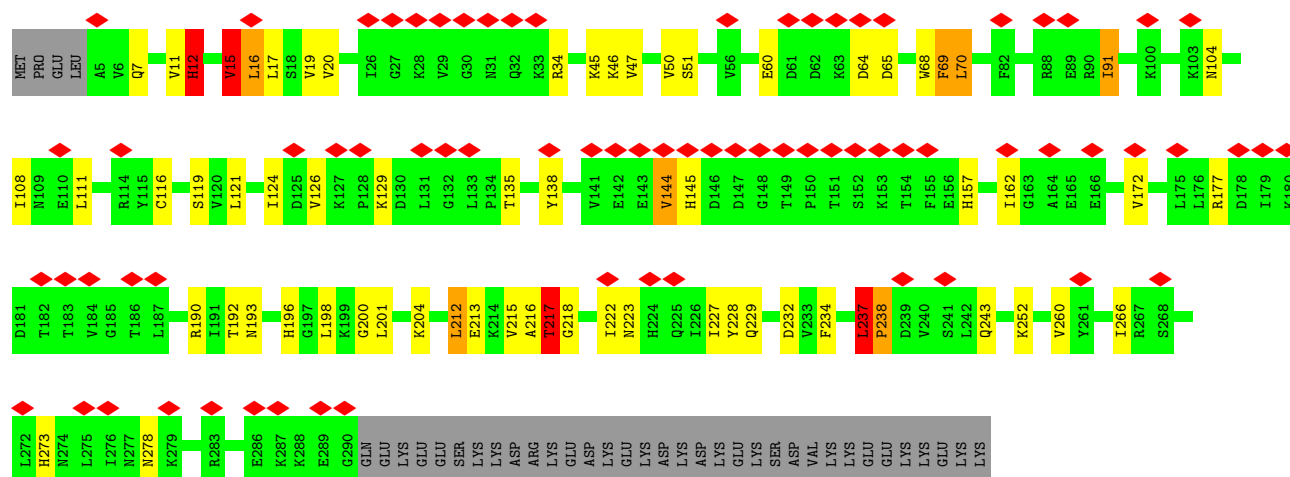


• Molecule 5: 26S proteasome non-ATPase regulatory subunit 6

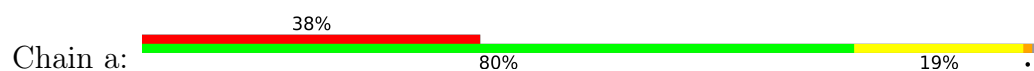




- Molecule 6: 26S proteasome non-ATPase regulatory subunit 7

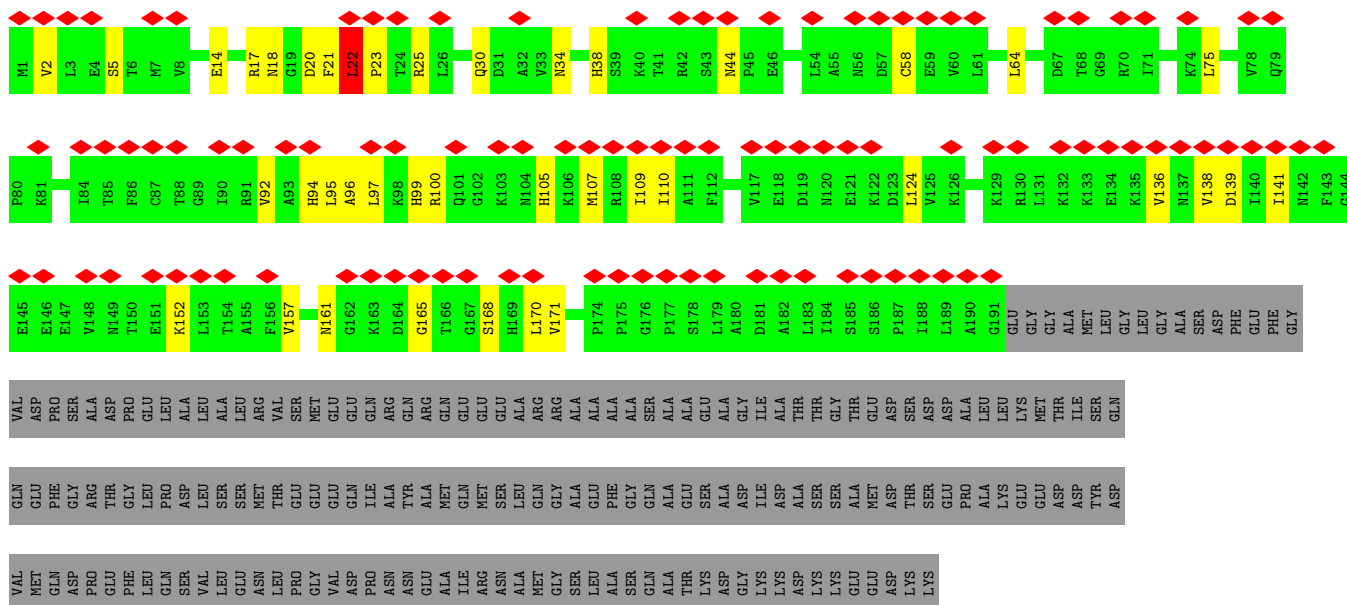


- Molecule 7: 26S proteasome non-ATPase regulatory subunit 13

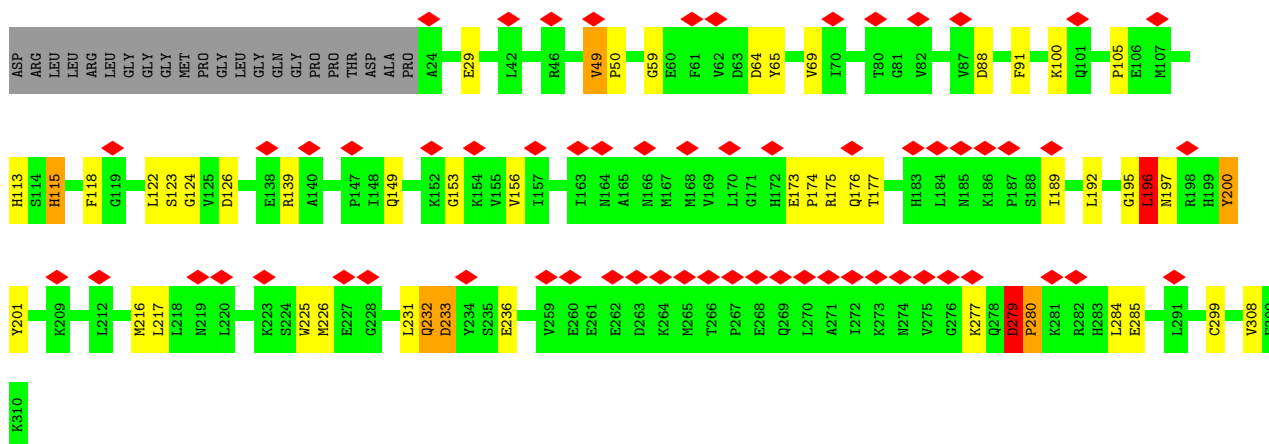
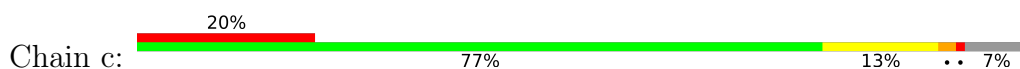


- Molecule 8: 26S proteasome non-ATPase regulatory subunit 4

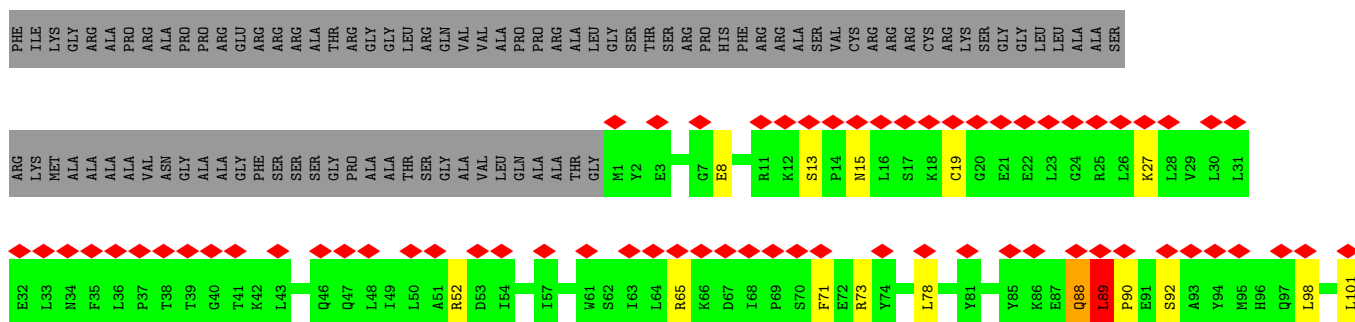
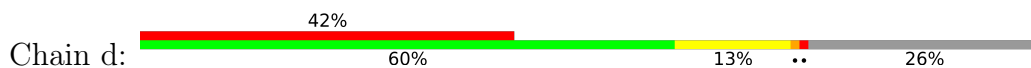


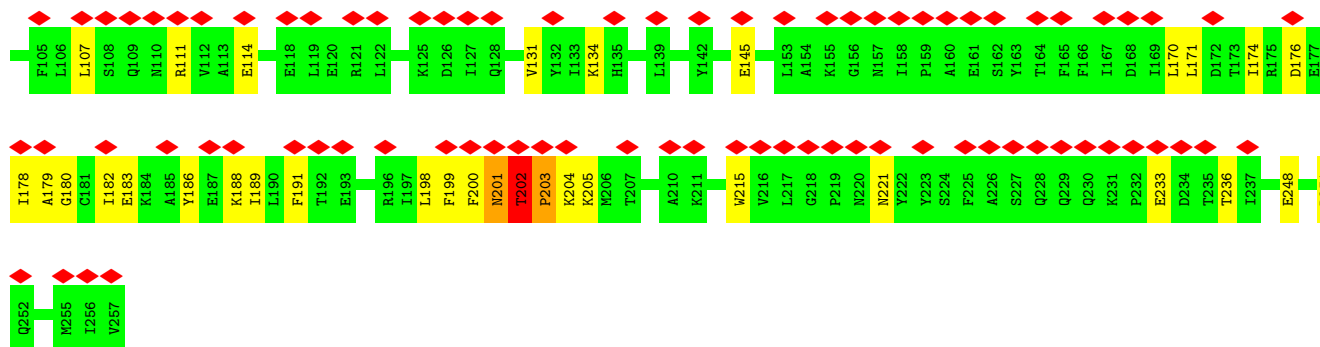


- Molecule 9: 26S proteasome non-ATPase regulatory subunit 14



- Molecule 10: 26S proteasome non-ATPase regulatory subunit 8

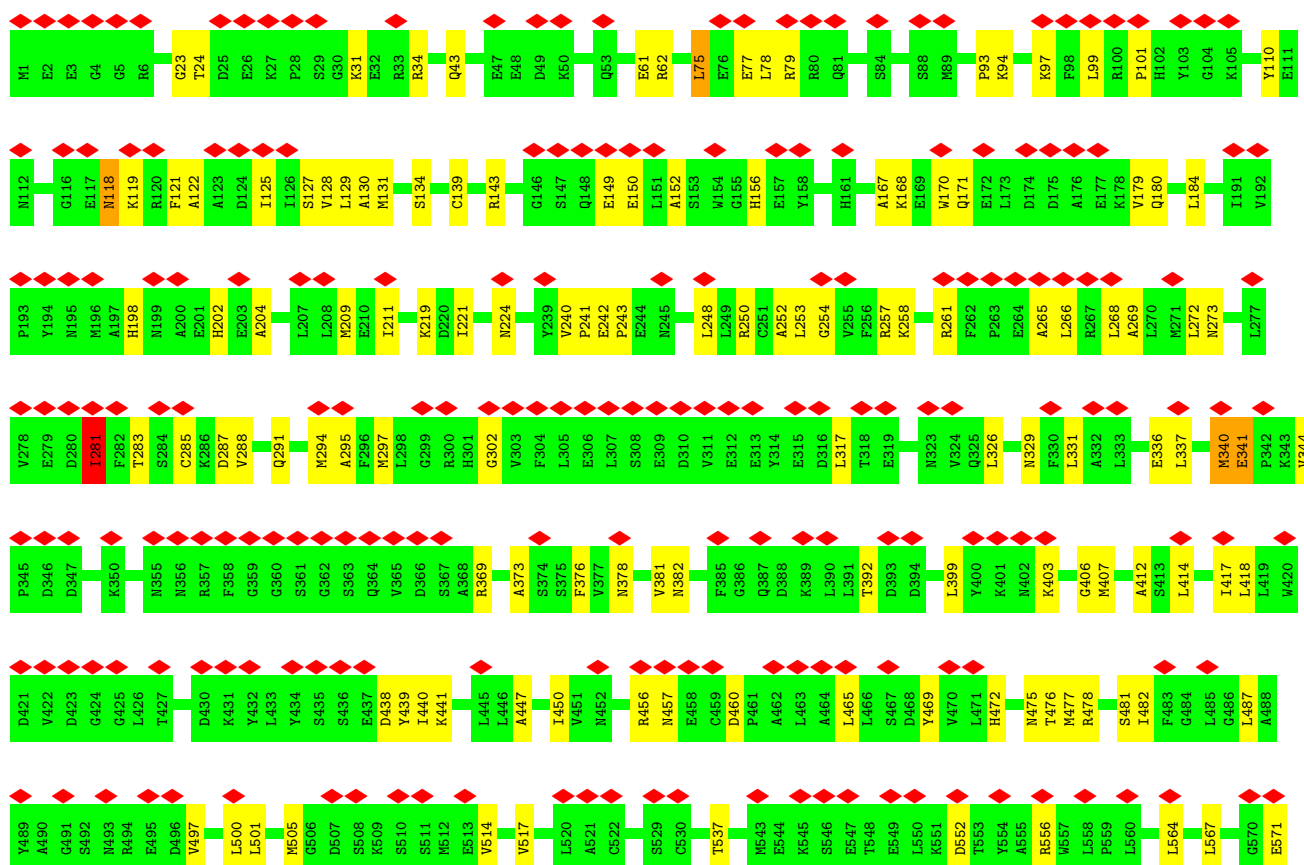


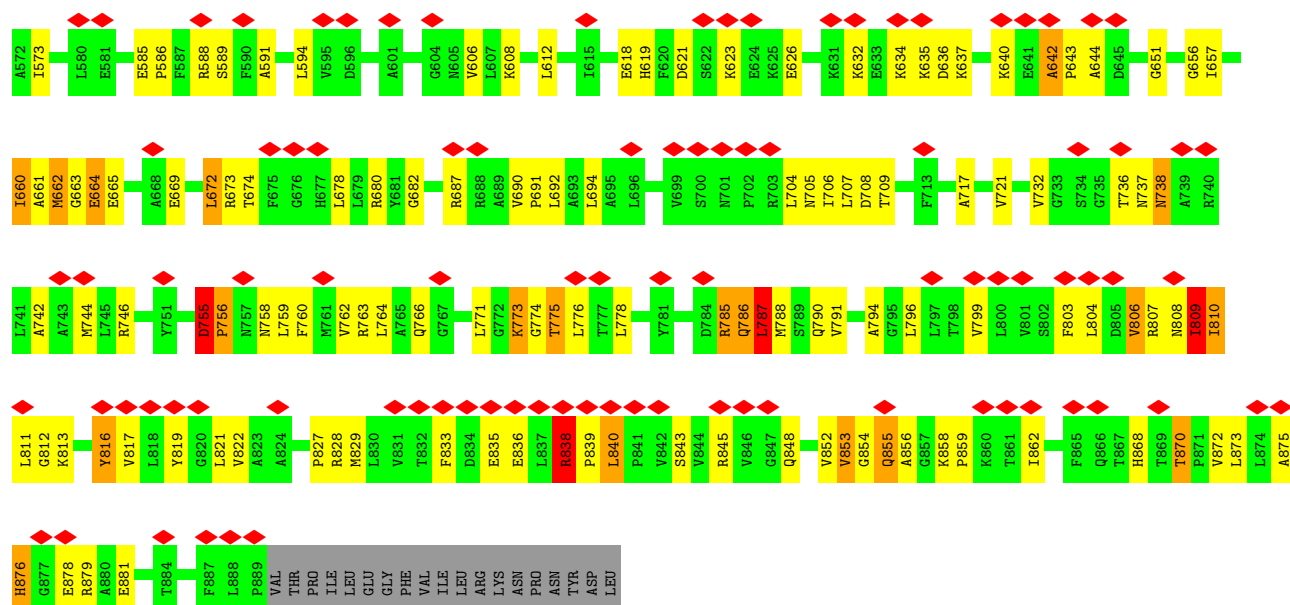


- Molecule 11: 26S proteasome complex subunit SEM1

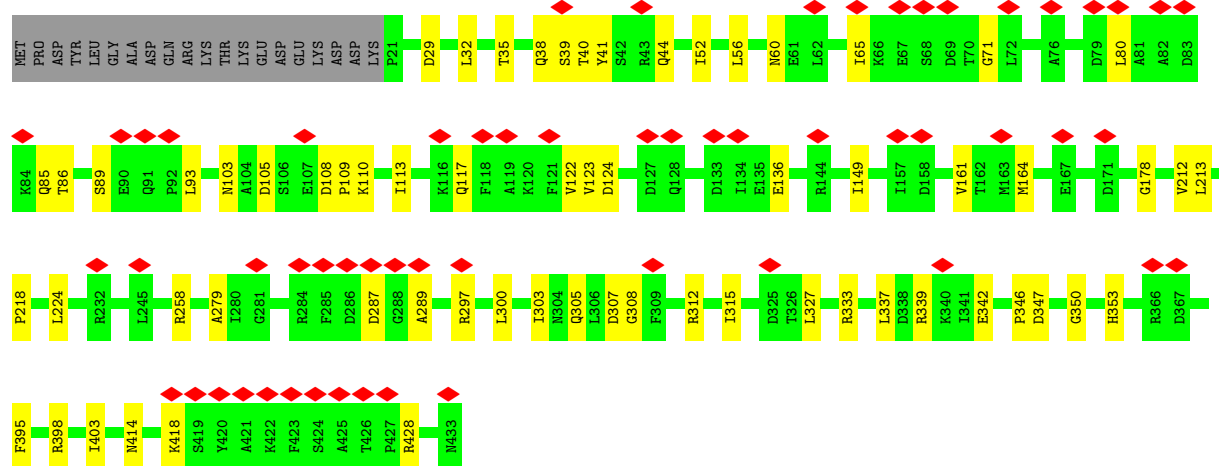
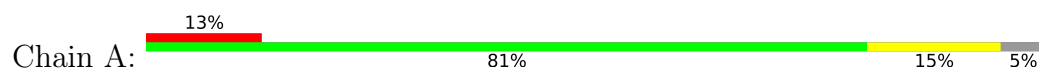


- Molecule 12: 26S proteasome non-ATPase regulatory subunit 2

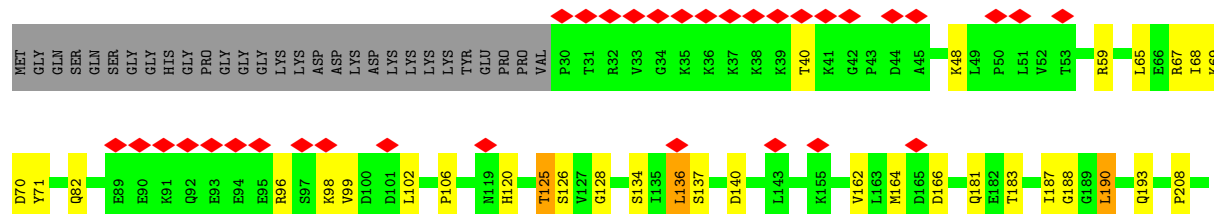
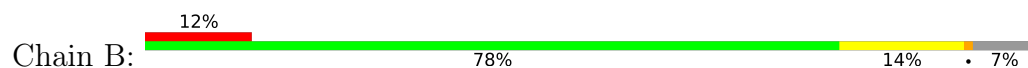


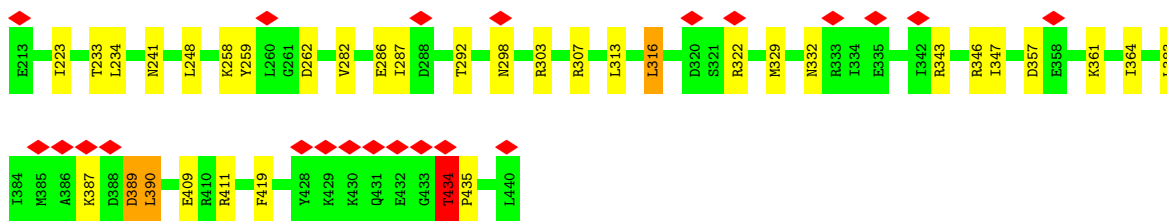


• Molecule 13: 26S protease regulatory subunit 7

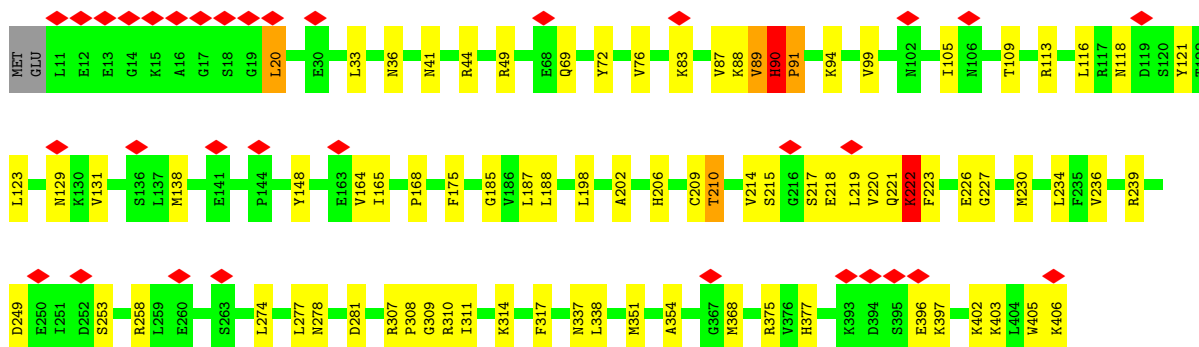
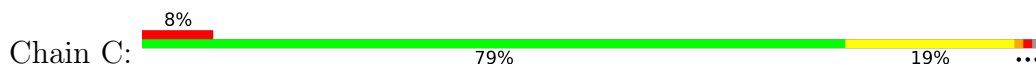


• Molecule 14: 26S protease regulatory subunit 4

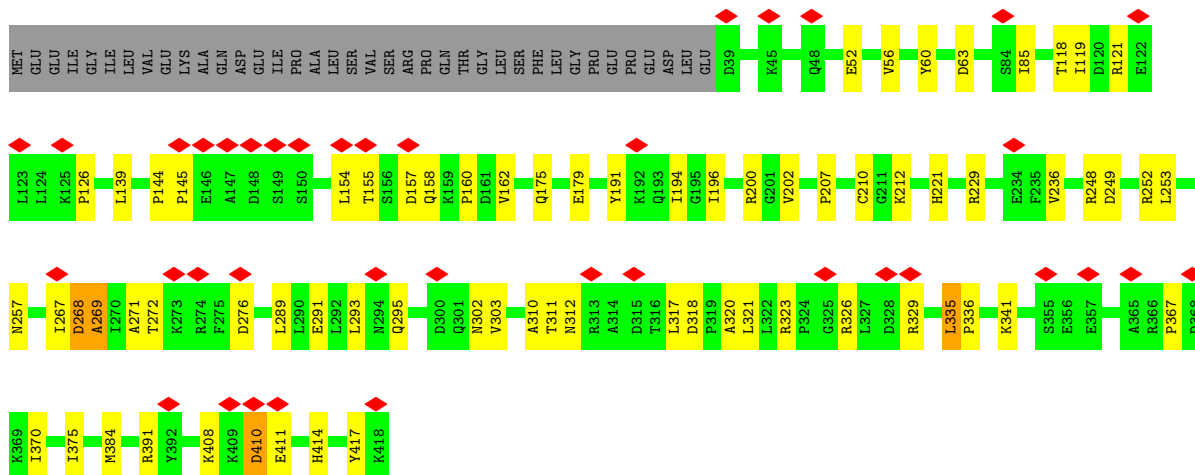
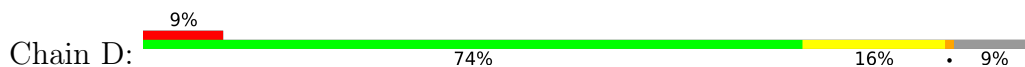




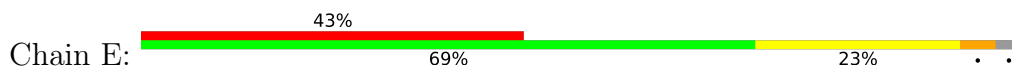
- Molecule 15: Isoform 2 of 26S proteasome regulatory subunit 8

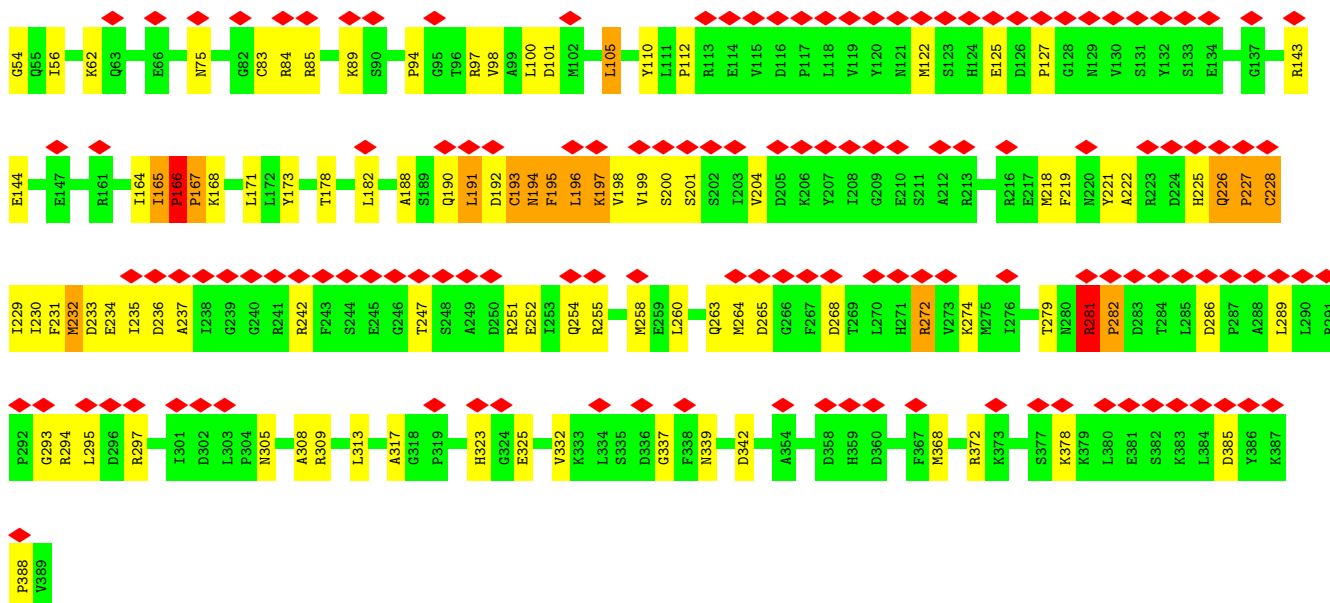


- Molecule 16: 26S protease regulatory subunit 6B



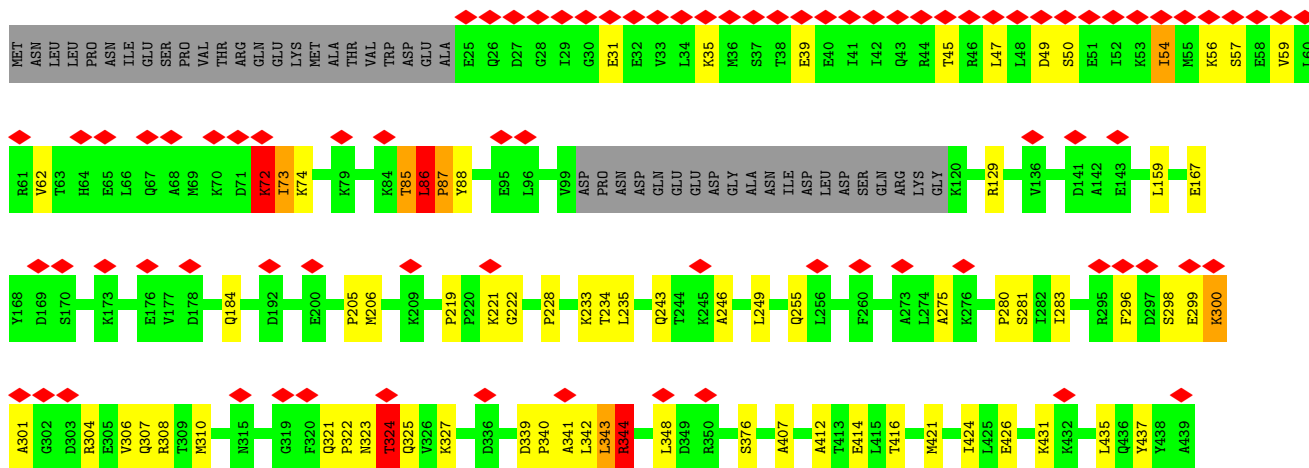
- Molecule 17: 26S proteasome regulatory subunit 10B





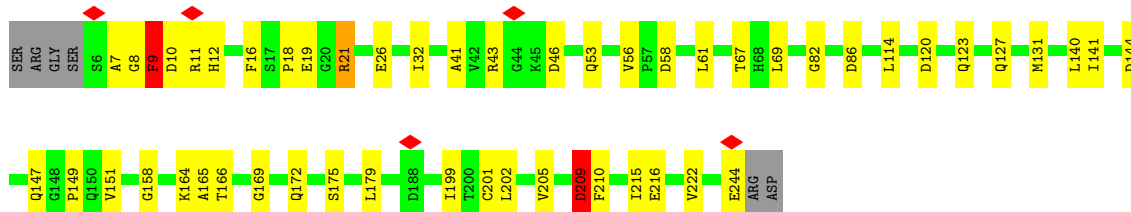
- Molecule 18: 26S protease regulatory subunit 6A

Chain F: 19% 73% 15% 10%



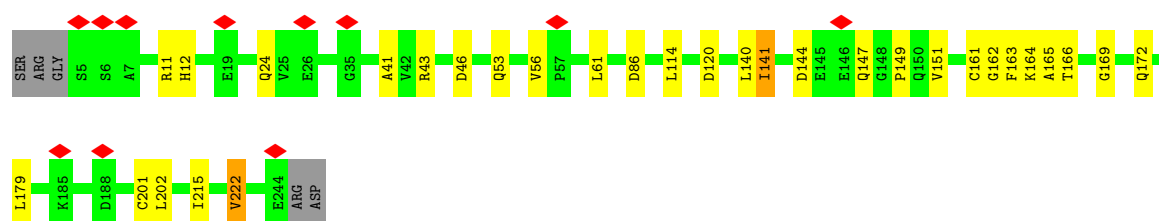
- Molecule 19: Proteasome subunit alpha type-6

Chain G: 76% 20%



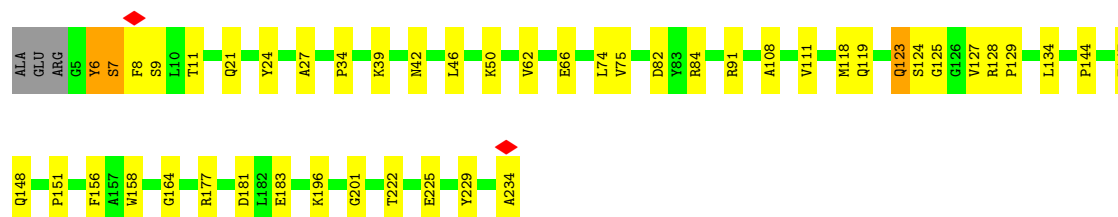
- Molecule 19: Proteasome subunit alpha type-6

Chain g:  85% 12% ..

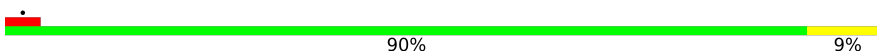


- Molecule 20: Proteasome subunit alpha type-2

Chain H:  79% 19% ..



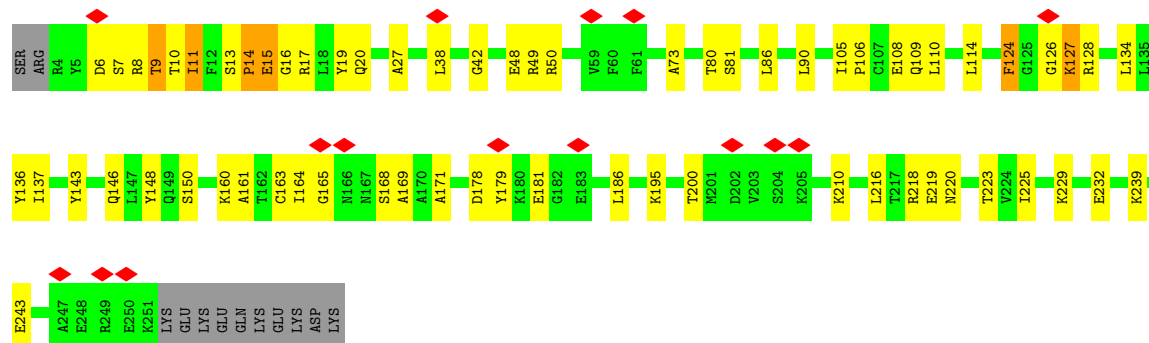
- Molecule 20: Proteasome subunit alpha type-2

Chain h:  90% 9%



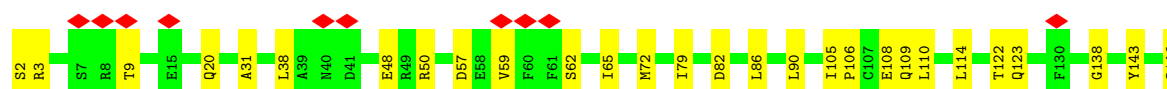
- Molecule 21: Proteasome subunit alpha type-4

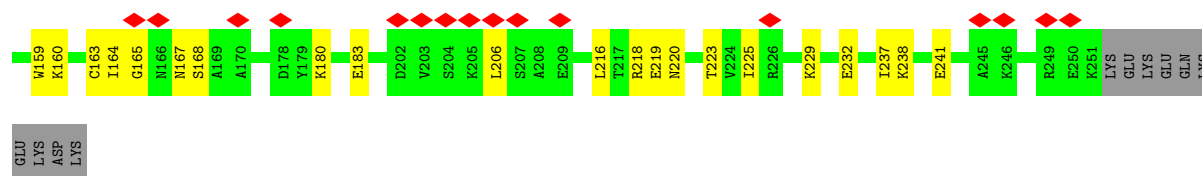
Chain I:  70% 23% 6% 5%



- Molecule 21: Proteasome subunit alpha type-4

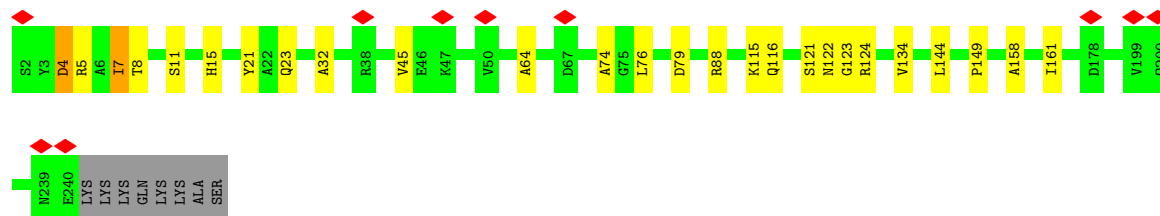
Chain i:  77% 19% 10%





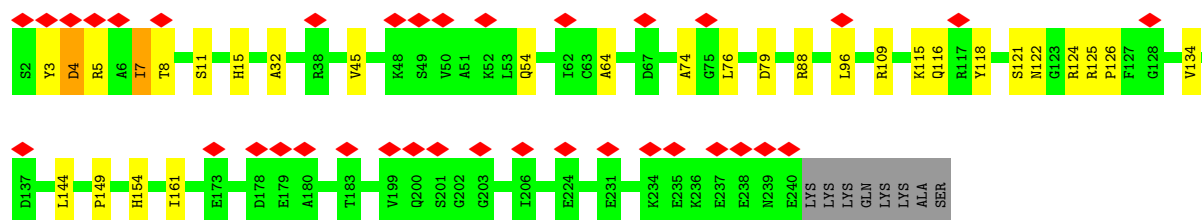
• Molecule 22: Proteasome subunit alpha type-7

Chain J: 86% 10% ..



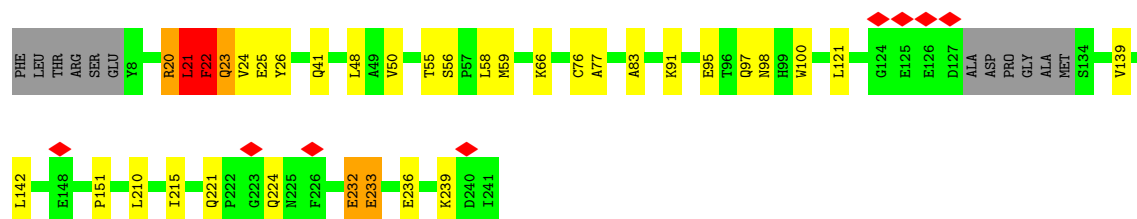
• Molecule 22: Proteasome subunit alpha type-7

Chain j: 15% 85% 11% ..



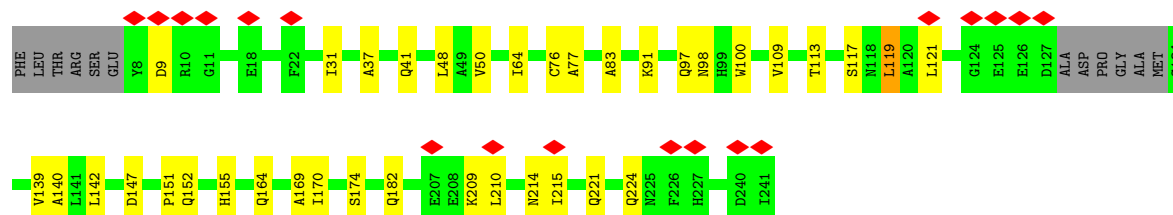
• Molecule 23: Proteasome subunit alpha type-5

Chain K: 80% 12% .. 5%

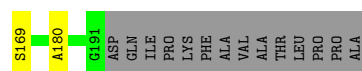


• Molecule 23: Proteasome subunit alpha type-5

Chain k: 8% 80% 15% 5%

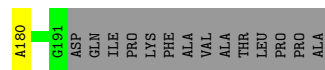
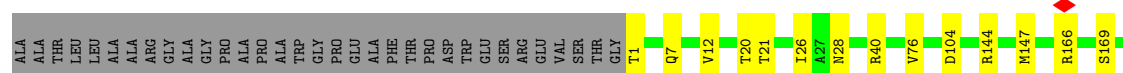


- | | | | | | | | | | | | | | | |
|-----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|
| ALA | ALA | THR | LEU | ALA | ARG | GLY | GLY | PRO | ALA | PRO | TRP | ALA | ALA | GLY | GLY | PRO | GLU | ALA | PHE | THR | PRO | ASP | TPR | GLU | SER | ANG | GLU | VAL | SER | THR | GLY | T1 | Q7 | V12 | T20 | T21 | I26 | A27 | N28 | D32 | R40 | V76 | D104 | R144 | M147 | R166 |
|-----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|



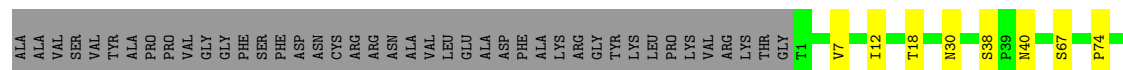
- Molecule 26: Proteasome subunit beta type-6

Chain n: 74% 6% 20%



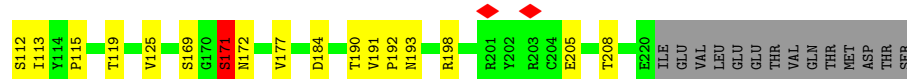
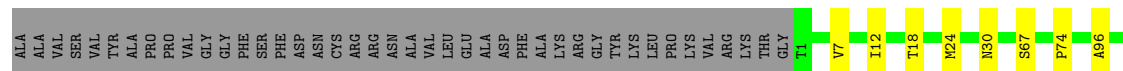
- Molecule 27: Proteasome subunit beta type-7

Chain O: 70% 9% 20%



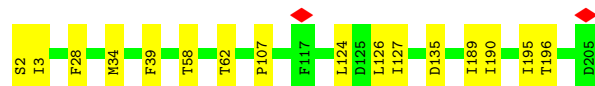
- Molecule 27: Proteasome subunit beta type-7

Chain o: 71% 9% 20%



- Molecule 28: Proteasome subunit beta type-3

Chain P: 92% 8%

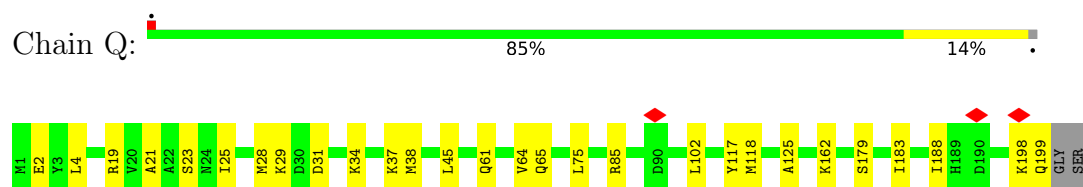


- Molecule 28: Proteasome subunit beta type-3

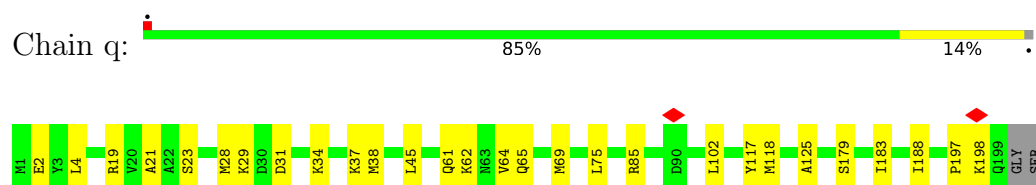
Chain p: 91% 9%



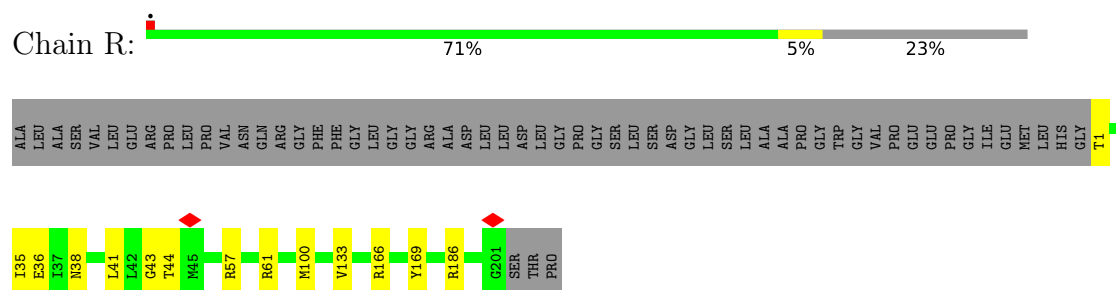
- Molecule 29: Proteasome subunit beta type-2



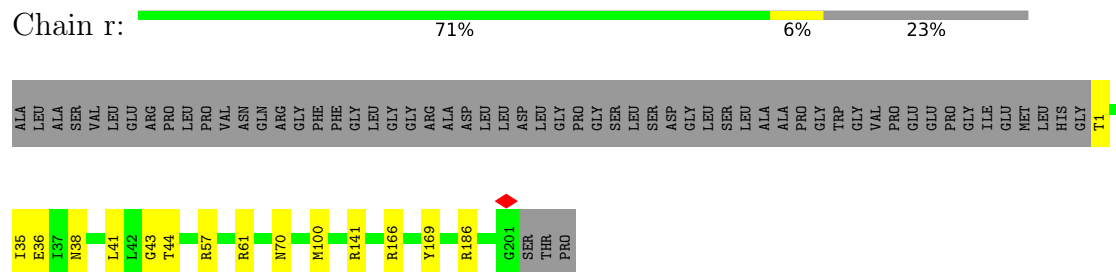
- Molecule 29: Proteasome subunit beta type-2



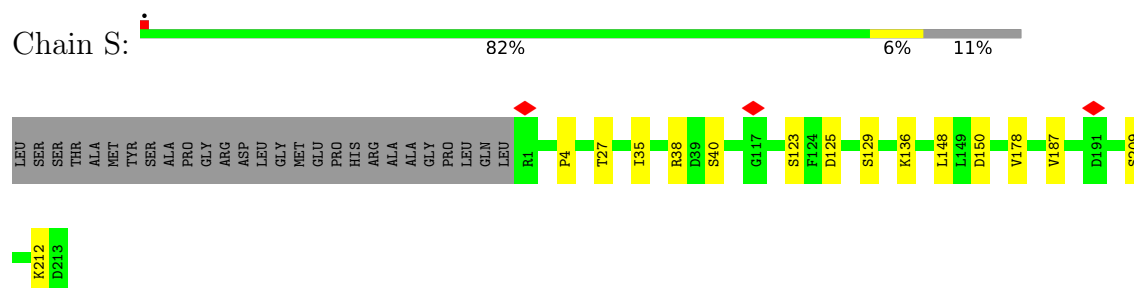
- Molecule 30: Proteasome subunit beta type-5



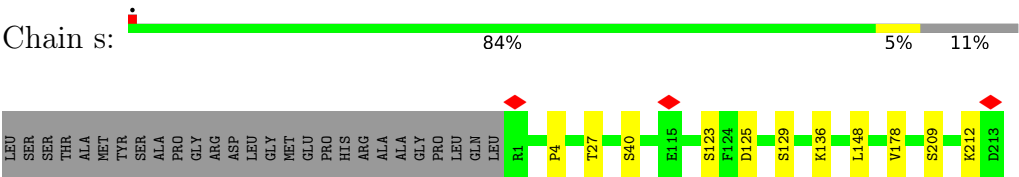
- Molecule 30: Proteasome subunit beta type-5



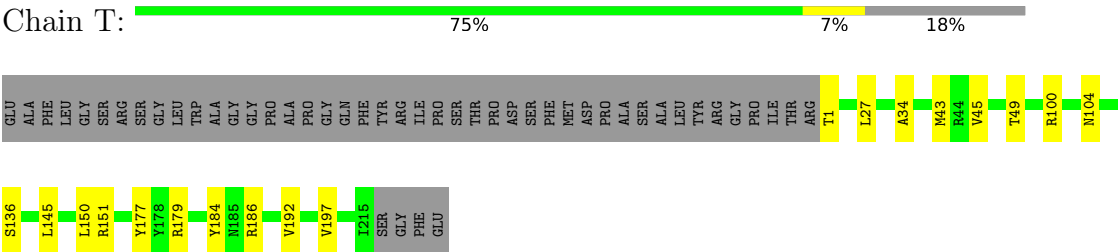
- Molecule 31: Proteasome subunit beta type-1



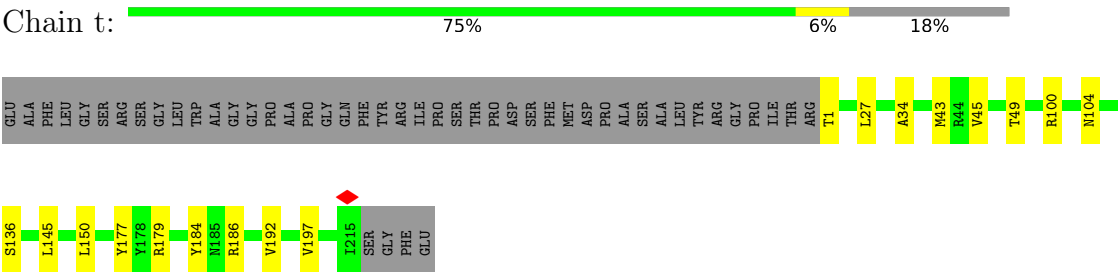
- Molecule 31: Proteasome subunit beta type-1



• Molecule 32: Proteasome subunit beta type-4



• Molecule 32: Proteasome subunit beta type-4



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	59461	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	46.6	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	0.117	Depositor
Minimum map value	-0.045	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.0337	Depositor
Map size (Å)	438.4, 438.4, 438.4	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.37, 1.37, 1.37	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	U	0.25	0/6945	0.71	17/9382 (0.2%)
2	V	0.26	0/3929	0.71	4/5309 (0.1%)
3	W	0.24	0/3751	0.65	3/5042 (0.1%)
4	X	0.21	0/3053	0.66	4/4115 (0.1%)
5	Y	0.24	0/3173	0.66	2/4273 (0.0%)
6	Z	0.24	0/2318	0.82	12/3142 (0.4%)
7	a	0.24	0/3053	0.65	2/4133 (0.0%)
8	b	0.22	0/1478	0.70	2/2001 (0.1%)
9	c	0.25	0/2302	0.80	5/3110 (0.2%)
10	d	0.25	0/2162	0.72	8/2919 (0.3%)
11	e	0.24	0/338	0.67	0/450
12	f	0.30	0/6980	0.87	15/9433 (0.2%)
13	A	0.22	0/3283	0.59	0/4433
14	B	0.27	0/3254	0.69	3/4388 (0.1%)
15	C	0.24	0/3146	0.65	2/4226 (0.0%)
16	D	0.26	0/3090	0.68	5/4168 (0.1%)
17	E	0.27	0/3145	0.98	13/4233 (0.3%)
18	F	0.22	0/3137	0.63	5/4223 (0.1%)
19	G	0.26	0/1853	0.62	3/2515 (0.1%)
19	g	0.19	0/1859	0.52	0/2523
20	H	0.22	0/1727	0.53	0/2350
20	h	0.18	0/1743	0.44	0/2372
21	I	0.24	0/1925	0.62	2/2606 (0.1%)
21	i	0.18	0/1942	0.54	0/2628
22	J	0.18	0/1728	0.52	2/2358 (0.1%)
22	j	0.19	0/1728	0.53	2/2358 (0.1%)
23	K	0.23	0/1755	0.62	4/2375 (0.2%)
23	k	0.19	0/1747	0.47	0/2364
24	L	0.27	0/1885	0.53	1/2552 (0.0%)
24	l	0.27	0/1885	0.53	1/2552 (0.0%)
25	M	0.20	0/1891	0.59	2/2552 (0.1%)
25	m	0.20	0/1891	0.59	2/2552 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
26	N	0.18	0/1454	0.49	0/1967
26	n	0.18	0/1454	0.49	0/1967
27	O	0.19	0/1670	0.48	1/2265 (0.0%)
27	o	0.19	0/1670	0.48	1/2265 (0.0%)
28	P	0.17	0/1614	0.40	0/2177
28	p	0.17	0/1614	0.40	0/2177
29	Q	0.18	0/1603	0.47	0/2174
29	q	0.18	0/1603	0.47	0/2174
30	R	0.16	0/1579	0.37	0/2134
30	r	0.16	0/1579	0.37	0/2134
31	S	0.17	0/1671	0.42	0/2253
31	s	0.17	0/1671	0.42	0/2253
32	T	0.18	0/1700	0.43	0/2305
32	t	0.18	0/1700	0.43	0/2305
All	All	0.23	0/106678	0.64	123/144187 (0.1%)

There are no bond length outliers.

The worst 5 of 123 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	E	165	ILE	CA-C-N	25.38	146.53	120.38
17	E	165	ILE	C-N-CA	25.38	146.53	120.38
12	f	755	ASP	CA-C-N	14.88	138.45	119.84
12	f	755	ASP	C-N-CA	14.88	138.45	119.84
12	f	774	GLY	N-CA-C	-14.36	97.61	114.69

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	U	6828	0	6884	154	0
2	V	3852	0	3893	120	0
3	W	3703	0	3822	125	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	X	3009	0	3112	64	0
5	Y	3115	0	3118	118	0
6	Z	2276	0	2307	104	0
7	a	2995	0	3012	51	0
8	b	1458	0	1505	43	0
9	c	2260	0	2276	52	0
10	d	2116	0	2146	33	0
11	e	334	0	294	27	0
12	f	6866	0	6866	277	0
13	A	3229	0	3260	55	0
14	B	3207	0	3278	53	0
15	C	3105	0	3218	76	0
16	D	3040	0	3075	64	0
17	E	3097	0	3174	145	0
18	F	3098	0	3186	117	0
19	G	1820	0	1791	67	0
19	g	1826	0	1796	29	0
20	H	1692	0	1586	53	0
20	h	1708	0	1594	13	0
21	I	1895	0	1833	56	0
21	i	1912	0	1851	28	0
22	J	1704	0	1517	18	0
22	j	1704	0	1517	24	0
23	K	1729	0	1680	41	0
23	k	1722	0	1673	24	0
24	L	1850	0	1822	11	0
24	l	1850	0	1822	16	0
25	M	1856	0	1814	38	0
25	m	1856	0	1814	36	0
26	N	1430	0	1398	8	0
26	n	1430	0	1398	8	0
27	O	1643	0	1644	21	0
27	o	1643	0	1644	20	0
28	P	1585	0	1598	10	0
28	p	1585	0	1598	12	0
29	Q	1570	0	1547	18	0
29	q	1570	0	1547	17	0
30	R	1548	0	1499	10	0
30	r	1548	0	1499	11	0
31	S	1641	0	1618	11	0
31	s	1641	0	1618	6	0
32	T	1667	0	1628	14	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
32	t	1667	0	1628	11	0
33	c	1	0	0	0	0
34	A	31	0	12	1	0
34	B	31	0	12	2	0
34	C	31	0	12	2	0
34	F	31	0	12	1	0
35	A	1	0	0	0	0
35	B	1	0	0	0	0
35	C	1	0	0	0	0
35	D	1	0	0	0	0
35	F	1	0	0	0	0
36	D	27	0	12	0	0
All	All	105037	0	104460	1986	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 1986 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)
17:E:164:ILE:CG1	17:E:166:PRO:HG2	1.39	1.52
8:b:22:LEU:HB3	8:b:23:PRO:CD	1.32	1.46
13:A:123:VAL:HA	18:F:86:LEU:CD1	1.42	1.46
12:f:281:ILE:HD11	12:f:285:CYS:SG	1.56	1.43
4:X:143:TYR:OH	5:Y:252:SER:CA	1.64	1.43

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	U	868/953 (91%)	782 (90%)	78 (9%)	8 (1%)	14	49
2	V	478/533 (90%)	413 (86%)	63 (13%)	2 (0%)	30	65
3	W	454/456 (100%)	409 (90%)	41 (9%)	4 (1%)	14	49
4	X	378/422 (90%)	355 (94%)	21 (6%)	2 (0%)	24	61
5	Y	376/389 (97%)	332 (88%)	41 (11%)	3 (1%)	16	52
6	Z	284/324 (88%)	243 (86%)	37 (13%)	4 (1%)	9	39
7	a	371/376 (99%)	334 (90%)	34 (9%)	3 (1%)	16	52
8	b	189/377 (50%)	164 (87%)	24 (13%)	1 (0%)	24	61
9	c	285/309 (92%)	246 (86%)	32 (11%)	7 (2%)	4	29
10	d	255/349 (73%)	214 (84%)	38 (15%)	3 (1%)	10	42
11	e	36/70 (51%)	29 (81%)	7 (19%)	0	100	100
12	f	887/908 (98%)	707 (80%)	164 (18%)	16 (2%)	6	35
13	A	411/433 (95%)	364 (89%)	46 (11%)	1 (0%)	43	76
14	B	409/440 (93%)	349 (85%)	57 (14%)	3 (1%)	18	55
15	C	394/398 (99%)	353 (90%)	36 (9%)	5 (1%)	9	41
16	D	378/418 (90%)	321 (85%)	53 (14%)	4 (1%)	11	44
17	E	387/403 (96%)	323 (84%)	54 (14%)	10 (3%)	4	28
18	F	391/439 (89%)	336 (86%)	47 (12%)	8 (2%)	6	33
19	G	237/245 (97%)	209 (88%)	25 (10%)	3 (1%)	9	41
19	g	238/245 (97%)	222 (93%)	16 (7%)	0	100	100
20	H	228/233 (98%)	215 (94%)	10 (4%)	3 (1%)	9	41
20	h	230/233 (99%)	222 (96%)	8 (4%)	0	100	100
21	I	246/260 (95%)	217 (88%)	26 (11%)	3 (1%)	10	42
21	i	248/260 (95%)	223 (90%)	25 (10%)	0	100	100
22	J	237/247 (96%)	216 (91%)	20 (8%)	1 (0%)	30	65
22	j	237/247 (96%)	216 (91%)	20 (8%)	1 (0%)	30	65
23	K	224/240 (93%)	208 (93%)	14 (6%)	2 (1%)	14	49
23	k	224/240 (93%)	210 (94%)	14 (6%)	0	100	100
24	L	236/268 (88%)	226 (96%)	10 (4%)	0	100	100
24	l	236/268 (88%)	226 (96%)	10 (4%)	0	100	100
25	M	238/254 (94%)	220 (92%)	18 (8%)	0	100	100
25	m	238/254 (94%)	219 (92%)	19 (8%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
26	N	189/238 (79%)	180 (95%)	9 (5%)	0	100	100
26	n	189/238 (79%)	180 (95%)	9 (5%)	0	100	100
27	O	218/276 (79%)	207 (95%)	10 (5%)	1 (0%)	24	61
27	o	218/276 (79%)	207 (95%)	10 (5%)	1 (0%)	24	61
28	P	202/204 (99%)	194 (96%)	8 (4%)	0	100	100
28	p	202/204 (99%)	194 (96%)	8 (4%)	0	100	100
29	Q	197/201 (98%)	186 (94%)	11 (6%)	0	100	100
29	q	197/201 (98%)	186 (94%)	11 (6%)	0	100	100
30	R	199/262 (76%)	193 (97%)	6 (3%)	0	100	100
30	r	199/262 (76%)	193 (97%)	6 (3%)	0	100	100
31	S	211/240 (88%)	205 (97%)	6 (3%)	0	100	100
31	s	211/240 (88%)	205 (97%)	6 (3%)	0	100	100
32	T	213/263 (81%)	202 (95%)	11 (5%)	0	100	100
32	t	213/263 (81%)	202 (95%)	11 (5%)	0	100	100
All	All	13386/14859 (90%)	12057 (90%)	1230 (9%)	99 (1%)	20	55

5 of 99 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	U	48	LEU
3	W	91	SER
3	W	176	SER
3	W	179	LYS
4	X	203	PRO

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	U	748/816 (92%)	736 (98%)	12 (2%)	55	69
2	V	414/459 (90%)	410 (99%)	4 (1%)	68	76

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	W	416/416 (100%)	408 (98%)	8 (2%)	50	67
4	X	327/362 (90%)	321 (98%)	6 (2%)	51	67
5	Y	334/344 (97%)	326 (98%)	8 (2%)	43	63
6	Z	256/295 (87%)	247 (96%)	9 (4%)	32	54
7	a	333/336 (99%)	330 (99%)	3 (1%)	70	76
8	b	167/312 (54%)	166 (99%)	1 (1%)	78	80
9	c	252/267 (94%)	246 (98%)	6 (2%)	43	63
10	d	231/293 (79%)	225 (97%)	6 (3%)	40	61
11	e	38/63 (60%)	37 (97%)	1 (3%)	40	61
12	f	745/763 (98%)	723 (97%)	22 (3%)	36	57
13	A	348/372 (94%)	347 (100%)	1 (0%)	86	85
14	B	357/385 (93%)	348 (98%)	9 (2%)	42	62
15	C	340/346 (98%)	333 (98%)	7 (2%)	47	65
16	D	333/366 (91%)	325 (98%)	8 (2%)	43	63
17	E	341/353 (97%)	333 (98%)	8 (2%)	44	64
18	F	340/379 (90%)	331 (97%)	9 (3%)	40	61
19	G	192/209 (92%)	189 (98%)	3 (2%)	55	69
19	g	193/209 (92%)	191 (99%)	2 (1%)	68	76
20	H	163/190 (86%)	161 (99%)	2 (1%)	63	73
20	h	164/190 (86%)	164 (100%)	0	100	100
21	I	191/220 (87%)	187 (98%)	4 (2%)	47	65
21	i	193/220 (88%)	193 (100%)	0	100	100
22	J	152/210 (72%)	151 (99%)	1 (1%)	76	79
22	j	152/210 (72%)	151 (99%)	1 (1%)	76	79
23	K	187/202 (93%)	181 (97%)	6 (3%)	34	56
23	k	186/202 (92%)	183 (98%)	3 (2%)	55	69
24	L	198/229 (86%)	196 (99%)	2 (1%)	68	76
24	l	198/229 (86%)	197 (100%)	1 (0%)	81	82
25	M	192/211 (91%)	190 (99%)	2 (1%)	68	76
25	m	192/211 (91%)	190 (99%)	2 (1%)	68	76
26	N	148/180 (82%)	148 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
26	n	148/180 (82%)	148 (100%)	0	100	100
27	O	177/227 (78%)	177 (100%)	0	100	100
27	o	177/227 (78%)	177 (100%)	0	100	100
28	P	172/173 (99%)	172 (100%)	0	100	100
28	p	172/173 (99%)	172 (100%)	0	100	100
29	Q	164/171 (96%)	164 (100%)	0	100	100
29	q	164/171 (96%)	164 (100%)	0	100	100
30	R	153/201 (76%)	153 (100%)	0	100	100
30	r	153/201 (76%)	153 (100%)	0	100	100
31	S	174/198 (88%)	174 (100%)	0	100	100
31	s	174/198 (88%)	174 (100%)	0	100	100
32	T	175/214 (82%)	175 (100%)	0	100	100
32	t	175/214 (82%)	175 (100%)	0	100	100
All	All	11199/12597 (89%)	11042 (99%)	157 (1%)	57	71

5 of 157 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
17	E	196	LEU
23	K	233	GLU
18	F	54	ILE
19	G	209	ASP
22	j	7	ILE

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

Mol	Chain	Res	Type
16	D	127	ASN
23	K	155	HIS
16	D	414	HIS
20	H	71	HIS
27	O	172	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 11 ligands modelled in this entry, 6 are monoatomic - leaving 5 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	$\# Z > 2$
34	ATP	C	501	35	32,33,33	1.31	4 (12%)	48,52,52	1.81	7 (14%)
34	ATP	B	501	35	32,33,33	1.32	4 (12%)	48,52,52	1.90	9 (18%)
34	ATP	F	501	35	32,33,33	1.35	5 (15%)	48,52,52	1.81	8 (16%)
36	ADP	D	501	35	28,29,29	1.37	4 (14%)	43,45,45	1.88	9 (20%)
34	ATP	A	501	35	32,33,33	1.32	4 (12%)	48,52,52	1.79	10 (20%)

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
34	ATP	C	501	35	-	5/22/38/38	0/3/3/3
34	ATP	B	501	35	-	3/22/38/38	0/3/3/3
34	ATP	F	501	35	-	4/22/38/38	0/3/3/3
36	ADP	D	501	35	-	1/16/32/32	0/3/3/3
34	ATP	A	501	35	-	3/22/38/38	0/3/3/3

The worst 5 of 21 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
34	B	501	ATP	C5-C4	4.65	1.47	1.39
34	C	501	ATP	C5-C4	4.64	1.47	1.39
34	F	501	ATP	C5-C4	4.60	1.47	1.39
34	A	501	ATP	C5-C4	4.53	1.47	1.39
36	D	501	ADP	C5-C4	4.52	1.47	1.39

The worst 5 of 43 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	B	501	ATP	C5-C4-N3	-6.55	117.70	126.72
34	C	501	ATP	C5-C4-N3	-6.34	117.99	126.72
34	A	501	ATP	C5-C4-N3	-5.92	118.57	126.72
34	F	501	ATP	C5-C4-N3	-5.76	118.78	126.72
36	D	501	ADP	C5-C4-N3	-5.74	118.81	126.72

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
34	A	501	ATP	C5'-O5'-PA-O1A
34	A	501	ATP	C5'-O5'-PA-O2A
34	A	501	ATP	C5'-O5'-PA-O3A
34	B	501	ATP	C5'-O5'-PA-O2A
34	B	501	ATP	C5'-O5'-PA-O3A

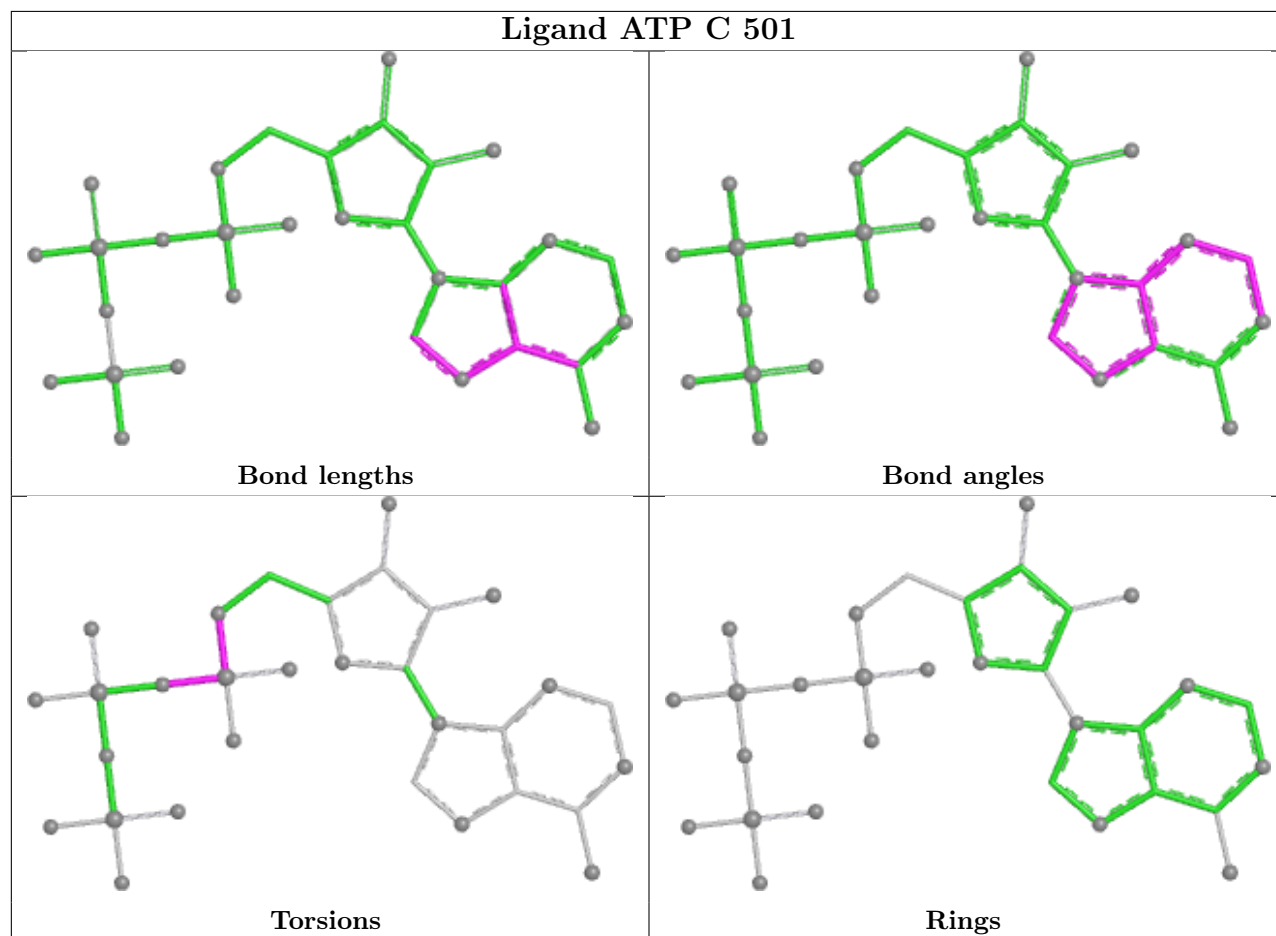
There are no ring outliers.

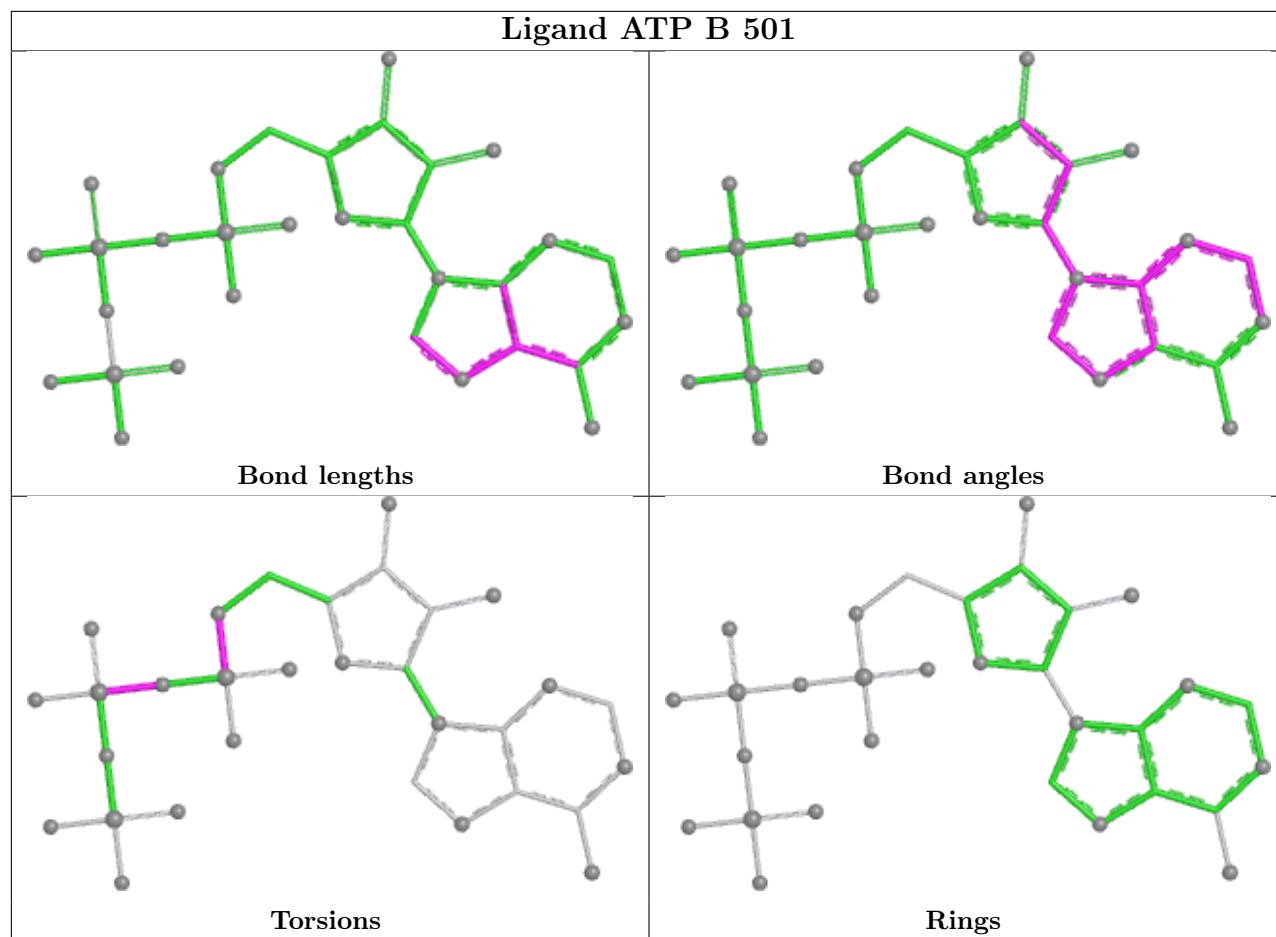
4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
34	C	501	ATP	2	0
34	B	501	ATP	2	0
34	F	501	ATP	1	0
34	A	501	ATP	1	0

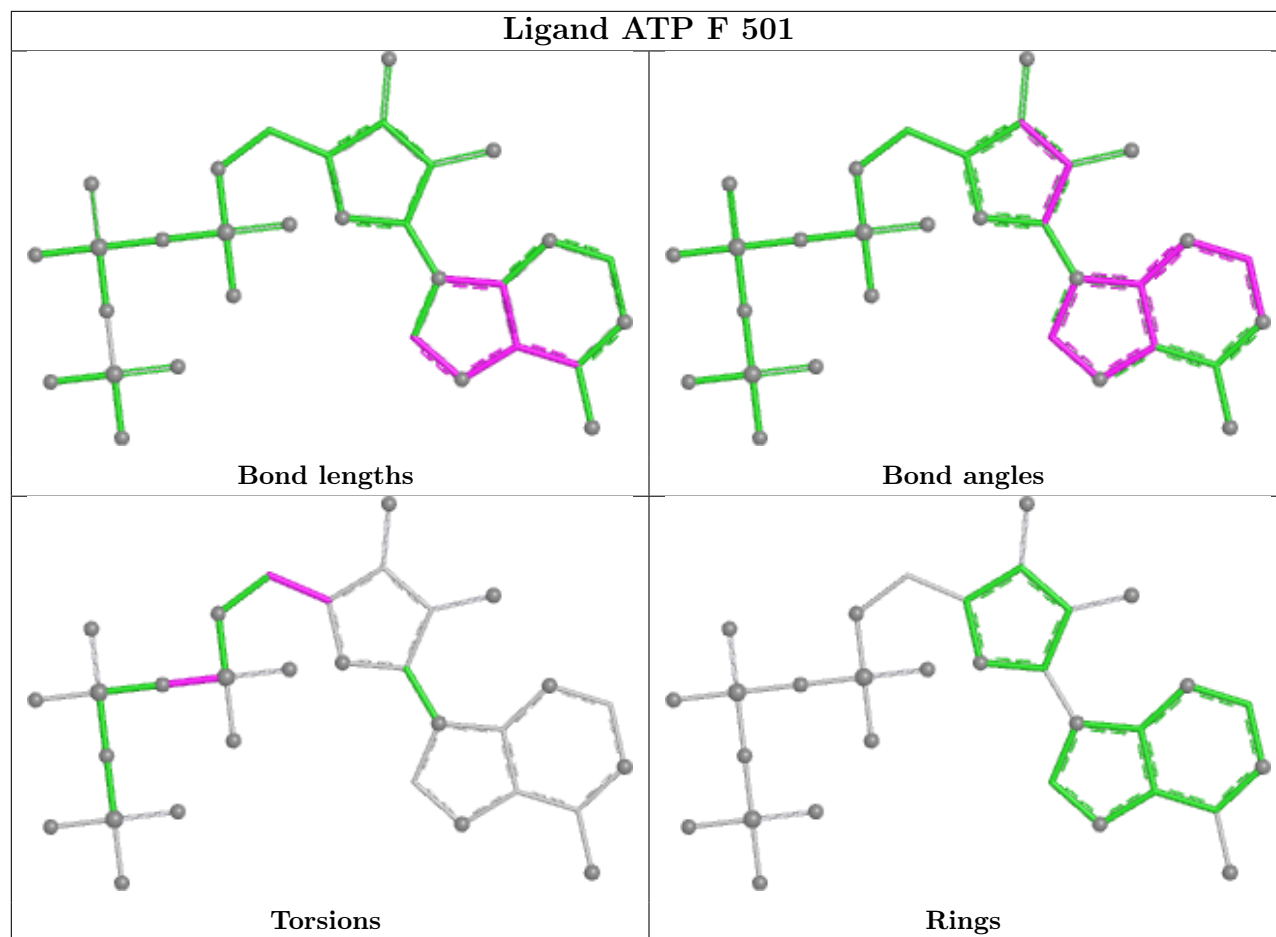
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

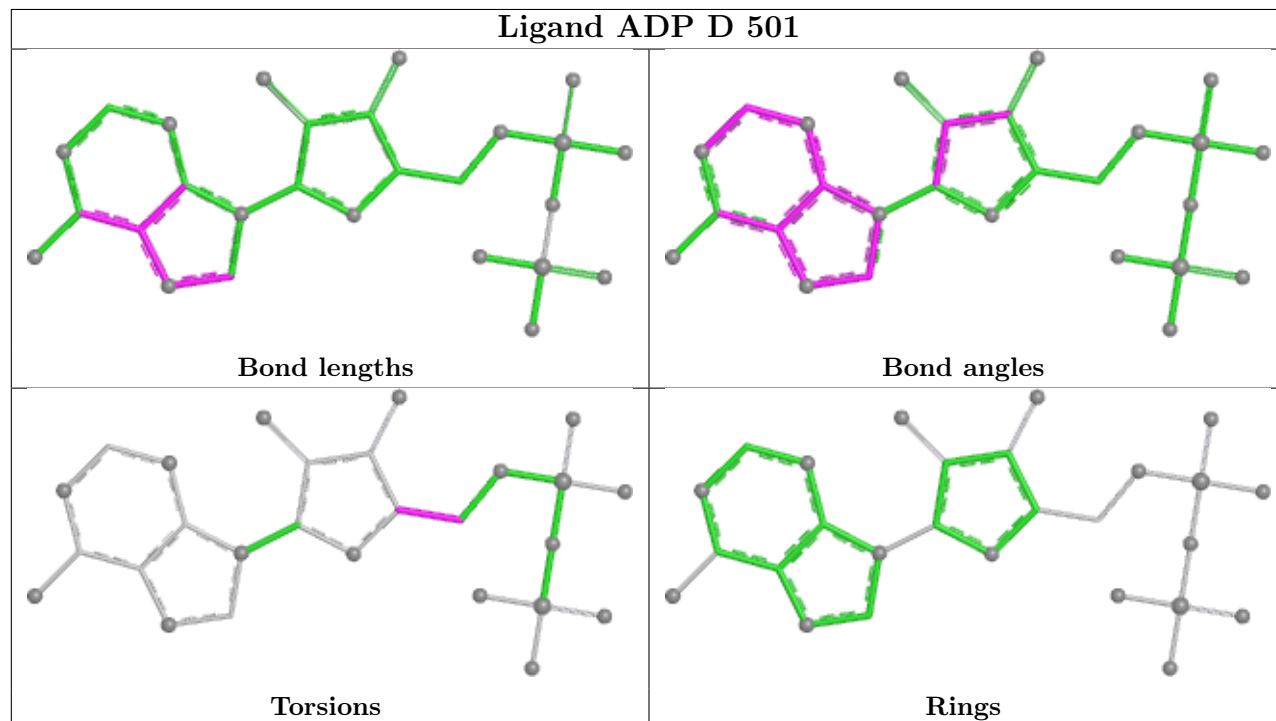


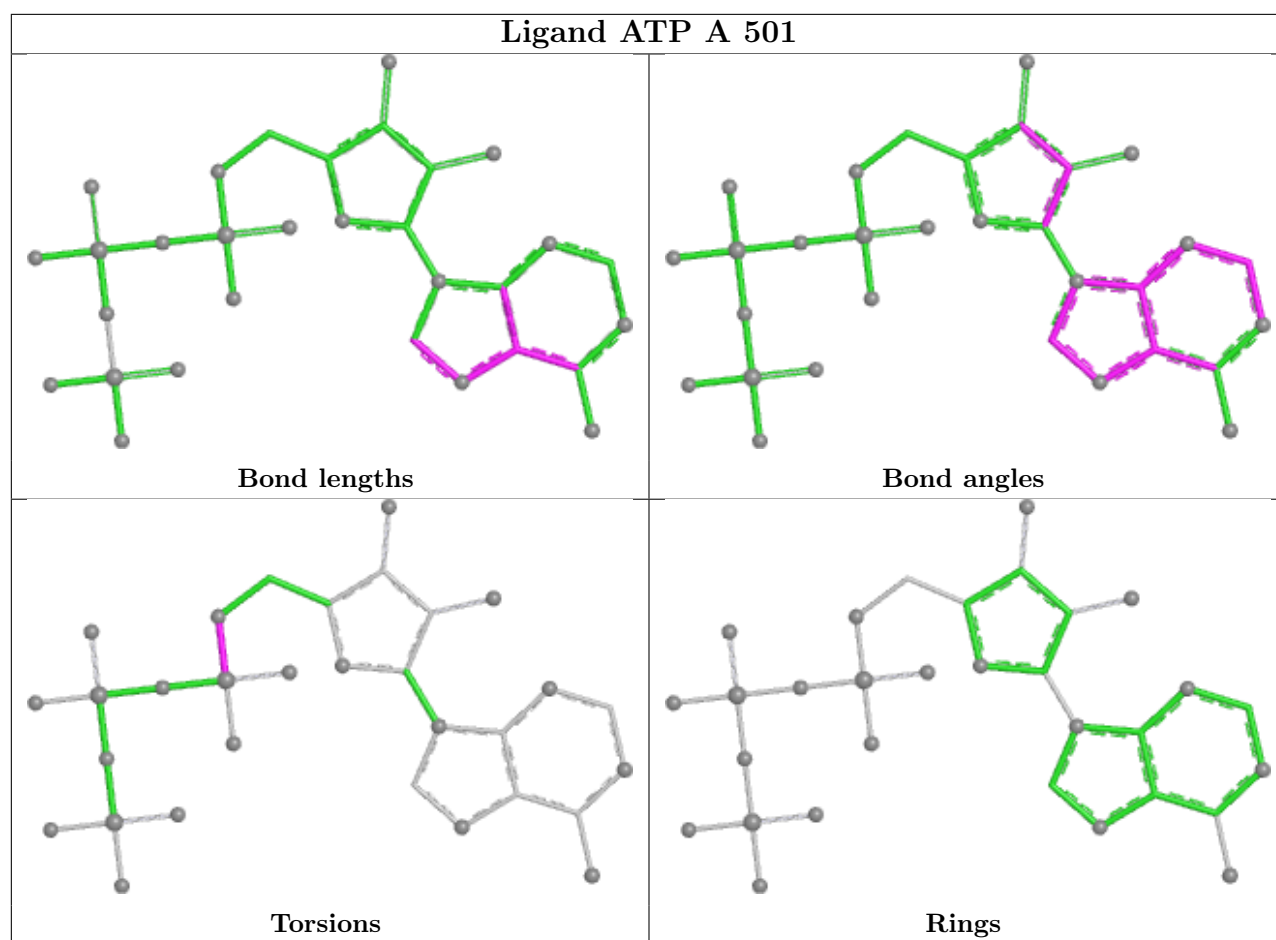


Ligand ATP F 501



Ligand ADP D 501





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.

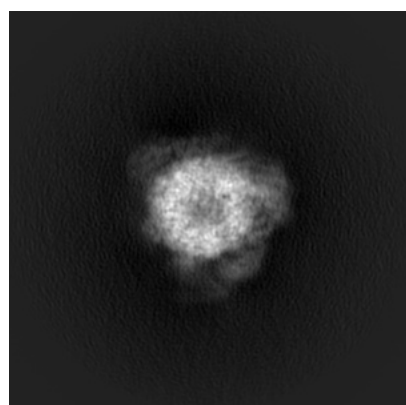
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-14211. These allow visual inspection of the internal detail of the map and identification of artifacts.

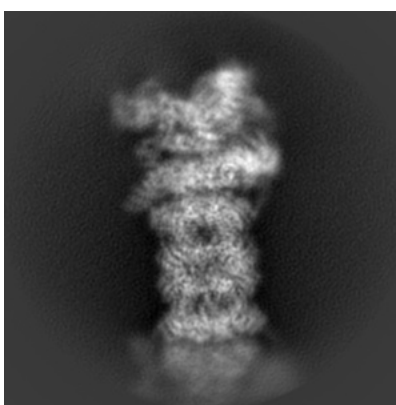
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

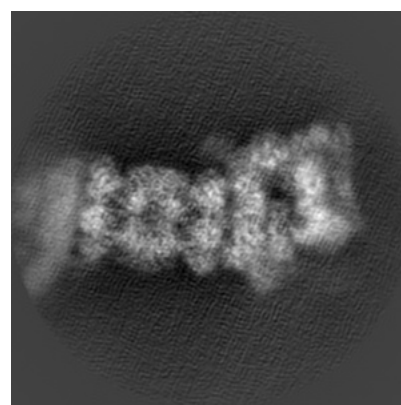
6.1.1 Primary map



X



Y

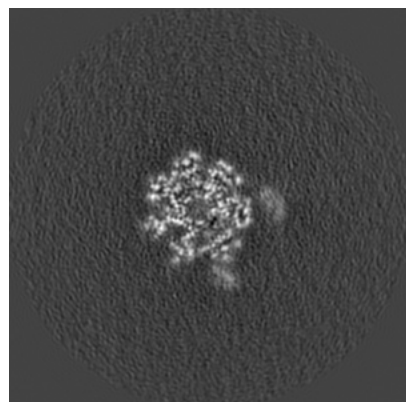


Z

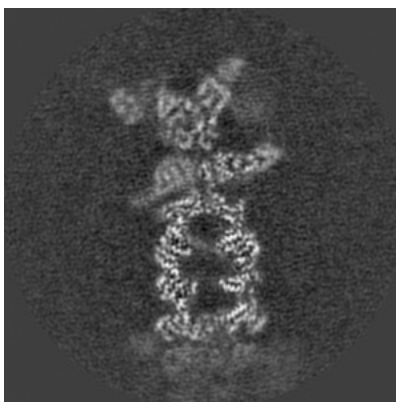
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

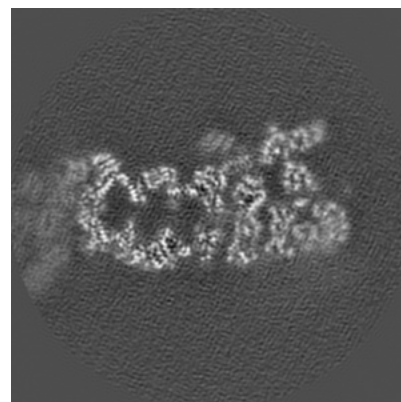
6.2.1 Primary map



X Index: 160



Y Index: 160

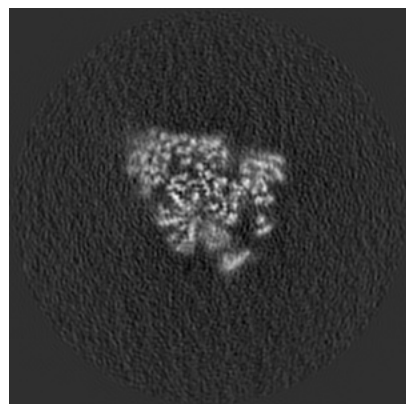


Z Index: 160

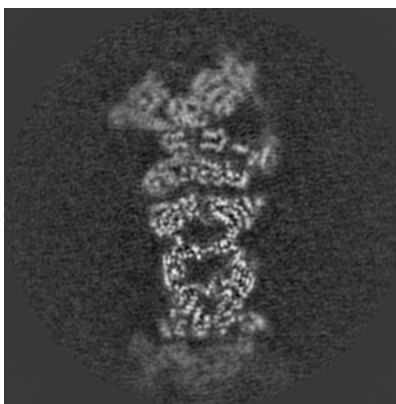
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

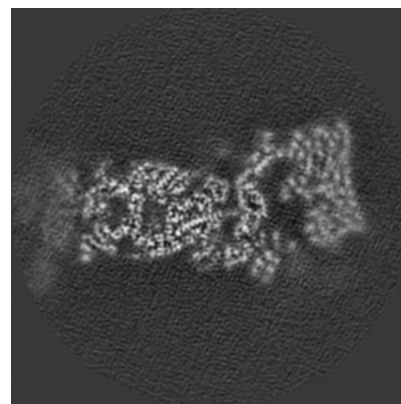
6.3.1 Primary map



X Index: 196



Y Index: 143

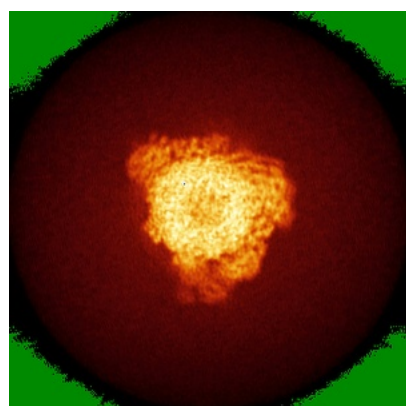


Z Index: 180

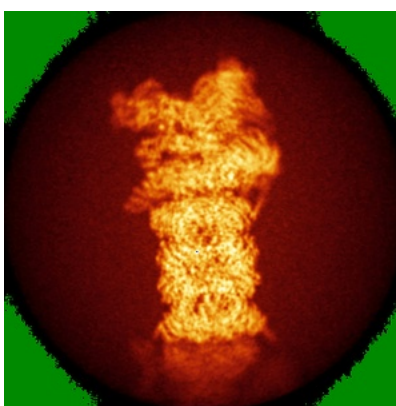
The images above show the largest variance slices of the map in three orthogonal directions.

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

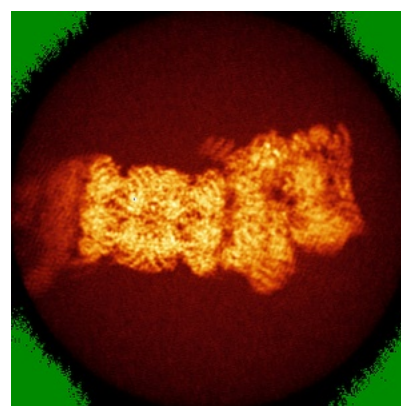
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views

This section was not generated.

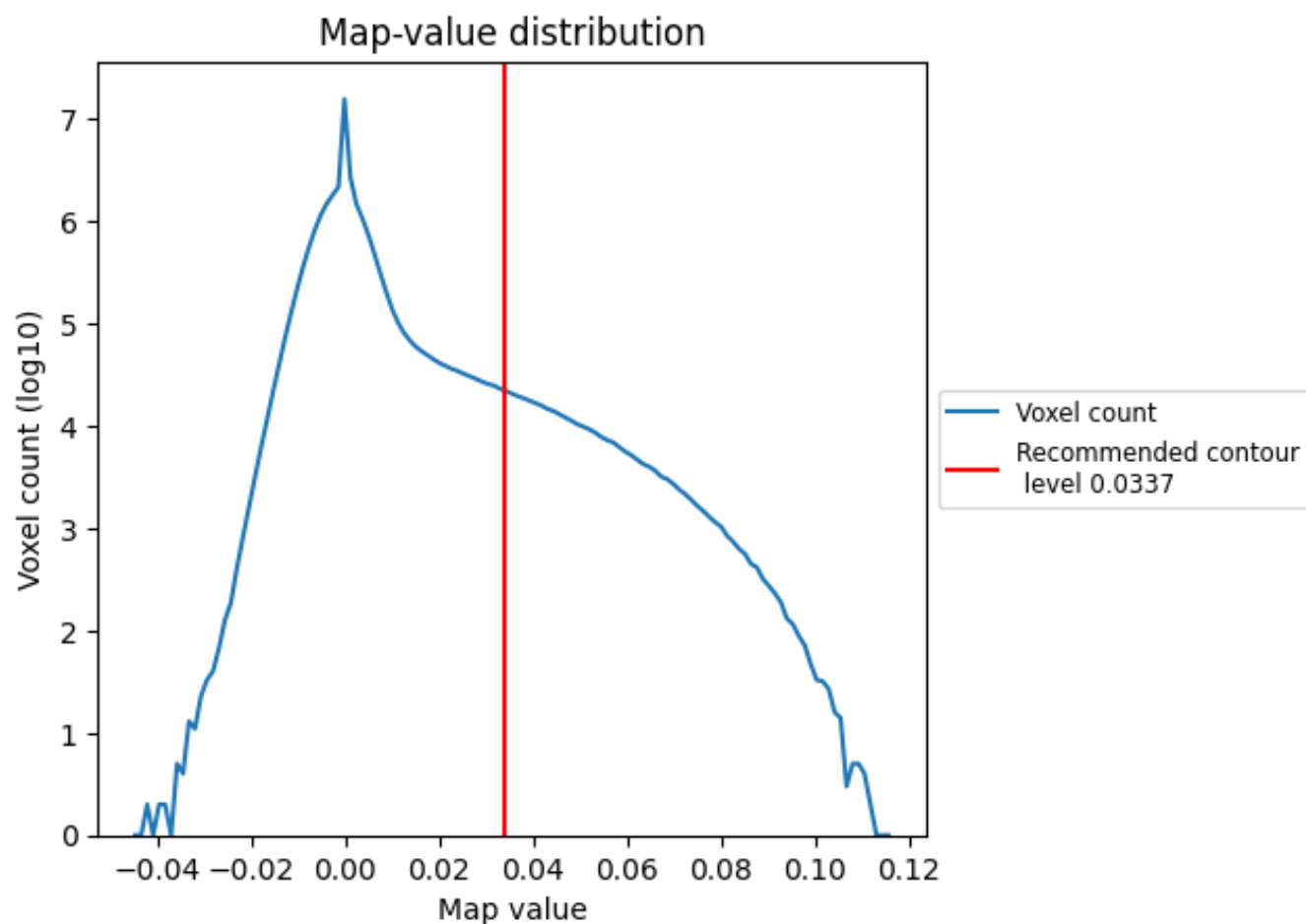
6.6 Mask visualisation

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

7 Map analysis [i](#)

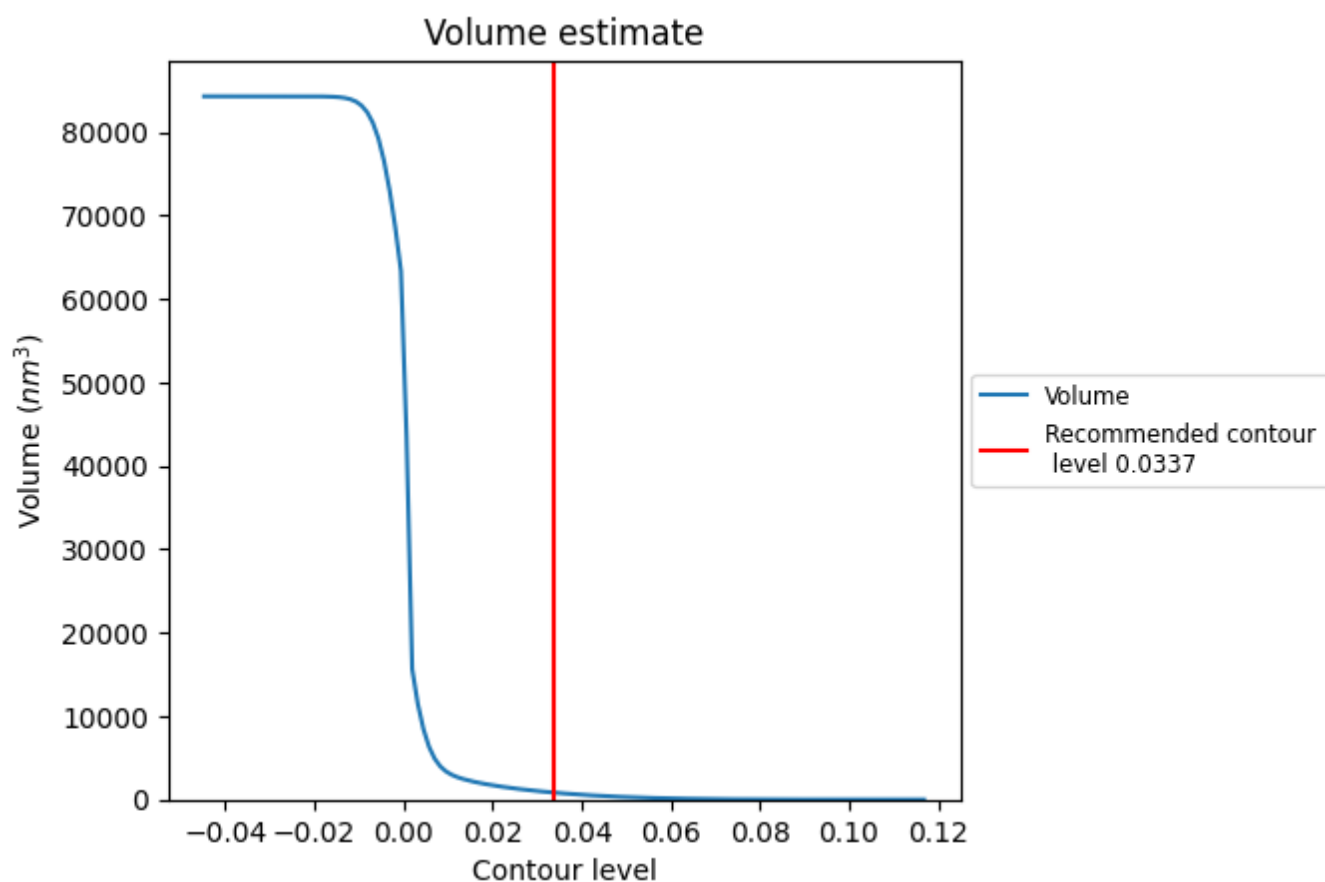
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

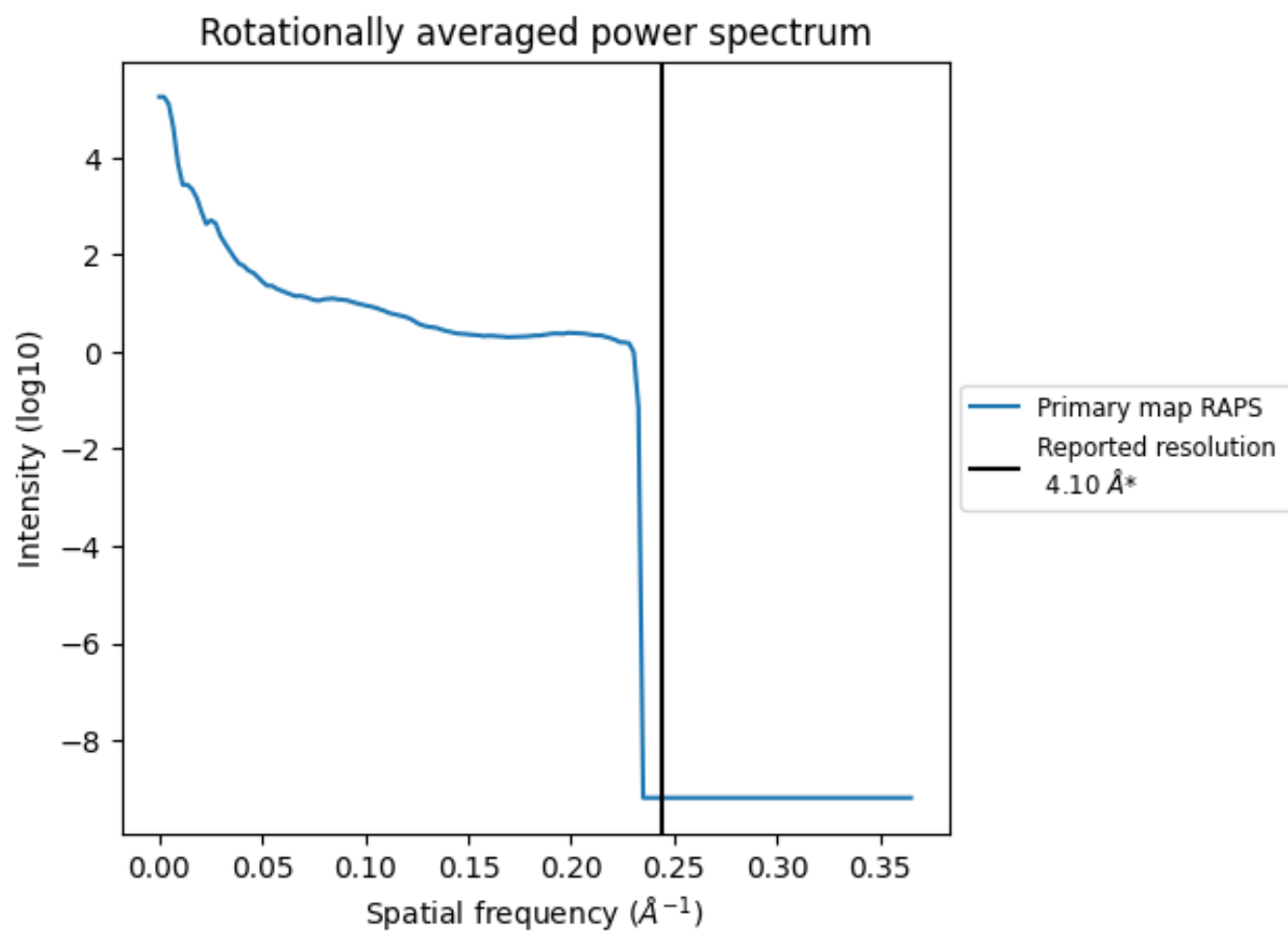
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 837 nm³; this corresponds to an approximate mass of 756 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 0.244 Å⁻¹

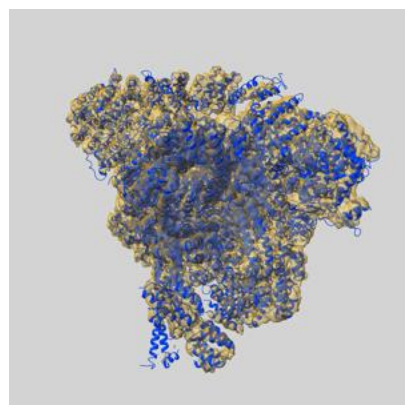
8 Fourier-Shell correlation ⓘ

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

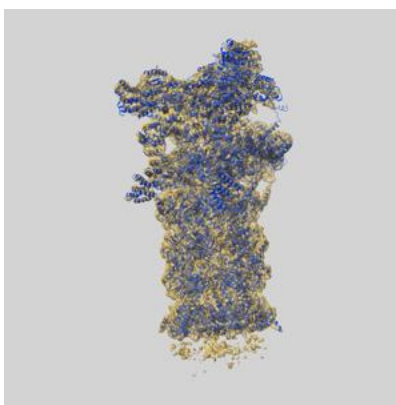
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-14211 and PDB model 7QYB. Per-residue inclusion information can be found in [section 3](#) on [page 12](#).

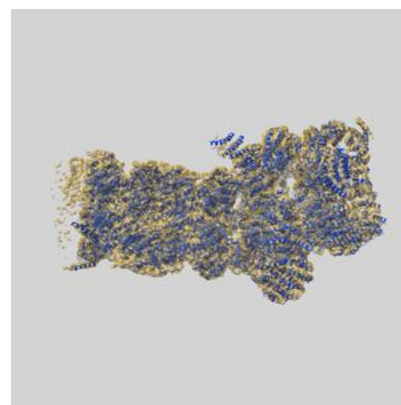
9.1 Map-model overlay [i](#)



X



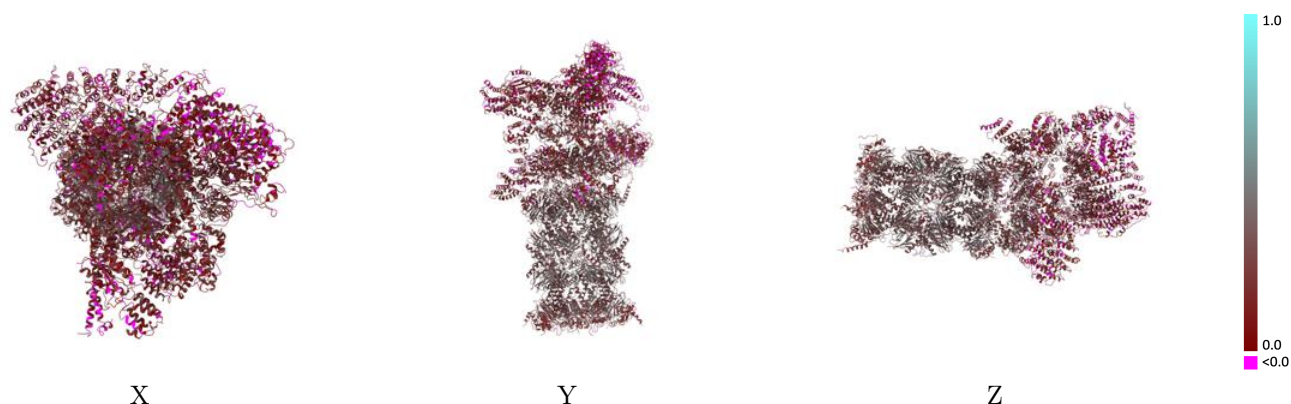
Y



Z

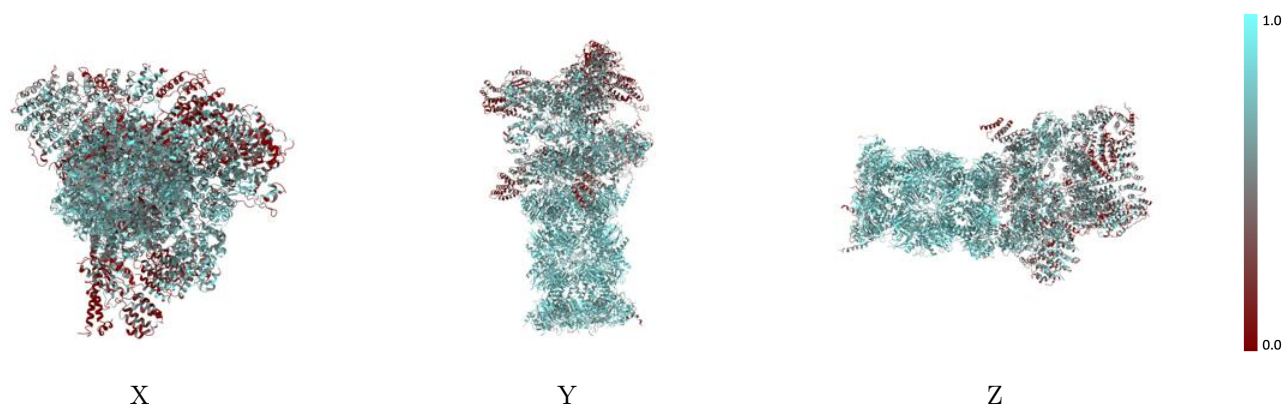
The images above show the 3D surface view of the map at the recommended contour level 0.0337 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



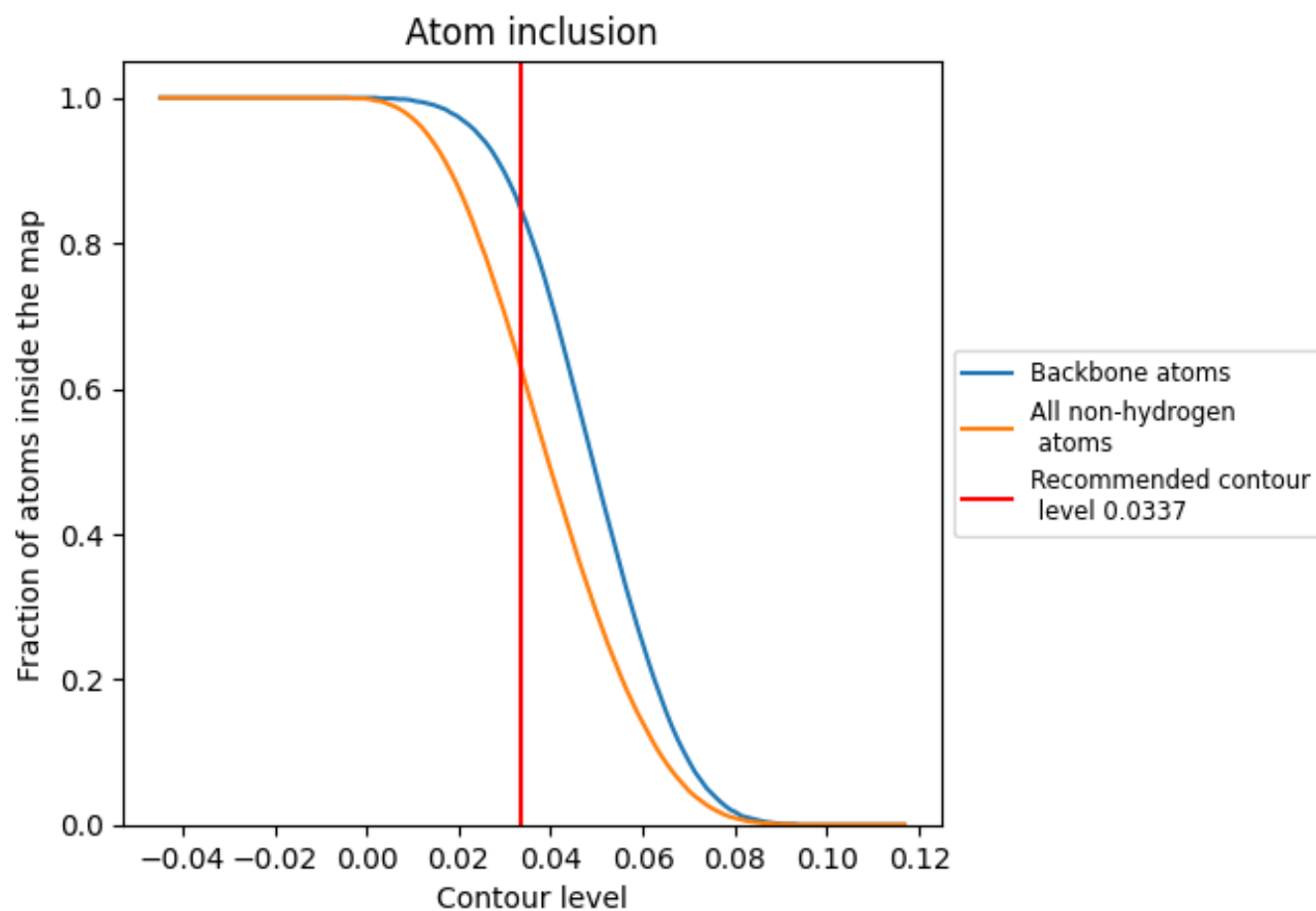
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0337).

9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 62% 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 (0.0337) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6250	 0.2520
A	 0.6340	 0.2650
B	 0.6170	 0.2700
C	 0.6530	 0.2820
D	 0.6330	 0.2600
E	 0.4120	 0.1550
F	 0.5660	 0.2320
G	 0.7440	 0.3140
H	 0.7590	 0.3370
I	 0.7170	 0.3180
J	 0.7540	 0.3340
K	 0.7340	 0.3320
L	 0.7760	 0.3470
M	 0.7380	 0.3150
N	 0.7930	 0.3520
O	 0.7890	 0.3540
P	 0.7890	 0.3590
Q	 0.7860	 0.3480
R	 0.8200	 0.3590
S	 0.7960	 0.3680
T	 0.7970	 0.3560
U	 0.4710	 0.1370
V	 0.3860	 0.1150
W	 0.4450	 0.1670
X	 0.5250	 0.2160
Y	 0.6590	 0.1970
Z	 0.5510	 0.2010
a	 0.4530	 0.1600
b	 0.3520	 0.1510
c	 0.5780	 0.2070
d	 0.3770	 0.1210
e	 0.5240	 0.1550
f	 0.5070	 0.1490
g	 0.7070	 0.3090
h	 0.7030	 0.3050



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
i	 0.6580	 0.2750
j	 0.6690	 0.2670
k	 0.6850	 0.2950
l	 0.7380	 0.3180
m	 0.7240	 0.2940
n	 0.7980	 0.3480
o	 0.7910	 0.3540
p	 0.7750	 0.3510
q	 0.7840	 0.3370
r	 0.8140	 0.3570
s	 0.7920	 0.3600
t	 0.7970	 0.3560