



Full wwPDB EM Validation Report ⓘ

Mar 4, 2026 – 10:19 PM UTC

PDB ID : 9O1A / pdb_00009o1a
EMDB ID : EMD-70000
Title : Pseudomonas aeruginosa ATPase State1 F1Fo focused
Authors : Stewart, A.G.; Sobti, M.
Deposited on : 2025-04-03
Resolution : 2.77 Å(reported)

This is a Full wwPDB EM Validation 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

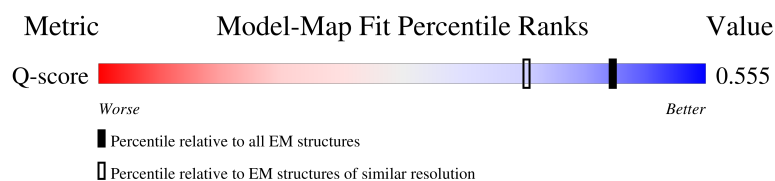
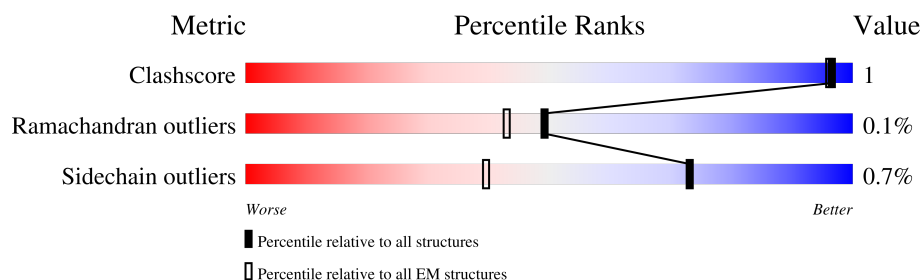
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

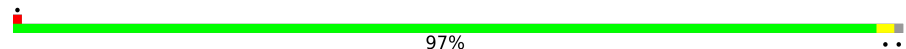
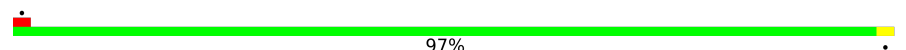
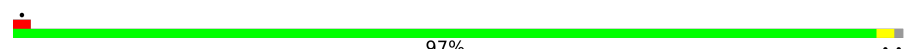
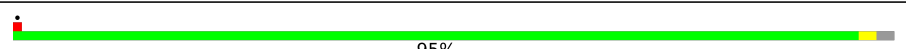
The reported resolution of this entry is 2.77 Å.

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.



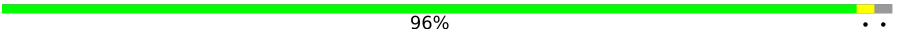
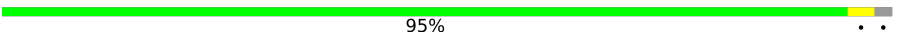
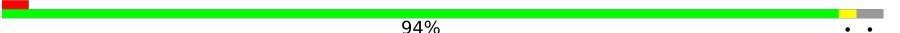
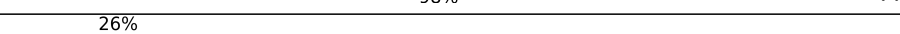
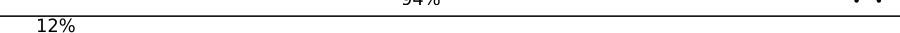
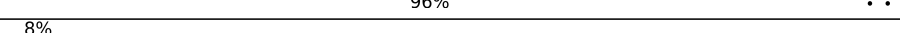
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10695 (2.27 - 3.27)

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	513	 97% ..
1	B	513	 97% .
1	C	513	 97% ..
2	D	464	 95% ..

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Mol	Chain	Length	Quality of chain
2	E	464	 96%
2	F	464	 95%
3	G	286	 94%
4	H	141	 12% 89% 9%
5	I	85	 91% 7%
5	J	85	 88% 9%
5	L	85	 91% 7%
5	M	85	 88% 9%
5	N	85	 94%
5	O	85	 93% 5%
5	P	85	 91% 6%
5	Q	85	 93% 5%
5	R	85	 93%
5	S	85	 93% 5%
6	W	178	 98%
7	X	156	 26% 94%
7	Y	156	 12% 96%
8	a	289	 8% 94%

2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 37051 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 ATP synthase subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	506	Total	C	N	O	S	0	0
			3823	2405	662	743	13		
1	B	512	Total	C	N	O	S	0	0
			3872	2435	671	752	14		
1	C	509	Total	C	N	O	S	0	0
			3846	2420	666	747	13		

- Molecule 2 is a protein called ATP synthase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	456	Total	C	N	O	S	0	0
			3464	2188	591	673	12		
2	E	456	Total	C	N	O	S	0	0
			3464	2188	591	673	12		
2	F	456	Total	C	N	O	S	0	0
			3464	2188	591	673	12		

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MET	-	expression tag	UNP A6VF32
D	2	HIS	-	expression tag	UNP A6VF32
D	3	HIS	-	expression tag	UNP A6VF32
D	4	HIS	-	expression tag	UNP A6VF32
D	5	HIS	-	expression tag	UNP A6VF32
D	6	HIS	-	expression tag	UNP A6VF32
D	7	HIS	-	expression tag	UNP A6VF32
E	1	MET	-	expression tag	UNP A6VF32
E	2	HIS	-	expression tag	UNP A6VF32
E	3	HIS	-	expression tag	UNP A6VF32
E	4	HIS	-	expression tag	UNP A6VF32
E	5	HIS	-	expression tag	UNP A6VF32
E	6	HIS	-	expression tag	UNP A6VF32

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Chain	Residue	Modelled	Actual	Comment	Reference
E	7	HIS	-	expression tag	UNP A6VF32
F	1	MET	-	expression tag	UNP A6VF32
F	2	HIS	-	expression tag	UNP A6VF32
F	3	HIS	-	expression tag	UNP A6VF32
F	4	HIS	-	expression tag	UNP A6VF32
F	5	HIS	-	expression tag	UNP A6VF32
F	6	HIS	-	expression tag	UNP A6VF32
F	7	HIS	-	expression tag	UNP A6VF32

- Molecule 3 is a protein called ATP synthase gamma chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	276	Total	C	N	O	S	0	0
			2132	1339	381	399	13		

- Molecule 4 is a protein called ATP synthase epsilon chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	128	Total	C	N	O	S	0	0
			933	590	159	181	3		

- Molecule 5 is a protein called ATP synthase subunit c.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	I	83	Total	C	N	O	S	0	0
			594	397	92	100	5		
5	J	83	Total	C	N	O	S	0	0
			594	397	92	100	5		
5	L	83	Total	C	N	O	S	0	0
			594	397	92	100	5		
5	M	83	Total	C	N	O	S	0	0
			594	397	92	100	5		
5	N	83	Total	C	N	O	S	0	0
			594	397	92	100	5		
5	O	83	Total	C	N	O	S	0	0
			594	397	92	100	5		
5	P	83	Total	C	N	O	S	0	0
			591	394	92	100	5		
5	Q	83	Total	C	N	O	S	0	0
			594	397	92	100	5		
5	R	83	Total	C	N	O	S	0	0
			594	397	92	100	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
5	S	83	Total	C	N	O	S	0	0
			594	397	92	100	5		

- Molecule 6 is a protein called ATP synthase subunit delta.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	W	177	Total	C	N	O	S	0	0
			1345	840	237	266	2		

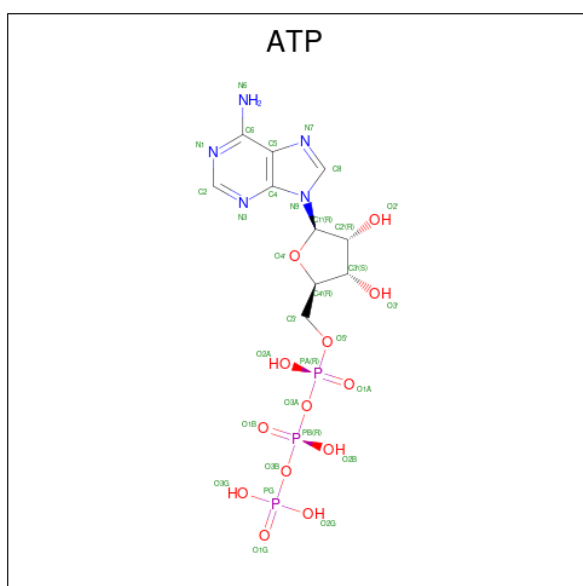
- Molecule 7 is a protein called ATP synthase subunit b.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	X	154	Total	C	N	O	S	0	0
			1174	731	217	223	3		
7	Y	154	Total	C	N	O	S	0	0
			1174	731	217	223	3		

- Molecule 8 is a protein called ATP synthase subunit a.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	a	286	Total	C	N	O	S	0	0
			2243	1503	354	377	9		

- Molecule 9 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).

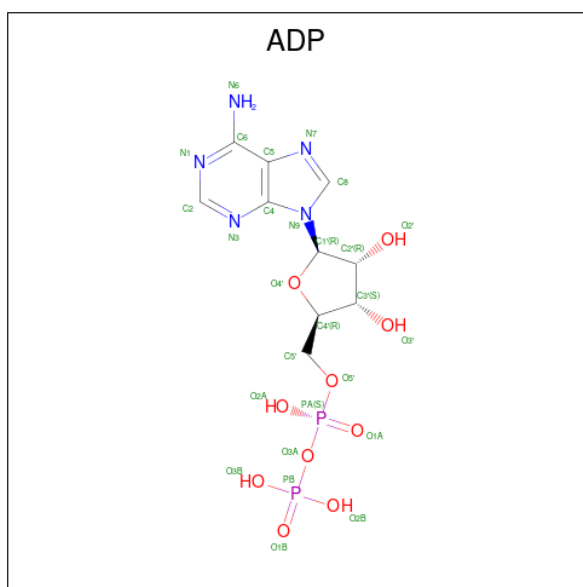


Mol	Chain	Residues	Atoms					AltConf
9	A	1	Total	C	N	O	P	0
			31	10	5	13	3	
9	B	1	Total	C	N	O	P	0
			31	10	5	13	3	
9	C	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
10	A	1	Total	Mg	0
			1	1	
10	B	1	Total	Mg	0
			1	1	
10	C	1	Total	Mg	0
			1	1	
10	D	1	Total	Mg	0
			1	1	
10	E	1	Total	Mg	0
			1	1	

- Molecule 11 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂) (labeled as "Ligand of Interest" by depositor).



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

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Mol	Chain	Residues	Atoms					AltConf
11	E	1	Total	C	N	O	P	0
			27	10	5	10	2	
11	F	1	Total	C	N	O	P	0
			27	10	5	10	2	

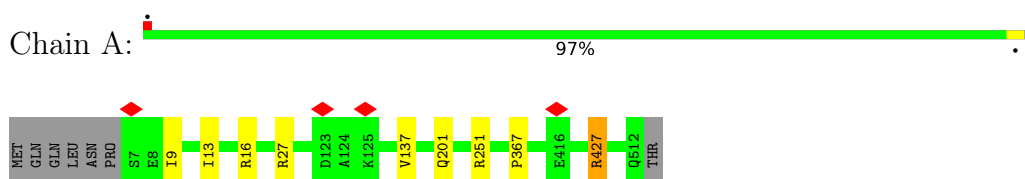
- Molecule 12 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
12	a	1	Total	Zn	0
			1	1	

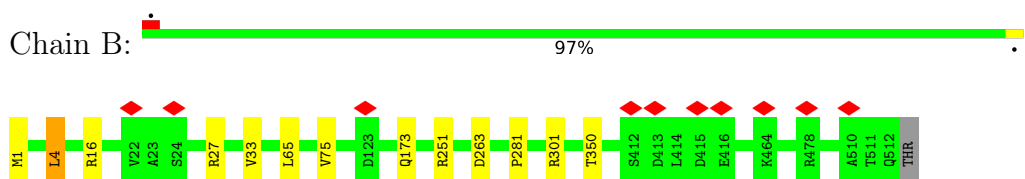
3 Residue-property plots [i](#)

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

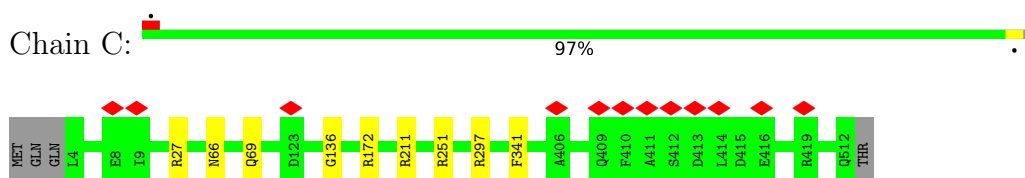
- Molecule 1: ATP synthase subunit alpha



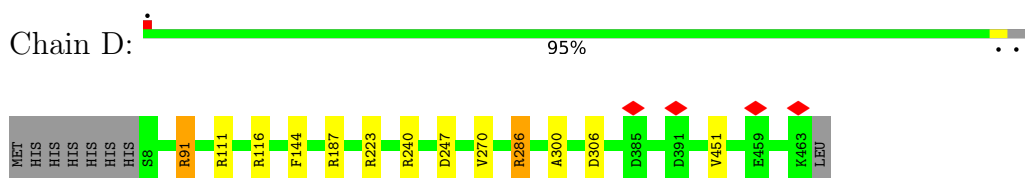
- Molecule 1: ATP synthase subunit alpha



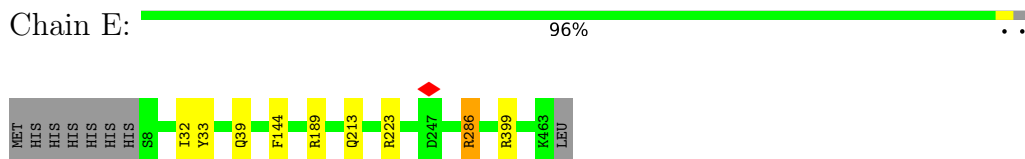
- Molecule 1: ATP synthase subunit alpha



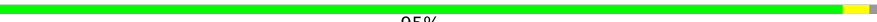
- Molecule 2: ATP synthase subunit beta

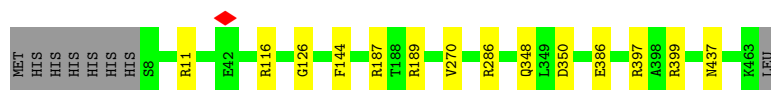


- Molecule 2: ATP synthase subunit beta



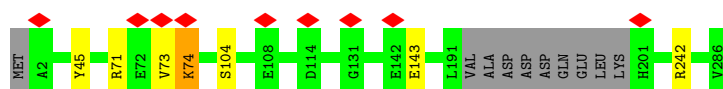
- Molecule 2: ATP synthase subunit beta

Chain F:  95%



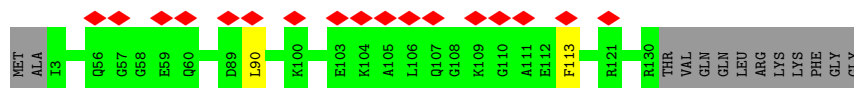
- Molecule 3: ATP synthase gamma chain

Chain G:  94%



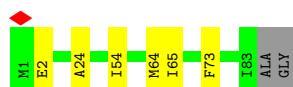
- Molecule 4: ATP synthase epsilon chain

Chain H:  12% 89% 9%



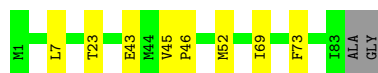
- Molecule 5: ATP synthase subunit c

Chain I:  91% 7%



- Molecule 5: ATP synthase subunit c

Chain J:  88% 9%



- Molecule 5: ATP synthase subunit c

Chain L:  91% 7%



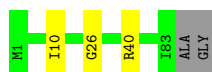
- Molecule 5: ATP synthase subunit c

Chain M:  88% 9%



- Molecule 5: ATP synthase subunit c

Chain N: 94% . .



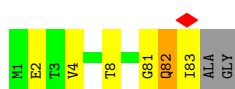
- Molecule 5: ATP synthase subunit c

Chain O: 93% 5% .



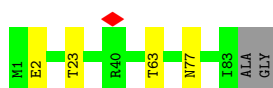
- Molecule 5: ATP synthase subunit c

Chain P: 91% 6% . .



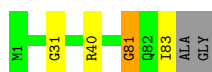
- Molecule 5: ATP synthase subunit c

Chain Q: 93% 5% .



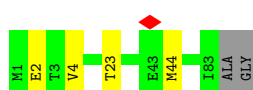
- Molecule 5: ATP synthase subunit c

Chain R: 93% . . .



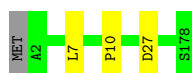
- Molecule 5: ATP synthase subunit c

Chain S: 93% 5% .



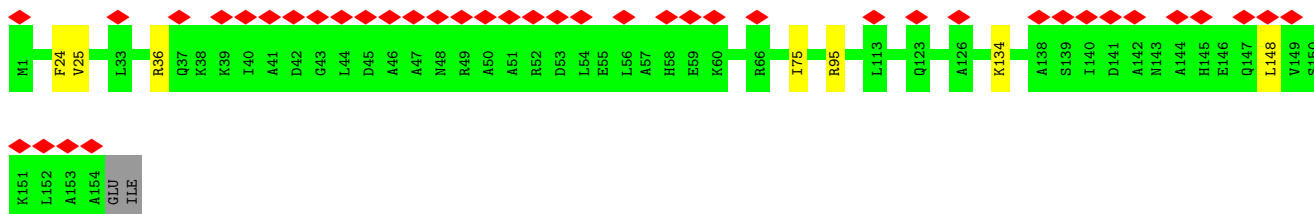
- Molecule 6: ATP synthase subunit delta

Chain W: 98% . .



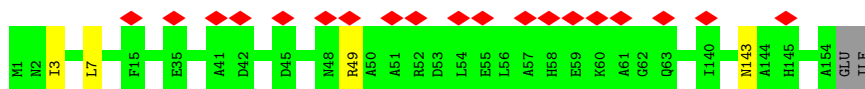
- Molecule 7: ATP synthase subunit b

Chain X: 26% 94%



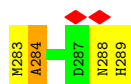
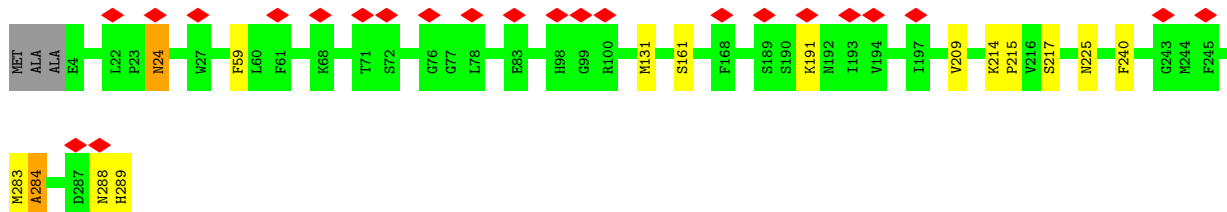
- Molecule 7: ATP synthase subunit b

Chain Y: 12% 96%



- Molecule 8: ATP synthase subunit a

Chain a: 8% 94%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	28761	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 ($e^-/\text{\AA}^2$)	62	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.350	Depositor
Minimum map value	-0.122	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.012	Depositor
Recommended contour level	0.06	Depositor
Map size ($\text{\AA}$)	332.0, 332.0, 332.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.83, 0.83, 0.83	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, ATP, 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	0.73	0/3875	1.31	6/5233 (0.1%)
1	B	0.74	0/3925	1.32	7/5301 (0.1%)
1	C	0.72	0/3899	1.35	7/5267 (0.1%)
2	D	0.75	0/3521	1.33	7/4767 (0.1%)
2	E	0.73	0/3521	1.32	8/4767 (0.2%)
2	F	0.73	0/3521	1.32	8/4767 (0.2%)
3	G	0.74	0/2161	1.34	2/2908 (0.1%)
4	H	0.72	0/946	1.38	4/1282 (0.3%)
5	I	0.68	0/603	1.27	1/816 (0.1%)
5	J	0.78	0/603	1.34	1/816 (0.1%)
5	L	0.66	0/603	1.36	1/816 (0.1%)
5	M	0.69	0/603	1.33	1/816 (0.1%)
5	N	0.68	0/603	1.36	1/816 (0.1%)
5	O	0.69	0/603	1.41	0/816
5	P	0.73	1/600 (0.2%)	1.41	1/812 (0.1%)
5	Q	0.73	0/603	1.34	1/816 (0.1%)
5	R	0.82	2/603 (0.3%)	1.42	2/816 (0.2%)
5	S	0.66	0/603	1.35	1/816 (0.1%)
6	W	0.72	0/1358	1.31	0/1833
7	X	0.74	0/1184	1.40	3/1593 (0.2%)
7	Y	0.77	0/1184	1.36	2/1593 (0.1%)
8	a	0.78	1/2306 (0.0%)	1.36	5/3141 (0.2%)
All	All	0.74	4/37428 (0.0%)	1.34	69/50608 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
2	D	0	4
2	E	0	1
2	F	0	1
3	G	0	1
5	P	0	1
5	R	0	1
7	X	0	1
8	a	0	1
All	All	0	12

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	R	81	GLY	C-O	-9.11	1.13	1.23
5	P	82	GLN	C-O	-5.30	1.14	1.24
8	a	161	SER	CA-CB	-5.29	1.44	1.53
5	R	83	ILE	C-O	-5.06	1.13	1.23

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	a	284	ALA	O-C-N	-26.17	89.63	122.34
5	R	81	GLY	O-C-N	-11.51	107.79	122.29
5	P	82	GLN	O-C-N	-9.40	110.25	122.00
8	a	284	ALA	CA-C-N	7.80	134.67	122.26
8	a	284	ALA	C-N-CA	7.80	134.67	122.26
2	D	451	VAL	N-CA-C	7.03	118.02	112.12
2	F	144	PHE	CA-CB-CG	6.61	120.41	113.80
7	X	25	VAL	N-CA-C	6.55	117.17	110.82
3	G	74	LYS	N-CA-C	6.47	120.42	111.55
2	E	39	GLN	N-CA-C	6.45	118.88	111.02
3	G	242	ARG	NE-CZ-NH2	6.29	124.86	119.20
1	C	27	ARG	NE-CZ-NH2	6.29	124.86	119.20
7	X	24	PHE	N-CA-C	6.21	118.97	111.40
1	A	367	PRO	N-CA-C	-6.16	100.83	110.50
5	L	73	PHE	CA-CB-CG	6.10	119.90	113.80
2	D	91	ARG	NE-CZ-NH2	6.03	124.63	119.20
8	a	24	ASN	CA-CB-CG	6.00	118.60	112.60
2	E	189	ARG	NE-CZ-NH2	5.95	124.56	119.20
2	D	223	ARG	NE-CZ-NH2	5.83	124.45	119.20
8	a	240	PHE	CA-CB-CG	5.78	119.58	113.80
2	E	213	GLN	N-CA-C	5.75	118.40	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	144	PHE	CA-CB-CG	5.74	119.54	113.80
2	D	286	ARG	NE-CZ-NH2	5.73	124.36	119.20
1	A	251	ARG	NE-CZ-NH2	5.69	124.32	119.20
2	E	286	ARG	NE-CZ-NH2	5.59	124.23	119.20
2	F	11	ARG	NE-CZ-NH2	5.59	124.23	119.20
1	B	251	ARG	NE-CZ-NH2	5.58	124.23	119.20
5	Q	77	ASN	CA-CB-CG	5.55	118.15	112.60
2	E	399	ARG	NE-CZ-NH2	5.50	124.15	119.20
1	B	27	ARG	NE-CZ-NH2	5.43	124.09	119.20
5	J	73	PHE	CA-CB-CG	5.41	119.21	113.80
1	A	427	ARG	NE-CZ-NH2	5.38	124.05	119.20
2	E	144	PHE	CA-CB-CG	5.38	119.18	113.80
4	H	90	LEU	CA-C-N	5.33	127.42	120.28
4	H	90	LEU	C-N-CA	5.33	127.42	120.28
2	E	223	ARG	NE-CZ-NH2	5.32	123.99	119.20
1	C	251	ARG	NE-CZ-NH2	5.32	123.98	119.20
7	Y	49	ARG	NE-CZ-NH2	5.31	123.98	119.20
2	F	350	ASP	CA-CB-CG	5.31	117.91	112.60
5	S	2	GLU	N-CA-C	5.31	117.48	111.11
1	B	301	ARG	NE-CZ-NH2	5.30	123.97	119.20
1	B	281	PRO	CA-C-N	5.25	125.45	120.31
1	B	281	PRO	C-N-CA	5.25	125.45	120.31
2	E	213	GLN	OE1-CD-NE2	-5.22	117.38	122.60
7	Y	143	ASN	OD1-CG-ND2	-5.21	117.39	122.60
2	F	397	ARG	NE-CZ-NH2	5.20	123.88	119.20
1	A	201	GLN	OE1-CD-NE2	-5.18	117.42	122.60
1	C	172	ARG	NE-CZ-NH2	5.17	123.86	119.20
7	X	95	ARG	NE-CZ-NH2	5.17	123.85	119.20
1	C	341	PHE	N-CA-C	5.17	116.60	110.97
2	F	399	ARG	NE-CZ-NH2	5.13	123.82	119.20
2	D	111	ARG	NE-CZ-NH2	5.12	123.81	119.20
2	F	286	ARG	NE-CZ-NH2	5.12	123.81	119.20
5	R	40	ARG	NE-CZ-NH2	5.12	123.81	119.20
2	D	240	ARG	NE-CZ-NH2	5.11	123.80	119.20
5	N	40	ARG	NE-CZ-NH2	5.11	123.80	119.20
1	C	136	GLY	CA-C-N	5.08	127.16	120.60
1	C	136	GLY	C-N-CA	5.08	127.16	120.60
5	I	73	PHE	CA-CB-CG	5.08	118.88	113.80
2	F	187	ARG	NE-CZ-NH2	5.07	123.76	119.20
1	A	27	ARG	NE-CZ-NH2	5.06	123.76	119.20
1	B	350	THR	N-CA-C	5.06	117.21	110.53
4	H	113	PHE	CA-C-N	5.06	127.31	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	H	113	PHE	C-N-CA	5.06	127.31	120.38
1	A	16	ARG	NE-CZ-NH2	5.05	123.75	119.20
5	M	73	PHE	CA-CB-CG	5.04	118.84	113.80
1	B	173	GLN	N-CA-C	5.04	118.69	112.24
2	F	189	ARG	NE-CZ-NH2	5.01	123.71	119.20
1	C	297	ARG	NE-CZ-NH2	5.01	123.71	119.20

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	16	ARG	Sidechain
2	D	116	ARG	Sidechain
2	D	187	ARG	Sidechain
2	D	286	ARG	Sidechain
2	D	91	ARG	Sidechain
2	E	286	ARG	Sidechain
2	F	116	ARG	Sidechain
3	G	45	TYR	Sidechain
5	P	82	GLN	Mainchain
5	R	81	GLY	Mainchain
7	X	36	ARG	Sidechain
8	a	284	ALA	Mainchain

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	A	3823	0	3872	1	0
1	B	3872	0	3924	7	0
1	C	3846	0	3896	2	0
2	D	3464	0	3482	1	0
2	E	3464	0	3482	1	0
2	F	3464	0	3482	2	0
3	G	2132	0	2171	1	0
4	H	933	0	940	0	0
5	I	594	0	641	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	J	594	0	641	5	0
5	L	594	0	641	5	0
5	M	594	0	641	6	0
5	N	594	0	641	2	0
5	O	594	0	641	3	0
5	P	591	0	632	4	0
5	Q	594	0	641	3	0
5	R	594	0	641	1	0
5	S	594	0	641	1	0
6	W	1345	0	1384	4	0
7	X	1174	0	1204	1	0
7	Y	1174	0	1204	8	0
8	a	2243	0	2290	17	0
9	A	31	0	12	0	0
9	B	31	0	12	0	0
9	C	31	0	12	0	0
10	A	1	0	0	0	0
10	B	1	0	0	0	0
10	C	1	0	0	0	0
10	D	1	0	0	0	0
10	E	1	0	0	0	0
11	D	27	0	12	0	0
11	E	27	0	12	0	0
11	F	27	0	12	0	0
12	a	1	0	0	0	0
All	All	37051	0	37804	56	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 1.

All (56) 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)
5:M:49:GLN:NE2	8:a:288:ASN:O	1.93	1.01
1:B:4:LEU:HD21	6:W:10:PRO:HG2	1.72	0.72
5:L:81:GLY:O	5:L:83:ILE:HG13	1.90	0.72
8:a:191:LYS:HD2	8:a:191:LYS:H	1.54	0.71
5:I:2:GLU:H	5:I:2:GLU:CD	1.99	0.70
7:Y:3:ILE:HD11	8:a:131:MET:CE	2.26	0.65
5:I:64:MET:HE2	8:a:217:SER:HA	1.79	0.65
5:M:60:ASP:HB3	5:M:64:MET:HE2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:Y:3:ILE:HG13	8:a:131:MET:HE1	1.83	0.60
7:Y:3:ILE:CG1	8:a:131:MET:HE1	2.32	0.59
5:I:54:ILE:HG12	8:a:283:MET:CE	2.33	0.58
1:B:65:LEU:HD12	1:B:75:VAL:HG21	1.87	0.56
5:I:54:ILE:HG12	8:a:283:MET:HE1	1.87	0.56
5:L:1:MET:H2	5:Q:2:GLU:CD	2.15	0.55
5:Q:23:THR:HG23	5:Q:63:THR:HG21	1.89	0.54
5:O:64:MET:HA	5:O:67:VAL:HG12	1.89	0.54
5:L:81:GLY:O	5:L:83:ILE:N	2.41	0.53
1:C:211:ARG:CZ	2:F:126:GLY:HA2	2.38	0.53
5:J:45:VAL:HB	5:J:46:PRO:HD3	1.91	0.53
8:a:191:LYS:HD2	8:a:191:LYS:N	2.23	0.52
5:M:60:ASP:O	5:M:64:MET:HG2	2.09	0.52
1:B:1:MET:O	1:B:1:MET:HG2	2.09	0.51
5:M:53:PHE:CE2	8:a:289:HIS:ND1	2.77	0.51
1:B:65:LEU:HD12	1:B:75:VAL:CG2	2.42	0.49
7:Y:3:ILE:HD11	8:a:131:MET:HE1	1.94	0.48
5:J:7:LEU:HD22	5:N:10:ILE:HD11	1.95	0.48
5:L:81:GLY:C	5:L:83:ILE:H	2.20	0.48
7:Y:3:ILE:HD11	8:a:131:MET:HE2	1.94	0.48
5:Q:23:THR:CG2	5:Q:63:THR:HG21	2.43	0.48
5:L:81:GLY:C	5:L:83:ILE:N	2.69	0.48
5:O:55:VAL:HG11	5:R:31:GLY:HA3	1.96	0.48
1:C:211:ARG:NH1	2:F:126:GLY:HA2	2.29	0.48
5:P:81:GLY:C	5:P:83:ILE:N	2.72	0.47
1:A:9:ILE:HG23	1:A:13:ILE:HG22	1.95	0.47
1:B:4:LEU:HD21	6:W:10:PRO:CG	2.42	0.46
7:Y:3:ILE:CD1	8:a:131:MET:HE1	2.45	0.46
5:I:65:ILE:HD11	8:a:209:VAL:HG13	1.98	0.46
7:X:148:LEU:HD22	7:X:148:LEU:H	1.80	0.46
1:B:1:MET:HG3	6:W:7:LEU:HD11	1.98	0.45
2:E:32:ILE:O	2:E:33:TYR:HB2	2.17	0.44
5:J:52:MET:HE2	5:J:52:MET:HB3	1.90	0.44
5:M:55:VAL:HG11	5:O:31:GLY:HA3	2.00	0.43
5:J:43:GLU:CD	5:J:43:GLU:H	2.27	0.43
7:Y:3:ILE:O	7:Y:7:LEU:HD23	2.19	0.43
5:P:4:VAL:O	5:P:8:THR:HG23	2.19	0.42
5:I:54:ILE:HG12	8:a:283:MET:HE2	2.02	0.41
2:D:247:ASP:O	2:D:300:ALA:HB3	2.20	0.41
5:P:81:GLY:C	5:P:83:ILE:H	2.26	0.41
7:Y:3:ILE:CD1	8:a:131:MET:CE	2.95	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:24:ALA:O	5:N:26:GLY:HA3	2.21	0.41
5:P:2:GLU:HG3	5:S:4:VAL:HG23	2.02	0.41
1:B:1:MET:CG	6:W:7:LEU:HD11	2.51	0.41
3:G:73:VAL:HG12	3:G:74:LYS:H	1.85	0.41
5:M:1:MET:HE2	5:M:1:MET:HB2	1.99	0.41
8:a:214:LYS:N	8:a:215:PRO:HD2	2.36	0.40
5:J:7:LEU:HD23	5:J:7:LEU:HA	1.93	0.40

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	504/513 (98%)	492 (98%)	12 (2%)	0	100	100
1	B	510/513 (99%)	491 (96%)	19 (4%)	0	100	100
1	C	507/513 (99%)	494 (97%)	13 (3%)	0	100	100
2	D	454/464 (98%)	444 (98%)	9 (2%)	1 (0%)	43	70
2	E	454/464 (98%)	441 (97%)	13 (3%)	0	100	100
2	F	454/464 (98%)	444 (98%)	9 (2%)	1 (0%)	43	70
3	G	272/286 (95%)	268 (98%)	4 (2%)	0	100	100
4	H	126/141 (89%)	122 (97%)	4 (3%)	0	100	100
5	I	81/85 (95%)	81 (100%)	0	0	100	100
5	J	81/85 (95%)	81 (100%)	0	0	100	100
5	L	81/85 (95%)	79 (98%)	1 (1%)	1 (1%)	10	30
5	M	81/85 (95%)	81 (100%)	0	0	100	100
5	N	81/85 (95%)	78 (96%)	3 (4%)	0	100	100
5	O	81/85 (95%)	80 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
5	P	81/85 (95%)	79 (98%)	2 (2%)	0	100	100
5	Q	81/85 (95%)	80 (99%)	1 (1%)	0	100	100
5	R	81/85 (95%)	80 (99%)	1 (1%)	0	100	100
5	S	81/85 (95%)	81 (100%)	0	0	100	100
6	W	175/178 (98%)	173 (99%)	2 (1%)	0	100	100
7	X	152/156 (97%)	151 (99%)	1 (1%)	0	100	100
7	Y	152/156 (97%)	149 (98%)	3 (2%)	0	100	100
8	a	284/289 (98%)	275 (97%)	9 (3%)	0	100	100
All	All	4854/4987 (97%)	4744 (98%)	107 (2%)	3 (0%)	49	75

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	L	82	GLN
2	F	270	VAL
2	D	270	VAL

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	398/405 (98%)	396 (100%)	2 (0%)	81	92
1	B	404/405 (100%)	401 (99%)	3 (1%)	76	90
1	C	401/405 (99%)	399 (100%)	2 (0%)	81	92
2	D	368/376 (98%)	367 (100%)	1 (0%)	86	94
2	E	368/376 (98%)	368 (100%)	0	100	100
2	F	368/376 (98%)	365 (99%)	3 (1%)	73	89
3	G	226/235 (96%)	223 (99%)	3 (1%)	61	83
4	H	91/101 (90%)	91 (100%)	0	100	100
5	I	59/59 (100%)	59 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
5	J	59/59 (100%)	57 (97%)	2 (3%)	32	65
5	L	59/59 (100%)	58 (98%)	1 (2%)	53	80
5	M	59/59 (100%)	58 (98%)	1 (2%)	53	80
5	N	59/59 (100%)	59 (100%)	0	100	100
5	O	59/59 (100%)	59 (100%)	0	100	100
5	P	58/59 (98%)	58 (100%)	0	100	100
5	Q	59/59 (100%)	59 (100%)	0	100	100
5	R	59/59 (100%)	59 (100%)	0	100	100
5	S	59/59 (100%)	57 (97%)	2 (3%)	32	65
6	W	141/142 (99%)	140 (99%)	1 (1%)	76	90
7	X	115/117 (98%)	113 (98%)	2 (2%)	53	80
7	Y	115/117 (98%)	115 (100%)	0	100	100
8	a	238/239 (100%)	235 (99%)	3 (1%)	61	83
All	All	3822/3884 (98%)	3796 (99%)	26 (1%)	73	90

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	VAL
1	A	427	ARG
1	B	4	LEU
1	B	33	VAL
1	B	263	ASP
1	C	66	ASN
1	C	69	GLN
2	D	306	ASP
2	F	348	GLN
2	F	386	GLU
2	F	437	ASN
3	G	71	ARG
3	G	104	SER
3	G	143	GLU
5	J	23	THR
5	J	69	ILE
5	L	80	VAL
5	M	2	GLU
5	S	23	THR

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Mol	Chain	Res	Type
5	S	44	MET
6	W	27	ASP
7	X	75	ILE
7	X	134	LYS
8	a	24	ASN
8	a	59	PHE
8	a	225	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (54) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	145	GLN
1	A	187	GLN
1	A	203	GLN
1	A	221	ASN
1	A	273	GLN
1	A	424	HIS
1	A	426	GLN
1	A	480	HIS
1	B	3	GLN
1	B	145	GLN
1	B	203	GLN
1	B	221	ASN
1	B	273	GLN
1	C	5	ASN
1	C	88	ASN
1	C	187	GLN
1	C	221	ASN
1	C	273	GLN
1	C	362	ASN
1	C	480	HIS
2	D	14	GLN
2	D	48	GLN
2	D	72	ASN
2	D	124	GLN
2	D	192	ASN
2	D	348	GLN
2	D	370	GLN
2	E	72	ASN
2	E	192	ASN
2	E	248	ASN

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Mol	Chain	Res	Type
2	E	370	GLN
2	F	192	ASN
2	F	284	GLN
2	F	299	GLN
2	F	370	GLN
3	G	51	GLN
3	G	187	GLN
3	G	226	GLN
5	J	82	GLN
5	L	49	GLN
5	L	82	GLN
5	N	41	GLN
5	P	82	GLN
5	Q	82	GLN
5	S	41	GLN
6	W	84	ASN
6	W	104	GLN
7	X	123	GLN
7	X	143	ASN
7	Y	37	GLN
7	Y	111	GLN
7	Y	143	ASN
8	a	11	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 6 are 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$
11	ADP	E	501	10	28,29,29	0.50	0	43,45,45	0.56	0
9	ATP	B	600	10	32,33,33	1.34	4 (12%)	48,52,52	1.16	5 (10%)
9	ATP	A	600	10	32,33,33	1.35	3 (9%)	48,52,52	1.28	6 (12%)
11	ADP	F	501	-	28,29,29	1.10	1 (3%)	43,45,45	1.18	4 (9%)
11	ADP	D	501	10	28,29,29	1.02	1 (3%)	43,45,45	1.24	5 (11%)
9	ATP	C	600	10	32,33,33	1.19	4 (12%)	48,52,52	1.27	6 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	ADP	E	501	10	-	5/16/32/32	0/3/3/3
9	ATP	B	600	10	-	1/22/38/38	0/3/3/3
9	ATP	A	600	10	-	1/22/38/38	0/3/3/3
11	ADP	F	501	-	-	0/16/32/32	0/3/3/3
11	ADP	D	501	10	-	0/16/32/32	0/3/3/3
9	ATP	C	600	10	-	0/22/38/38	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	600	ATP	PB-O3A	-3.42	1.55	1.59
9	A	600	ATP	PB-O3A	-3.26	1.56	1.59
9	A	600	ATP	PA-O3A	-3.25	1.56	1.59
9	B	600	ATP	PA-O3A	-3.01	1.56	1.59
9	B	600	ATP	PB-O3B	-2.83	1.56	1.59
9	C	600	ATP	PB-O3B	-2.63	1.56	1.59
9	A	600	ATP	PB-O3B	-2.54	1.56	1.59
9	C	600	ATP	PB-O3A	-2.23	1.57	1.59
11	F	501	ADP	C5-C4	-2.17	1.35	1.39
9	C	600	ATP	PA-O3A	-2.15	1.57	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	C	600	ATP	C5-C4	-2.09	1.35	1.39
11	D	501	ADP	C5-C4	-2.04	1.35	1.39
9	B	600	ATP	C5-C4	-2.04	1.35	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	D	501	ADP	O2A-PA-O3A	3.34	116.31	107.27
9	A	600	ATP	C4-C5-N7	3.23	114.28	110.58
9	A	600	ATP	C5-C4-N3	-3.05	122.52	126.72
11	F	501	ADP	C4-C5-N7	2.98	113.99	110.58
9	B	600	ATP	C4-C5-N7	2.98	113.98	110.58
11	D	501	ADP	C4-C5-N7	2.90	113.90	110.58
9	C	600	ATP	C4-C5-N7	2.84	113.83	110.58
9	C	600	ATP	C5-C4-N3	-2.83	122.81	126.72
9	C	600	ATP	O2B-PB-O3B	2.82	114.88	107.27
9	A	600	ATP	O2B-PB-O3B	2.66	114.45	107.27
9	B	600	ATP	O2A-PA-O3A	2.65	114.45	107.27
9	B	600	ATP	C5-C4-N3	-2.61	123.12	126.72
11	F	501	ADP	C5-C4-N3	-2.61	123.13	126.72
11	D	501	ADP	C5-C4-N3	-2.60	123.14	126.72
9	A	600	ATP	N3-C4-N9	2.41	131.26	127.17
9	C	600	ATP	O2B-PB-O3A	2.33	113.58	107.27
11	F	501	ADP	O2B-PB-O3A	2.31	112.37	104.64
9	C	600	ATP	N3-C4-N9	2.24	130.98	127.17
11	D	501	ADP	C2'-C3'-C4'	2.17	106.80	102.61
9	B	600	ATP	O2B-PB-O3B	2.14	113.05	107.27
9	A	600	ATP	C2-N1-C6	-2.05	115.36	118.73
9	C	600	ATP	O4'-C1'-C2'	-2.03	102.26	106.62
9	B	600	ATP	N3-C4-N9	2.03	130.61	127.17
11	F	501	ADP	C5-C6-N1	2.02	122.65	117.51
9	A	600	ATP	C5-C6-N1	2.01	122.63	117.51
11	D	501	ADP	C2-N1-C6	-2.01	115.43	118.73

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	E	501	ADP	C5'-O5'-PA-O1A
11	E	501	ADP	C5'-O5'-PA-O2A
11	E	501	ADP	C5'-O5'-PA-O3A
11	E	501	ADP	C4'-C5'-O5'-PA

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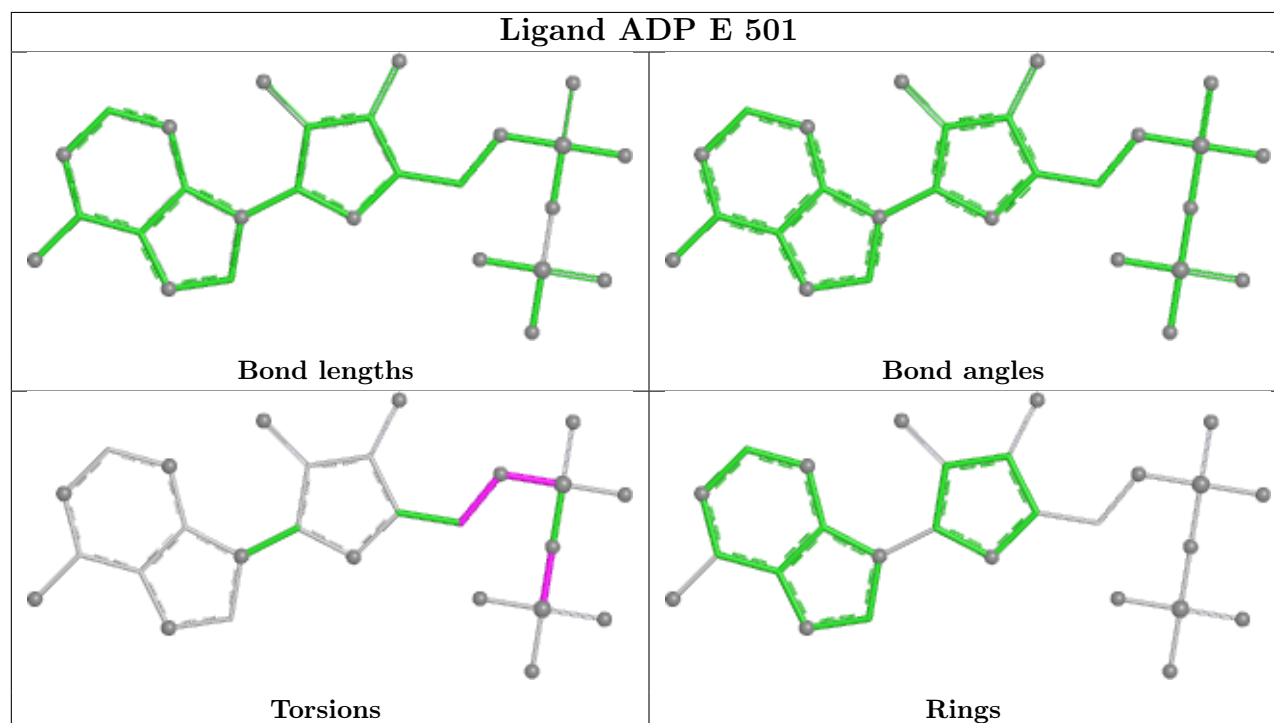
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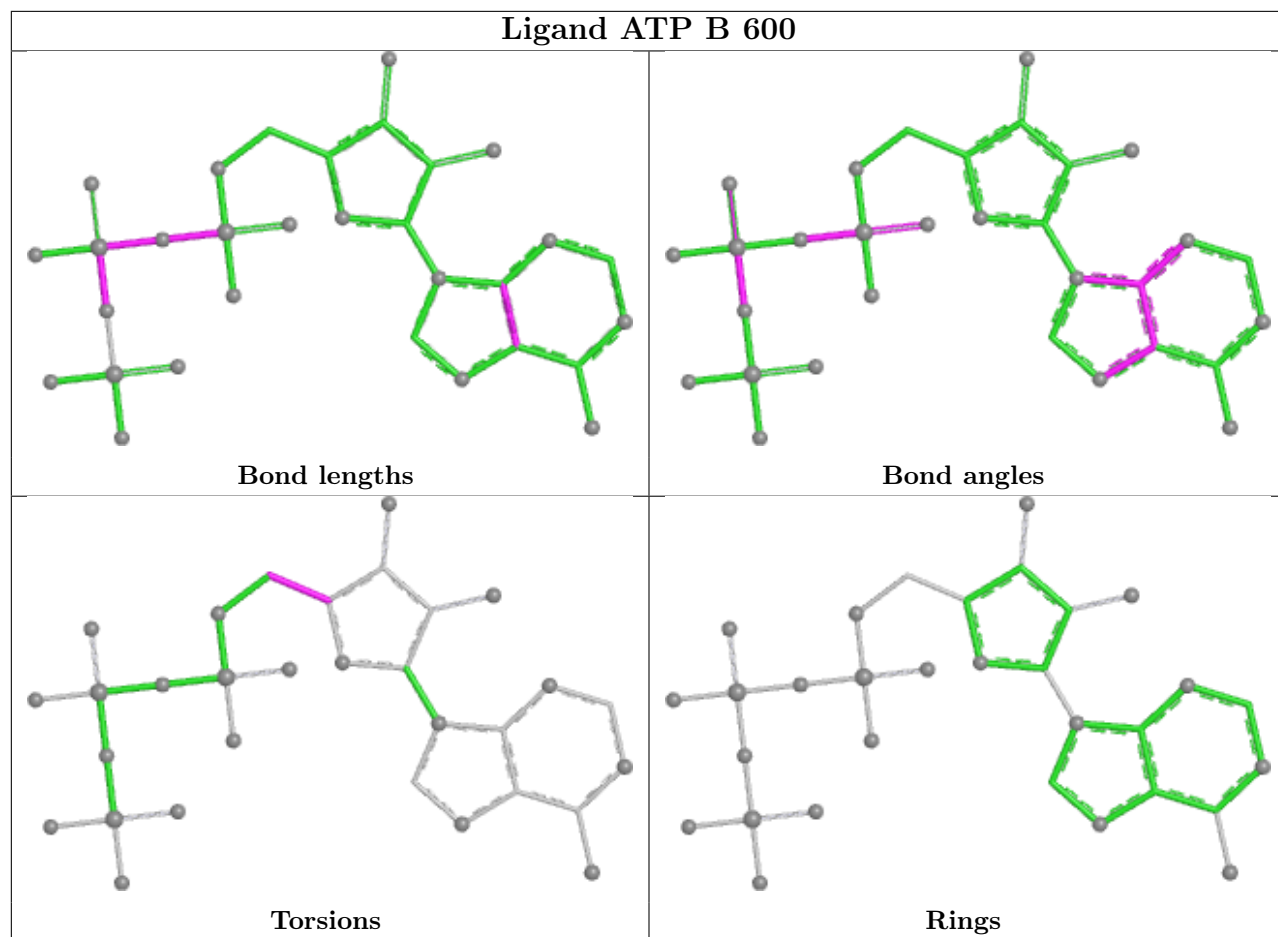
Mol	Chain	Res	Type	Atoms
9	B	600	ATP	O4'-C4'-C5'-O5'
11	E	501	ADP	PA-O3A-PB-O3B
9	A	600	ATP	O4'-C1'-N9-C8

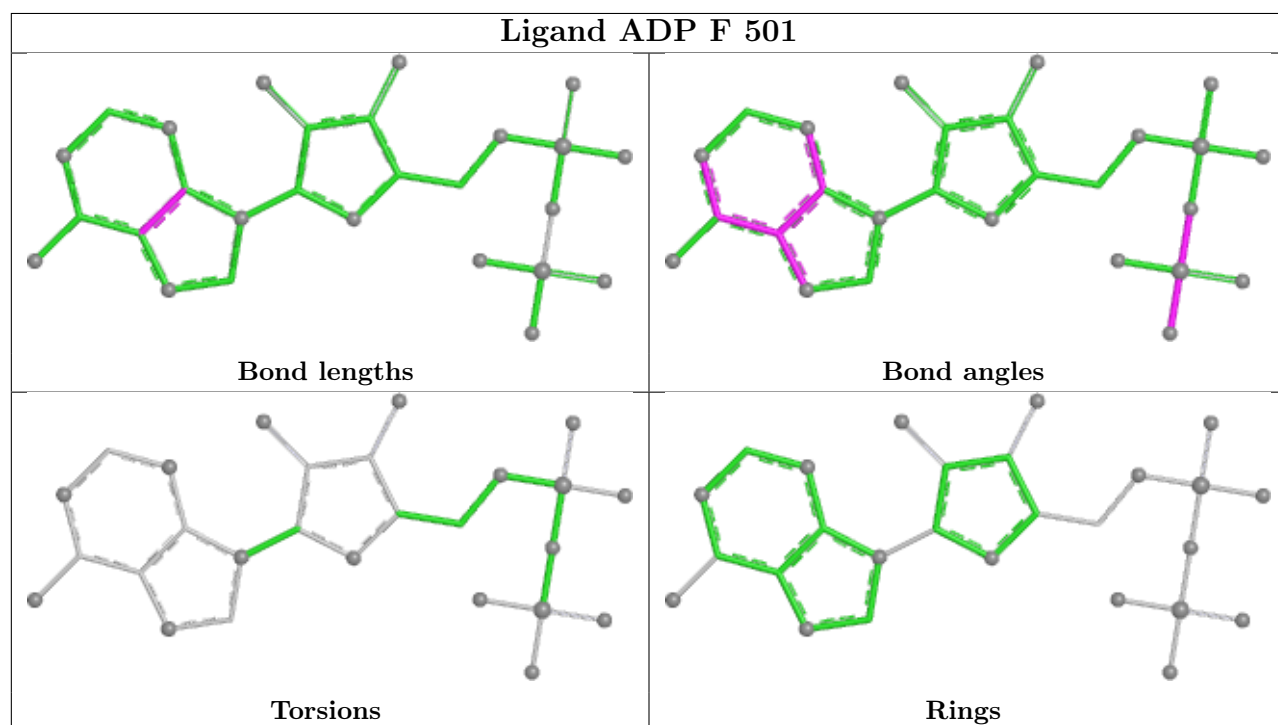
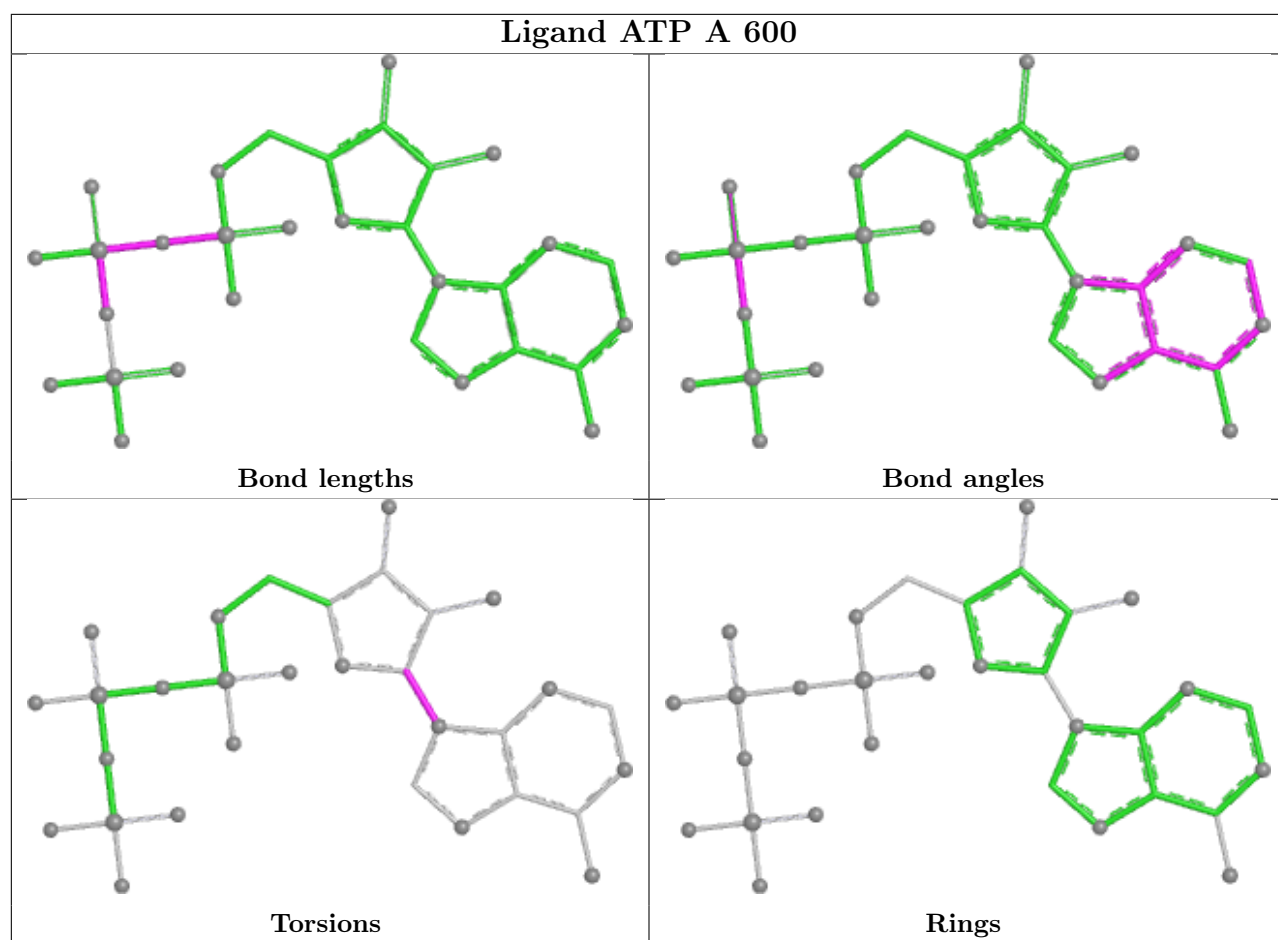
There are no ring outliers.

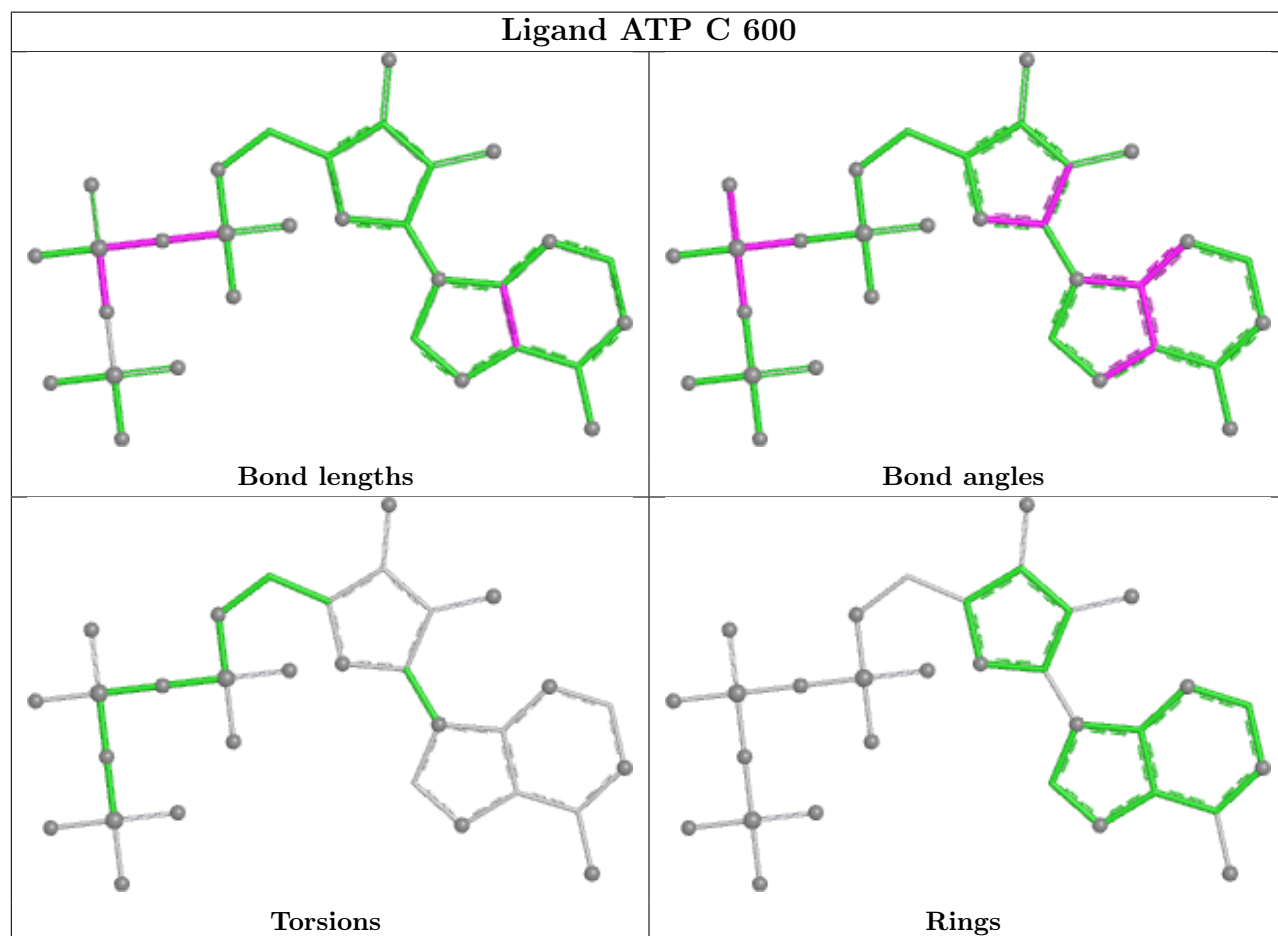
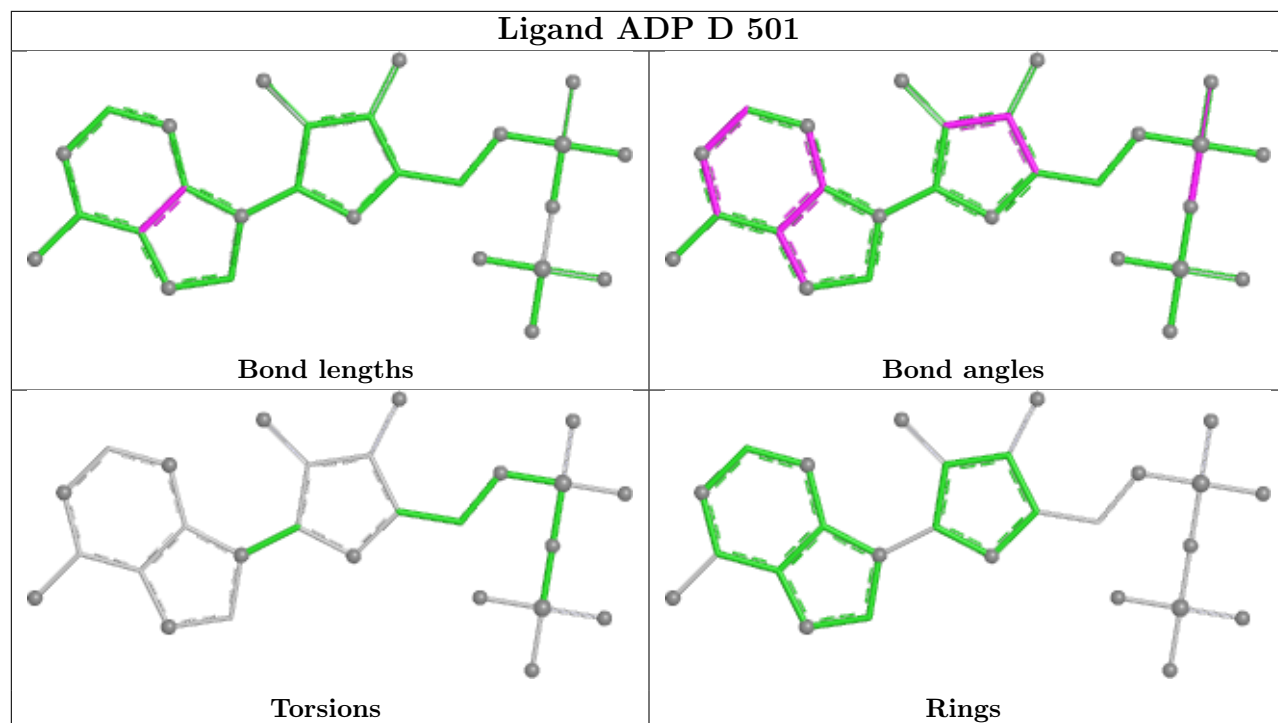
No monomer is involved in short contacts.

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.

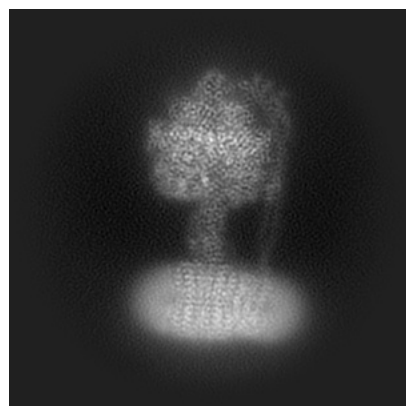
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-70000. These allow visual inspection of the internal detail of the map and identification of artifacts.

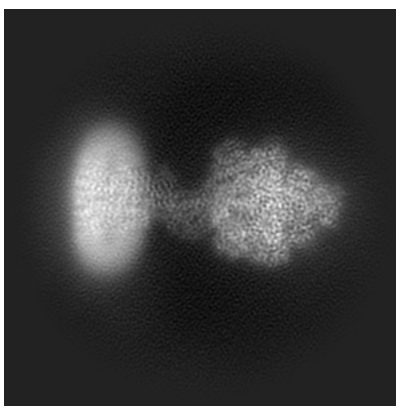
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

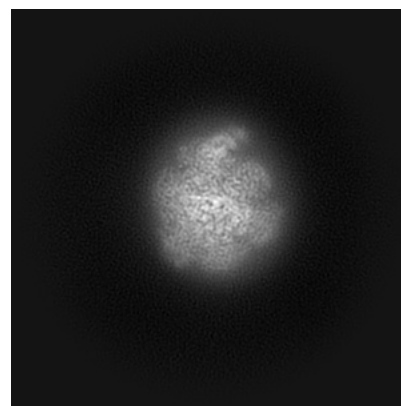
6.1.1 Primary map



X

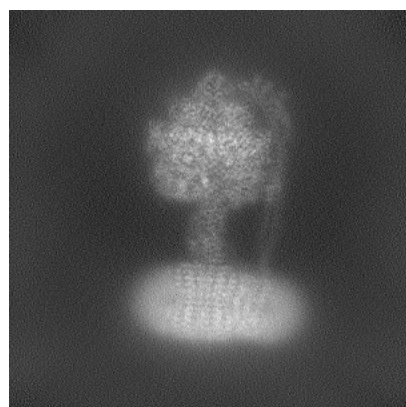


Y

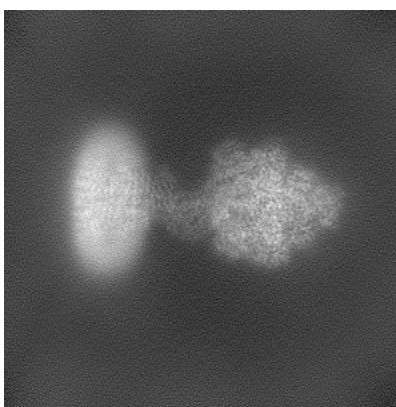


Z

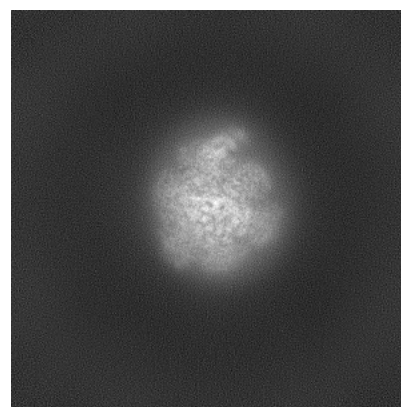
6.1.2 Raw map



X



Y

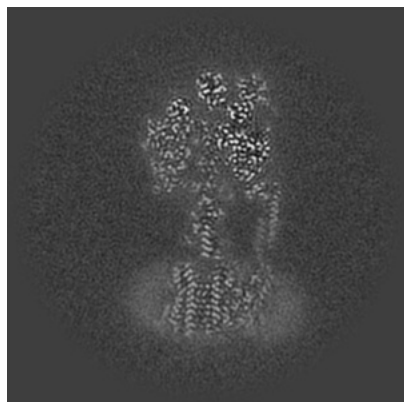


Z

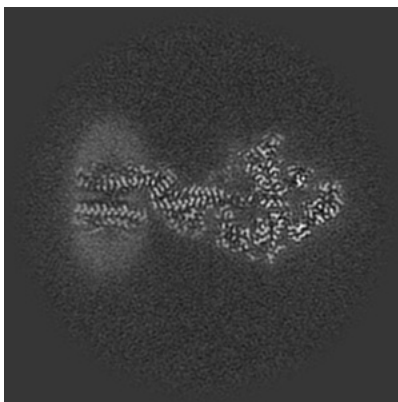
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

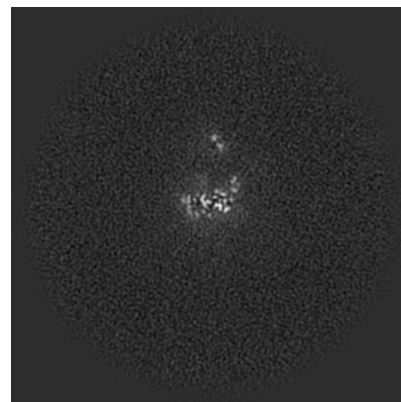
6.2.1 Primary map



X Index: 200

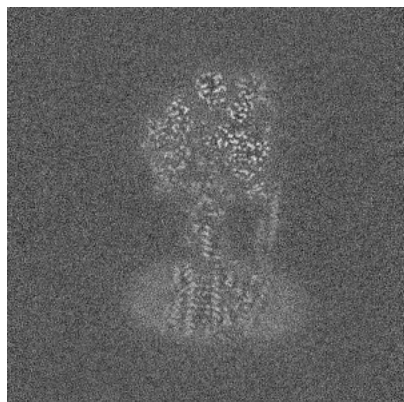


Y Index: 200

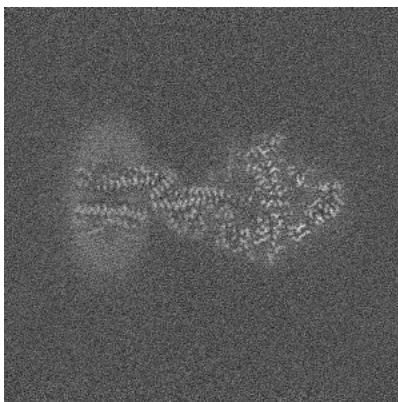


Z Index: 200

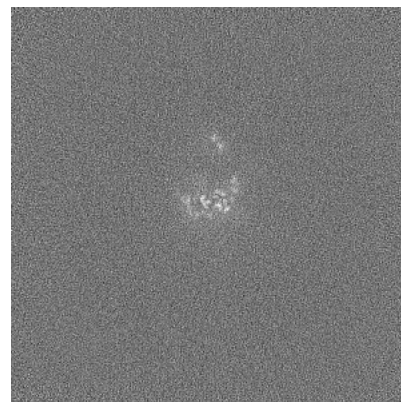
6.2.2 Raw map



X Index: 200



Y Index: 200

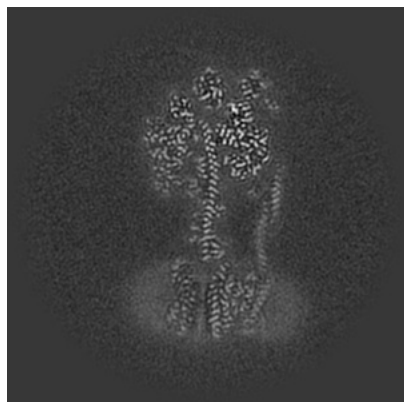


Z Index: 200

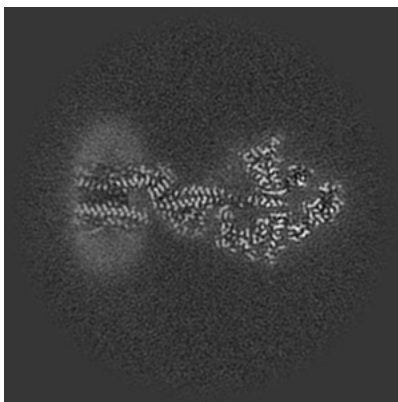
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

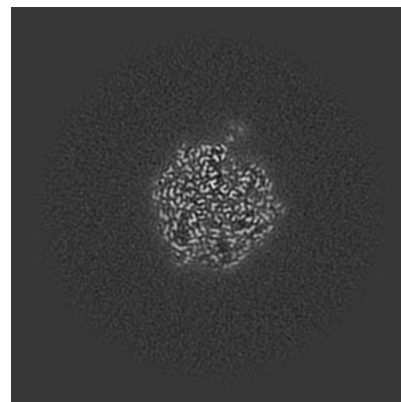
6.3.1 Primary map



X Index: 204

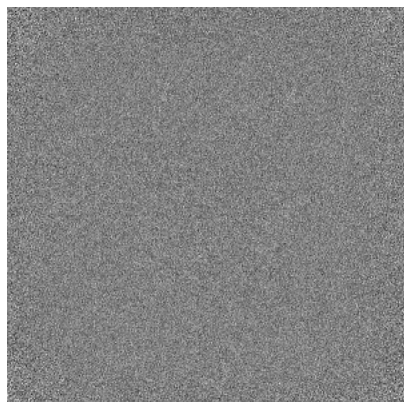


Y Index: 201

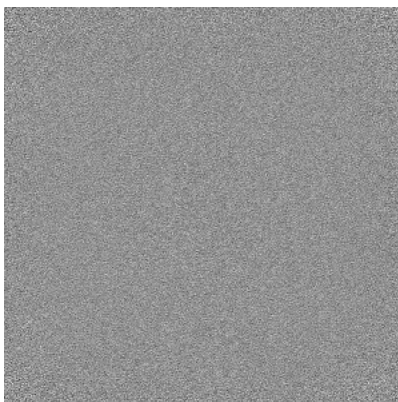


Z Index: 270

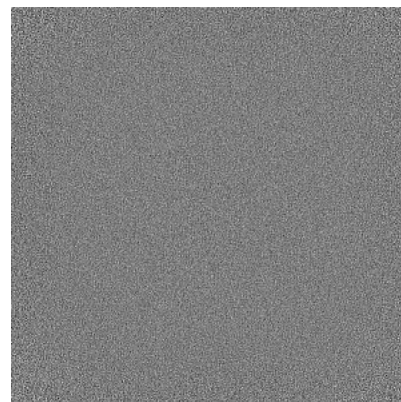
6.3.2 Raw map



X Index: 0



Y Index: 0

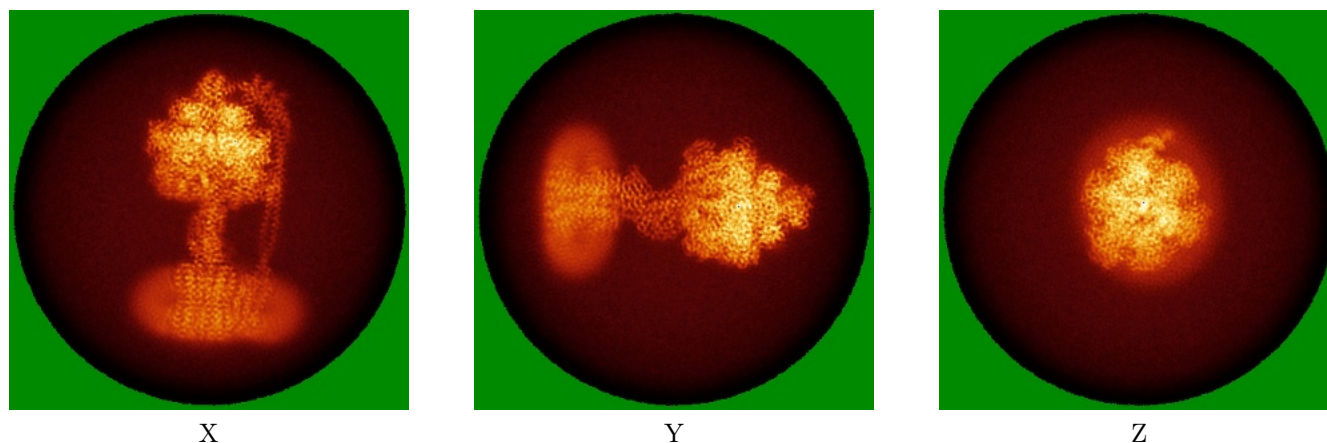


Z Index: 0

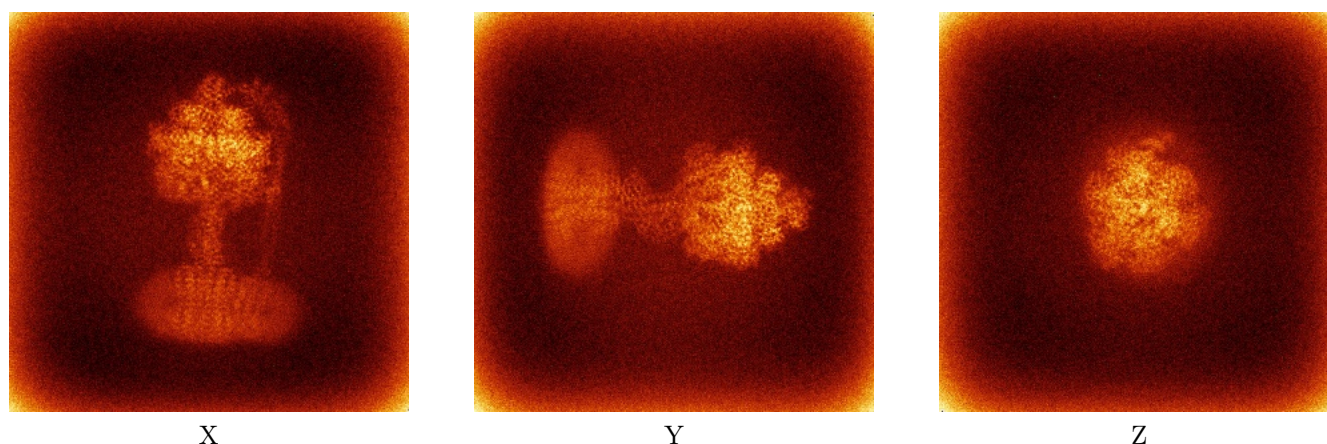
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

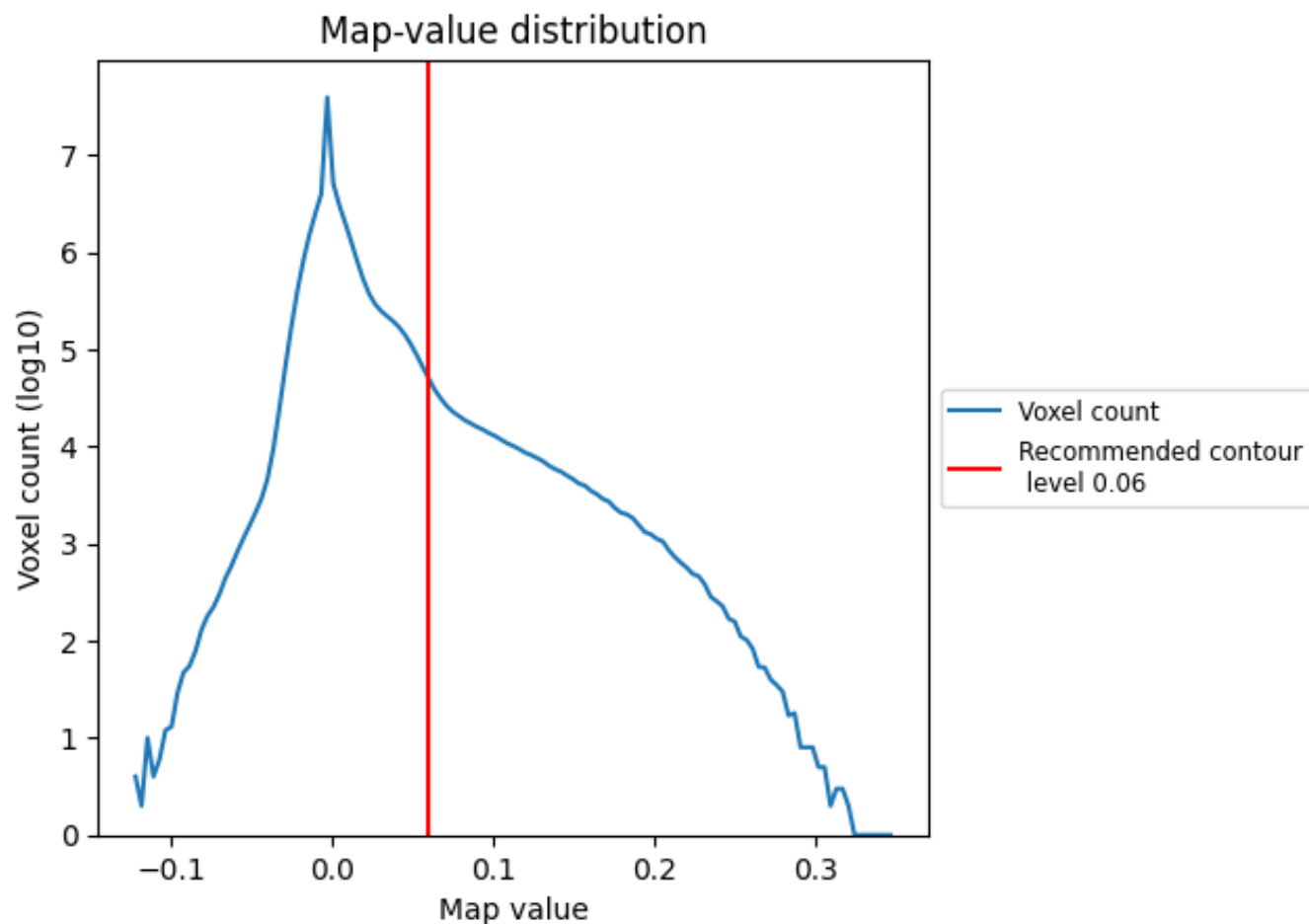
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

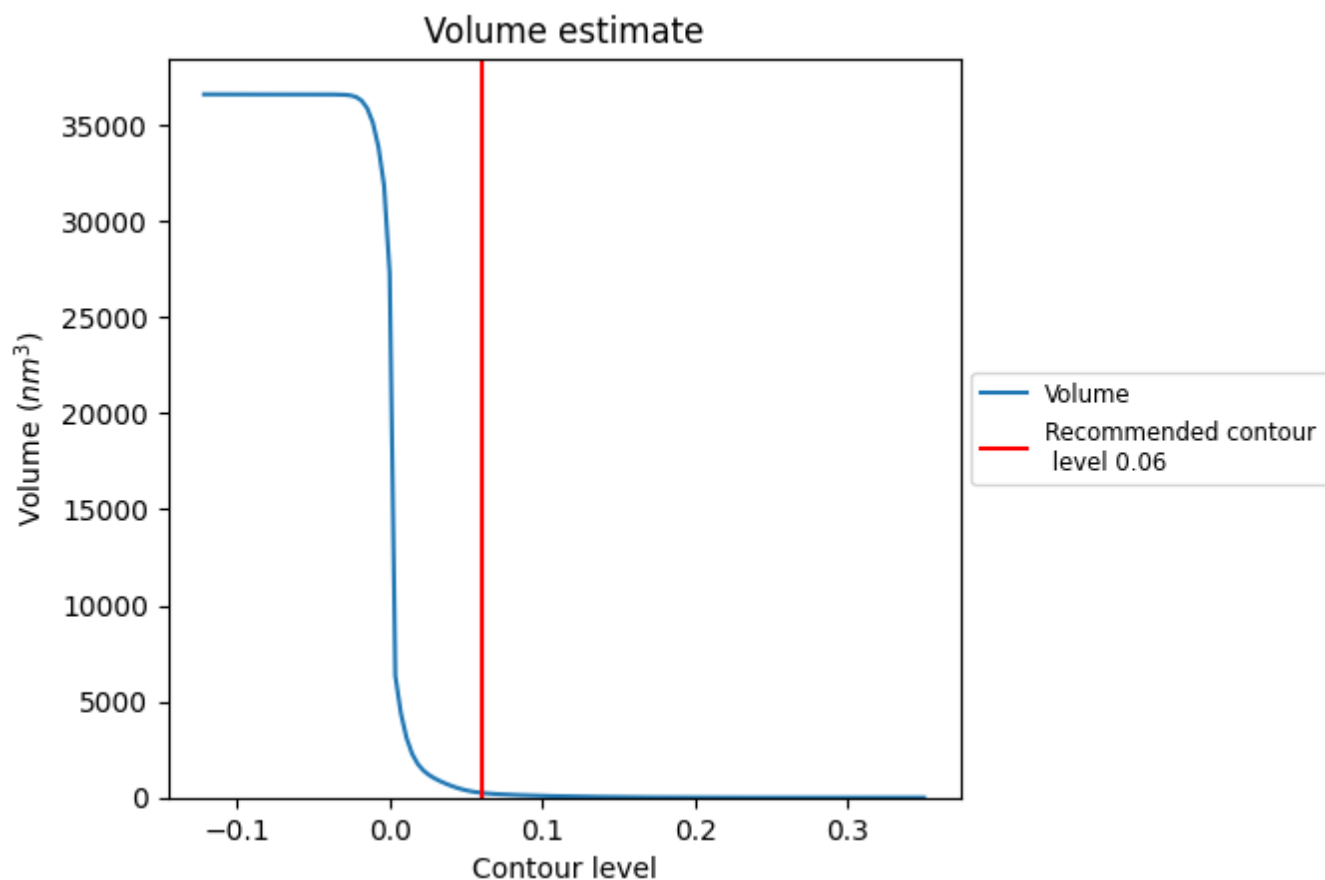
This section contains the results of statistical analysis of the map.

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.

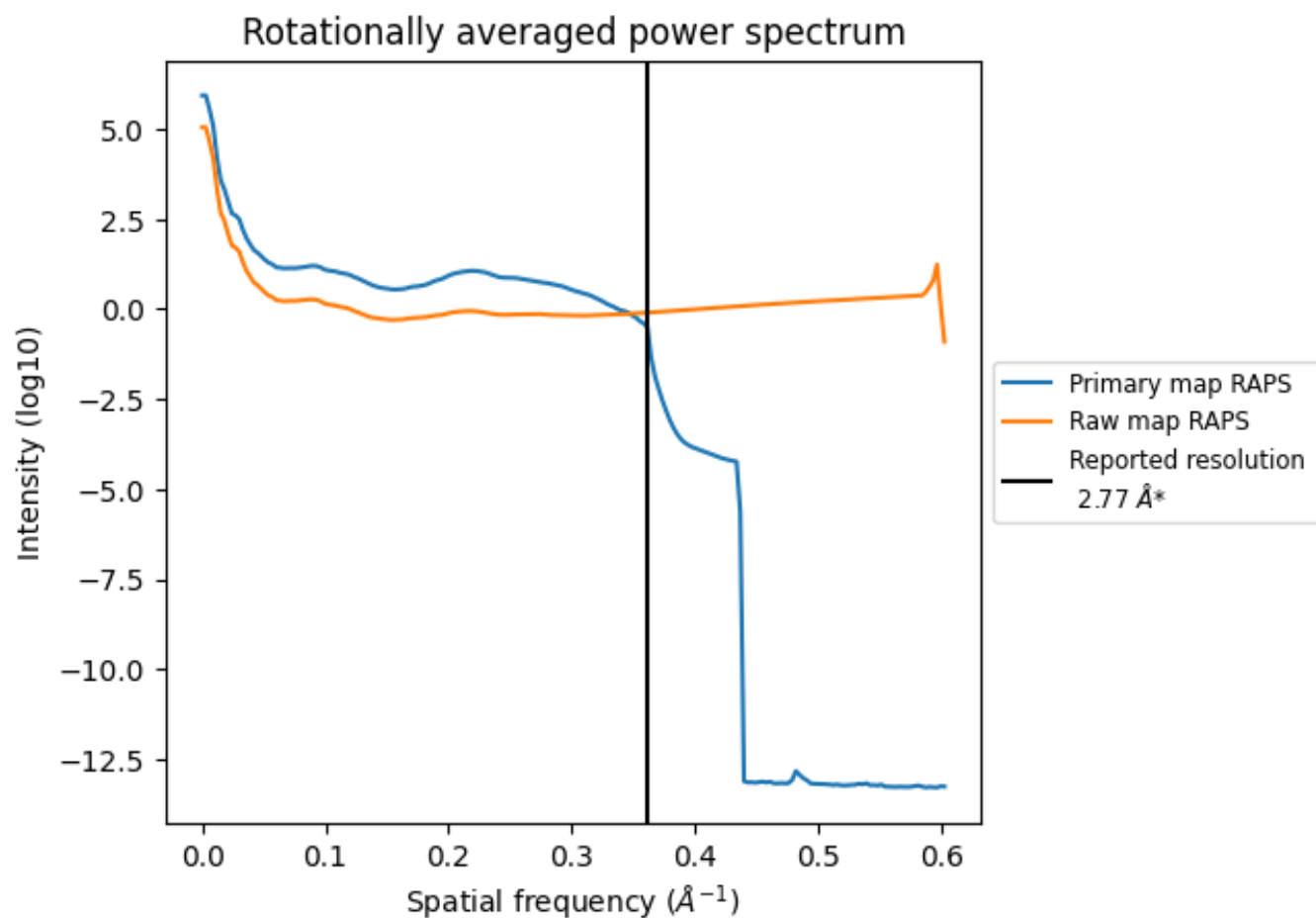
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 245 nm³; this corresponds to an approximate mass of 222 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 ⓘ

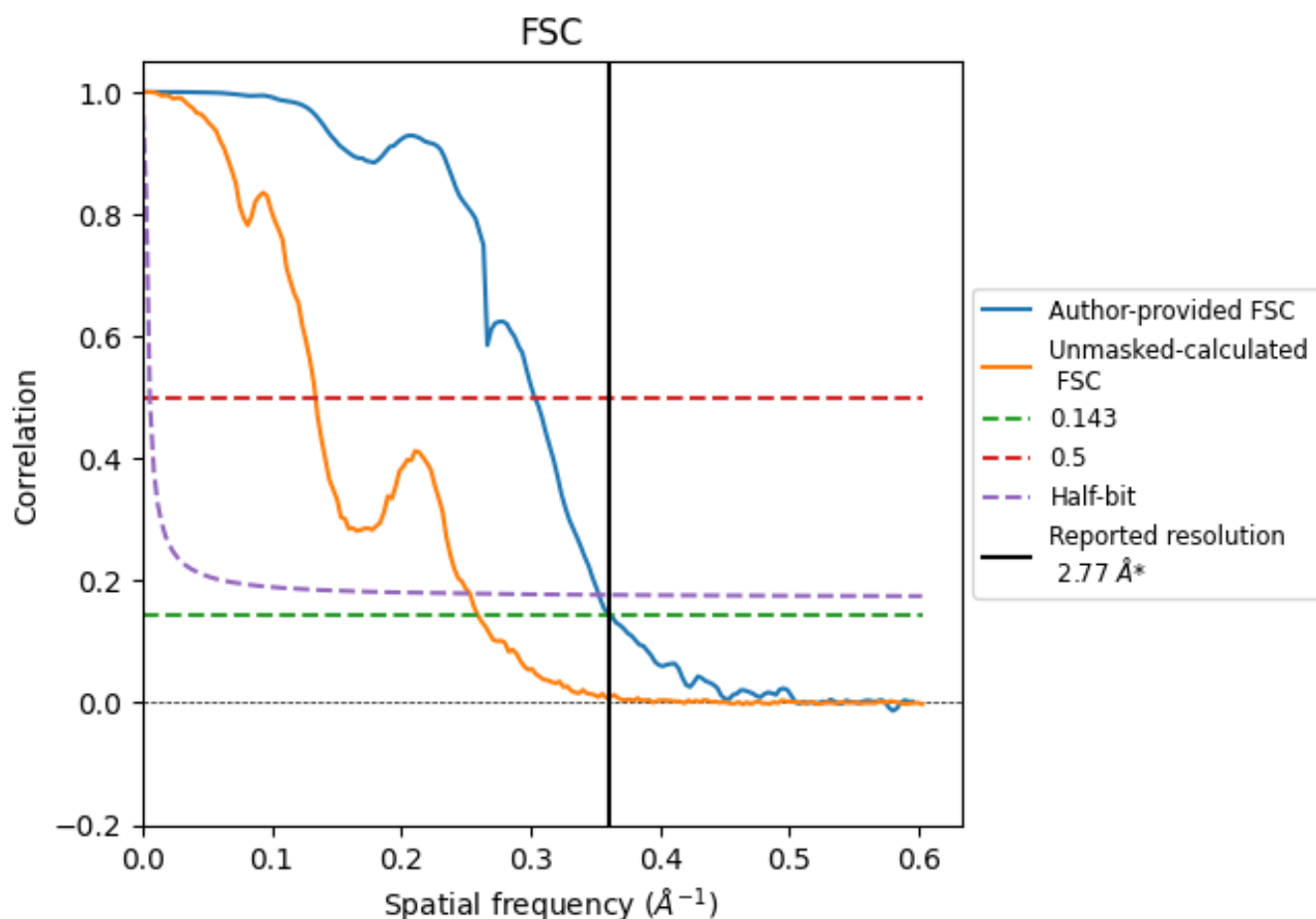


*Reported resolution corresponds to spatial frequency of 0.361 Å⁻¹

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 0.361 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.77	-	-
Author-provided FSC curve	2.77	3.30	2.84
Unmasked-calculated*	3.86	7.47	3.95

*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 3.86 differs from the reported value 2.77 by more than 10 %

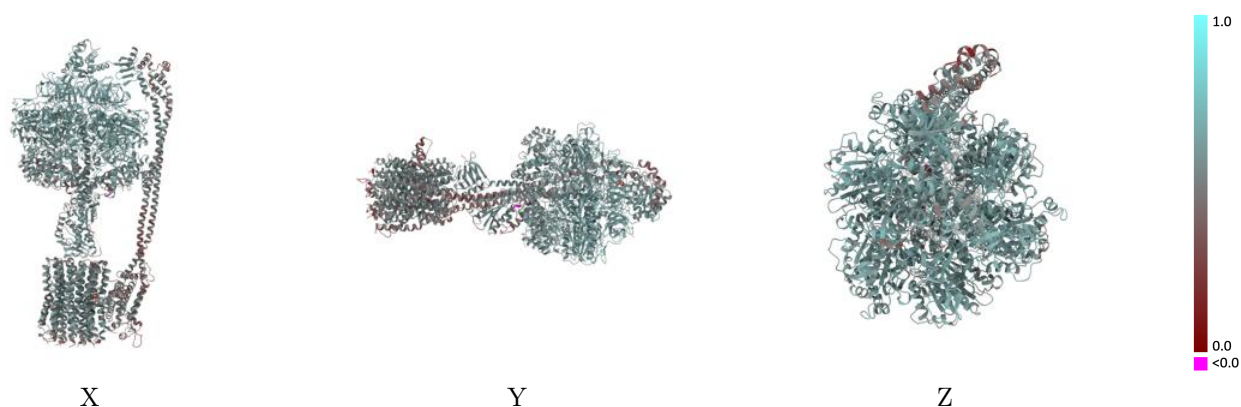
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-70000 and PDB model 9O1A. Per-residue inclusion information can be found in section 3 on page 9.

9.1 Map-model overlay [i](#)

This section was not generated.

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

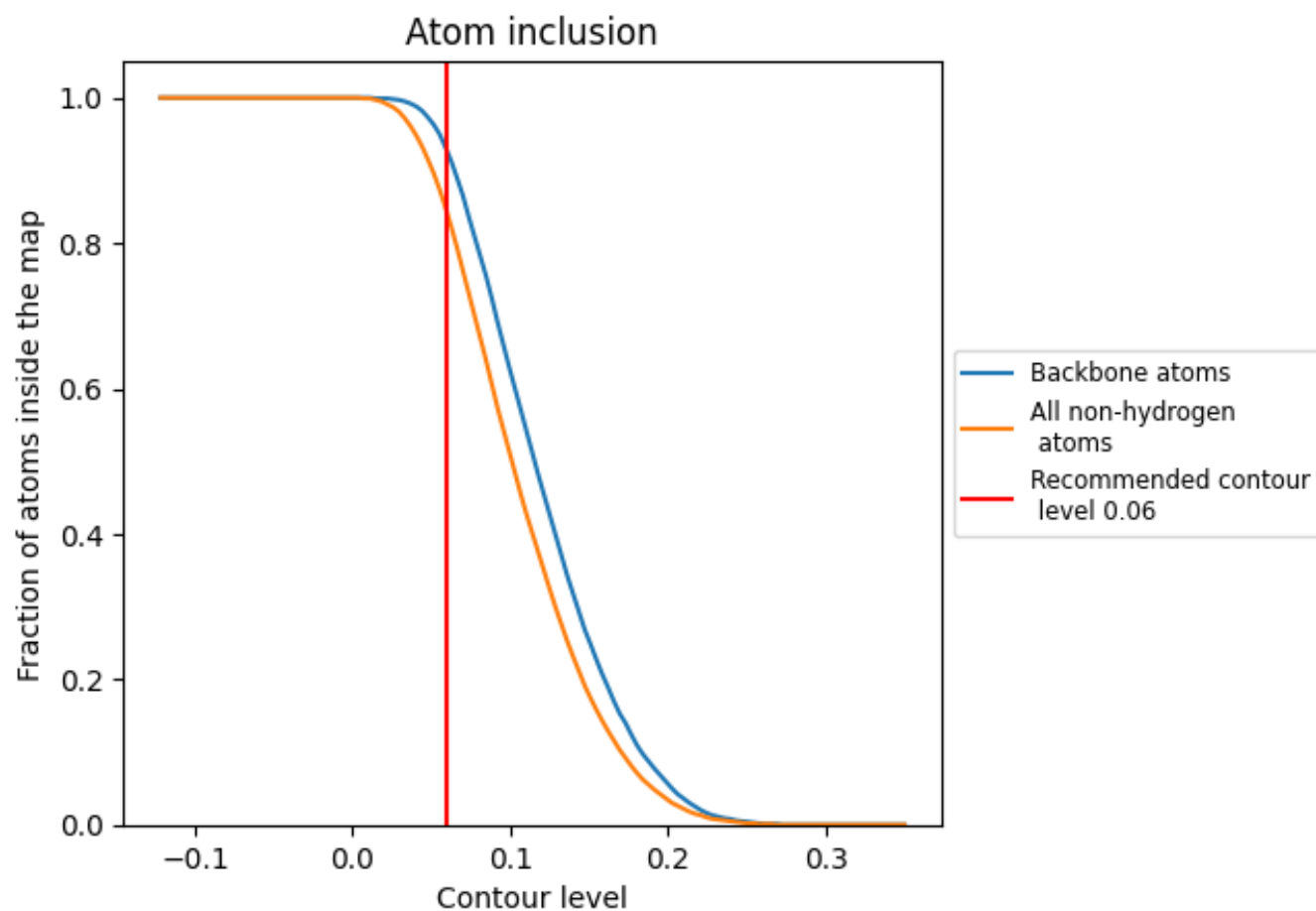


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 84% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8420	 0.5550
A	 0.8880	 0.5850
B	 0.8820	 0.5810
C	 0.9040	 0.5960
D	 0.8750	 0.5790
E	 0.9170	 0.6030
F	 0.9130	 0.6000
G	 0.8230	 0.5390
H	 0.6950	 0.4990
I	 0.8240	 0.5280
J	 0.7870	 0.5200
L	 0.7840	 0.5100
M	 0.8280	 0.5180
N	 0.8400	 0.5370
O	 0.8480	 0.5440
P	 0.8010	 0.5090
Q	 0.7800	 0.5010
R	 0.8290	 0.5150
S	 0.8290	 0.5190
W	 0.8810	 0.5810
X	 0.5460	 0.4370
Y	 0.6340	 0.4470
a	 0.6920	 0.4420

