



wwPDB EM Validation Summary Report ⓘ

Mar 26, 2026 – 06:14 PM UTC

PDB ID : 7NVN / pdb_00007nvn
EMDB ID : EMD-12607
Title : Human TRiC complex in closed state with nanobody and tubulin bound
Authors : Kelly, J.J.; Chi, G.; Bulawa, C.; Paavilainen, V.O.; Bountra, C.; Huiskonen, J.T.; Yue, W.
Deposited on : 2021-03-15
Resolution : 3.00 Å(reported)

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

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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

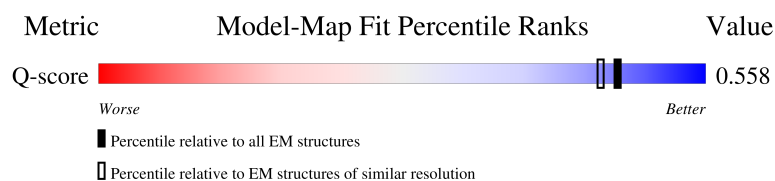
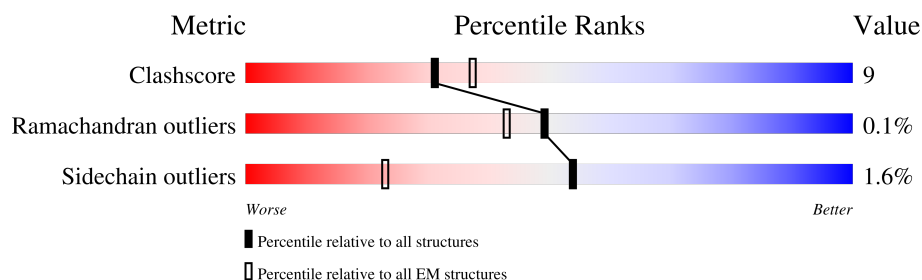
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.00 Å.

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



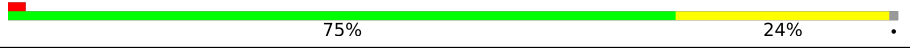
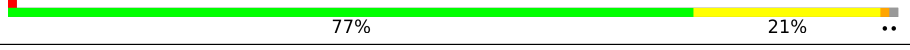
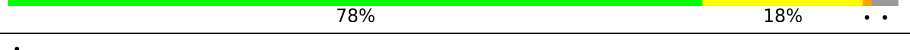
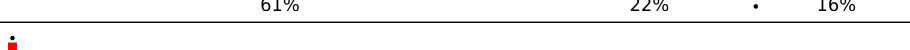
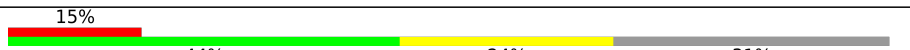
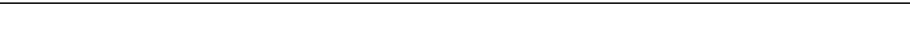
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14081 (2.50 - 3.50)

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

Mol	Chain	Length	Quality of chain
1	A	556	
1	a	556	
2	B	535	
2	b	535	

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Mol	Chain	Length	Quality of chain
3	D	539	
3	d	539	
4	E	541	
4	e	541	
5	G	545	
5	g	545	
6	H	543	
6	h	543	
7	N	129	
7	n	129	
8	Q	548	
8	q	548	
9	Z	531	
9	z	531	
10	T	445	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
13	AF3	B	603	-	-	X	-
13	AF3	G	603	-	-	X	-
13	AF3	H	603	-	-	X	-
13	AF3	Q	603	-	-	X	-
13	AF3	Z	603	-	-	X	-
13	AF3	a	603	-	-	X	-
13	AF3	b	603	-	-	X	-
13	AF3	e	603	-	-	X	-
13	AF3	q	603	-	-	X	-
13	AF3	z	603	-	-	X	-

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 69139 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 T-complex protein 1 subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	535	Total	C	N	O	S	0	0
			4063	2545	710	785	23		
1	a	536	Total	C	N	O	S	0	0
			4071	2550	711	786	24		

- Molecule 2 is a protein called T-complex protein 1 subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	526	Total	C	N	O	S	0	0
			3952	2473	696	764	19		
2	b	527	Total	C	N	O	S	0	0
			3960	2478	697	765	20		

- Molecule 3 is a protein called T-complex protein 1 subunit delta.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	519	Total	C	N	O	S	0	0
			3919	2451	682	763	23		
3	d	519	Total	C	N	O	S	0	0
			3919	2451	682	763	23		

- Molecule 4 is a protein called T-complex protein 1 subunit epsilon.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	537	Total	C	N	O	S	1	0
			4142	2596	720	795	31		
4	e	538	Total	C	N	O	S	1	0
			4148	2599	721	797	31		

- Molecule 5 is a protein called T-complex protein 1 subunit gamma.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	G	523	Total	C	N	O	S	0	0
			4064	2534	718	782	30		
5	g	523	Total	C	N	O	S	0	0
			4064	2534	718	782	30		

- Molecule 6 is a protein called T-complex protein 1 subunit eta.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	H	524	Total	C	N	O	S	0	0
			4026	2545	695	761	25		
6	h	524	Total	C	N	O	S	0	0
			4026	2545	695	761	25		

- Molecule 7 is a protein called Nanobody.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	N	109	Total	C	N	O	S	0	0
			859	535	158	162	4		
7	n	109	Total	C	N	O	S	0	0
			854	532	157	161	4		

- Molecule 8 is a protein called T-complex protein 1 subunit theta.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	Q	530	Total	C	N	O	S	0	0
			4033	2545	686	775	27		
8	q	529	Total	C	N	O	S	0	0
			4029	2543	685	774	27		

- Molecule 9 is a protein called T-complex protein 1 subunit zeta.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	Z	529	Total	C	N	O	S	0	0
			4051	2546	709	775	21		
9	z	526	Total	C	N	O	S	0	0
			4028	2531	705	771	21		

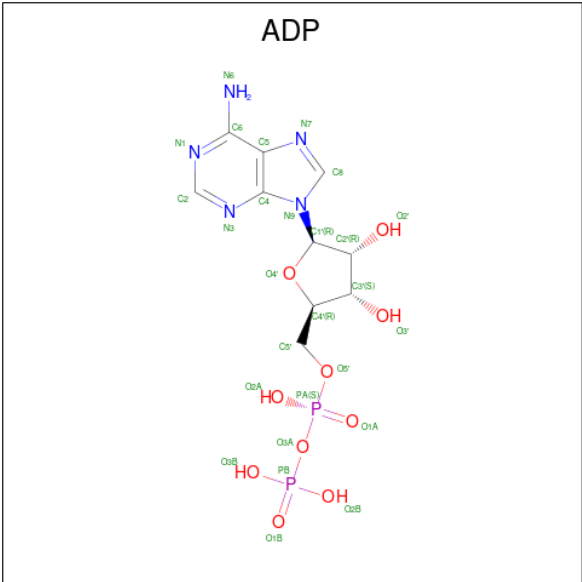
- Molecule 10 is a protein called Tubulin beta-2A chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	T	307	Total	C	N	O	S	0	0
			2400	1507	402	475	16		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
T	170	VAL	MET	conflict	UNP Q13885

- Molecule 11 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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Mol	Chain	Residues	Atoms					AltConf
11	d	1	Total	C	N	O	P	0
			27	10	5	10	2	
11	e	1	Total	C	N	O	P	0
			27	10	5	10	2	
11	g	1	Total	C	N	O	P	0
			27	10	5	10	2	
11	h	1	Total	C	N	O	P	0
			27	10	5	10	2	
11	q	1	Total	C	N	O	P	0
			27	10	5	10	2	
11	z	1	Total	C	N	O	P	0
			27	10	5	10	2	

- Molecule 12 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

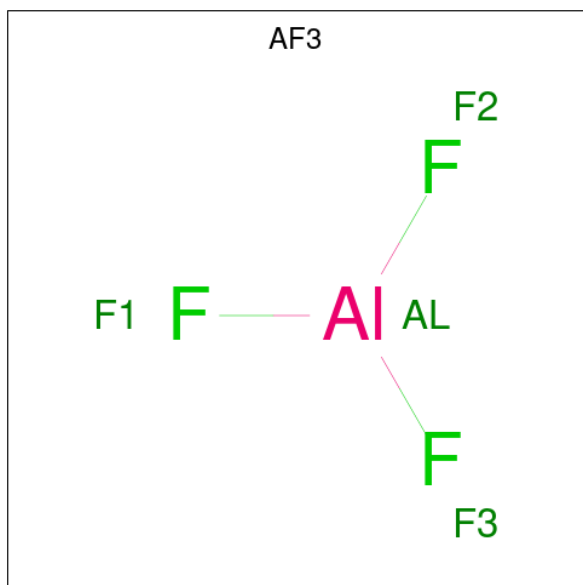
Mol	Chain	Residues	Atoms		AltConf
12	A	1	Total	Mg	0
			1	1	
12	B	1	Total	Mg	0
			1	1	
12	D	1	Total	Mg	0
			1	1	
12	E	1	Total	Mg	0
			1	1	
12	G	1	Total	Mg	0
			1	1	
12	H	1	Total	Mg	0
			1	1	
12	Q	1	Total	Mg	0
			1	1	
12	Z	1	Total	Mg	0
			1	1	
12	a	1	Total	Mg	0
			1	1	
12	b	1	Total	Mg	0
			1	1	
12	d	1	Total	Mg	0
			1	1	
12	e	1	Total	Mg	0
			1	1	
12	g	1	Total	Mg	0
			1	1	

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Mol	Chain	Residues	Atoms		AltConf
12	h	1	Total 1	Mg 1	0
12	q	1	Total 1	Mg 1	0
12	z	1	Total 1	Mg 1	0

- Molecule 13 is ALUMINUM FLUORIDE (CCD ID: AF3) (formula: AlF_3).



Mol	Chain	Residues	Atoms			AltConf
13	A	1	Total 4	Al 1	F 3	0
13	B	1	Total 4	Al 1	F 3	0
13	D	1	Total 4	Al 1	F 3	0
13	E	1	Total 4	Al 1	F 3	0
13	G	1	Total 4	Al 1	F 3	0
13	H	1	Total 4	Al 1	F 3	0
13	Q	1	Total 4	Al 1	F 3	0
13	Z	1	Total 4	Al 1	F 3	0

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Mol	Chain	Residues	Atoms			AltConf
13	a	1	Total 4	Al 1	F 3	0
13	b	1	Total 4	Al 1	F 3	0
13	d	1	Total 4	Al 1	F 3	0
13	e	1	Total 4	Al 1	F 3	0
13	g	1	Total 4	Al 1	F 3	0
13	h	1	Total 4	Al 1	F 3	0
13	q	1	Total 4	Al 1	F 3	0
13	z	1	Total 4	Al 1	F 3	0

- Molecule 14 is water.

Mol	Chain	Residues	Atoms		AltConf
14	A	2	Total 2	O 2	0
14	B	2	Total 2	O 2	0
14	D	1	Total 1	O 1	0
14	E	1	Total 1	O 1	0
14	G	1	Total 1	O 1	0
14	H	1	Total 1	O 1	0
14	Q	1	Total 1	O 1	0
14	Z	1	Total 1	O 1	0
14	a	2	Total 2	O 2	0
14	b	1	Total 1	O 1	0
14	d	1	Total 1	O 1	0

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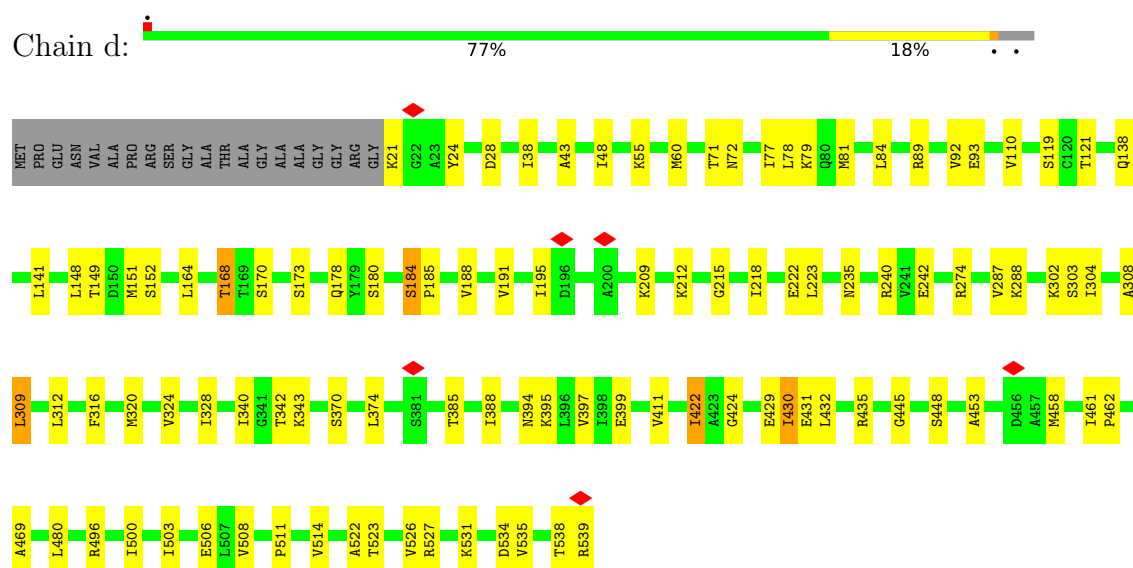
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Mol	Chain	Residues	Atoms		AltConf
14	e	1	Total 1	O 1	0
14	g	1	Total 1	O 1	0
14	h	1	Total 1	O 1	0
14	q	1	Total 1	O 1	0
14	z	1	Total 1	O 1	0

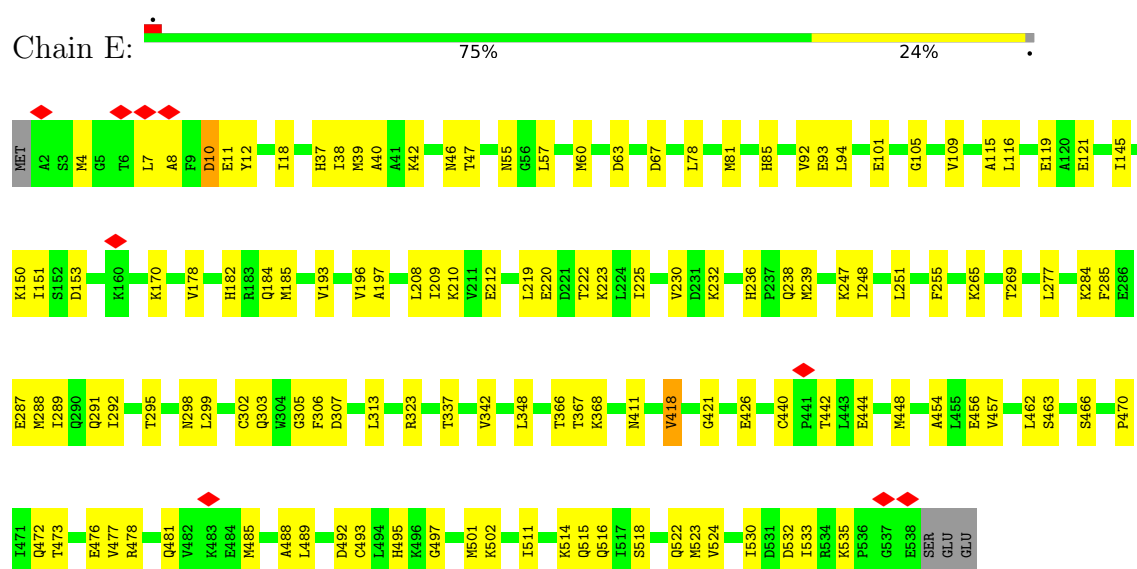




• Molecule 3: T-complex protein 1 subunit delta



• Molecule 4: T-complex protein 1 subunit epsilon



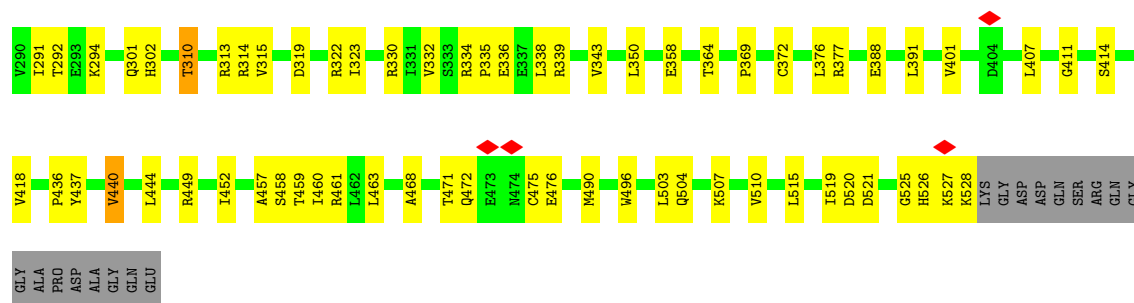
[illegible][illegible]

Chain g:

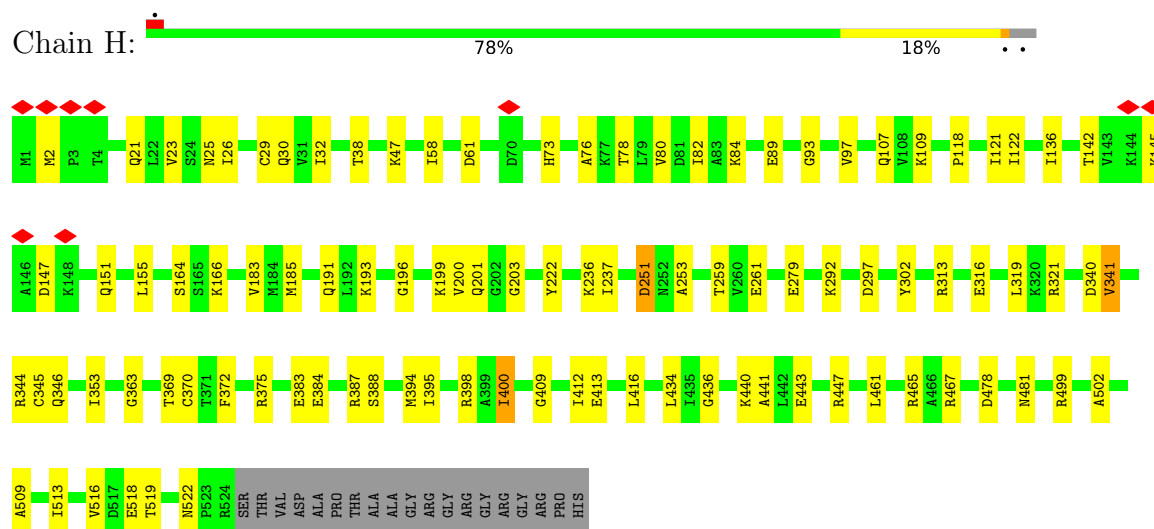
73% 22%

0.15
0.10
0.05
0.00

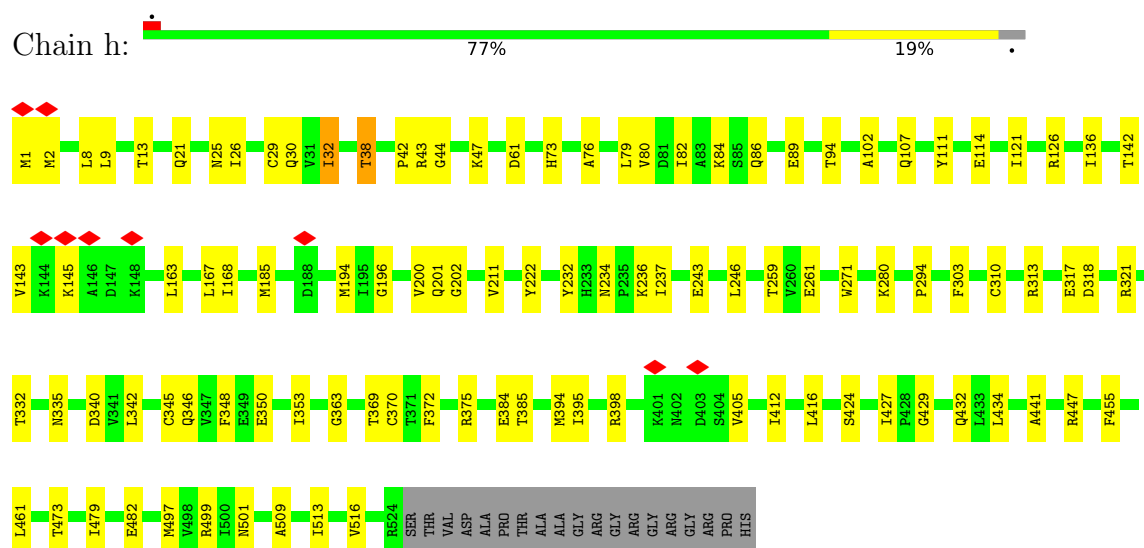
MET MET GLY HIS ARG P6 V7 Q12 M13 D35 R38 L41 G42 P43 K44 S45 K48 M54 I66 Q73 P75 S79 T83 S84 E89 E90 T96 E104 M105 V108 P119 Y126 R127 L130 D131 M132 M133 L137 D145



- Molecule 6: T-complex protein 1 subunit eta

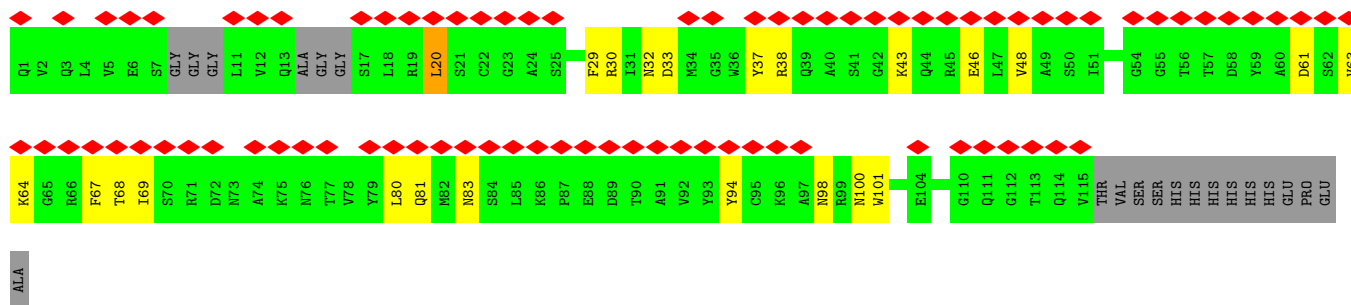


- Molecule 6: T-complex protein 1 subunit eta



- Molecule 7: Nanobody

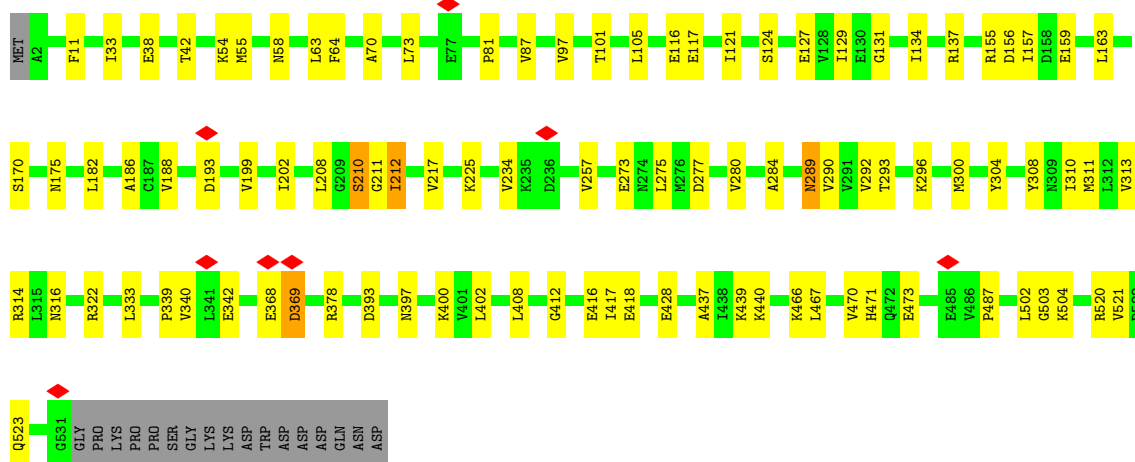
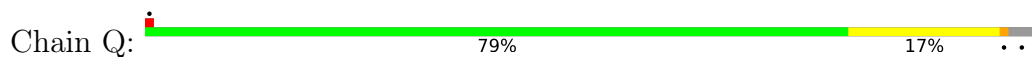




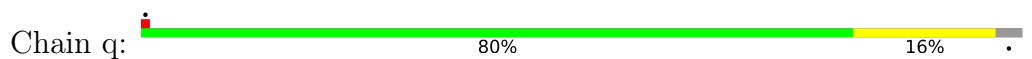
• Molecule 7: Nanobody

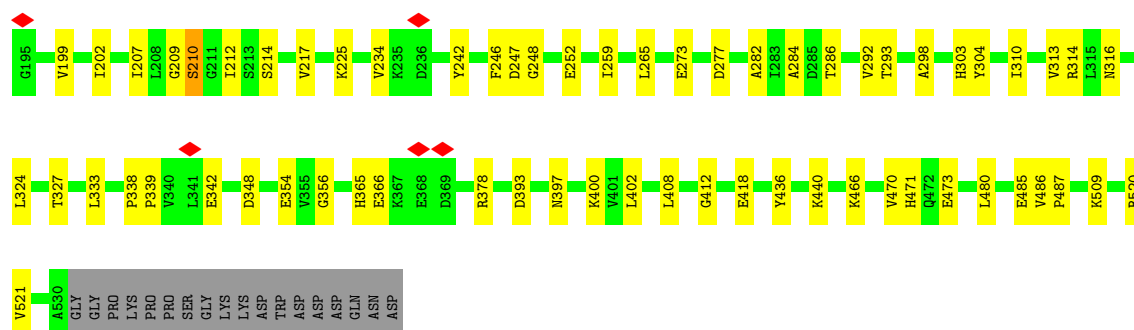


• Molecule 8: T-complex protein 1 subunit theta



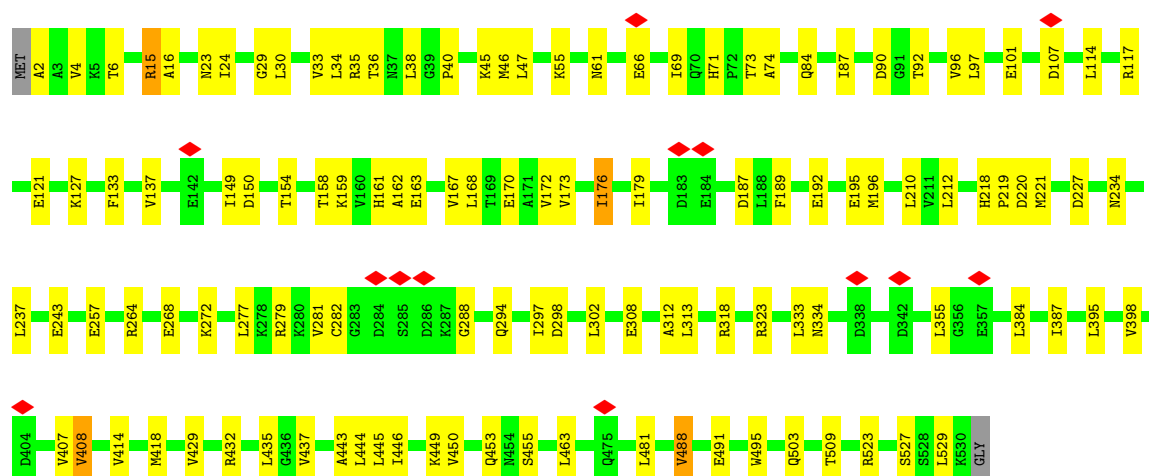
• Molecule 8: T-complex protein 1 subunit theta





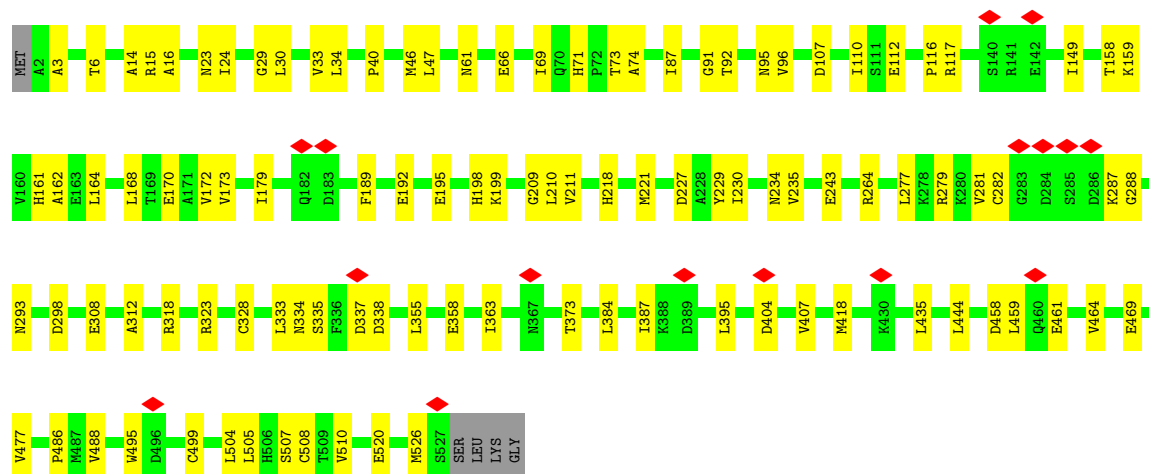
• Molecule 9: T-complex protein 1 subunit zeta

Chain Z: 77% 22%



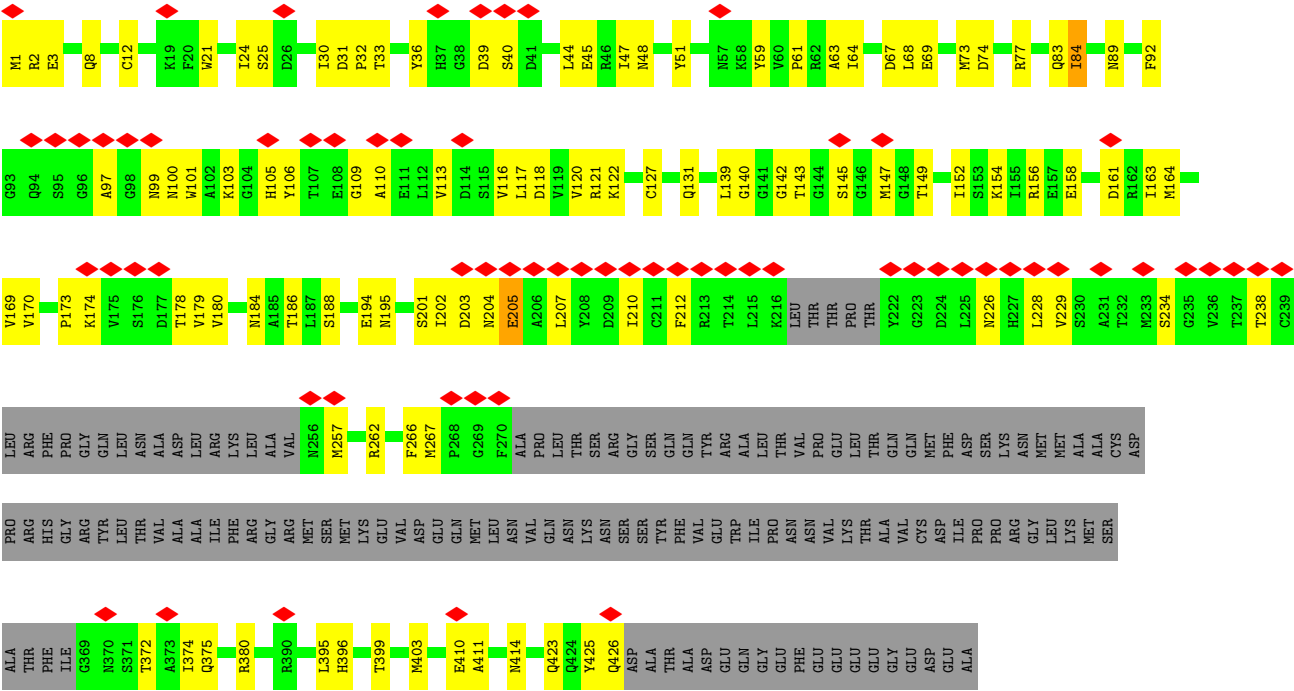
• Molecule 9: T-complex protein 1 subunit zeta

Chain z: 79% 20%



• Molecule 10: Tubulin beta-2A chain

Chain T: 15% 44% 24% 31%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	93758	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	43	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.187	Depositor
Minimum map value	-0.089	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.025	Depositor
Map size (Å)	423.99997, 423.99997, 423.99997	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.06, 1.06, 1.06	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AF3, MG, ADP

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	A	0.17	0/4103	0.36	0/5540
1	a	0.17	0/4111	0.36	0/5550
2	B	0.16	0/3995	0.35	0/5386
2	b	0.16	0/4003	0.35	0/5396
3	D	0.16	0/3951	0.33	0/5333
3	d	0.16	0/3951	0.32	0/5333
4	E	0.19	0/4196	0.38	0/5652
4	e	0.17	0/4202	0.36	0/5660
5	G	0.17	0/4110	0.35	0/5544
5	g	0.16	0/4110	0.35	0/5544
6	H	0.17	0/4083	0.33	0/5511
6	h	0.18	0/4083	0.34	0/5511
7	N	0.17	0/874	0.42	0/1179
7	n	0.13	0/869	0.43	0/1172
8	Q	0.17	0/4091	0.35	0/5530
8	q	0.19	0/4087	0.35	0/5525
9	Z	0.18	0/4098	0.36	0/5524
9	z	0.17	0/4075	0.35	0/5494
10	T	0.24	0/2453	0.53	0/3320
All	All	0.17	0/69445	0.36	0/93704

There are no bond length outliers.

There are no bond angle outliers.

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	A	4063	0	4219	99	0
1	a	4071	0	4231	86	0
2	B	3952	0	4070	84	0
2	b	3960	0	4082	76	0
3	D	3919	0	4128	84	0
3	d	3919	0	4128	71	0
4	E	4142	0	4256	93	0
4	e	4148	0	4261	82	0
5	G	4064	0	4202	87	0
5	g	4064	0	4202	90	0
6	H	4026	0	4135	70	0
6	h	4026	0	4135	67	0
7	N	859	0	817	17	0
7	n	854	0	812	21	0
8	Q	4033	0	4102	63	0
8	q	4029	0	4099	62	0
9	Z	4051	0	4195	76	0
9	z	4028	0	4166	68	0
10	T	2400	0	2245	77	0
11	A	27	0	12	1	0
11	B	27	0	12	5	0
11	D	27	0	12	0	0
11	E	27	0	12	0	0
11	G	27	0	12	2	0
11	H	27	0	12	4	0
11	Q	27	0	12	2	0
11	Z	27	0	12	1	0
11	a	27	0	12	3	0
11	b	27	0	12	3	0
11	d	27	0	12	0	0
11	e	27	0	12	1	0
11	g	27	0	12	3	0
11	h	27	0	12	1	0
11	q	27	0	12	2	0
11	z	27	0	12	2	0
12	A	1	0	0	0	0
12	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	D	1	0	0	0	0
12	E	1	0	0	0	0
12	G	1	0	0	0	0
12	H	1	0	0	0	0
12	Q	1	0	0	0	0
12	Z	1	0	0	0	0
12	a	1	0	0	0	0
12	b	1	0	0	0	0
12	d	1	0	0	0	0
12	e	1	0	0	0	0
12	g	1	0	0	0	0
12	h	1	0	0	0	0
12	q	1	0	0	0	0
12	z	1	0	0	0	0
13	A	4	0	0	1	0
13	B	4	0	0	2	0
13	D	4	0	0	0	0
13	E	4	0	0	0	0
13	G	4	0	0	2	0
13	H	4	0	0	2	0
13	Q	4	0	0	2	0
13	Z	4	0	0	2	0
13	a	4	0	0	3	0
13	b	4	0	0	4	0
13	d	4	0	0	0	0
13	e	4	0	0	2	0
13	g	4	0	0	0	0
13	h	4	0	0	1	0
13	q	4	0	0	2	0
13	z	4	0	0	3	0
14	A	2	0	0	0	0
14	B	2	0	0	0	0
14	D	1	0	0	0	0
14	E	1	0	0	0	0
14	G	1	0	0	0	0
14	H	1	0	0	1	0
14	Q	1	0	0	0	0
14	Z	1	0	0	0	0
14	a	2	0	0	0	0
14	b	1	0	0	0	0
14	d	1	0	0	0	0
14	e	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	g	1	0	0	0	0
14	h	1	0	0	1	0
14	q	1	0	0	0	0
14	z	1	0	0	0	0
All	All	69139	0	70677	1230	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 1230 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)
10:T:203:ASP:O	10:T:207:LEU:HG	1.45	1.13
10:T:202:ILE:HG22	10:T:207:LEU:HD11	1.41	0.99
10:T:161:ASP:OD1	10:T:161:ASP:O	1.91	0.89
6:H:340:ASP:OD1	6:H:341:VAL:HG13	1.79	0.82
9:Z:154:THR:HG21	9:Z:495:TRP:H	1.44	0.82

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	A	533/556 (96%)	512 (96%)	20 (4%)	1 (0%)	43	76
1	a	534/556 (96%)	514 (96%)	19 (4%)	1 (0%)	43	76
2	B	524/535 (98%)	504 (96%)	20 (4%)	0	100	100
2	b	525/535 (98%)	506 (96%)	19 (4%)	0	100	100
3	D	517/539 (96%)	505 (98%)	12 (2%)	0	100	100
3	d	517/539 (96%)	505 (98%)	12 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	E	536/541 (99%)	518 (97%)	17 (3%)	1 (0%)	43	76
4	e	537/541 (99%)	517 (96%)	19 (4%)	1 (0%)	43	76
5	G	521/545 (96%)	501 (96%)	20 (4%)	0	100	100
5	g	521/545 (96%)	500 (96%)	21 (4%)	0	100	100
6	H	522/543 (96%)	510 (98%)	12 (2%)	0	100	100
6	h	522/543 (96%)	506 (97%)	16 (3%)	0	100	100
7	N	103/129 (80%)	100 (97%)	2 (2%)	1 (1%)	12	45
7	n	103/129 (80%)	101 (98%)	1 (1%)	1 (1%)	12	45
8	Q	528/548 (96%)	511 (97%)	15 (3%)	2 (0%)	30	65
8	q	527/548 (96%)	513 (97%)	13 (2%)	1 (0%)	43	76
9	Z	527/531 (99%)	510 (97%)	16 (3%)	1 (0%)	43	76
9	z	524/531 (99%)	507 (97%)	16 (3%)	1 (0%)	43	76
10	T	299/445 (67%)	282 (94%)	17 (6%)	0	100	100
All	All	8920/9379 (95%)	8622 (97%)	287 (3%)	11 (0%)	49	80

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	CYS
4	E	10	ASP
8	Q	210	SER
9	Z	16	ALA
1	a	357	CYS

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	A	446/463 (96%)	438 (98%)	8 (2%)	51	77
1	a	447/463 (96%)	442 (99%)	5 (1%)	65	83
2	B	418/427 (98%)	409 (98%)	9 (2%)	45	74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	b	419/427 (98%)	410 (98%)	9 (2%)	47	75
3	D	442/452 (98%)	432 (98%)	10 (2%)	44	74
3	d	442/452 (98%)	432 (98%)	10 (2%)	44	74
4	E	453/456 (99%)	450 (99%)	3 (1%)	76	86
4	e	454/456 (100%)	448 (99%)	6 (1%)	61	81
5	G	454/469 (97%)	449 (99%)	5 (1%)	65	83
5	g	454/469 (97%)	447 (98%)	7 (2%)	57	80
6	H	431/443 (97%)	426 (99%)	5 (1%)	63	82
6	h	431/443 (97%)	424 (98%)	7 (2%)	55	79
7	N	88/105 (84%)	86 (98%)	2 (2%)	44	74
7	n	87/105 (83%)	85 (98%)	2 (2%)	44	74
8	Q	436/452 (96%)	429 (98%)	7 (2%)	55	79
8	q	436/452 (96%)	430 (99%)	6 (1%)	59	80
9	Z	441/442 (100%)	431 (98%)	10 (2%)	44	74
9	z	438/442 (99%)	433 (99%)	5 (1%)	65	83
10	T	260/382 (68%)	253 (97%)	7 (3%)	39	71
All	All	7477/7800 (96%)	7354 (98%)	123 (2%)	54	79

5 of 123 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
9	Z	408	VAL
8	q	480	LEU
2	b	446	LEU
8	q	234	VAL
10	T	84	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 171 such sidechains are listed below:

Mol	Chain	Res	Type
4	e	465	ASN
8	q	34	GLN
5	g	13	ASN
5	g	472	GLN
8	q	461	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 48 ligands modelled in this entry, 16 are monoatomic - leaving 32 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$
13	AF3	h	603	-	0,3,3	-	-	-		
11	ADP	H	601	12	28,29,29	1.39	5 (17%)	43,45,45	1.82	9 (20%)
11	ADP	Q	601	12	28,29,29	1.36	5 (17%)	43,45,45	1.83	10 (23%)
11	ADP	Z	601	12	28,29,29	1.40	4 (14%)	43,45,45	1.85	9 (20%)
13	AF3	z	603	-	0,3,3	-	-	-		
11	ADP	G	601	12	28,29,29	1.38	5 (17%)	43,45,45	1.80	9 (20%)
13	AF3	g	603	-	0,3,3	-	-	-		
11	ADP	A	601	12	28,29,29	1.41	4 (14%)	43,45,45	1.83	9 (20%)
13	AF3	q	603	-	0,3,3	-	-	-		
11	ADP	d	601	12	28,29,29	1.37	5 (17%)	43,45,45	1.83	9 (20%)
11	ADP	g	601	12	28,29,29	1.38	5 (17%)	43,45,45	1.80	9 (20%)
13	AF3	A	603	-	0,3,3	-	-	-		
11	ADP	h	601	12	28,29,29	1.41	5 (17%)	43,45,45	1.84	9 (20%)
13	AF3	D	603	-	0,3,3	-	-	-		
13	AF3	E	603	-	0,3,3	-	-	-		
11	ADP	z	601	12	28,29,29	1.39	4 (14%)	43,45,45	1.83	9 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
13	AF3	b	603	-	0,3,3	-	-	-		
11	ADP	B	601	12	28,29,29	1.39	5 (17%)	43,45,45	1.81	9 (20%)
13	AF3	e	603	-	0,3,3	-	-	-		
11	ADP	e	601	12	28,29,29	1.39	4 (14%)	43,45,45	1.81	9 (20%)
11	ADP	E	601	12	28,29,29	1.38	4 (14%)	43,45,45	1.81	9 (20%)
11	ADP	D	601	12	28,29,29	1.38	5 (17%)	43,45,45	1.80	9 (20%)
13	AF3	B	603	-	0,3,3	-	-	-		
13	AF3	Q	603	-	0,3,3	-	-	-		
11	ADP	q	601	12	28,29,29	1.37	5 (17%)	43,45,45	1.84	9 (20%)
13	AF3	G	603	-	0,3,3	-	-	-		
13	AF3	a	603	-	0,3,3	-	-	-		
11	ADP	b	601	12	28,29,29	1.38	4 (14%)	43,45,45	1.83	9 (20%)
13	AF3	d	603	-	0,3,3	-	-	-		
13	AF3	H	603	-	0,3,3	-	-	-		
13	AF3	Z	603	-	0,3,3	-	-	-		
11	ADP	a	601	12	28,29,29	1.41	4 (14%)	43,45,45	1.81	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
11	ADP	B	601	12	-	0/16/32/32	0/3/3/3
11	ADP	H	601	12	-	3/16/32/32	0/3/3/3
11	ADP	Q	601	12	-	3/16/32/32	0/3/3/3
11	ADP	A	601	12	-	7/16/32/32	0/3/3/3
11	ADP	E	601	12	-	3/16/32/32	0/3/3/3
11	ADP	D	601	12	-	1/16/32/32	0/3/3/3
11	ADP	Z	601	12	-	1/16/32/32	0/3/3/3
11	ADP	e	601	12	-	5/16/32/32	0/3/3/3
11	ADP	z	601	12	-	2/16/32/32	0/3/3/3
11	ADP	d	601	12	-	3/16/32/32	0/3/3/3
11	ADP	g	601	12	-	5/16/32/32	0/3/3/3
11	ADP	q	601	12	-	0/16/32/32	0/3/3/3
11	ADP	h	601	12	-	6/16/32/32	0/3/3/3
11	ADP	b	601	12	-	4/16/32/32	0/3/3/3
11	ADP	G	601	12	-	6/16/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	ADP	a	601	12	-	7/16/32/32	0/3/3/3

The worst 5 of 73 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	Z	601	ADP	C5-C4	4.60	1.47	1.39
11	e	601	ADP	C5-C4	4.58	1.47	1.39
11	b	601	ADP	C5-C4	4.56	1.47	1.39
11	A	601	ADP	C5-C4	4.55	1.47	1.39
11	B	601	ADP	C5-C4	4.53	1.47	1.39

The worst 5 of 145 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	Z	601	ADP	C5-C4-N3	-6.09	118.33	126.72
11	A	601	ADP	C5-C4-N3	-5.99	118.46	126.72
11	z	601	ADP	C5-C4-N3	-5.99	118.47	126.72
11	h	601	ADP	C5-C4-N3	-5.84	118.67	126.72
11	a	601	ADP	C5-C4-N3	-5.82	118.70	126.72

There are no chirality outliers.

5 of 56 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	E	601	ADP	C5'-O5'-PA-O2A
11	E	601	ADP	C5'-O5'-PA-O3A
11	G	601	ADP	C5'-O5'-PA-O1A
11	H	601	ADP	C5'-O5'-PA-O3A
11	a	601	ADP	C5'-O5'-PA-O1A

There are no ring outliers.

25 monomers are involved in 43 short contacts:

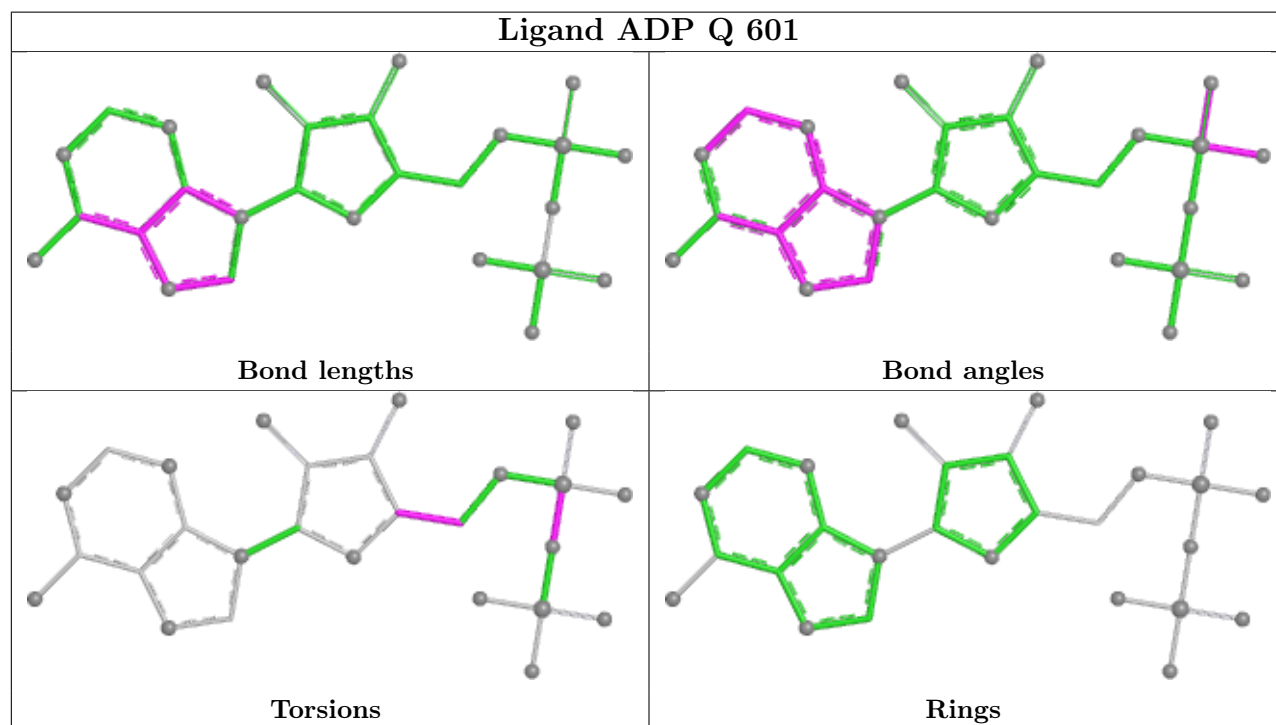
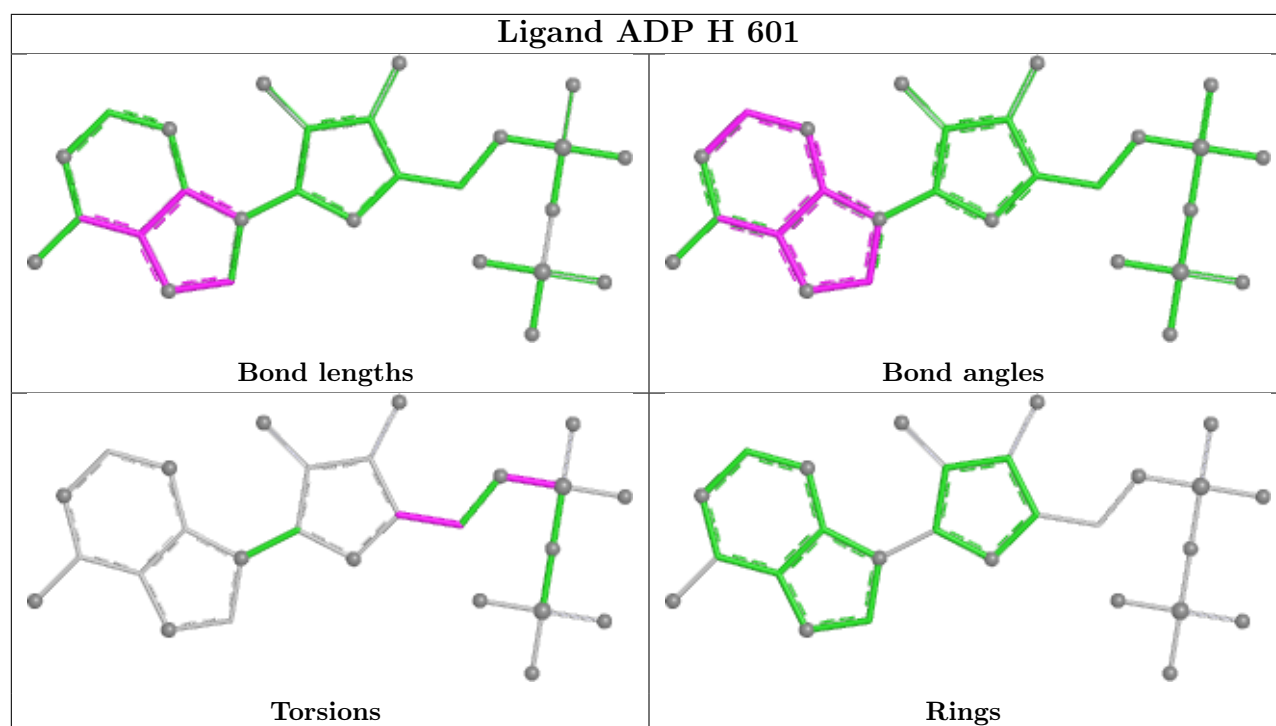
Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	h	603	AF3	1	0
11	H	601	ADP	4	0
11	Q	601	ADP	2	0
11	Z	601	ADP	1	0
13	z	603	AF3	3	0
11	G	601	ADP	2	0
11	A	601	ADP	1	0

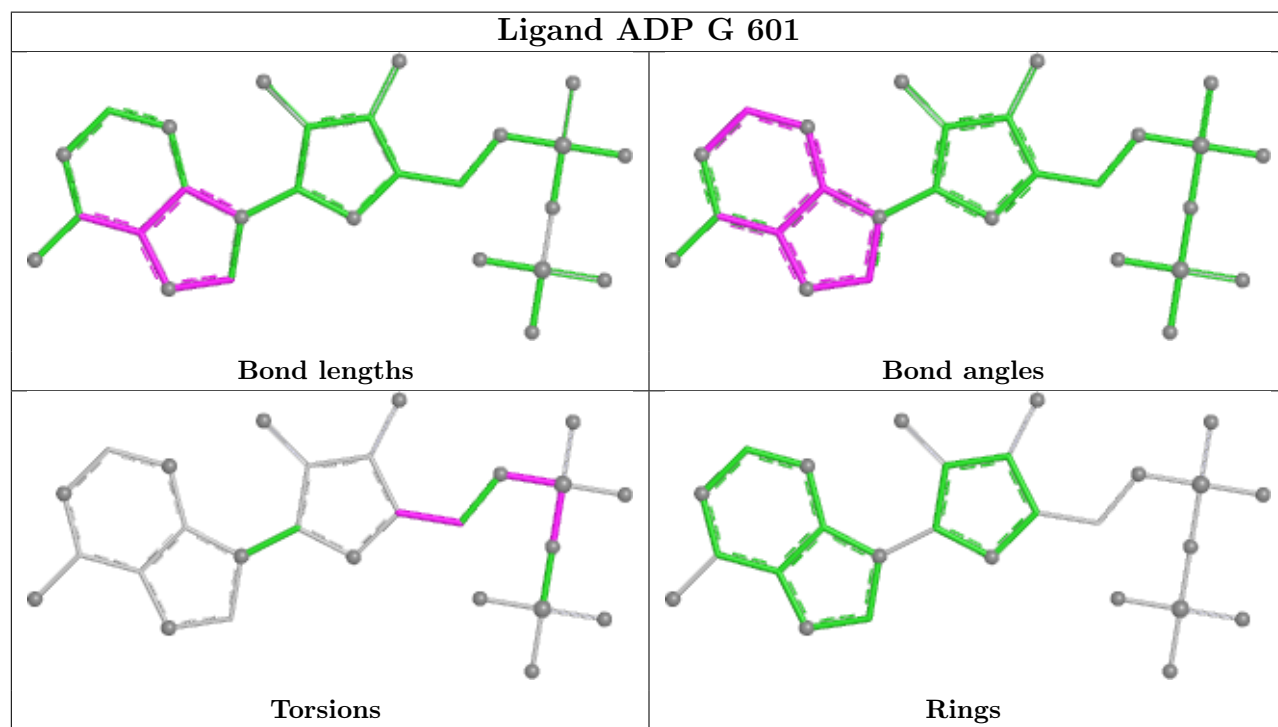
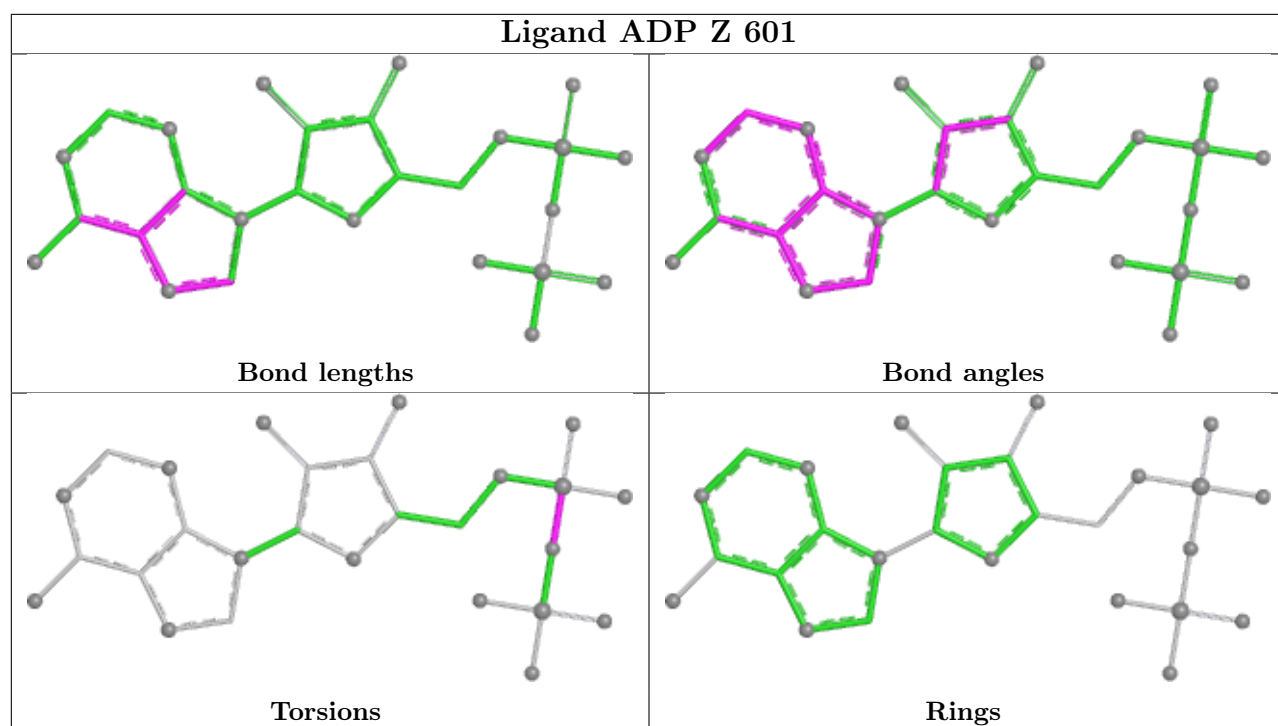
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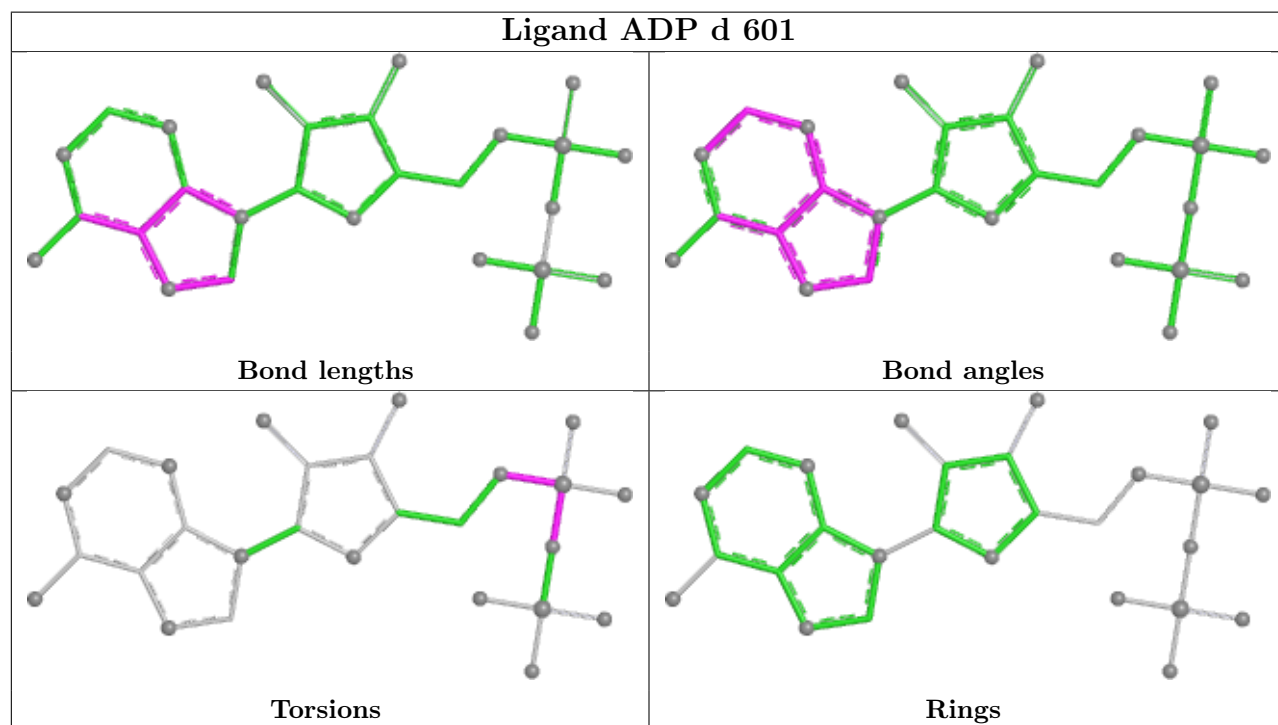
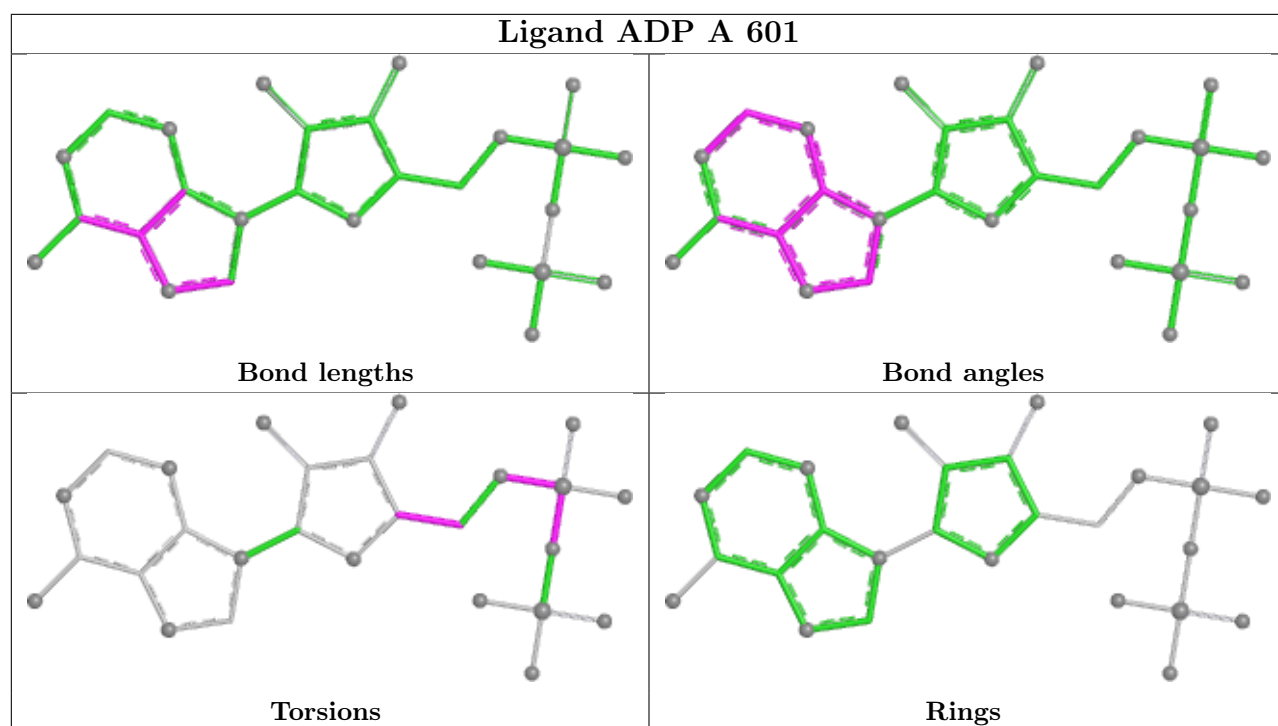
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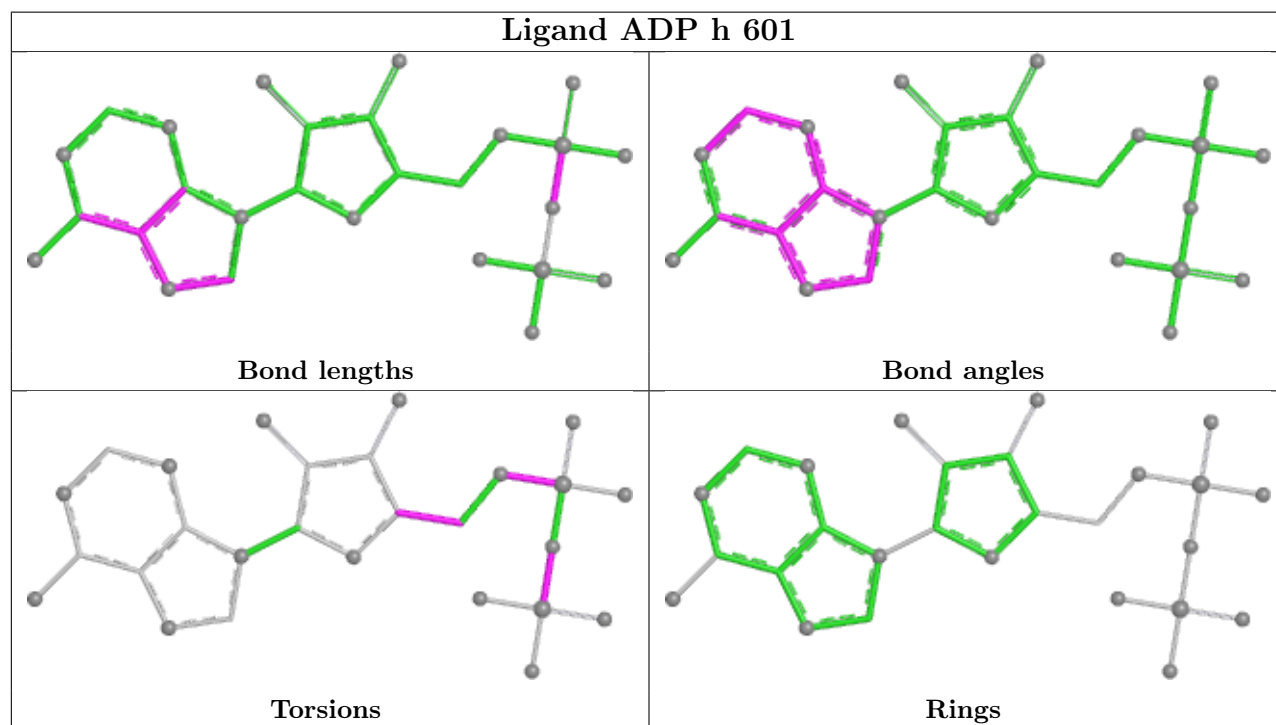
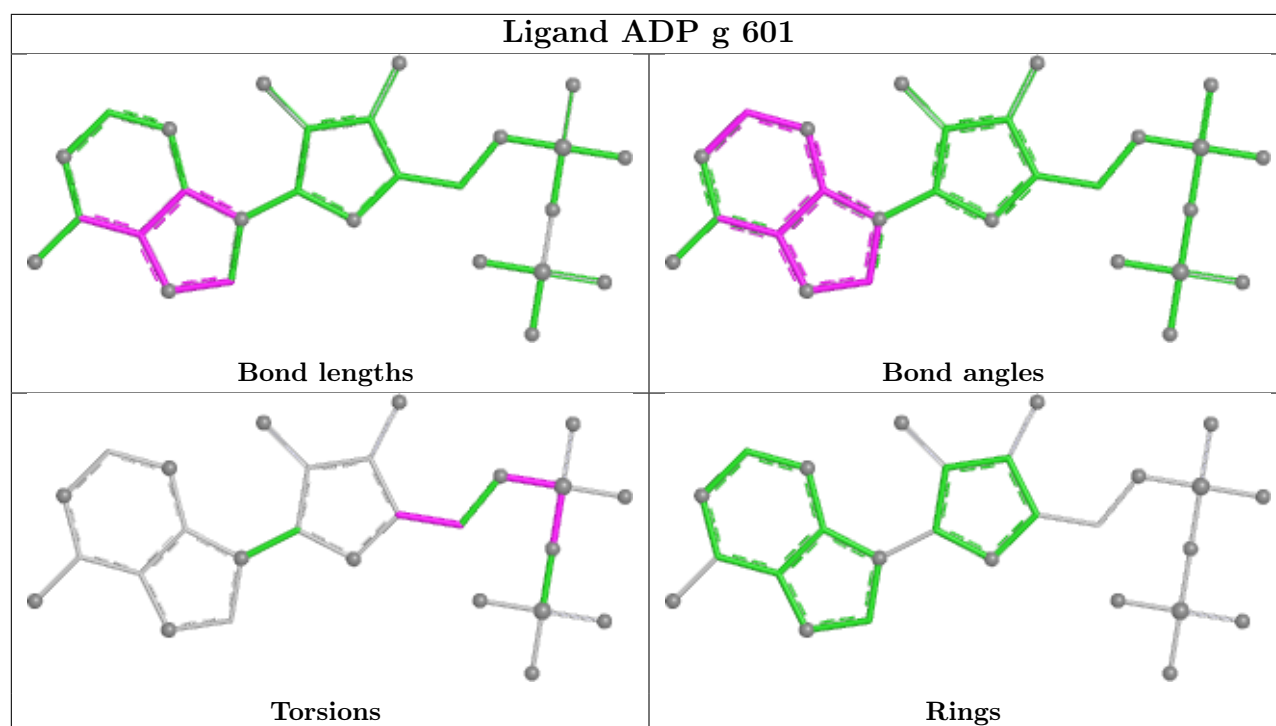
Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	q	603	AF3	2	0
11	g	601	ADP	3	0
13	A	603	AF3	1	0
11	h	601	ADP	1	0
11	z	601	ADP	2	0
13	b	603	AF3	4	0
11	B	601	ADP	5	0
13	e	603	AF3	2	0
11	e	601	ADP	1	0
13	B	603	AF3	2	0
13	Q	603	AF3	2	0
11	q	601	ADP	2	0
13	G	603	AF3	2	0
13	a	603	AF3	3	0
11	b	601	ADP	3	0
13	H	603	AF3	2	0
13	Z	603	AF3	2	0
11	a	601	ADP	3	0

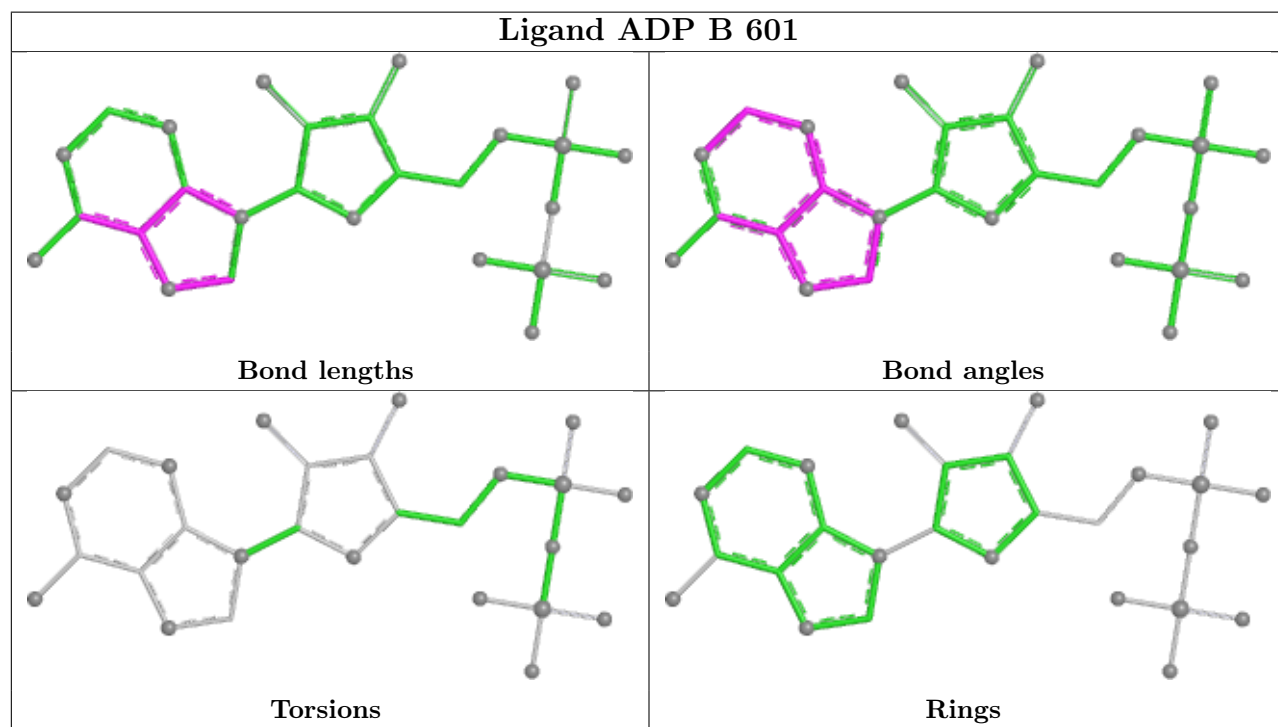
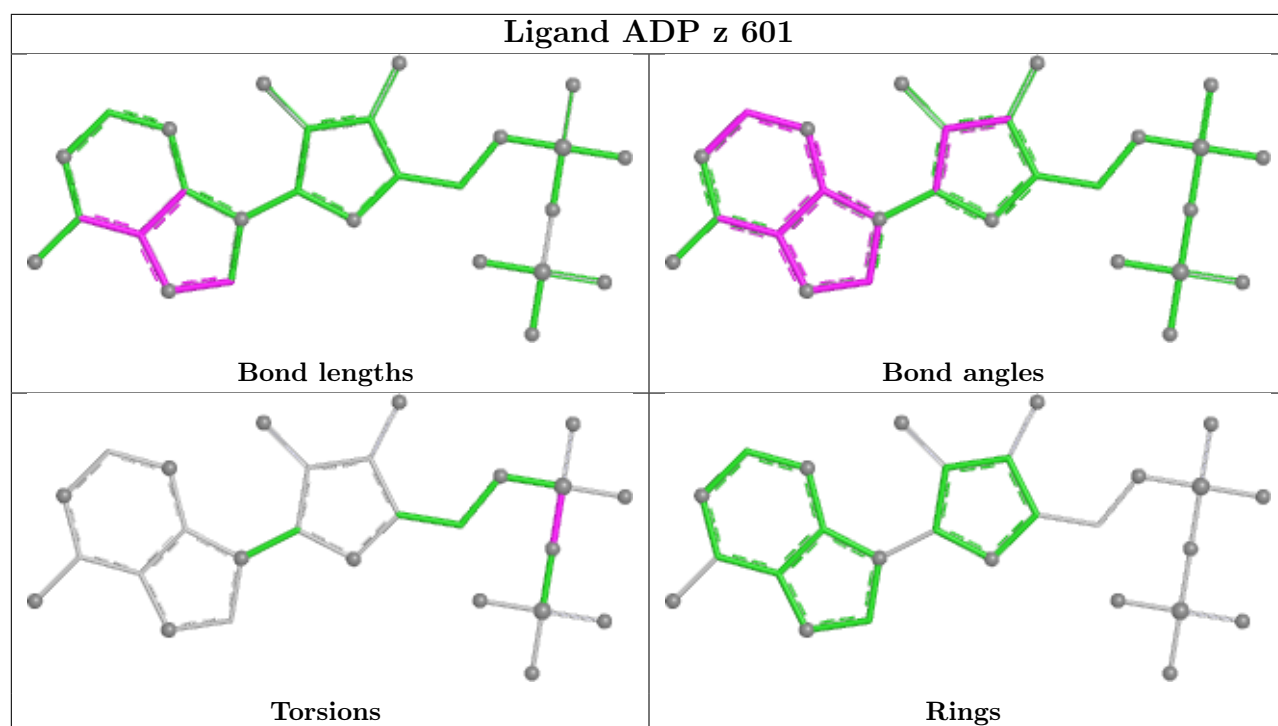
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

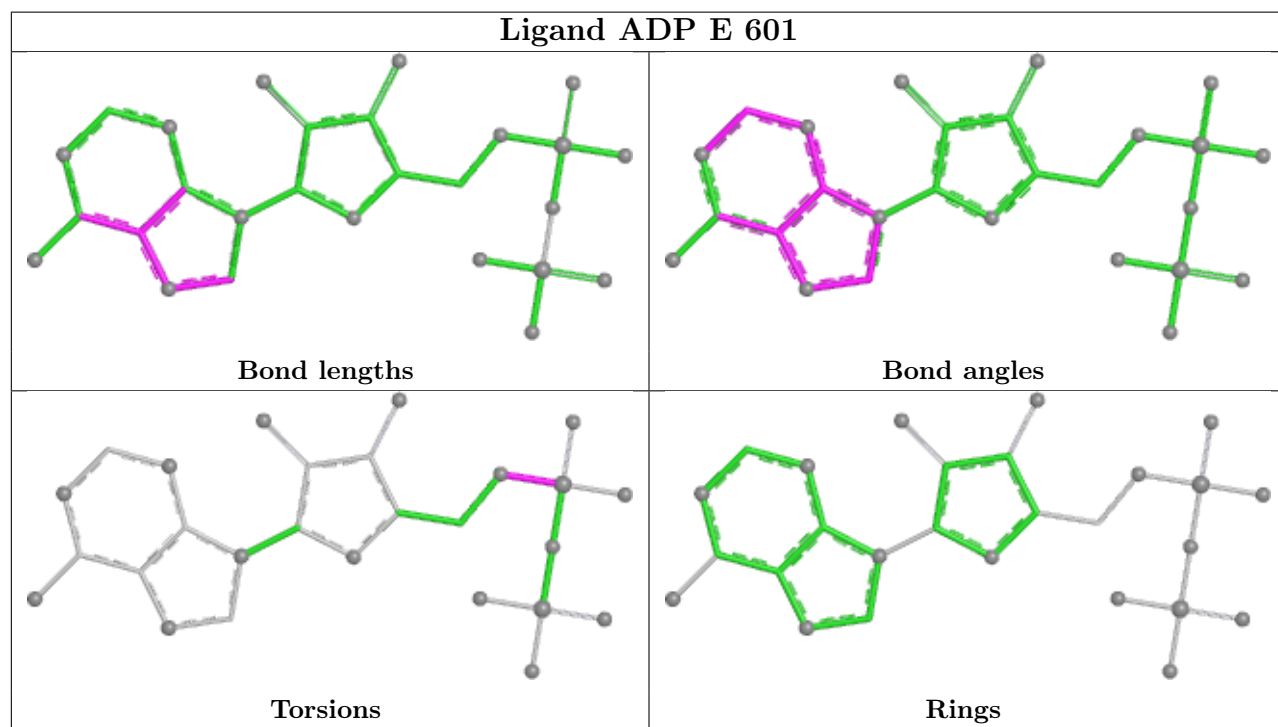
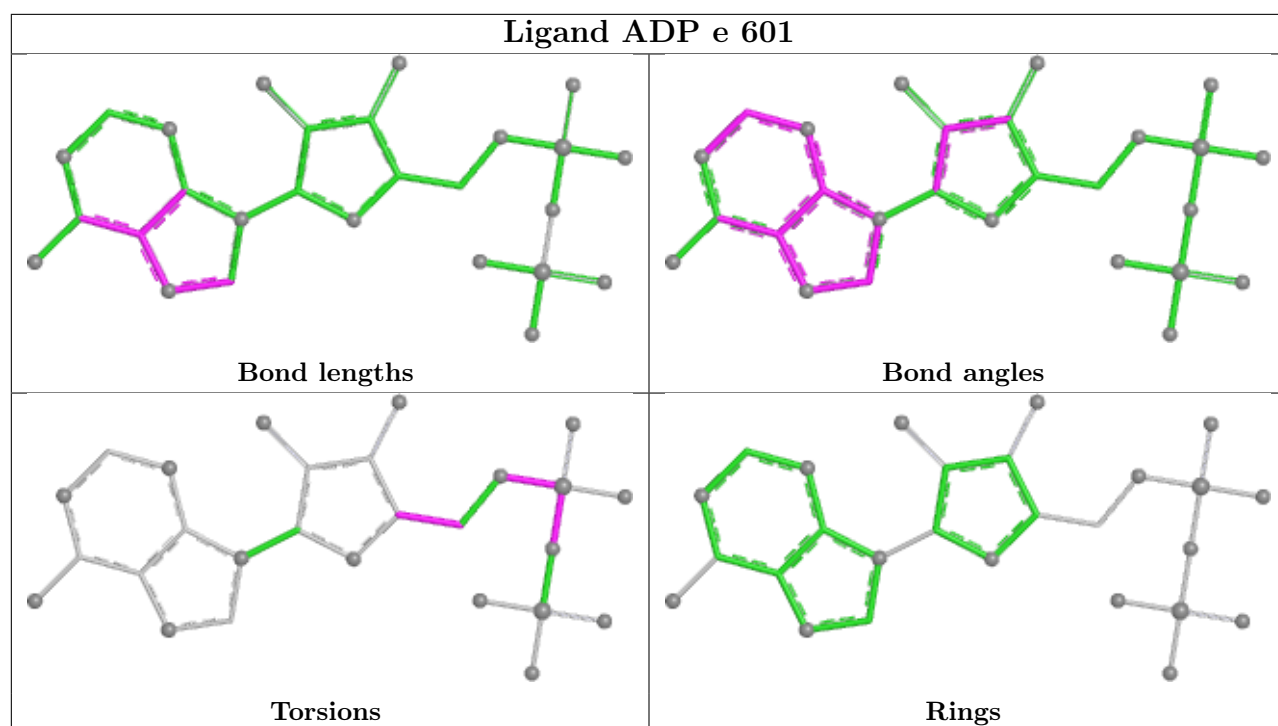


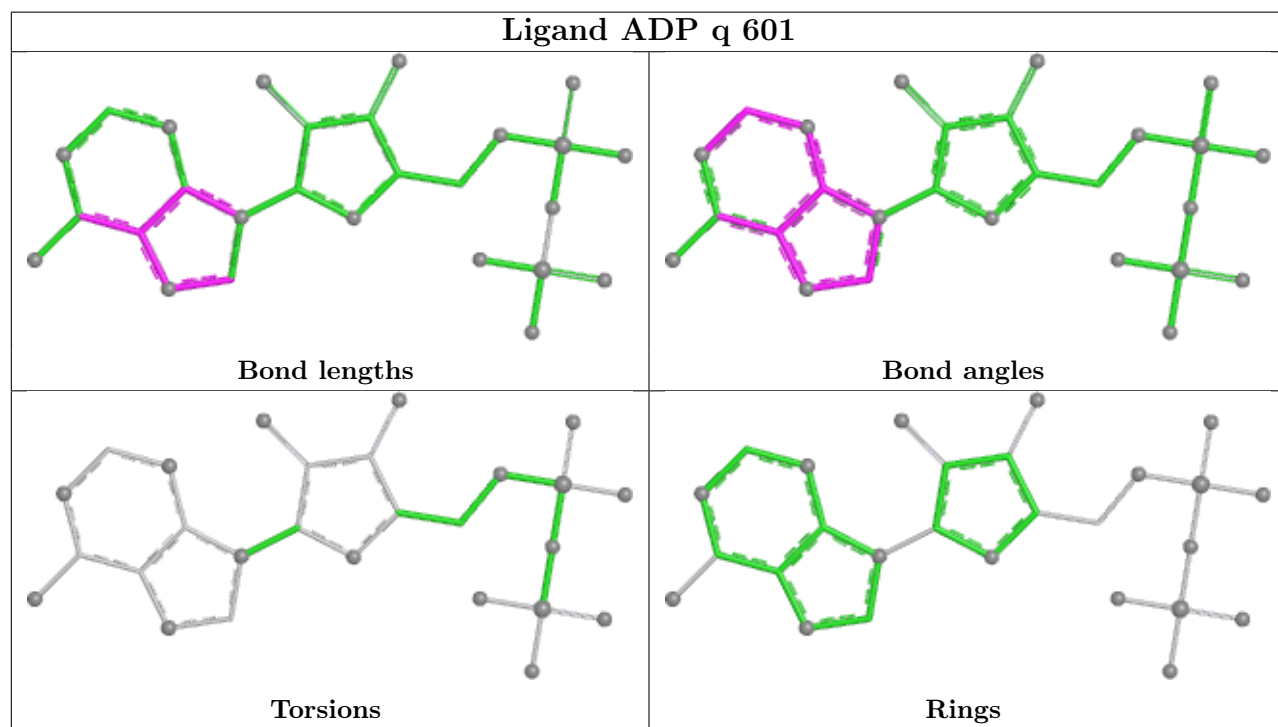
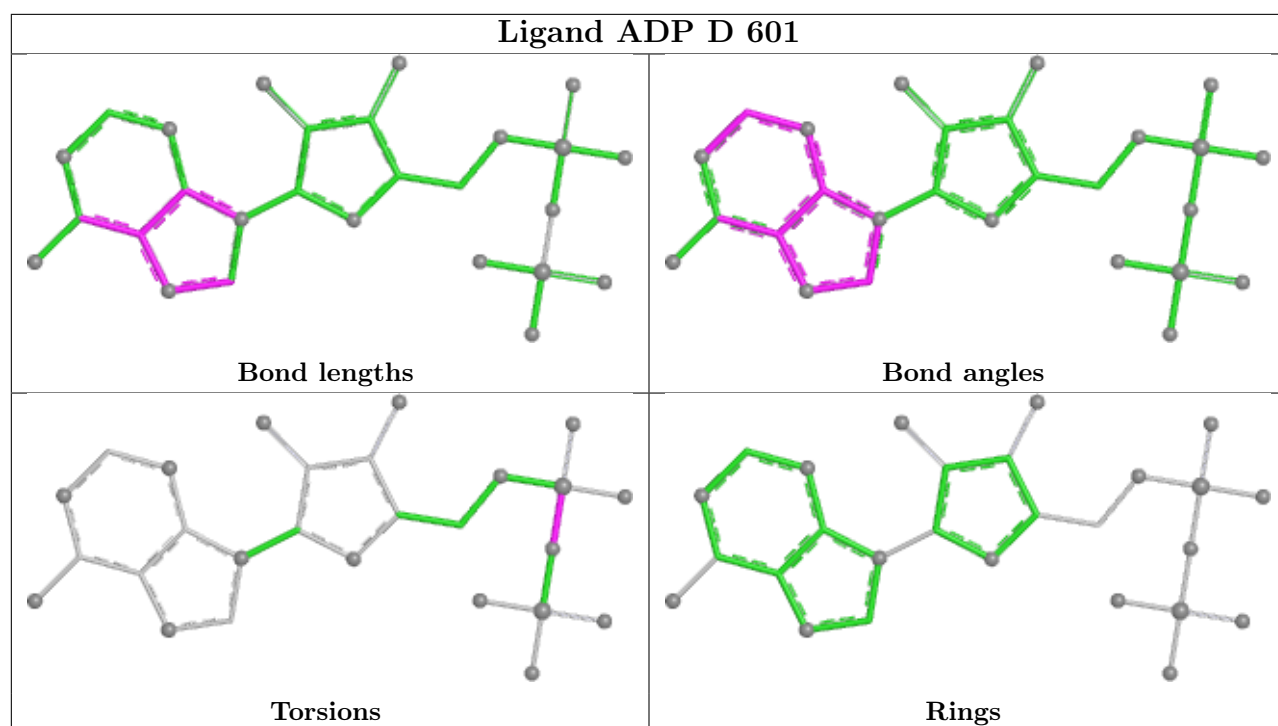


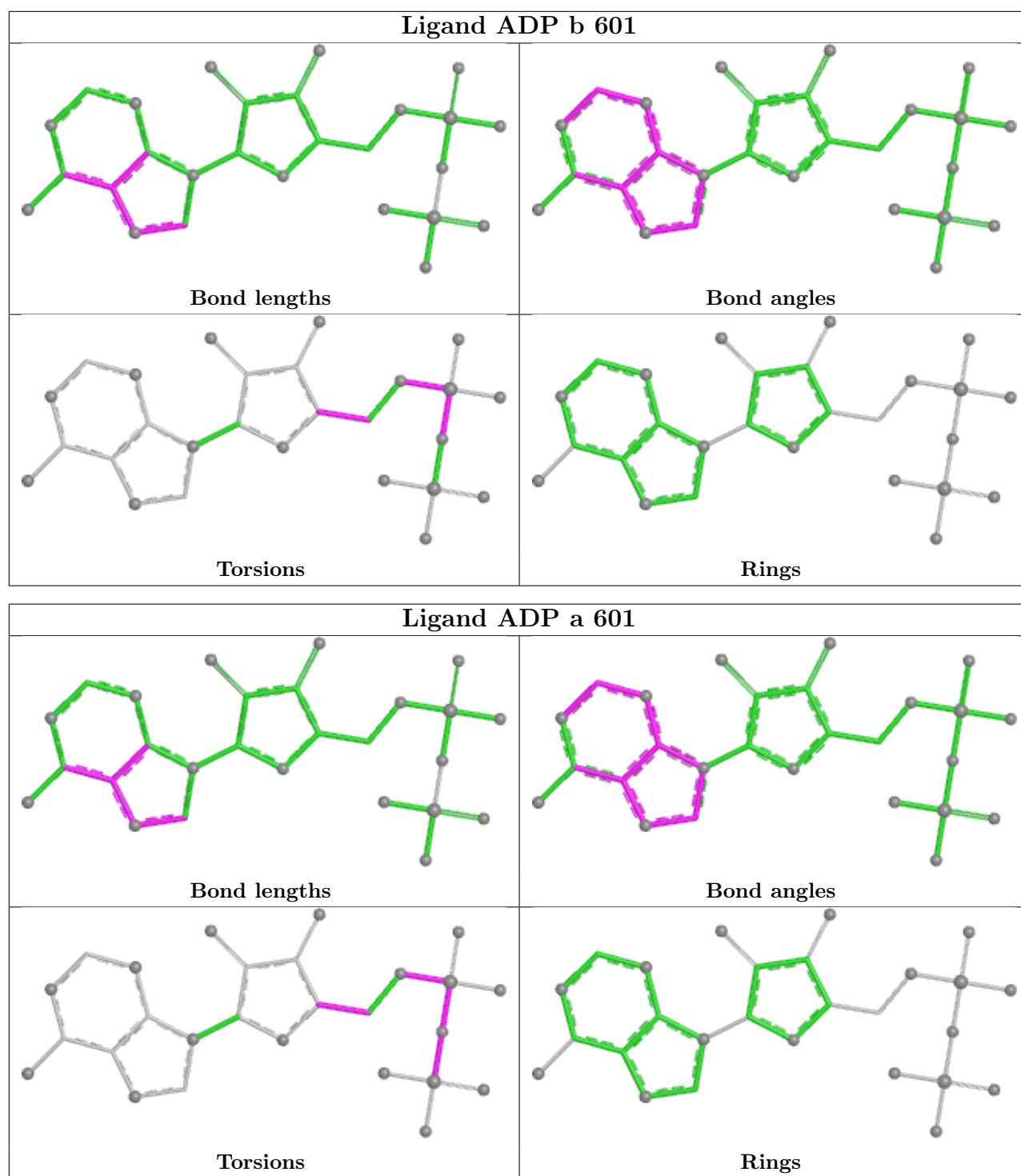












5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

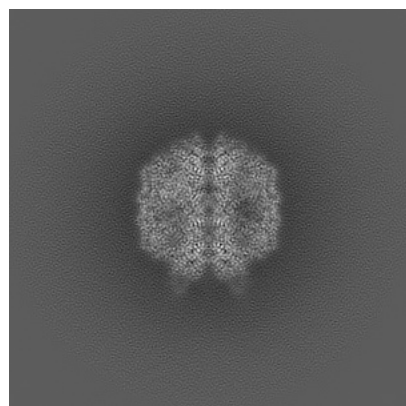
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12607. These allow visual inspection of the internal detail of the map and identification of artifacts.

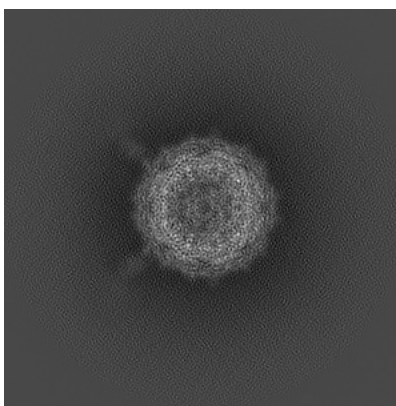
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

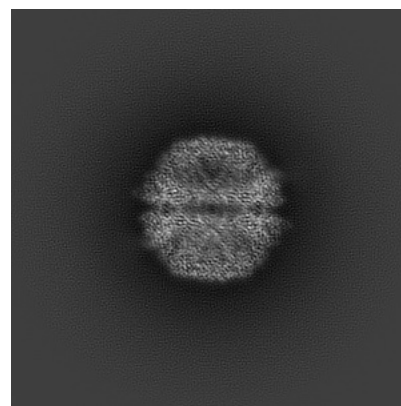
6.1.1 Primary map



X

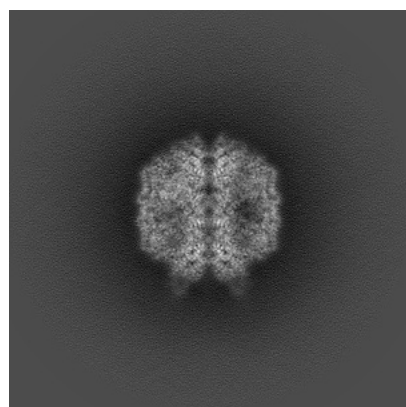


Y

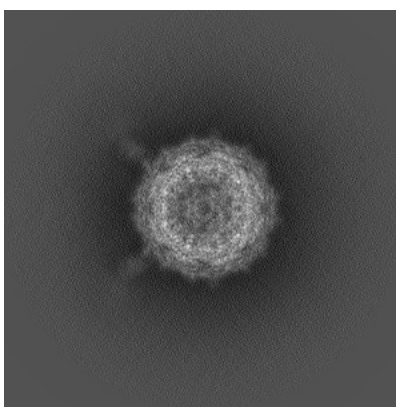


Z

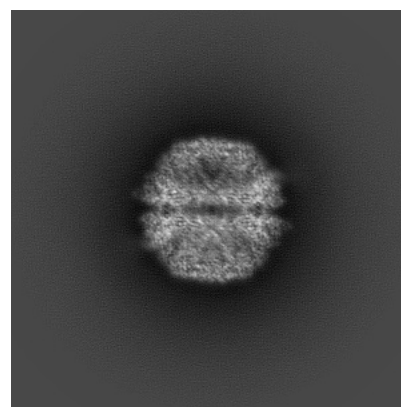
6.1.2 Raw map



X



Y

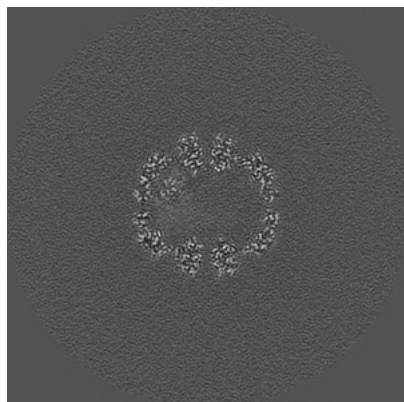


Z

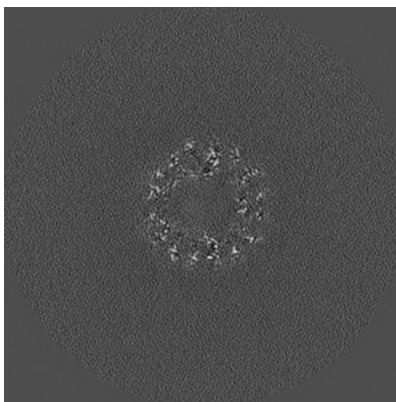
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

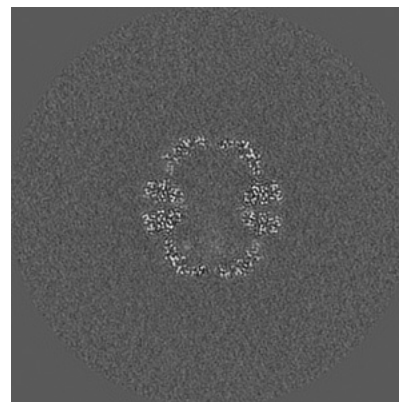
6.2.1 Primary map



X Index: 200

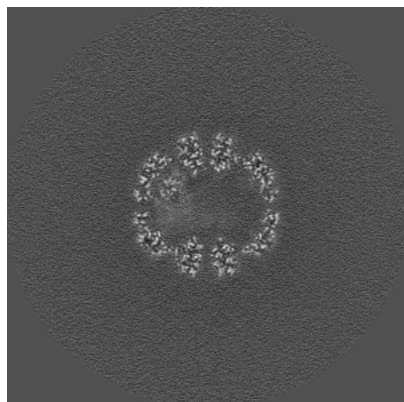


Y Index: 200

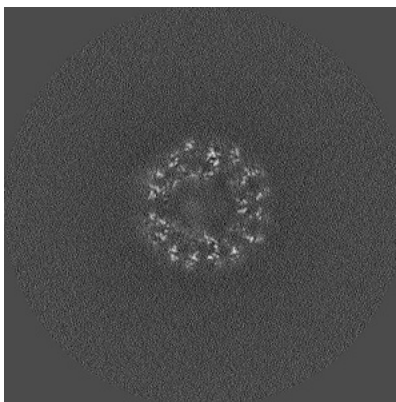


Z Index: 200

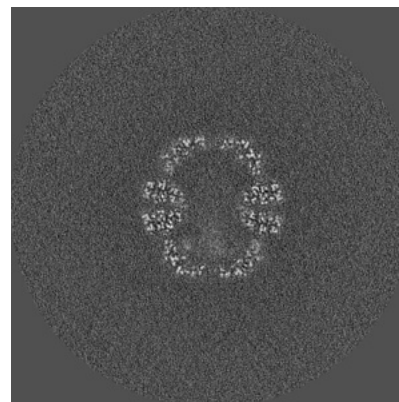
6.2.2 Raw map



X Index: 200



Y Index: 200

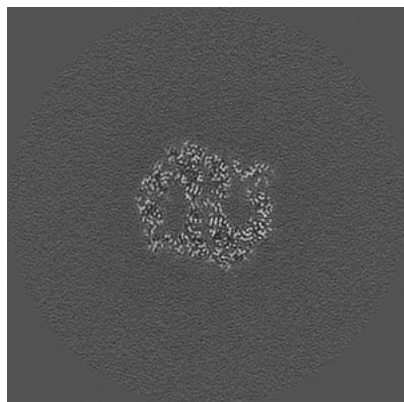


Z Index: 200

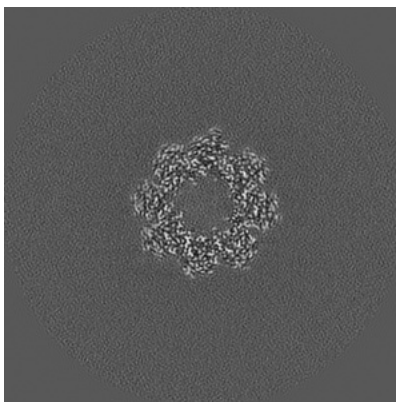
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

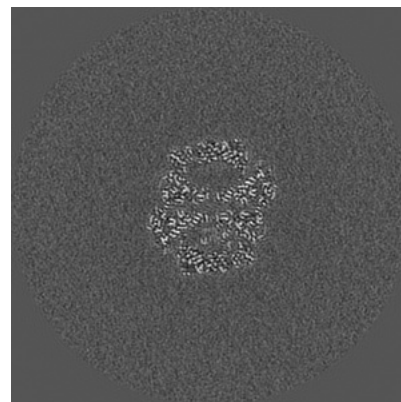
6.3.1 Primary map



X Index: 233

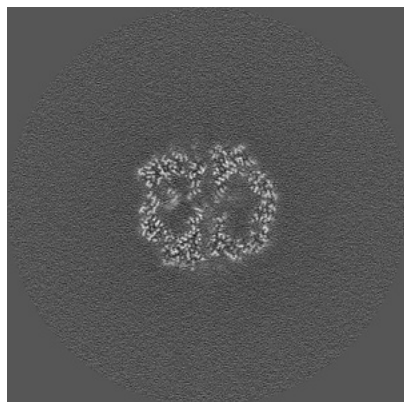


Y Index: 186

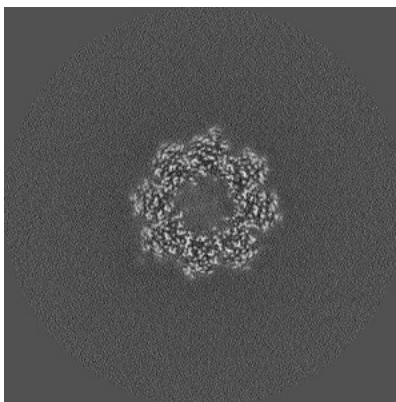


Z Index: 231

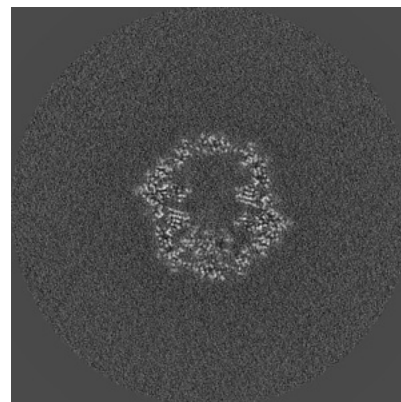
6.3.2 Raw map



X Index: 174



Y Index: 186

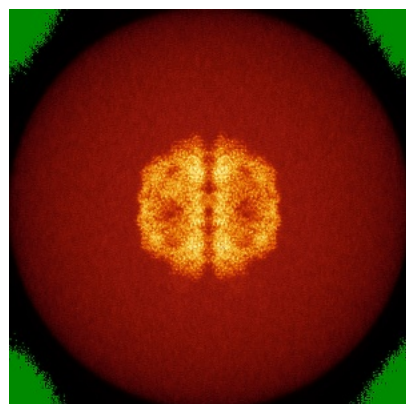


Z Index: 214

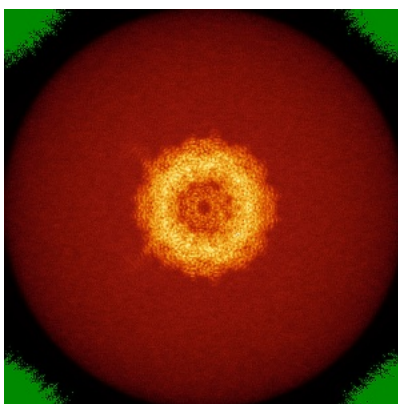
The images above show the largest variance slices of the map in three orthogonal directions.

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

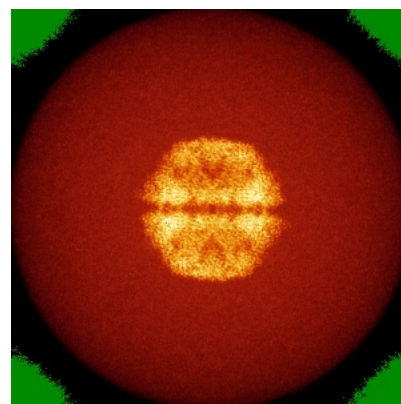
6.4.1 Primary map



X

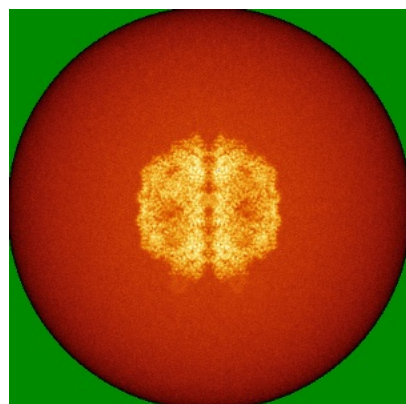


Y

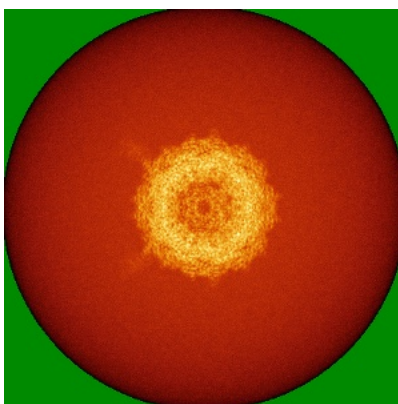


Z

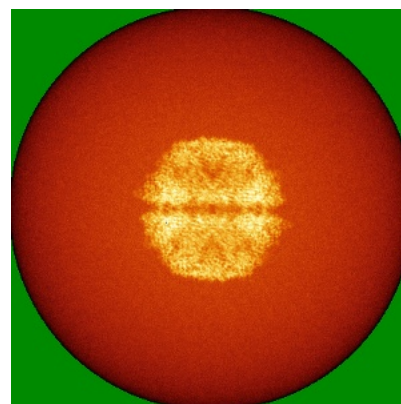
6.4.2 Raw map



X



Y

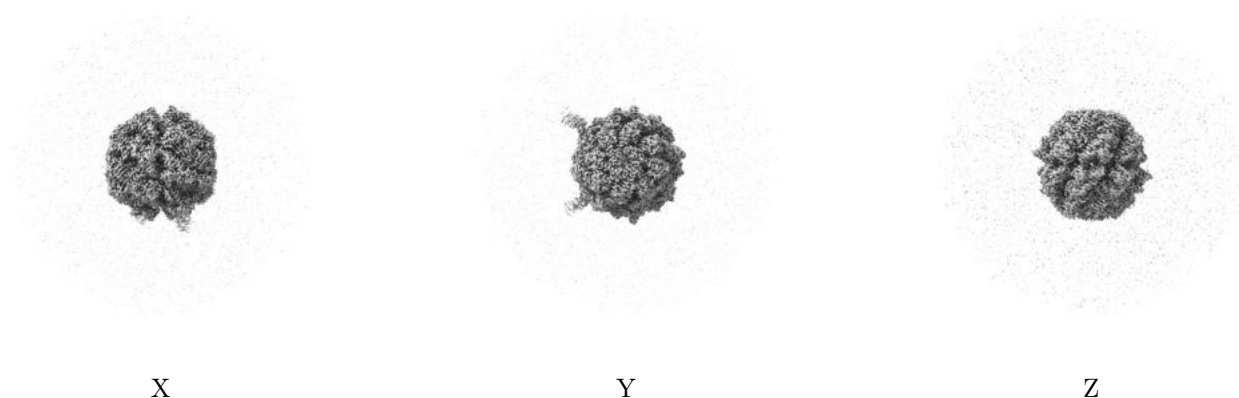


Z

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.

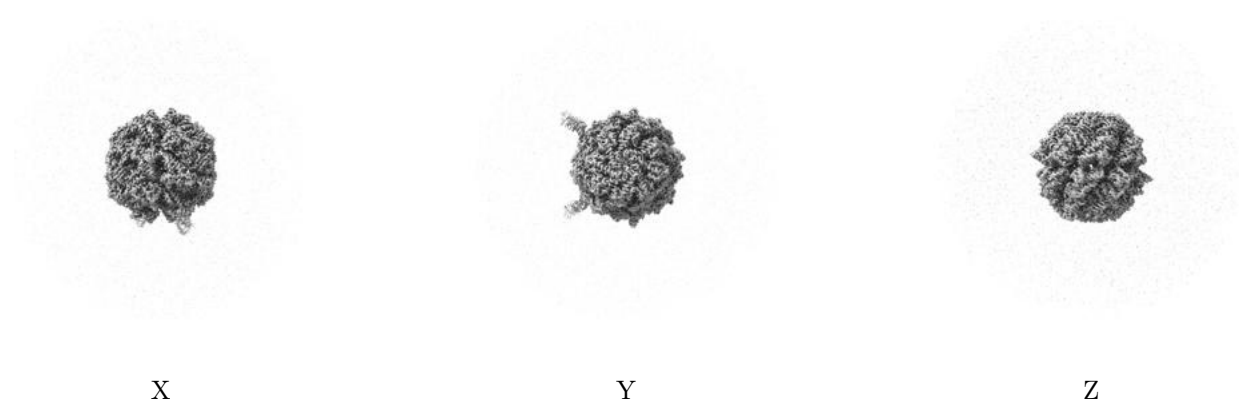
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.025. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

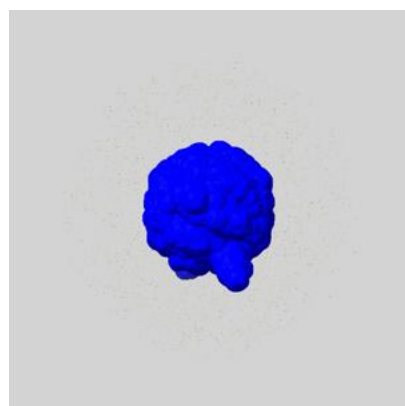
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

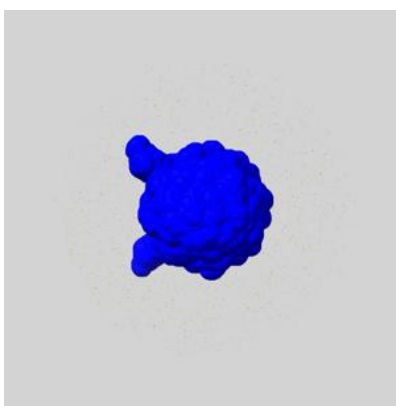
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

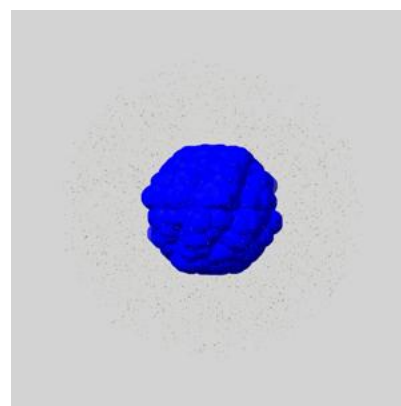
6.6.1 emd_12607_msk_1.map [i](#)



X



Y

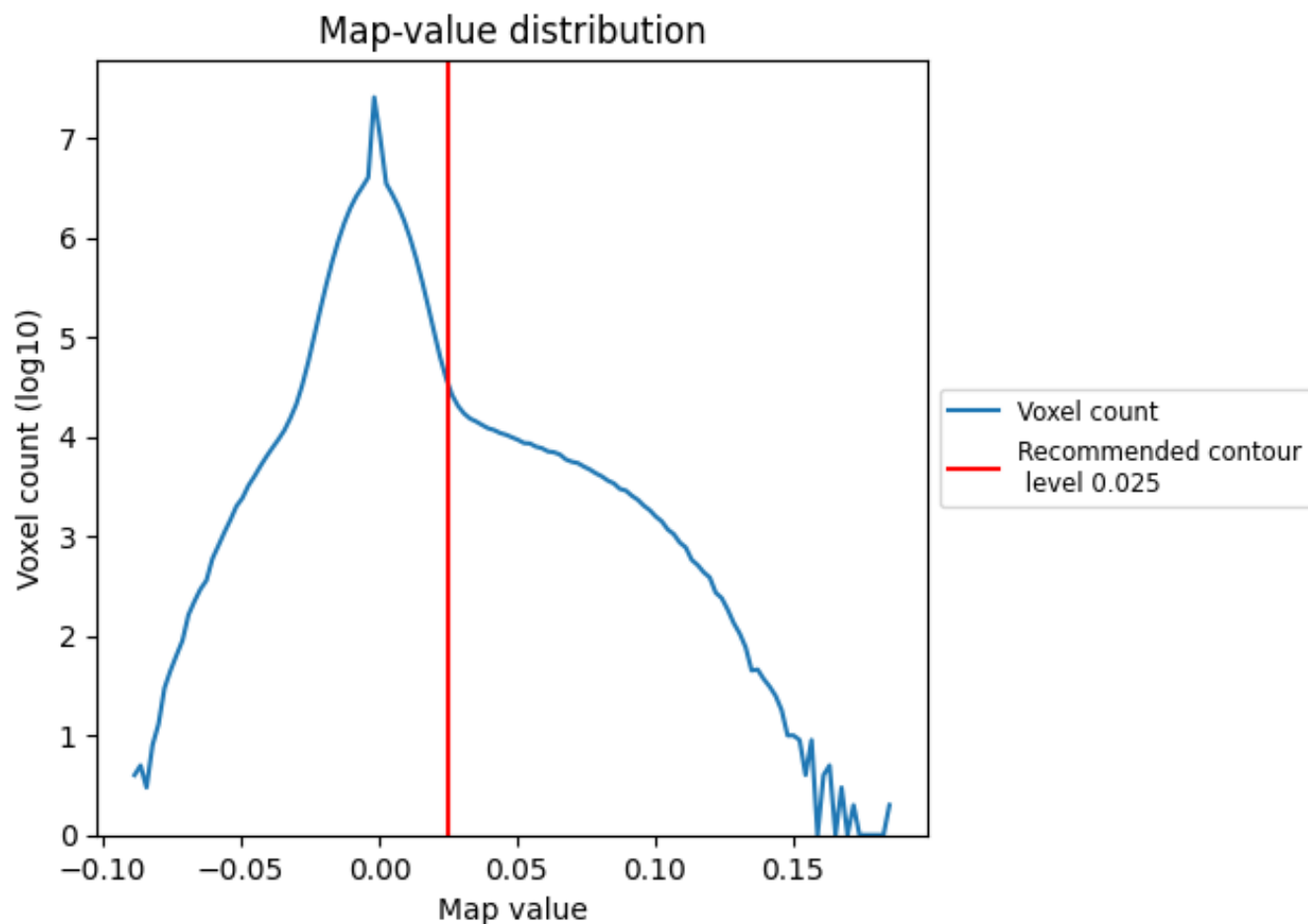


Z

7 Map analysis [i](#)

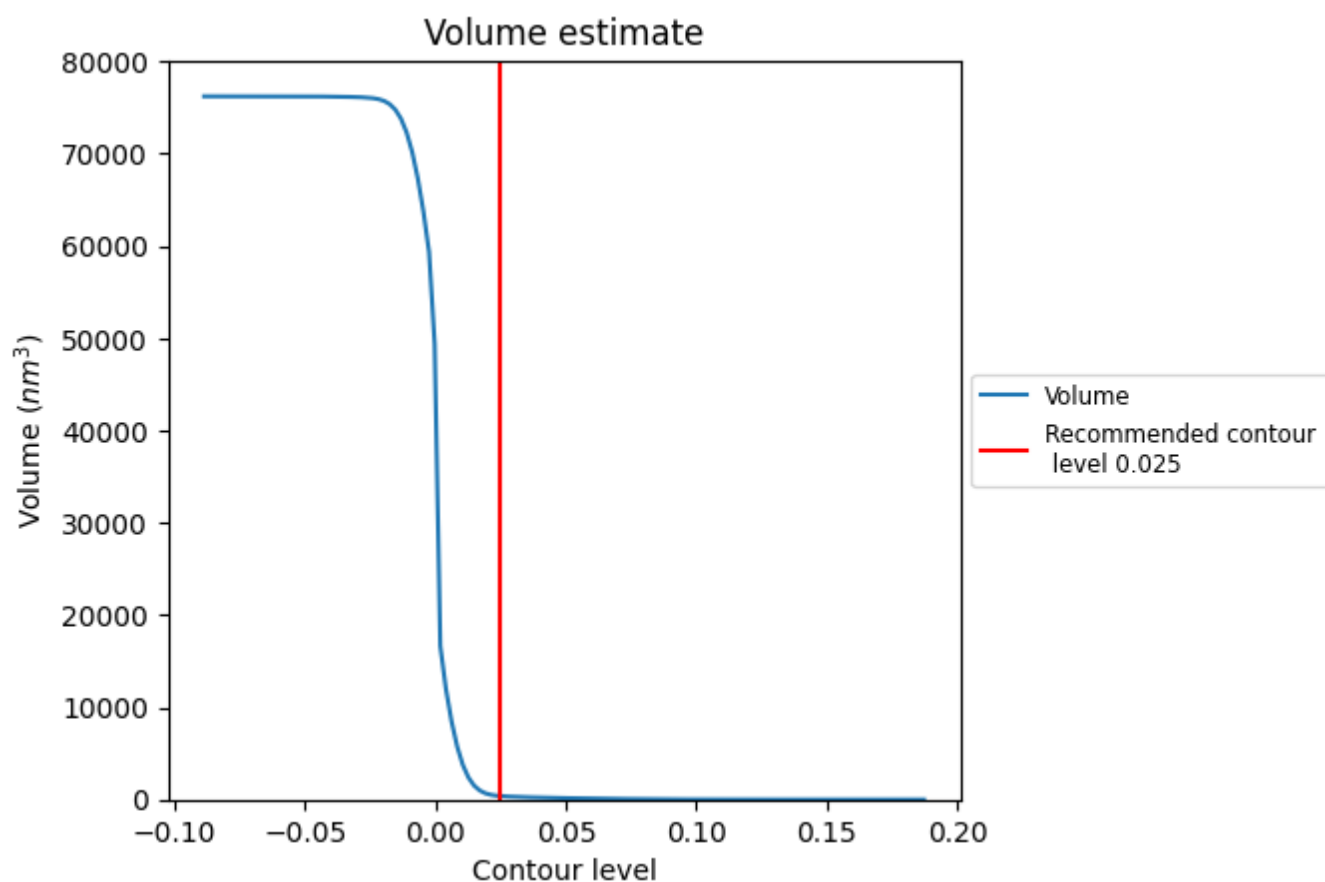
This section contains the results of statistical analysis of the map.

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.

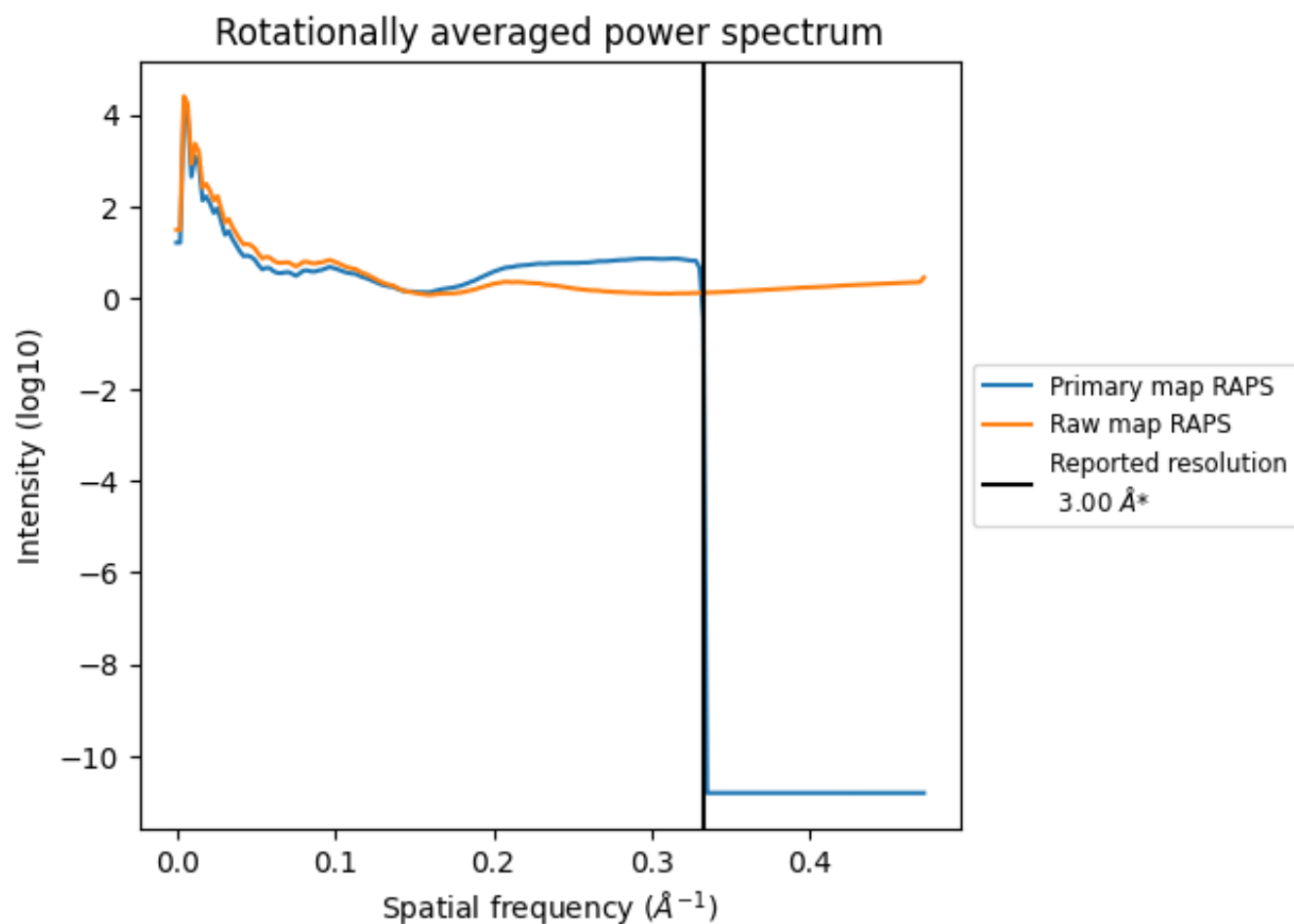
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 380 nm³; this corresponds to an approximate mass of 343 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 ⓘ

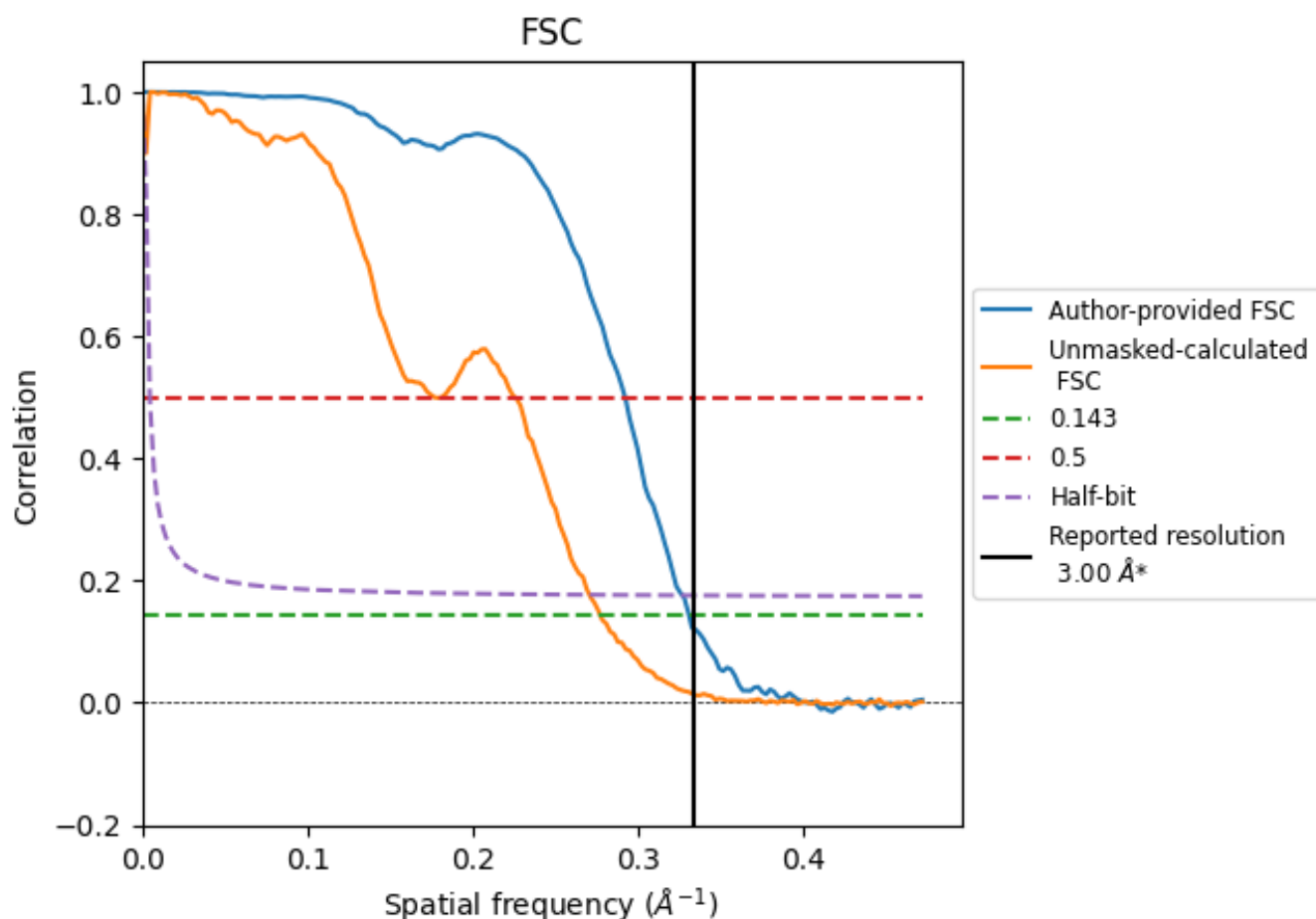


*Reported resolution corresponds to spatial frequency of 0.333 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.333 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

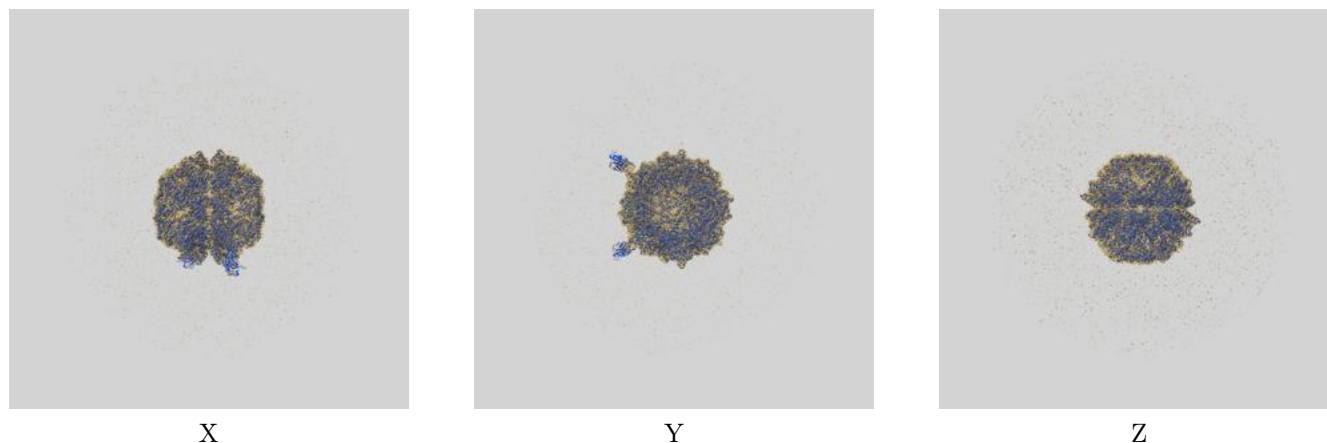
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	3.02	3.42	3.06
Unmasked-calculated*	3.61	5.66	3.70

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.61 differs from the reported value 3.0 by more than 10 %

9 Map-model fit [i](#)

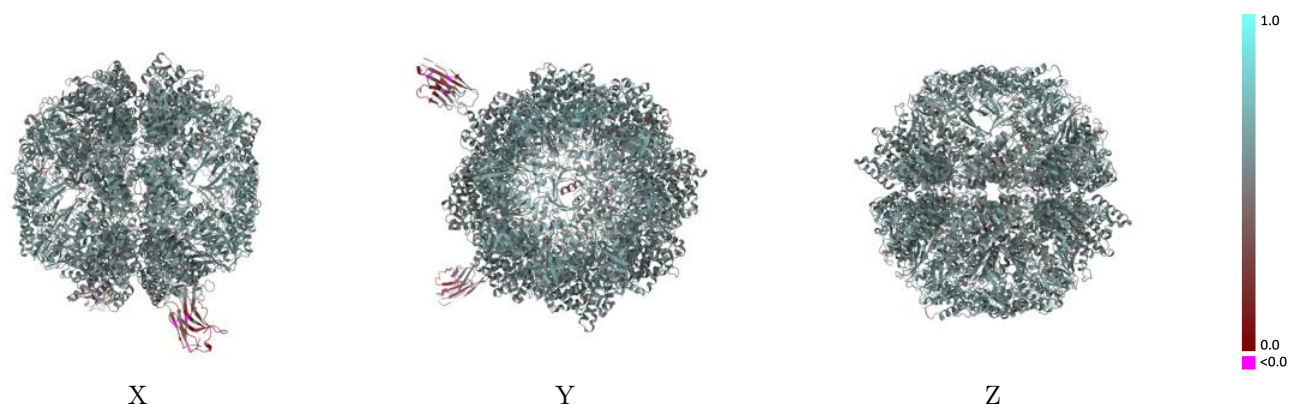
This section contains information regarding the fit between EMDB map EMD-12607 and PDB model 7NVN. Per-residue inclusion information can be found in [section 3](#) on [page 11](#).

9.1 Map-model overlay [i](#)



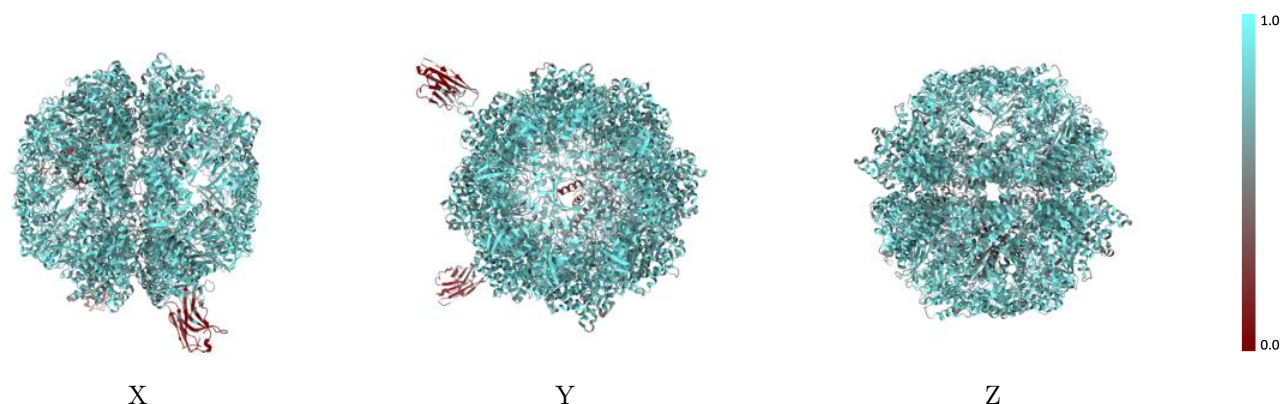
The images above show the 3D surface view of the map at the recommended contour level 0.025 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



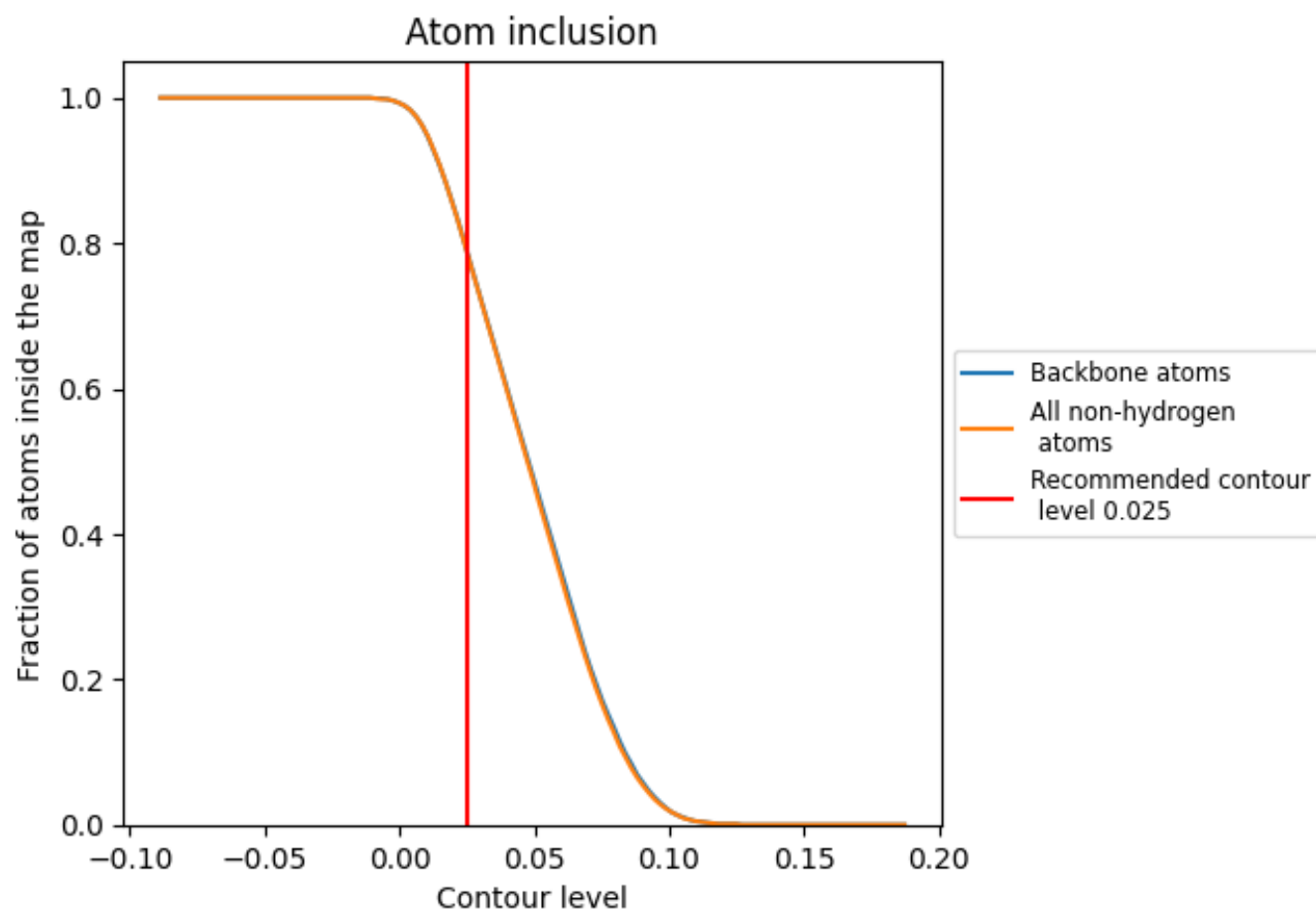
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.025).

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 79% 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 (0.025) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7880	 0.5580
A	 0.8040	 0.5630
B	 0.8020	 0.5530
D	 0.8140	 0.5660
E	 0.8140	 0.5650
G	 0.8110	 0.5630
H	 0.8310	 0.5860
N	 0.2410	 0.3180
Q	 0.8190	 0.5760
T	 0.6240	 0.5150
Z	 0.8070	 0.5600
a	 0.8030	 0.5590
b	 0.8080	 0.5550
d	 0.8070	 0.5620
e	 0.8170	 0.5660
g	 0.8040	 0.5590
h	 0.8420	 0.5860
n	 0.2510	 0.3100
q	 0.8210	 0.5750
z	 0.7990	 0.5530

