



wwPDB EM Validation Summary Report ⓘ

Mar 28, 2026 – 08:11 PM UTC

PDB ID : 7W3I / pdb_00007w3i
EMDB ID : EMD-32281
Title : Structure of USP14-bound human 26S proteasome in substrate-inhibited state
SB_USP14
Authors : Zhang, S.; Zou, S.; Yin, D.; Wu, Z.; Mao, Y.
Deposited on : 2021-11-25
Resolution : 3.50 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

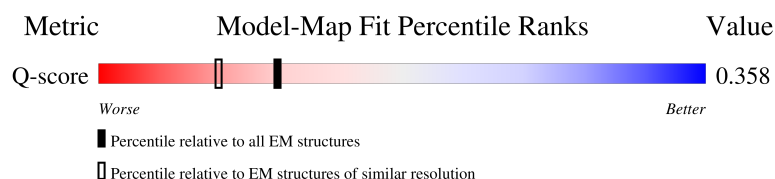
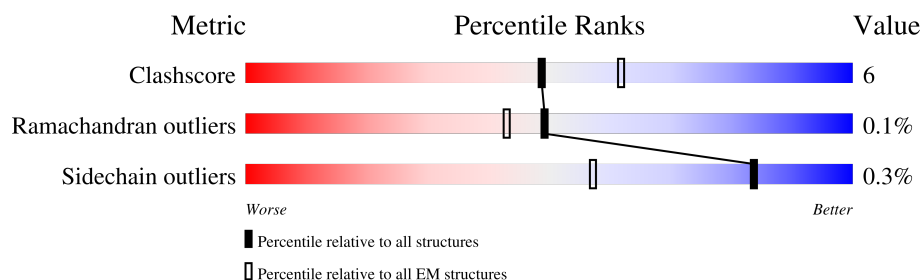
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.50 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.



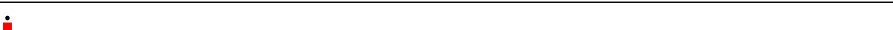
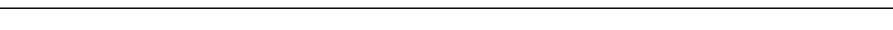
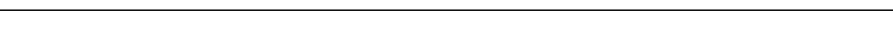
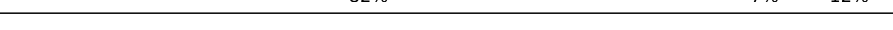
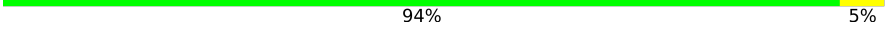
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13950 (3.00 - 4.00)

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	x	494	
2	A	433	
3	B	440	
4	C	398	

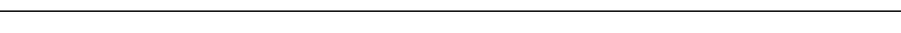
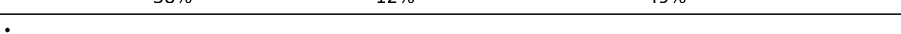
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Mol	Chain	Length	Quality of chain
5	D	418	 70%21%9%
6	E	403	 79%17%.
7	F	439	 76%15%9%
8	G	246	 91%9%.
8	g	246	 91%8%.
9	H	234	 91%9%.
9	h	234	 94%6%.
10	I	261	 90%6%.
10	i	261	 90%6%.
11	J	248	 88%8%.
11	j	248	 87%9%.
12	K	241	 87%10%.
12	k	241	 93%. .
13	L	269	 82%7%12%
13	l	269	 82%7%12%
14	M	255	 87%7%6%
14	m	255	 86%8%6%
15	N	239	 83%.15%
15	n	239	 79%5%15%
16	O	277	 73%7%21%
16	o	277	 71%8%21%
17	P	205	 94%5%
17	p	205	 87%13%
18	Q	201	 87%12%.
18	q	201	 89%10%.

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Mol	Chain	Length	Quality of chain
19	R	263	
19	r	263	
20	S	241	
20	s	241	
21	T	264	
21	t	264	
22	y	76	
23	U	953	
24	V	534	
25	W	456	
26	X	422	
27	Y	389	
28	Z	324	
29	a	376	
30	b	377	
31	c	310	
32	d	350	
33	f	908	
34	e	70	

2 Entry composition

There are 37 unique types of molecules in this entry. The entry contains 110481 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 Ubiquitin carboxyl-terminal hydrolase 14.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	x	494	Total	C	N	O	S	0	0
			3929	2485	647	769	28		

- Molecule 2 is a protein called 26S protease regulatory subunit 7.

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

- Molecule 3 is a protein called 26S protease regulatory subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	399	Total	C	N	O	S	0	0
			3112	1961	533	603	15		

- Molecule 4 is a protein called Isoform 2 of 26S proteasome regulatory subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	C	386	Total	C	N	O	S	0	0
			3017	1895	543	563	16		

- Molecule 5 is a protein called 26S protease regulatory subunit 6B.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	D	380	Total	C	N	O	S	0	0
			3039	1923	524	579	13		

- Molecule 6 is a protein called 26S proteasome regulatory subunit 10B.

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

- Molecule 7 is a protein called 26S protease regulatory subunit 6A.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	F	400	Total	C	N	O	S	0	0
			3133	1972	540	603	18		

- Molecule 8 is a protein called Proteasome subunit alpha type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	G	244	Total	C	N	O	S	0	0
			1889	1198	316	362	13		
8	g	244	Total	C	N	O	S	0	0
			1880	1193	318	356	13		

- Molecule 9 is a protein called Proteasome subunit alpha type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	H	232	Total	C	N	O	S	0	0
			1805	1152	305	342	6		
9	h	232	Total	C	N	O	S	0	0
			1805	1154	307	338	6		

- Molecule 10 is a protein called Proteasome subunit alpha type-4.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	I	250	Total	C	N	O	S	1	0
			1958	1236	336	376	10		
10	i	250	Total	C	N	O	S	0	0
			1955	1234	336	375	10		

- Molecule 11 is a protein called Proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	J	239	Total	C	N	O	S	0	0
			1880	1179	333	363	5		
11	j	239	Total	C	N	O	S	0	0
			1861	1168	332	356	5		

- Molecule 12 is a protein called Proteasome subunit alpha type-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	K	234	Total	C	N	O	S	0	0
			1777	1117	295	354	11		

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Mol	Chain	Residues	Atoms					AltConf	Trace
12	k	234	Total	C	N	O	S	0	0
			1782	1119	295	357	11		

- Molecule 13 is a protein called Isoform Long of Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	L	238	Total	C	N	O	S	0	0
			1866	1169	336	350	11		
13	l	238	Total	C	N	O	S	0	0
			1861	1165	335	350	11		

- Molecule 14 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	M	240	Total	C	N	O	S	0	0
			1876	1191	321	353	11		
14	m	240	Total	C	N	O	S	0	0
			1881	1193	321	356	11		

- Molecule 15 is a protein called Proteasome subunit beta type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	N	202	Total	C	N	O	S	0	0
			1514	949	258	295	12		
15	n	202	Total	C	N	O	S	0	0
			1510	947	258	293	12		

- Molecule 16 is a protein called Proteasome subunit beta type-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	O	220	Total	C	N	O	S	0	0
			1649	1038	279	320	12		
16	o	220	Total	C	N	O	S	0	0
			1659	1044	283	320	12		

- Molecule 17 is a protein called Proteasome subunit beta type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	P	204	Total	C	N	O	S	0	0
			1587	1010	264	294	19		
17	p	204	Total	C	N	O	S	0	0
			1591	1013	265	294	19		

- Molecule 18 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	Q	199	Total	C	N	O	S	0	0
			1588	1017	270	292	9		
18	q	199	Total	C	N	O	S	0	0
			1588	1017	270	292	9		

- Molecule 19 is a protein called Proteasome subunit beta type-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	R	201	Total	C	N	O	S	0	0
			1559	982	274	294	9		
19	r	201	Total	C	N	O	S	0	0
			1559	982	274	294	9		

- Molecule 20 is a protein called Proteasome subunit beta type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	S	213	Total	C	N	O	S	0	0
			1641	1041	281	309	10		
20	s	213	Total	C	N	O	S	0	0
			1654	1047	284	313	10		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
21	T	216	Total	C	N	O	S	0	0
			1683	1062	291	318	12		
21	t	216	Total	C	N	O	S	0	0
			1687	1064	291	320	12		

- Molecule 22 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	y	76	Total	C	N	O	S	0	0
			601	378	105	117	1		

- Molecule 23 is a protein called 26S proteasome non-ATPase regulatory subunit 1.

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

- Molecule 24 is a protein called 26S proteasome non-ATPase regulatory subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	V	444	Total	C	N	O	S	0	0
			3612	2301	645	653	13		

- Molecule 25 is a protein called 26S proteasome non-ATPase regulatory subunit 12.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	W	446	Total	C	N	O	S	0	0
			3635	2302	622	687	24		

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

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

- Molecule 27 is a protein called 26S proteasome non-ATPase regulatory subunit 6.

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

- Molecule 28 is a protein called 26S proteasome non-ATPase regulatory subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	Z	286	Total	C	N	O	S	0	0
			2281	1457	392	427	5		

- Molecule 29 is a protein called 26S proteasome non-ATPase regulatory subunit 13.

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

- Molecule 30 is a protein called 26S proteasome non-ATPase regulatory subunit 4.

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

- Molecule 31 is a protein called 26S proteasome non-ATPase regulatory subunit 14.

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

- Molecule 32 is a protein called 26S proteasome non-ATPase regulatory subunit 8.

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

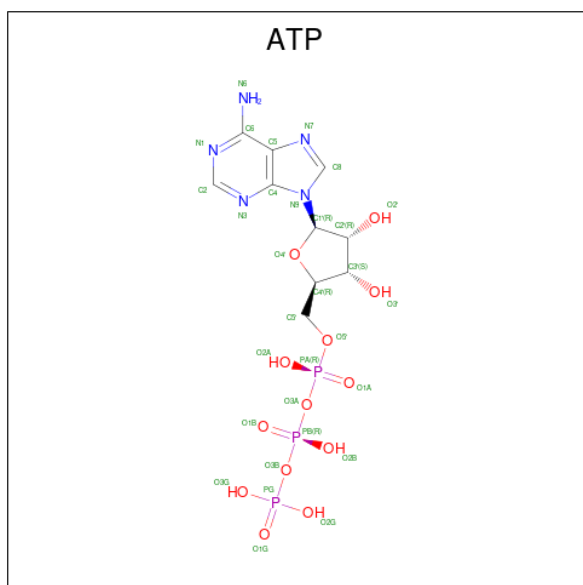
- Molecule 33 is a protein called 26S proteasome non-ATPase regulatory subunit 2.

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

- Molecule 34 is a protein called 26S proteasome complex subunit DSS1.

Mol	Chain	Residues	Atoms				AltConf	Trace
34	e	50	Total	C	N	O	0	0
			425	260	65	100		

- Molecule 35 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).



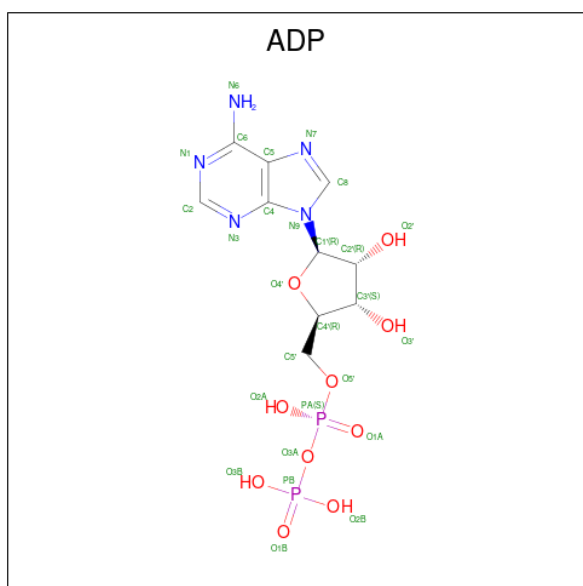
Mol	Chain	Residues	Atoms					AltConf
35	A	1	Total	C	N	O	P	0
			31	10	5	13	3	

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

- Molecule 36 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



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

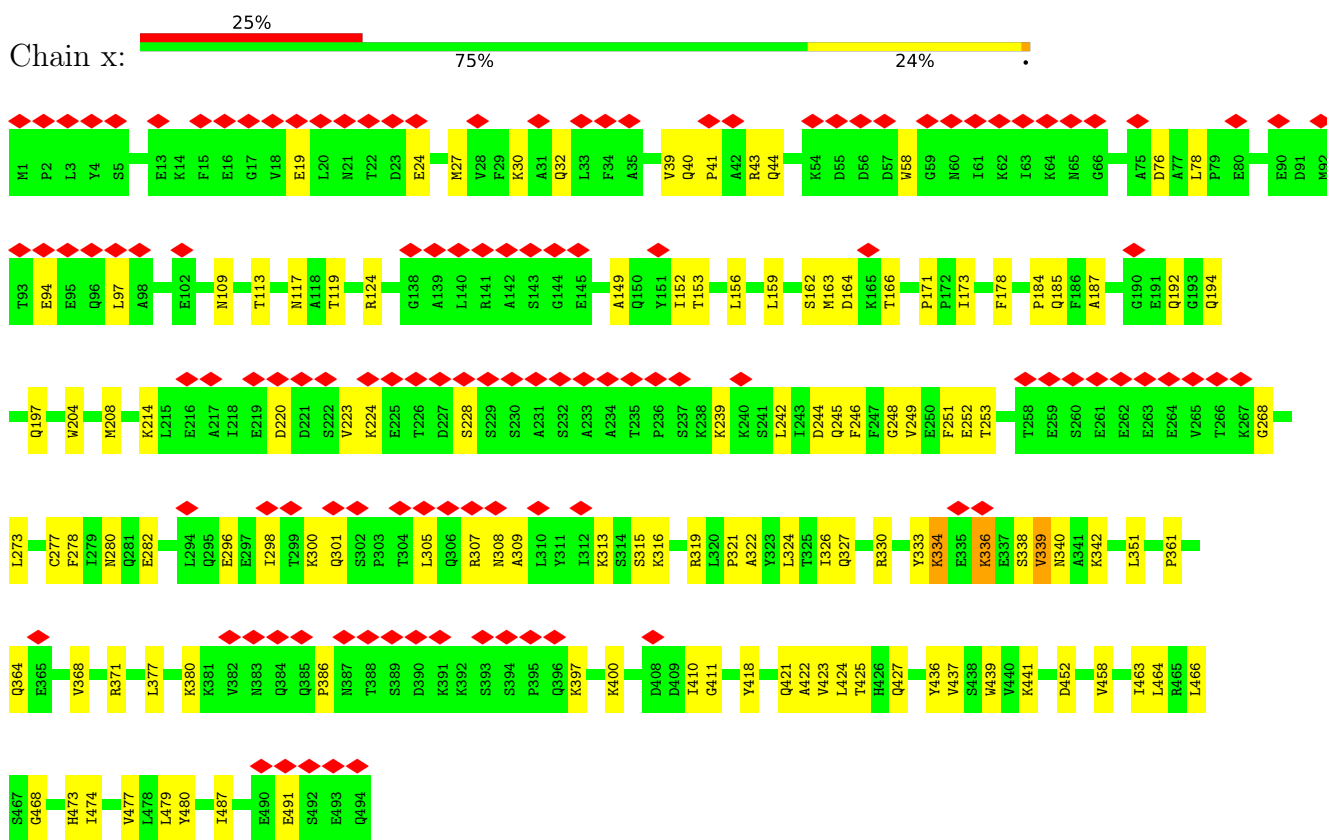
- Molecule 37 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

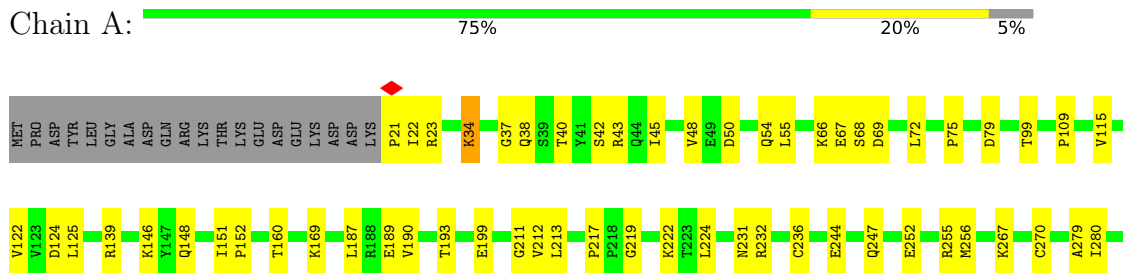
3 Residue-property plots

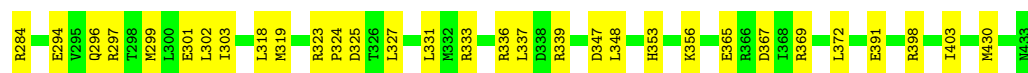
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Ubiquitin carboxyl-terminal hydrolase 14

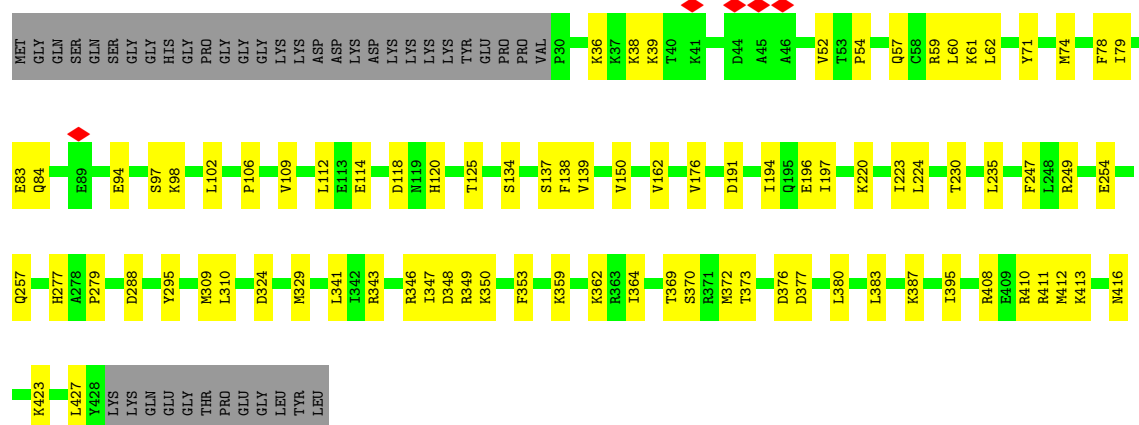


• Molecule 2: 26S protease regulatory subunit 7

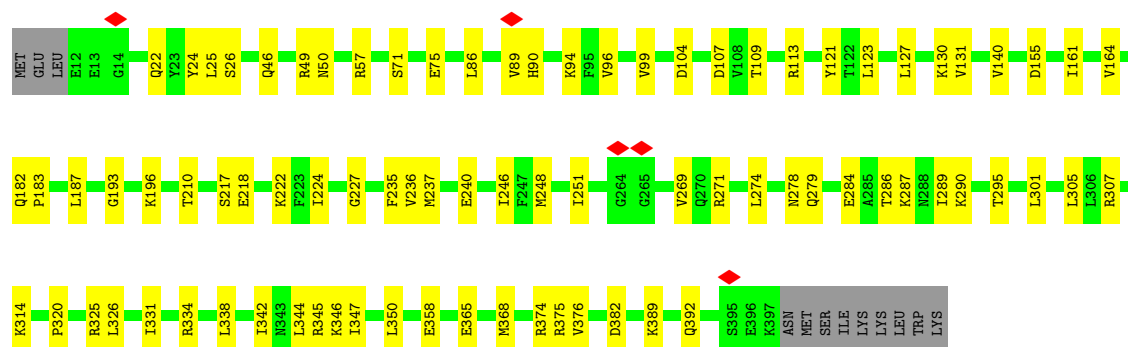
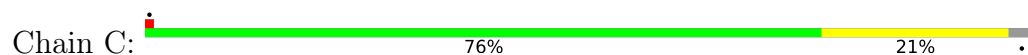




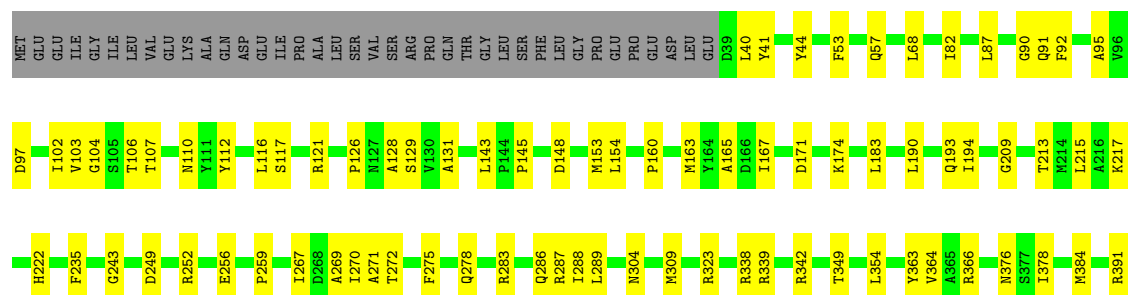
• Molecule 3: 26S protease regulatory subunit 4



• Molecule 4: Isoform 2 of 26S proteasome regulatory subunit 8



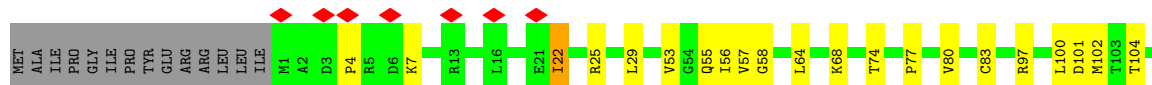
• Molecule 5: 26S protease regulatory subunit 6B





- Molecule 6: 26S proteasome regulatory subunit 10B

Chain E: 79% 17%



- Molecule 7: 26S protease regulatory subunit 6A

Chain F: 76% 15% 9%



- Molecule 8: Proteasome subunit alpha type-6

Chain G: 91% 9%



- Molecule 8: Proteasome subunit alpha type-6

Chain g: 91% 8%



- Molecule 9: Proteasome subunit alpha type-2

Chain H: 91% 9%



- Molecule 9: Proteasome subunit alpha type-2

Chain h: 94% 6%



- Molecule 10: Proteasome subunit alpha type-4

Chain I: 90% 6%



- Molecule 10: Proteasome subunit alpha type-4

Chain i: 90% 6%



- Molecule 11: Proteasome subunit alpha type-7

Chain J: 88% 8%



- Molecule 11: Proteasome subunit alpha type-7

Chain j: 87% 9%



- Molecule 12: Proteasome subunit alpha type-5

Chain K: 87% 10%



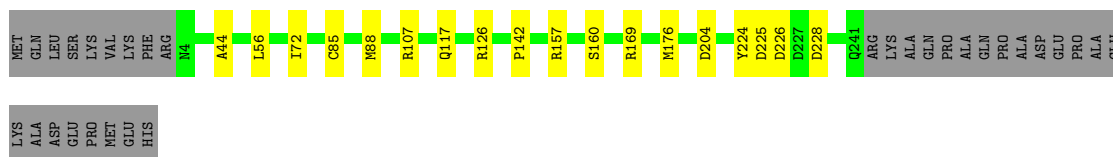
- Molecule 12: Proteasome subunit alpha type-5

Chain k: 93% 6%



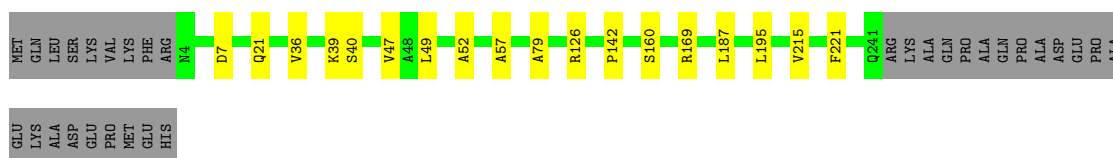
- Molecule 13: Isoform Long of Proteasome subunit alpha type-1

Chain L: 82% 7% 12%



- Molecule 13: Isoform Long of Proteasome subunit alpha type-1

Chain I: 82% 7% 12%



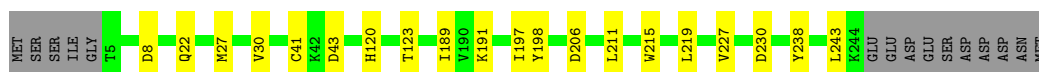
- Molecule 14: Proteasome subunit alpha type-3

Chain M: 87% 7% 6%



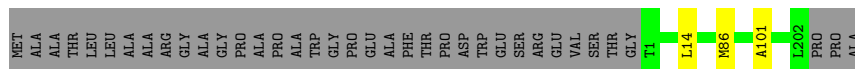
- Molecule 14: Proteasome subunit alpha type-3

Chain m: 86% 8% 6%



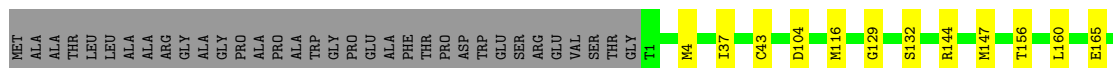
- Molecule 15: Proteasome subunit beta type-6

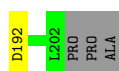
Chain N: 83% 15%



- Molecule 15: Proteasome subunit beta type-6

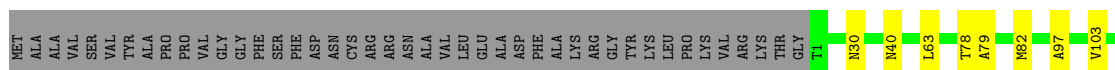
Chain n: 79% 5% 15%





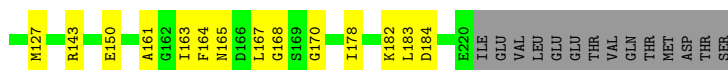
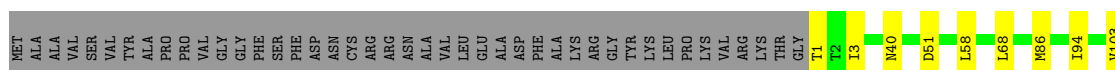
- Molecule 16: Proteasome subunit beta type-7

Chain O: 73% 7% 21%



- Molecule 16: Proteasome subunit beta type-7

Chain o: 71% 8% 21%



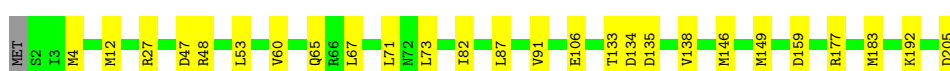
- Molecule 17: Proteasome subunit beta type-3

Chain P: 94% 5%



- Molecule 17: Proteasome subunit beta type-3

Chain p: 87% 13%



- Molecule 18: Proteasome subunit beta type-2

Chain Q: 87% 12%



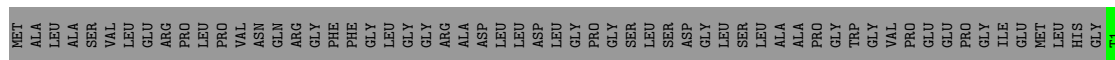
- Molecule 18: Proteasome subunit beta type-2

Chain q: 89% 10%



- Molecule 19: Proteasome subunit beta type-5

Chain R: 72% 5% 24%



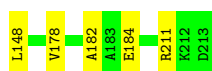
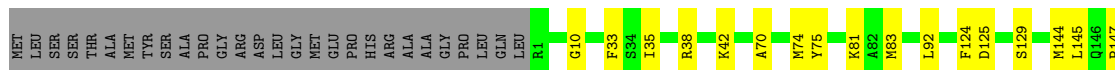
- Molecule 19: Proteasome subunit beta type-5

Chain r: 72% 5% 24%



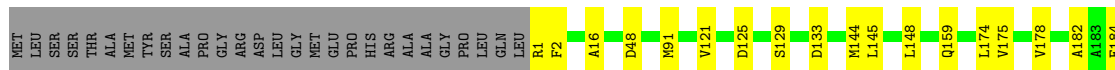
- Molecule 20: Proteasome subunit beta type-1

Chain S: 79% 9% 12%



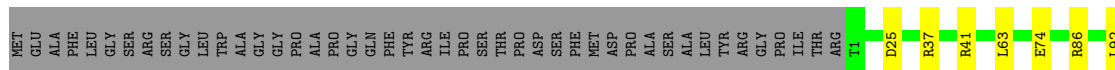
- Molecule 20: Proteasome subunit beta type-1

Chain s: 80% 9% 12%



- Molecule 21: Proteasome subunit beta type-4

Chain T: 75% 6% 18%





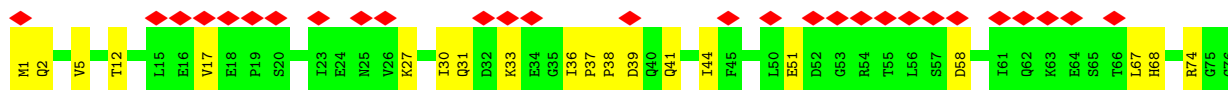
• Molecule 21: Proteasome subunit beta type-4

Chain t: 76% 6% 18%



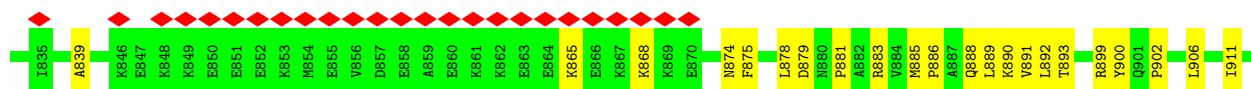
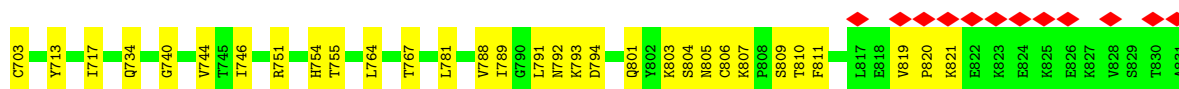
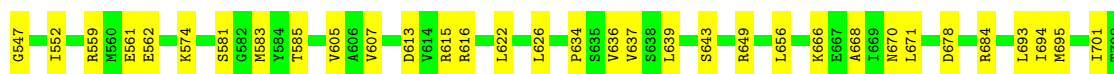
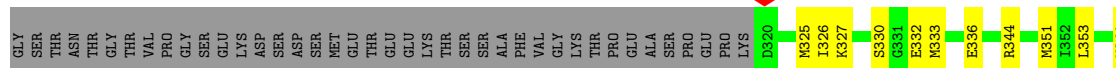
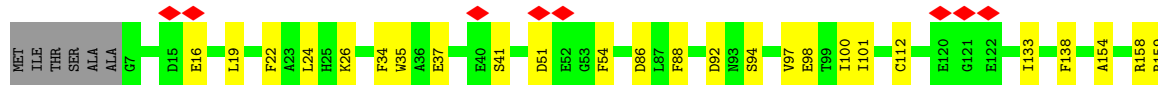
• Molecule 22: Ubiquitin

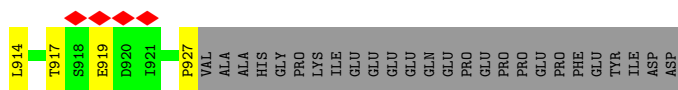
Chain y: 37% 74% 26%



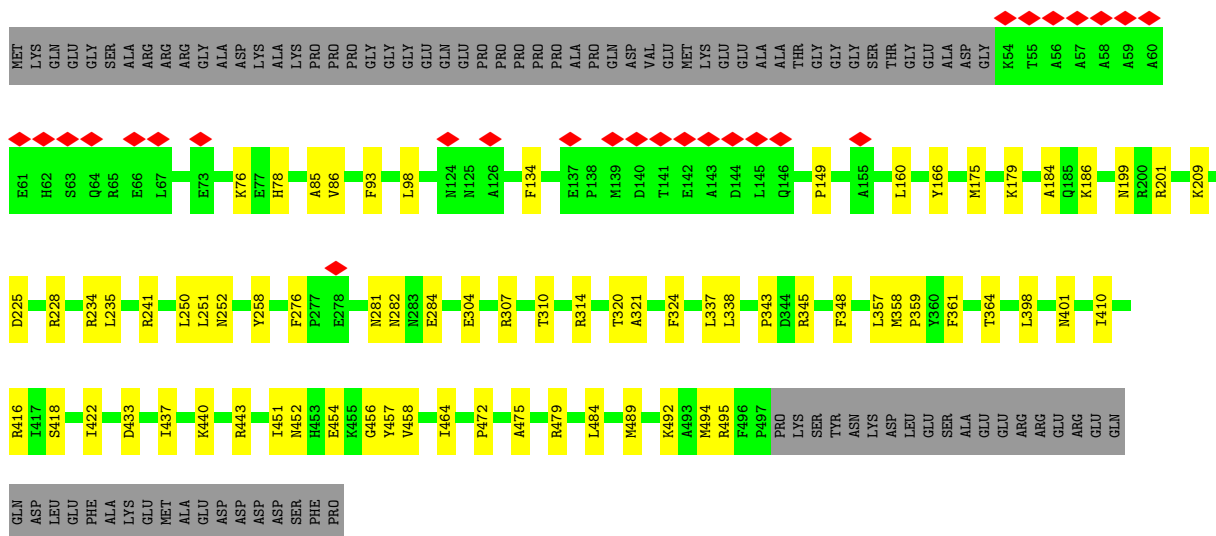
• Molecule 23: 26S proteasome non-ATPase regulatory subunit 1

Chain U: 5% 72% 20% 8%

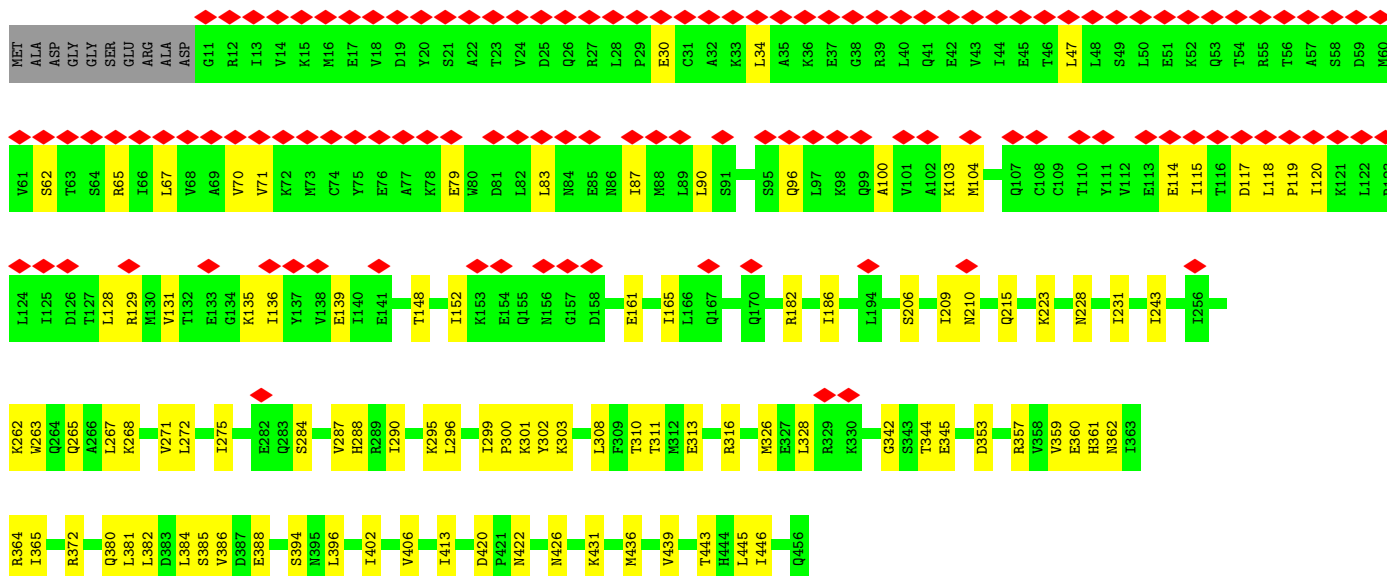
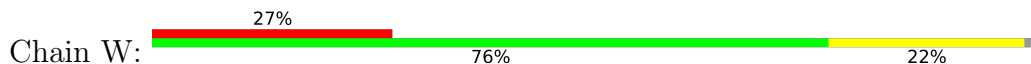




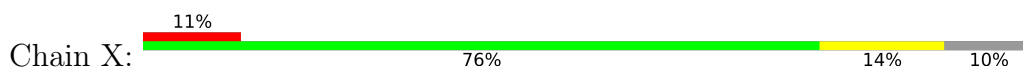
- Molecule 24: 26S proteasome non-ATPase regulatory subunit 3

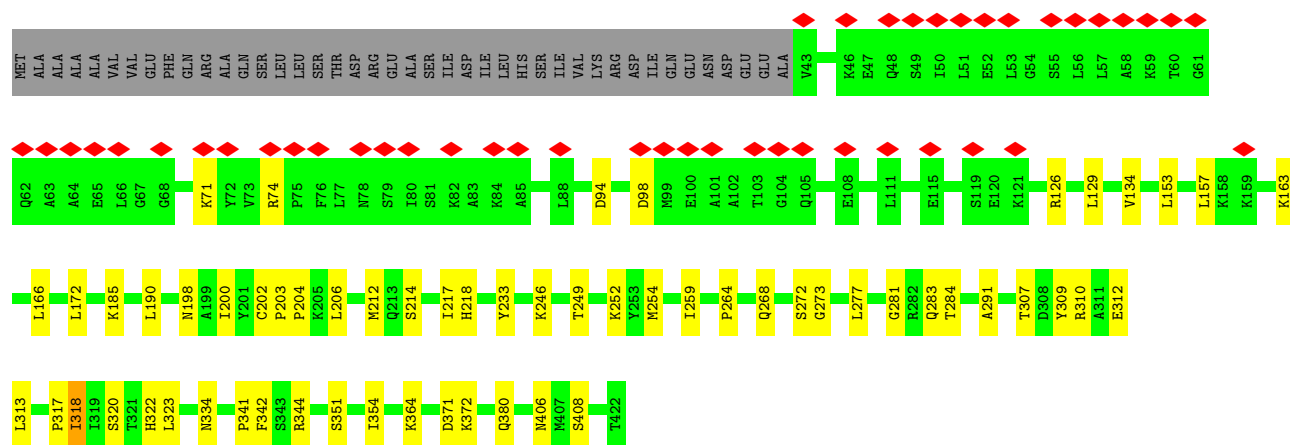


- Molecule 25: 26S proteasome non-ATPase regulatory subunit 12

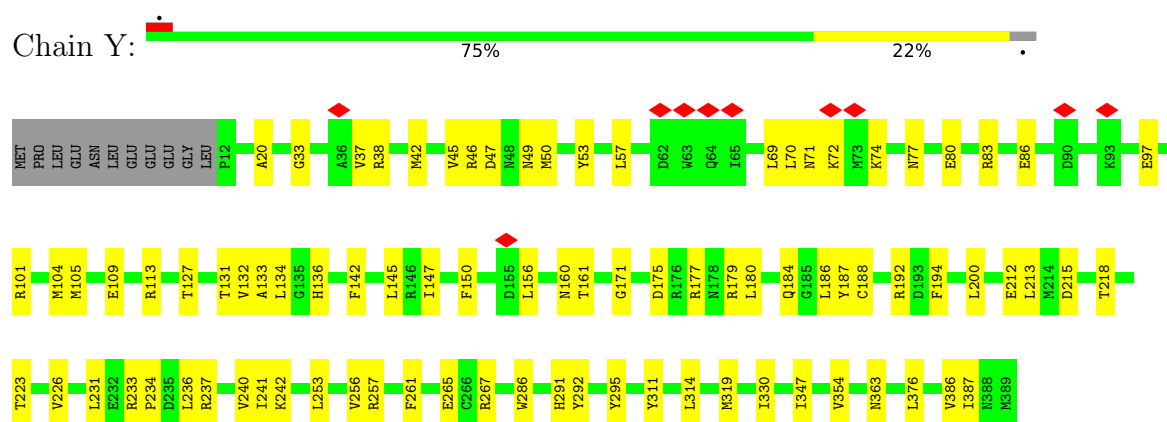


- Molecule 26: 26S proteasome non-ATPase regulatory subunit 11

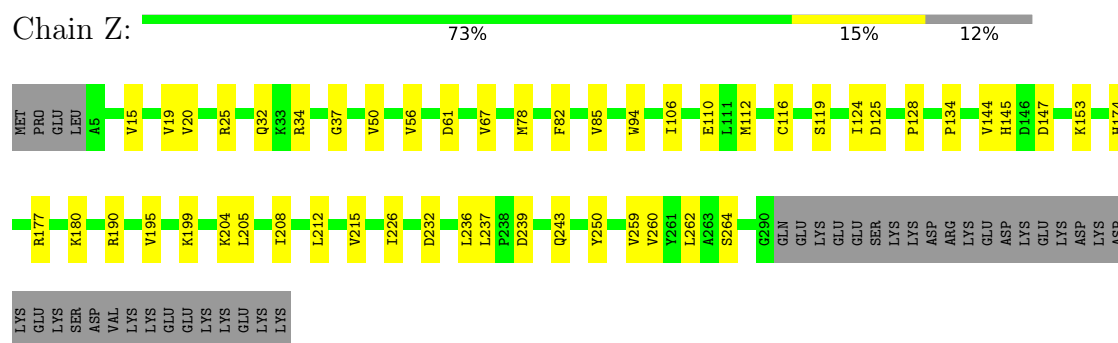




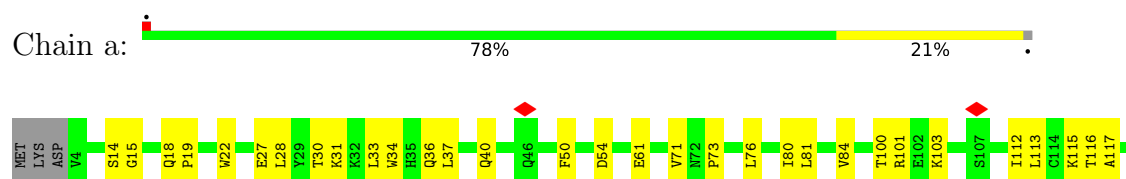
- Molecule 27: 26S proteasome non-ATPase regulatory subunit 6

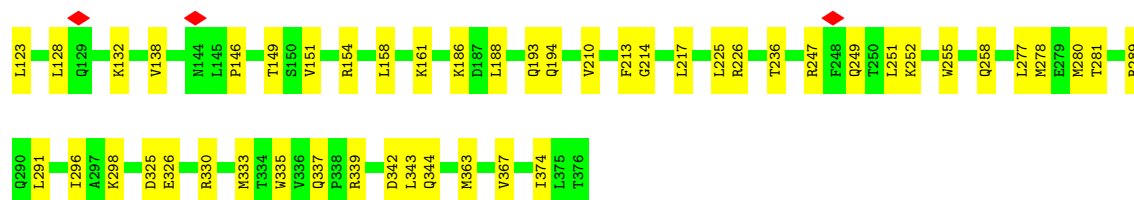


- Molecule 28: 26S proteasome non-ATPase regulatory subunit 7



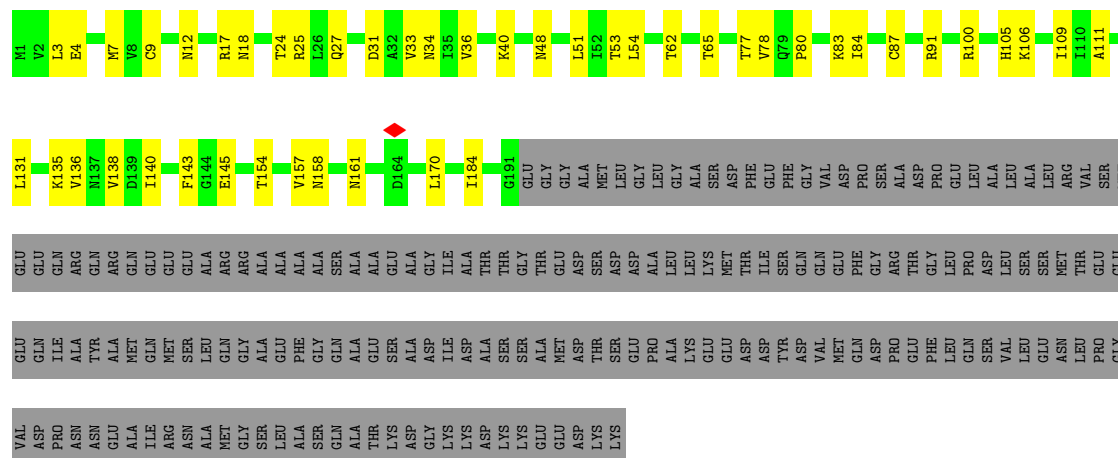
- Molecule 29: 26S proteasome non-ATPase regulatory subunit 13





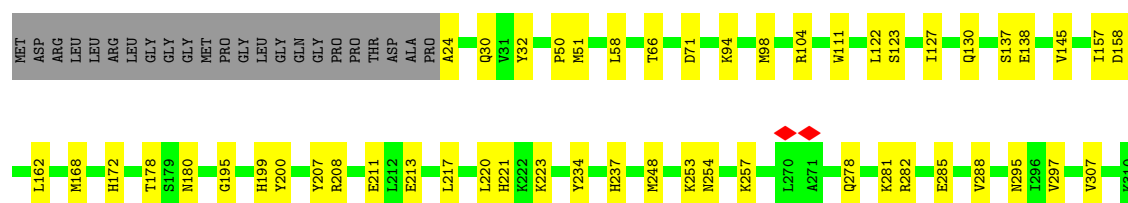
• Molecule 30: 26S proteasome non-ATPase regulatory subunit 4

Chain b: 38% 12% 49%



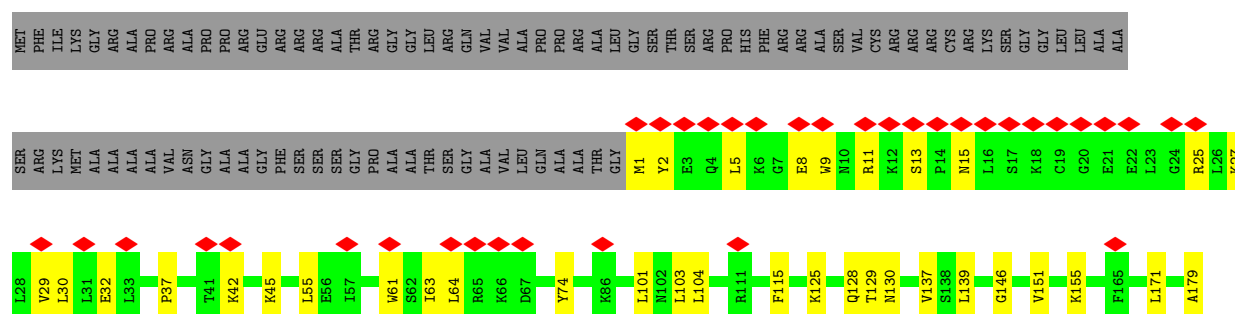
• Molecule 31: 26S proteasome non-ATPase regulatory subunit 14

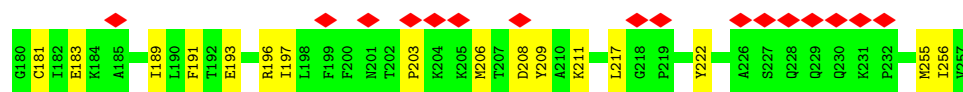
Chain c: 76% 16% 7%



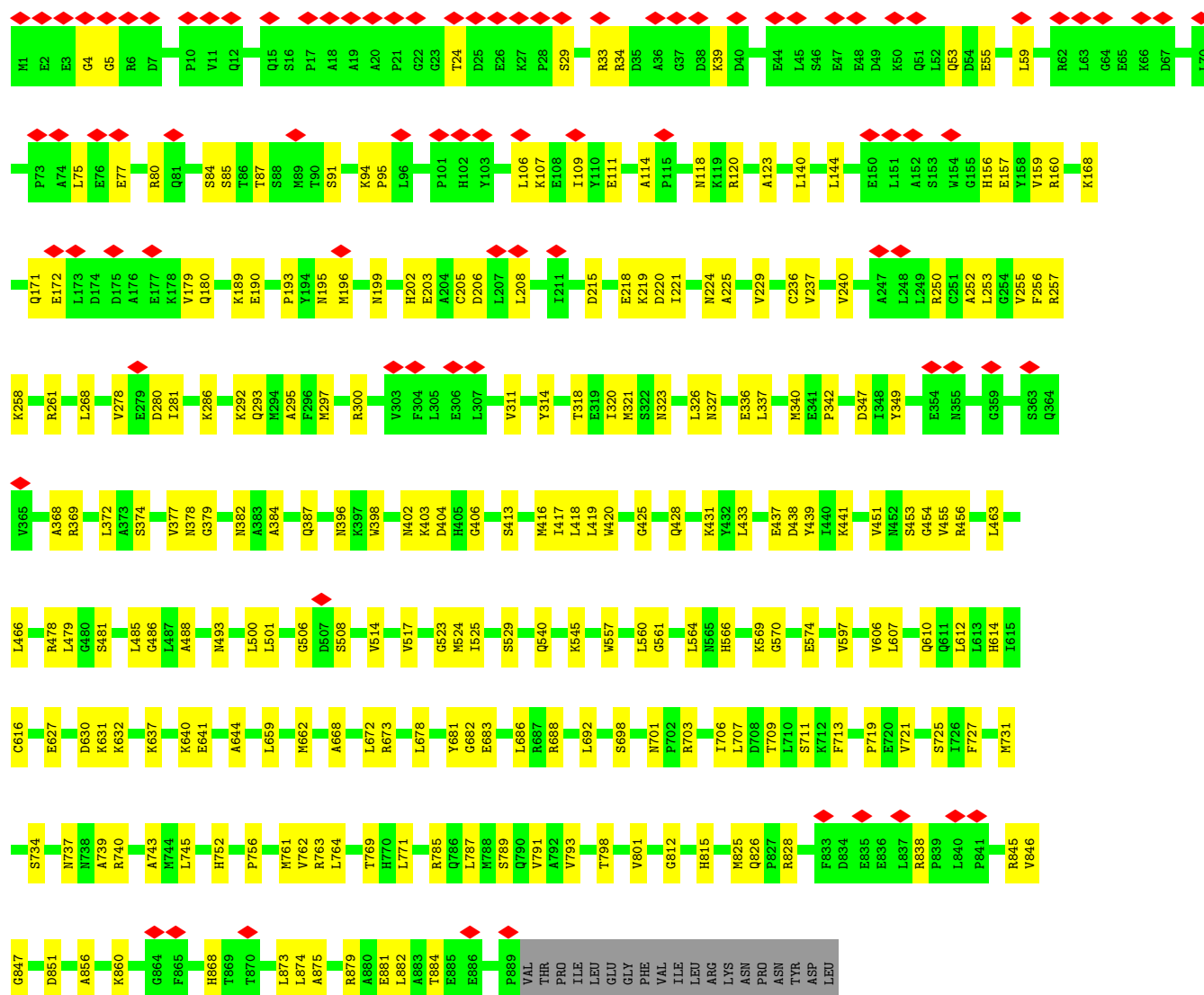
• Molecule 32: 26S proteasome non-ATPase regulatory subunit 8

Chain d: 15% 59% 15% 27%

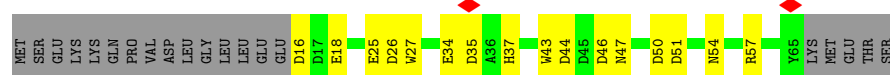




- Molecule 33: 26S proteasome non-ATPase regulatory subunit 2



- Molecule 34: 26S proteasome complex subunit DSS1



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	53225	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	400	Depositor
Maximum defocus (nm)	5000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.020	Depositor
Minimum map value	-0.005	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.005	Depositor
Map size ($\text{\AA}$)	438.4, 438.4, 438.4	wwPDB
Map dimensions	640, 640, 640	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.685, 0.685, 0.685	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: ADP, ATP, ZN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	x	0.16	0/4002	0.42	0/5390
2	A	0.21	0/3283	0.50	2/4433 (0.0%)
3	B	0.21	0/3157	0.52	0/4258
4	C	0.18	0/3056	0.46	0/4108
5	D	0.18	0/3089	0.47	0/4168
6	E	0.17	0/3145	0.42	0/4233
7	F	0.19	0/3173	0.47	0/4273
8	G	0.25	0/1923	0.38	0/2601
8	g	0.24	0/1914	0.39	0/2590
9	H	0.26	0/1844	0.40	0/2499
9	h	0.25	0/1844	0.38	0/2497
10	I	0.24	0/1991	0.42	0/2685
10	i	0.23	0/1985	0.41	0/2677
11	J	0.20	0/1906	0.41	0/2573
11	j	0.21	0/1887	0.39	0/2549
12	K	0.24	0/1804	0.41	0/2436
12	k	0.22	0/1809	0.35	0/2444
13	L	0.25	0/1901	0.36	0/2570
13	l	0.25	0/1896	0.37	0/2565
14	M	0.25	0/1911	0.39	0/2573
14	m	0.24	0/1916	0.41	0/2580
15	N	0.24	0/1540	0.31	0/2085
15	n	0.24	0/1536	0.35	0/2080
16	O	0.25	0/1676	0.42	0/2271
16	o	0.25	0/1686	0.40	0/2282
17	P	0.26	0/1616	0.42	0/2180
17	p	0.27	0/1620	0.39	0/2184
18	Q	0.25	0/1621	0.37	0/2194
18	q	0.26	0/1621	0.39	0/2194
19	R	0.24	0/1590	0.35	0/2147
19	r	0.24	0/1590	0.36	0/2147
20	S	0.26	0/1671	0.39	0/2252

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
20	s	0.25	0/1684	0.36	0/2268
21	T	0.26	0/1716	0.40	0/2323
21	t	0.25	0/1720	0.38	0/2328
22	y	0.20	0/607	0.49	0/816
23	U	0.19	0/6945	0.50	0/9382
24	V	0.17	0/3681	0.44	0/4969
25	W	0.25	0/3683	0.51	0/4952
26	X	0.17	0/3053	0.46	2/4115 (0.0%)
27	Y	0.19	0/3173	0.52	0/4273
28	Z	0.21	0/2324	0.53	0/3150
29	a	0.18	0/3053	0.52	0/4133
30	b	0.20	0/1478	0.51	0/2001
31	c	0.22	0/2302	0.54	0/3110
32	d	0.21	0/2162	0.52	0/2919
33	f	0.20	0/6980	0.52	0/9433
34	e	0.22	0/437	0.53	0/595
All	All	0.22	0/112201	0.45	4/151485 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	48	VAL	N-CA-C	-5.54	107.42	112.96
2	A	37	GLY	N-CA-C	-5.26	104.95	114.46
26	X	317	PRO	CA-C-N	5.19	131.31	121.97
26	X	317	PRO	C-N-CA	5.19	131.31	121.97

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	x	3929	0	3915	72	0
2	A	3229	0	3261	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	3112	0	3181	63	0
4	C	3017	0	3117	56	0
5	D	3039	0	3076	63	0
6	E	3097	0	3174	50	0
7	F	3133	0	3220	46	0
8	G	1889	0	1885	13	0
8	g	1880	0	1875	12	0
9	H	1805	0	1784	13	0
9	h	1805	0	1798	10	0
10	I	1958	0	1960	10	0
10	i	1955	0	1955	10	0
11	J	1880	0	1892	14	0
11	j	1861	0	1865	15	0
12	K	1777	0	1762	16	0
12	k	1782	0	1766	7	0
13	L	1866	0	1852	14	0
13	l	1861	0	1839	11	0
14	M	1876	0	1861	12	0
14	m	1881	0	1868	12	0
15	N	1514	0	1487	2	0
15	n	1510	0	1483	10	0
16	O	1649	0	1659	13	0
16	o	1659	0	1681	13	0
17	P	1587	0	1598	8	0
17	p	1591	0	1609	18	0
18	Q	1588	0	1584	17	0
18	q	1588	0	1584	16	0
19	R	1559	0	1523	10	0
19	r	1559	0	1523	7	0
20	S	1641	0	1639	17	0
20	s	1654	0	1656	14	0
21	T	1683	0	1662	13	0
21	t	1687	0	1666	10	0
22	y	601	0	629	14	0
23	U	6828	0	6884	115	0
24	V	3612	0	3682	47	0
25	W	3635	0	3762	61	0
26	X	3009	0	3113	38	0
27	Y	3115	0	3120	53	0
28	Z	2281	0	2312	34	0
29	a	2995	0	3012	52	0
30	b	1458	0	1505	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
31	c	2260	0	2276	31	0
32	d	2116	0	2146	35	0
33	f	6866	0	6866	160	0
34	e	425	0	328	9	0
35	A	31	0	12	2	0
35	D	31	0	12	3	0
35	E	31	0	12	1	0
35	F	31	0	12	1	0
36	B	27	0	12	1	0
36	C	27	0	12	2	0
37	c	1	0	0	0	0
All	All	110481	0	110967	1277	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 6.

The worst 5 of 1277 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)
12:K:195:ILE:HD11	12:K:217:LEU:HD11	1.62	0.81
33:f:286:LYS:HG3	33:f:873:LEU:HD12	1.64	0.79
5:D:267:ILE:HG12	5:D:309:MET:HE2	1.69	0.74
23:U:479:LEU:HG	23:U:514:LEU:HD11	1.72	0.72
31:c:50:PRO:HG2	31:c:51:MET:HE2	1.71	0.72

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	x	492/494 (100%)	463 (94%)	29 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	411/433 (95%)	365 (89%)	46 (11%)	0	100	100
3	B	397/440 (90%)	352 (89%)	45 (11%)	0	100	100
4	C	384/398 (96%)	341 (89%)	42 (11%)	1 (0%)	36	67
5	D	378/418 (90%)	344 (91%)	34 (9%)	0	100	100
6	E	387/403 (96%)	353 (91%)	33 (8%)	1 (0%)	36	67
7	F	396/439 (90%)	356 (90%)	40 (10%)	0	100	100
8	G	242/246 (98%)	231 (96%)	11 (4%)	0	100	100
8	g	242/246 (98%)	226 (93%)	16 (7%)	0	100	100
9	H	230/234 (98%)	220 (96%)	10 (4%)	0	100	100
9	h	230/234 (98%)	217 (94%)	13 (6%)	0	100	100
10	I	249/261 (95%)	240 (96%)	9 (4%)	0	100	100
10	i	248/261 (95%)	240 (97%)	8 (3%)	0	100	100
11	J	237/248 (96%)	223 (94%)	14 (6%)	0	100	100
11	j	237/248 (96%)	221 (93%)	16 (7%)	0	100	100
12	K	232/241 (96%)	215 (93%)	16 (7%)	1 (0%)	30	62
12	k	232/241 (96%)	221 (95%)	11 (5%)	0	100	100
13	L	236/269 (88%)	231 (98%)	5 (2%)	0	100	100
13	l	236/269 (88%)	227 (96%)	9 (4%)	0	100	100
14	M	238/255 (93%)	231 (97%)	7 (3%)	0	100	100
14	m	238/255 (93%)	232 (98%)	6 (2%)	0	100	100
15	N	200/239 (84%)	197 (98%)	3 (2%)	0	100	100
15	n	200/239 (84%)	190 (95%)	10 (5%)	0	100	100
16	O	218/277 (79%)	209 (96%)	9 (4%)	0	100	100
16	o	218/277 (79%)	210 (96%)	8 (4%)	0	100	100
17	P	202/205 (98%)	189 (94%)	13 (6%)	0	100	100
17	p	202/205 (98%)	191 (95%)	11 (5%)	0	100	100
18	Q	197/201 (98%)	190 (96%)	7 (4%)	0	100	100
18	q	197/201 (98%)	191 (97%)	6 (3%)	0	100	100
19	R	199/263 (76%)	193 (97%)	6 (3%)	0	100	100
19	r	199/263 (76%)	193 (97%)	6 (3%)	0	100	100
20	S	211/241 (88%)	203 (96%)	8 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
20	s	211/241 (88%)	203 (96%)	8 (4%)	0	100	100
21	T	214/264 (81%)	209 (98%)	5 (2%)	0	100	100
21	t	214/264 (81%)	206 (96%)	8 (4%)	0	100	100
22	y	74/76 (97%)	71 (96%)	3 (4%)	0	100	100
23	U	868/953 (91%)	806 (93%)	62 (7%)	0	100	100
24	V	442/534 (83%)	428 (97%)	14 (3%)	0	100	100
25	W	444/456 (97%)	417 (94%)	27 (6%)	0	100	100
26	X	378/422 (90%)	366 (97%)	11 (3%)	1 (0%)	36	67
27	Y	376/389 (97%)	338 (90%)	37 (10%)	1 (0%)	36	67
28	Z	284/324 (88%)	248 (87%)	34 (12%)	2 (1%)	18	51
29	a	371/376 (99%)	339 (91%)	31 (8%)	1 (0%)	36	67
30	b	189/377 (50%)	172 (91%)	17 (9%)	0	100	100
31	c	285/310 (92%)	248 (87%)	37 (13%)	0	100	100
32	d	255/350 (73%)	218 (86%)	37 (14%)	0	100	100
33	f	887/908 (98%)	792 (89%)	95 (11%)	0	100	100
34	e	48/70 (69%)	43 (90%)	5 (10%)	0	100	100
All	All	13955/15458 (90%)	13009 (93%)	938 (7%)	8 (0%)	49	79

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
26	X	318	ILE
28	Z	134	PRO
29	a	343	LEU
27	Y	292	TYR
28	Z	145	HIS

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	x	439/439 (100%)	431 (98%)	8 (2%)	51	69
2	A	348/372 (94%)	344 (99%)	4 (1%)	65	74
3	B	347/385 (90%)	345 (99%)	2 (1%)	78	79
4	C	330/346 (95%)	328 (99%)	2 (1%)	78	79
5	D	333/366 (91%)	333 (100%)	0	100	100
6	E	341/353 (97%)	341 (100%)	0	100	100
7	F	344/379 (91%)	343 (100%)	1 (0%)	86	83
8	G	205/210 (98%)	205 (100%)	0	100	100
8	g	202/210 (96%)	202 (100%)	0	100	100
9	H	188/191 (98%)	188 (100%)	0	100	100
9	h	188/191 (98%)	188 (100%)	0	100	100
10	I	207/221 (94%)	205 (99%)	2 (1%)	68	75
10	i	206/221 (93%)	206 (100%)	0	100	100
11	J	201/211 (95%)	201 (100%)	0	100	100
11	j	196/211 (93%)	196 (100%)	0	100	100
12	K	193/203 (95%)	193 (100%)	0	100	100
12	k	195/203 (96%)	195 (100%)	0	100	100
13	L	202/230 (88%)	202 (100%)	0	100	100
13	l	201/230 (87%)	201 (100%)	0	100	100
14	M	196/212 (92%)	196 (100%)	0	100	100
14	m	198/212 (93%)	198 (100%)	0	100	100
15	N	157/181 (87%)	157 (100%)	0	100	100
15	n	156/181 (86%)	156 (100%)	0	100	100
16	O	179/228 (78%)	179 (100%)	0	100	100
16	o	181/228 (79%)	181 (100%)	0	100	100
17	P	172/174 (99%)	172 (100%)	0	100	100
17	p	173/174 (99%)	173 (100%)	0	100	100
18	Q	168/171 (98%)	168 (100%)	0	100	100
18	q	168/171 (98%)	168 (100%)	0	100	100
19	R	156/202 (77%)	156 (100%)	0	100	100
19	r	156/202 (77%)	156 (100%)	0	100	100
20	S	175/199 (88%)	175 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
20	s	178/199 (89%)	178 (100%)	0	100	100
21	T	178/215 (83%)	178 (100%)	0	100	100
21	t	179/215 (83%)	179 (100%)	0	100	100
22	y	68/68 (100%)	68 (100%)	0	100	100
23	U	748/816 (92%)	748 (100%)	0	100	100
24	V	390/460 (85%)	390 (100%)	0	100	100
25	W	410/416 (99%)	408 (100%)	2 (0%)	81	80
26	X	327/362 (90%)	327 (100%)	0	100	100
27	Y	334/344 (97%)	334 (100%)	0	100	100
28	Z	257/295 (87%)	255 (99%)	2 (1%)	73	77
29	a	333/336 (99%)	333 (100%)	0	100	100
30	b	167/312 (54%)	166 (99%)	1 (1%)	78	79
31	c	252/268 (94%)	249 (99%)	3 (1%)	63	73
32	d	231/294 (79%)	231 (100%)	0	100	100
33	f	745/763 (98%)	744 (100%)	1 (0%)	88	87
34	e	44/63 (70%)	40 (91%)	4 (9%)	9	32
All	All	11942/13133 (91%)	11910 (100%)	32 (0%)	84	83

5 of 32 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
34	e	25	GLU
34	e	26	ASP
3	B	102	LEU
2	A	403	ILE
34	e	27	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 129 such sidechains are listed below:

Mol	Chain	Res	Type
31	c	166	ASN
33	f	323	ASN
19	R	119	ASN
19	R	62	GLN
33	f	382	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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$
35	ATP	D	501	-	32,33,33	0.32	0	48,52,52	0.30	0
36	ADP	C	501	-	28,29,29	1.37	3 (10%)	43,45,45	1.96	9 (20%)
35	ATP	F	501	-	32,33,33	0.63	1 (3%)	48,52,52	0.36	0
35	ATP	E	501	-	32,33,33	0.46	0	48,52,52	0.32	0
35	ATP	A	501	-	32,33,33	0.60	1 (3%)	48,52,52	0.34	0
36	ADP	B	501	-	28,29,29	1.37	4 (14%)	43,45,45	1.94	9 (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
35	ATP	D	501	-	-	5/22/38/38	0/3/3/3
36	ADP	C	501	-	-	3/16/32/32	0/3/3/3
35	ATP	F	501	-	-	6/22/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
35	ATP	E	501	-	-	2/22/38/38	0/3/3/3
35	ATP	A	501	-	-	6/22/38/38	0/3/3/3
36	ADP	B	501	-	-	3/16/32/32	0/3/3/3

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
36	B	501	ADP	C5-C4	4.67	1.47	1.39
36	C	501	ADP	C5-C4	4.51	1.47	1.39
36	C	501	ADP	C5-N7	-2.62	1.34	1.39
35	F	501	ATP	PA-O3A	-2.59	1.56	1.59
36	C	501	ADP	C5-C6	2.52	1.48	1.41

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	B	501	ADP	C5-C4-N3	-6.61	117.61	126.72
36	C	501	ADP	C5-C4-N3	-6.27	118.08	126.72
36	B	501	ADP	N3-C4-N9	5.55	136.60	127.17
36	C	501	ADP	N3-C4-N9	5.27	136.14	127.17
36	B	501	ADP	C2-N3-C4	3.92	121.42	111.83

There are no chirality outliers.

5 of 25 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
35	A	501	ATP	C5'-O5'-PA-O1A
35	A	501	ATP	C5'-O5'-PA-O3A
35	D	501	ATP	C5'-O5'-PA-O3A
35	E	501	ATP	PB-O3B-PG-O2G
35	F	501	ATP	C5'-O5'-PA-O1A

There are no ring outliers.

6 monomers are involved in 10 short contacts:

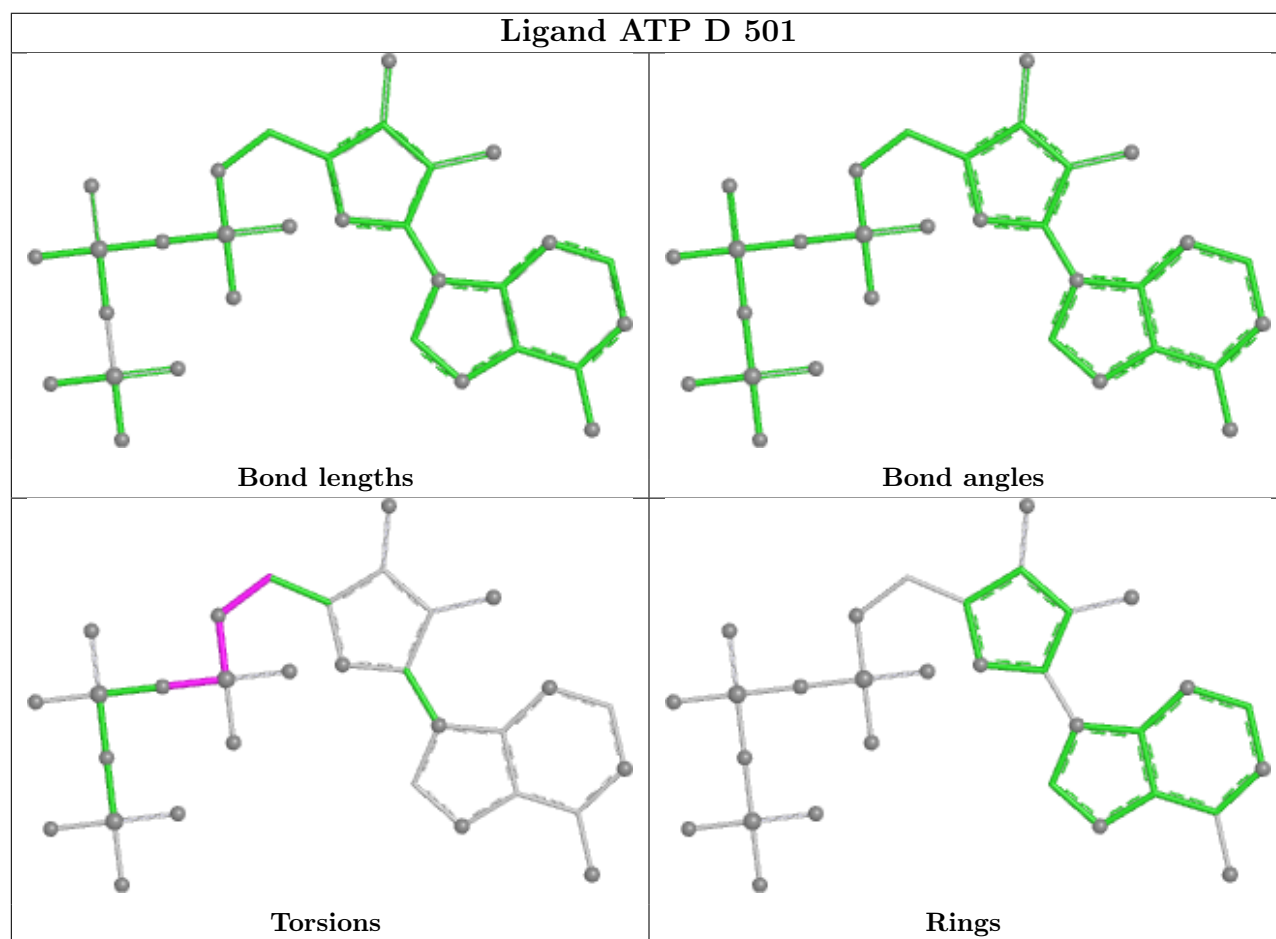
Mol	Chain	Res	Type	Clashes	Symm-Clashes
35	D	501	ATP	3	0
36	C	501	ADP	2	0
35	F	501	ATP	1	0
35	E	501	ATP	1	0

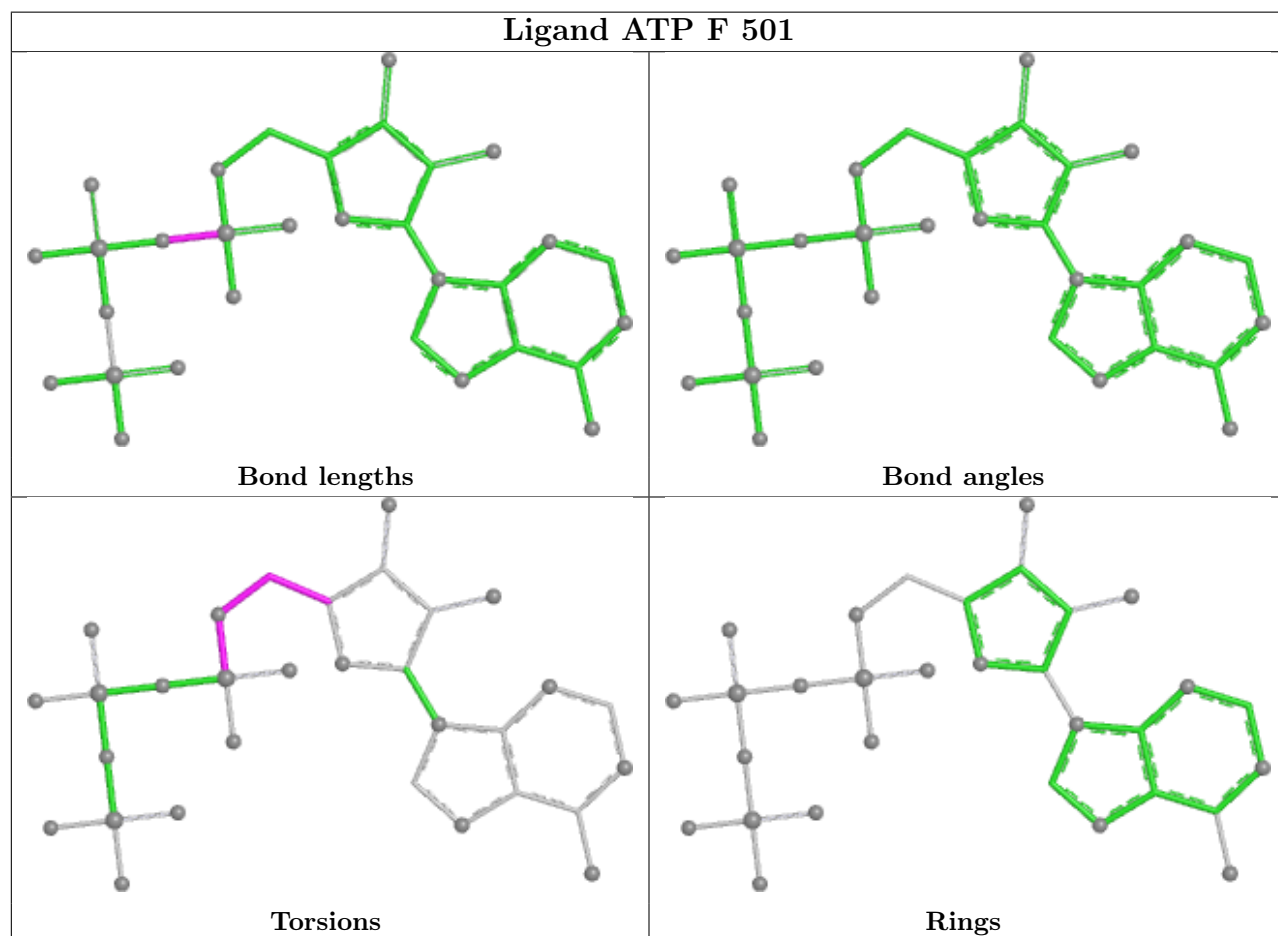
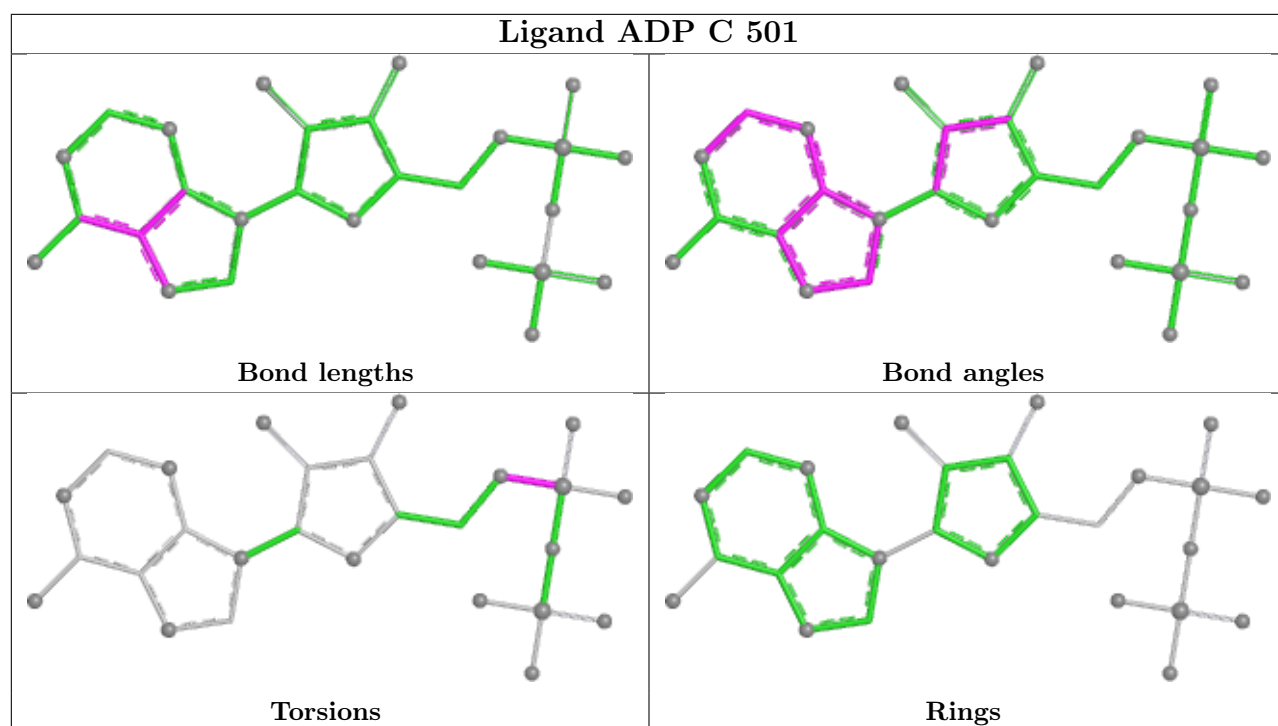
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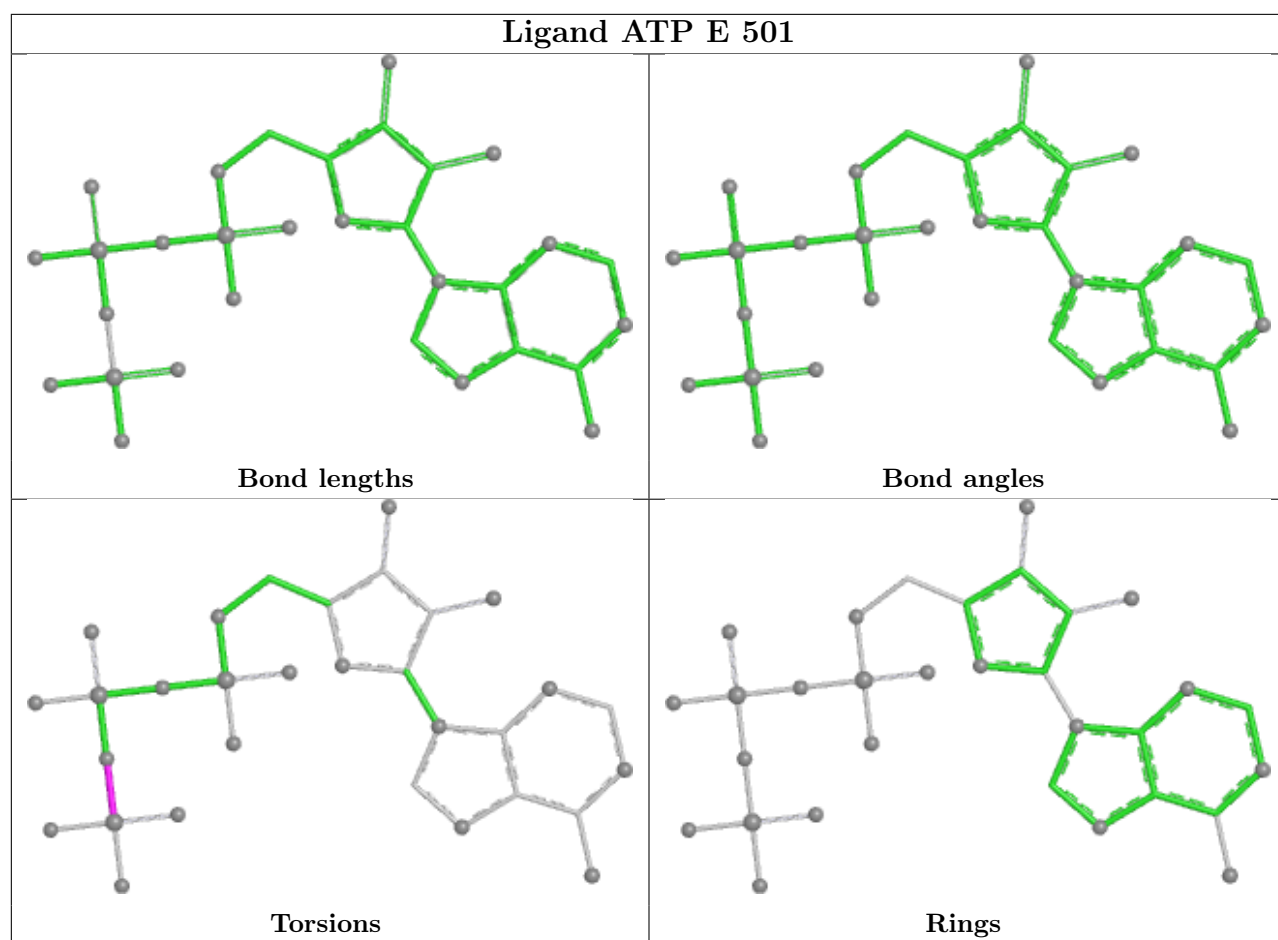
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
35	A	501	ATP	2	0
36	B	501	ADP	1	0

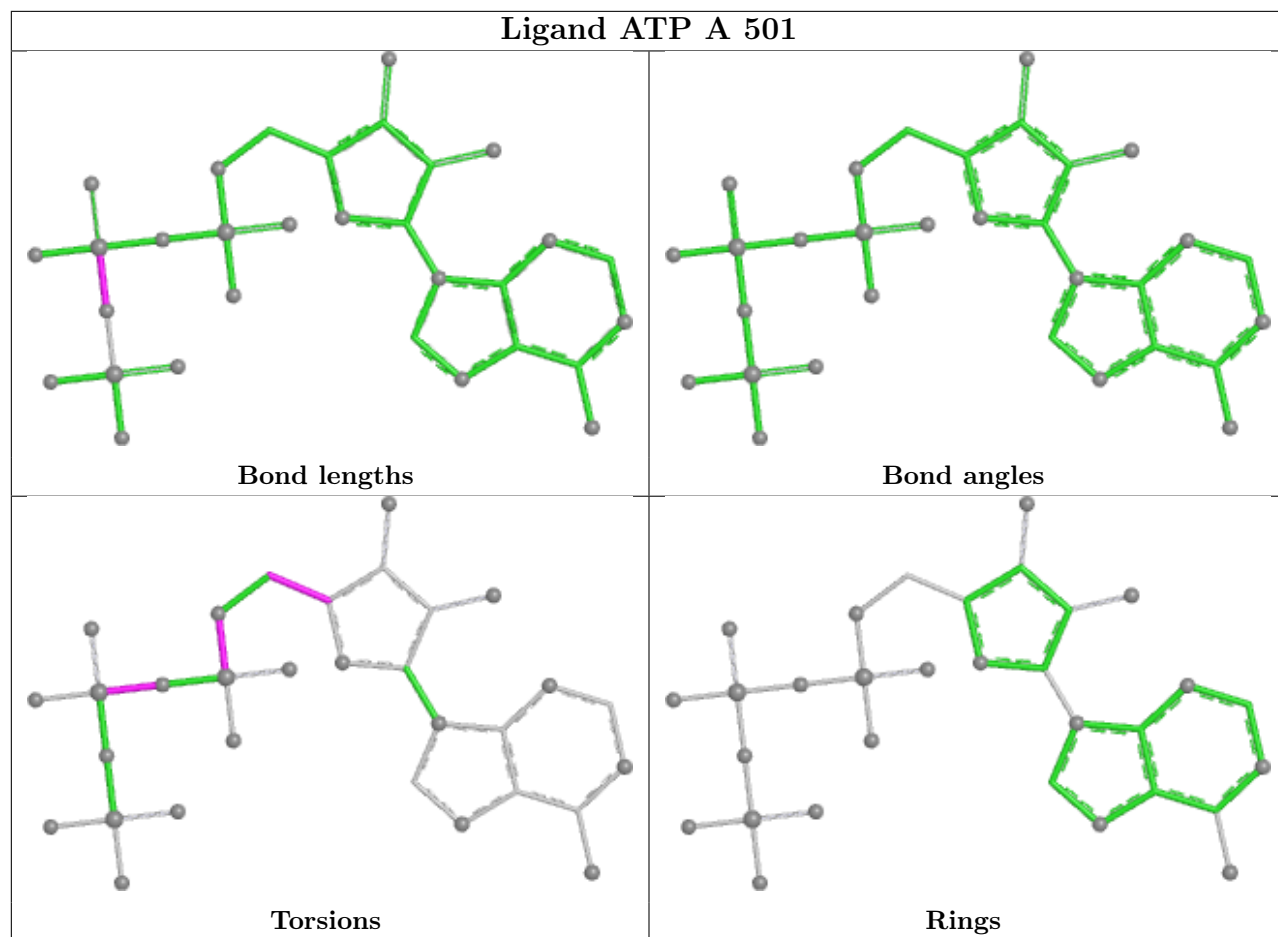
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



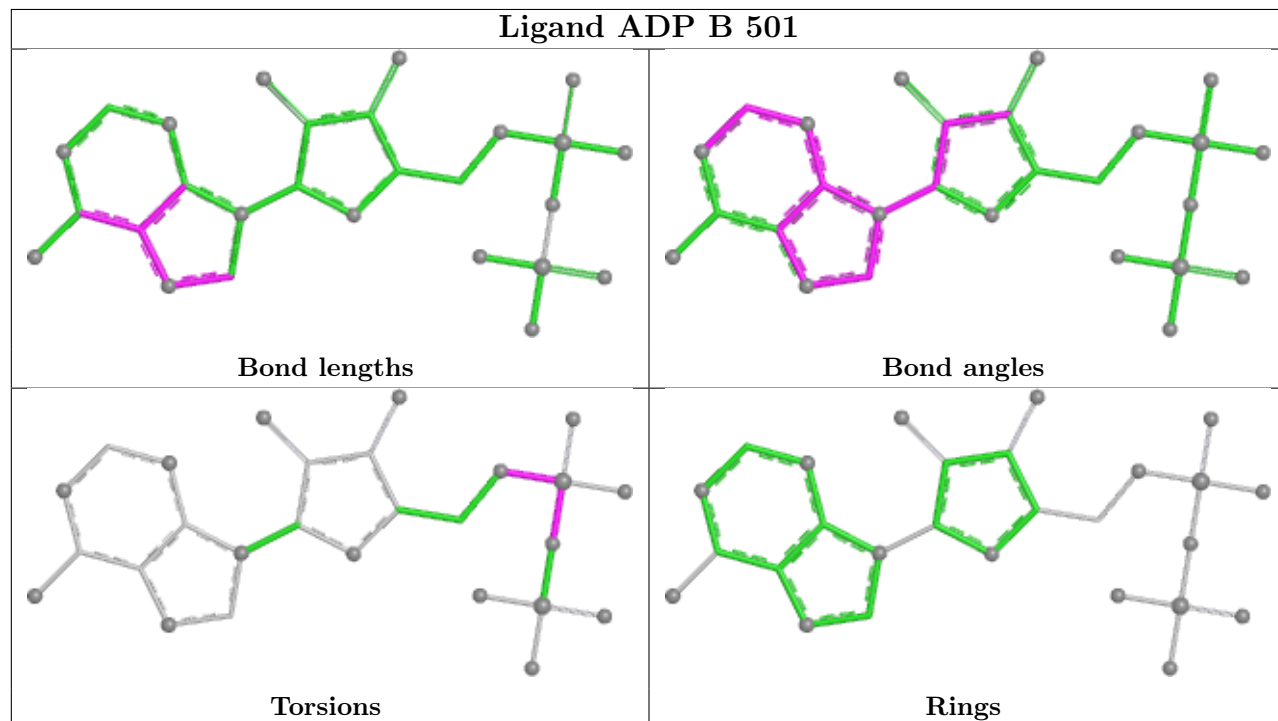




Ligand ATP A 501



Ligand ADP B 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.

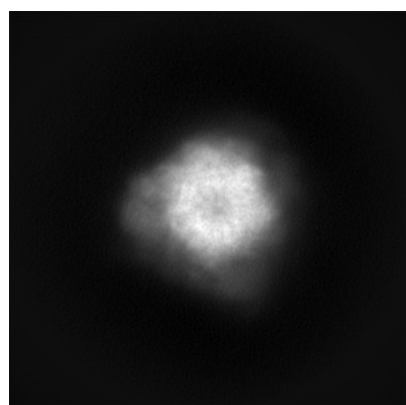
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32281. 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](#)

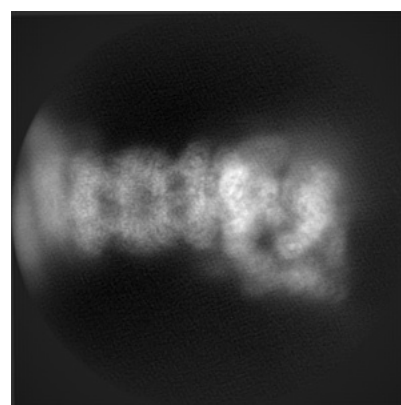
6.1.1 Primary map



X



Y

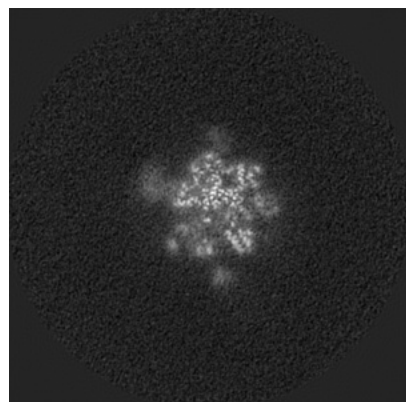


Z

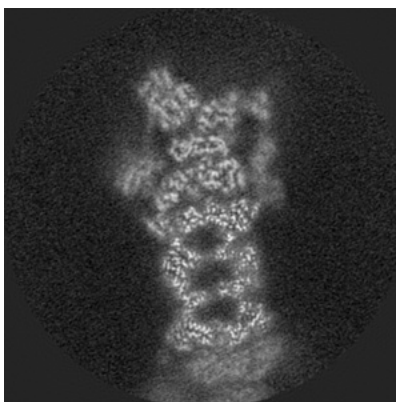
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

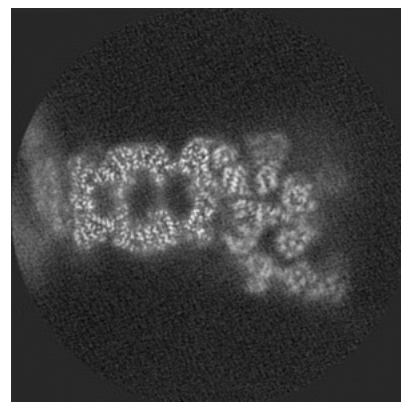
6.2.1 Primary map



X Index: 320



Y Index: 320

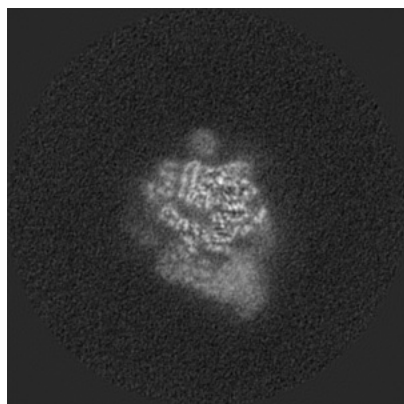


Z Index: 320

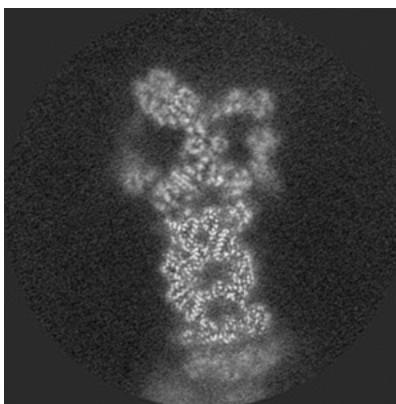
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

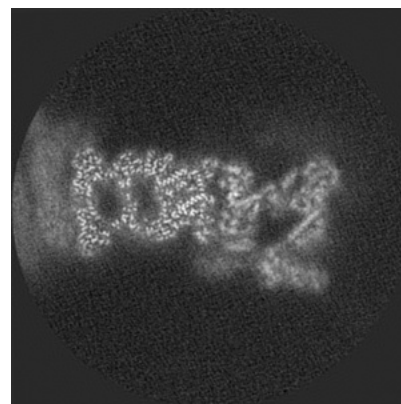
6.3.1 Primary map



X Index: 359



Y Index: 302

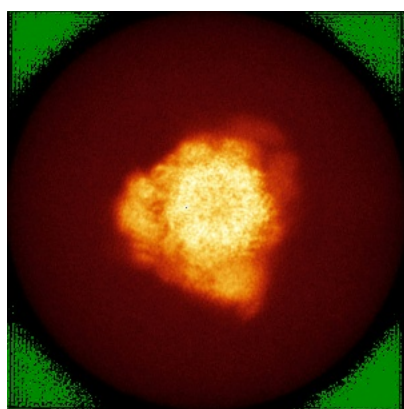


Z Index: 360

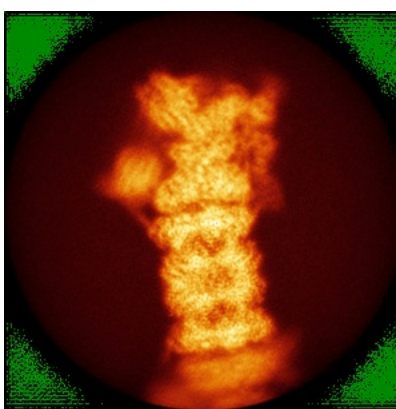
The images above show the largest variance slices of the map in three orthogonal directions.

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

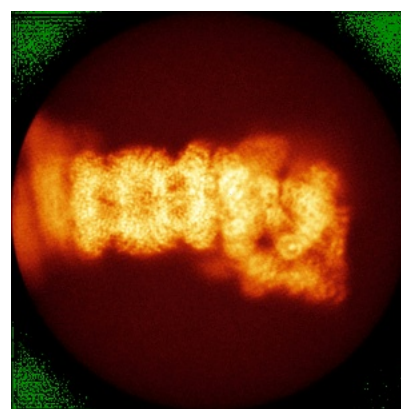
6.4.1 Primary map



X



Y

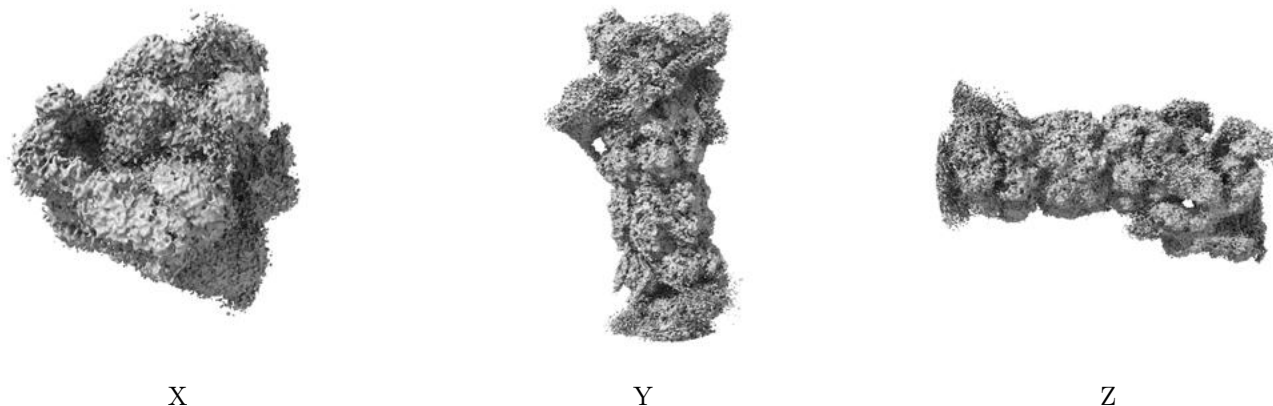


Z

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

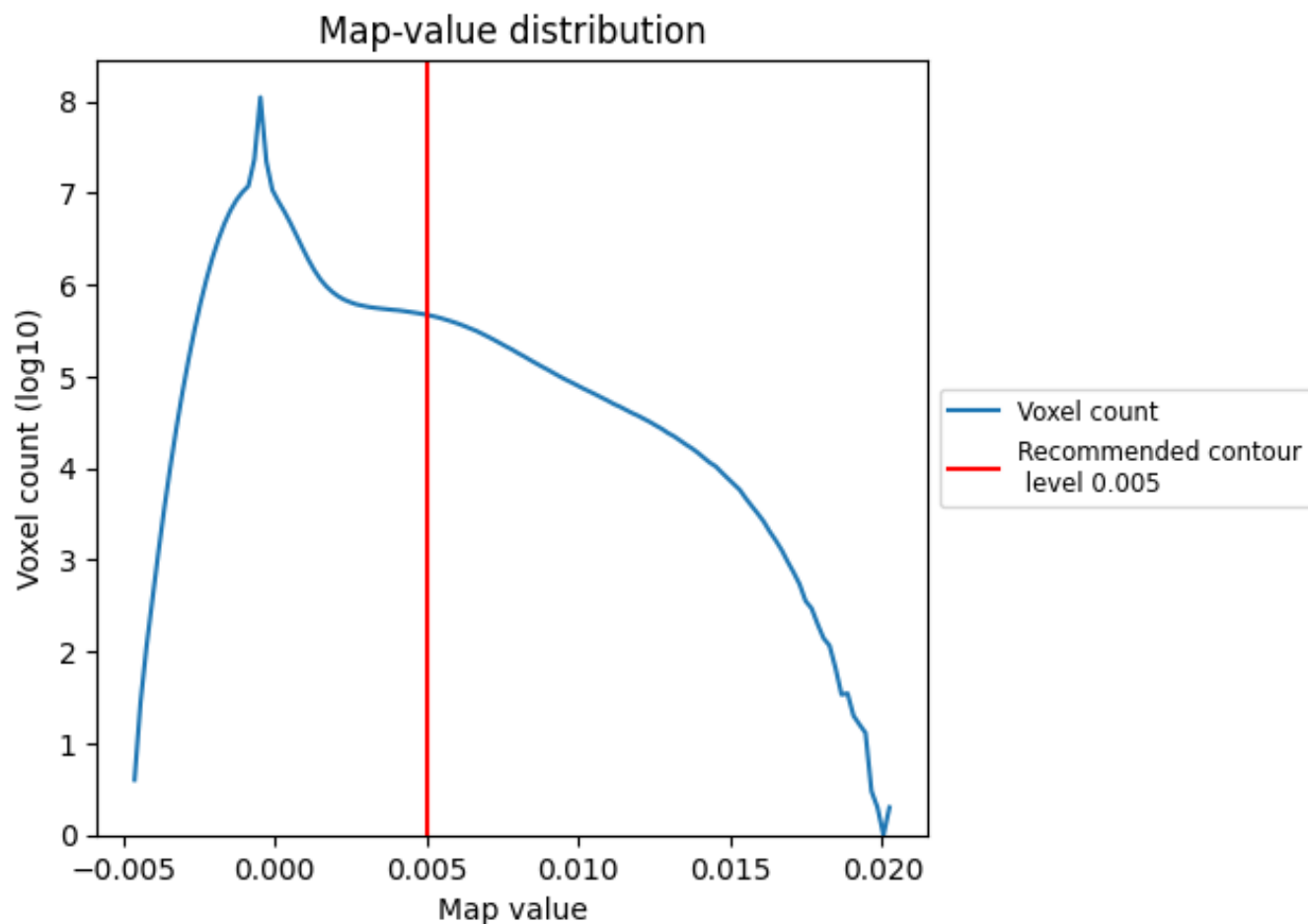
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

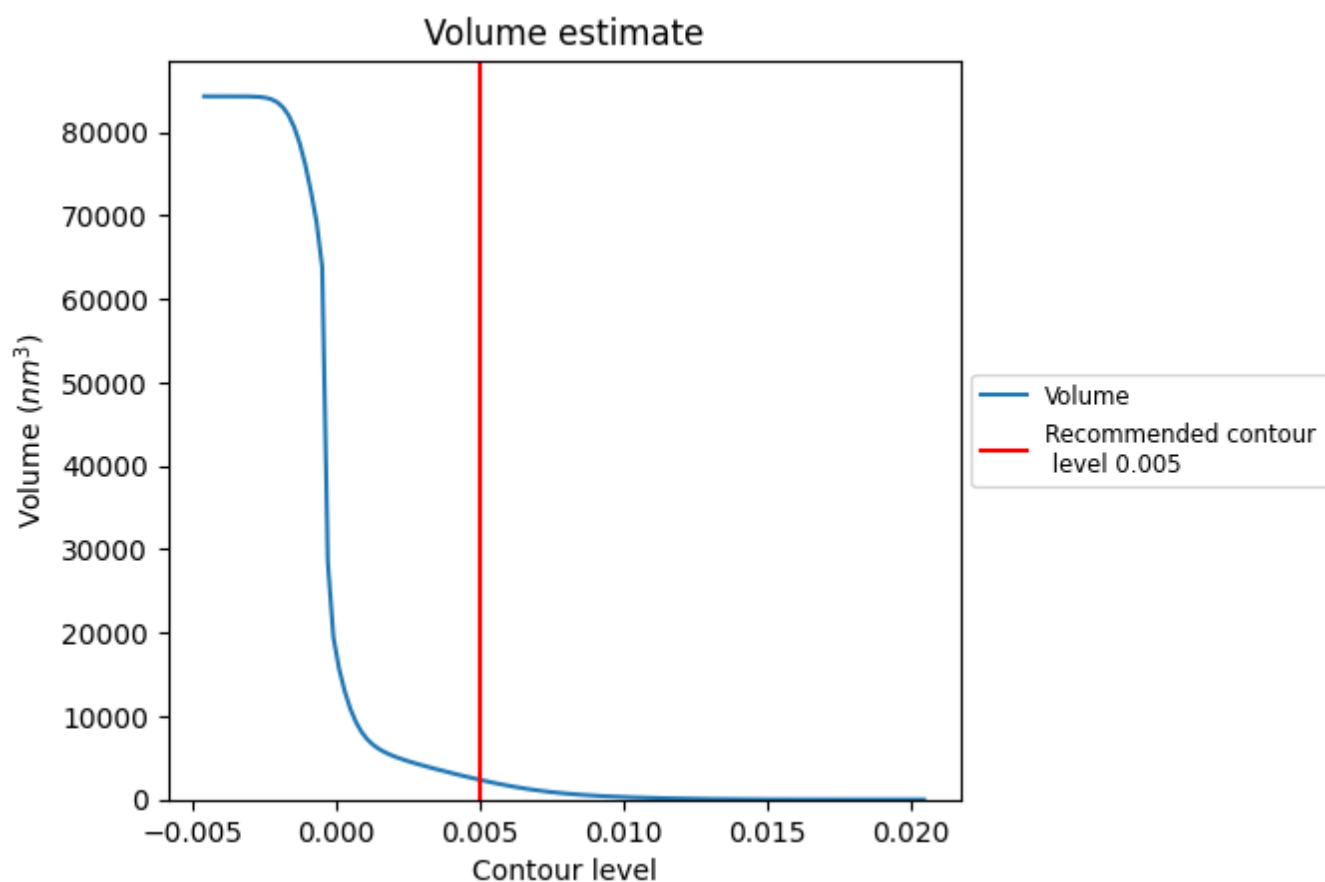
This section contains the results of statistical analysis of the map.

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.

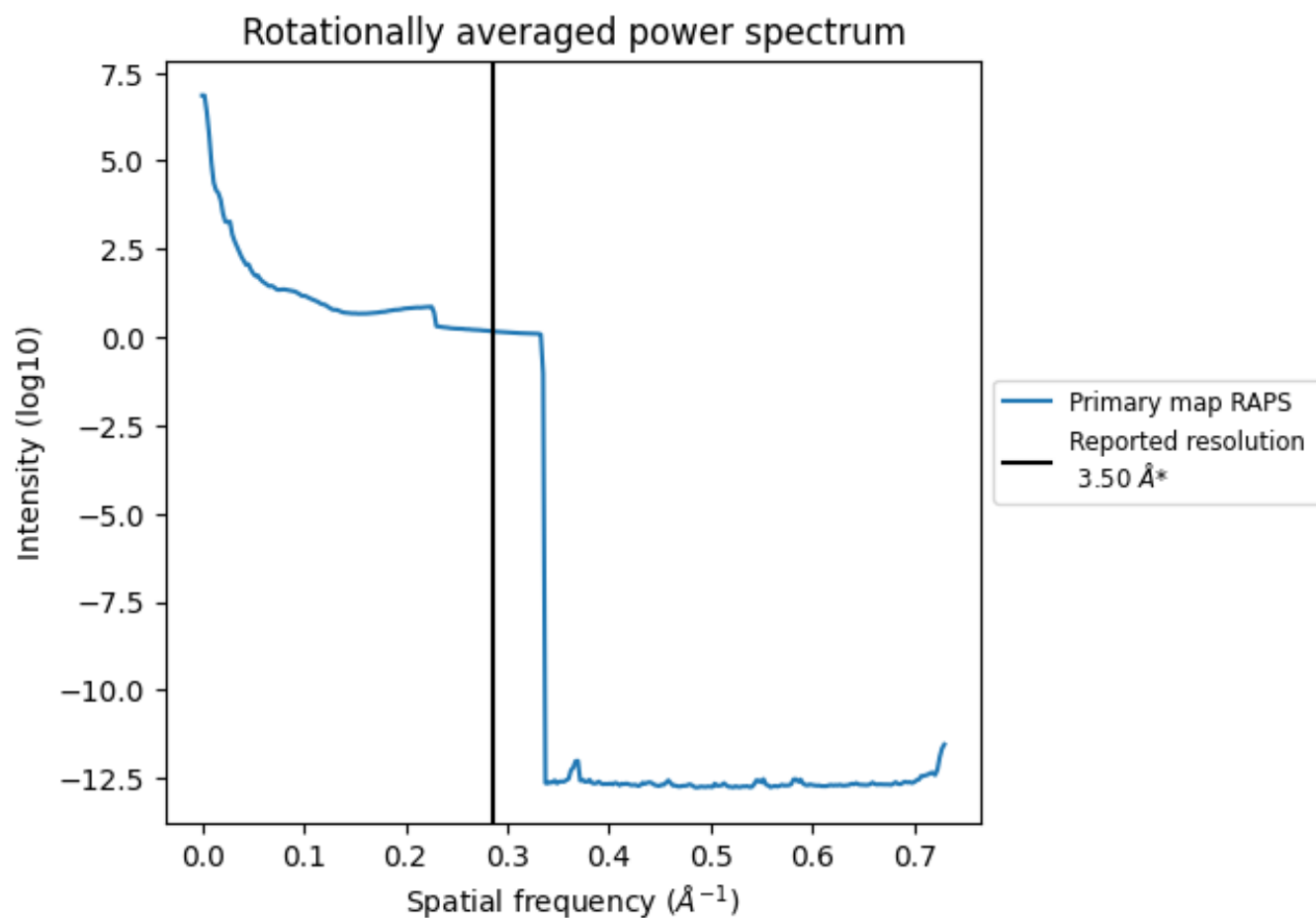
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2347 nm³; this corresponds to an approximate mass of 2121 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 0.286 $\text{\AA}^{-1}$

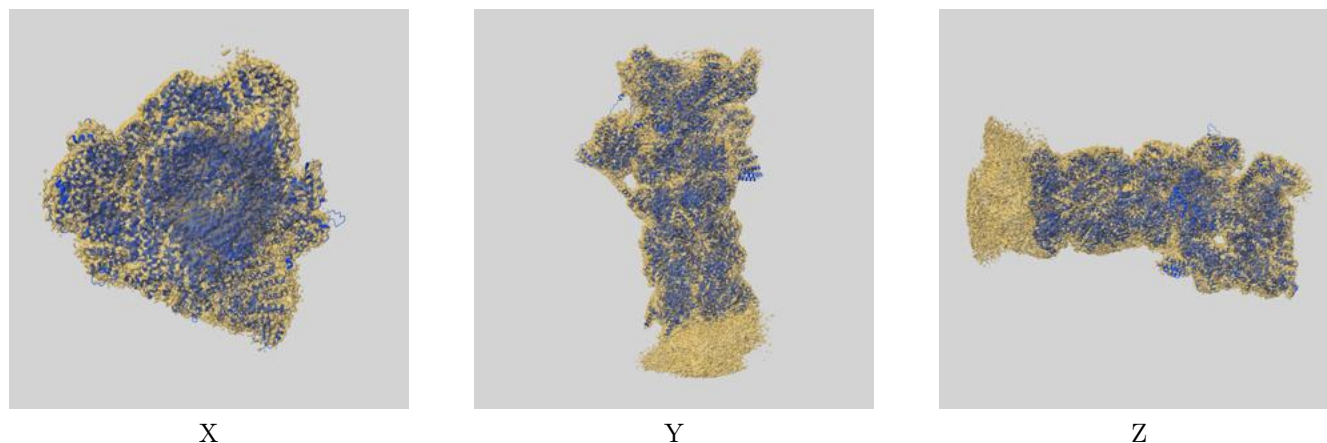
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

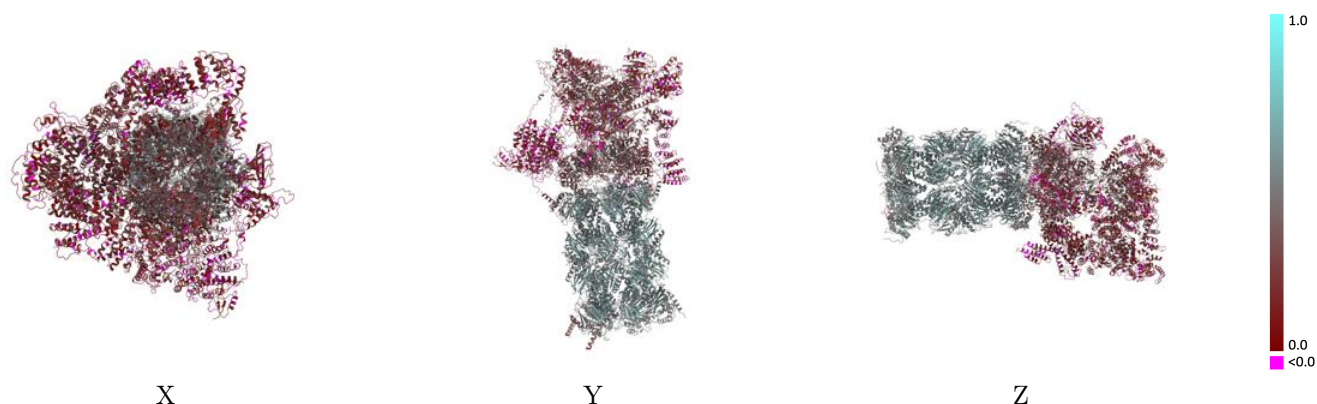
This section contains information regarding the fit between EMDB map EMD-32281 and PDB model 7W3I. Per-residue inclusion information can be found in section [3](#) on page [12](#).

9.1 Map-model overlay [i](#)



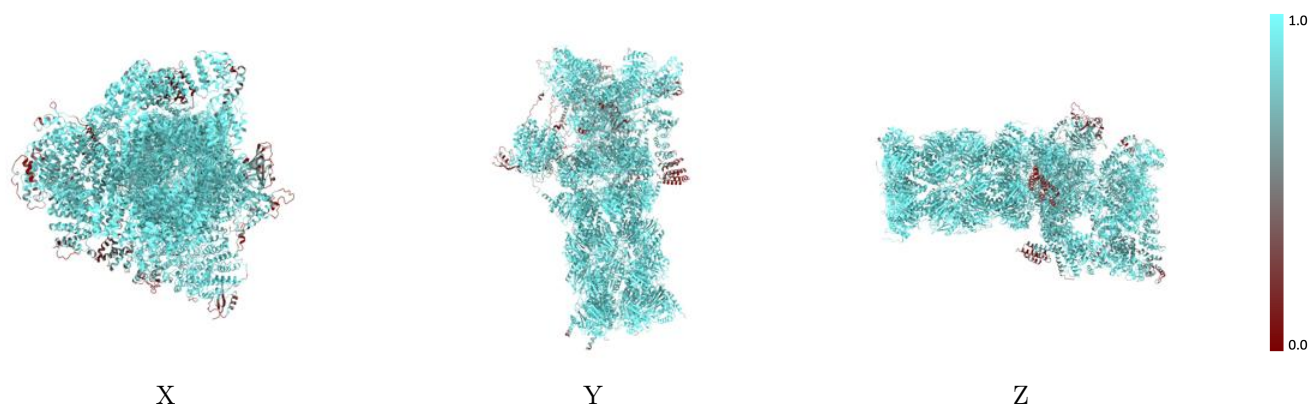
The images above show the 3D surface view of the map at the recommended contour level 0.005 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



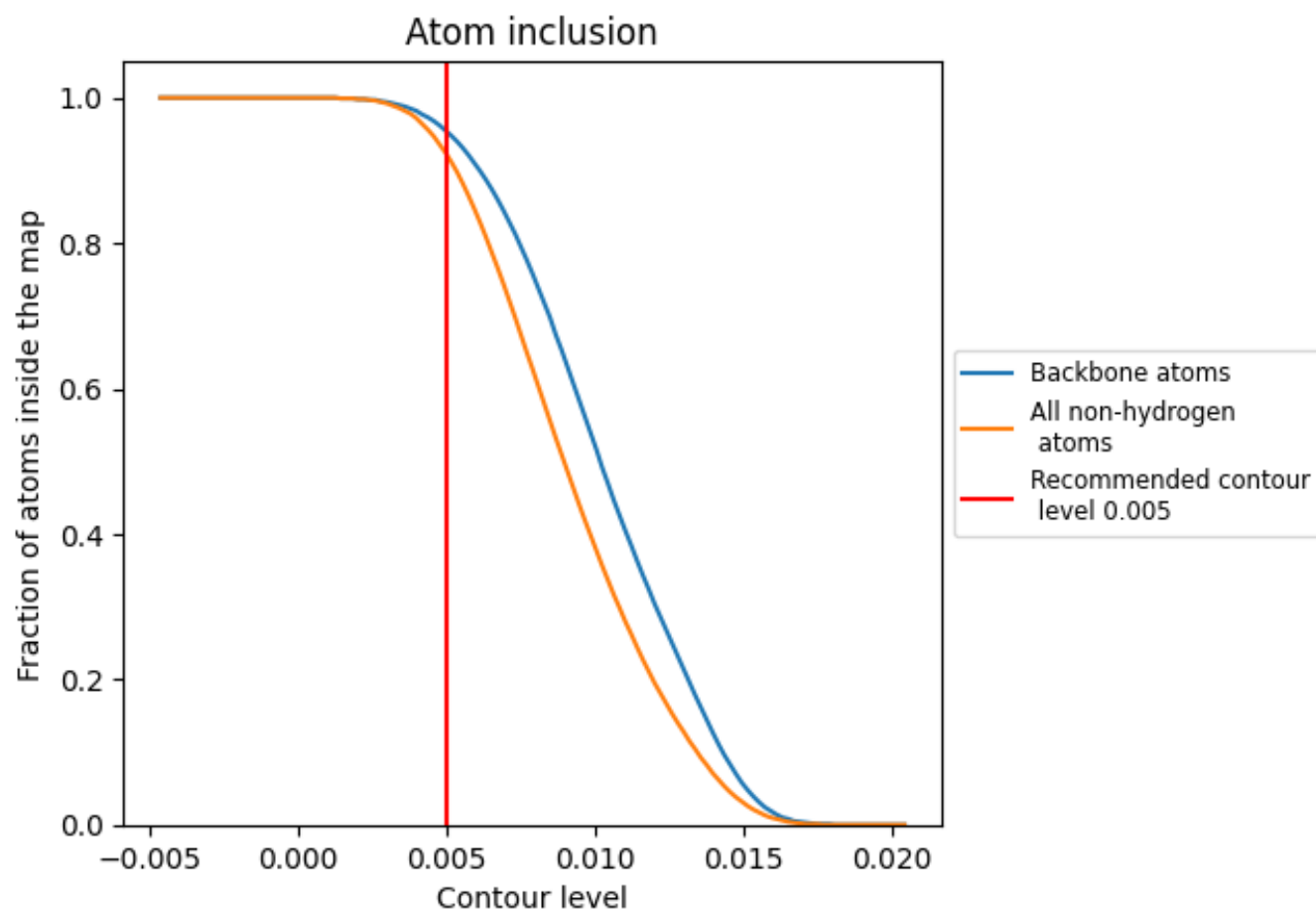
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.005).

9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 92% 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 (0.005) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9230	 0.3580
A	 0.9590	 0.3330
B	 0.9540	 0.2960
C	 0.9510	 0.2920
D	 0.9810	 0.3000
E	 0.9450	 0.3040
F	 0.9370	 0.3270
G	 0.9890	 0.5060
H	 0.9930	 0.5170
I	 0.9810	 0.4850
J	 0.9730	 0.4510
K	 0.9780	 0.4970
L	 0.9910	 0.5310
M	 0.9830	 0.5140
N	 0.9910	 0.5430
O	 0.9960	 0.5360
P	 0.9970	 0.5420
Q	 0.9920	 0.5350
R	 0.9900	 0.5390
S	 0.9840	 0.5330
T	 0.9840	 0.5470
U	 0.8980	 0.2470
V	 0.8670	 0.2150
W	 0.6990	 0.1590
X	 0.8180	 0.2020
Y	 0.9320	 0.1970
Z	 0.9730	 0.2790
a	 0.9130	 0.1860
b	 0.9240	 0.2210
c	 0.9630	 0.2800
d	 0.7080	 0.1640
e	 0.8970	 0.2030
f	 0.8490	 0.1320
g	 0.9710	 0.5090
h	 0.9850	 0.5240



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Chain	Atom inclusion	Q-score
i	 0.9540	 0.4950
j	 0.9500	 0.4460
k	 0.9790	 0.5130
l	 0.9910	 0.5330
m	 0.9820	 0.5140
n	 0.9930	 0.5450
o	 0.9820	 0.5410
p	 0.9870	 0.5420
q	 0.9880	 0.5400
r	 0.9920	 0.5370
s	 0.9930	 0.5380
t	 0.9920	 0.5490
x	 0.6640	 0.1620
y	 0.4920	 0.1750