



Full wwPDB EM Validation Report ⓘ

May 13, 2025 – 09:26 PM EDT

PDB ID : 6OQW / pdb_00006oqw
EMDB ID : EMD-20172
Title : E. coli ATP synthase State 3a
Authors : Stewart, A.G.; Sobti, M.; Walshe, J.L.
Deposited on : 2019-04-29
Resolution : 3.10 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.dev118
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0rc1
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) : 2.43.1

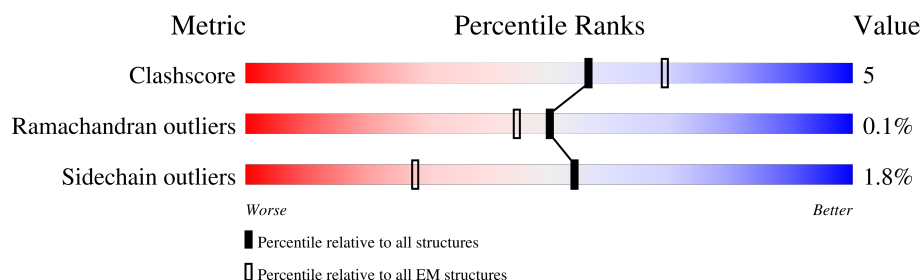
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.10 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
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 ≥ 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	W	177	
2	A	513	
2	B	513	
2	C	513	
3	X	156	
3	Y	156	
4	H	139	
5	G	287	

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Mol	Chain	Length	Quality of chain
6	D	471	
6	E	471	
6	F	471	
7	I	79	
7	J	79	
7	L	79	
7	M	79	
7	N	79	
7	O	79	
7	P	79	
7	Q	79	
7	R	79	
7	S	79	
8	a	271	

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 36901 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 delta.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	W	173	Total	C	N	O	S	0	0
			1317	826	233	252	6		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
W	64	ALA	CYS	conflict	UNP V0ZA15
W	140	ALA	CYS	conflict	UNP V0ZA15

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	510	Total	C	N	O	S	0	0
			3849	2419	679	742	9		
2	B	513	Total	C	N	O	S	0	0
			3877	2438	683	746	10		
2	A	510	Total	C	N	O	S	0	0
			3852	2422	679	742	9		

- Molecule 3 is a protein called ATP synthase subunit b.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	X	153	Total	C	N	O	S	0	0
			1179	737	213	226	3		
3	Y	155	Total	C	N	O	S	0	0
			1171	733	212	223	3		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	21	ALA	CYS	conflict	UNP A0A073FPT7
Y	21	ALA	CYS	conflict	UNP A0A073FPT7

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

Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	136	Total	C	N	O	S	0	0
			1034	649	180	201	4		

- Molecule 5 is a protein called ATP synthase gamma chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	G	284	Total	C	N	O	S	0	0
			2186	1373	382	419	12		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	87	ALA	CYS	conflict	UNP J7RYJ3
G	112	ALA	CYS	conflict	UNP J7RYJ3

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

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	458	Total	C	N	O	S	0	0
			3520	2219	602	685	14		
6	E	460	Total	C	N	O	S	0	0
			3533	2227	604	687	15		
6	D	460	Total	C	N	O	S	0	0
			3533	2227	604	687	15		

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-11	MET	-	initiating methionine	UNP A0A0F6CB56
F	-10	ARG	-	expression tag	UNP A0A0F6CB56
F	-9	GLY	-	expression tag	UNP A0A0F6CB56
F	-8	SER	-	expression tag	UNP A0A0F6CB56
F	-7	HIS	-	expression tag	UNP A0A0F6CB56
F	-6	HIS	-	expression tag	UNP A0A0F6CB56
F	-5	HIS	-	expression tag	UNP A0A0F6CB56
F	-4	HIS	-	expression tag	UNP A0A0F6CB56
F	-3	HIS	-	expression tag	UNP A0A0F6CB56
F	-2	HIS	-	expression tag	UNP A0A0F6CB56
F	-1	GLY	-	expression tag	UNP A0A0F6CB56
F	137	ALA	CYS	conflict	UNP A0A0F6CB56
E	-11	MET	-	initiating methionine	UNP A0A0F6CB56

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-10	ARG	-	expression tag	UNP A0A0F6CB56
E	-9	GLY	-	expression tag	UNP A0A0F6CB56
E	-8	SER	-	expression tag	UNP A0A0F6CB56
E	-7	HIS	-	expression tag	UNP A0A0F6CB56
E	-6	HIS	-	expression tag	UNP A0A0F6CB56
E	-5	HIS	-	expression tag	UNP A0A0F6CB56
E	-4	HIS	-	expression tag	UNP A0A0F6CB56
E	-3	HIS	-	expression tag	UNP A0A0F6CB56
E	-2	HIS	-	expression tag	UNP A0A0F6CB56
E	-1	GLY	-	expression tag	UNP A0A0F6CB56
E	137	ALA	CYS	conflict	UNP A0A0F6CB56
D	-11	MET	-	initiating methionine	UNP A0A0F6CB56
D	-10	ARG	-	expression tag	UNP A0A0F6CB56
D	-9	GLY	-	expression tag	UNP A0A0F6CB56
D	-8	SER	-	expression tag	UNP A0A0F6CB56
D	-7	HIS	-	expression tag	UNP A0A0F6CB56
D	-6	HIS	-	expression tag	UNP A0A0F6CB56
D	-5	HIS	-	expression tag	UNP A0A0F6CB56
D	-4	HIS	-	expression tag	UNP A0A0F6CB56
D	-3	HIS	-	expression tag	UNP A0A0F6CB56
D	-2	HIS	-	expression tag	UNP A0A0F6CB56
D	-1	GLY	-	expression tag	UNP A0A0F6CB56
D	137	ALA	CYS	conflict	UNP A0A0F6CB56

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

Mol	Chain	Residues	Atoms					AltConf	Trace
7	I	77	Total	C	N	O	S	0	0
			557	370	88	92	7		
7	J	77	Total	C	N	O	S	0	0
			557	370	88	92	7		
7	L	77	Total	C	N	O	S	0	0
			557	370	88	92	7		
7	M	77	Total	C	N	O	S	0	0
			557	370	88	92	7		
7	N	77	Total	C	N	O	S	0	0
			557	370	88	92	7		
7	O	77	Total	C	N	O	S	0	0
			554	369	88	90	7		
7	P	77	Total	C	N	O	S	0	0
			557	370	88	92	7		
7	Q	77	Total	C	N	O	S	0	0
			557	370	88	92	7		

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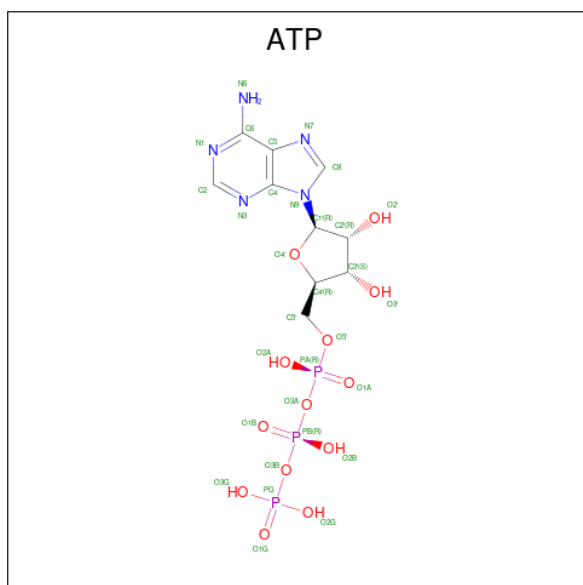
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Mol	Chain	Residues	Atoms					AltConf	Trace
7	R	76	Total	C	N	O	S	0	0
			552	367	87	91	7		
7	S	77	Total	C	N	O	S	0	0
			557	370	88	92	7		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
8	a	266	Total	C	N	O	S	0	0
			2104	1419	326	349	10		

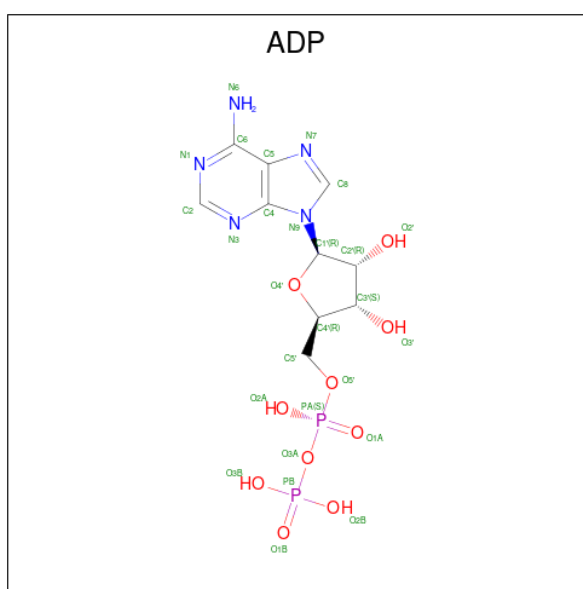
- Molecule 9 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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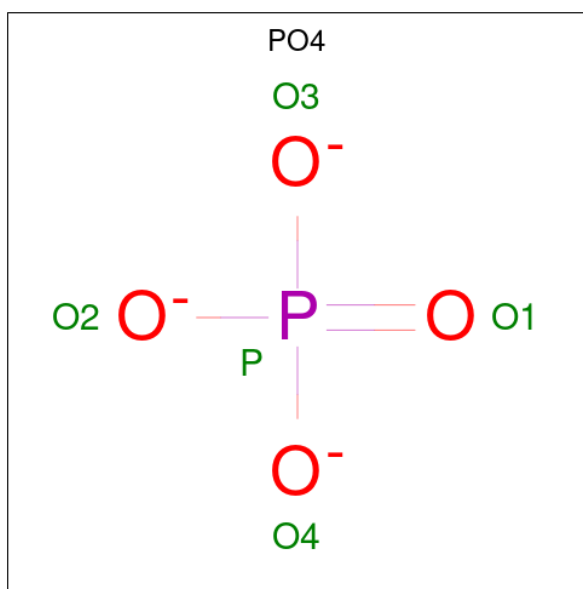
Mol	Chain	Residues	Atoms		AltConf
10	B	1	Total	Mg	0
			1	1	
10	A	1	Total	Mg	0
			1	1	
10	F	1	Total	Mg	0
			1	1	
10	D	1	Total	Mg	0
			1	1	

- Molecule 11 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

- Molecule 12 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).




Mol	Chain	Residues	Atoms			AltConf
			Total	O	P	
12	D	1	5	4	1	0

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.

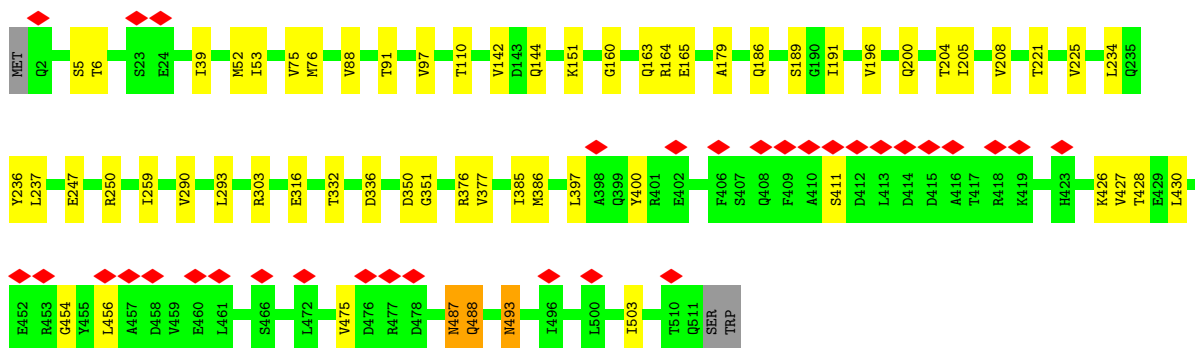
- Molecule 1: ATP synthase subunit delta

Chain W: 



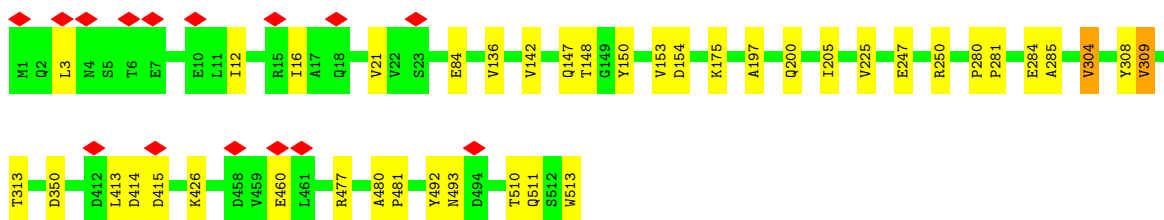
- Molecule 2: ATP synthase subunit alpha

Chain C: 



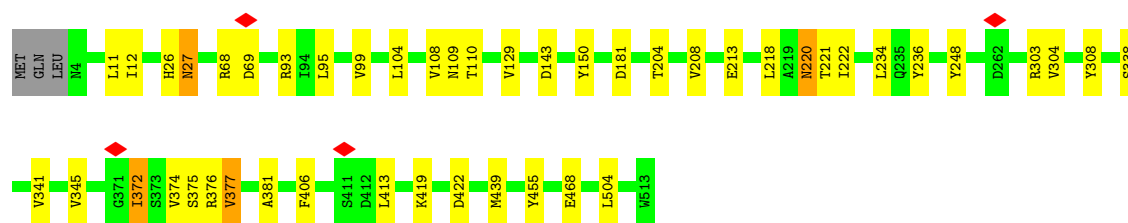
- Molecule 2: ATP synthase subunit alpha

Chain B: 

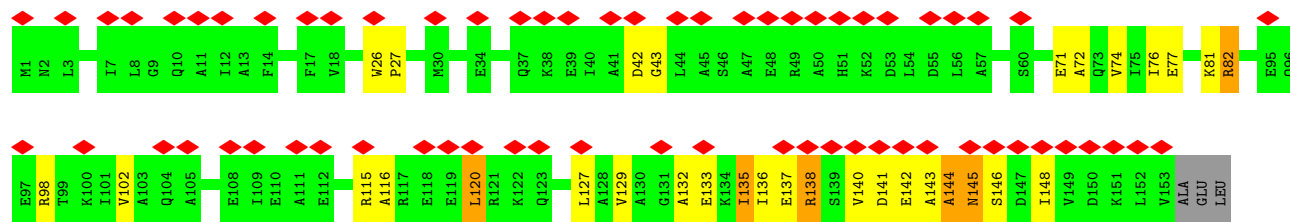
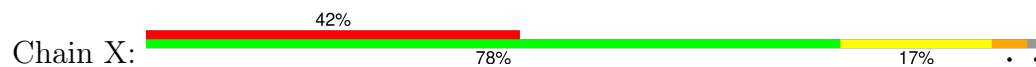


- Molecule 2: ATP synthase subunit alpha

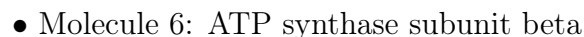
Chain A: 



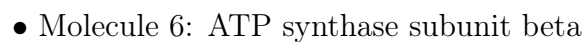
• Molecule 3: ATP synthase subunit b



87% 10% .



90% 7% .



Response	Percentage
Doing a good job	89%
Not doing a good job	8%



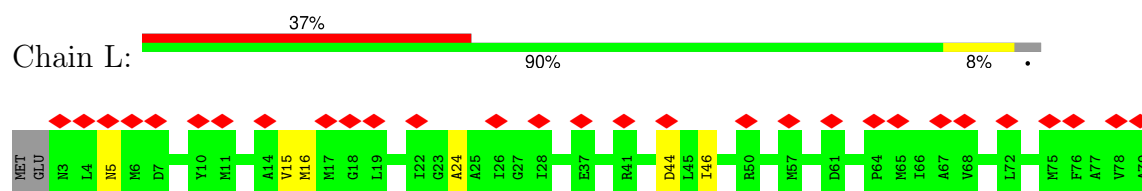
94%



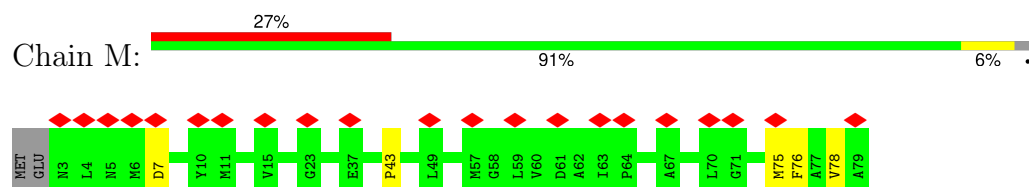
77% 16% ...



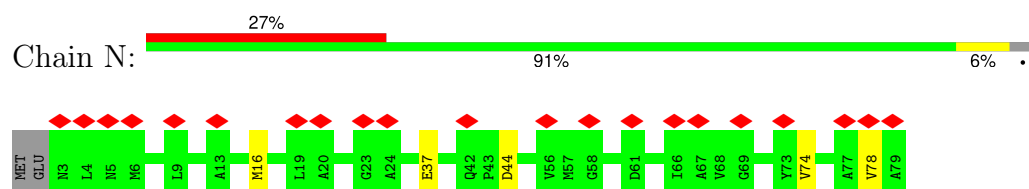
- Molecule 7: ATP synthase subunit c



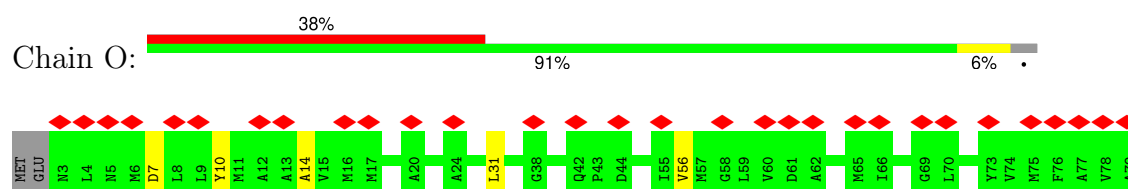
- Molecule 7: ATP synthase subunit c



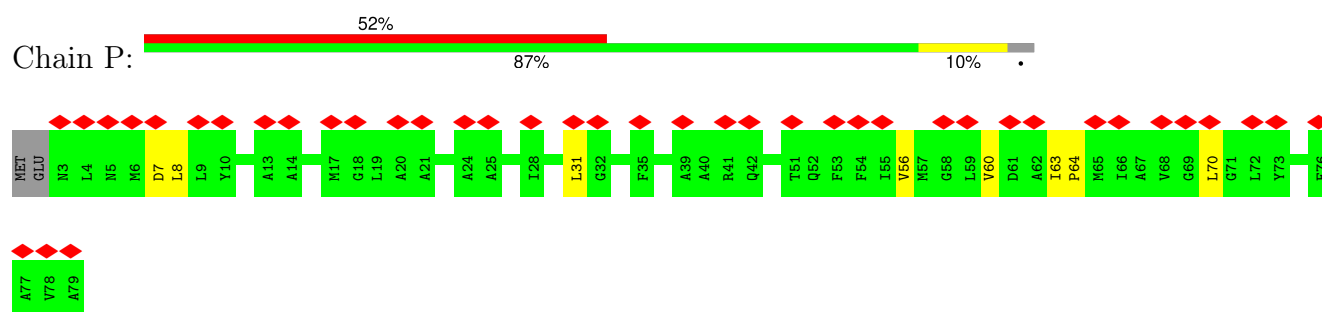
- Molecule 7: ATP synthase subunit c



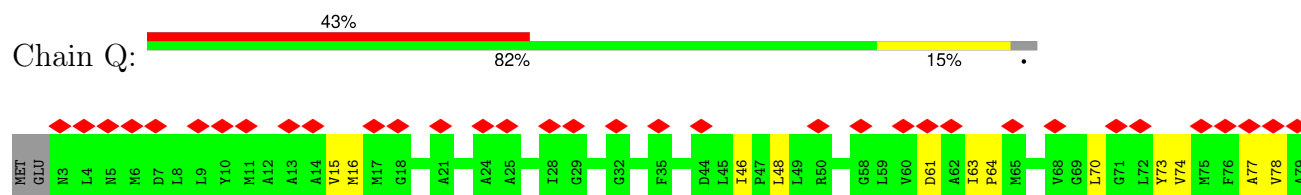
- Molecule 7: ATP synthase subunit c



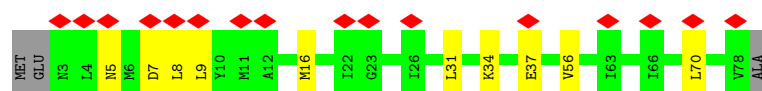
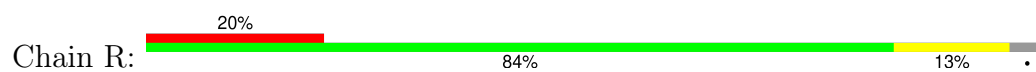
- Molecule 7: ATP synthase subunit c



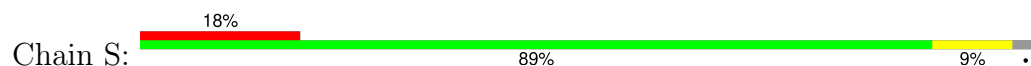
- Molecule 7: ATP synthase subunit c



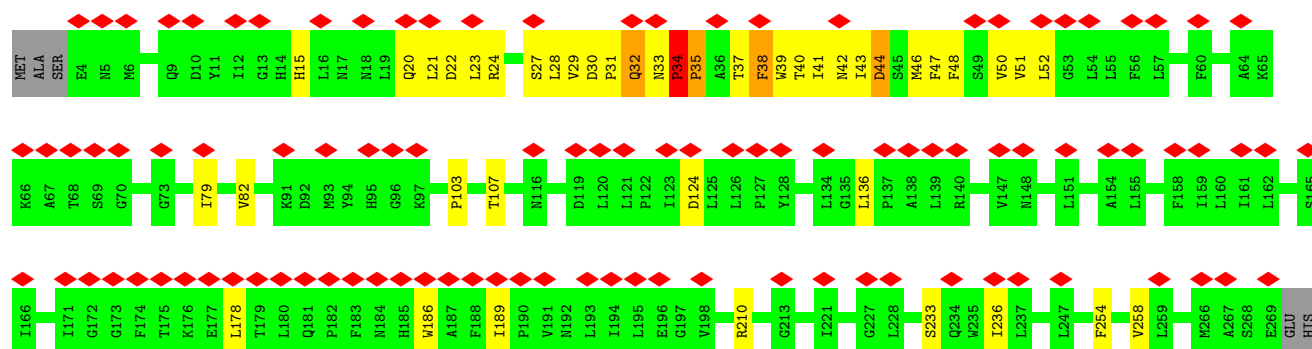
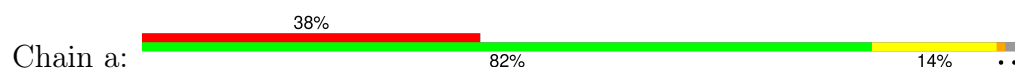
- Molecule 7: ATP synthase subunit c



• Molecule 7: ATP synthase subunit c



• Molecule 8: ATP synthase subunit a



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	78283	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	48	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.092	Depositor
Minimum map value	-0.033	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0171	Depositor
Map size (\AA)	377.65, 377.65, 377.65	wwPDB
Map dimensions	350, 350, 350	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.079, 1.079, 1.079	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	W	0.19	0/1330	0.41	0/1794
2	A	0.33	1/3908 (0.0%)	0.58	1/5288 (0.0%)
2	B	0.25	0/3933	0.46	2/5321 (0.0%)
2	C	0.27	0/3903	0.53	1/5280 (0.0%)
3	X	0.35	0/1187	0.74	6/1596 (0.4%)
3	Y	0.17	0/1179	0.39	0/1589
4	H	0.18	0/1049	0.40	0/1415
5	G	0.37	0/2217	0.66	4/2990 (0.1%)
6	D	0.20	0/3590	0.41	0/4858
6	E	0.17	0/3590	0.37	0/4858
6	F	0.29	0/3577	0.48	3/4841 (0.1%)
7	I	0.19	0/565	0.45	0/764
7	J	0.40	0/565	0.75	3/764 (0.4%)
7	L	0.16	0/565	0.36	0/764
7	M	0.17	0/565	0.36	0/764
7	N	0.18	0/565	0.37	0/764
7	O	0.18	0/562	0.40	0/760
7	P	0.19	0/565	0.39	0/764
7	Q	0.17	0/565	0.38	0/764
7	R	0.19	0/560	0.41	0/757
7	S	0.18	0/565	0.37	0/764
8	a	0.33	0/2164	0.57	4/2955 (0.1%)
All	All	0.26	1/37269 (0.0%)	0.50	24/50414 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	375	SER	CA-CB	-5.45	1.45	1.53

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	377	VAL	N-CA-C	-19.27	95.24	111.81
6	F	264	ALA	N-CA-C	9.04	120.74	111.07
3	X	144	ALA	N-CA-C	-7.50	104.02	113.02
7	J	43	PRO	N-CA-C	-6.00	102.41	111.41
8	a	35	PRO	N-CA-C	-5.96	101.14	110.50
5	G	210	ASP	CA-C-N	5.86	125.40	119.19
5	G	210	ASP	C-N-CA	5.86	125.40	119.19
7	J	47	PRO	CB-CA-C	-5.65	102.24	111.56
3	X	82	ARG	N-CA-C	-5.62	105.06	111.07
5	G	202	SER	N-CA-C	-5.54	106.55	113.15
2	B	280	PRO	CA-C-N	5.43	126.62	119.84
2	B	280	PRO	C-N-CA	5.43	126.62	119.84
8	a	34	PRO	CB-CA-C	-5.37	104.37	110.92
8	a	34	PRO	N-CA-CB	-5.33	97.91	103.08
7	J	47	PRO	N-CA-CB	-5.26	97.73	103.25
2	C	316	GLU	N-CA-C	5.25	120.04	111.37
6	F	262	PRO	CA-C-O	-5.22	115.30	121.67
6	F	264	ALA	CA-C-O	-5.20	115.36	120.82
5	G	191	PRO	CB-CA-C	-5.20	104.14	110.95
3	X	77	GLU	N-CA-C	-5.16	105.83	111.82
3	X	76	ILE	CA-C-N	-5.09	112.57	120.31
3	X	76	ILE	C-N-CA	-5.09	112.57	120.31
3	X	145	ASN	N-CA-C	-5.04	107.15	113.15
8	a	178	LEU	N-CA-C	5.04	119.06	113.01

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	W	1317	0	1343	11	0
2	A	3852	0	3891	27	0
2	B	3877	0	3923	26	0
2	C	3849	0	3895	41	0
3	X	1179	0	1223	36	0
3	Y	1171	0	1198	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	1034	0	1041	10	0
5	G	2186	0	2238	31	0
6	D	3533	0	3548	21	0
6	E	3533	0	3548	17	0
6	F	3520	0	3530	34	0
7	I	557	0	598	2	0
7	J	557	0	598	17	0
7	L	557	0	598	6	0
7	M	557	0	598	5	0
7	N	557	0	598	3	0
7	O	554	0	596	3	0
7	P	557	0	598	6	0
7	Q	557	0	598	9	0
7	R	552	0	593	10	0
7	S	557	0	598	7	0
8	a	2104	0	2196	56	0
9	A	31	0	12	0	0
9	B	31	0	12	1	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	F	1	0	0	0	0
11	D	27	0	12	1	0
11	E	27	0	12	1	0
11	F	27	0	12	3	0
12	D	5	0	0	0	0
All	All	36901	0	37619	353	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 5.

All (353) 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)
8:a:30:ASP:OD2	8:a:31:PRO:HD2	1.41	1.19
3:X:145:ASN:HB3	3:X:148:ILE:HG12	1.38	1.00
3:X:138:ARG:HA	3:X:138:ARG:CZ	1.95	0.96
5:G:210:ASP:OD2	5:G:211:PRO:HD2	1.67	0.94
8:a:30:ASP:OD2	8:a:31:PRO:CD	2.17	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:X:136:ILE:HG23	3:Y:136:ILE:HD12	1.55	0.87
8:a:46:MET:HE3	8:a:46:MET:HA	1.58	0.86
8:a:46:MET:HE3	8:a:46:MET:CA	2.04	0.85
7:J:46:ILE:N	7:J:47:PRO:HD2	1.93	0.84
8:a:38:PHE:HD1	8:a:38:PHE:H	1.27	0.82
8:a:23:LEU:HD21	8:a:41:ILE:HD11	1.60	0.81
3:X:136:ILE:HD12	3:X:138:ARG:HB3	1.62	0.80
3:X:145:ASN:HA	3:X:148:ILE:HG23	1.64	0.79
3:X:145:ASN:CB	3:X:148:ILE:HG12	2.14	0.77
5:G:210:ASP:OD2	5:G:211:PRO:CD	2.34	0.75
8:a:21:LEU:HD12	8:a:27:SER:O	1.87	0.75
6:F:363:ILE:HD11	6:F:396:ILE:HD12	1.69	0.75
8:a:30:ASP:CG	8:a:31:PRO:HD2	2.12	0.75
8:a:46:MET:O	8:a:50:VAL:HG22	1.87	0.74
3:Y:88:ASP:OD2	3:Y:89:GLU:N	2.21	0.73
3:X:133:GLU:HB3	3:X:140:VAL:HG23	1.71	0.72
8:a:32:GLN:N	8:a:32:GLN:HE21	1.89	0.71
1:W:79:MET:HE1	1:W:88:LEU:HD21	1.73	0.71
2:A:218:LEU:HA	2:A:221:THR:HG22	1.73	0.70
8:a:46:MET:HE1	8:a:50:VAL:HG13	1.73	0.70
6:F:136:MET:HE1	6:F:361:GLN:HG3	1.74	0.69
7:P:63:ILE:HG23	7:P:64:PRO:HD3	1.75	0.69
8:a:37:THR:HG22	8:a:39:TRP:HD1	1.58	0.69
8:a:35:PRO:HB2	8:a:40:THR:HG21	1.74	0.68
6:F:265:VAL:HG12	6:F:265:VAL:O	1.93	0.68
8:a:233:SER:O	8:a:236:ILE:HG22	1.93	0.68
2:B:247:GLU:OE2	2:B:250:ARG:NH2	2.27	0.68
2:A:376:ARG:HH21	6:D:151:ALA:HB1	1.58	0.67
2:C:151:LYS:NZ	2:C:430:LEU:O	2.24	0.67
8:a:21:LEU:O	8:a:40:THR:HA	1.95	0.67
2:C:376:ARG:HG2	11:F:500:ADP:O3'	1.95	0.66
2:B:492:TYR:O	2:B:493:ASN:ND2	2.29	0.65
2:C:205:ILE:HD13	2:C:225:VAL:HG23	1.78	0.64
8:a:22:ASP:HA	8:a:40:THR:HG22	1.80	0.64
8:a:103:PRO:O	8:a:107:THR:HG23	1.97	0.64
8:a:46:MET:CE	8:a:50:VAL:HG13	2.28	0.63
3:X:136:ILE:HG12	3:Y:136:ILE:HD11	1.79	0.63
3:X:132:ALA:O	3:X:135:ILE:HG22	1.99	0.62
1:W:33:PHE:O	1:W:37:VAL:HG13	2.00	0.62
7:R:70:LEU:HD23	7:S:75:MET:SD	2.40	0.61
2:C:487:ASN:OD1	2:C:487:ASN:C	2.43	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:200:GLN:HB2	2:C:205:ILE:HD11	1.82	0.61
6:F:23:PRO:O	6:F:45:GLN:NE2	2.33	0.61
5:G:167:LEU:HD12	5:G:168:TYR:N	2.15	0.61
6:F:83:THR:OG1	6:F:88:MET:SD	2.59	0.60
2:B:148:THR:HG21	2:B:153:VAL:HB	1.82	0.60
6:D:345:ASP:O	6:D:349:VAL:HG22	2.02	0.60
8:a:46:MET:HA	8:a:46:MET:CE	2.28	0.60
3:X:138:ARG:HA	3:X:138:ARG:NE	2.17	0.60
6:D:20:ASP:N	6:D:20:ASP:OD1	2.35	0.60
8:a:44:ASP:OD2	8:a:44:ASP:N	2.31	0.59
8:a:30:ASP:HB3	8:a:33:ASN:HB2	1.84	0.59
7:R:70:LEU:HD22	7:S:17:MET:SD	2.43	0.59
8:a:46:MET:CA	8:a:46:MET:CE	2.79	0.59
1:W:79:MET:HE1	1:W:88:LEU:CD2	2.33	0.59
3:X:129:VAL:HG23	3:X:140:VAL:HG11	1.85	0.58
6:F:239:LEU:HD23	6:F:282:ILE:HD12	1.84	0.58
3:Y:141:ASP:OD2	3:Y:144:ALA:N	2.36	0.58
5:G:167:LEU:HD12	5:G:167:LEU:C	2.29	0.58
2:A:218:LEU:HD12	2:A:221:THR:CG2	2.34	0.58
8:a:21:LEU:HD22	8:a:43:ILE:HD11	1.84	0.57
8:a:28:LEU:HD13	8:a:136:LEU:HD22	1.86	0.57
6:F:363:ILE:HD11	6:F:396:ILE:CD1	2.34	0.57
7:J:46:ILE:N	7:J:47:PRO:CD	2.66	0.57
7:J:5:ASN:ND2	7:M:7:ASP:OD2	2.37	0.57
6:D:23:PRO:O	6:D:45:GLN:NE2	2.33	0.57
2:C:493:ASN:HD22	2:C:493:ASN:H	1.52	0.56
8:a:47:PHE:O	8:a:51:VAL:HG23	2.06	0.56
6:F:136:MET:HE1	6:F:361:GLN:CG	2.35	0.56
6:D:233:GLU:OE2	6:D:235:ARG:NH1	2.38	0.56
2:B:413:LEU:H	2:B:413:LEU:HD23	1.69	0.56
6:F:15:VAL:HG11	6:F:68:VAL:HG21	1.88	0.56
7:J:45:LEU:CD1	7:M:43:PRO:HG3	2.36	0.56
3:X:136:ILE:CG2	3:Y:136:ILE:HD12	2.34	0.56
7:J:43:PRO:O	7:J:44:ASP:HB2	2.05	0.56
7:J:45:LEU:C	7:J:47:PRO:HD2	2.31	0.55
5:G:66:TYR:HB3	5:G:188:LEU:HD22	1.87	0.55
2:C:163:GLN:HG3	2:C:377:VAL:HG21	1.88	0.55
5:G:273:THR:HB	6:E:264:ALA:O	2.07	0.55
8:a:28:LEU:CD1	8:a:136:LEU:HD22	2.37	0.55
6:F:25:VAL:HG12	6:F:25:VAL:O	2.07	0.55
2:C:163:GLN:NE2	2:C:165:GLU:OE2	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:154:GLY:HA2	11:F:500:ADP:O2A	2.07	0.54
2:B:148:THR:HG22	2:B:154:ASP:OD1	2.08	0.54
3:X:133:GLU:CB	3:X:140:VAL:HG23	2.37	0.53
2:B:281:PRO:HG2	5:G:276:LEU:CD1	2.38	0.53
8:a:21:LEU:HB2	8:a:43:ILE:HD11	1.91	0.53
1:W:3:PHE:HB2	1:W:6:VAL:HG22	1.90	0.53
2:C:411:SER:O	2:C:411:SER:OG	2.22	0.53
7:J:14:ALA:HB1	7:L:16:MET:HG3	1.91	0.53
3:X:145:ASN:HA	3:X:148:ILE:CG2	2.36	0.53
8:a:34:PRO:CG	8:a:34:PRO:O	2.56	0.53
8:a:23:LEU:HD12	8:a:39:TRP:CA	2.39	0.53
5:G:54:HIS:HD2	5:G:200:HIS:ND1	2.07	0.52
7:R:5:ASN:ND2	7:R:9:LEU:HD13	2.24	0.52
6:F:80:GLY:O	6:F:83:THR:HG22	2.10	0.52
6:E:96:ASP:OD1	6:E:97:MET:N	2.43	0.52
8:a:21:LEU:CD2	8:a:43:ILE:HD11	2.39	0.52
7:R:70:LEU:HD21	7:S:72:LEU:HD23	1.90	0.52
2:B:175:LYS:NZ	9:B:600:ATP:O1B	2.36	0.52
6:D:298:VAL:HG23	6:D:298:VAL:O	2.08	0.52
2:C:493:ASN:N	2:C:493:ASN:ND2	2.57	0.52
2:C:247:GLU:OE2	2:C:250:ARG:NH2	2.43	0.52
2:B:284:GLU:O	2:B:285:ALA:HB3	2.10	0.52
8:a:47:PHE:C	8:a:47:PHE:CD2	2.87	0.51
2:B:477:ARG:HH12	2:B:511:GLN:HE22	1.58	0.51
5:G:178:MET:SD	5:G:178:MET:N	2.83	0.51
8:a:32:GLN:N	8:a:32:GLN:NE2	2.57	0.51
6:F:263:SER:OG	6:F:264:ALA:N	2.42	0.51
7:J:7:ASP:OD1	7:L:5:ASN:ND2	2.35	0.51
2:B:205:ILE:HD13	2:B:225:VAL:HG13	1.92	0.51
2:A:26:HIS:O	2:A:27:ASN:ND2	2.33	0.51
6:D:46:LEU:HD21	6:D:52:ARG:HB2	1.93	0.51
8:a:38:PHE:CD1	8:a:38:PHE:N	2.64	0.51
8:a:29:VAL:HG23	8:a:35:PRO:HD3	1.91	0.51
2:B:281:PRO:HG3	5:G:279:ILE:HG21	1.92	0.51
3:Y:49:ARG:O	3:Y:52:LYS:HG3	2.11	0.51
2:B:426:LYS:HE3	2:B:460:GLU:HA	1.93	0.50
3:X:116:ALA:O	3:X:120:LEU:HD22	2.11	0.50
6:D:308:PRO:HA	6:D:311:THR:HG22	1.93	0.50
8:a:34:PRO:O	8:a:34:PRO:CD	2.58	0.50
7:O:31:LEU:HD21	7:O:56:VAL:HG13	1.92	0.50
8:a:30:ASP:CG	8:a:31:PRO:CD	2.78	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:156:THR:HB	11:F:500:ADP:O1A	2.11	0.50
7:S:65:MET:HE3	8:a:210:ARG:HG3	1.93	0.50
2:C:332:THR:OG1	2:C:336:ASP:O	2.28	0.49
8:a:46:MET:HE3	8:a:46:MET:C	2.37	0.49
3:X:142:GLU:O	3:X:143:ALA:HB3	2.12	0.49
4:H:88:ASP:OD1	4:H:88:ASP:N	2.43	0.49
2:C:76:MET:HE3	2:C:237:LEU:HD12	1.94	0.49
3:X:72:ALA:HB1	3:Y:79:ALA:HB2	1.94	0.49
6:D:191:HIS:HA	6:D:194:THR:HG22	1.94	0.49
6:D:180:GLY:O	6:D:246:ARG:NE	2.44	0.49
8:a:21:LEU:CB	8:a:43:ILE:HD11	2.42	0.49
2:A:377:VAL:HG11	2:A:381:ALA:HB2	1.95	0.49
2:A:220:ASN:CG	2:A:220:ASN:O	2.56	0.49
6:F:307:SER:HB3	6:F:308:PRO:HD3	1.95	0.49
2:C:76:MET:HE2	2:C:234:LEU:HD23	1.95	0.49
2:C:189:SER:OG	2:C:191:ILE:HG22	2.13	0.49
2:A:204:THR:O	2:A:208:VAL:HG13	2.12	0.48
8:a:30:ASP:OD2	8:a:31:PRO:N	2.47	0.48
6:E:224:LEU:HD21	6:E:281:ARG:HB2	1.94	0.48
3:Y:10:GLN:NE2	8:a:124:ASP:OD2	2.37	0.48
7:P:63:ILE:CG2	7:P:64:PRO:HD3	2.43	0.48
2:A:372:ILE:O	2:A:372:ILE:CG2	2.61	0.48
7:J:45:LEU:HD12	7:M:43:PRO:HG3	1.96	0.48
3:X:136:ILE:O	3:X:137:GLU:HB2	2.12	0.48
6:E:265:VAL:HG12	6:E:265:VAL:O	2.13	0.48
2:A:455:TYR:CD2	2:A:504:LEU:HD23	2.48	0.48
3:X:135:ILE:HG23	3:X:136:ILE:N	2.28	0.48
7:P:56:VAL:O	7:P:60:VAL:HG23	2.14	0.48
7:R:34:LYS:HA	7:R:37:GLU:HG2	1.95	0.48
3:X:129:VAL:CG2	3:X:140:VAL:HG11	2.42	0.48
3:X:135:ILE:C	3:X:135:ILE:HD13	2.39	0.48
1:W:112:SER:HB2	1:W:150:VAL:HG12	1.95	0.47
2:A:110:THR:HG21	2:A:234:LEU:HB3	1.96	0.47
7:P:7:ASP:OD2	7:P:8:LEU:N	2.47	0.47
3:Y:99:THR:HA	3:Y:102:VAL:HG22	1.96	0.47
2:B:414:ASP:OD1	2:B:415:ASP:N	2.47	0.47
8:a:46:MET:HE3	8:a:46:MET:O	2.14	0.47
1:W:37:VAL:HG12	2:A:12:ILE:HD11	1.95	0.47
6:E:42:VAL:HA	6:E:53:THR:HG22	1.96	0.47
7:J:46:ILE:C	7:J:48:LEU:H	2.23	0.47
8:a:21:LEU:HB2	8:a:43:ILE:CD1	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:G:210:ASP:OD2	5:G:211:PRO:N	2.48	0.47
6:E:42:VAL:HG13	6:E:51:VAL:CG1	2.44	0.47
6:E:152:GLY:O	11:E:500:ADP:H5'2	2.15	0.47
2:C:204:THR:O	2:C:208:VAL:HG13	2.15	0.47
3:X:133:GLU:HB3	3:X:140:VAL:CG2	2.42	0.47
2:B:281:PRO:HG2	5:G:276:LEU:HD11	1.97	0.47
6:F:144:LYS:NZ	6:F:279:GLN:O	2.42	0.46
7:Q:70:LEU:O	7:Q:74:VAL:HG23	2.15	0.46
6:F:219:VAL:HA	6:F:222:THR:HG22	1.97	0.46
6:E:450:GLU:O	6:E:453:VAL:HG22	2.14	0.46
4:H:40:PRO:O	5:G:206:LEU:N	2.45	0.46
8:a:32:GLN:NE2	8:a:32:GLN:H	2.13	0.46
7:S:8:LEU:HD13	7:S:8:LEU:O	2.16	0.46
5:G:5:ASP:OD1	5:G:5:ASP:N	2.48	0.46
8:a:21:LEU:O	8:a:40:THR:CA	2.62	0.46
2:B:304:VAL:HG21	2:B:308:TYR:CD2	2.50	0.46
5:G:188:LEU:HA	5:G:189:PRO:C	2.41	0.46
2:A:304:VAL:HG11	2:A:308:TYR:CD2	2.50	0.46
3:X:71:GLU:O	3:X:74:VAL:HG22	2.16	0.46
4:H:13:GLN:N	4:H:13:GLN:OE1	2.49	0.46
5:G:105:TRP:O	5:G:110:VAL:HG22	2.15	0.46
5:G:269:GLN:O	5:G:273:THR:HG22	2.15	0.46
6:F:282:ILE:HD11	6:F:292:SER:CB	2.44	0.46
2:A:419:LYS:O	2:A:422:ASP:OD2	2.34	0.46
8:a:23:LEU:HD12	8:a:39:TRP:HA	1.97	0.46
2:C:5:SER:O	2:C:6:THR:OG1	2.30	0.46
2:A:104:LEU:HA	2:A:222:ILE:HG12	1.97	0.46
7:L:44:ASP:OD2	7:L:44:ASP:N	2.45	0.46
6:F:27:ASP:O	6:F:42:VAL:HG13	2.16	0.45
1:W:28:GLN:OE1	1:W:99:ARG:NH1	2.48	0.45
1:W:166:ARG:NH1	2:B:21:VAL:HG12	2.30	0.45
2:B:12:ILE:HG21	3:X:127:LEU:HD13	1.97	0.45
5:G:274:GLN:O	5:G:278:GLU:HG2	2.16	0.45
7:J:45:LEU:HD23	7:J:45:LEU:HA	1.79	0.45
5:G:253:GLY:O	5:G:257:ILE:HG12	2.16	0.45
2:C:163:GLN:HG2	2:C:164:ARG:H	1.82	0.45
2:B:136:VAL:HG22	6:E:183:THR:HG23	1.98	0.45
3:X:136:ILE:HG12	3:Y:136:ILE:CD1	2.46	0.45
2:C:163:GLN:HG3	2:C:377:VAL:CG2	2.46	0.45
2:C:385:ILE:HG23	2:C:386:MET:H	1.82	0.45
3:Y:26:TRP:N	3:Y:27:PRO:CD	2.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:a:48:PHE:O	8:a:52:LEU:HG	2.17	0.45
6:D:379:MET:HE1	6:D:390:VAL:HG11	1.98	0.45
7:J:46:ILE:C	7:J:48:LEU:N	2.75	0.45
8:a:186:TRP:HA	8:a:189:ILE:HG13	1.98	0.45
2:A:95:LEU:HG	2:A:129:VAL:HG21	1.99	0.45
5:G:60:LEU:HD12	5:G:60:LEU:O	2.17	0.45
7:N:74:VAL:HA	7:N:78:VAL:HG22	1.99	0.45
2:A:218:LEU:HA	2:A:221:THR:CG2	2.45	0.45
6:E:191:HIS:HA	6:E:194:THR:HG22	1.99	0.45
6:D:189:PHE:CE1	6:D:204:LEU:HD21	2.52	0.44
7:J:41:ARG:HG2	7:J:42:GLN:HG3	1.99	0.44
1:W:75:LEU:HG	1:W:79:MET:HE2	1.98	0.44
2:C:397:LEU:HA	2:C:400:TYR:HB3	1.99	0.44
3:X:136:ILE:HG23	3:Y:136:ILE:CD1	2.38	0.44
2:A:108:VAL:HG12	2:A:109:ASN:O	2.18	0.44
2:A:143:ASP:OD1	2:A:303:ARG:NH1	2.44	0.44
2:A:439:MET:HE1	2:A:468:GLU:OE2	2.17	0.44
7:R:56:VAL:HG23	7:S:53:PHE:CZ	2.53	0.44
2:C:385:ILE:HG23	2:C:386:MET:N	2.32	0.44
2:A:99:VAL:HG21	2:A:248:TYR:HB2	2.00	0.44
3:Y:88:ASP:OD2	3:Y:88:ASP:C	2.60	0.44
2:A:11:LEU:C	2:A:11:LEU:HD23	2.43	0.44
5:G:222:TYR:O	5:G:223:VAL:C	2.59	0.44
8:a:20:GLN:H	8:a:29:VAL:HG12	1.83	0.44
8:a:47:PHE:CE2	8:a:51:VAL:HG21	2.52	0.44
6:D:159:MET:HE3	6:D:240:PHE:HB3	2.00	0.44
5:G:60:LEU:HD12	5:G:60:LEU:C	2.42	0.44
2:C:144:GLN:OE1	2:C:303:ARG:NH1	2.51	0.44
6:F:265:VAL:O	6:F:265:VAL:CG1	2.62	0.44
6:F:400:LEU:HD12	6:F:427:PHE:CZ	2.53	0.43
2:B:84:GLU:OE1	6:D:19:GLN:NE2	2.51	0.43
6:E:172:GLY:HA3	6:E:236:ASP:O	2.18	0.43
6:E:285:THR:HG22	6:E:286:LYS:N	2.33	0.43
7:J:63:ILE:HB	7:J:64:PRO:HD3	2.00	0.43
5:G:227:VAL:O	5:G:228:TYR:C	2.59	0.43
2:C:290:VAL:O	2:C:293:LEU:HB3	2.18	0.43
6:F:298:VAL:O	6:F:298:VAL:HG23	2.17	0.43
7:J:43:PRO:O	7:J:44:ASP:CB	2.66	0.43
8:a:79:ILE:O	8:a:82:VAL:HG12	2.19	0.43
2:A:406:PHE:HB3	2:A:413:LEU:HD11	1.99	0.43
7:R:7:ASP:OD1	7:R:8:LEU:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:D:372:ASP:OD2	6:D:373:ILE:N	2.51	0.43
7:R:5:ASN:HD21	7:R:9:LEU:HD13	1.82	0.43
6:E:307:SER:HB3	6:E:308:PRO:HD3	2.01	0.43
2:C:53:ILE:CG2	2:C:88:VAL:CG1	2.96	0.43
7:Q:73:TYR:CD1	7:Q:77:ALA:HB3	2.54	0.43
2:C:165:GLU:OE2	2:C:351:GLY:HA3	2.18	0.42
3:X:26:TRP:HB3	3:X:27:PRO:HD3	2.01	0.42
5:G:256:LEU:HD22	6:F:376:ILE:HD11	2.01	0.42
7:R:16:MET:HG3	7:S:14:ALA:HB1	2.01	0.42
7:R:31:LEU:HD21	7:R:56:VAL:HG13	2.01	0.42
6:D:207:GLY:HA3	6:D:219:VAL:HG21	2.01	0.42
8:a:30:ASP:HB3	8:a:33:ASN:CB	2.48	0.42
2:C:427:VAL:HG13	2:C:428:THR:N	2.34	0.42
2:C:454:GLY:C	2:C:456:LEU:H	2.28	0.42
8:a:34:PRO:O	8:a:34:PRO:HG2	2.20	0.42
2:C:191:ILE:HG23	2:C:191:ILE:O	2.20	0.42
2:C:488:GLN:HA	2:C:488:GLN:NE2	2.34	0.42
3:X:138:ARG:NE	3:X:138:ARG:CA	2.80	0.42
6:F:239:LEU:HD23	6:F:282:ILE:CD1	2.49	0.42
6:D:166:ILE:HG22	6:D:166:ILE:O	2.20	0.42
5:G:60:LEU:O	5:G:60:LEU:CG	2.66	0.42
6:E:90:VAL:HG21	6:E:215:ASN:HA	2.00	0.42
8:a:24:ARG:HA	8:a:24:ARG:HD3	1.75	0.42
2:C:186:GLN:HB3	2:C:191:ILE:HG23	2.02	0.42
2:A:341:VAL:O	2:A:345:VAL:HG23	2.19	0.42
5:G:227:VAL:O	5:G:230:GLY:N	2.52	0.42
6:F:163:ILE:O	6:F:167:ALA:HB3	2.20	0.42
6:F:261:MET:HE2	6:F:261:MET:HB3	1.85	0.42
2:C:179:ALA:HB1	2:C:259:ILE:HG21	2.02	0.42
2:A:213:GLU:OE1	2:A:213:GLU:HA	2.19	0.42
3:X:81:LYS:O	3:X:82:ARG:C	2.62	0.42
3:X:98:ARG:O	3:X:102:VAL:HG22	2.19	0.42
6:D:224:LEU:HD21	6:D:282:ILE:HG12	2.02	0.42
2:B:510:THR:C	2:B:511:GLN:HG2	2.44	0.42
3:X:140:VAL:HG12	3:X:141:ASP:N	2.34	0.42
3:X:145:ASN:O	3:X:148:ILE:HG13	2.20	0.42
2:C:488:GLN:CA	2:C:488:GLN:HE21	2.32	0.42
6:F:138:PRO:O	6:F:341:SER:OG	2.28	0.42
6:F:146:GLY:HA3	6:F:315:LEU:HD13	2.00	0.42
7:L:46:ILE:HD12	7:Q:48:LEU:HD21	2.00	0.42
7:Q:63:ILE:CG2	7:Q:64:PRO:HD3	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:493:ASN:H	2:C:493:ASN:ND2	2.13	0.42
7:I:75:MET:HE1	7:P:70:LEU:HD11	2.02	0.42
7:J:51:THR:O	7:J:55:ILE:HG12	2.20	0.42
7:L:15:VAL:HG13	7:L:16:MET:N	2.35	0.42
2:B:309:VAL:O	2:B:313:THR:HG22	2.20	0.41
2:A:150:TYR:OH	2:A:181:ASP:OD2	2.34	0.41
5:G:42:ARG:N	5:G:43:PRO:CD	2.83	0.41
2:B:12:ILE:O	2:B:16:ILE:HG12	2.21	0.41
2:C:350:ASP:O	2:C:376:ARG:HG3	2.21	0.41
2:C:487:ASN:OD1	2:C:487:ASN:O	2.39	0.41
2:A:376:ARG:HG2	11:D:500:ADP:H5'1	2.01	0.41
2:C:144:GLN:O	2:C:160:GLY:HA2	2.21	0.41
2:A:68:ARG:O	2:A:69:ASP:HB2	2.21	0.41
6:F:86:ARG:NH2	6:F:99:GLY:O	2.49	0.41
7:M:78:VAL:O	7:M:78:VAL:HG12	2.20	0.41
2:C:475:VAL:HG12	2:C:503:ILE:HD11	2.02	0.41
6:F:83:THR:O	6:F:83:THR:HG23	2.20	0.41
6:D:79:VAL:HB	6:D:229:LYS:HG2	2.01	0.41
6:D:455:LYS:O	6:D:459:LEU:N	2.45	0.41
7:N:44:ASP:OD2	7:N:44:ASP:N	2.48	0.41
4:H:23:ILE:HG12	4:H:34:ILE:HB	2.03	0.41
7:Q:78:VAL:HG12	7:Q:78:VAL:O	2.21	0.41
1:W:117:SER:C	1:W:119:GLN:H	2.28	0.41
2:C:164:ARG:HH11	2:C:164:ARG:HG3	1.85	0.41
4:H:130:VAL:HG13	4:H:131:ILE:H	1.86	0.41
4:H:136:LYS:NZ	5:G:259:GLU:OE2	2.49	0.41
6:E:153:VAL:HG23	6:E:321:LEU:HD23	2.03	0.41
6:D:265:VAL:HG12	6:D:265:VAL:O	2.20	0.41
8:a:20:GLN:HG2	8:a:42:ASN:OD1	2.21	0.41
8:a:254:PHE:CE2	8:a:258:VAL:HG21	2.56	0.41
2:B:197:ALA:HB1	2:B:200:GLN:CG	2.50	0.41
3:X:144:ALA:C	3:X:146:SER:H	2.29	0.41
7:Q:15:VAL:HG23	7:Q:16:MET:N	2.36	0.41
4:H:72:GLN:HB3	4:H:73:PRO:HD2	2.03	0.40
6:F:219:VAL:HA	6:F:222:THR:CG2	2.51	0.40
6:E:282:ILE:HG21	6:E:292:SER:HB2	2.03	0.40
7:I:48:LEU:HD21	7:Q:46:ILE:HD12	2.03	0.40
7:N:16:MET:HG3	7:O:14:ALA:HB1	2.02	0.40
2:C:52:MET:O	2:C:91:THR:OG1	2.34	0.40
5:G:75:VAL:HG12	5:G:76:GLY:N	2.36	0.40
6:F:400:LEU:HD12	6:F:427:PHE:CE1	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:148:THR:HG23	2:B:150:TYR:H	1.85	0.40
3:X:135:ILE:HD13	3:X:135:ILE:O	2.21	0.40
4:H:41:LEU:HB3	4:H:71:VAL:CG1	2.52	0.40
5:G:98:LEU:C	5:G:98:LEU:HD23	2.46	0.40
6:F:420:LEU:HA	6:F:423:THR:HG22	2.04	0.40
7:Q:61:ASP:C	7:Q:64:PRO:HD2	2.47	0.40
1:W:8:ARG:HB3	1:W:9:PRO:HD3	2.02	0.40
2:C:39:ILE:HD12	2:C:75:VAL:CG2	2.51	0.40
3:X:42:ASP:OD2	3:X:43:GLY:N	2.54	0.40
4:H:137:ALA:O	4:H:138:MET:HG2	2.21	0.40
6:E:169:GLU:HG3	6:E:417:TYR:CD2	2.57	0.40
7:J:45:LEU:O	7:J:48:LEU:HB3	2.21	0.40
7:O:7:ASP:HA	7:O:10:TYR:HD1	1.87	0.40
2:B:3:LEU:HD22	3:X:115:ARG:HD3	2.03	0.40
2:B:480:ALA:N	2:B:481:PRO:CD	2.85	0.40
4:H:42:LEU:HD22	5:G:214:LEU:HD21	2.04	0.40
6:F:242:ASP:O	6:F:295:ALA:HB3	2.21	0.40
7:L:24:ALA:HB3	7:Q:63:ILE:HG21	2.03	0.40
7:M:75:MET:HE2	7:M:76:PHE:CE2	2.57	0.40
7:P:31:LEU:HD11	7:P:56:VAL:HG21	2.02	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	W	171/177 (97%)	166 (97%)	5 (3%)	0	100	100
2	A	508/513 (99%)	497 (98%)	11 (2%)	0	100	100
2	B	511/513 (100%)	492 (96%)	19 (4%)	0	100	100
2	C	508/513 (99%)	487 (96%)	21 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	X	151/156 (97%)	145 (96%)	6 (4%)	0	100	100
3	Y	153/156 (98%)	151 (99%)	2 (1%)	0	100	100
4	H	134/139 (96%)	128 (96%)	6 (4%)	0	100	100
5	G	282/287 (98%)	274 (97%)	8 (3%)	0	100	100
6	D	458/471 (97%)	442 (96%)	16 (4%)	0	100	100
6	E	458/471 (97%)	441 (96%)	17 (4%)	0	100	100
6	F	456/471 (97%)	438 (96%)	18 (4%)	0	100	100
7	I	75/79 (95%)	73 (97%)	2 (3%)	0	100	100
7	J	75/79 (95%)	71 (95%)	2 (3%)	2 (3%)	4	21
7	L	75/79 (95%)	73 (97%)	2 (3%)	0	100	100
7	M	75/79 (95%)	71 (95%)	4 (5%)	0	100	100
7	N	75/79 (95%)	73 (97%)	2 (3%)	0	100	100
7	O	75/79 (95%)	73 (97%)	2 (3%)	0	100	100
7	P	75/79 (95%)	71 (95%)	4 (5%)	0	100	100
7	Q	75/79 (95%)	72 (96%)	3 (4%)	0	100	100
7	R	74/79 (94%)	73 (99%)	1 (1%)	0	100	100
7	S	75/79 (95%)	74 (99%)	1 (1%)	0	100	100
8	a	264/271 (97%)	256 (97%)	7 (3%)	1 (0%)	30	63
All	All	4803/4928 (98%)	4641 (97%)	159 (3%)	3 (0%)	50	79

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
8	a	34	PRO
7	J	44	ASP
7	J	47	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	W	135/140 (96%)	132 (98%)	3 (2%)	47	71
2	A	400/407 (98%)	393 (98%)	7 (2%)	54	76
2	B	403/407 (99%)	397 (98%)	6 (2%)	60	80
2	C	400/407 (98%)	390 (98%)	10 (2%)	42	69
3	X	117/121 (97%)	114 (97%)	3 (3%)	41	68
3	Y	112/121 (93%)	109 (97%)	3 (3%)	40	67
4	H	108/110 (98%)	105 (97%)	3 (3%)	38	66
5	G	236/238 (99%)	227 (96%)	9 (4%)	28	59
6	D	380/389 (98%)	373 (98%)	7 (2%)	54	76
6	E	380/389 (98%)	375 (99%)	5 (1%)	65	82
6	F	379/389 (97%)	375 (99%)	4 (1%)	70	84
7	I	54/56 (96%)	53 (98%)	1 (2%)	52	75
7	J	54/56 (96%)	52 (96%)	2 (4%)	29	59
7	L	54/56 (96%)	54 (100%)	0	100	100
7	M	54/56 (96%)	54 (100%)	0	100	100
7	N	54/56 (96%)	53 (98%)	1 (2%)	52	75
7	O	53/56 (95%)	53 (100%)	0	100	100
7	P	54/56 (96%)	54 (100%)	0	100	100
7	Q	54/56 (96%)	54 (100%)	0	100	100
7	R	54/56 (96%)	54 (100%)	0	100	100
7	S	54/56 (96%)	54 (100%)	0	100	100
8	a	233/237 (98%)	229 (98%)	4 (2%)	56	78
All	All	3822/3915 (98%)	3754 (98%)	68 (2%)	54	76

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	W	3	PHE
1	W	6	VAL
1	W	37	VAL
2	C	97	VAL
2	C	110	THR
2	C	142	VAL
2	C	196	VAL
2	C	221	THR

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Mol	Chain	Res	Type
2	C	236	TYR
2	C	426	LYS
2	C	487	ASN
2	C	488	GLN
2	C	493	ASN
2	B	142	VAL
2	B	147	GLN
2	B	304	VAL
2	B	309	VAL
2	B	350	ASP
2	B	513	TRP
2	A	27	ASN
2	A	93	ARG
2	A	220	ASN
2	A	236	TYR
2	A	338	SER
2	A	372	ILE
2	A	374	VAL
3	X	120	LEU
3	X	135	ILE
3	X	138	ARG
4	H	42	LEU
4	H	102	GLU
4	H	132	GLU
5	G	5	ASP
5	G	60	LEU
5	G	110	VAL
5	G	167	LEU
5	G	177	THR
5	G	190	LEU
5	G	193	SER
5	G	196	ASP
5	G	226	GLN
6	F	42	VAL
6	F	79	VAL
6	F	95	VAL
6	F	224	LEU
6	E	15	VAL
6	E	54	ILE
6	E	318	THR
6	E	389	VAL
6	E	438	LEU

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Mol	Chain	Res	Type
6	D	19	GLN
6	D	20	ASP
6	D	51	VAL
6	D	103	GLU
6	D	181	GLU
6	D	277	VAL
6	D	451	GLU
7	I	44	ASP
7	J	37	GLU
7	J	47	PRO
7	N	37	GLU
3	Y	19	LEU
3	Y	46	SER
3	Y	83	ARG
8	a	15	HIS
8	a	32	GLN
8	a	38	PHE
8	a	44	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	W	22	GLN
1	W	82	ASN
2	C	147	GLN
2	C	186	GLN
2	C	200	GLN
2	C	215	HIS
2	C	443	GLN
2	C	488	GLN
2	C	493	ASN
2	B	140	GLN
2	B	147	GLN
2	B	186	GLN
2	B	215	HIS
2	B	399	GLN
2	B	493	ASN
2	B	511	GLN
2	A	58	ASN
2	A	144	GLN
2	A	333	GLN
2	A	344	ASN

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Mol	Chain	Res	Type
2	A	420	GLN
2	A	484	GLN
3	X	37	GLN
3	X	73	GLN
4	H	104	HIS
4	H	127	GLN
5	G	54	HIS
5	G	91	ASN
5	G	226	GLN
5	G	252	ASN
5	G	265	ASN
5	G	269	GLN
5	G	274	GLN
6	E	243	ASN
6	E	441	GLN
6	D	19	GLN
6	D	33	ASN
6	D	43	GLN
6	D	158	ASN
6	D	279	GLN
6	D	294	GLN
6	D	324	GLN
6	D	343	GLN
7	O	3	ASN
7	R	5	ASN
7	S	3	ASN
8	a	32	GLN
8	a	95	HIS
8	a	184	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 12 ligands modelled in this entry, 5 are monoatomic - leaving 7 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	F	500	10	24,29,29	0.84	0	29,45,45	1.42	3 (10%)
9	ATP	B	600	10	28,33,33	0.88	0	34,52,52	1.20	2 (5%)
11	ADP	E	500	-	24,29,29	0.76	0	29,45,45	0.93	2 (6%)
12	PO4	D	501	10	4,4,4	1.05	0	6,6,6	0.43	0
9	ATP	C	600	10	28,33,33	0.89	0	34,52,52	1.23	2 (5%)
11	ADP	D	500	10	24,29,29	0.89	0	29,45,45	1.17	2 (6%)
9	ATP	A	600	10	28,33,33	0.87	0	34,52,52	1.23	2 (5%)

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	F	500	10	-	1/12/32/32	0/3/3/3
9	ATP	B	600	10	-	2/18/38/38	0/3/3/3
11	ADP	E	500	-	-	1/12/32/32	0/3/3/3
9	ATP	C	600	10	-	4/18/38/38	0/3/3/3
11	ADP	D	500	10	-	6/12/32/32	0/3/3/3
9	ATP	A	600	10	-	1/18/38/38	0/3/3/3

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	F	500	ADP	C4'-O4'-C1'	-5.98	104.45	109.92
9	C	600	ATP	N3-C2-N1	-3.72	123.62	128.67
9	B	600	ATP	N3-C2-N1	-3.70	123.65	128.67
9	A	600	ATP	N3-C2-N1	-3.66	123.70	128.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	D	500	ADP	N3-C2-N1	-3.61	123.77	128.67
9	A	600	ATP	C4-C5-N7	-2.80	106.38	109.34
9	C	600	ATP	C4-C5-N7	-2.74	106.44	109.34
9	B	600	ATP	C4-C5-N7	-2.71	106.48	109.34
11	D	500	ADP	C4-C5-N7	-2.67	106.52	109.34
11	F	500	ADP	O4'-C4'-C3'	-2.40	100.39	105.15
11	E	500	ADP	C5-C6-N6	2.30	123.81	120.31
11	F	500	ADP	C5-C6-N6	2.28	123.78	120.31
11	E	500	ADP	O3A-PB-O1B	-2.18	99.57	111.04

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	B	600	ATP	PB-O3B-PG-O2G
11	D	500	ADP	PA-O3A-PB-O2B
11	D	500	ADP	PA-O3A-PB-O3B
11	D	500	ADP	C5'-O5'-PA-O2A
11	D	500	ADP	O4'-C4'-C5'-O5'
11	D	500	ADP	C3'-C4'-C5'-O5'
9	C	600	ATP	PA-O3A-PB-O2B
11	D	500	ADP	C5'-O5'-PA-O3A
9	A	600	ATP	PA-O3A-PB-O2B
9	B	600	ATP	PB-O3B-PG-O1G
9	C	600	ATP	PB-O3B-PG-O2G
11	F	500	ADP	C4'-C5'-O5'-PA
11	E	500	ADP	O4'-C4'-C5'-O5'
9	C	600	ATP	PB-O3B-PG-O1G
9	C	600	ATP	PA-O3A-PB-O1B

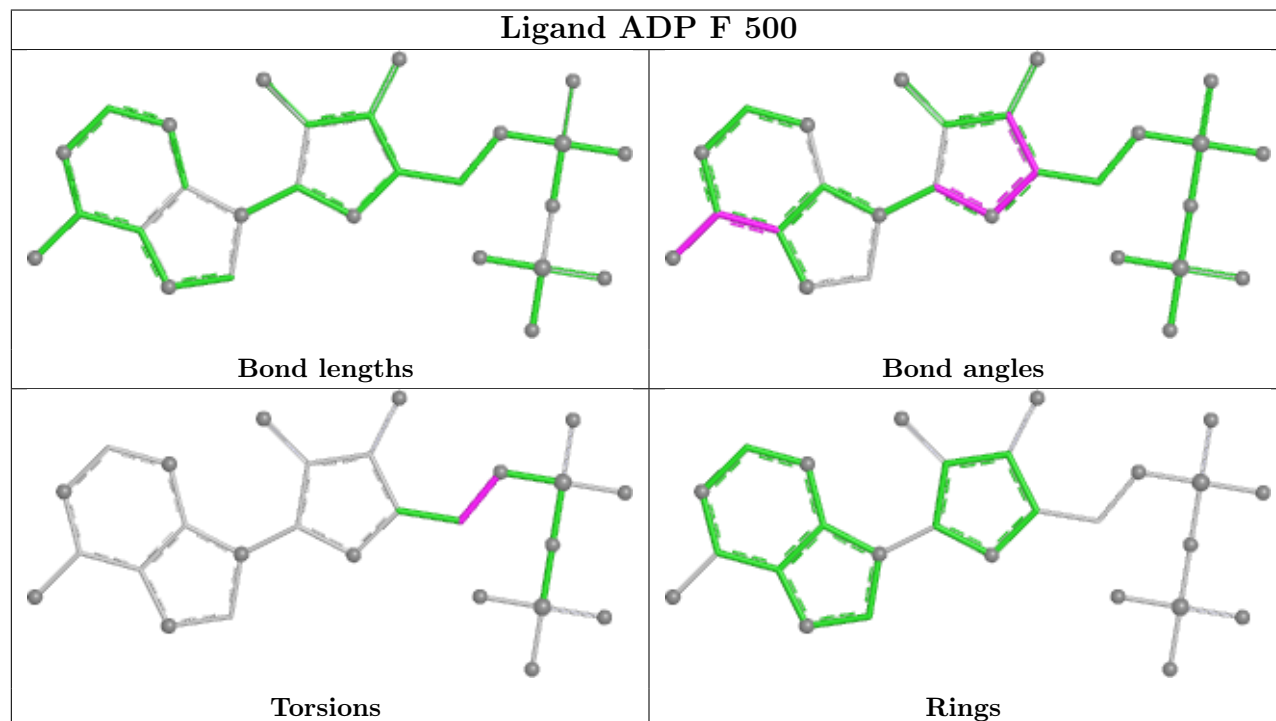
There are no ring outliers.

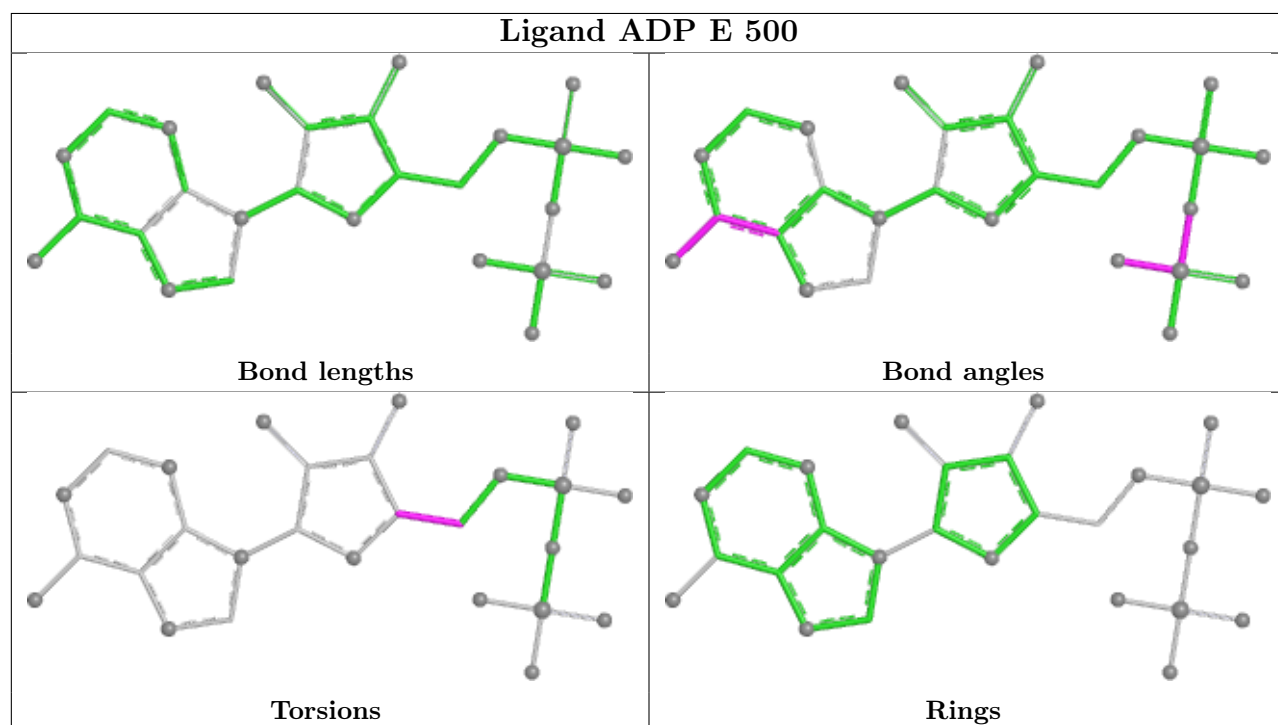
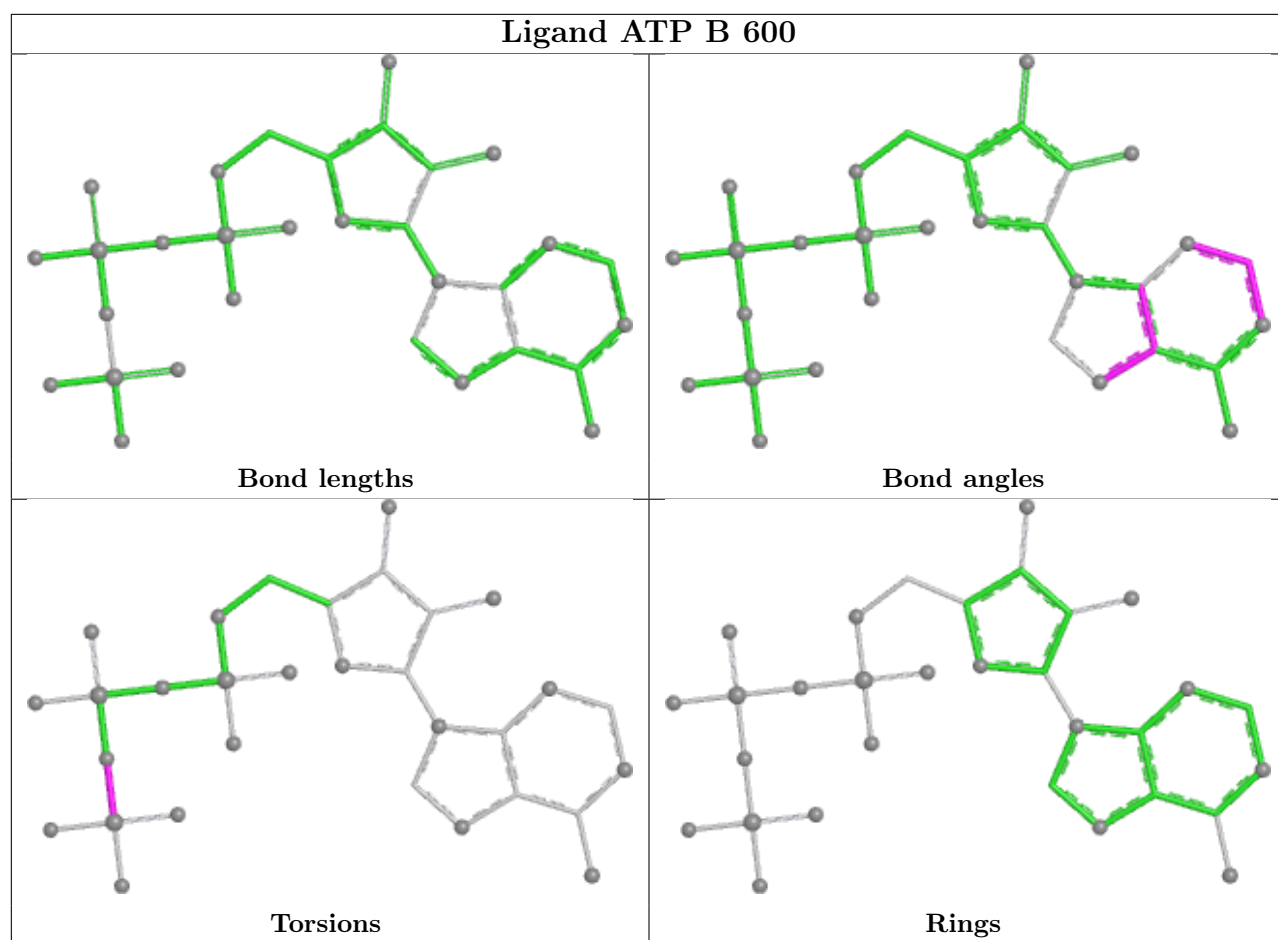
4 monomers are involved in 6 short contacts:

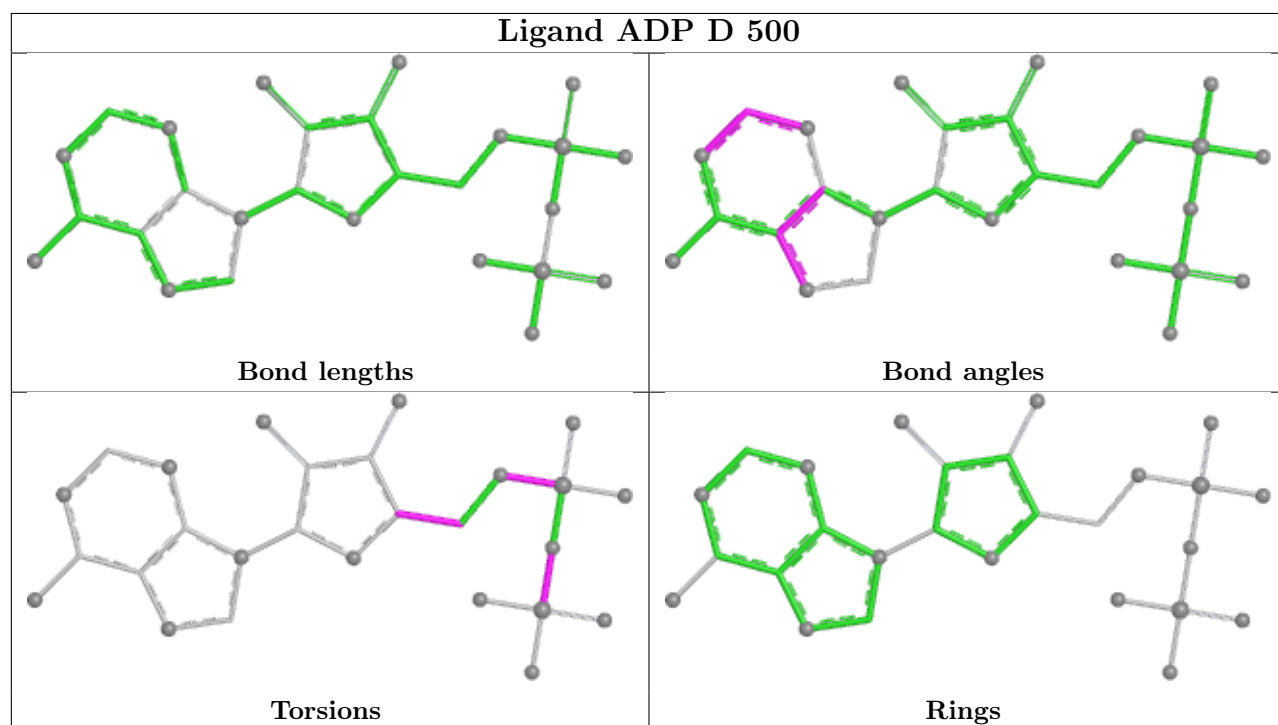
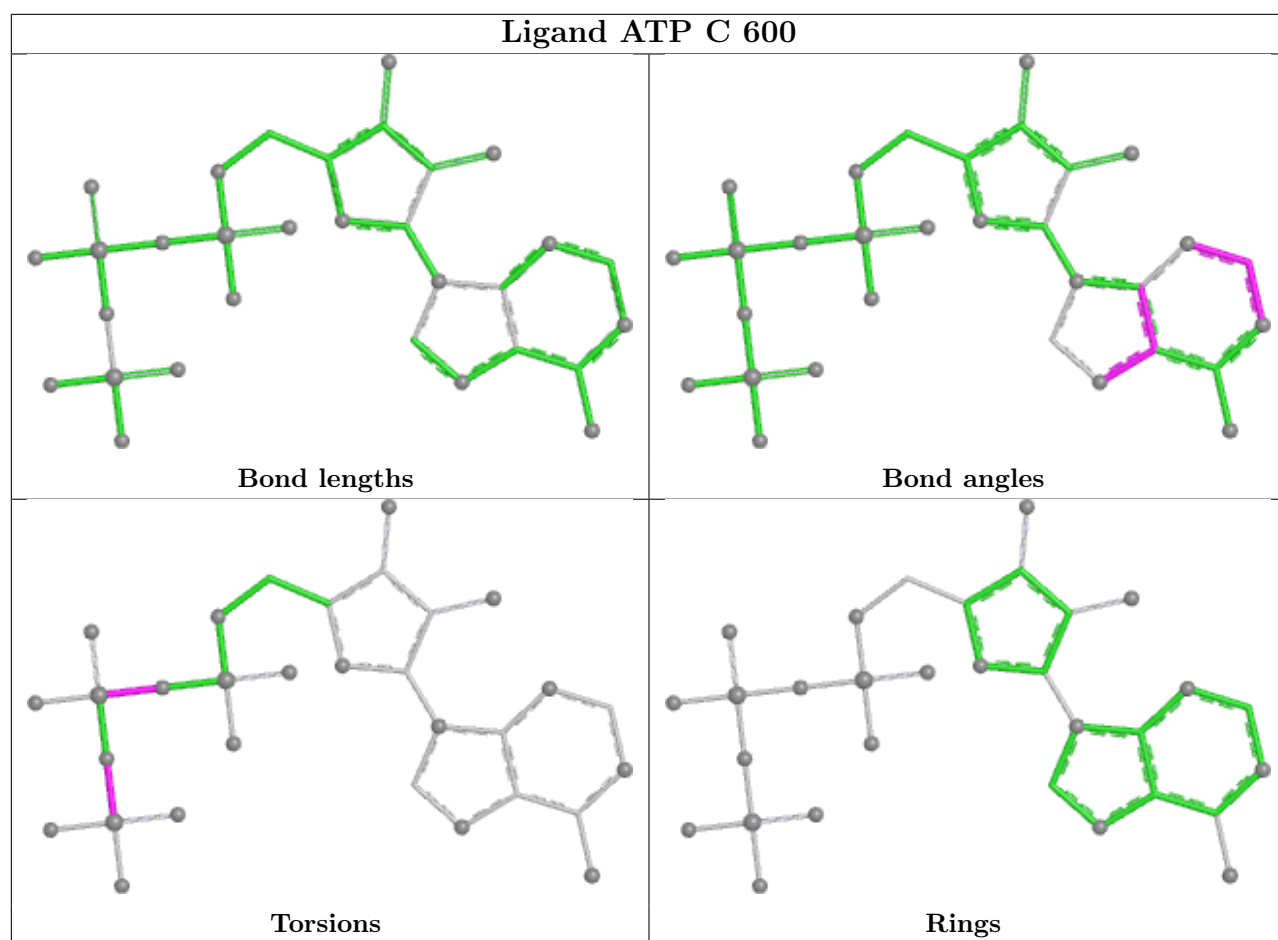
Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	F	500	ADP	3	0
9	B	600	ATP	1	0
11	E	500	ADP	1	0
11	D	500	ADP	1	0

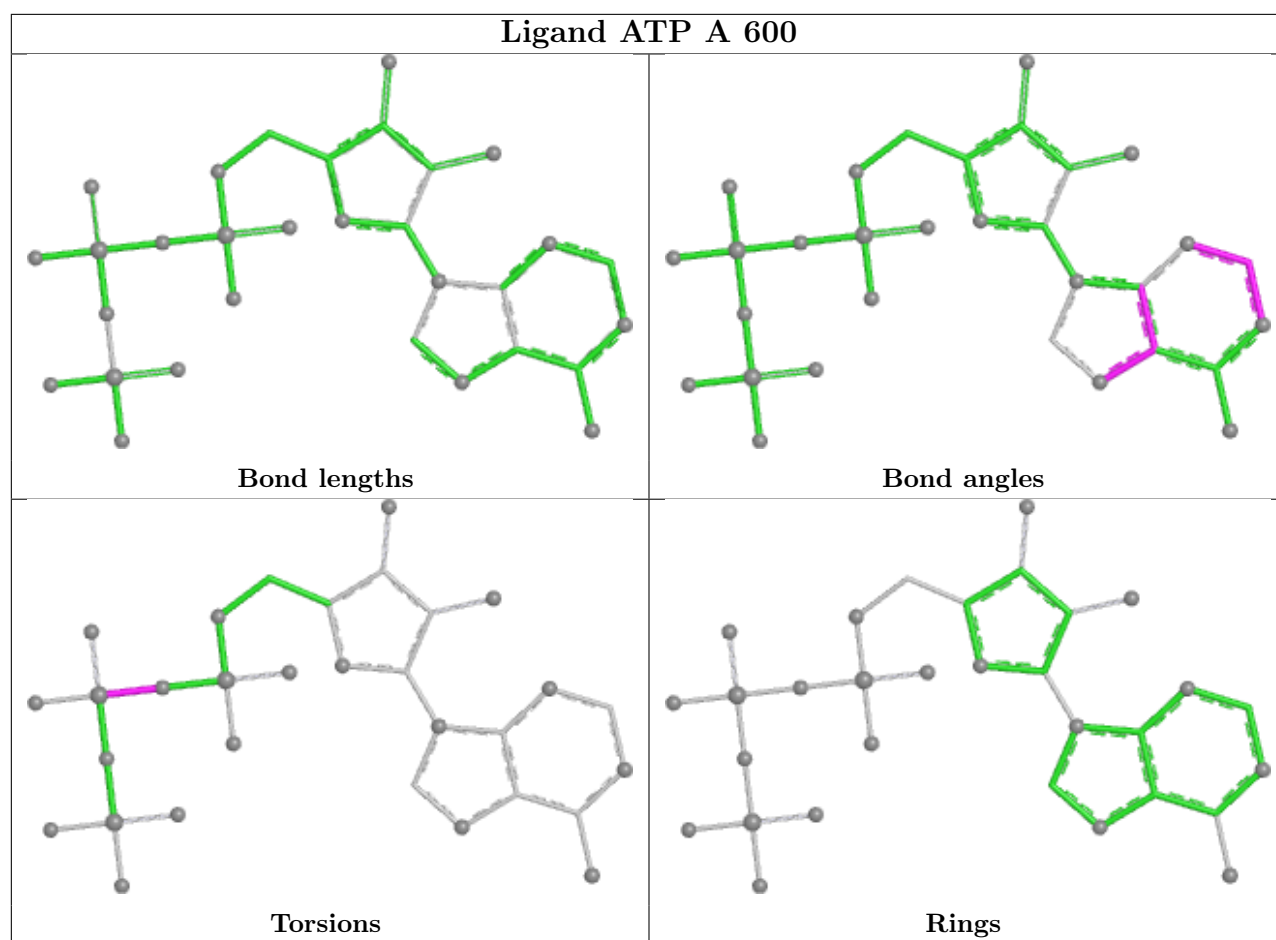
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

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.

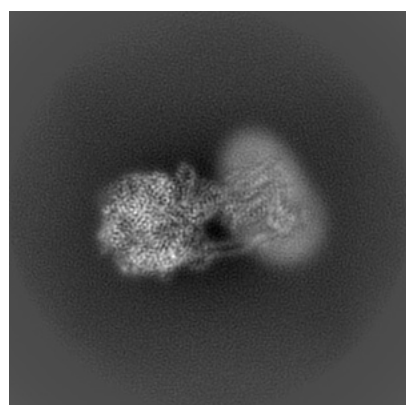
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20172. These allow visual inspection of the internal detail of the map and identification of artifacts.

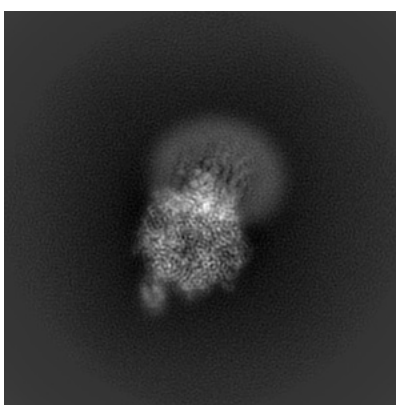
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

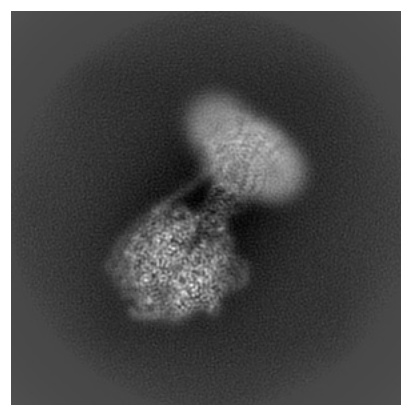
6.1.1 Primary map



X



Y

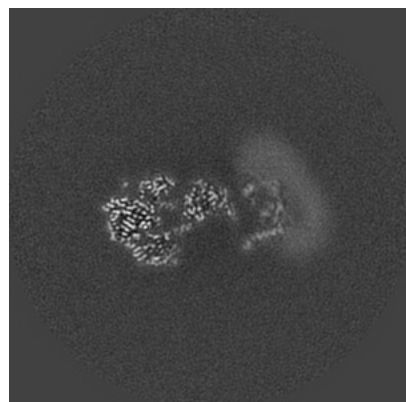


Z

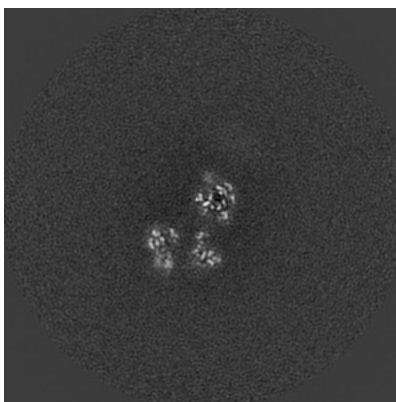
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

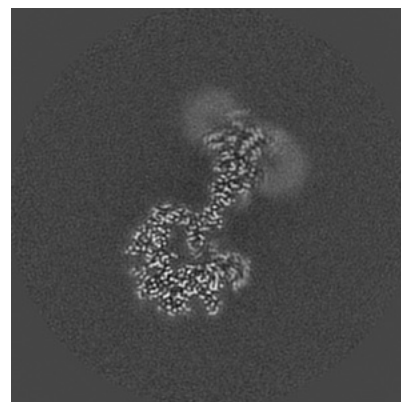
6.2.1 Primary map



X Index: 175



Y Index: 175

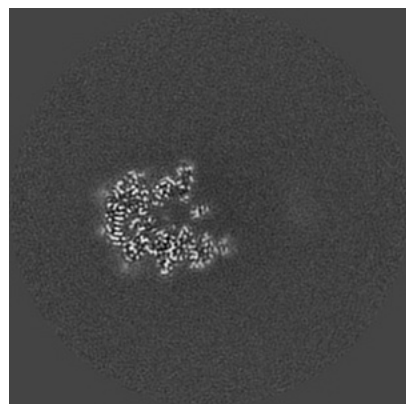


Z Index: 175

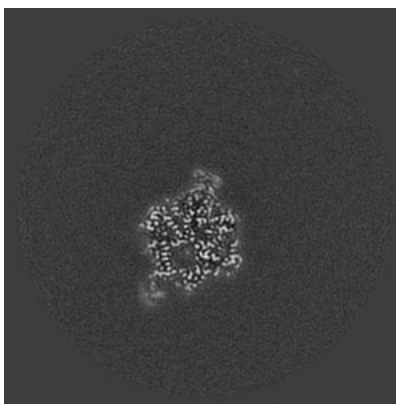
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

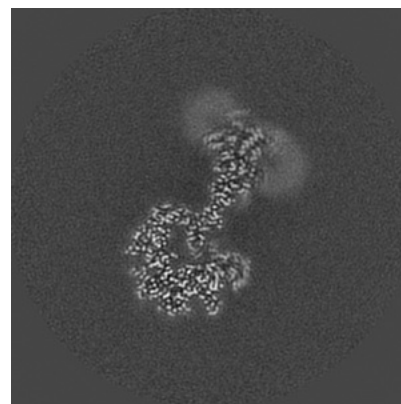
6.3.1 Primary map



X Index: 146



Y Index: 113

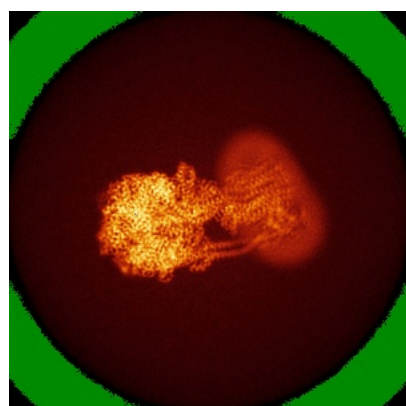


Z Index: 175

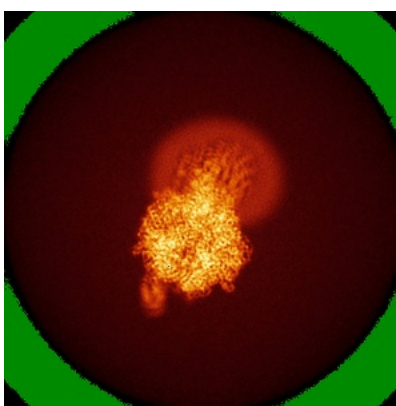
The images above show the largest variance slices of the map in three orthogonal directions.

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

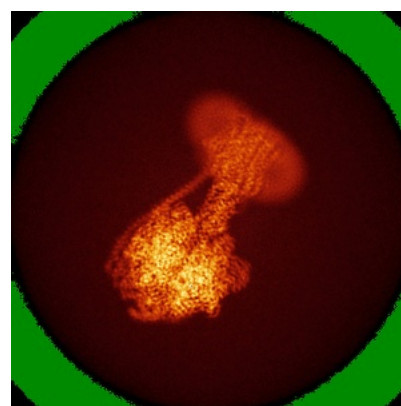
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0171. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

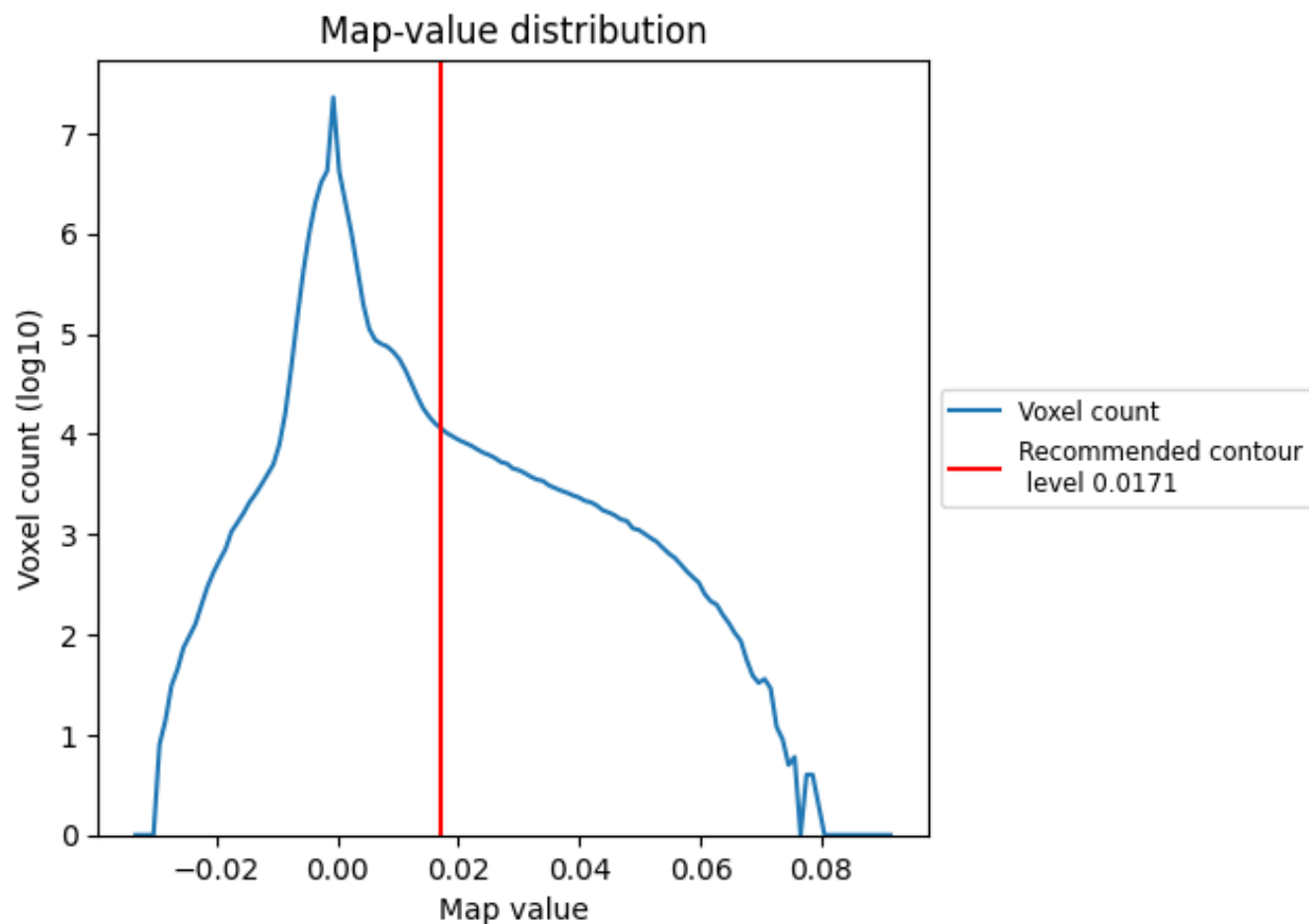
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

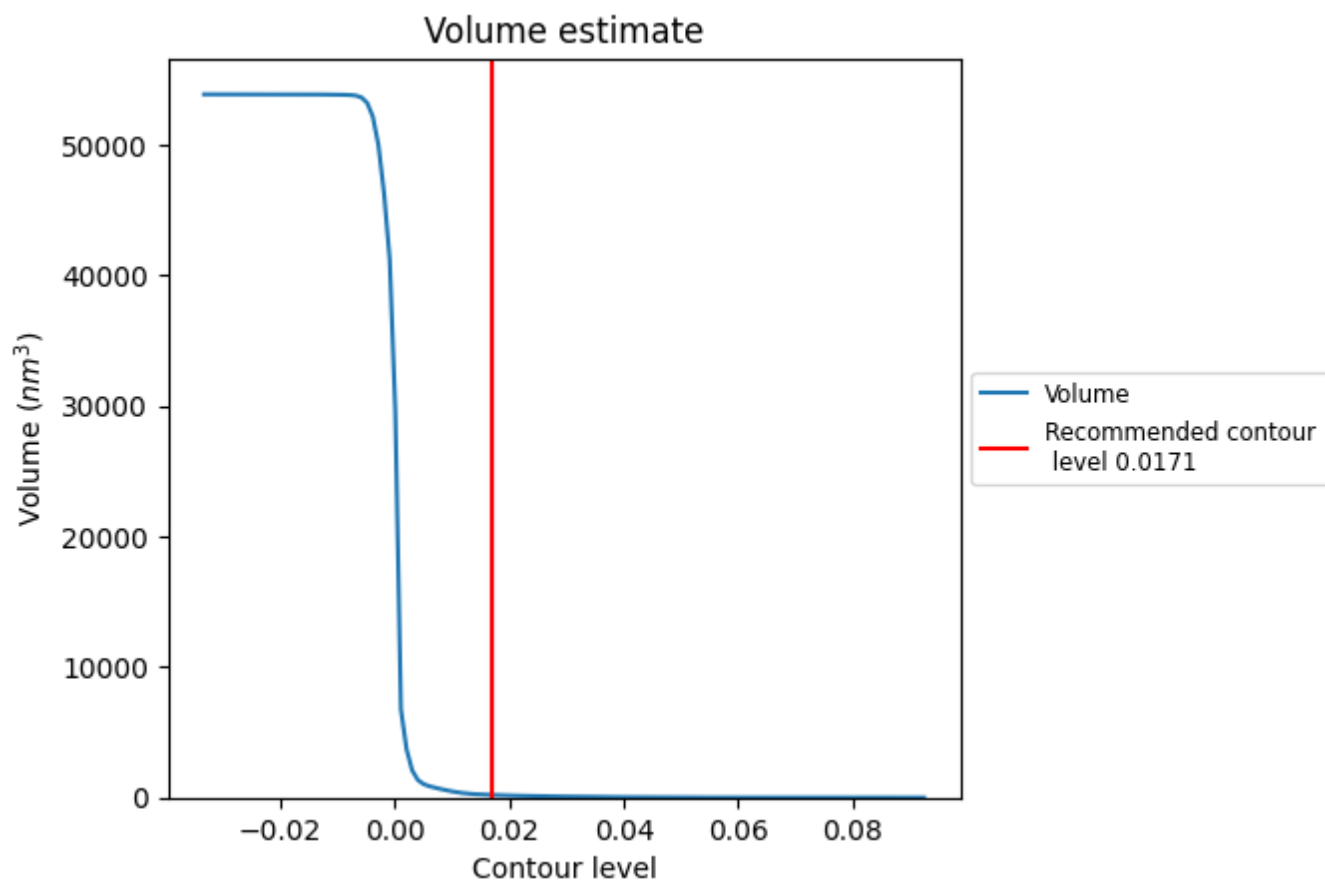
This section contains the results of statistical analysis of the map.

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.

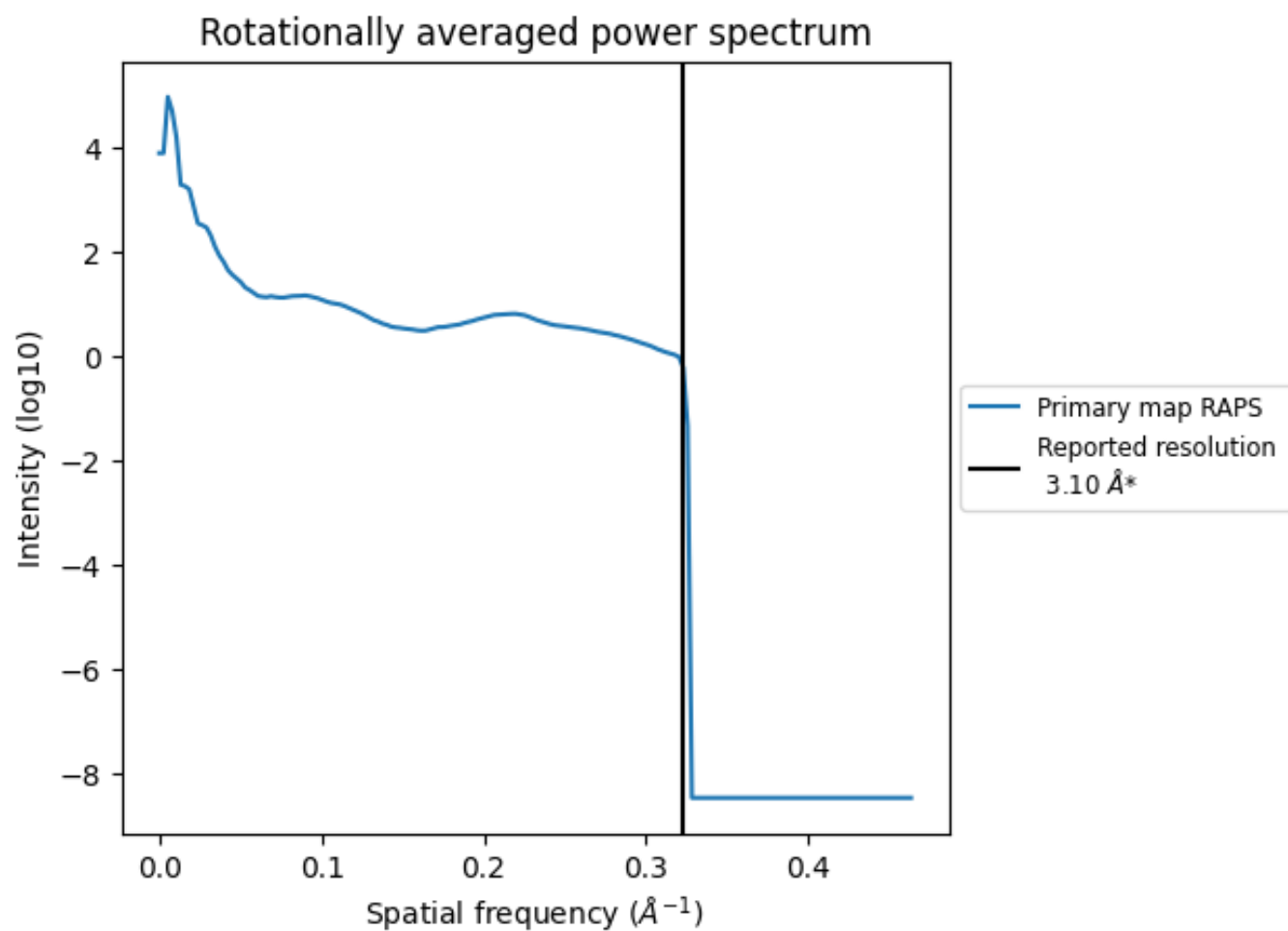
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 196 nm³; this corresponds to an approximate mass of 177 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 0.323 Å⁻¹

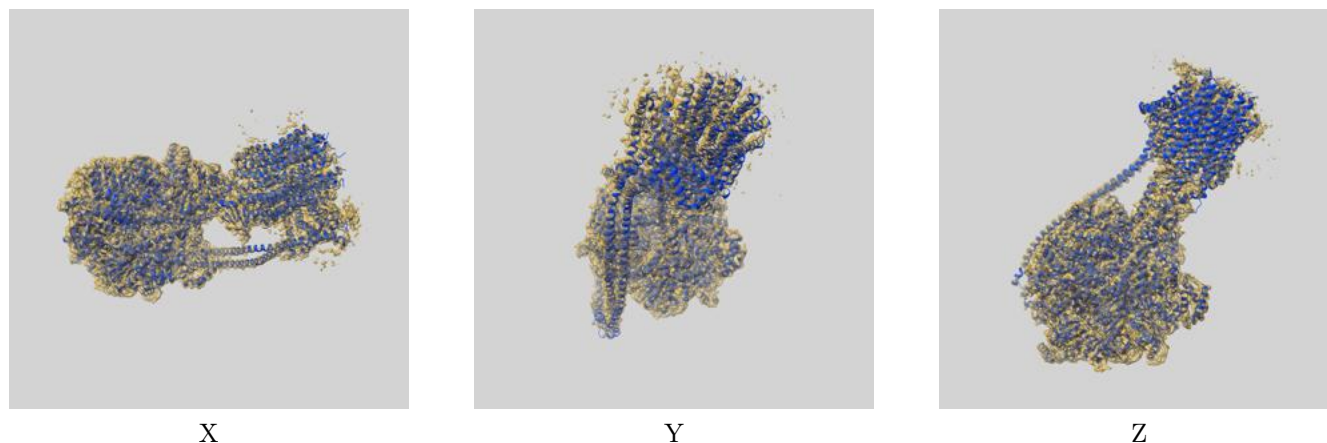
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

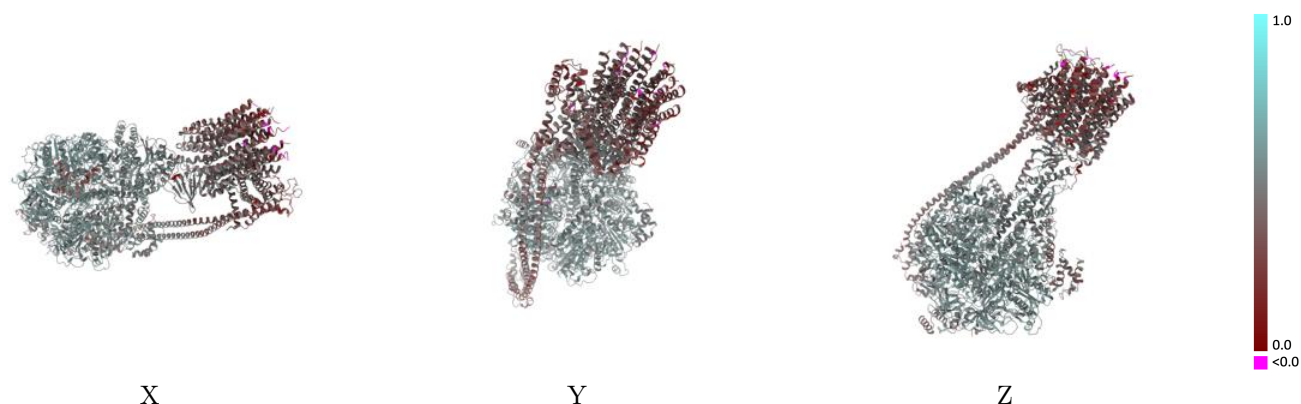
This section contains information regarding the fit between EMDB map EMD-20172 and PDB model 6OQW. Per-residue inclusion information can be found in section 3 on page 10.

9.1 Map-model overlay [i](#)



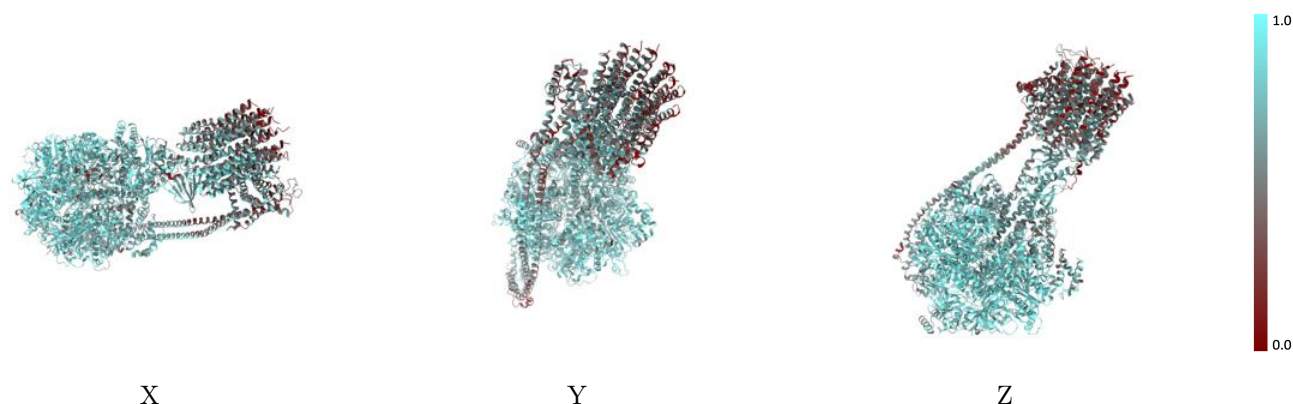
The images above show the 3D surface view of the map at the recommended contour level 0.0171 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



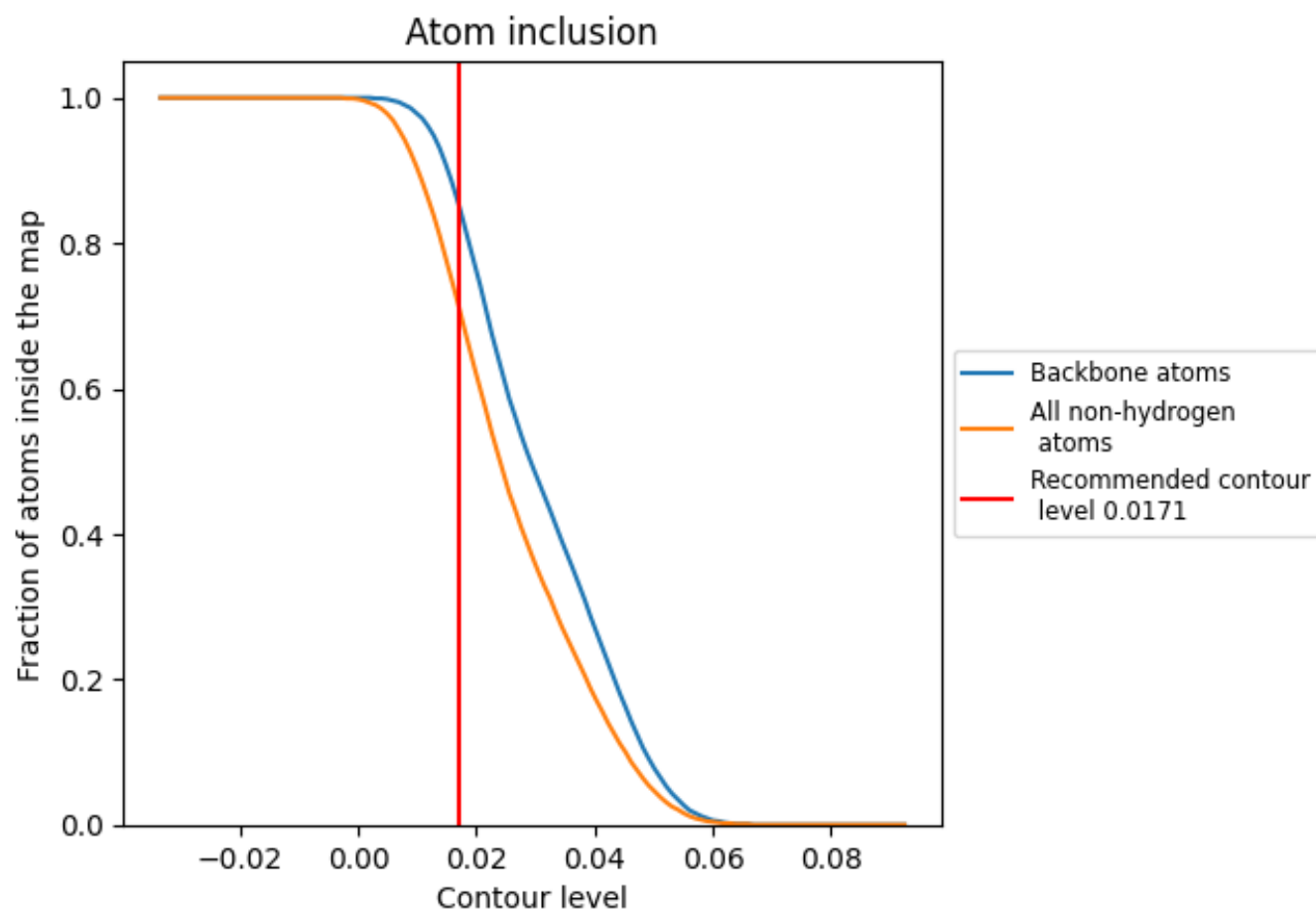
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.0171).















































9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 71% 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 (0.0171) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7120	 0.4870
A	 0.8340	 0.5440
B	 0.8050	 0.5370
C	 0.7960	 0.5260
D	 0.8120	 0.5540
E	 0.7960	 0.5320
F	 0.8500	 0.5580
G	 0.7090	 0.5070
H	 0.6550	 0.4970
I	 0.4100	 0.3230
J	 0.5320	 0.3980
L	 0.4830	 0.3740
M	 0.5570	 0.4020
N	 0.5440	 0.3710
O	 0.4600	 0.3180
P	 0.4060	 0.3020
Q	 0.4320	 0.3420
R	 0.6230	 0.3970
S	 0.5990	 0.3980
W	 0.7010	 0.4520
X	 0.4420	 0.3660
Y	 0.5310	 0.3700
a	 0.4620	 0.3550

