



## wwPDB EM Validation Summary Report ⓘ

Mar 31, 2025 – 02:24 PM JST

PDB ID : 3G37 / pdb\_00003g37  
EMDB ID : EMD-1674  
Title : Cryo-EM structure of actin filament in the presence of phosphate  
Authors : Wakabayshi, T.; Murakami, K.; Yasunaga, T.; Noguchi, T.Q.; Uyeda, T.Q.  
Deposited on : 2009-02-02  
Resolution : 6.00 Å(reported)

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We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev117  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
MolProbity : 4.02b-467  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.42

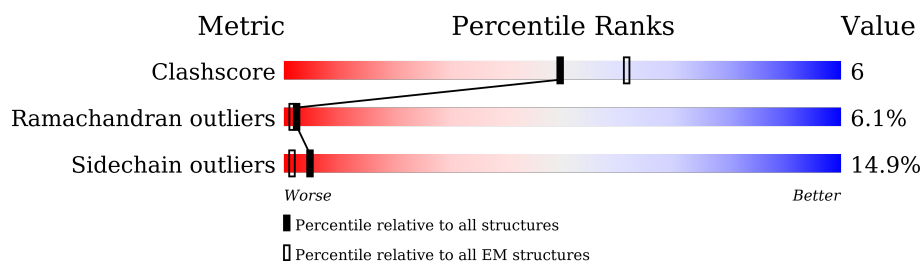
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*





The reported resolution of this entry is 6.00 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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  $\geq 3$ , 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	O	376	
1	P	376	
1	Q	376	
1	R	376	
1	S	376	
1	T	376	
1	U	376	
1	V	376	

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Mol	Chain	Length	Quality of chain
1	W	376	
1	X	376	
1	Y	376	
1	Z	376	

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
3	PO4	O	803	-	X	-	-
3	PO4	P	803	-	X	-	-
3	PO4	R	803	-	X	-	-
3	PO4	S	803	-	X	-	-
3	PO4	T	803	-	X	-	-
3	PO4	U	802	-	X	-	-
3	PO4	V	803	-	X	-	-
3	PO4	X	803	-	X	-	-
3	PO4	Y	801	-	X	X	-
3	PO4	Y	803	-	X	-	-
3	PO4	Z	802	-	X	-	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 35808 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 Actin, alpha skeletal muscle.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	O	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	P	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	Q	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	R	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	S	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	T	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	U	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	V	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	W	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	X	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	Y	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		
1	Z	376	Total	C	N	O	S	0	0
			2936	1857	493	565	21		

There are 12 discrepancies between the modelled and reference sequences:

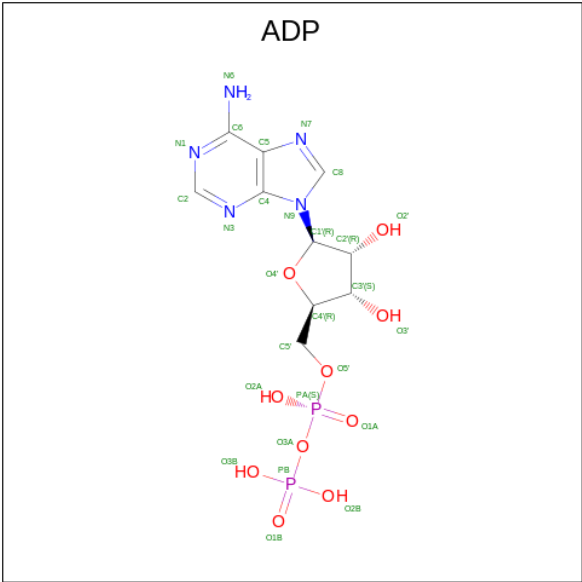
Chain	Residue	Modelled	Actual	Comment	Reference
O	0	ACE	-	acetylation	UNP P68135
P	0	ACE	-	acetylation	UNP P68135
Q	0	ACE	-	acetylation	UNP P68135
R	0	ACE	-	acetylation	UNP P68135
S	0	ACE	-	acetylation	UNP P68135
T	0	ACE	-	acetylation	UNP P68135

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Chain	Residue	Modelled	Actual	Comment	Reference
U	0	ACE	-	acetylation	UNP P68135
V	0	ACE	-	acetylation	UNP P68135
W	0	ACE	-	acetylation	UNP P68135
X	0	ACE	-	acetylation	UNP P68135
Y	0	ACE	-	acetylation	UNP P68135
Z	0	ACE	-	acetylation	UNP P68135

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>).



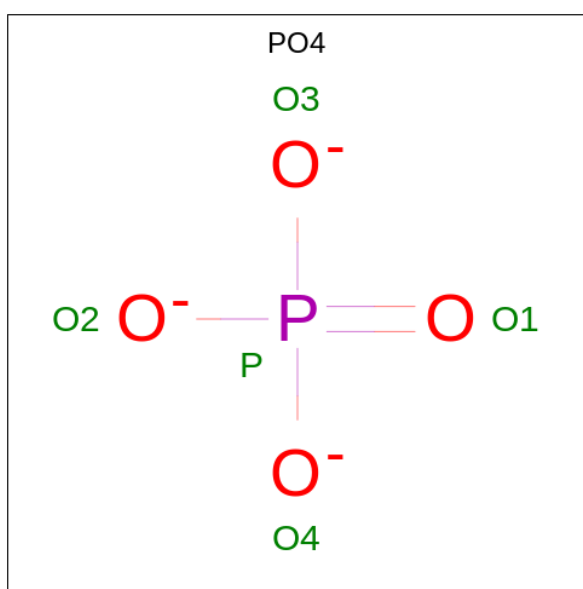
Mol	Chain	Residues	Atoms					AltConf
2	O	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	P	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	Q	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	R	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	S	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	T	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	U	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	V	1	Total	C	N	O	P	0
			27	10	5	10	2	

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Mol	Chain	Residues	Atoms					AltConf
2	W	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	X	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	Y	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	Z	1	Total	C	N	O	P	0
			27	10	5	10	2	

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			AltConf
3	O	1	Total	O	P	0
			5	4	1	
3	O	1	Total	O	P	0
			5	4	1	
3	O	1	Total	O	P	0
			5	4	1	
3	P	1	Total	O	P	0
			5	4	1	
3	P	1	Total	O	P	0
			5	4	1	
3	P	1	Total	O	P	0
			5	4	1	
3	Q	1	Total	O	P	0
			5	4	1	

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Mol	Chain	Residues	Atoms			AltConf
3	Q	1	Total 5	O 4	P 1	0
3	Q	1	Total 5	O 4	P 1	0
3	R	1	Total 5	O 4	P 1	0
3	R	1	Total 5	O 4	P 1	0
3	R	1	Total 5	O 4	P 1	0
3	S	1	Total 5	O 4	P 1	0
3	S	1	Total 5	O 4	P 1	0
3	S	1	Total 5	O 4	P 1	0
3	T	1	Total 5	O 4	P 1	0
3	T	1	Total 5	O 4	P 1	0
3	T	1	Total 5	O 4	P 1	0
3	T	1	Total 5	O 4	P 1	0
3	T	1	Total 5	O 4	P 1	0
3	U	1	Total 5	O 4	P 1	0
3	U	1	Total 5	O 4	P 1	0
3	U	1	Total 5	O 4	P 1	0
3	V	1	Total 5	O 4	P 1	0
3	V	1	Total 5	O 4	P 1	0
3	W	1	Total 5	O 4	P 1	0
3	W	1	Total 5	O 4	P 1	0
3	W	1	Total 5	O 4	P 1	0
3	X	1	Total 5	O 4	P 1	0

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Mol	Chain	Residues	Atoms			AltConf
3	X	1	Total	O	P	0
			5	4	1	
3	X	1	Total	O	P	0
			5	4	1	
3	Y	1	Total	O	P	0
			5	4	1	
3	Y	1	Total	O	P	0
			5	4	1	
3	Y	1	Total	O	P	0
			5	4	1	
3	Z	1	Total	O	P	0
			5	4	1	
3	Z	1	Total	O	P	0
			5	4	1	
3	Z	1	Total	O	P	0
			5	4	1	

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
4	O	6	Total	Mg	0
			6	6	
4	P	5	Total	Mg	0
			5	5	
4	Q	7	Total	Mg	0
			7	7	
4	R	6	Total	Mg	0
			6	6	
4	S	6	Total	Mg	0
			6	6	
4	T	6	Total	Mg	0
			6	6	
4	U	6	Total	Mg	0
			6	6	
4	V	6	Total	Mg	0
			6	6	
4	W	6	Total	Mg	0
			6	6	
4	X	6	Total	Mg	0
			6	6	
4	Y	6	Total	Mg	0
			6	6	

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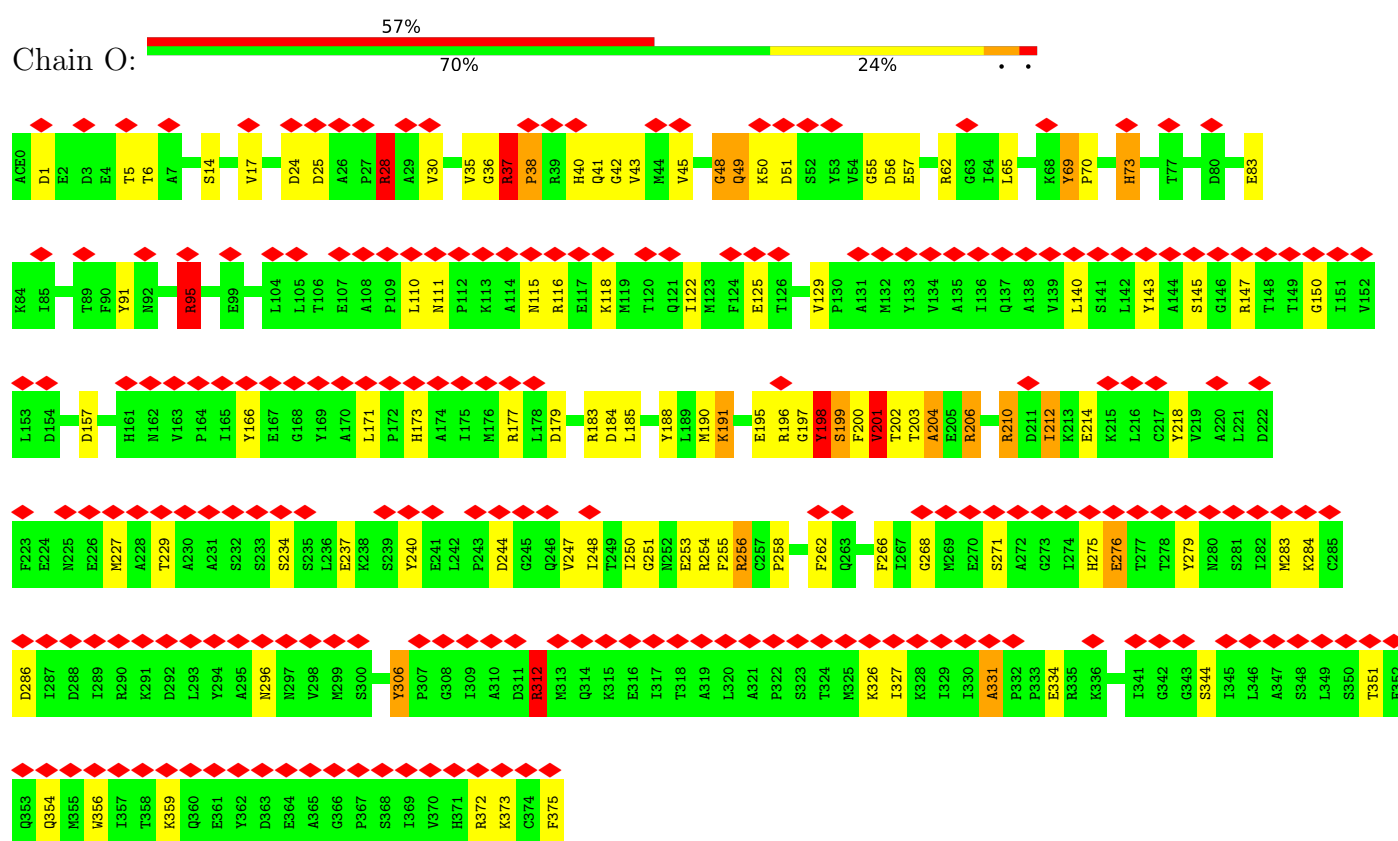
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Mol	Chain	Residues	Atoms		AltConf
			Total	Mg	
4	Z	6	6	6	0

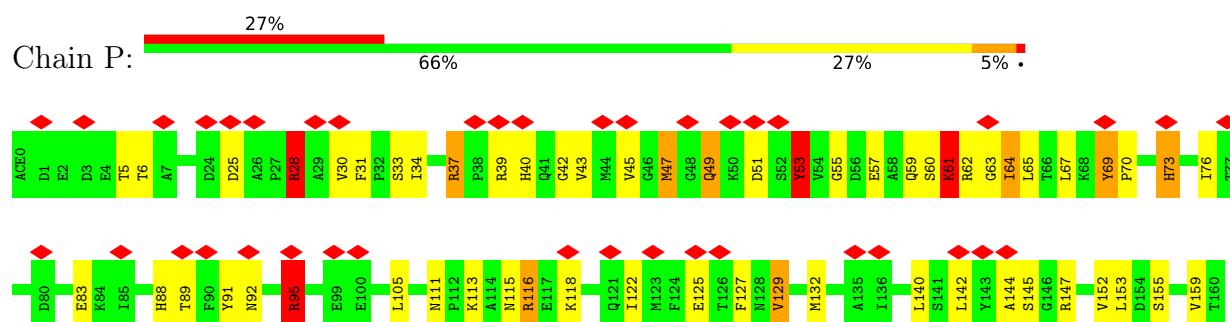
### 3 Residue-property plots

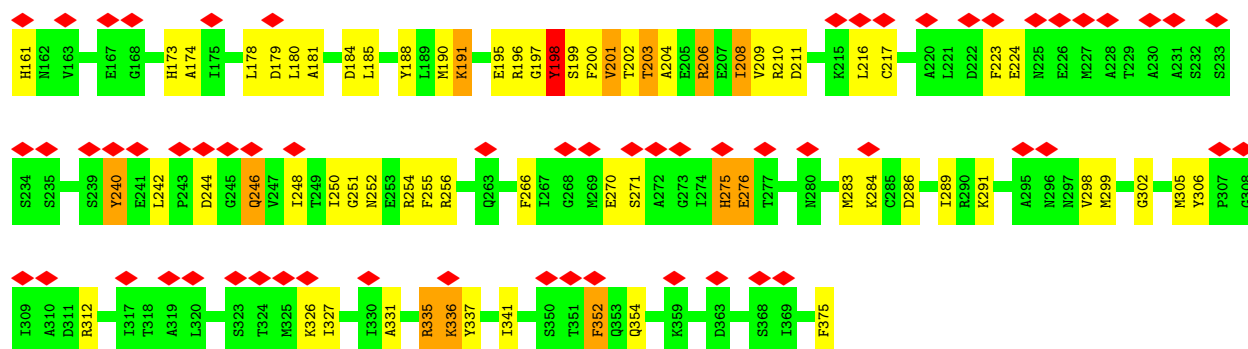
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

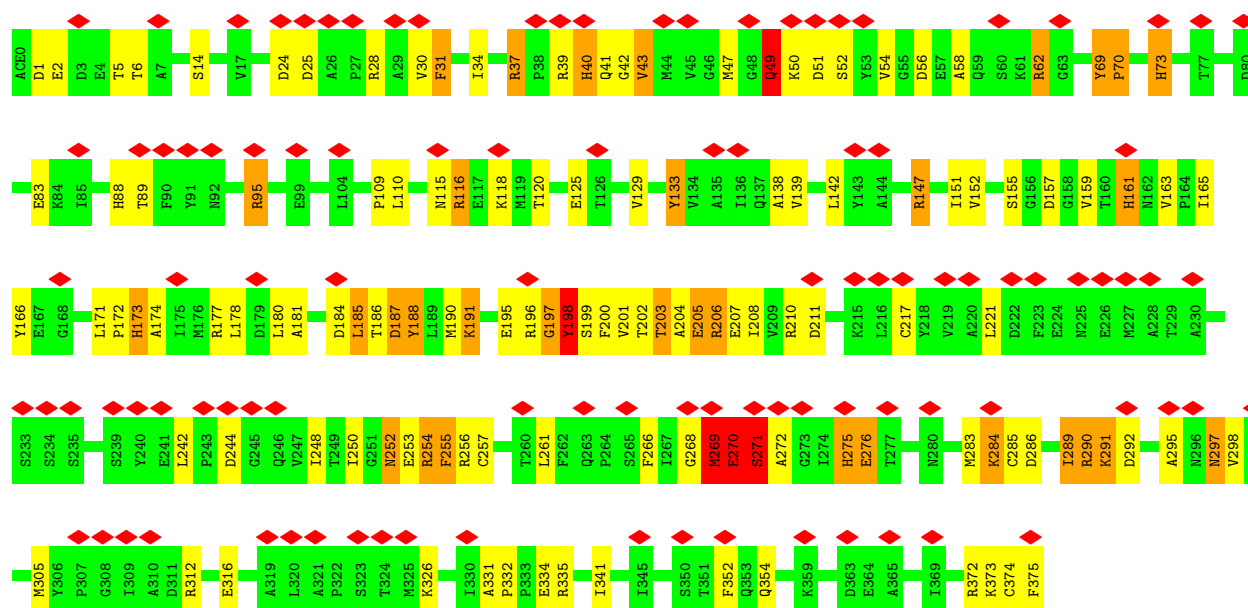


- Molecule 1: Actin, alpha skeletal muscle

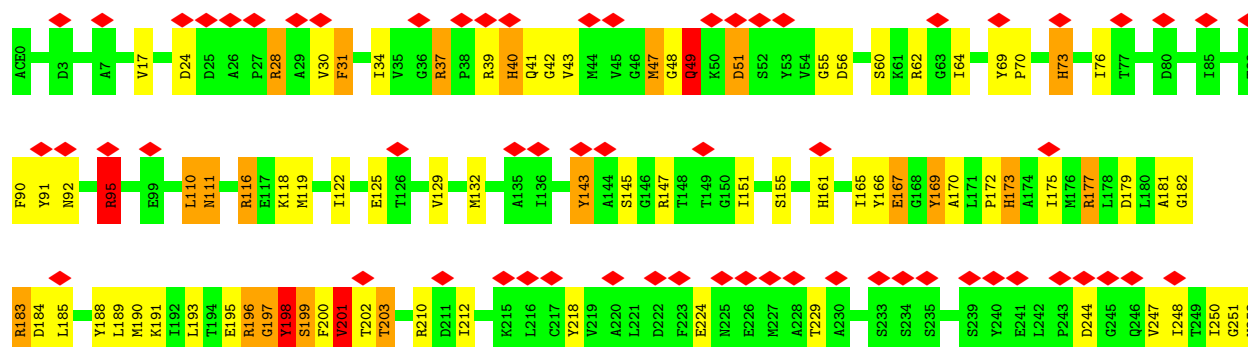


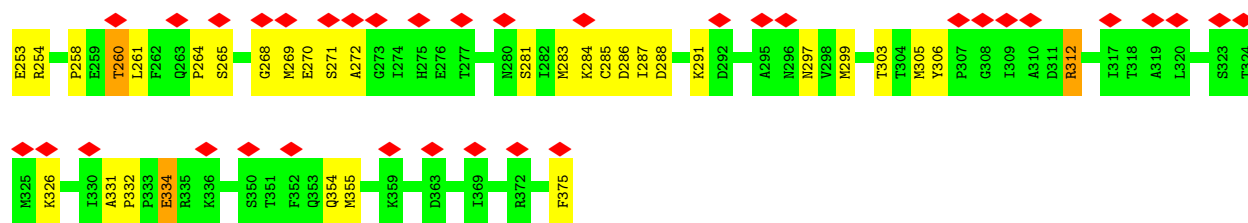


• Molecule 1: Actin, alpha skeletal muscle

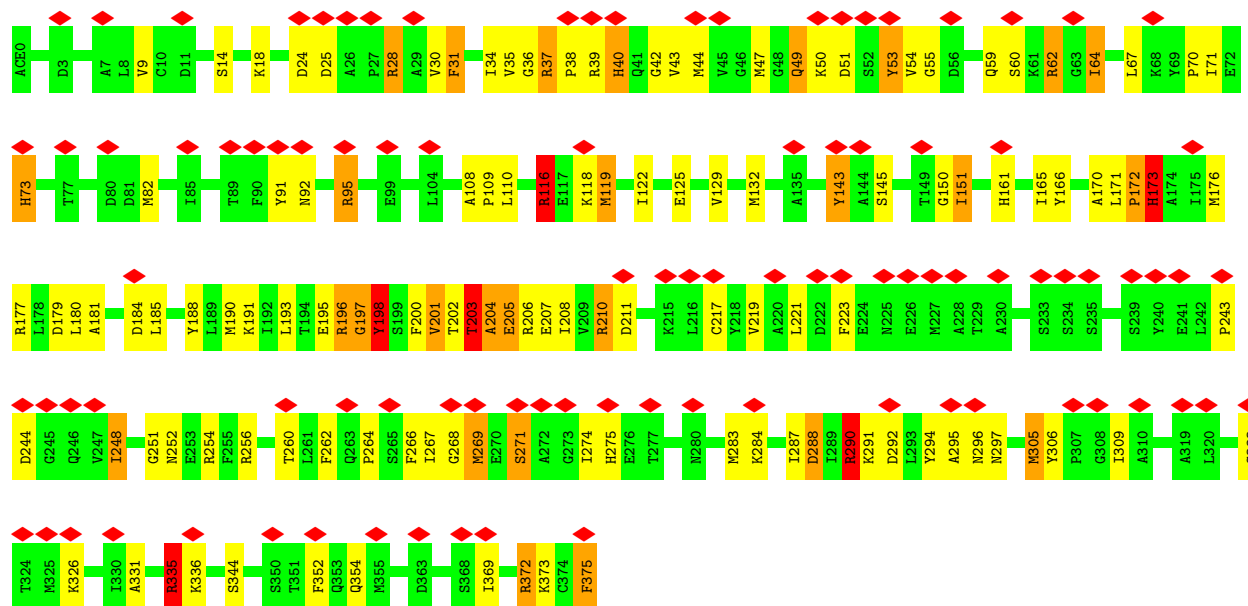


• Molecule 1: Actin, alpha skeletal muscle

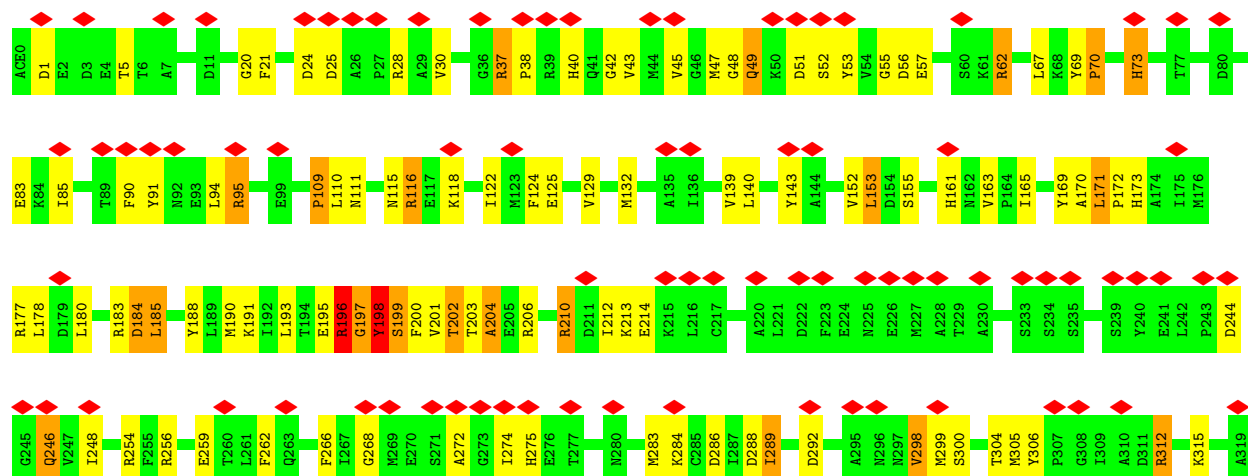


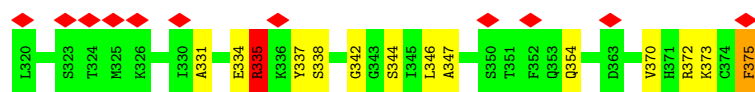


• Molecule 1: Actin, alpha skeletal muscle

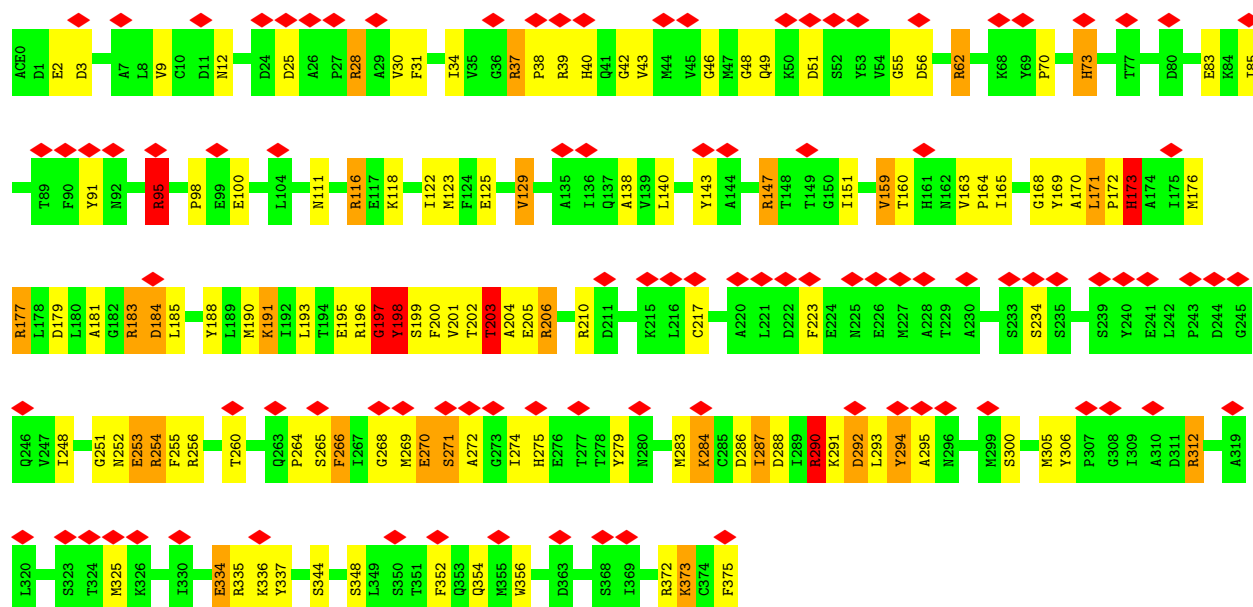


• Molecule 1: Actin, alpha skeletal muscle

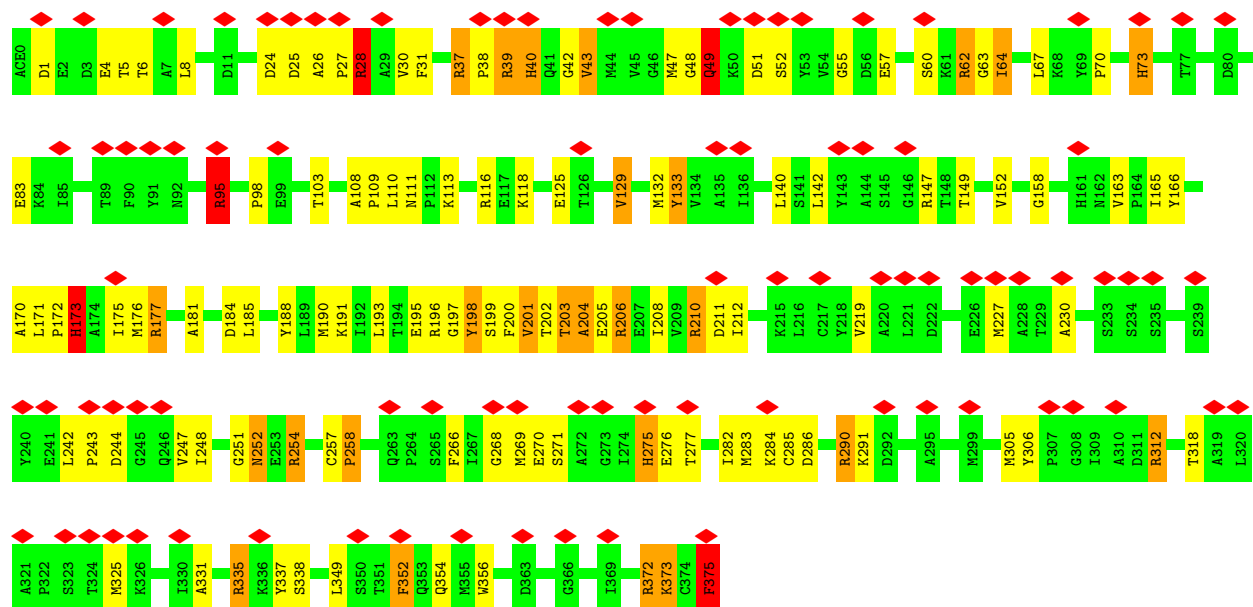




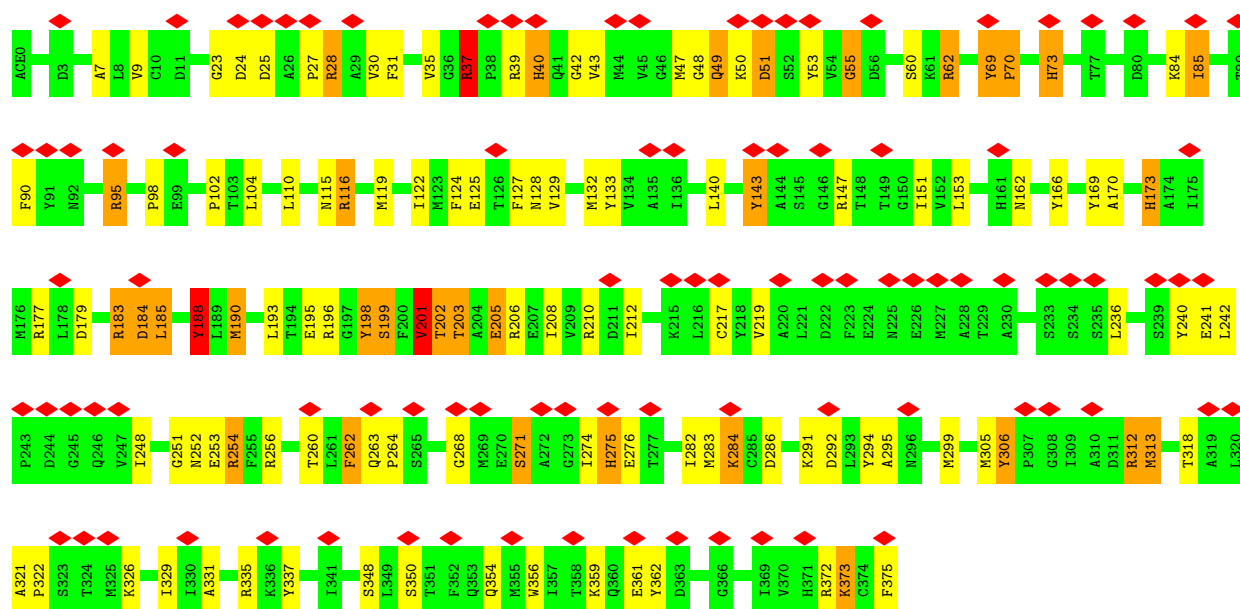
- Molecule 1: Actin, alpha skeletal muscle



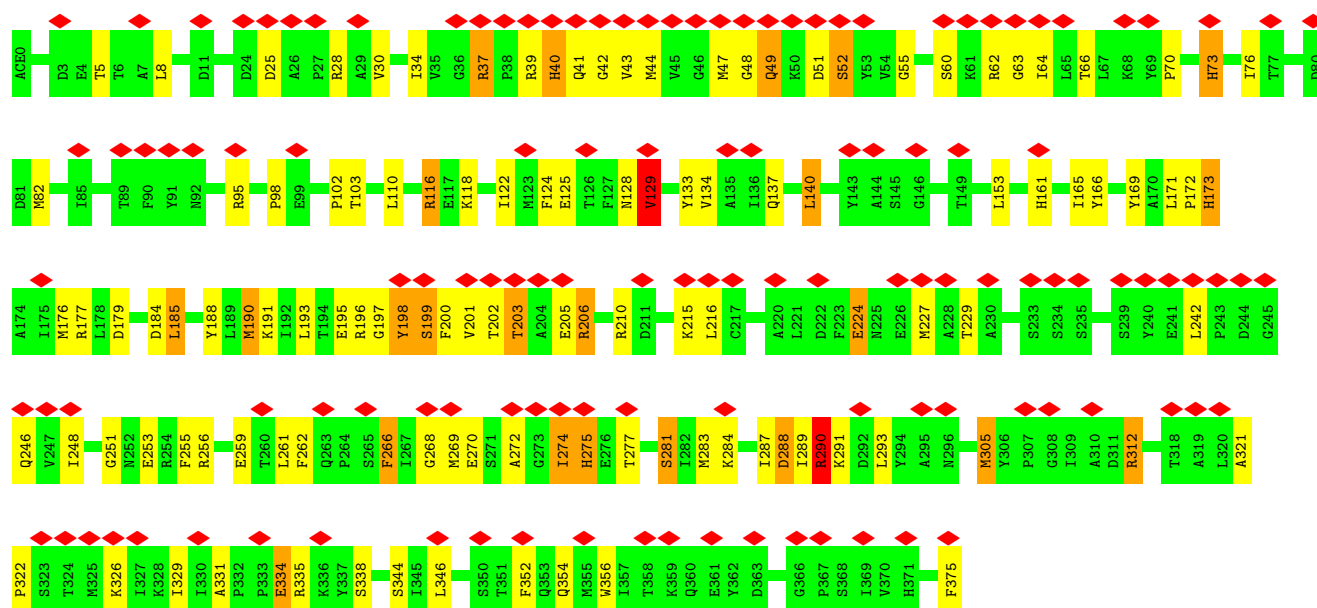
- Molecule 1: Actin, alpha skeletal muscle



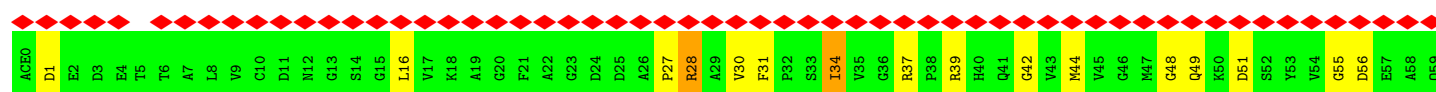
- Molecule 1: Actin, alpha skeletal muscle

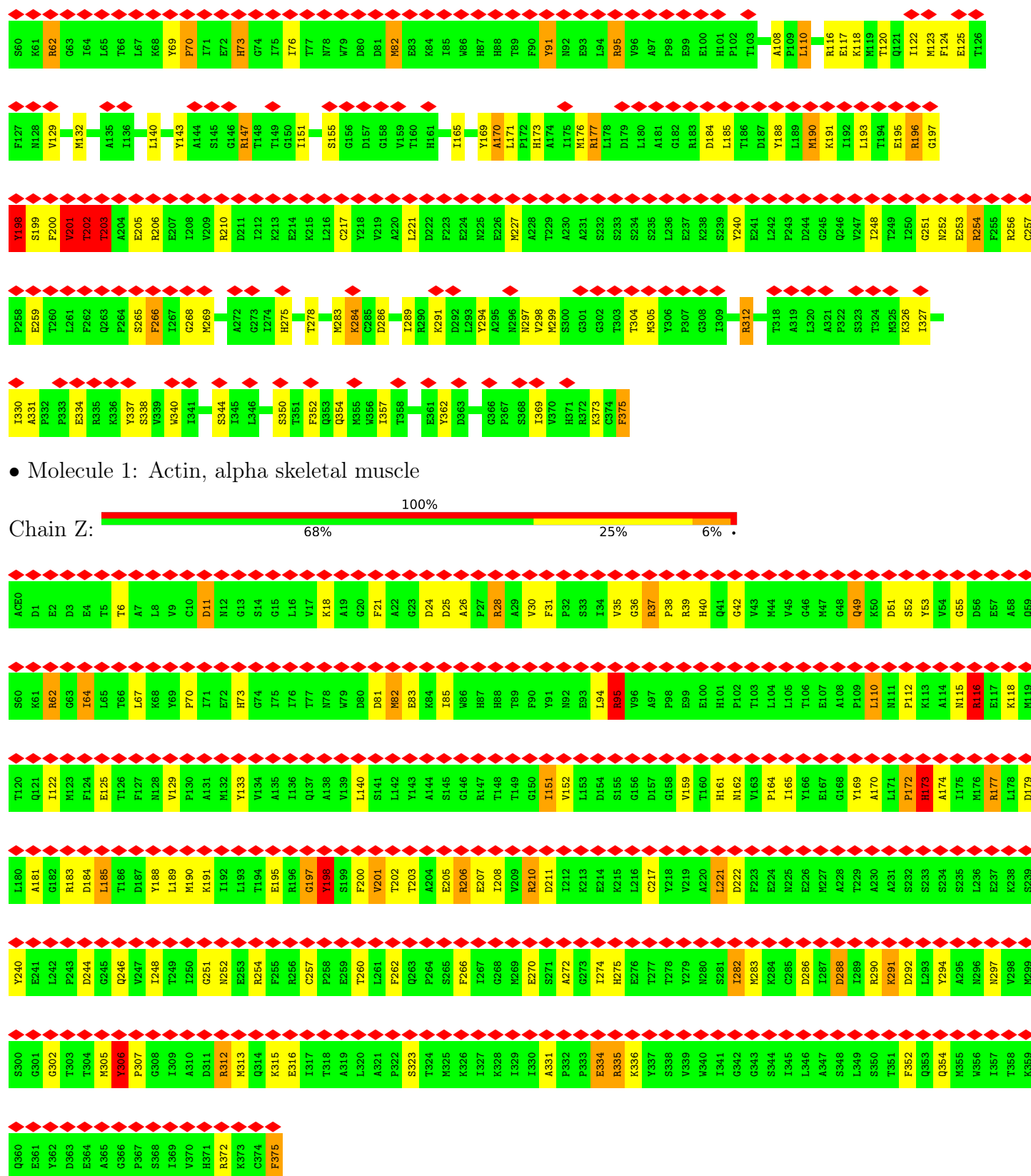


• Molecule 1: Actin, alpha skeletal muscle



• Molecule 1: Actin, alpha skeletal muscle





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	HELICAL, twist=Not provided°, rise=Not provided Å, axial sym=Not provided	Depositor
Number of particles used	8000	Depositor
Resolution determination method	Not provided	
CTF correction method	FSC at 0.143 cut-off	Depositor
Microscope	HITACHI EF2000	Depositor
Voltage (kV)	200	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	15	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	100000	Depositor
Image detector	GENERIC CCD	Depositor
Maximum map value	105.405	Depositor
Minimum map value	-66.216	Depositor
Average map value	0.454	Depositor
Map value standard deviation	12.051	Depositor
Recommended contour level	33	Depositor
Map size (Å)	275.275, 275.275, 275.275	wwPDB
Map dimensions	121, 121, 121	wwPDB
Map angles (°)	90, 90, 90	wwPDB
Pixel spacing (Å)	2.275, 2.275, 2.275	Depositor



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HIC, ACE, PO4, ADP, MG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	O	0.73	0/2984	1.47	30/4042 (0.7%)
1	P	0.91	9/2984 (0.3%)	1.48	36/4042 (0.9%)
1	Q	0.80	4/2984 (0.1%)	1.52	31/4042 (0.8%)
1	R	0.87	3/2984 (0.1%)	1.48	27/4042 (0.7%)
1	S	0.74	1/2984 (0.0%)	1.49	30/4042 (0.7%)
1	T	0.74	0/2984	1.46	23/4042 (0.6%)
1	U	0.73	0/2984	1.44	29/4042 (0.7%)
1	V	0.71	0/2984	1.45	34/4042 (0.8%)
1	W	0.73	0/2984	1.48	34/4042 (0.8%)
1	X	0.72	0/2984	1.45	29/4042 (0.7%)
1	Y	0.72	0/2984	1.40	27/4042 (0.7%)
1	Z	0.72	0/2984	1.44	27/4042 (0.7%)
All	All	0.76	17/35808 (0.0%)	1.46	357/48504 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	O	0	18
1	P	0	17
1	Q	0	25
1	R	0	16
1	S	0	21
1	T	0	19
1	U	0	25
1	V	0	19
1	W	0	24
1	X	0	14
1	Y	0	22

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Mol	Chain	#Chirality outliers	#Planarity outliers
1	Z	0	23
All	All	0	243

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	R	167	GLU	CD-OE1	-17.54	1.06	1.25
1	R	167	GLU	CD-OE2	12.87	1.39	1.25
1	P	62	ARG	C-N	12.41	1.55	1.33
1	P	61	LYS	CG-CD	10.97	1.89	1.52
1	Q	187	ASP	CG-OD2	-9.47	1.03	1.25

The worst 5 of 357 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	269	MET	CG-SD-CE	-16.11	74.43	100.20
1	S	288	ASP	CB-CG-OD1	14.03	130.92	118.30
1	R	166	TYR	CB-CG-CD1	-13.80	112.72	121.00
1	Q	187	ASP	CB-CG-OD2	12.64	129.68	118.30
1	R	166	TYR	CB-CG-CD2	12.48	128.49	121.00

There are no chirality outliers.

5 of 243 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	O	28	ARG	Sidechain
1	O	38	PRO	Peptide
1	O	50	LYS	Peptide
1	O	69	TYR	Sidechain
1	O	91	TYR	Sidechain

## 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	O	2936	0	2896	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	P	2936	0	2895	38	0
1	Q	2936	0	2896	52	0
1	R	2936	0	2896	35	0
1	S	2936	0	2896	40	0
1	T	2936	0	2896	42	0
1	U	2936	0	2896	37	0
1	V	2936	0	2896	48	0
1	W	2936	0	2896	41	0
1	X	2936	0	2896	37	0
1	Y	2936	0	2896	33	0
1	Z	2936	0	2896	23	0
2	O	27	0	12	1	0
2	P	27	0	12	0	0
2	Q	27	0	12	0	0
2	R	27	0	12	0	0
2	S	27	0	12	0	0
2	T	27	0	12	0	0
2	U	27	0	12	1	0
2	V	27	0	12	0	0
2	W	27	0	12	0	0
2	X	27	0	12	0	0
2	Y	27	0	12	0	0
2	Z	27	0	12	0	0
3	O	15	0	0	0	0
3	P	15	0	0	1	0
3	Q	15	0	0	0	0
3	R	15	0	0	1	0
3	S	15	0	0	0	0
3	T	20	0	0	1	0
3	U	15	0	0	0	0
3	V	10	0	0	0	0
3	W	15	0	0	0	0
3	X	15	0	0	0	0
3	Y	15	0	0	2	0
3	Z	15	0	0	0	0
4	O	6	0	0	0	0
4	P	5	0	0	0	0
4	Q	7	0	0	0	0
4	R	6	0	0	0	0
4	S	6	0	0	0	0
4	T	6	0	0	0	0
4	U	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	V	6	0	0	0	0
4	W	6	0	0	0	0
4	X	6	0	0	0	0
4	Y	6	0	0	0	0
4	Z	6	0	0	0	0
All	All	35808	0	34895	391	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 391 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)
1:P:61:LYS:CD	1:P:61:LYS:CG	1.89	1.50
1:Q:202:THR:HG23	1:Q:203:THR:H	1.48	0.78
1:P:202:THR:HG23	1:P:203:THR:H	1.50	0.76
1:O:201:VAL:HG13	1:O:202:THR:HG22	1.68	0.74
1:V:275:HIS:CD2	1:V:276:GLU:H	2.05	0.74

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	O	373/376 (99%)	283 (76%)	66 (18%)	24 (6%)	1	13
1	P	373/376 (99%)	298 (80%)	54 (14%)	21 (6%)	1	14
1	Q	373/376 (99%)	281 (75%)	66 (18%)	26 (7%)	1	11
1	R	373/376 (99%)	295 (79%)	52 (14%)	26 (7%)	1	11
1	S	373/376 (99%)	286 (77%)	67 (18%)	20 (5%)	1	15

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	T	373/376 (99%)	275 (74%)	73 (20%)	25 (7%)	1	12
1	U	373/376 (99%)	284 (76%)	66 (18%)	23 (6%)	1	13
1	V	373/376 (99%)	293 (79%)	56 (15%)	24 (6%)	1	13
1	W	373/376 (99%)	288 (77%)	60 (16%)	25 (7%)	1	12
1	X	373/376 (99%)	292 (78%)	63 (17%)	18 (5%)	2	16
1	Y	373/376 (99%)	292 (78%)	64 (17%)	17 (5%)	2	17
1	Z	373/376 (99%)	289 (78%)	59 (16%)	25 (7%)	1	12
All	All	4476/4512 (99%)	3456 (77%)	746 (17%)	274 (6%)	2	13

5 of 274 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	O	42	GLY
1	O	51	ASP
1	O	129	VAL
1	O	254	ARG
1	P	42	GLY

### 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	O	317/317 (100%)	272 (86%)	45 (14%)	2	12
1	P	317/317 (100%)	269 (85%)	48 (15%)	2	11
1	Q	317/317 (100%)	269 (85%)	48 (15%)	2	11
1	R	317/317 (100%)	275 (87%)	42 (13%)	3	14
1	S	317/317 (100%)	265 (84%)	52 (16%)	2	10
1	T	317/317 (100%)	275 (87%)	42 (13%)	3	14
1	U	317/317 (100%)	265 (84%)	52 (16%)	2	10
1	V	317/317 (100%)	269 (85%)	48 (15%)	2	11
1	W	317/317 (100%)	279 (88%)	38 (12%)	4	16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	X	317/317 (100%)	266 (84%)	51 (16%)	2	10
1	Y	317/317 (100%)	276 (87%)	41 (13%)	3	14
1	Z	317/317 (100%)	259 (82%)	58 (18%)	1	8
All	All	3804/3804 (100%)	3239 (85%)	565 (15%)	5	11

5 of 565 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	344	SER
1	Y	116	ARG
1	X	338	SER
1	Z	85	ILE
1	S	59	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
1	Z	111	ASN
1	Z	246	GLN
1	V	101	HIS
1	U	162	ASN
1	Z	252	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

12 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	HIC	W	73	1	8,11,12	2.84	1 (12%)	6,14,16	3.79	2 (33%)
1	HIC	Y	73	1	8,11,12	2.65	1 (12%)	6,14,16	2.72	2 (33%)
1	HIC	S	73	1	8,11,12	2.90	1 (12%)	6,14,16	3.13	2 (33%)
1	HIC	V	73	1	8,11,12	2.74	1 (12%)	6,14,16	2.53	2 (33%)
1	HIC	T	73	1	8,11,12	2.42	1 (12%)	6,14,16	2.22	2 (33%)
1	HIC	X	73	1	8,11,12	2.83	1 (12%)	6,14,16	2.21	3 (50%)
1	HIC	Z	73	1	8,11,12	2.91	1 (12%)	6,14,16	4.00	4 (66%)
1	HIC	P	73	1	8,11,12	2.46	1 (12%)	6,14,16	2.39	2 (33%)
1	HIC	O	73	1	8,11,12	1.47	1 (12%)	6,14,16	1.05	0
1	HIC	R	73	1	8,11,12	2.59	1 (12%)	6,14,16	2.67	2 (33%)
1	HIC	Q	73	1	8,11,12	2.60	1 (12%)	6,14,16	2.49	2 (33%)
1	HIC	U	73	1	8,11,12	2.54	1 (12%)	6,14,16	1.51	2 (33%)

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
1	HIC	W	73	1	-	2/5/6/8	0/1/1/1
1	HIC	Y	73	1	-	1/5/6/8	0/1/1/1
1	HIC	S	73	1	-	0/5/6/8	0/1/1/1
1	HIC	V	73	1	-	2/5/6/8	0/1/1/1
1	HIC	T	73	1	-	0/5/6/8	0/1/1/1
1	HIC	X	73	1	-	1/5/6/8	0/1/1/1
1	HIC	Z	73	1	-	1/5/6/8	0/1/1/1
1	HIC	P	73	1	-	1/5/6/8	0/1/1/1
1	HIC	O	73	1	-	1/5/6/8	0/1/1/1
1	HIC	R	73	1	-	1/5/6/8	0/1/1/1
1	HIC	Q	73	1	-	1/5/6/8	0/1/1/1
1	HIC	U	73	1	-	0/5/6/8	0/1/1/1

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	S	73	HIC	CZ-NE2	-7.80	1.26	1.48
1	Z	73	HIC	CZ-NE2	-7.78	1.27	1.48
1	W	73	HIC	CZ-NE2	-7.71	1.27	1.48
1	X	73	HIC	CZ-NE2	-7.64	1.27	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	V	73	HIC	CZ-NE2	-7.48	1.27	1.48

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	W	73	HIC	CZ-NE2-CD2	7.38	151.91	126.31
1	Z	73	HIC	CZ-NE2-CD2	6.95	150.43	126.31
1	S	73	HIC	CZ-NE2-CD2	5.97	147.05	126.31
1	Y	73	HIC	CZ-NE2-CD2	5.31	144.73	126.31
1	W	73	HIC	CZ-NE2-CE1	-5.11	100.52	125.48

There are no chirality outliers.

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	P	73	HIC	CA-CB-CG-ND1
1	Q	73	HIC	CA-CB-CG-ND1
1	R	73	HIC	CA-CB-CG-ND1
1	V	73	HIC	O-C-CA-CB
1	V	73	HIC	CA-CB-CG-ND1

There are no ring outliers.

11 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	W	73	HIC	3	0
1	Y	73	HIC	2	0
1	S	73	HIC	3	0
1	V	73	HIC	3	0
1	T	73	HIC	2	0
1	X	73	HIC	3	0
1	P	73	HIC	1	0
1	O	73	HIC	2	0
1	R	73	HIC	3	0
1	Q	73	HIC	3	0
1	U	73	HIC	2	0

## 5.5 Carbohydrates

There are no oligosaccharides in this entry.



## 5.6 Ligand geometry ⓘ

Of 120 ligands modelled in this entry, 72 are monoatomic - leaving 48 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$
3	PO4	S	802	-	4,4,4	1.84	1 (25%)	6,6,6	1.94	1 (16%)
3	PO4	S	803	-	4,4,4	1.89	1 (25%)	6,6,6	3.09	3 (50%)
3	PO4	R	801	-	4,4,4	1.87	1 (25%)	6,6,6	2.74	2 (33%)
3	PO4	W	801	-	4,4,4	1.69	1 (25%)	6,6,6	3.13	2 (33%)
3	PO4	O	802	-	4,4,4	1.74	1 (25%)	6,6,6	2.35	2 (33%)
3	PO4	P	801	-	4,4,4	1.94	1 (25%)	6,6,6	3.10	2 (33%)
2	ADP	Y	800	4	24,29,29	0.88	0	29,45,45	1.50	4 (13%)
2	ADP	V	800	4	24,29,29	0.94	1 (4%)	29,45,45	1.54	3 (10%)
3	PO4	T	801	4	4,4,4	1.73	1 (25%)	6,6,6	3.20	2 (33%)
3	PO4	T	376	-	4,4,4	1.66	1 (25%)	6,6,6	3.18	2 (33%)
3	PO4	V	802	-	4,4,4	1.81	1 (25%)	6,6,6	2.53	2 (33%)
3	PO4	W	802	-	4,4,4	1.60	1 (25%)	6,6,6	2.07	2 (33%)
3	PO4	O	803	-	4,4,4	2.03	1 (25%)	6,6,6	3.05	4 (66%)
3	PO4	Q	802	-	4,4,4	1.80	1 (25%)	6,6,6	2.14	2 (33%)
3	PO4	R	802	-	4,4,4	1.49	1 (25%)	6,6,6	2.25	1 (16%)
3	PO4	R	803	-	4,4,4	2.01	1 (25%)	6,6,6	5.65	5 (83%)
3	PO4	U	802	-	4,4,4	1.38	1 (25%)	6,6,6	2.44	3 (50%)
2	ADP	R	800	4	24,29,29	1.02	1 (4%)	29,45,45	2.00	6 (20%)
3	PO4	Y	801	4	4,4,4	1.98	1 (25%)	6,6,6	2.02	3 (50%)
3	PO4	P	803	-	4,4,4	3.04	1 (25%)	6,6,6	6.67	5 (83%)
3	PO4	Z	803	-	4,4,4	1.73	1 (25%)	6,6,6	2.73	2 (33%)
2	ADP	P	800	4	24,29,29	0.97	1 (4%)	29,45,45	1.39	4 (13%)
3	PO4	P	802	-	4,4,4	1.51	1 (25%)	6,6,6	1.81	2 (33%)
3	PO4	Z	802	-	4,4,4	1.88	1 (25%)	6,6,6	3.19	3 (50%)
3	PO4	U	803	-	4,4,4	1.75	1 (25%)	6,6,6	2.08	2 (33%)
2	ADP	U	800	4	24,29,29	0.84	0	29,45,45	1.85	5 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	PO4	Y	802	-	4,4,4	1.85	1 (25%)	6,6,6	1.93	2 (33%)
3	PO4	Q	801	4	4,4,4	1.78	1 (25%)	6,6,6	4.15	2 (33%)
3	PO4	T	803	-	4,4,4	1.85	1 (25%)	6,6,6	3.91	3 (50%)
3	PO4	V	803	-	4,4,4	1.80	1 (25%)	6,6,6	3.12	3 (50%)
3	PO4	Y	803	-	4,4,4	1.89	1 (25%)	6,6,6	3.91	4 (66%)
2	ADP	Z	800	4	24,29,29	0.89	0	29,45,45	1.53	5 (17%)
2	ADP	X	800	4	24,29,29	0.94	1 (4%)	29,45,45	1.81	5 (17%)
3	PO4	W	803	-	4,4,4	1.86	1 (25%)	6,6,6	3.33	2 (33%)
2	ADP	W	800	4	24,29,29	0.93	1 (4%)	29,45,45	1.56	5 (17%)
2	ADP	S	800	4	24,29,29	0.93	0	29,45,45	1.67	4 (13%)
3	PO4	Z	801	-	4,4,4	1.84	1 (25%)	6,6,6	3.26	2 (33%)
3	PO4	U	801	-	4,4,4	3.14	1 (25%)	6,6,6	5.60	2 (33%)
2	ADP	O	800	4	24,29,29	1.02	1 (4%)	29,45,45	1.75	6 (20%)
3	PO4	X	803	-	4,4,4	1.76	1 (25%)	6,6,6	4.22	5 (83%)
3	PO4	X	801	-	4,4,4	1.52	1 (25%)	6,6,6	1.97	2 (33%)
2	ADP	Q	800	4	24,29,29	0.95	0	29,45,45	1.74	5 (17%)
2	ADP	T	800	4	24,29,29	0.84	0	29,45,45	1.96	6 (20%)
3	PO4	Q	803	-	4,4,4	1.74	1 (25%)	6,6,6	2.93	2 (33%)
3	PO4	O	801	-	4,4,4	1.69	1 (25%)	6,6,6	3.05	2 (33%)
3	PO4	T	802	-	4,4,4	1.83	1 (25%)	6,6,6	2.03	1 (16%)
3	PO4	X	802	-	4,4,4	1.71	1 (25%)	6,6,6	2.40	2 (33%)
3	PO4	S	801	-	4,4,4	1.37	1 (25%)	6,6,6	2.13	2 (33%)

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
2	ADP	R	800	4	-	0/12/32/32	0/3/3/3
2	ADP	O	800	4	-	0/12/32/32	0/3/3/3
2	ADP	Z	800	4	-	0/12/32/32	0/3/3/3
2	ADP	X	800	4	-	0/12/32/32	0/3/3/3
2	ADP	Q	800	4	-	2/12/32/32	0/3/3/3
2	ADP	P	800	4	-	0/12/32/32	0/3/3/3
2	ADP	T	800	4	-	0/12/32/32	0/3/3/3
2	ADP	W	800	4	-	1/12/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	Y	800	4	-	0/12/32/32	0/3/3/3
2	ADP	S	800	4	-	0/12/32/32	0/3/3/3
2	ADP	V	800	4	-	1/12/32/32	0/3/3/3
2	ADP	U	800	4	-	0/12/32/32	0/3/3/3

The worst 5 of 42 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	U	801	PO4	P-O1	5.67	1.64	1.50
3	P	803	PO4	P-O1	5.50	1.63	1.50
3	O	803	PO4	P-O4	3.50	1.65	1.54
3	Z	802	PO4	P-O4	3.49	1.65	1.54
3	P	801	PO4	P-O4	3.47	1.65	1.54

The worst 5 of 146 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	803	PO4	O4-P-O1	-12.86	63.84	110.89
3	U	801	PO4	O4-P-O1	-11.29	69.57	110.89
3	R	803	PO4	O4-P-O1	-9.65	75.57	110.89
3	U	801	PO4	O3-P-O2	7.70	132.67	107.97
3	X	803	PO4	O4-P-O1	-7.55	83.28	110.89

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	Q	800	ADP	C5'-O5'-PA-O3A
2	Q	800	ADP	C5'-O5'-PA-O2A
2	W	800	ADP	C3'-C4'-C5'-O5'
2	V	800	ADP	C5'-O5'-PA-O3A

There are no ring outliers.

6 monomers are involved in 7 short contacts:

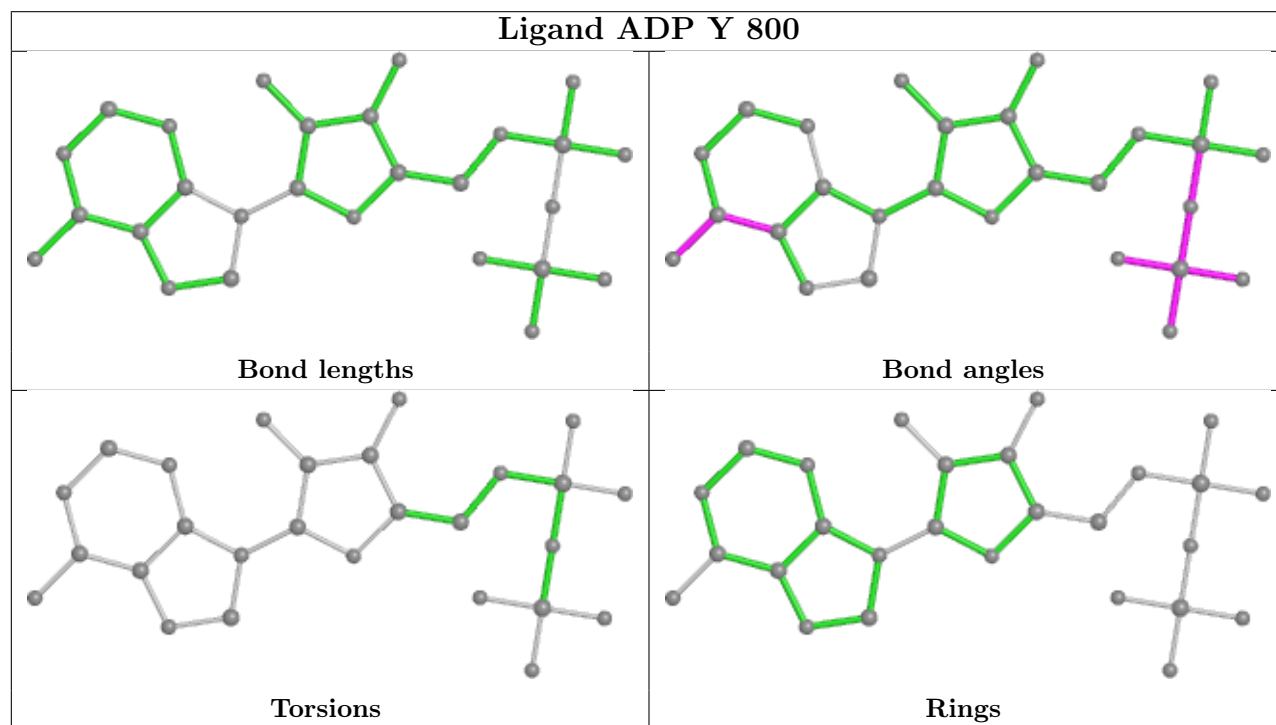
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	R	801	PO4	1	0
3	Y	801	PO4	2	0
3	P	803	PO4	1	0
2	U	800	ADP	1	0
3	T	803	PO4	1	0

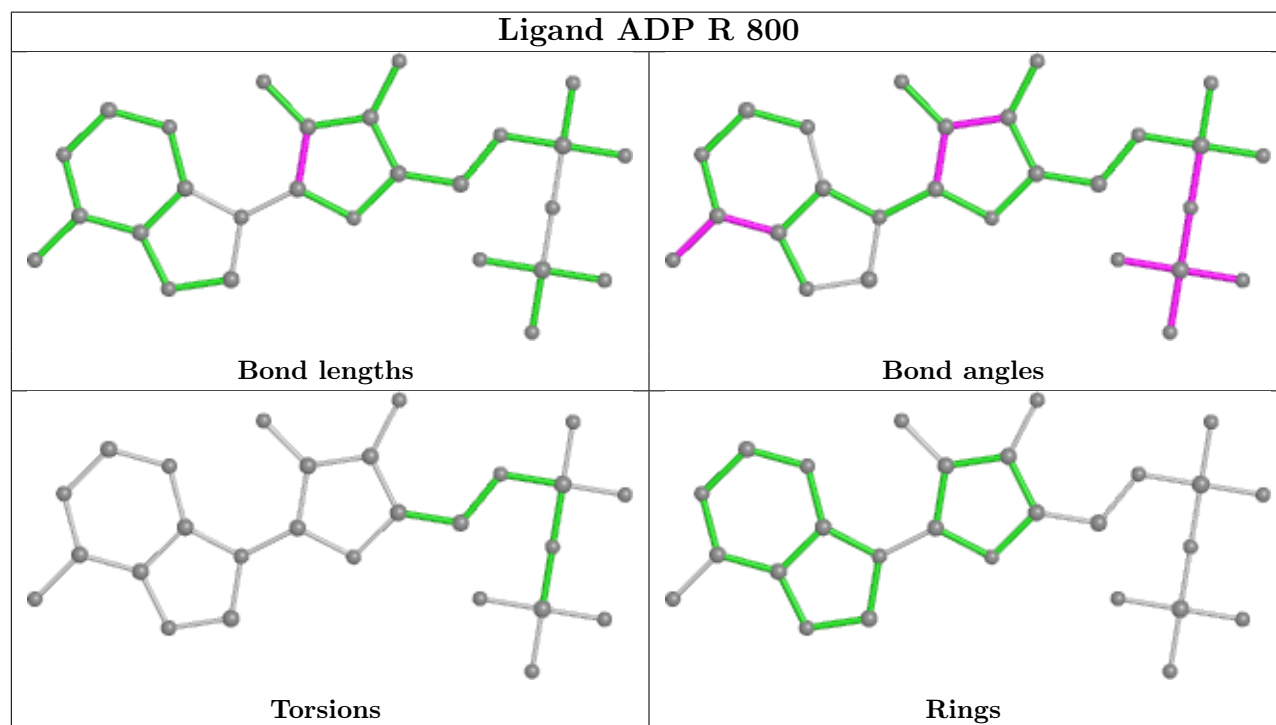
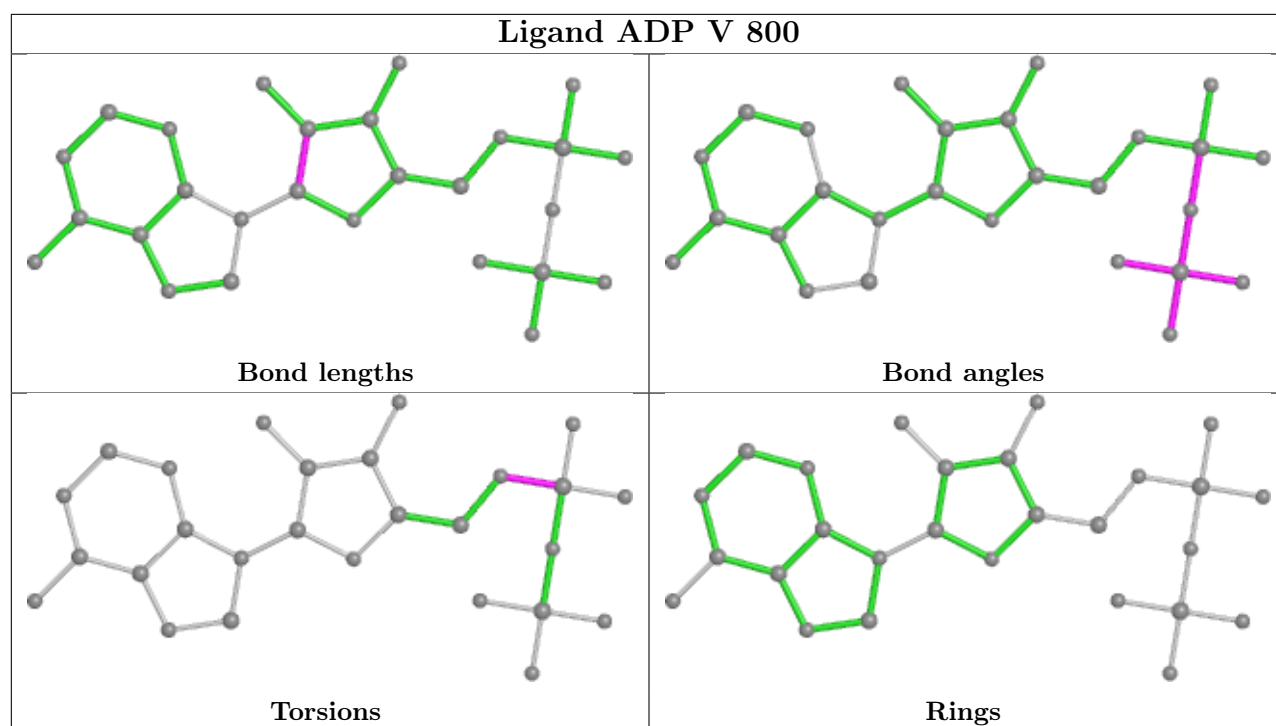
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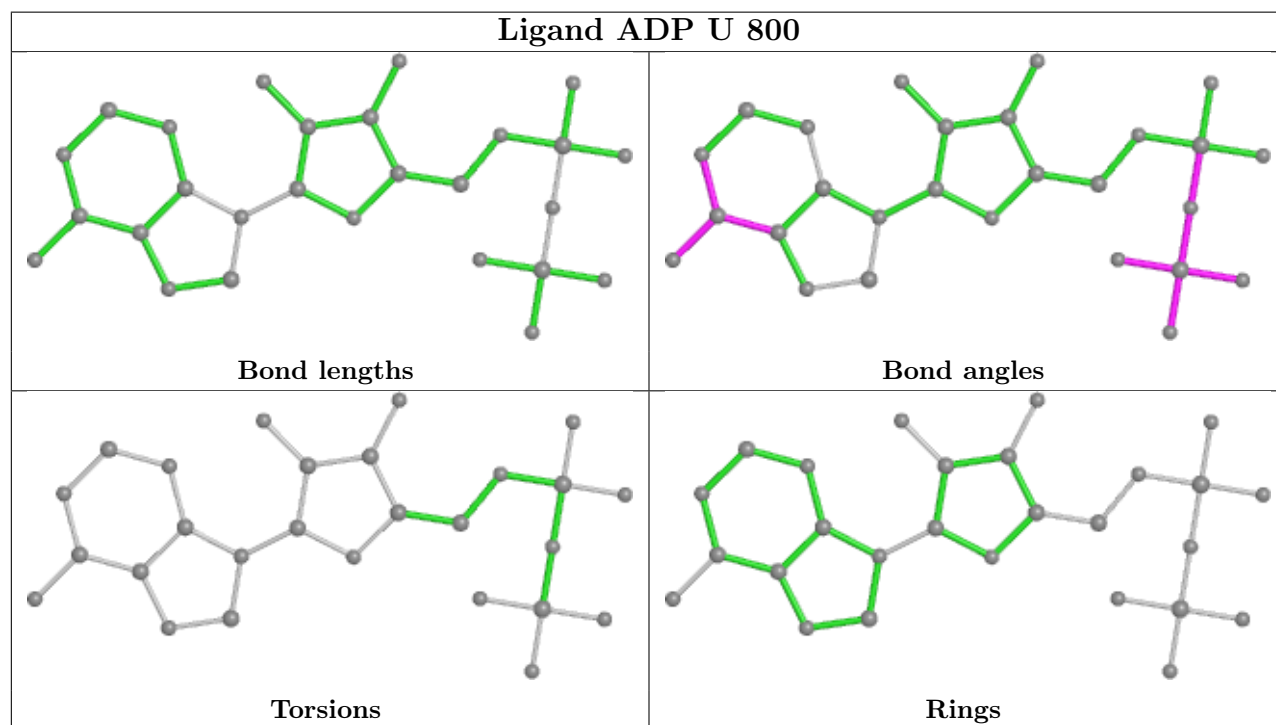
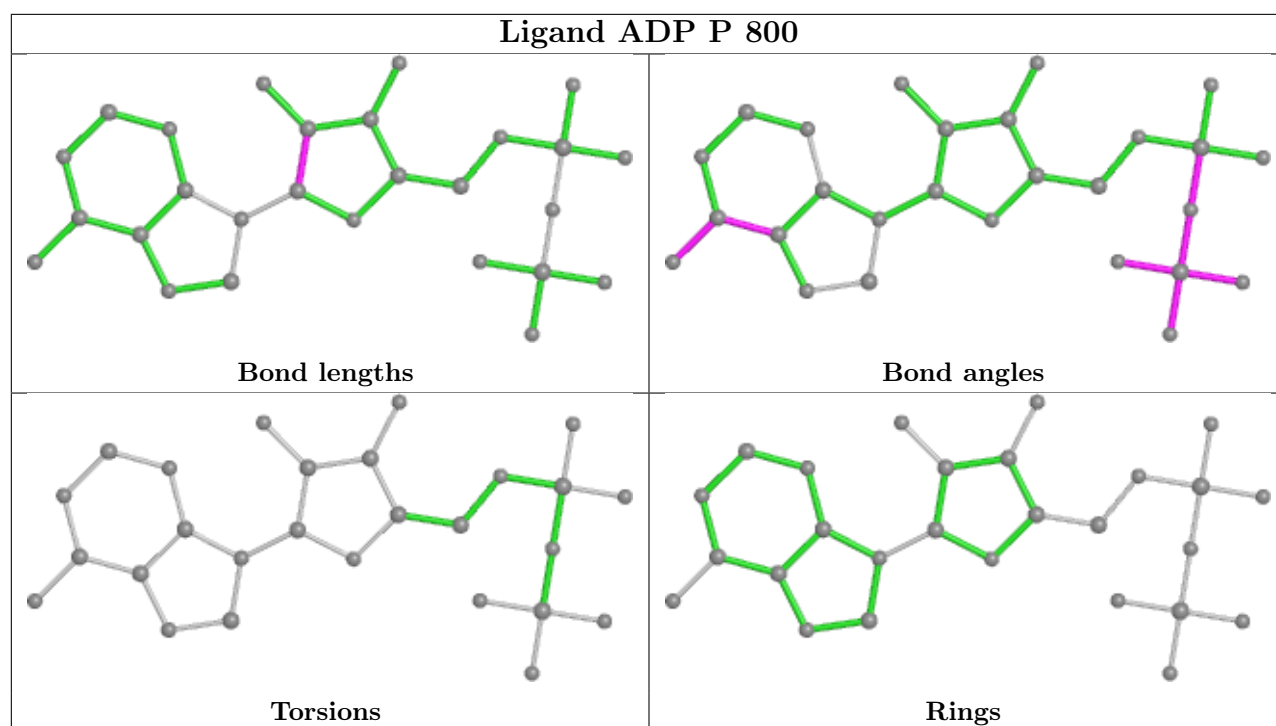
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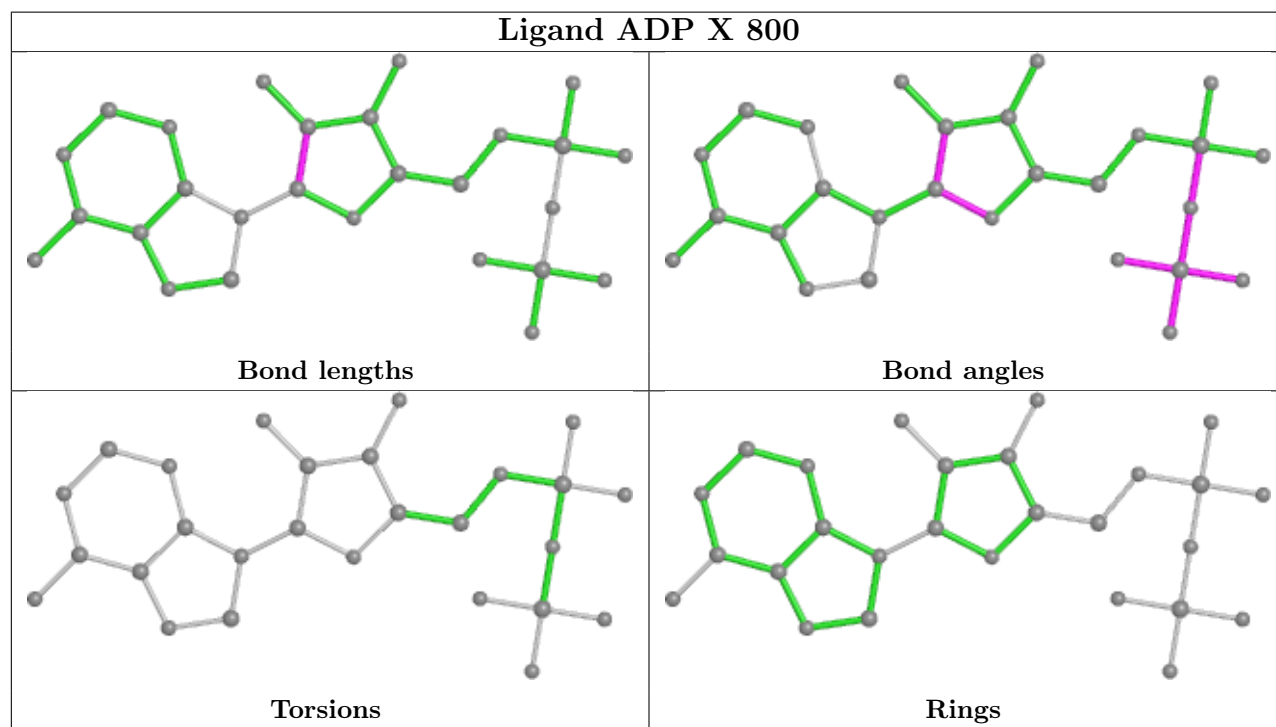
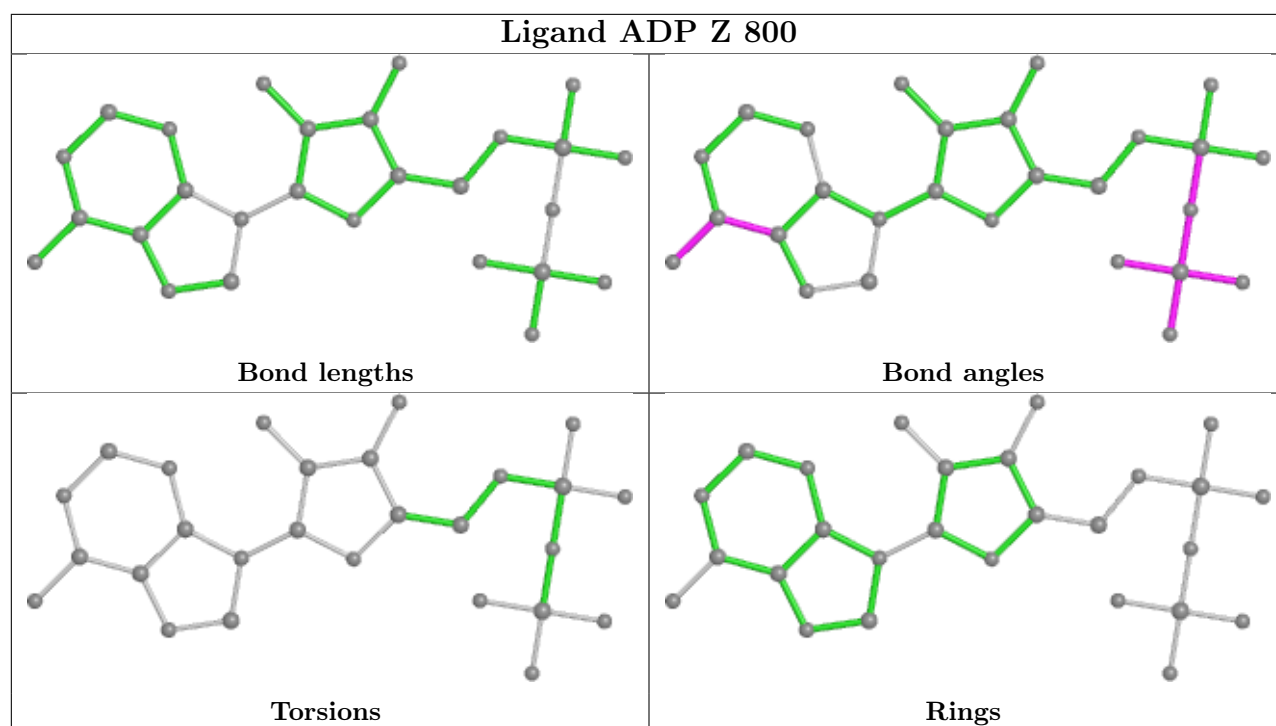
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	O	800	ADP	1	0

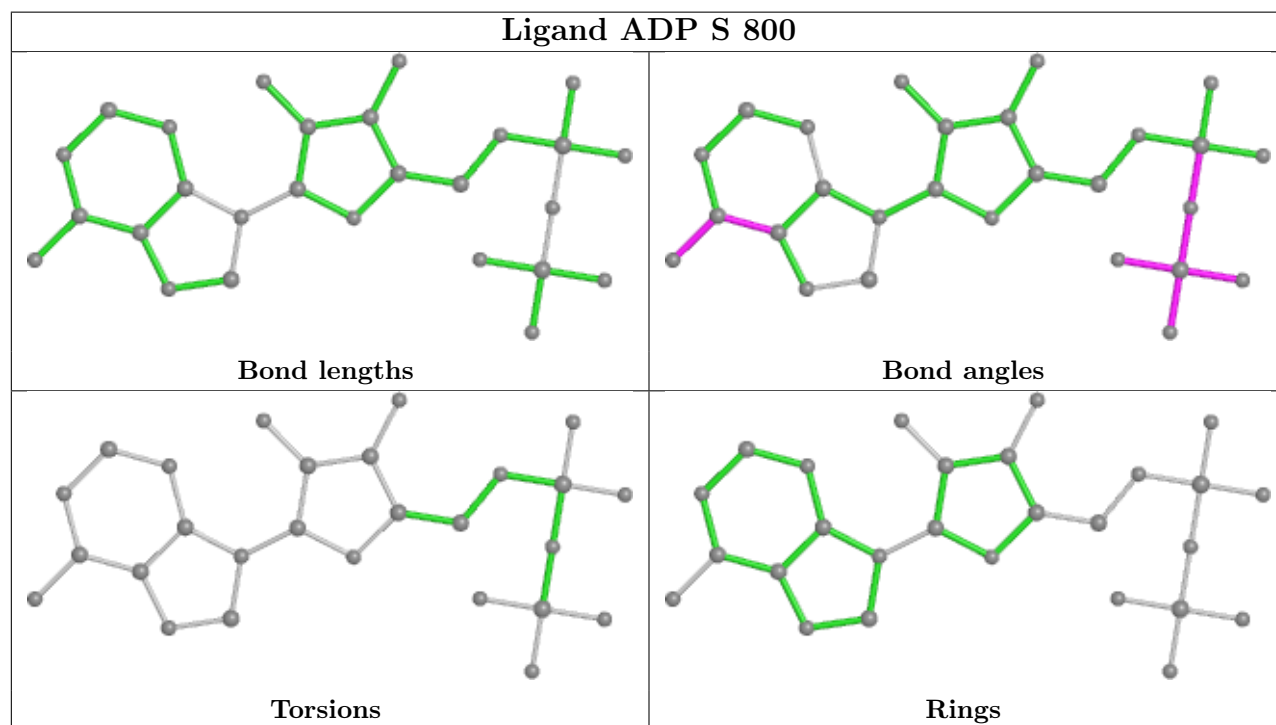
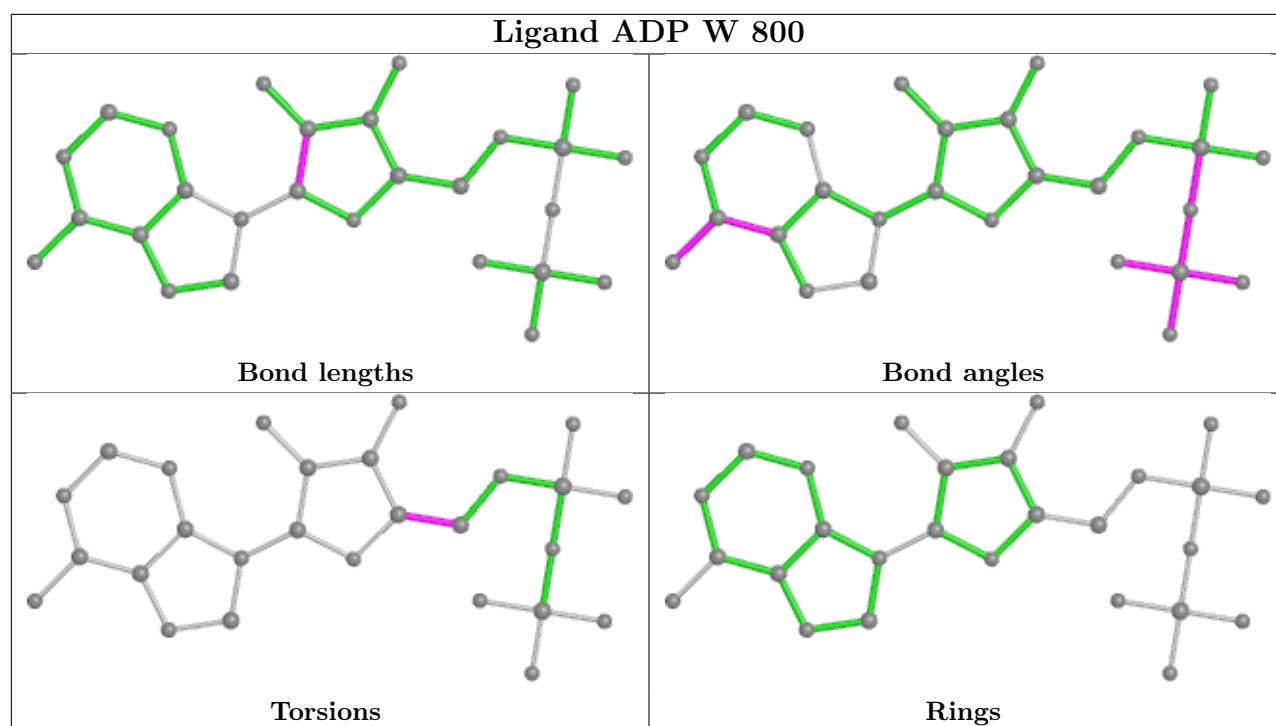
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



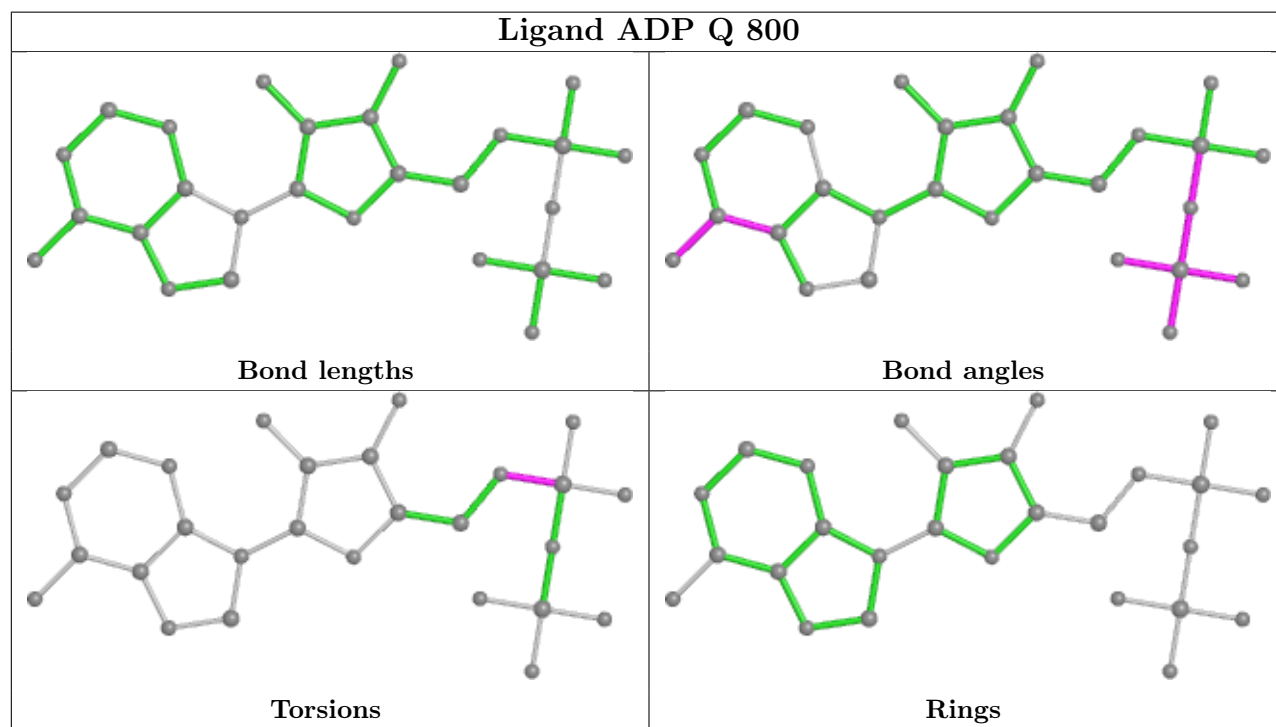
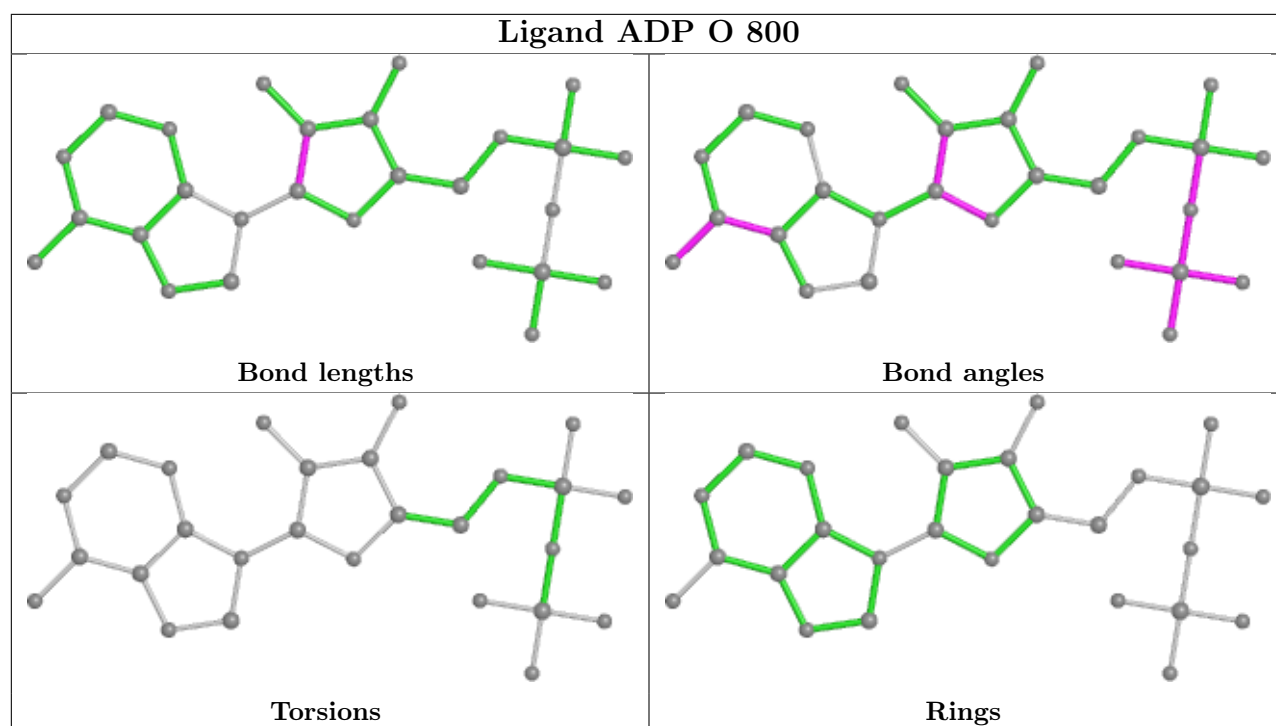


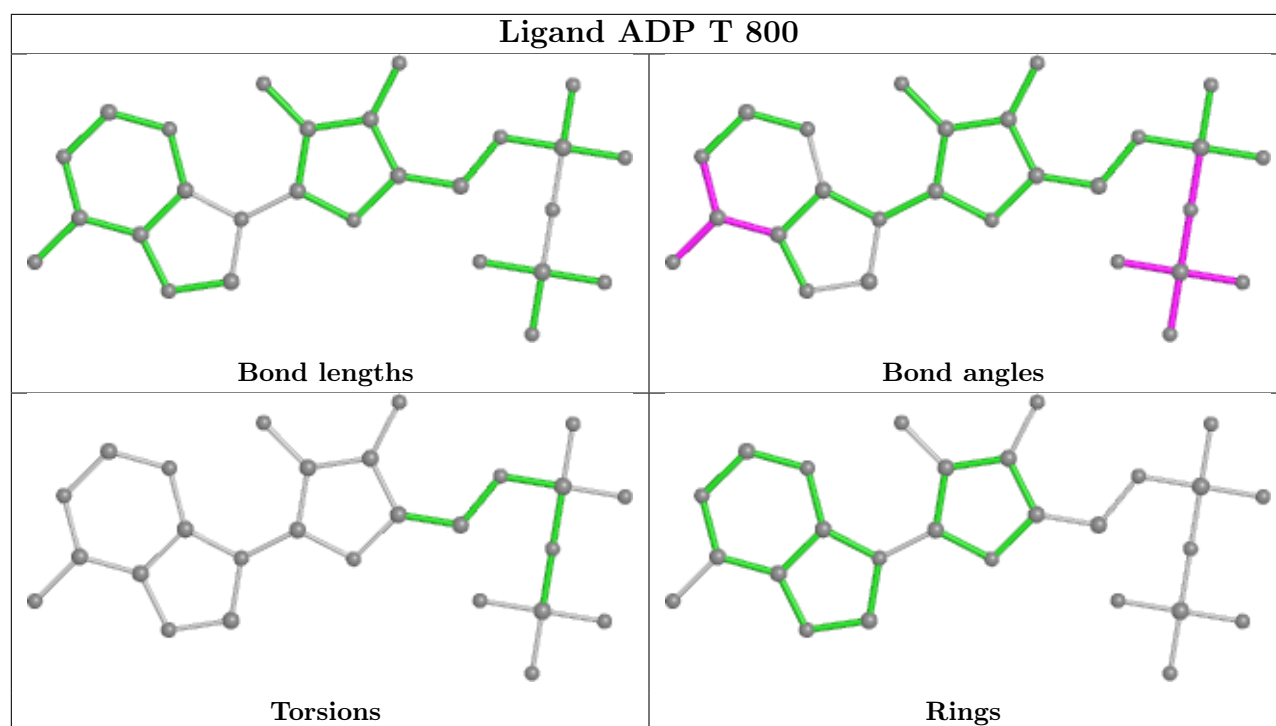












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

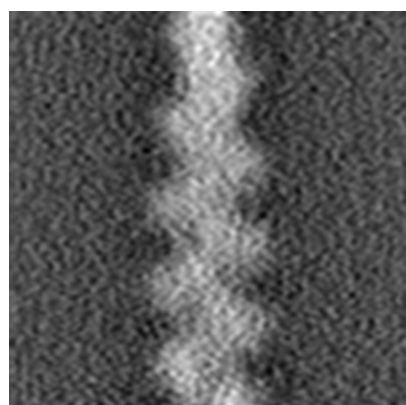
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-1674. These allow visual inspection of the internal detail of the map and identification of artifacts.

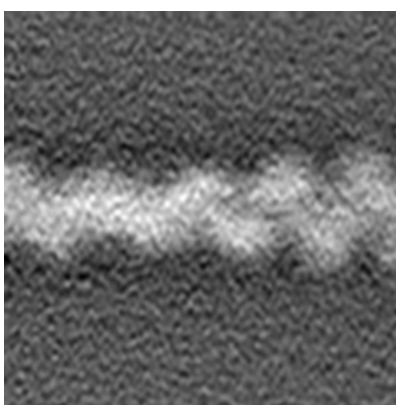
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

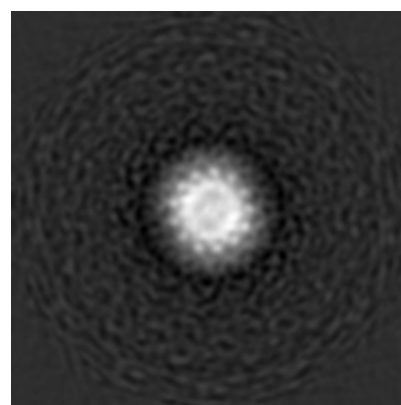
#### 6.1.1 Primary map



X



Y

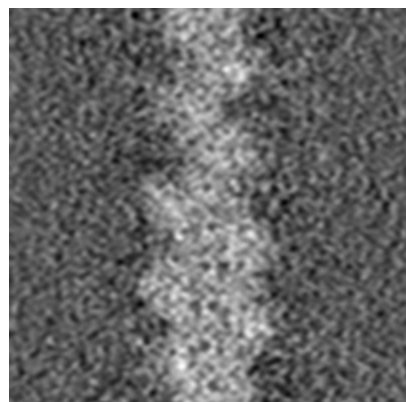


Z

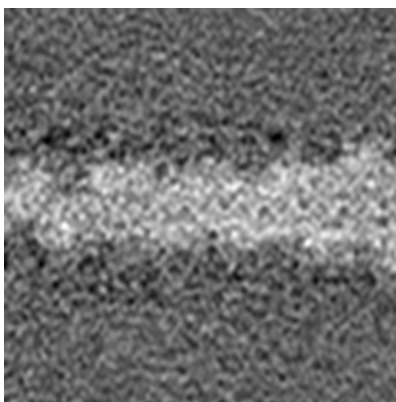
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

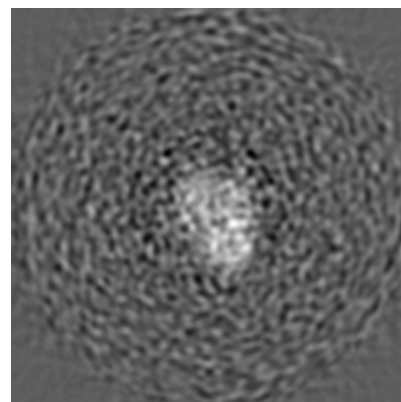
#### 6.2.1 Primary map



X Index: 60



Y Index: 60

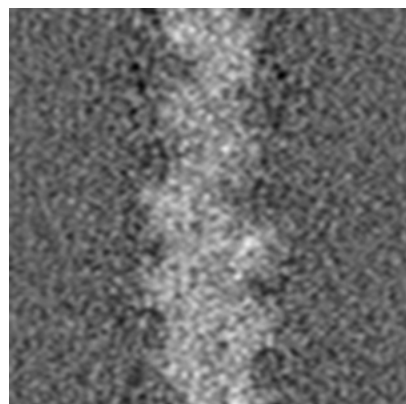


Z Index: 60

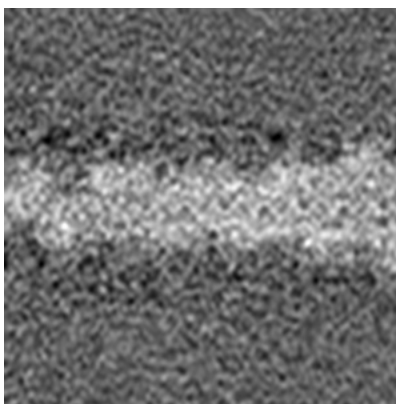
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

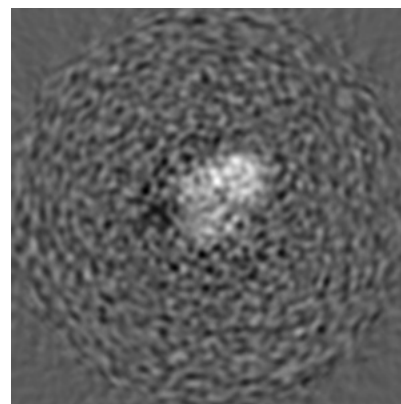
### 6.3.1 Primary map



X Index: 62



Y Index: 60

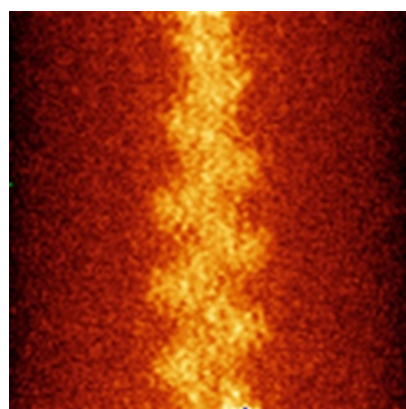


Z Index: 0

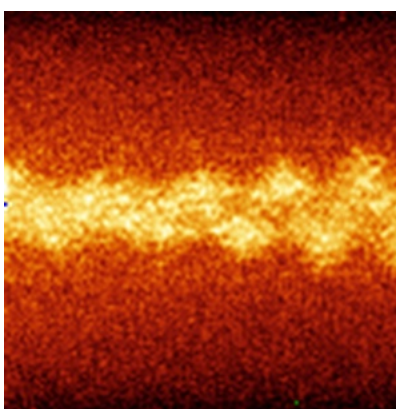
The images above show the largest variance slices of the map in three orthogonal directions.

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

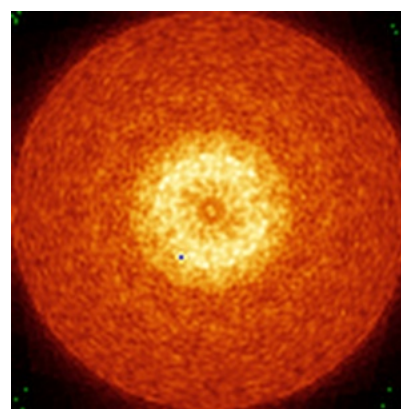
### 6.4.1 Primary map



X



Y

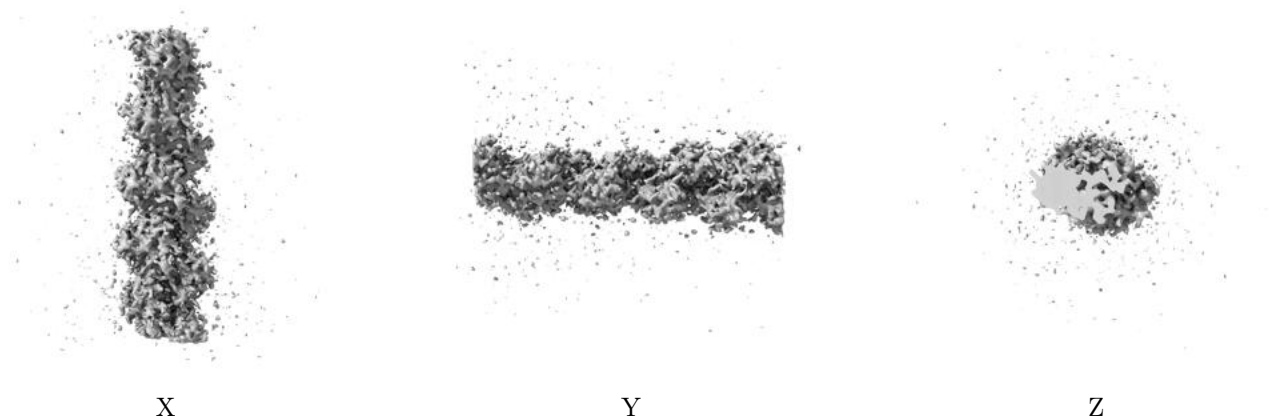


Z

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

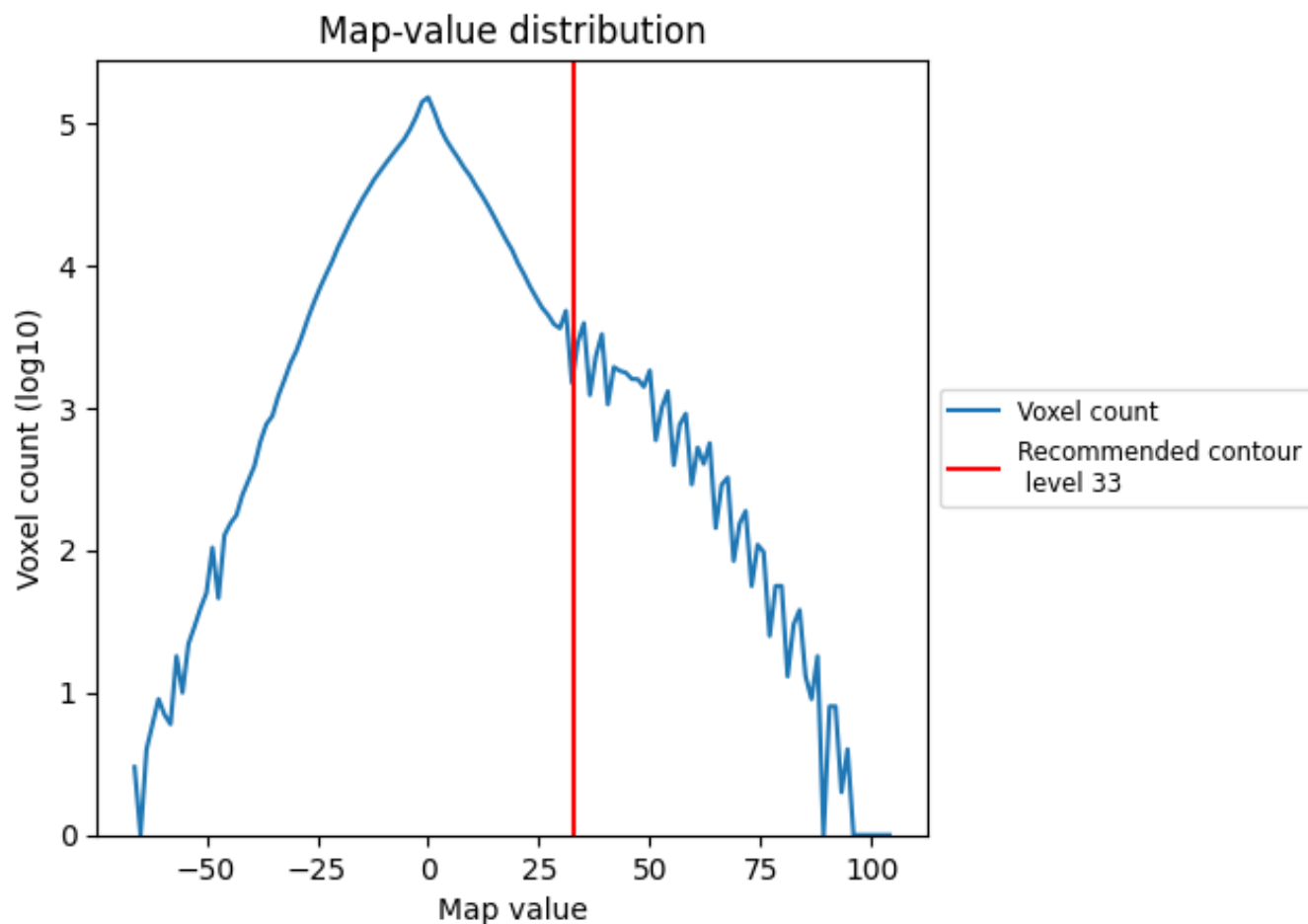
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

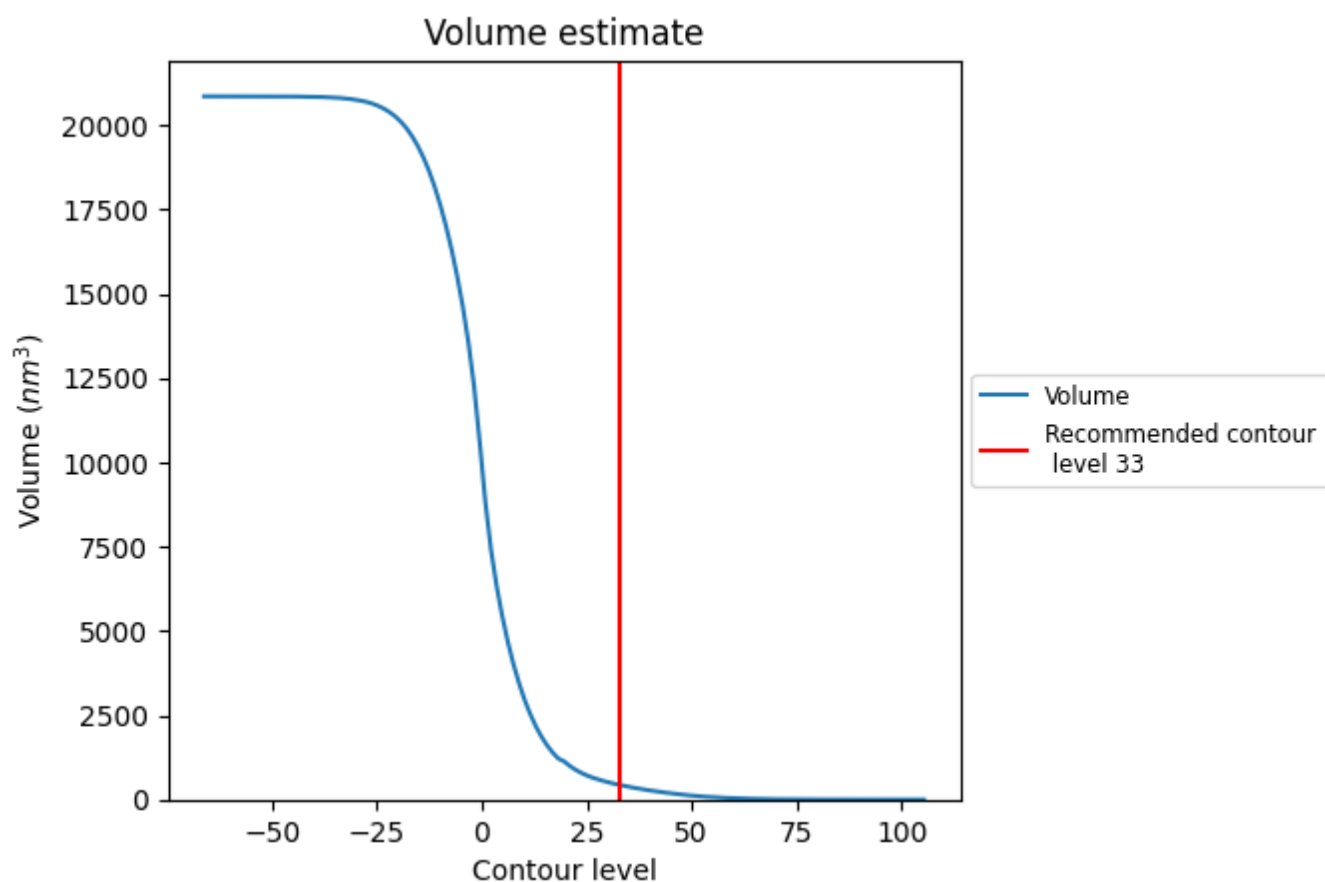
This section contains the results of statistical analysis of the map.

### 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.

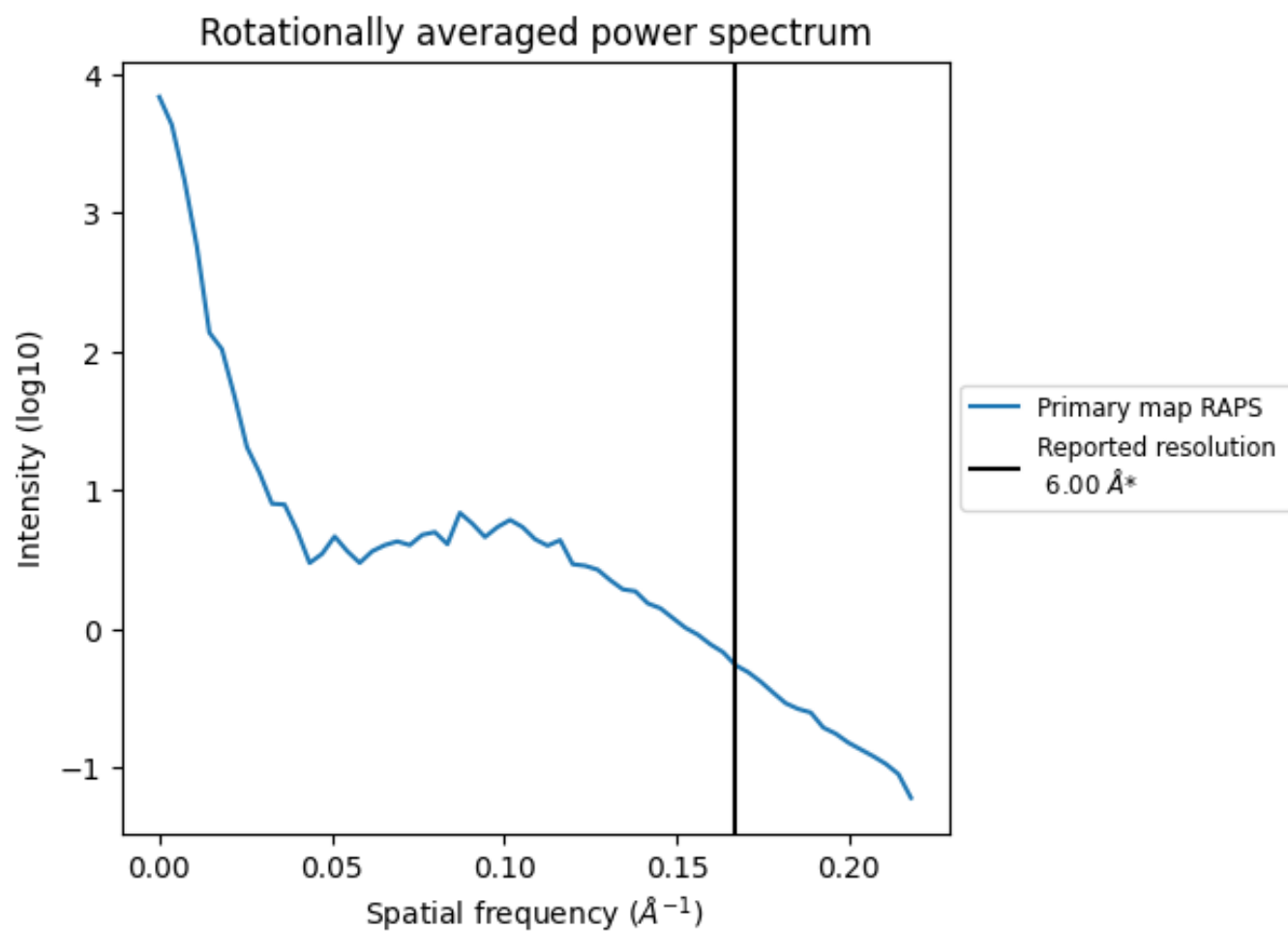
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 430 nm<sup>3</sup>; this corresponds to an approximate mass of 388 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 0.167 Å<sup>-1</sup>



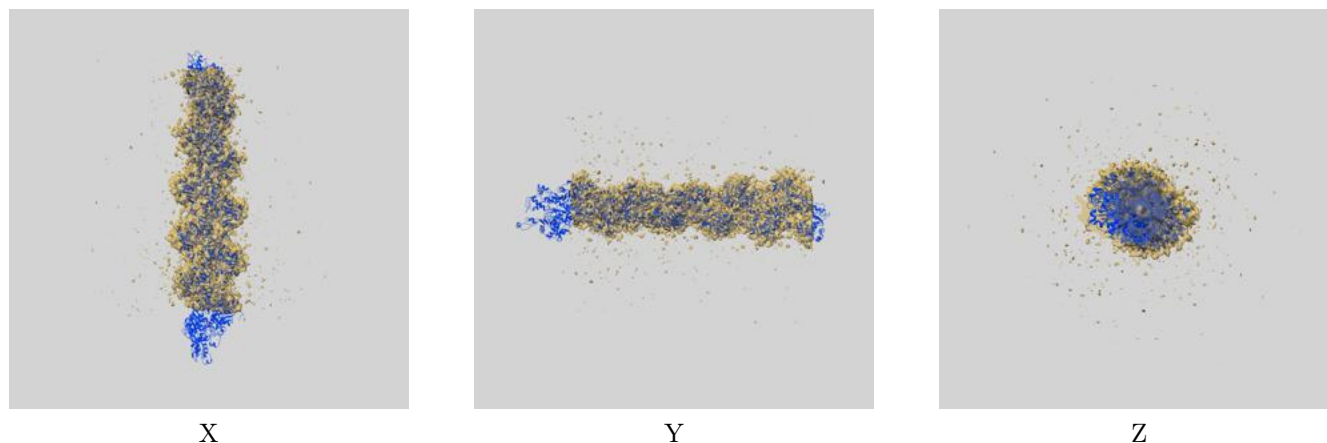
## 8 Fourier-Shell correlation

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

## 9 Map-model fit [i](#)

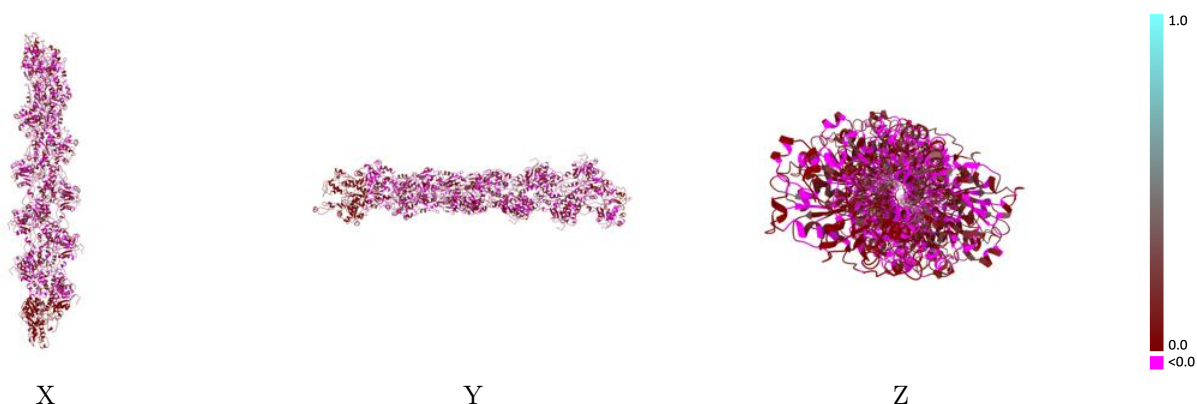
This section contains information regarding the fit between EMDB map EMD-1674 and PDB model 3G37. Per-residue inclusion information can be found in [section 3](#) on [page 10](#).

### 9.1 Map-model overlay [i](#)



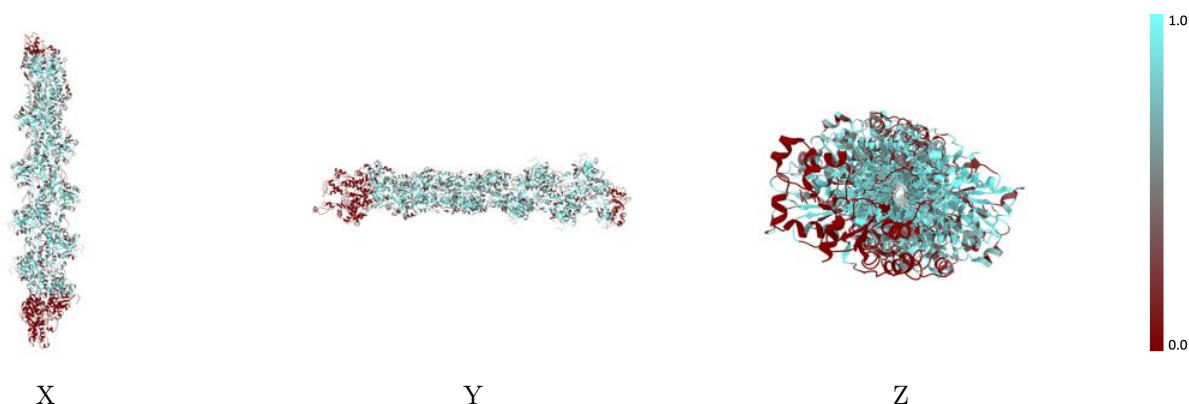
The images above show the 3D surface view of the map at the recommended contour level 33.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



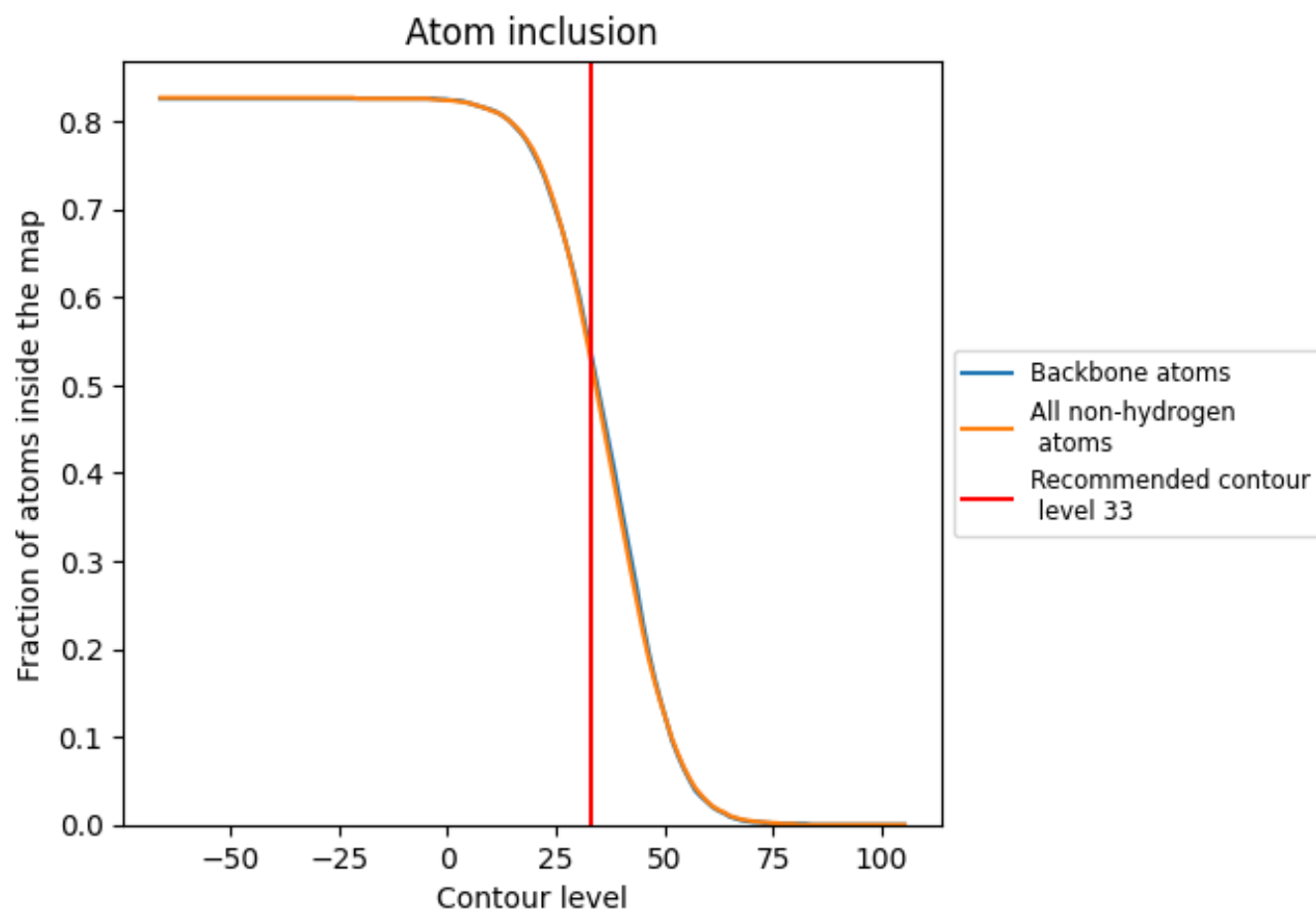
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (33).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 54% of all backbone atoms, 53% 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 (33) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.5300	<div></div> 0.0660
O	<div></div> 0.3890	<div></div> 0.0510
P	<div></div> 0.6370	<div></div> 0.0790
Q	<div></div> 0.6350	<div></div> 0.0770
R	<div></div> 0.6400	<div></div> 0.0840
S	<div></div> 0.6460	<div></div> 0.0830
T	<div></div> 0.6410	<div></div> 0.0810
U	<div></div> 0.6410	<div></div> 0.0860
V	<div></div> 0.6450	<div></div> 0.0830
W	<div></div> 0.6350	<div></div> 0.0820
X	<div></div> 0.5890	<div></div> 0.0640
Y	<div></div> 0.2590	<div></div> 0.0200
Z	<div></div> 0.0000	<div></div> -0.0000

