



Full wwPDB EM Validation Report ⓘ

Mar 25, 2026 – 04:57 AM UTC

PDB ID : 6J5I / pdb_00006j5i
EMDB ID : EMD-0677
Title : Cryo-EM structure of the mammalian DP-state ATP synthase
Authors : Gu, J.; Zhang, L.; Yi, J.; Yang, M.
Deposited on : 2019-01-11
Resolution : 3.34 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:

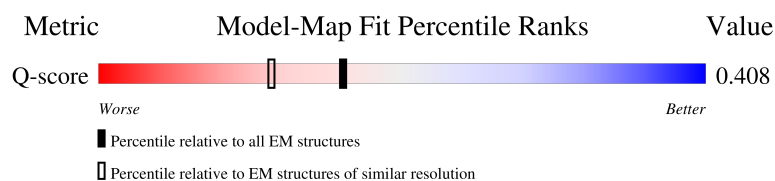
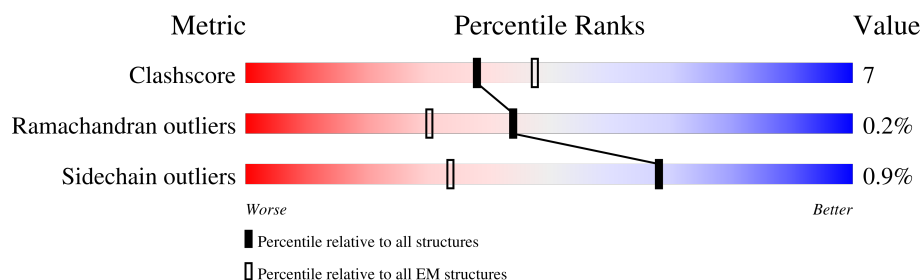
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 3.34 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	14446 (2.84 - 3.84)

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	509	
1	B	509	
1	C	509	
2	D	480	

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Mol	Chain	Length	Quality of chain
2	E	480	83% 13% ..
2	F	480	89% 8%
3	J	83	10% 55% 10% 34%
4	G	272	83% 16%
5	H	132	14% 80% 19%
6	I	48	6% 83% 17%
7	S	187	84% 16%
8	b	209	54% 85% 14%
9	c	70	6% 94% 6%
10	d	147	52% 97%
11	e	63	100% 79% 21%
12	f	87	90% 79% 13% 8%
13	g	84	98% 93% 7%
14	i	42	93% 93% 7%
15	k	29	100% 72% 28%
16	8	67	75% 54% 24% 16%
17	a	226	86% 73% 25%
18	K	72	82% 86% 14%
18	L	72	75% 89% 11%
18	M	72	67% 79% 19%
18	N	72	71% 79% 21%
18	O	72	82% 86% 14%
18	P	72	83% 83% 17%
18	Q	72	74% 86% 14%
18	R	72	71% 86% 14%

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Mol	Chain	Length	Quality of chain
19	u	42	100% 64% 36%

2 Entry composition

There are 22 unique types of molecules in this entry. The entry contains 38144 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, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	509	Total	C	N	O	S	0	0
			3873	2438	682	741	12		
1	B	481	Total	C	N	O	S	0	0
			3669	2312	649	696	12		
1	C	502	Total	C	N	O	S	0	0
			3821	2407	674	728	12		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	471	Total	C	N	O	S	0	0
			3564	2258	605	689	12		
2	E	467	Total	C	N	O	S	0	0
			3532	2240	599	681	12		
2	F	468	Total	C	N	O	S	0	0
			3536	2242	600	682	12		

- Molecule 3 is a protein called ATPase inhibitor, mitochondrial.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	J	55	Total	C	N	O	0	0
			411	246	83	82		

- Molecule 4 is a protein called ATP synthase subunit gamma.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	272	Total	C	N	O	S	0	0
			2112	1332	366	407	7		

- Molecule 5 is a protein called ATP synthase subunit delta, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	H	132	Total	C	N	O	S	0	0
			973	610	165	196	2		

- Molecule 6 is a protein called ATP synthase F1 subunit epsilon.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	I	48	Total	C	N	O	S	0	0
			377	241	69	66	1		

- Molecule 7 is a protein called ATP synthase subunit O, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	S	187	Total	C	N	O	S	0	0
			1435	913	247	266	9		

- Molecule 8 is a protein called ATP synthase peripheral stalk-membrane subunit b.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	b	209	Total	C	N	O	S	0	0
			1476	915	278	278	5		

- Molecule 9 is a protein called ATP synthase-coupling factor 6, mitochondrial.

Mol	Chain	Residues	Atoms				AltConf	Trace
9	c	70	Total	C	N	O	0	0
			428	263	84	81		

- Molecule 10 is a protein called ATP synthase subunit d, mitochondrial.

Mol	Chain	Residues	Atoms				AltConf	Trace
10	d	147	Total	C	N	O	0	0
			741	446	148	147		

- Molecule 11 is a protein called ATP synthase subunit e.

Mol	Chain	Residues	Atoms				AltConf	Trace
11	e	63	Total	C	N	O	0	0
			315	189	63	63		

- Molecule 12 is a protein called ATP synthase subunit f, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	f	80	Total	C	N	O	S	0	0
			495	306	96	91	2		

- Molecule 13 is a protein called ATP synthase subunit g.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	g	84	Total	C	N	O		0	0
			420	252	84	84			

- Molecule 14 is a protein called ATP synthase membrane subunit DAPIT.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	i	42	Total	C	N	O	S	0	0
			303	199	49	53	2		

- Molecule 15 is a protein called subunit k analog.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	k	29	Total	C	N	O		0	0
			145	87	29	29			

- Molecule 16 is a protein called ATP synthase protein 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	8	56	Total	C	N	O	S	0	0
			374	242	60	70	2		

- Molecule 17 is a protein called ATP synthase subunit a.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	a	223	Total	C	N	O	S	0	0
			1687	1119	266	291	11		

- Molecule 18 is a protein called Mitochondrial H⁺ transporting ATP synthase subunit c isoform 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	K	72	Total	C	N	O	S	0	0
			513	340	80	89	4		
18	L	72	Total	C	N	O	S	0	0
			513	340	80	89	4		

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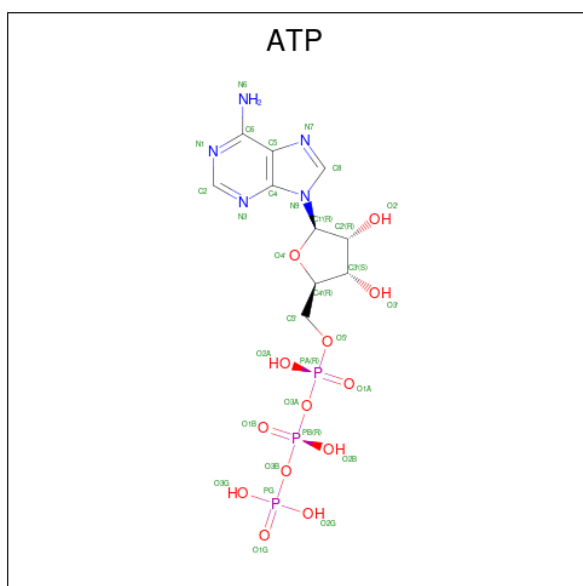
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Mol	Chain	Residues	Atoms					AltConf	Trace
18	M	72	Total	C	N	O	S	0	0
			513	340	80	89	4		
18	N	72	Total	C	N	O	S	0	0
			510	338	80	89	3		
18	O	72	Total	C	N	O	S	0	0
			510	338	80	89	3		
18	P	72	Total	C	N	O	S	0	0
			513	340	80	89	4		
18	Q	72	Total	C	N	O	S	0	0
			513	340	80	89	4		
18	R	72	Total	C	N	O	S	0	0
			510	337	80	89	4		

- Molecule 19 is a protein called ATP synthase membrane subunit 6.8PL.

Mol	Chain	Residues	Atoms				AltConf	Trace
19	u	42	Total	C	N	O	0	0
			210	126	42	42		

- Molecule 20 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
20	A	1	Total	C	N	O	P	0
			31	10	5	13	3	
20	B	1	Total	C	N	O	P	0
			31	10	5	13	3	

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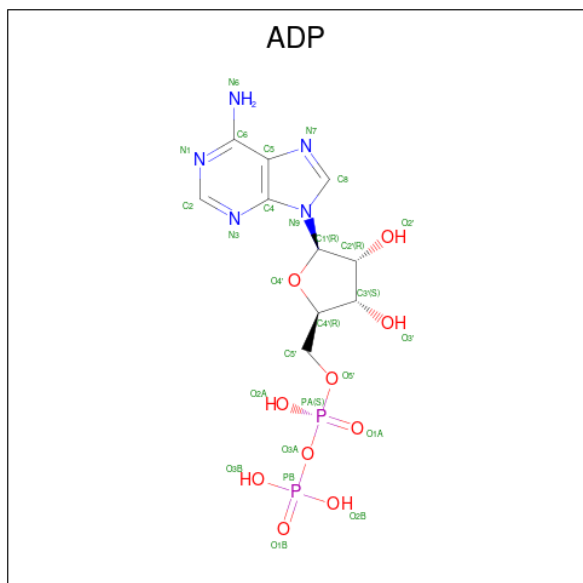
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Mol	Chain	Residues	Atoms					AltConf
20	C	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 21 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 22 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂) (labeled as "Ligand of Interest" by depositor).

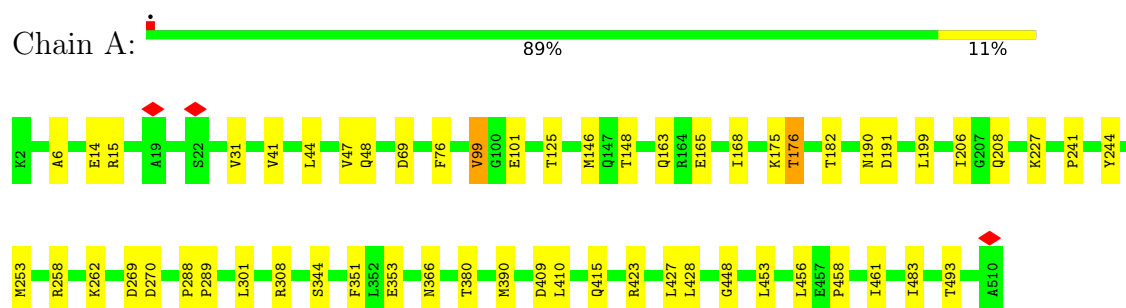


Mol	Chain	Residues	Atoms					AltConf
22	D	1	Total	C	N	O	P	0
			27	10	5	10	2	
22	F	1	Total	C	N	O	P	0
			27	10	5	10	2	

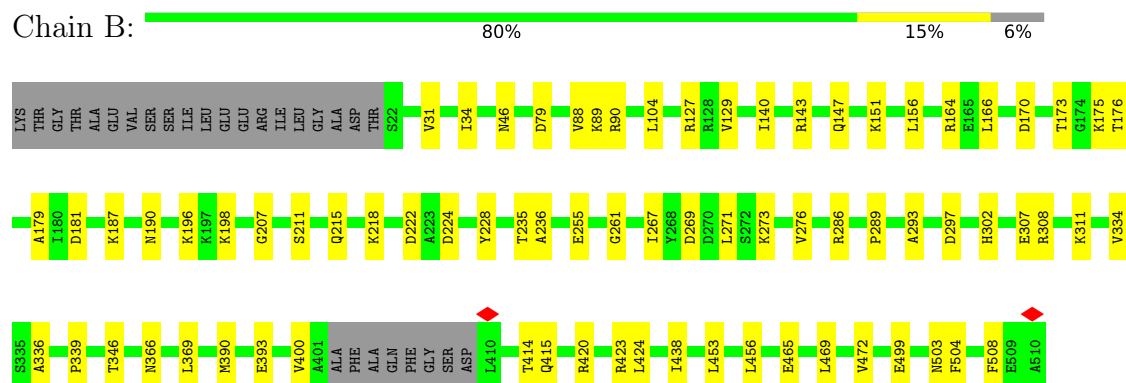
3 Residue-property plots

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

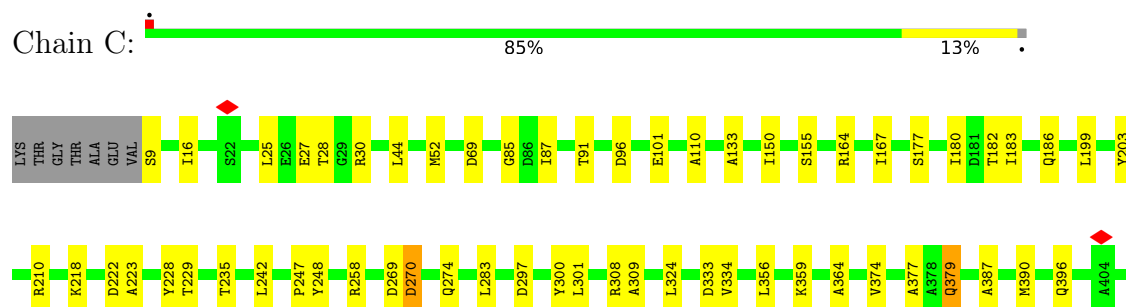
- Molecule 1: ATP synthase subunit alpha, mitochondrial



- Molecule 1: ATP synthase subunit alpha, mitochondrial



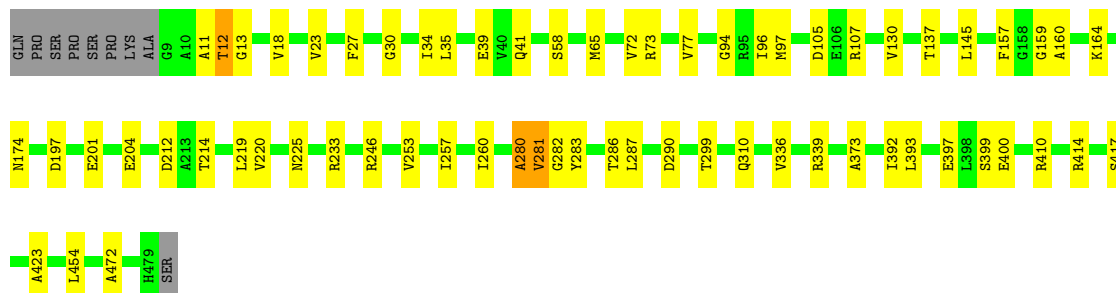
- Molecule 1: ATP synthase subunit alpha, mitochondrial





- Molecule 2: ATP synthase subunit beta

Chain D: 85% 13% ..



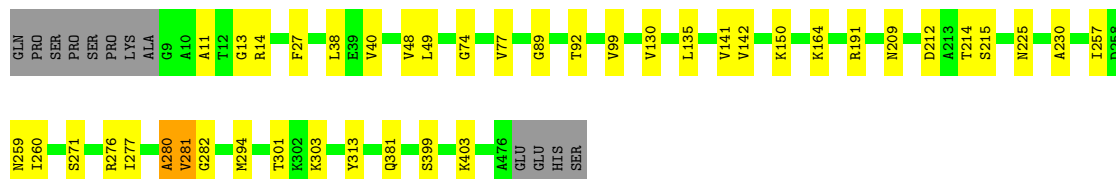
- Molecule 2: ATP synthase subunit beta

Chain E: 83% 13% ..



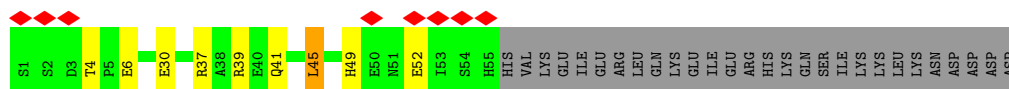
- Molecule 2: ATP synthase subunit beta

Chain F: 89% 8% .



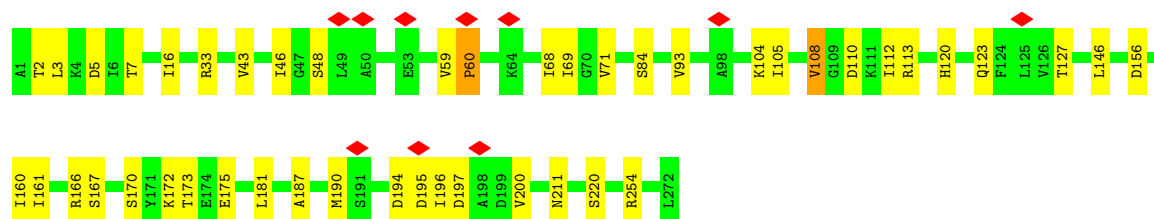
- Molecule 3: ATPase inhibitor, mitochondrial

Chain J: 10% 55% 10% 34%



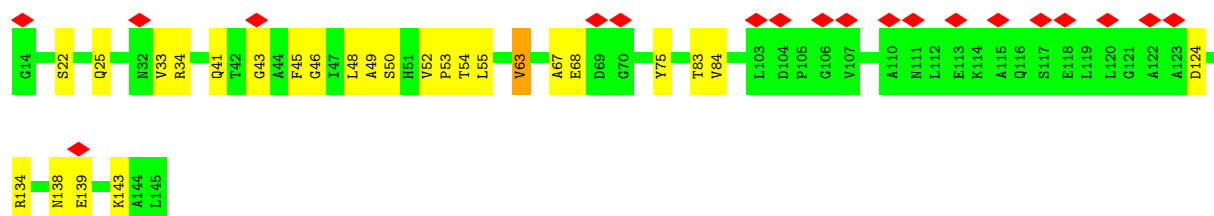
- Molecule 4: ATP synthase subunit gamma

Chain G: 



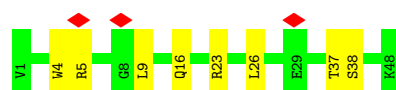
- Molecule 5: ATP synthase subunit delta, mitochondrial

Chain H: 



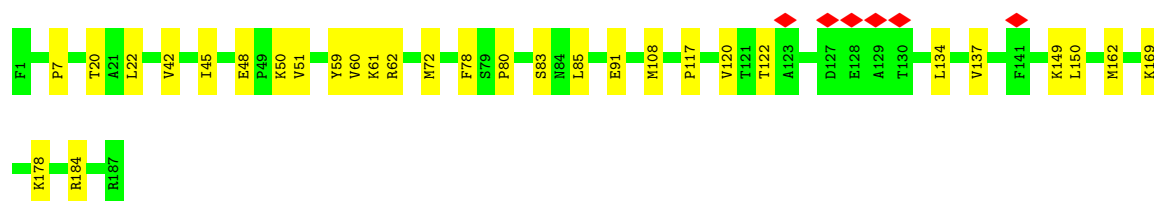
- Molecule 6: ATP synthase F1 subunit epsilon

Chain I: 



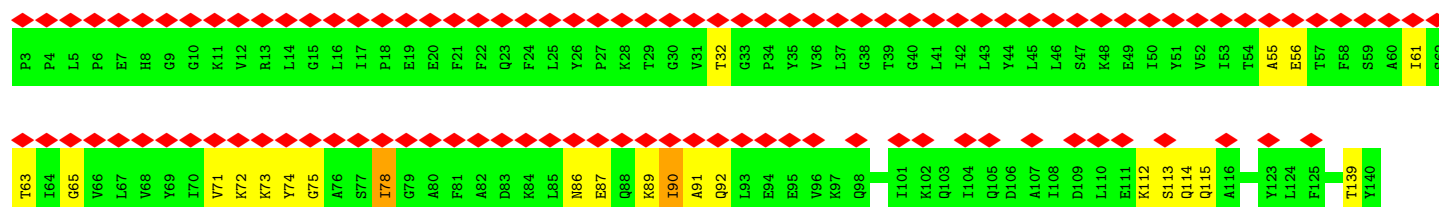
- Molecule 7: ATP synthase subunit O, mitochondrial

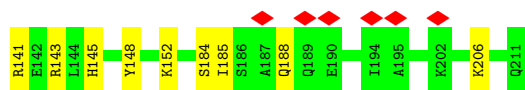
Chain S: 



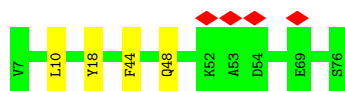
- Molecule 8: ATP synthase peripheral stalk-membrane subunit b

Chain b: 

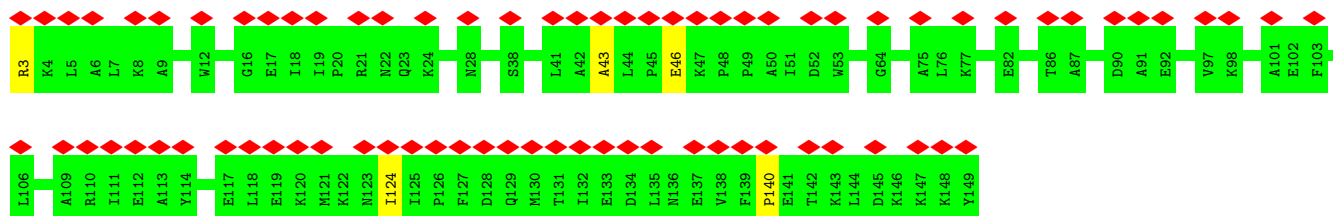




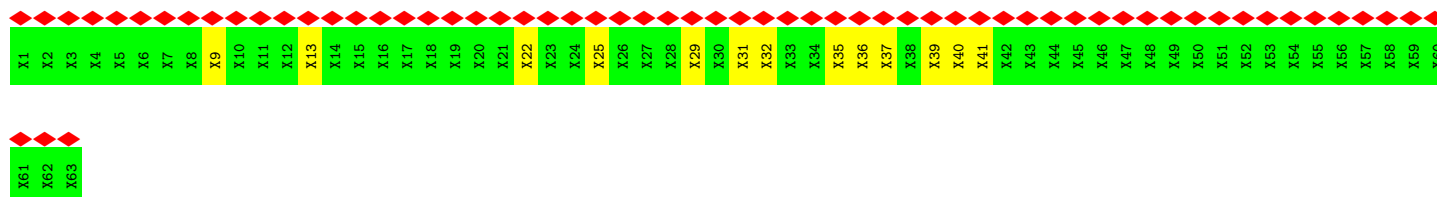
- Molecule 9: ATP synthase-coupling factor 6, mitochondrial



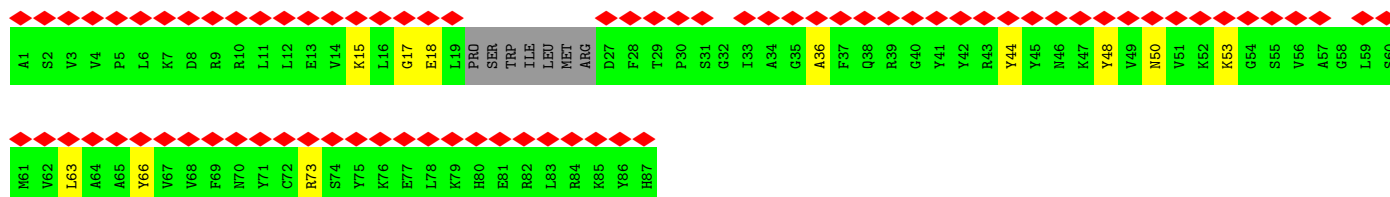
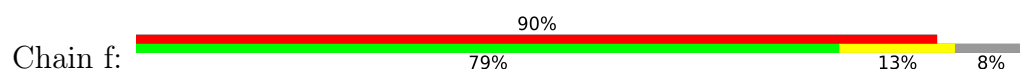
- Molecule 10: ATP synthase subunit d, mitochondrial



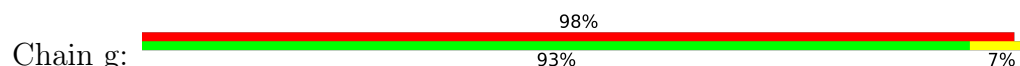
- Molecule 11: ATP synthase subunit e

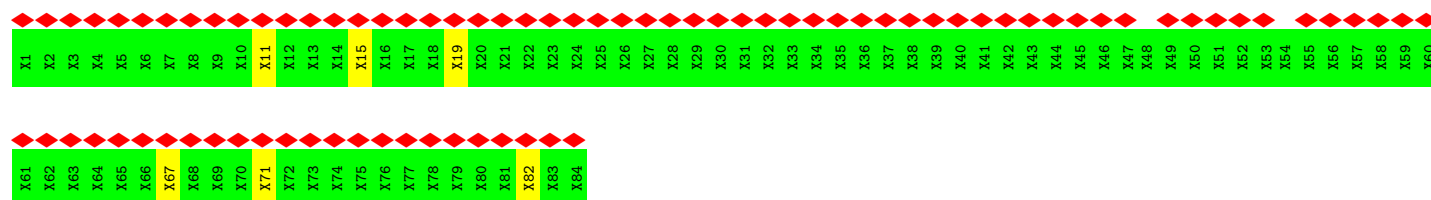


- Molecule 12: ATP synthase subunit f, mitochondrial

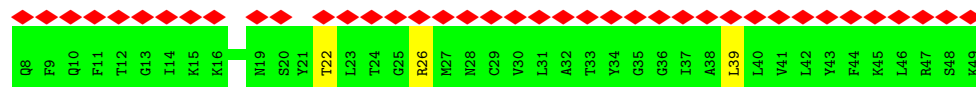


- Molecule 13: ATP synthase subunit g

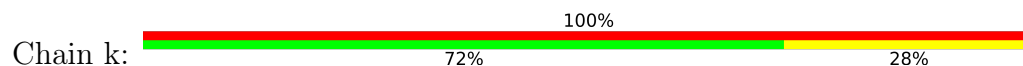




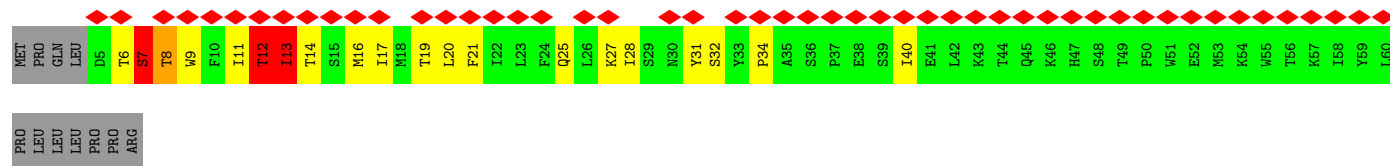
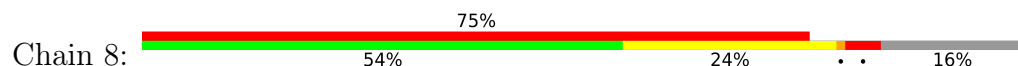
• Molecule 14: ATP synthase membrane subunit DAPIT



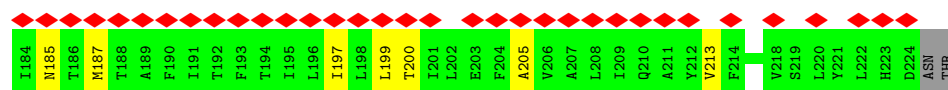
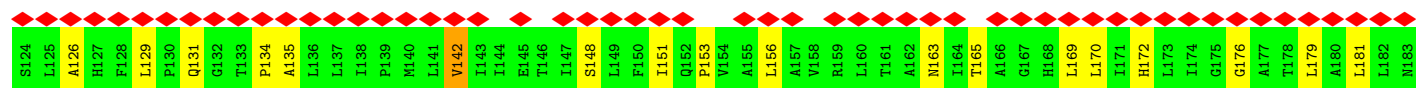
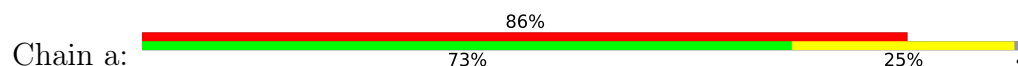
• Molecule 15: subunit k analog



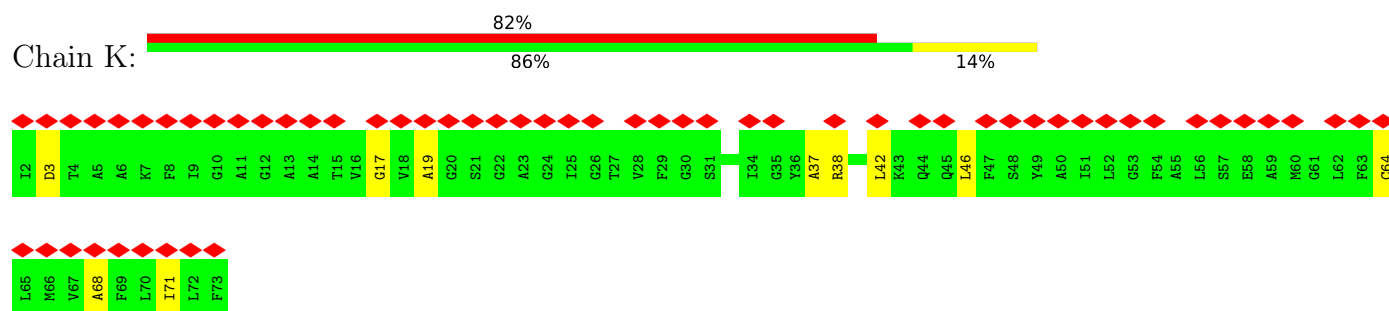
• Molecule 16: ATP synthase protein 8



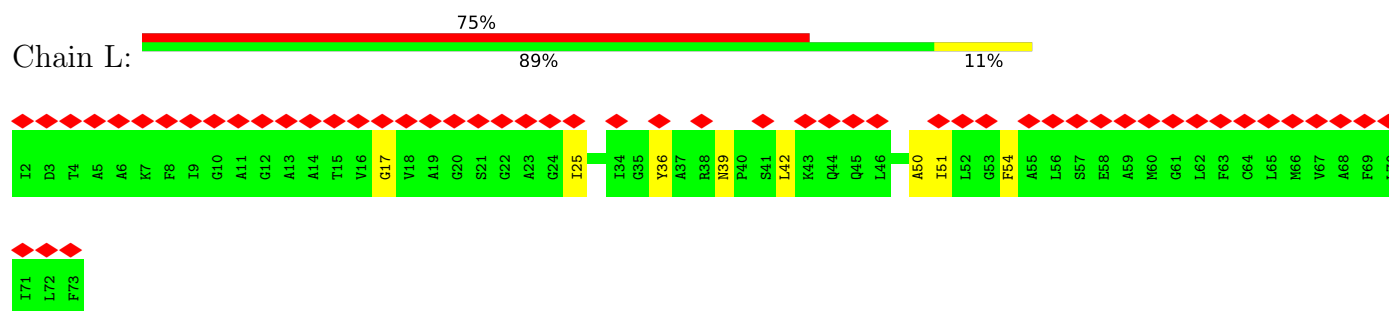
• Molecule 17: ATP synthase subunit a



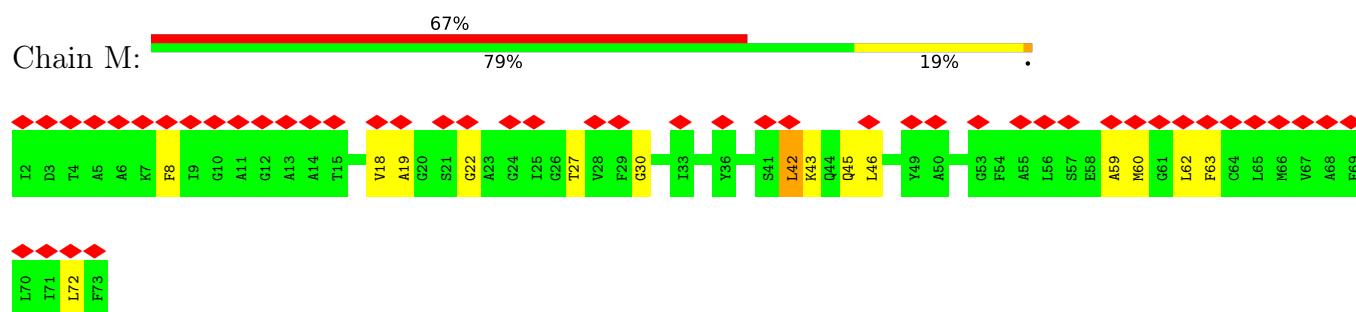
- Molecule 18: Mitochondrial H⁺ transporting ATP synthase subunit c isoform 1



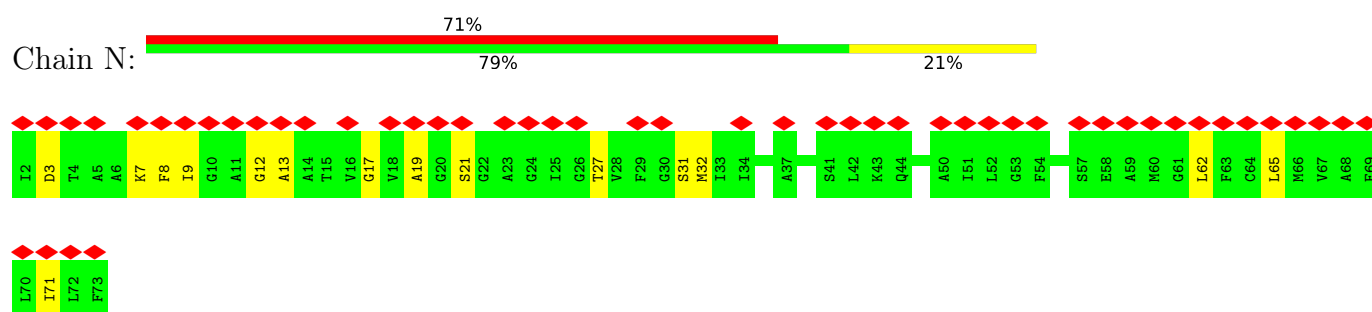
- Molecule 18: Mitochondrial H⁺ transporting ATP synthase subunit c isoform 1



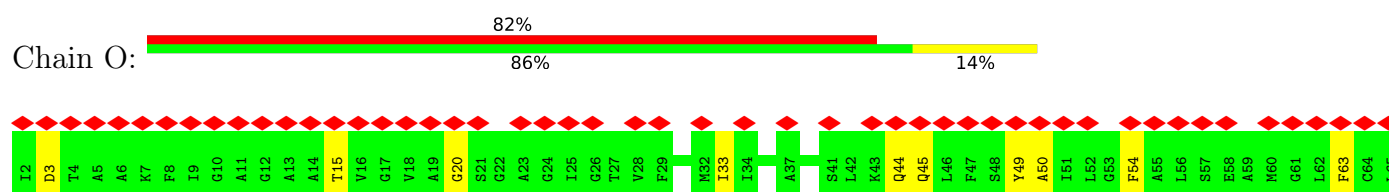
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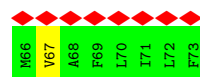
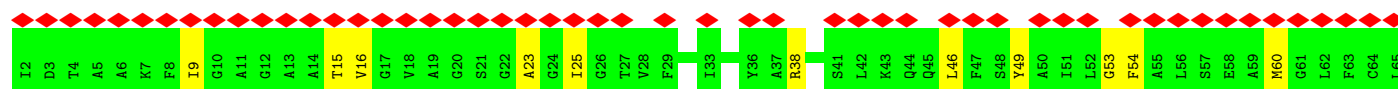
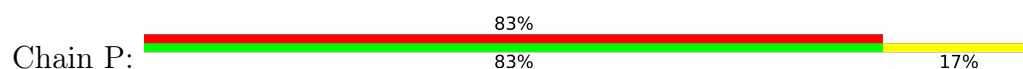


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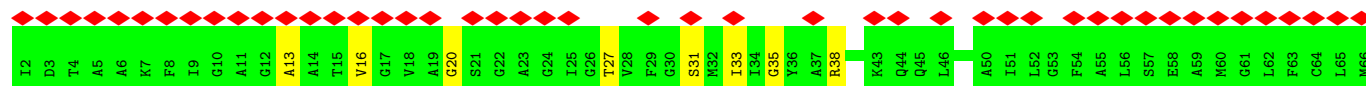
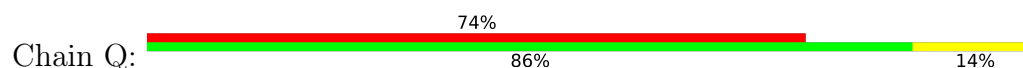




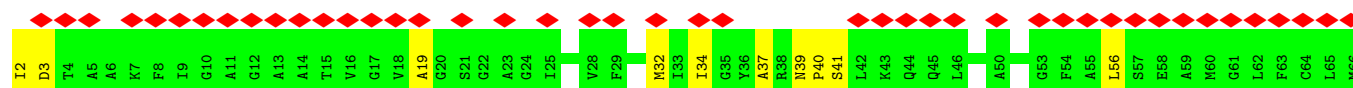
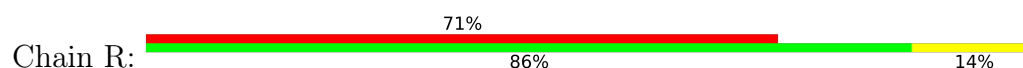
- Molecule 18: Mitochondrial H⁺ transporting ATP synthase subunit c isoform 1



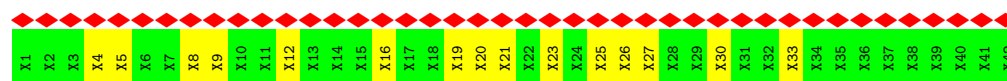
- Molecule 18: Mitochondrial H⁺ transporting ATP synthase subunit c isoform 1



- Molecule 18: Mitochondrial H⁺ transporting ATP synthase subunit c isoform 1



- Molecule 19: ATP synthase membrane subunit 6.8PL



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	291128	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1.56	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.215	Depositor
Minimum map value	-0.126	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.005	Depositor
Recommended contour level	0.0188	Depositor
Map size ($\text{\AA}$)	331.2, 331.2, 331.2	wwPDB
Map dimensions	240, 240, 240	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.38, 1.38, 1.38	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, ATP, 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	A	0.38	0/3924	0.62	0/5295
1	B	0.37	0/3717	0.62	1/5014 (0.0%)
1	C	0.44	0/3872	0.69	2/5225 (0.0%)
2	D	0.41	0/3622	0.64	1/4912 (0.0%)
2	E	0.35	0/3589	0.59	1/4868 (0.0%)
2	F	0.44	0/3593	0.62	1/4873 (0.0%)
3	J	0.37	0/416	0.75	0/555
4	G	0.30	0/2138	0.63	0/2874
5	H	0.27	0/986	0.65	0/1342
6	I	0.24	0/382	0.61	0/511
7	S	0.29	0/1453	0.61	0/1954
8	b	0.28	0/1487	0.79	3/2006 (0.1%)
9	c	0.20	0/431	0.58	0/587
10	d	0.23	0/742	0.75	1/1036 (0.1%)
12	f	0.30	0/499	0.85	0/678
14	i	0.33	0/308	0.73	0/417
16	8	0.45	0/380	1.11	4/522 (0.8%)
17	a	0.44	0/1724	0.82	3/2362 (0.1%)
18	K	0.36	0/522	0.67	0/703
18	L	0.30	0/522	0.75	0/703
18	M	0.38	0/522	0.78	0/703
18	N	0.37	0/519	0.68	0/700
18	O	0.33	0/519	0.74	0/700
18	P	0.28	0/522	0.66	0/703
18	Q	0.29	0/522	0.62	0/703
18	R	0.34	0/519	0.71	0/699
All	All	0.37	0/37430	0.67	17/50645 (0.0%)

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	C	0	1
2	D	0	1
2	E	0	1
2	F	0	1
5	H	0	1
12	f	0	1
16	8	0	4
All	All	0	10

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	b	86	ASN	N-CA-C	-7.51	103.17	111.36
10	d	124	ILE	N-CA-C	7.22	117.98	110.62
1	B	334	VAL	N-CA-C	-6.83	104.67	111.77
16	8	7	SER	CA-C-N	6.10	133.19	121.54
16	8	7	SER	C-N-CA	6.10	133.19	121.54
8	b	90	ILE	N-CA-C	-5.94	104.72	110.72
2	D	281	VAL	N-CA-C	5.85	121.51	109.34
2	E	281	VAL	N-CA-C	5.63	121.05	109.34
17	a	135	ALA	N-CA-C	5.58	118.55	111.69
8	b	32	THR	N-CA-C	5.45	117.22	111.28
2	F	281	VAL	N-CA-C	5.24	120.23	109.34
17	a	16	GLY	CA-C-N	-5.22	116.74	123.16
17	a	16	GLY	C-N-CA	-5.22	116.74	123.16
16	8	12	THR	CA-C-N	5.10	131.15	121.97
16	8	12	THR	C-N-CA	5.10	131.15	121.97
1	C	379	GLN	CA-C-N	5.04	131.16	121.54
1	C	379	GLN	C-N-CA	5.04	131.16	121.54

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
16	8	12	THR	Peptide
16	8	6	THR	Peptide
16	8	7	SER	Peptide
16	8	8	THR	Peptide
1	C	27	GLU	Peptide

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Mol	Chain	Res	Type	Group
2	D	280	ALA	Peptide
2	E	280	ALA	Peptide
2	F	280	ALA	Peptide
5	H	41	GLN	Peptide
12	f	44	TYR	Peptide

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	3873	0	3978	32	0
1	B	3669	0	3781	47	0
1	C	3821	0	3924	43	0
2	D	3564	0	3611	44	0
2	E	3532	0	3590	42	0
2	F	3536	0	3593	32	0
3	J	411	0	384	11	0
4	G	2112	0	2189	33	0
5	H	973	0	971	35	0
6	I	377	0	400	5	0
7	S	1435	0	1546	21	0
8	b	1476	0	1340	30	0
9	c	428	0	305	4	0
10	d	741	0	348	5	0
11	e	315	0	65	12	0
12	f	495	0	371	12	0
13	g	420	0	87	6	0
14	i	303	0	284	21	0
15	k	145	0	31	17	0
16	8	374	0	299	48	0
17	a	1687	0	1757	76	0
18	K	513	0	530	11	0
18	L	513	0	530	6	0
18	M	513	0	530	29	0
18	N	510	0	523	16	0
18	O	510	0	523	8	0
18	P	513	0	530	24	0
18	Q	513	0	530	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	R	510	0	521	27	0
19	u	210	0	45	32	0
20	A	31	0	12	1	0
20	B	31	0	12	2	0
20	C	31	0	12	1	0
21	A	1	0	0	0	0
21	B	1	0	0	0	0
21	C	1	0	0	0	0
21	D	1	0	0	0	0
21	F	1	0	0	0	0
22	D	27	0	11	1	0
22	F	27	0	12	1	0
All	All	38144	0	37175	554	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 7.

All (554) 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:H:48:LEU:HD21	18:R:41:SER:CB	1.23	1.66
16:8:28:ILE:HA	16:8:31:TYR:CE1	1.41	1.54
5:H:48:LEU:CD2	18:R:41:SER:HB3	1.30	1.53
5:H:48:LEU:CD1	18:R:39:ASN:HD21	1.27	1.48
5:H:48:LEU:HD13	18:R:39:ASN:ND2	1.11	1.40
15:k:12:UNK:CB	17:a:129:LEU:HD13	1.58	1.33
14:i:39:LEU:HD21	16:8:11:ILE:CG1	1.60	1.30
8:b:89:LYS:O	8:b:92:GLN:HG3	1.32	1.30
18:P:9:ILE:HD13	19:u:27:UNK:CB	1.63	1.28
18:Q:16:VAL:HG21	19:u:19:UNK:CB	1.67	1.25
10:d:140:PRO:N	16:8:40:ILE:CB	2.07	1.18
3:J:49:HIS:O	3:J:52:GLU:HG3	1.45	1.16
18:P:9:ILE:CD1	19:u:27:UNK:CB	2.23	1.15
5:H:48:LEU:CD2	18:R:41:SER:CB	2.03	1.13
14:i:39:LEU:HD21	16:8:11:ILE:HG12	1.30	1.13
8:b:71:VAL:HG11	12:f:48:TYR:HA	1.30	1.11
17:a:151:ILE:HD12	18:M:62:LEU:HD13	1.17	1.10
17:a:151:ILE:HG21	18:M:62:LEU:HD21	1.35	1.08
16:8:28:ILE:CA	16:8:31:TYR:HE1	1.64	1.08
17:a:151:ILE:HD12	18:M:62:LEU:CD1	1.85	1.06
14:i:39:LEU:HD21	16:8:11:ILE:HG13	1.33	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:a:151:ILE:CD1	18:M:62:LEU:HD13	1.85	1.05
14:i:39:LEU:CD2	16:8:11:ILE:CG1	2.36	1.03
16:8:28:ILE:CA	16:8:31:TYR:CE1	2.37	1.03
17:a:151:ILE:CD1	18:M:62:LEU:CD1	2.37	1.03
15:k:12:UNK:CB	17:a:129:LEU:CD1	2.38	1.01
17:a:151:ILE:HG13	18:M:62:LEU:HD11	1.42	1.01
2:E:469:VAL:O	2:E:473:ASP:OD1	1.79	1.00
17:a:151:ILE:CG2	18:M:62:LEU:HD21	1.92	0.98
18:P:23:ALA:HB1	19:u:9:UNK:CB	1.94	0.98
18:P:9:ILE:HD11	19:u:23:UNK:O	1.65	0.96
15:k:10:UNK:N	17:a:126:ALA:HB1	1.80	0.95
16:8:28:ILE:O	16:8:31:TYR:HD1	1.50	0.94
16:8:28:ILE:HA	16:8:31:TYR:CD1	2.02	0.94
8:b:78:ILE:HG22	12:f:53:LYS:H	1.33	0.94
14:i:39:LEU:HD11	16:8:11:ILE:HG12	1.49	0.93
12:f:53:LYS:O	17:a:41:ARG:O	1.88	0.92
17:a:151:ILE:CG1	18:M:62:LEU:HD11	2.01	0.90
14:i:39:LEU:CD2	16:8:11:ILE:HG12	1.97	0.90
18:Q:16:VAL:CG2	19:u:19:UNK:CB	2.51	0.88
5:H:48:LEU:HD13	18:R:39:ASN:CG	1.99	0.88
4:G:194:ASP:OD1	18:K:38:ARG:HG2	1.74	0.87
12:f:18:GLU:CB	17:a:40:ASN:HB2	2.07	0.85
18:Q:16:VAL:HG12	19:u:16:UNK:HA	1.59	0.84
18:R:2:ILE:HD13	19:u:33:UNK:CB	2.07	0.84
16:8:28:ILE:HG13	16:8:31:TYR:CE1	2.12	0.84
5:H:43:GLY:HA3	18:P:38:ARG:O	1.77	0.83
3:J:49:HIS:O	3:J:52:GLU:CG	2.27	0.83
16:8:16:MET:HG2	17:a:21:THR:HB	1.61	0.82
14:i:39:LEU:CD1	16:8:11:ILE:HG12	2.10	0.81
15:k:10:UNK:N	17:a:126:ALA:CB	2.43	0.80
14:i:39:LEU:HD22	16:8:11:ILE:HD11	1.65	0.79
2:D:472:ALA:O	3:J:45:LEU:HD12	1.81	0.78
2:D:11:ALA:O	2:D:27:PHE:CD1	2.36	0.78
16:8:28:ILE:O	16:8:31:TYR:CD1	2.37	0.78
4:G:2:THR:OG1	4:G:5:ASP:OD2	2.02	0.78
8:b:89:LYS:O	8:b:92:GLN:CG	2.25	0.78
11:e:36:UNK:O	11:e:40:UNK:CB	2.33	0.76
8:b:87:GLU:O	8:b:91:ALA:N	2.19	0.76
18:P:16:VAL:CG1	19:u:16:UNK:O	2.35	0.75
15:k:7:UNK:HA	17:a:129:LEU:HD21	1.68	0.74
5:H:43:GLY:CA	18:P:38:ARG:O	2.35	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:45:PHE:HB3	18:Q:38:ARG:HG3	1.69	0.74
2:D:11:ALA:O	2:D:27:PHE:HD1	1.71	0.73
11:e:22:UNK:CB	13:g:82:UNK:O	2.36	0.73
11:e:35:UNK:O	11:e:39:UNK:N	2.21	0.73
5:H:48:LEU:CD1	18:R:39:ASN:ND2	2.07	0.73
2:D:472:ALA:O	3:J:45:LEU:CD1	2.37	0.73
2:E:10:ALA:HB3	2:E:79:ASP:OD2	1.88	0.73
18:Q:16:VAL:CG1	19:u:16:UNK:HA	2.18	0.73
11:e:22:UNK:HA	13:g:82:UNK:O	1.89	0.72
5:H:48:LEU:HD22	18:R:41:SER:CB	2.17	0.72
17:a:151:ILE:CD1	18:M:62:LEU:HD11	2.17	0.72
18:P:23:ALA:CB	19:u:9:UNK:CB	2.68	0.71
17:a:151:ILE:CG2	18:M:62:LEU:CD2	2.69	0.71
8:b:55:ALA:O	17:a:179:LEU:HD11	1.90	0.71
18:Q:13:ALA:HB2	19:u:23:UNK:CB	2.21	0.71
11:e:31:UNK:O	11:e:35:UNK:N	2.24	0.70
16:8:28:ILE:C	16:8:31:TYR:HD1	2.00	0.70
18:P:23:ALA:O	19:u:9:UNK:CB	2.39	0.69
5:H:48:LEU:CD2	18:R:41:SER:OG	2.41	0.69
14:i:39:LEU:CD2	16:8:9:TRP:O	2.41	0.69
5:H:46:GLY:N	18:Q:38:ARG:O	2.18	0.69
14:i:39:LEU:HD11	16:8:11:ILE:CG1	2.21	0.69
15:k:9:UNK:CB	17:a:126:ALA:HB1	2.22	0.69
18:P:9:ILE:HD11	19:u:23:UNK:C	2.22	0.69
8:b:71:VAL:CG1	12:f:48:TYR:HA	2.19	0.67
1:C:182:THR:O	1:C:186:GLN:NE2	2.27	0.67
18:P:9:ILE:CD1	19:u:23:UNK:O	2.39	0.67
14:i:39:LEU:HD23	16:8:9:TRP:O	1.95	0.67
18:Q:20:GLY:HA2	19:u:12:UNK:CB	2.25	0.67
1:A:353:GLU:OE2	1:A:366:ASN:ND2	2.27	0.67
18:P:25:ILE:HD11	18:P:54:PHE:HA	1.76	0.67
18:P:16:VAL:HG13	19:u:16:UNK:O	1.94	0.66
8:b:56:GLU:HA	17:a:179:LEU:HG	1.77	0.65
2:E:279:SER:OG	2:E:280:ALA:N	2.29	0.65
16:8:21:PHE:O	16:8:25:GLN:NE2	2.29	0.65
5:H:48:LEU:HD21	18:R:41:SER:CA	2.18	0.65
1:C:410:LEU:HB3	1:C:413:ALA:HB3	1.79	0.65
4:G:161:ILE:HG12	4:G:175:GLU:HG3	1.79	0.64
2:E:465:ILE:O	2:E:469:VAL:HG12	1.97	0.64
2:F:11:ALA:HB1	2:F:27:PHE:CD2	2.33	0.64
8:b:56:GLU:HA	17:a:179:LEU:CD2	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:194:ASP:OD2	18:K:38:ARG:CD	2.47	0.63
14:i:39:LEU:CG	16:8:11:ILE:HG12	2.28	0.63
5:H:48:LEU:CD2	18:R:41:SER:CA	2.77	0.63
8:b:78:ILE:HG22	12:f:53:LYS:N	2.10	0.63
16:8:28:ILE:C	16:8:31:TYR:CD1	2.77	0.63
2:E:427:THR:HG23	2:E:429:HIS:H	1.63	0.62
3:J:4:THR:HG22	3:J:6:GLU:H	1.64	0.62
4:G:194:ASP:OD2	18:K:38:ARG:HD3	1.99	0.62
1:C:270:ASP:N	1:C:270:ASP:OD1	2.32	0.62
5:H:139:GLU:HG2	5:H:143:LYS:HE2	1.80	0.62
16:8:27:LYS:NZ	17:a:31:PHE:O	2.25	0.62
11:e:22:UNK:CA	13:g:82:UNK:O	2.47	0.62
2:E:473:ASP:OD1	2:E:473:ASP:N	2.33	0.61
11:e:37:UNK:O	11:e:41:UNK:CB	2.49	0.61
2:F:280:ALA:O	2:F:282:GLY:N	2.31	0.61
2:D:65:MET:HE1	2:D:233:ARG:HG3	1.82	0.61
14:i:39:LEU:CD2	16:8:11:ILE:HG13	2.16	0.61
1:A:483:ILE:HD11	1:A:493:THR:HG21	1.83	0.60
14:i:39:LEU:CD1	16:8:11:ILE:CD1	2.79	0.60
2:D:159:GLY:O	2:D:164:LYS:NZ	2.35	0.60
2:E:347:TYR:HB2	2:E:461:MET:HE1	1.84	0.60
4:G:146:LEU:HD22	6:I:16:GLN:HE22	1.67	0.60
5:H:45:PHE:CB	18:Q:38:ARG:HG3	2.30	0.60
8:b:89:LYS:C	8:b:92:GLN:HG3	2.21	0.60
1:C:186:GLN:HG2	1:C:199:LEU:HB3	1.83	0.60
2:D:137:THR:O	2:D:174:ASN:ND2	2.34	0.60
1:C:297:ASP:HB2	1:C:300:TYR:HB3	1.84	0.59
1:B:499:GLU:O	1:B:503:ASN:ND2	2.35	0.59
12:f:15:LYS:O	12:f:36:ALA:HB3	2.02	0.59
1:B:140:ILE:HD13	1:B:143:ARG:HH22	1.67	0.59
14:i:39:LEU:CD2	16:8:11:ILE:CD1	2.78	0.59
16:8:28:ILE:CA	16:8:31:TYR:CD1	2.74	0.59
17:a:163:ASN:HB2	18:N:62:LEU:CD1	2.33	0.59
2:D:246:ARG:NH1	2:D:299:THR:O	2.35	0.59
1:C:258:ARG:NH1	1:C:308:ARG:O	2.35	0.59
1:B:179:ALA:HB1	1:B:267:ILE:HD13	1.83	0.59
2:F:209:ASN:H	2:F:215:SER:HG	1.51	0.59
1:B:390:MET:HG3	1:B:424:LEU:HD13	1.85	0.58
16:8:17:ILE:HD12	17:a:100:MET:HB2	1.84	0.58
1:A:390:MET:HE1	1:A:428:LEU:HD21	1.85	0.58
14:i:39:LEU:CD2	16:8:11:ILE:HD11	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:196:LYS:NZ	1:B:261:GLY:O	2.33	0.58
2:D:423:ALA:HB2	22:D:501:ADP:HN61	1.69	0.58
8:b:74:TYR:CB	12:f:50:ASN:CB	2.81	0.58
2:D:392:ILE:HD11	4:G:16:ILE:HG23	1.86	0.58
7:S:78:PHE:HB2	7:S:80:PRO:HD3	1.85	0.58
16:8:28:ILE:CG1	16:8:31:TYR:CE1	2.86	0.58
17:a:123:THR:HA	17:a:126:ALA:HB3	1.85	0.58
1:C:420:ARG:NH1	1:C:449:VAL:O	2.37	0.57
17:a:151:ILE:HG23	18:M:62:LEU:CD2	2.34	0.57
2:D:34:ILE:HG22	2:D:35:LEU:HG	1.85	0.57
1:A:423:ARG:NH2	1:A:456:LEU:O	2.37	0.57
2:F:301:THR:HG23	2:F:303:LYS:H	1.69	0.57
18:K:17:GLY:HA2	18:R:19:ALA:HB2	1.86	0.57
2:D:11:ALA:O	2:D:27:PHE:CE1	2.57	0.57
15:k:7:UNK:C	17:a:129:LEU:HD21	2.35	0.57
17:a:170:LEU:HD11	18:O:63:PHE:CE1	2.40	0.57
18:K:42:LEU:HA	18:K:46:LEU:HD12	1.86	0.57
15:k:13:UNK:CB	17:a:142:VAL:HG11	2.35	0.57
2:E:172:ILE:HG21	2:E:217:VAL:HG22	1.87	0.56
18:P:23:ALA:HB1	19:u:9:UNK:CA	2.35	0.56
18:K:38:ARG:HH21	18:R:34:ILE:HD11	1.69	0.56
2:E:57:GLU:OE2	7:S:62:ARG:NH1	2.39	0.56
14:i:39:LEU:CD1	16:8:11:ILE:CG1	2.82	0.56
6:I:5:ARG:NH1	6:I:9:LEU:O	2.39	0.56
18:N:8:PHE:O	18:N:12:GLY:N	2.37	0.56
1:B:307:GLU:HG2	2:F:225:ASN:HD22	1.70	0.56
2:D:280:ALA:O	2:D:282:GLY:N	2.37	0.56
8:b:78:ILE:CG2	12:f:53:LYS:H	2.15	0.55
14:i:39:LEU:HD21	16:8:9:TRP:O	2.07	0.55
15:k:7:UNK:O	17:a:129:LEU:CD2	2.55	0.55
7:S:122:THR:HB	7:S:162:MET:HB3	1.87	0.55
17:a:71:MET:HE1	17:a:108:LEU:HD21	1.88	0.55
17:a:93:THR:HG1	17:a:96:THR:HG1	1.52	0.55
7:S:120:VAL:HG23	7:S:150:LEU:HD11	1.88	0.55
15:k:7:UNK:CA	17:a:129:LEU:HD21	2.34	0.55
18:Q:27:THR:HB	19:u:5:UNK:CB	2.37	0.55
14:i:39:LEU:HD22	16:8:11:ILE:CD1	2.37	0.55
2:F:11:ALA:HB1	2:F:27:PHE:HD2	1.72	0.55
14:i:39:LEU:HD13	16:8:11:ILE:CD1	2.37	0.55
1:B:307:GLU:HG2	2:F:225:ASN:HB3	1.88	0.54
1:C:25:LEU:HG	8:b:206:LYS:HD3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:248:TYR:OH	1:C:301:LEU:O	2.25	0.54
2:E:246:ARG:HD3	2:E:306:ILE:HG13	1.89	0.54
3:J:37:ARG:O	3:J:41:GLN:HG3	2.07	0.54
2:E:318:ASP:OD2	4:G:254:ARG:NH2	2.40	0.54
2:F:135:LEU:HB2	2:F:150:LYS:HG3	1.89	0.54
4:G:48:SER:OG	4:G:211:ASN:ND2	2.41	0.54
5:H:48:LEU:CD1	18:R:41:SER:OG	2.56	0.54
1:B:190:ASN:HA	1:B:198:LYS:HG2	1.90	0.54
1:C:359:LYS:HG3	2:F:381:GLN:HG2	1.89	0.54
1:B:218:LYS:NZ	1:B:222:ASP:OD2	2.41	0.54
4:G:166:ARG:HH21	4:G:170:SER:HB2	1.71	0.54
17:a:163:ASN:CB	18:N:62:LEU:HD12	2.38	0.54
18:L:50:ALA:O	18:L:54:PHE:N	2.41	0.54
1:A:15:ARG:NH2	7:S:91:GLU:OE1	2.41	0.53
1:B:46:ASN:O	1:B:90:ARG:NE	2.41	0.53
1:C:110:ALA:HB3	1:C:242:LEU:HD22	1.90	0.53
15:k:7:UNK:C	17:a:129:LEU:CD2	2.86	0.53
1:A:270:ASP:OD1	1:A:270:ASP:N	2.40	0.53
17:a:172:HIS:O	17:a:176:GLY:N	2.38	0.53
1:A:344:SER:O	2:E:191:ARG:NH2	2.41	0.53
18:M:42:LEU:HA	18:M:46:LEU:HB3	1.89	0.53
2:E:474:LYS:HD3	2:E:474:LYS:N	2.23	0.53
1:C:52:MET:O	1:C:91:THR:OG1	2.21	0.53
10:d:140:PRO:CB	16:8:40:ILE:CB	2.86	0.53
2:E:99:VAL:HG13	2:E:100:ILE:HG23	1.91	0.53
2:E:300:THR:HG23	2:E:305:SER:HB3	1.89	0.53
1:C:133:ALA:HB3	2:D:225:ASN:HD22	1.73	0.53
4:G:110:ASP:OD1	4:G:113:ARG:NH2	2.42	0.53
8:b:141:ARG:O	8:b:145:HIS:N	2.41	0.53
1:A:208:GLN:NE2	1:A:269:ASP:OD2	2.42	0.53
5:H:48:LEU:HD22	18:R:41:SER:OG	2.08	0.53
11:e:32:UNK:O	11:e:36:UNK:N	2.42	0.53
1:B:423:ARG:NH1	1:B:456:LEU:O	2.42	0.53
4:G:68:ILE:HB	4:G:105:ILE:HG22	1.91	0.53
2:D:454:LEU:O	3:J:41:GLN:OE1	2.26	0.52
1:A:191:ASP:OD1	1:A:227:LYS:NZ	2.42	0.52
1:B:175:LYS:NZ	20:B:601:ATP:O1B	2.37	0.52
2:D:160:ALA:O	2:D:339:ARG:NH2	2.36	0.52
2:E:173:ASN:ND2	2:E:421:GLN:OE1	2.42	0.52
8:b:61:ILE:O	8:b:65:GLY:N	2.43	0.52
2:F:40:VAL:HG22	2:F:77:VAL:HG22	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:209:ASN:ND2	2:F:214:THR:OG1	2.41	0.52
5:H:53:PRO:HA	5:H:84:VAL:HG13	1.92	0.52
5:H:34:ARG:HD3	5:H:68:GLU:HG3	1.90	0.52
18:P:16:VAL:HG11	19:u:16:UNK:O	2.10	0.52
2:D:58:SER:OG	2:D:58:SER:O	2.27	0.52
4:G:84:SER:HG	4:G:173:THR:HG1	1.58	0.51
18:M:30:GLY:O	18:N:31:SER:OG	2.27	0.51
1:A:423:ARG:HH12	1:A:458:PRO:HG3	1.75	0.51
2:E:15:ILE:CG2	2:E:16:VAL:N	2.73	0.51
2:F:212:ASP:OD1	2:F:212:ASP:N	2.44	0.51
18:L:25:ILE:HD11	18:L:54:PHE:HD1	1.75	0.51
1:A:69:ASP:N	1:A:69:ASP:OD1	2.43	0.51
1:B:297:ASP:OD1	1:B:297:ASP:N	2.41	0.51
2:D:410:ARG:NH1	3:J:30:GLU:OE1	2.43	0.51
2:F:271:SER:OG	2:F:276:ARG:NH1	2.43	0.51
6:I:37:THR:OG1	6:I:38:SER:N	2.43	0.51
13:g:67:UNK:O	13:g:71:UNK:N	2.44	0.51
1:C:16:ILE:HD12	7:S:184:ARG:HD3	1.91	0.51
2:F:399:SER:O	2:F:403:LYS:NZ	2.39	0.51
7:S:178:LYS:HD2	9:c:10:LEU:HD11	1.91	0.51
17:a:83:ASN:HB3	17:a:205:ALA:HB1	1.92	0.51
17:a:148:SER:OG	18:M:59:ALA:HB2	2.09	0.51
2:F:209:ASN:ND2	2:F:212:ASP:OD1	2.43	0.51
1:B:151:LYS:NZ	1:B:465:GLU:OE2	2.44	0.51
2:D:96:ILE:HD12	2:D:105:ASP:HB3	1.92	0.51
2:D:414:ARG:O	2:D:417:SER:OG	2.28	0.51
1:A:244:TYR:HE1	1:A:301:LEU:HD11	1.76	0.50
8:b:56:GLU:HA	17:a:179:LEU:CG	2.41	0.50
17:a:151:ILE:HD11	18:M:62:LEU:CD1	2.39	0.50
2:D:257:ILE:HG21	2:D:260:ILE:HD13	1.93	0.50
1:C:155:SER:O	1:C:379:GLN:NE2	2.44	0.50
1:C:396:GLN:HB3	1:C:417:LEU:HD21	1.94	0.50
2:F:257:ILE:HG21	2:F:260:ILE:HD13	1.93	0.50
8:b:141:ARG:NH1	10:d:43:ALA:O	2.45	0.50
19:u:21:UNK:O	19:u:25:UNK:N	2.44	0.50
1:C:9:SER:N	9:c:18:TYR:OH	2.44	0.50
19:u:26:UNK:O	19:u:30:UNK:N	2.44	0.50
1:B:393:GLU:OE2	1:B:420:ARG:NH2	2.45	0.50
1:C:411:ASP:HA	1:C:415:GLN:HB2	1.94	0.50
17:a:131:GLN:HE22	17:a:134:PRO:HD2	1.75	0.50
18:Q:31:SER:O	18:Q:35:GLY:N	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:387:ALA:HA	1:C:390:MET:HB3	1.93	0.50
17:a:109:TRP:CD2	17:a:153:PRO:HG3	2.47	0.50
18:Q:27:THR:HG21	19:u:5:UNK:CB	2.41	0.50
1:B:31:VAL:O	7:S:59:TYR:OH	2.25	0.50
5:H:50:SER:HB2	18:R:40:PRO:HB2	1.94	0.50
17:a:181:LEU:HD21	17:a:187:MET:HG3	1.94	0.50
5:H:33:VAL:HG13	5:H:67:ALA:HB2	1.93	0.49
17:a:163:ASN:CB	18:N:62:LEU:CD1	2.89	0.49
1:C:9:SER:N	9:c:18:TYR:HH	2.11	0.49
2:D:39:GLU:O	2:D:41:GLN:NE2	2.44	0.49
13:g:15:UNK:O	13:g:19:UNK:N	2.45	0.49
1:A:190:ASN:ND2	1:A:199:LEU:O	2.46	0.49
1:C:44:LEU:O	2:D:73:ARG:NH2	2.46	0.49
2:D:257:ILE:HB	2:D:310:GLN:HG2	1.94	0.49
2:E:15:ILE:HG22	2:E:16:VAL:N	2.27	0.49
11:e:25:UNK:O	11:e:29:UNK:N	2.46	0.49
2:E:107:ARG:HH11	2:E:210:LEU:HD23	1.78	0.49
16:8:12:THR:OG1	16:8:13:ILE:N	2.44	0.49
17:a:148:SER:OG	18:M:59:ALA:CB	2.61	0.49
17:a:163:ASN:HB2	18:N:62:LEU:HD12	1.95	0.49
18:M:22:GLY:O	18:N:21:SER:OG	2.28	0.49
1:A:453:LEU:HD13	1:A:461:ILE:HD12	1.94	0.49
1:C:28:THR:HG23	1:C:87:ILE:HG23	1.94	0.49
2:E:156:LEU:HD12	2:E:335:THR:HB	1.95	0.49
7:S:7:PRO:HG3	7:S:20:THR:HG22	1.95	0.49
10:d:140:PRO:CA	16:8:40:ILE:CB	2.90	0.49
17:a:96:THR:HG23	17:a:165:THR:HG22	1.95	0.49
18:P:23:ALA:HB1	19:u:9:UNK:HA	1.95	0.49
4:G:105:ILE:N	4:G:123:GLN:O	2.44	0.48
7:S:48:GLU:HB3	7:S:51:VAL:HB	1.95	0.48
18:K:19:ALA:HB2	18:L:17:GLY:HA2	1.93	0.48
19:u:4:UNK:O	19:u:8:UNK:N	2.46	0.48
1:C:164:ARG:HD2	1:C:309:ALA:HB3	1.95	0.48
2:E:48:VAL:HG21	2:E:100:ILE:HG21	1.95	0.48
7:S:134:LEU:HA	7:S:137:VAL:HG22	1.95	0.48
11:e:36:UNK:O	11:e:40:UNK:N	2.46	0.48
5:H:22:SER:OG	5:H:25:GLN:O	2.31	0.48
5:H:48:LEU:HD22	18:R:41:SER:N	2.28	0.48
18:K:37:ALA:O	18:L:39:ASN:ND2	2.38	0.48
2:D:11:ALA:CB	2:D:30:GLY:O	2.61	0.48
7:S:42:VAL:HA	7:S:45:ILE:HG12	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:421:GLY:O	1:C:425:THR:N	2.43	0.48
1:B:504:PHE:O	1:B:508:PHE:N	2.45	0.48
18:O:33:ILE:HG23	18:P:46:LEU:HD11	1.96	0.48
2:D:212:ASP:OD2	2:D:214:THR:OG1	2.31	0.48
8:b:148:TYR:O	8:b:152:LYS:NZ	2.46	0.48
17:a:111:ALA:O	17:a:115:THR:OG1	2.32	0.48
1:C:177:SER:HA	1:C:180:ILE:HG22	1.96	0.47
4:G:172:LYS:NZ	4:G:173:THR:O	2.43	0.47
1:C:413:ALA:HA	3:J:39:ARG:HH12	1.78	0.47
2:E:31:LEU:HD12	2:E:32:PRO:HD2	1.96	0.47
17:a:197:ILE:O	17:a:200:THR:OG1	2.24	0.47
1:A:410:LEU:HB2	1:A:415:GLN:HE21	1.78	0.47
1:C:218:LYS:HD2	2:F:130:VAL:HG21	1.95	0.47
8:b:55:ALA:O	17:a:179:LEU:CD1	2.60	0.47
16:8:27:LYS:NZ	17:a:28:SER:O	2.29	0.47
17:a:76:ILE:HB	17:a:213:VAL:HG21	1.95	0.47
2:D:286:THR:OG1	2:D:290:ASP:OD2	2.31	0.47
1:A:101:GLU:OE2	1:A:262:LYS:NZ	2.45	0.47
2:D:11:ALA:O	2:D:13:GLY:N	2.48	0.47
5:H:134:ARG:HH21	5:H:138:ASN:HD21	1.63	0.47
2:D:96:ILE:HG12	2:D:219:LEU:HD12	1.97	0.47
5:H:48:LEU:CG	18:R:41:SER:HB3	2.27	0.47
18:M:27:THR:HA	18:N:27:THR:HG21	1.96	0.47
2:F:11:ALA:CB	2:F:27:PHE:CE2	2.98	0.47
7:S:80:PRO:O	7:S:83:SER:OG	2.33	0.47
17:a:68:TRP:HD1	17:a:72:LEU:HD11	1.80	0.47
18:K:68:ALA:HA	18:K:71:ILE:HD12	1.96	0.47
18:M:8:PHE:HD1	18:N:71:ILE:HG23	1.80	0.47
18:M:18:VAL:O	18:M:22:GLY:N	2.47	0.47
2:F:11:ALA:HB1	2:F:27:PHE:CE2	2.50	0.47
15:k:10:UNK:CA	17:a:126:ALA:HB2	2.45	0.47
15:k:9:UNK:O	15:k:11:UNK:N	2.49	0.46
17:a:163:ASN:HB2	18:N:62:LEU:HD13	1.98	0.46
1:B:34:ILE:HG22	2:E:54:HIS:HB2	1.96	0.46
5:H:43:GLY:HA2	18:P:38:ARG:O	2.14	0.46
1:C:247:PRO:HG2	1:C:274:GLN:HG3	1.97	0.46
2:E:280:ALA:O	2:E:282:GLY:N	2.40	0.46
18:R:2:ILE:CD1	19:u:33:UNK:CB	2.89	0.46
5:H:55:LEU:HD22	5:H:83:THR:HG23	1.97	0.46
7:S:117:PRO:HB3	7:S:149:LYS:HG3	1.95	0.46
5:H:48:LEU:CD2	18:R:41:SER:N	2.78	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:a:109:TRP:O	17:a:112:THR:OG1	2.30	0.46
7:S:72:MET:HE3	7:S:72:MET:HB2	1.80	0.46
18:M:42:LEU:HD21	18:M:43:LYS:HZ3	1.81	0.46
18:R:3:ASP:OD1	18:R:3:ASP:N	2.41	0.46
1:C:96:ASP:OD1	1:C:96:ASP:N	2.49	0.46
5:H:52:VAL:HG21	18:R:37:ALA:O	2.16	0.46
10:d:3:ARG:HA	10:d:46:GLU:HA	1.98	0.46
17:a:64:LYS:HD3	17:a:64:LYS:HA	1.74	0.46
18:O:3:ASP:OD1	18:O:3:ASP:N	2.45	0.46
2:F:141:VAL:HG23	2:F:142:VAL:HG23	1.99	0.45
17:a:163:ASN:HB3	18:N:62:LEU:HD12	1.98	0.45
1:C:333:ASP:OD2	4:G:2:THR:HG23	2.16	0.45
8:b:63:THR:HA	17:a:199:LEU:HD23	1.97	0.45
18:K:64:CYS:O	18:K:68:ALA:N	2.48	0.45
1:C:25:LEU:HD22	7:S:169:LYS:HD2	1.97	0.45
17:a:151:ILE:HD13	17:a:151:ILE:HA	1.86	0.45
1:B:196:LYS:HA	1:B:196:LYS:HD2	1.80	0.45
1:B:207:GLY:HA3	1:B:273:LYS:HD3	1.97	0.45
1:C:431:GLY:O	20:C:601:ATP:N6	2.49	0.45
7:S:61:LYS:HD2	7:S:61:LYS:HA	1.64	0.45
15:k:9:UNK:N	17:a:129:LEU:HD23	2.32	0.45
2:E:410:ARG:HE	2:E:456:GLU:HG2	1.82	0.45
2:F:38:LEU:HB2	2:F:49:LEU:HB2	1.98	0.45
2:F:209:ASN:N	2:F:215:SER:OG	2.38	0.45
1:A:6:ALA:HB3	1:A:69:ASP:HB2	1.97	0.45
1:A:44:LEU:HD13	1:A:47:VAL:HG21	1.99	0.45
1:C:167:ILE:HD11	1:C:324:LEU:HD22	1.97	0.45
7:S:22:LEU:HD13	7:S:85:LEU:HD22	1.98	0.45
18:N:3:ASP:O	18:N:7:LYS:N	2.47	0.45
2:D:257:ILE:HD12	2:D:310:GLN:HE21	1.82	0.45
7:S:108:MET:HE3	7:S:108:MET:HB2	1.80	0.45
8:b:139:THR:HG22	8:b:143:ARG:HE	1.81	0.45
1:B:176:THR:OG1	1:B:269:ASP:OD2	2.35	0.45
2:E:13:GLY:O	2:E:77:VAL:HG12	2.16	0.45
2:E:376:VAL:HG11	2:E:443:PHE:HD2	1.82	0.45
18:M:60:MET:HB2	18:M:60:MET:HE2	1.64	0.45
18:Q:13:ALA:CB	19:u:23:UNK:CB	2.92	0.45
1:B:366:ASN:HD22	1:B:369:LEU:HB2	1.81	0.44
2:E:291:MET:HE2	2:E:291:MET:HB2	1.81	0.44
2:E:377:GLN:O	2:E:381:GLN:NE2	2.50	0.44
4:G:3:LEU:O	4:G:7:THR:OG1	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:O:15:THR:HG21	18:P:67:VAL:HG21	1.99	0.44
8:b:75:GLY:O	8:b:78:ILE:HD11	2.17	0.44
16:8:20:LEU:HD13	17:a:24:ILE:HG22	2.00	0.44
2:D:157:PHE:HB2	2:D:336:VAL:HG22	1.99	0.44
18:P:60:MET:HE2	18:P:60:MET:HB3	1.87	0.44
1:B:156:LEU:HD21	1:B:390:MET:HE3	1.99	0.44
1:B:469:LEU:HA	1:B:472:VAL:HG12	1.99	0.44
2:D:18:VAL:HG13	2:D:23:VAL:HG22	1.99	0.44
18:M:45:GLN:H	18:M:45:GLN:HG2	1.58	0.44
18:O:44:GLN:HG2	18:O:45:GLN:H	1.83	0.44
18:P:16:VAL:HG21	19:u:20:UNK:CB	2.48	0.44
5:H:63:VAL:HG13	5:H:75:TYR:HB2	1.99	0.44
16:8:13:ILE:H	16:8:13:ILE:HG13	1.42	0.44
18:Q:33:ILE:HG21	18:R:32:MET:HA	1.99	0.44
1:C:356:LEU:HB2	1:C:364:ALA:HB1	1.98	0.44
8:b:72:LYS:HD2	8:b:73:LYS:HG3	2.00	0.44
8:b:112:LYS:HA	8:b:115:GLN:HB3	1.99	0.44
16:8:12:THR:HG21	17:a:17:LEU:HD11	1.99	0.44
8:b:75:GLY:HA2	8:b:78:ILE:CD1	2.47	0.44
7:S:60:VAL:HG23	7:S:61:LYS:H	1.82	0.44
18:O:50:ALA:O	18:O:54:PHE:N	2.43	0.44
1:B:79:ASP:OD1	1:B:79:ASP:N	2.49	0.43
2:F:259:ASN:HD21	2:F:313:TYR:HB2	1.83	0.43
8:b:113:SER:OG	8:b:114:GLN:OE1	2.36	0.43
17:a:79:ILE:HG22	17:a:95:THR:HG21	2.00	0.43
17:a:151:ILE:CG1	18:M:62:LEU:CD1	2.74	0.43
1:B:286:ARG:NH2	2:E:275:GLY:O	2.51	0.43
5:H:54:THR:H	5:H:84:VAL:HG13	1.83	0.43
18:L:36:TYR:HD2	18:M:46:LEU:HD13	1.82	0.43
1:A:99:VAL:HG22	1:A:253:MET:HA	1.99	0.43
1:A:148:THR:HG23	1:A:182:THR:HG23	1.99	0.43
18:Q:27:THR:CB	19:u:5:UNK:CB	2.97	0.43
2:D:11:ALA:HB1	2:D:27:PHE:CD1	2.53	0.43
2:D:27:PHE:O	2:D:58:SER:OG	2.24	0.43
1:A:168:ILE:HG23	1:A:351:PHE:HD1	1.83	0.43
1:B:453:LEU:HB2	1:B:456:LEU:HD22	2.00	0.43
2:F:164:LYS:NZ	22:F:501:ADP:O1B	2.51	0.43
6:I:23:ARG:HA	6:I:26:LEU:HD13	2.01	0.43
2:D:197:ASP:O	2:D:201:GLU:N	2.40	0.43
1:A:175:LYS:HB2	1:A:175:LYS:HE3	1.80	0.43
1:B:289:PRO:HB2	1:B:293:ALA:HA	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:145:LEU:HD21	2:E:376:VAL:HG23	2.01	0.43
4:G:156:ASP:HB3	4:G:181:LEU:HD22	2.01	0.43
7:S:48:GLU:HG3	7:S:50:LYS:H	1.83	0.43
18:R:56:LEU:HA	18:R:56:LEU:HD12	1.78	0.43
1:B:89:LYS:HE3	1:B:89:LYS:HB2	1.86	0.43
1:B:129:VAL:O	1:B:308:ARG:NH2	2.48	0.43
1:B:187:LYS:HE2	1:B:224:ASP:HB3	2.00	0.43
2:F:294:MET:HB3	2:F:294:MET:HE3	1.82	0.43
13:g:11:UNK:O	13:g:15:UNK:N	2.51	0.43
16:8:19:THR:OG1	17:a:25:MET:HE1	2.19	0.43
18:L:51:ILE:HD13	18:L:51:ILE:HA	1.83	0.43
18:Q:69:PHE:O	18:Q:73:PHE:N	2.44	0.43
1:A:409:ASP:OD2	4:G:33:ARG:NH2	2.42	0.42
1:C:203:TYR:OH	1:C:269:ASP:OD2	2.33	0.42
2:E:96:ILE:HG12	2:E:219:LEU:HB2	2.01	0.42
2:E:321:ASP:HB3	2:E:324:PRO:HD2	2.01	0.42
2:F:14:ARG:HB3	2:F:74:GLY:HA2	2.02	0.42
4:G:43:VAL:HA	4:G:46:ILE:HD12	2.01	0.42
17:a:50:ILE:HD13	17:a:50:ILE:HA	1.88	0.42
18:P:46:LEU:HD12	18:P:46:LEU:HA	1.85	0.42
1:C:183:ILE:O	1:C:228:TYR:OH	2.25	0.42
2:D:12:THR:HA	2:D:77:VAL:O	2.18	0.42
2:D:283:TYR:HB3	2:D:287:LEU:HD13	2.00	0.42
14:i:22:THR:HB	14:i:26:ARG:HB2	2.00	0.42
18:R:32:MET:HE3	18:R:32:MET:HB3	1.92	0.42
1:A:288:PRO:HA	1:A:289:PRO:HD3	1.82	0.42
1:C:283:LEU:HD22	2:F:277:ILE:HB	2.01	0.42
2:D:94:GLY:O	2:D:107:ARG:NH1	2.53	0.42
2:E:236:LEU:O	2:E:240:THR:OG1	2.35	0.42
4:G:197:ASP:H	4:G:200:VAL:HB	1.84	0.42
1:A:176:THR:OG1	20:A:601:ATP:O2B	2.38	0.42
1:B:166:LEU:HB2	1:B:346:THR:HG21	2.01	0.42
1:B:415:GLN:O	1:B:415:GLN:NE2	2.53	0.42
1:C:222:ASP:OD1	1:C:223:ALA:N	2.52	0.42
2:E:336:VAL:HG23	2:E:351:ASP:HB3	2.01	0.42
4:G:59:VAL:HA	4:G:60:PRO:HD3	1.89	0.42
4:G:113:ARG:HG3	4:G:127:THR:HG21	2.01	0.42
4:G:156:ASP:OD1	4:G:156:ASP:N	2.40	0.42
1:A:163:GLN:NE2	1:A:165:GLU:OE1	2.52	0.42
2:F:99:VAL:HG21	2:F:230:ALA:HB1	2.01	0.42
1:A:427:LEU:HD11	1:A:448:GLY:HA3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:k:3:UNK:HA	15:k:6:UNK:HA	2.01	0.42
18:P:49:TYR:O	18:P:53:GLY:N	2.51	0.42
1:A:76:PHE:CE1	1:A:241:PRO:HB2	2.55	0.42
1:B:140:ILE:HG22	1:B:311:LYS:HG2	2.01	0.42
5:H:124:ASP:OD1	5:H:124:ASP:N	2.49	0.42
12:f:66:TYR:HE2	17:a:91:SER:HB3	1.84	0.42
1:B:215:GLN:OE1	2:E:132:GLN:NE2	2.53	0.42
18:M:72:LEU:HD12	18:M:72:LEU:HA	1.88	0.42
1:B:181:ASP:OD1	1:B:181:ASP:N	2.52	0.42
1:B:336:ALA:HB3	1:B:339:PRO:HD2	2.02	0.42
3:J:45:LEU:HD23	3:J:45:LEU:HA	1.76	0.42
4:G:167:SER:OG	4:G:170:SER:OG	2.30	0.42
18:Q:27:THR:CG2	19:u:5:UNK:CB	2.98	0.42
1:B:164:ARG:NH2	2:F:191:ARG:HD3	2.35	0.42
16:8:12:THR:O	16:8:14:THR:N	2.53	0.42
1:B:400:VAL:HB	1:B:414:THR:HG23	2.02	0.41
2:E:138:GLY:HA3	2:E:433:LEU:HD11	2.02	0.41
16:8:32:SER:OG	17:a:47:GLN:NE2	2.53	0.41
1:B:273:LYS:HA	1:B:276:VAL:HG22	2.01	0.41
4:G:93:VAL:HG21	4:G:120:HIS:HE2	1.85	0.41
8:b:89:LYS:HA	8:b:92:GLN:HG2	2.02	0.41
1:B:104:LEU:HD13	1:B:228:TYR:HA	2.02	0.41
1:B:147:GLN:HG3	1:B:438:ILE:HD13	2.01	0.41
2:D:145:LEU:HD23	2:D:373:ALA:HB1	2.02	0.41
2:F:13:GLY:O	2:F:77:VAL:N	2.53	0.41
4:G:187:ALA:H	4:G:190:MET:HG2	1.85	0.41
8:b:185:ILE:HD12	8:b:185:ILE:HA	1.94	0.41
11:e:32:UNK:O	11:e:36:UNK:CB	2.68	0.41
12:f:63:LEU:HD23	17:a:88:LEU:HD11	2.01	0.41
12:f:73:ARG:H	12:f:73:ARG:HG2	1.69	0.41
1:C:210:ARG:HG3	1:C:235:THR:HG21	2.02	0.41
4:G:71:VAL:HA	4:G:108:VAL:HG13	2.02	0.41
16:8:34:PRO:HG2	17:a:51:GLN:CD	2.46	0.41
1:B:170:ASP:O	1:B:173:THR:OG1	2.31	0.41
2:D:399:SER:OG	2:D:400:GLU:N	2.54	0.41
8:b:184:SER:O	8:b:188:GLN:NE2	2.53	0.41
18:M:59:ALA:O	18:M:63:PHE:HB2	2.21	0.41
1:A:31:VAL:HG22	1:A:41:VAL:HG22	2.03	0.41
1:A:258:ARG:NH1	1:A:308:ARG:O	2.54	0.41
4:G:69:ILE:HB	4:G:160:ILE:HG13	2.03	0.41
15:k:7:UNK:O	17:a:129:LEU:HD21	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:97:MET:HG2	2:D:220:VAL:HG22	2.02	0.41
5:H:52:VAL:CG2	18:R:37:ALA:O	2.69	0.41
6:I:4:TRP:H	6:I:4:TRP:CD1	2.38	0.41
18:N:32:MET:HE1	18:O:49:TYR:CG	2.55	0.41
1:A:146:MET:HE3	1:A:146:MET:HB3	1.81	0.41
1:B:127:ARG:NH2	1:B:255:GLU:OE1	2.54	0.41
17:a:156:LEU:HA	17:a:156:LEU:HD12	1.84	0.41
18:M:46:LEU:HD12	18:M:46:LEU:HA	1.88	0.41
18:P:9:ILE:HD12	19:u:27:UNK:CB	2.36	0.41
1:B:211:SER:OG	1:B:215:GLN:NE2	2.54	0.41
1:B:271:LEU:HD13	1:B:302:HIS:CD2	2.55	0.41
1:C:374:VAL:HG13	1:C:377:ALA:H	1.85	0.41
2:D:286:THR:OG1	2:D:286:THR:O	2.36	0.41
2:E:95:ARG:NH2	2:E:108:GLY:O	2.44	0.41
2:F:11:ALA:CB	2:F:27:PHE:HE2	2.34	0.41
4:G:112:ILE:HD13	4:G:112:ILE:HA	1.88	0.41
4:G:220:SER:O	4:G:220:SER:OG	2.37	0.41
16:8:20:LEU:HD12	16:8:20:LEU:HA	1.93	0.41
18:K:3:ASP:N	18:K:3:ASP:OD1	2.54	0.41
1:C:101:GLU:H	1:C:101:GLU:HG2	1.70	0.41
2:E:336:VAL:HG21	2:E:354:ASP:HB3	2.03	0.41
2:F:89:GLY:O	2:F:92:THR:OG1	2.35	0.41
18:M:19:ALA:HB2	18:N:17:GLY:HA2	2.03	0.41
18:N:19:ALA:HB1	18:O:20:GLY:HA3	2.03	0.41
1:B:235:THR:OG1	1:B:236:ALA:N	2.52	0.40
2:E:12:THR:O	2:E:12:THR:HG23	2.21	0.40
3:J:37:ARG:HA	3:J:37:ARG:HD3	1.90	0.40
4:G:194:ASP:O	4:G:196:ILE:N	2.54	0.40
15:k:10:UNK:N	17:a:126:ALA:HB2	2.32	0.40
19:u:19:UNK:O	19:u:23:UNK:N	2.54	0.40
1:A:48:GLN:HB3	2:E:70:GLY:HA2	2.04	0.40
1:C:30:ARG:HD2	1:C:85:GLY:HA2	2.02	0.40
2:D:393:LEU:HD22	2:D:397:GLU:HG2	2.02	0.40
9:c:44:PHE:O	9:c:48:GLN:N	2.54	0.40
11:e:9:UNK:O	11:e:13:UNK:N	2.54	0.40
1:C:438:ILE:HD12	1:C:438:ILE:HA	1.93	0.40
2:D:204:GLU:H	2:D:204:GLU:HG3	1.75	0.40
4:G:104:LYS:HA	4:G:123:GLN:HB3	2.03	0.40
20:B:601:ATP:O2G	2:E:358:ARG:NH2	2.36	0.40
17:a:169:LEU:HD13	17:a:169:LEU:HA	1.94	0.40
18:N:9:ILE:O	18:N:13:ALA:N	2.49	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	507/509 (100%)	479 (94%)	28 (6%)	0	100	100
1	B	477/509 (94%)	455 (95%)	22 (5%)	0	100	100
1	C	500/509 (98%)	464 (93%)	36 (7%)	0	100	100
2	D	469/480 (98%)	422 (90%)	45 (10%)	2 (0%)	30	59
2	E	465/480 (97%)	422 (91%)	42 (9%)	1 (0%)	43	71
2	F	466/480 (97%)	424 (91%)	41 (9%)	1 (0%)	43	71
3	J	53/83 (64%)	51 (96%)	2 (4%)	0	100	100
4	G	270/272 (99%)	254 (94%)	14 (5%)	2 (1%)	18	49
5	H	130/132 (98%)	117 (90%)	12 (9%)	1 (1%)	16	46
6	I	46/48 (96%)	39 (85%)	7 (15%)	0	100	100
7	S	185/187 (99%)	163 (88%)	22 (12%)	0	100	100
8	b	207/209 (99%)	200 (97%)	7 (3%)	0	100	100
9	c	68/70 (97%)	61 (90%)	7 (10%)	0	100	100
10	d	145/147 (99%)	126 (87%)	19 (13%)	0	100	100
12	f	76/87 (87%)	63 (83%)	12 (16%)	1 (1%)	9	35
14	i	40/42 (95%)	39 (98%)	1 (2%)	0	100	100
16	8	54/67 (81%)	46 (85%)	4 (7%)	4 (7%)	1	6
17	a	221/226 (98%)	194 (88%)	27 (12%)	0	100	100
18	K	70/72 (97%)	67 (96%)	3 (4%)	0	100	100
18	L	70/72 (97%)	66 (94%)	4 (6%)	0	100	100
18	M	70/72 (97%)	65 (93%)	5 (7%)	0	100	100
18	N	70/72 (97%)	67 (96%)	3 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
18	O	70/72 (97%)	66 (94%)	4 (6%)	0	100	100
18	P	70/72 (97%)	67 (96%)	3 (4%)	0	100	100
18	Q	70/72 (97%)	66 (94%)	4 (6%)	0	100	100
18	R	70/72 (97%)	65 (93%)	5 (7%)	0	100	100
All	All	4939/5113 (97%)	4548 (92%)	379 (8%)	12 (0%)	44	71

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	12	THR
2	D	281	VAL
2	E	281	VAL
2	F	281	VAL
4	G	195	ASP
16	8	13	ILE
16	8	7	SER
16	8	8	THR
16	8	12	THR
5	H	49	ALA
12	f	17	GLY
4	G	60	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	A	410/410 (100%)	404 (98%)	6 (2%)	57	71
1	B	389/410 (95%)	388 (100%)	1 (0%)	86	85
1	C	404/410 (98%)	399 (99%)	5 (1%)	63	74
2	D	378/386 (98%)	375 (99%)	3 (1%)	73	78
2	E	375/386 (97%)	371 (99%)	4 (1%)	65	75
2	F	375/386 (97%)	374 (100%)	1 (0%)	86	85

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	J	37/68 (54%)	36 (97%)	1 (3%)	39	63
4	G	230/230 (100%)	229 (100%)	1 (0%)	84	83
5	H	105/105 (100%)	104 (99%)	1 (1%)	68	75
6	I	38/38 (100%)	38 (100%)	0	100	100
7	S	163/163 (100%)	163 (100%)	0	100	100
8	b	123/182 (68%)	121 (98%)	2 (2%)	55	71
9	c	22/63 (35%)	22 (100%)	0	100	100
10	d	1/127 (1%)	1 (100%)	0	100	100
12	f	27/75 (36%)	27 (100%)	0	100	100
14	i	28/36 (78%)	28 (100%)	0	100	100
16	8	29/66 (44%)	28 (97%)	1 (3%)	32	59
17	a	184/199 (92%)	181 (98%)	3 (2%)	55	71
18	K	49/49 (100%)	49 (100%)	0	100	100
18	L	49/49 (100%)	48 (98%)	1 (2%)	48	67
18	M	49/49 (100%)	48 (98%)	1 (2%)	48	67
18	N	48/49 (98%)	47 (98%)	1 (2%)	47	66
18	O	48/49 (98%)	48 (100%)	0	100	100
18	P	49/49 (100%)	48 (98%)	1 (2%)	48	67
18	Q	49/49 (100%)	49 (100%)	0	100	100
18	R	48/49 (98%)	48 (100%)	0	100	100
All	All	3707/4132 (90%)	3674 (99%)	33 (1%)	68	77

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLU
1	A	99	VAL
1	A	125	THR
1	A	176	THR
1	A	206	ILE
1	A	380	THR
1	B	88	VAL
1	C	69	ASP
1	C	150	ILE
1	C	229	THR

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Mol	Chain	Res	Type
1	C	270	ASP
1	C	334	VAL
2	D	72	VAL
2	D	130	VAL
2	D	253	VAL
2	E	45	THR
2	E	192	THR
2	E	469	VAL
2	E	473	ASP
2	F	48	VAL
3	J	45	LEU
4	G	108	VAL
5	H	63	VAL
8	b	78	ILE
8	b	90	ILE
16	8	13	ILE
17	a	121	THR
17	a	142	VAL
17	a	185	ASN
18	L	42	LEU
18	M	42	LEU
18	N	65	LEU
18	P	15	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	HIS
1	A	405	GLN
1	A	476	HIS
1	B	186	GLN
1	B	208	GLN
1	B	215	GLN
1	B	366	ASN
1	B	385	GLN
1	B	396	GLN
1	B	466	ASN
1	B	503	ASN
1	C	70	ASN
1	C	208	GLN
1	C	330	GLN
1	C	341	ASN

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Mol	Chain	Res	Type
1	C	441	GLN
2	D	363	ASN
2	D	369	HIS
2	E	119	HIS
2	E	132	GLN
2	E	179	HIS
2	E	209	ASN
2	E	369	HIS
2	F	225	ASN
2	F	295	GLN
4	G	82	HIS
4	G	88	GLN
4	G	211	ASN
4	G	234	ASN
5	H	35	GLN
6	I	16	GLN
6	I	41	ASN
8	b	145	HIS
8	b	188	GLN
9	c	23	GLN
17	a	39	ASN
17	a	47	GLN
17	a	63	GLN
17	a	83	ASN
17	a	152	GLN
18	L	44	GLN
18	P	39	ASN
18	R	45	GLN

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 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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
22	ADP	D	501	21	28,29,29	1.49	6 (21%)	43,45,45	2.27	9 (20%)
20	ATP	C	601	21	32,33,33	1.33	5 (15%)	48,52,52	1.79	7 (14%)
20	ATP	A	601	21	32,33,33	1.39	5 (15%)	48,52,52	1.80	7 (14%)
20	ATP	B	601	21	32,33,33	1.29	4 (12%)	48,52,52	1.82	10 (20%)
22	ADP	F	501	21	28,29,29	1.38	4 (14%)	43,45,45	1.80	8 (18%)

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
22	ADP	D	501	21	-	9/16/32/32	0/3/3/3
20	ATP	C	601	21	-	0/22/38/38	0/3/3/3
20	ATP	A	601	21	-	5/22/38/38	0/3/3/3
20	ATP	B	601	21	-	8/22/38/38	0/3/3/3
22	ADP	F	501	21	-	5/16/32/32	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	F	501	ADP	C5-C4	4.34	1.46	1.39
20	B	601	ATP	C5-C4	4.29	1.46	1.39
20	C	601	ATP	C5-C4	4.23	1.46	1.39
22	D	501	ADP	C5-C4	4.18	1.46	1.39
20	A	601	ATP	C5-C4	3.86	1.46	1.39
20	A	601	ATP	C5-N7	-3.12	1.33	1.39
22	F	501	ADP	C5-N7	-2.85	1.33	1.39
22	D	501	ADP	C5-N7	-2.70	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
20	A	601	ATP	C8-N9	-2.64	1.33	1.37
20	B	601	ATP	C5-C6	2.57	1.48	1.41
20	C	601	ATP	C5-N7	-2.52	1.34	1.39
22	D	501	ADP	PA-O3A	2.38	1.62	1.59
22	D	501	ADP	C4-N9	-2.34	1.32	1.37
20	B	601	ATP	C5-N7	-2.30	1.34	1.39
20	A	601	ATP	C4-N9	-2.30	1.32	1.37
20	C	601	ATP	C5-C6	2.26	1.47	1.41
22	D	501	ADP	C8-N7	2.24	1.36	1.31
22	F	501	ADP	C5-C6	2.20	1.47	1.41
22	D	501	ADP	C5-C6	2.17	1.47	1.41
20	B	601	ATP	C8-N7	2.15	1.35	1.31
20	A	601	ATP	C5-C6	2.14	1.47	1.41
20	C	601	ATP	C4-N9	-2.12	1.33	1.37
22	F	501	ADP	C4-N9	-2.04	1.33	1.37
20	C	601	ATP	C8-N7	2.00	1.35	1.31

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	D	501	ADP	N3-C2-N1	-7.91	116.61	128.58
20	A	601	ATP	C5-C4-N3	-6.17	118.22	126.72
20	B	601	ATP	C5-C4-N3	-6.05	118.38	126.72
22	F	501	ADP	C5-C4-N3	-5.88	118.62	126.72
20	C	601	ATP	C5-C4-N3	-5.63	118.96	126.72
22	D	501	ADP	C5-C4-N3	-5.46	119.19	126.72
22	D	501	ADP	C2-N3-C4	5.17	124.44	111.83
20	B	601	ATP	N3-C4-N9	4.97	135.61	127.17
20	A	601	ATP	N3-C4-N9	4.73	135.21	127.17
20	C	601	ATP	N3-C4-N9	4.70	135.17	127.17
22	F	501	ADP	N3-C4-N9	4.69	135.14	127.17
22	D	501	ADP	C2-N1-C6	4.38	125.93	118.73
22	D	501	ADP	N3-C4-N9	3.99	133.95	127.17
20	B	601	ATP	C2-N3-C4	3.87	121.29	111.83
20	A	601	ATP	C2-N3-C4	3.77	121.03	111.83
22	F	501	ADP	C2-N3-C4	3.71	120.90	111.83
20	C	601	ATP	C2-N3-C4	3.68	120.83	111.83
20	A	601	ATP	C4-C5-N7	-3.60	106.46	110.58
22	F	501	ADP	N3-C2-N1	-3.41	123.42	128.58
20	C	601	ATP	N3-C2-N1	-3.40	123.43	128.58
20	B	601	ATP	C4-C5-N7	-3.36	106.73	110.58
20	B	601	ATP	N3-C2-N1	-3.33	123.54	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
20	A	601	ATP	N3-C2-N1	-3.22	123.71	128.58
20	B	601	ATP	C4-N9-C8	3.05	108.94	105.74
22	F	501	ADP	C4-C5-N7	-2.98	107.17	110.58
22	D	501	ADP	C4-C5-N7	-2.97	107.18	110.58
20	C	601	ATP	C4-C5-N7	-2.95	107.21	110.58
20	C	601	ATP	C4-N9-C8	2.83	108.71	105.74
22	D	501	ADP	C2'-C1'-N9	-2.66	106.71	113.30
22	D	501	ADP	C6-C5-N7	2.63	137.17	132.09
20	B	601	ATP	C5-N7-C8	2.54	107.44	103.45
20	A	601	ATP	C5-N7-C8	2.51	107.40	103.45
20	A	601	ATP	C4-N9-C8	2.43	108.28	105.74
20	C	601	ATP	C5-N7-C8	2.20	106.91	103.45
20	B	601	ATP	N9-C8-N7	-2.15	110.88	113.94
20	B	601	ATP	C3'-C2'-C1'	2.12	105.47	101.46
22	F	501	ADP	C5-N7-C8	2.12	106.78	103.45
20	B	601	ATP	C6-C5-N7	2.08	136.10	132.09
22	D	501	ADP	C3'-C2'-C1'	2.07	105.39	101.46
22	F	501	ADP	C3'-C2'-C1'	2.05	105.35	101.46
22	F	501	ADP	C4-N9-C8	2.04	107.88	105.74

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
20	A	601	ATP	C5'-O5'-PA-O1A
20	A	601	ATP	C5'-O5'-PA-O3A
20	B	601	ATP	C5'-O5'-PA-O1A
20	B	601	ATP	C5'-O5'-PA-O2A
20	B	601	ATP	C5'-O5'-PA-O3A
22	D	501	ADP	PA-O3A-PB-O3B
22	D	501	ADP	C5'-O5'-PA-O2A
22	D	501	ADP	C5'-O5'-PA-O3A
22	F	501	ADP	PB-O3A-PA-O5'
22	F	501	ADP	C5'-O5'-PA-O1A
22	F	501	ADP	C5'-O5'-PA-O2A
22	F	501	ADP	C5'-O5'-PA-O3A
20	A	601	ATP	O4'-C4'-C5'-O5'
20	B	601	ATP	C3'-C4'-C5'-O5'
20	B	601	ATP	O4'-C4'-C5'-O5'
20	B	601	ATP	PB-O3B-PG-O1G
20	A	601	ATP	C3'-C4'-C5'-O5'
22	D	501	ADP	O4'-C1'-N9-C4

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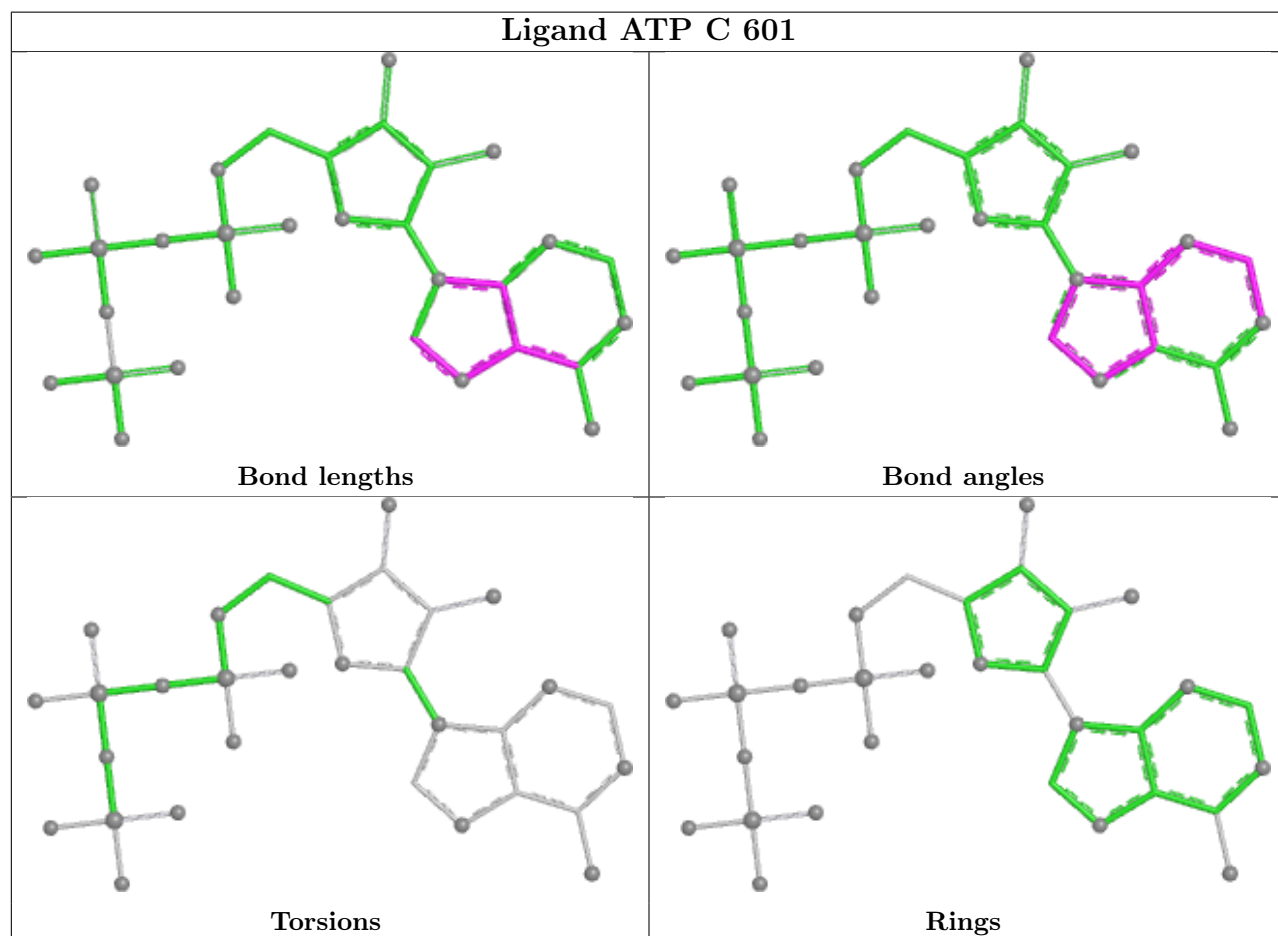
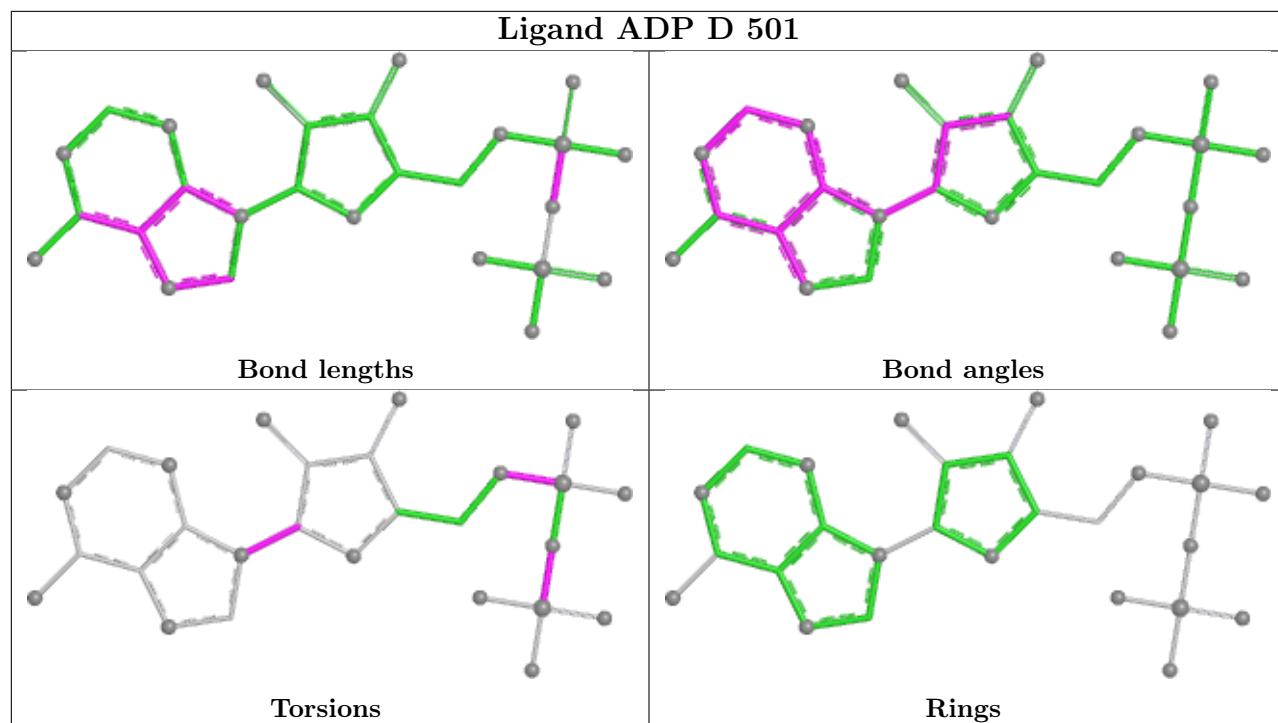
Mol	Chain	Res	Type	Atoms
22	D	501	ADP	O4'-C1'-N9-C8
20	A	601	ATP	C5'-O5'-PA-O2A
22	D	501	ADP	C5'-O5'-PA-O1A
22	F	501	ADP	PA-O3A-PB-O1B
20	B	601	ATP	PA-O3A-PB-O2B
22	D	501	ADP	C2'-C1'-N9-C8
22	D	501	ADP	PA-O3A-PB-O1B
22	D	501	ADP	PA-O3A-PB-O2B
20	B	601	ATP	PA-O3A-PB-O1B

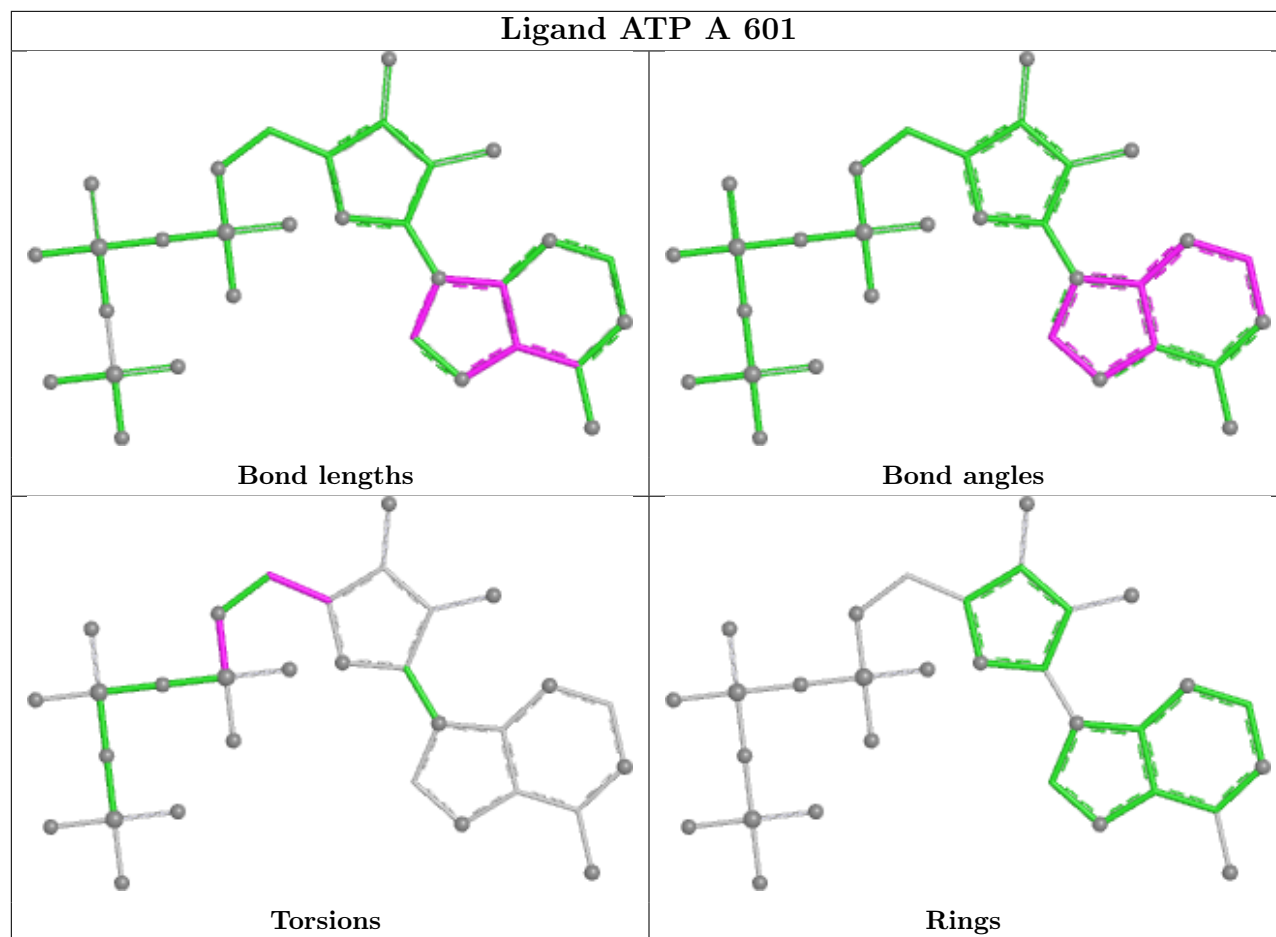
There are no ring outliers.

5 monomers are involved in 6 short contacts:

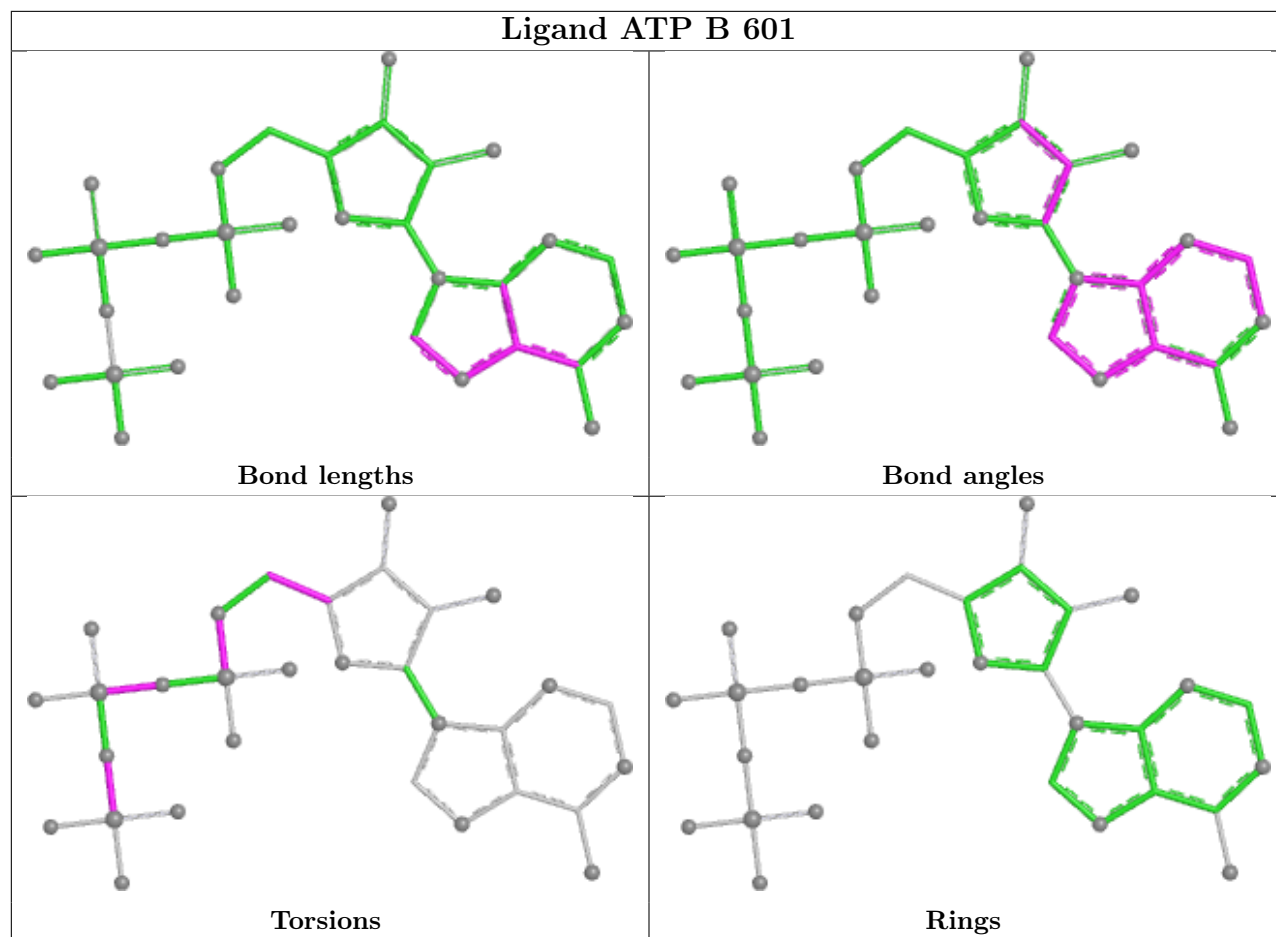
Mol	Chain	Res	Type	Clashes	Symm-Clashes
22	D	501	ADP	1	0
20	C	601	ATP	1	0
20	A	601	ATP	1	0
20	B	601	ATP	2	0
22	F	501	ADP	1	0

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

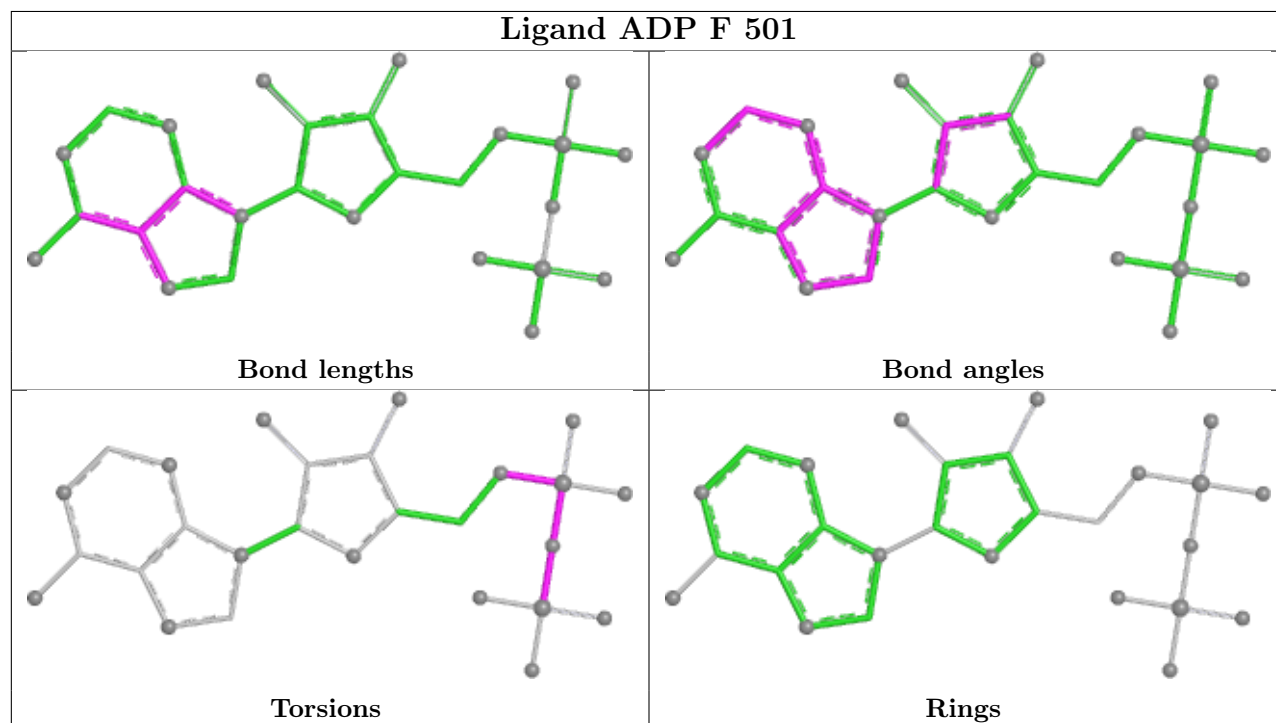




Ligand ATP B 601



Ligand ADP F 501



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

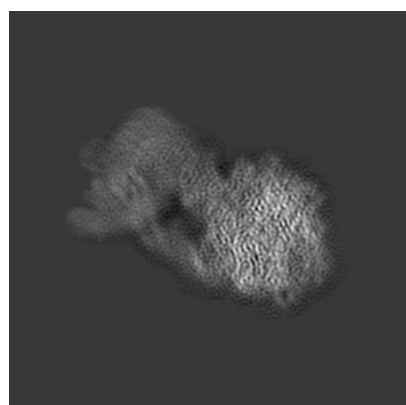
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-0677. These allow visual inspection of the internal detail of the map and identification of artifacts.

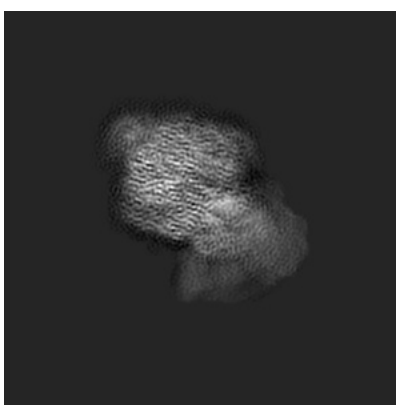
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

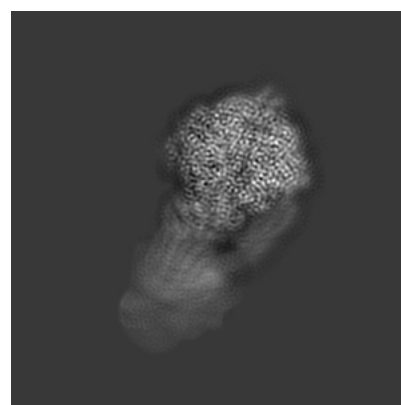
6.1.1 Primary map



X



Y

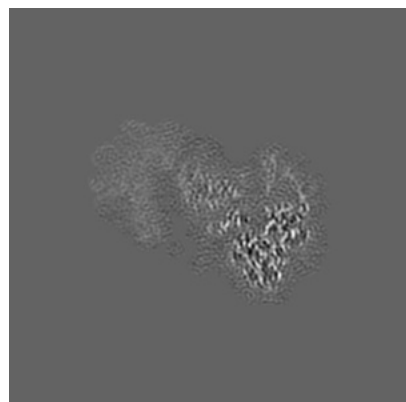


Z

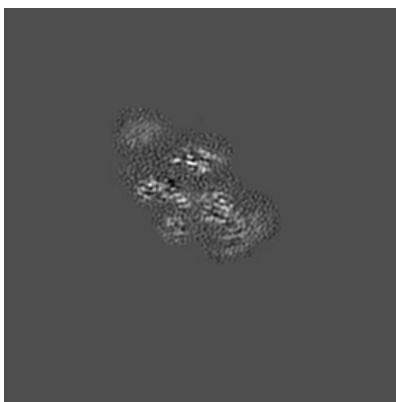
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 120



Y Index: 120

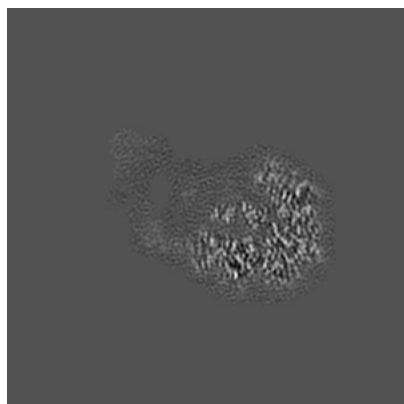


Z Index: 120

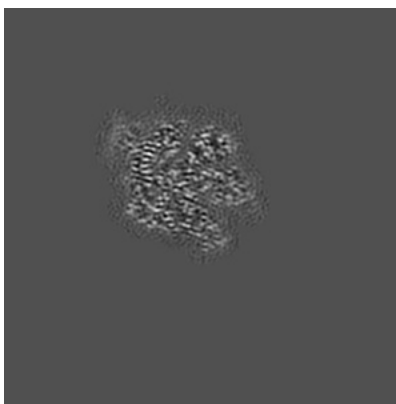
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

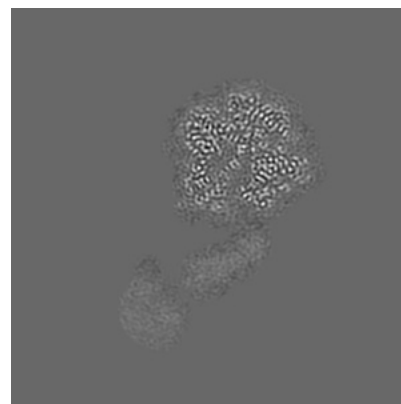
6.3.1 Primary map



X Index: 134



Y Index: 161

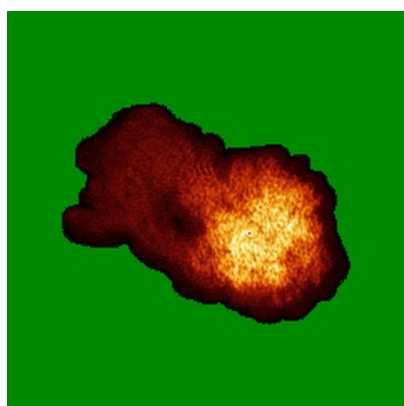


Z Index: 108

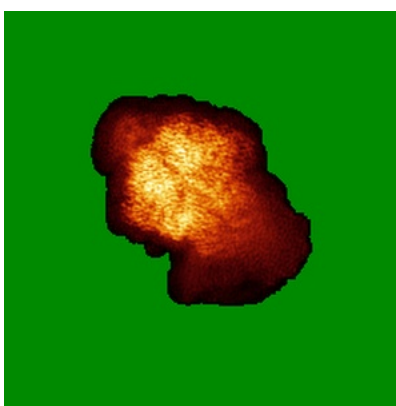
The images above show the largest variance slices of the map in three orthogonal directions.

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

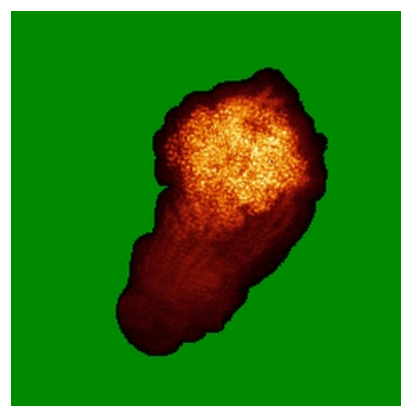
6.4.1 Primary map



X



Y

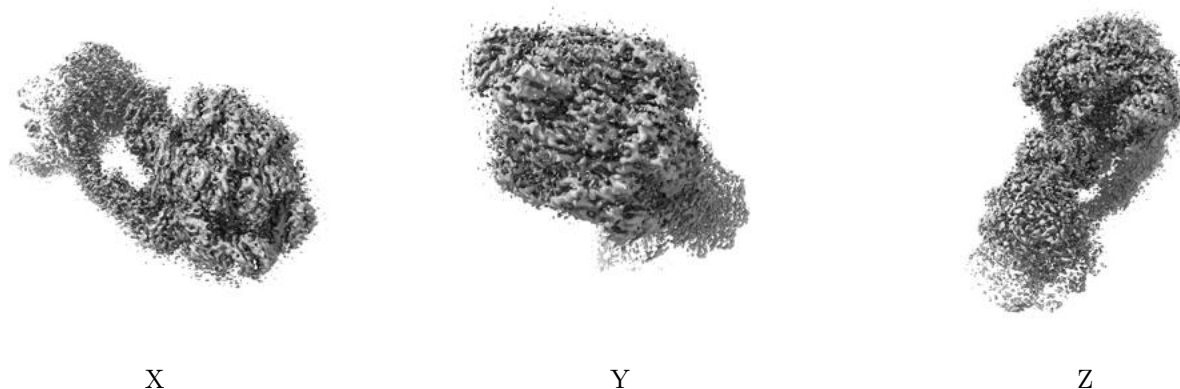


Z

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 0.0188. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

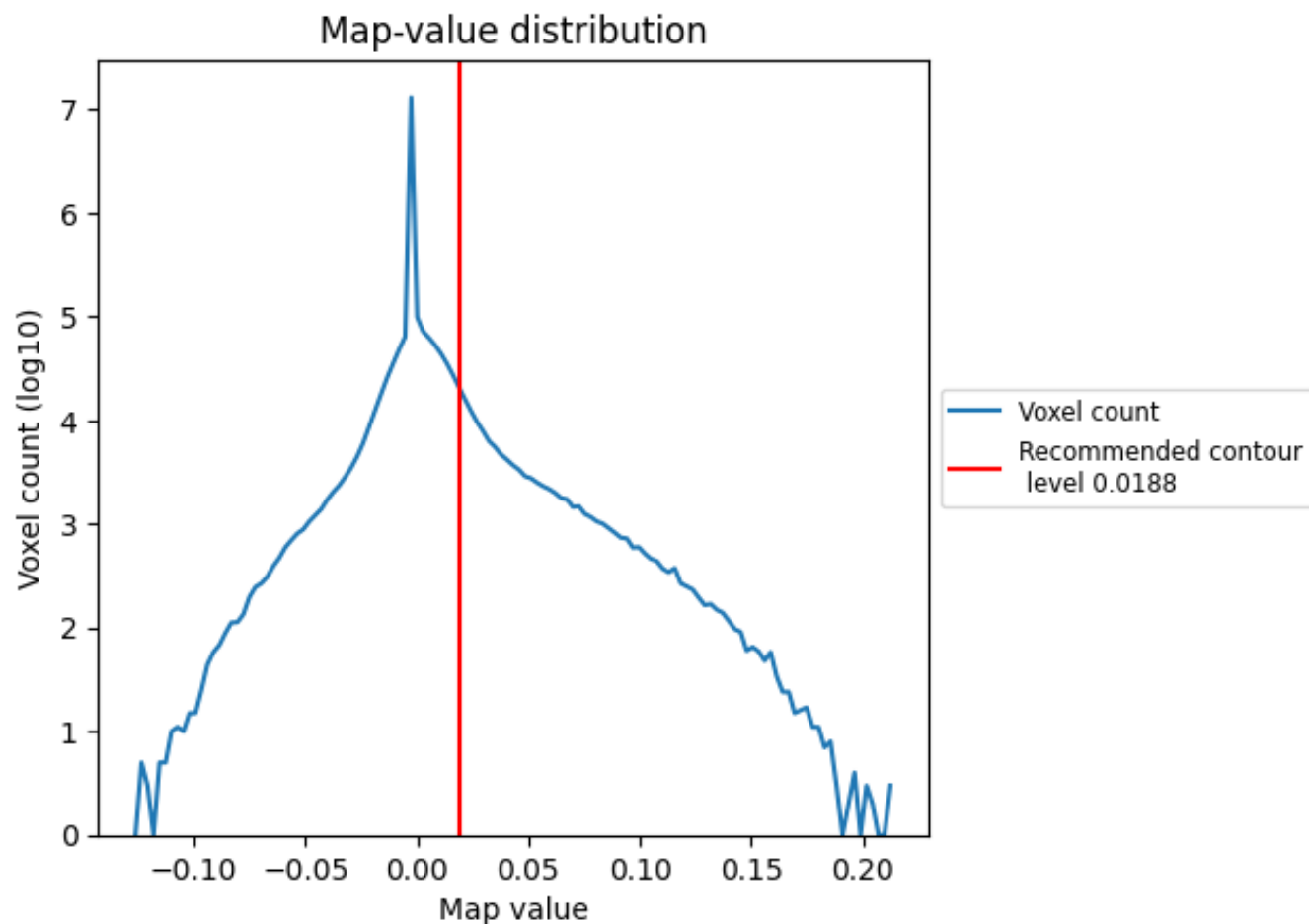
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

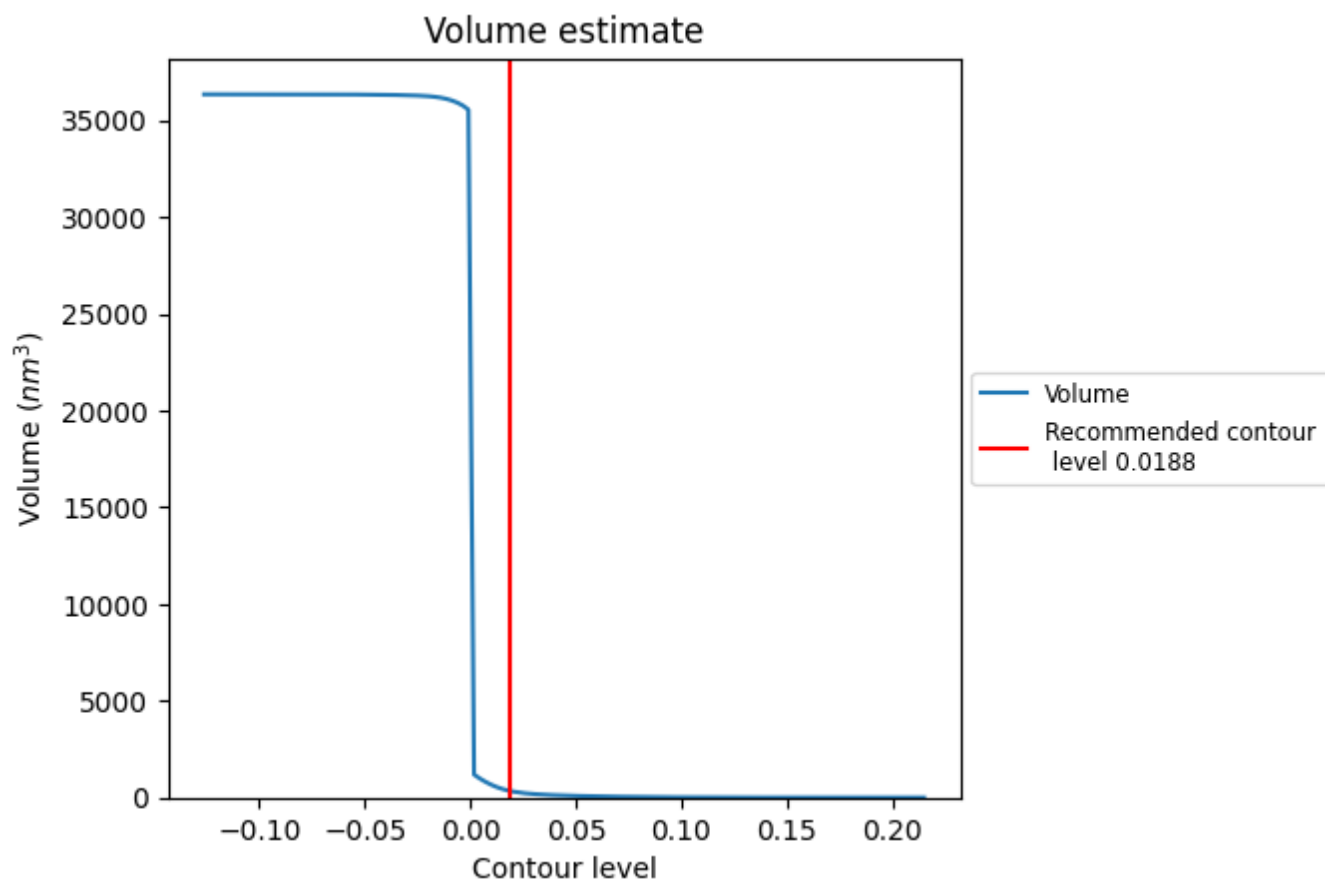
This section contains the results of statistical analysis of the map.

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.

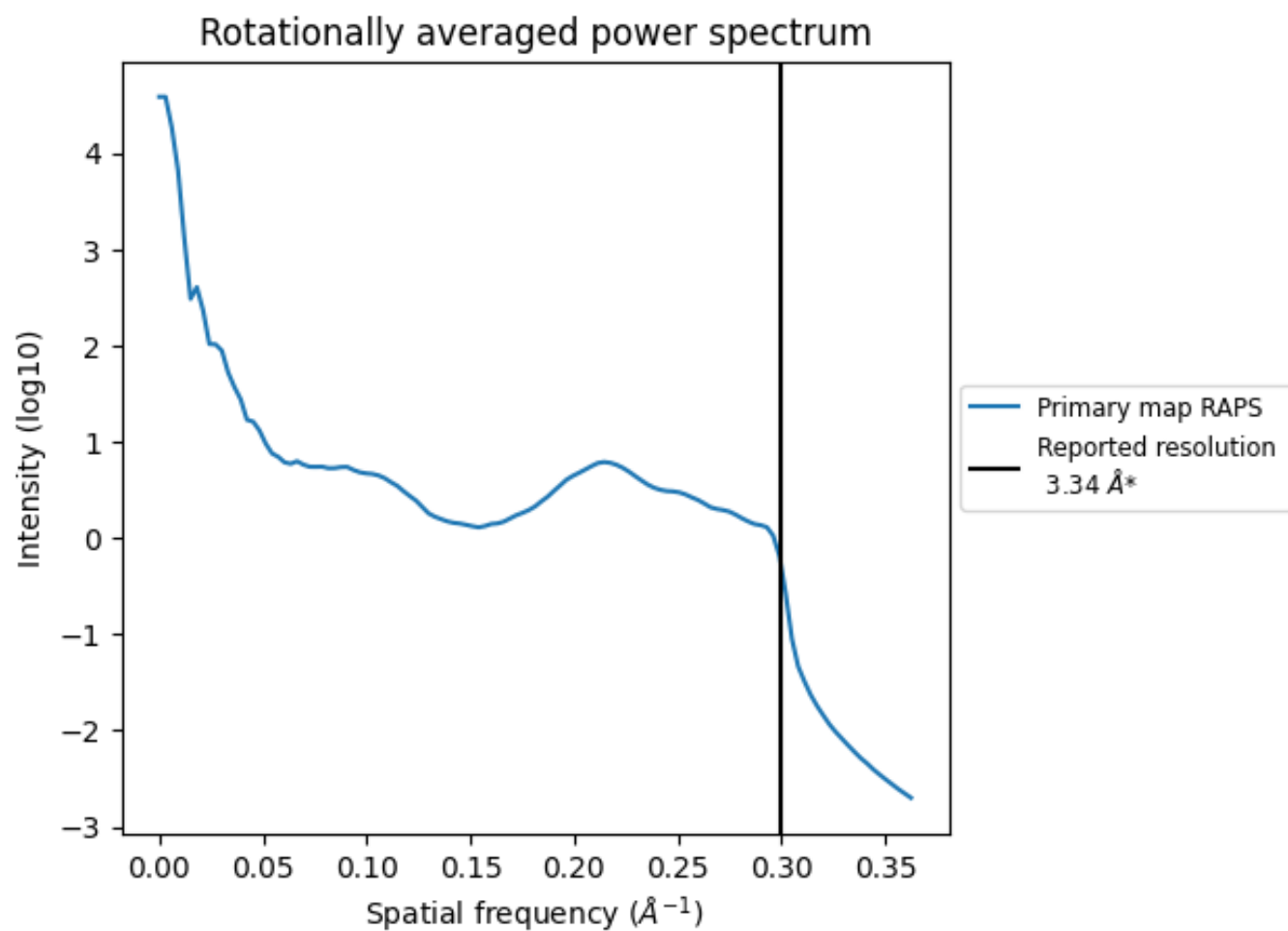
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 343 nm^3 ; this corresponds to an approximate mass of 310 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 0.299 Å⁻¹

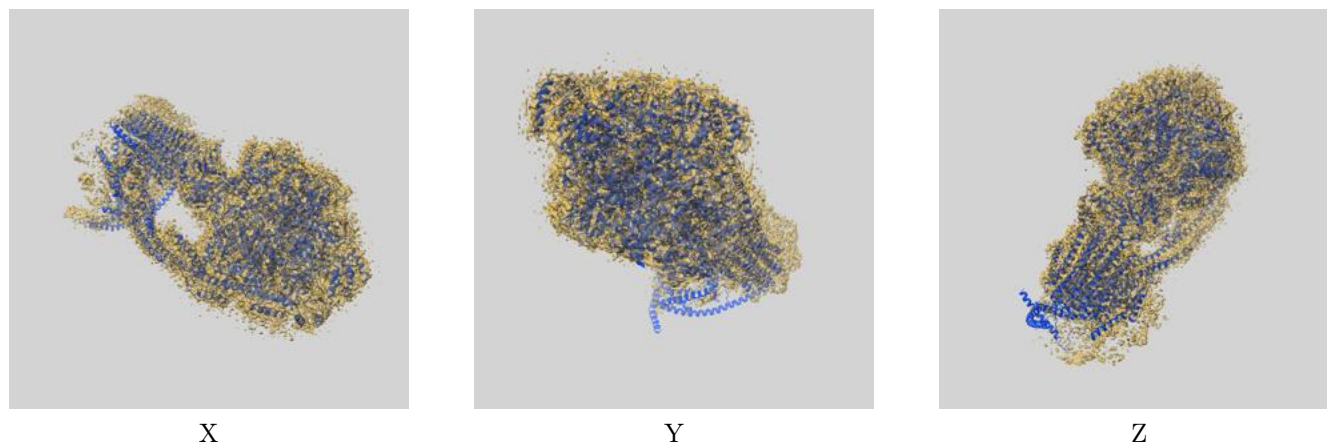
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

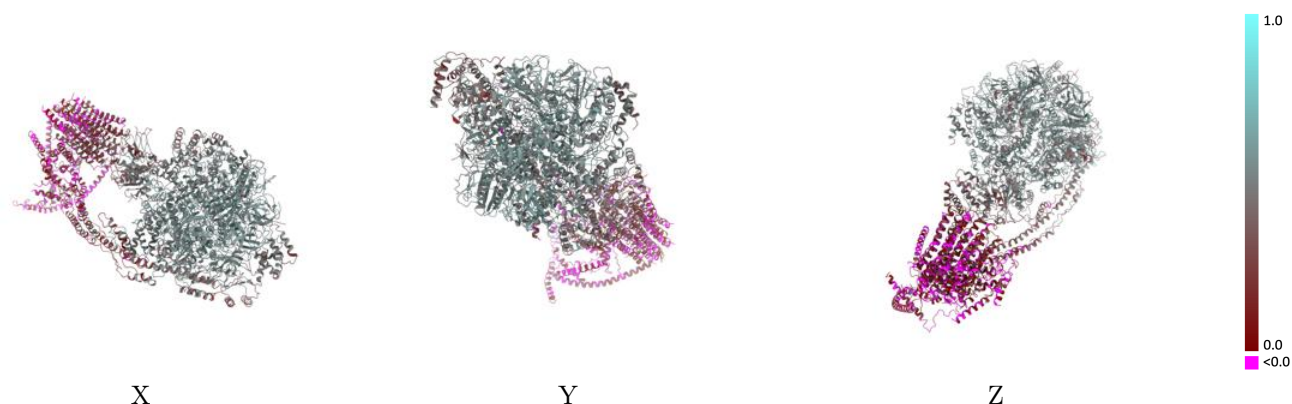
This section contains information regarding the fit between EMDB map EMD-0677 and PDB model 6J5I. Per-residue inclusion information can be found in section [3](#) on page [10](#).

9.1 Map-model overlay [i](#)



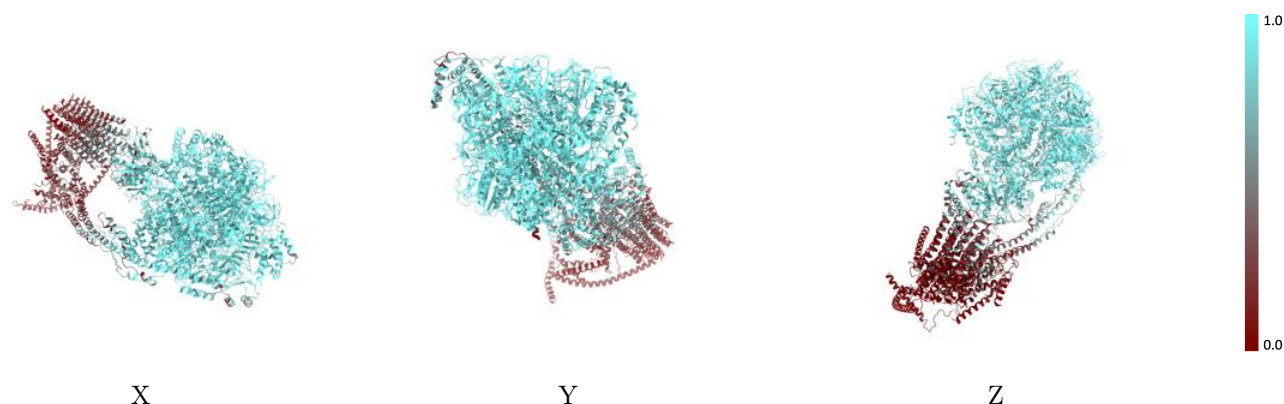
The images above show the 3D surface view of the map at the recommended contour level 0.0188 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



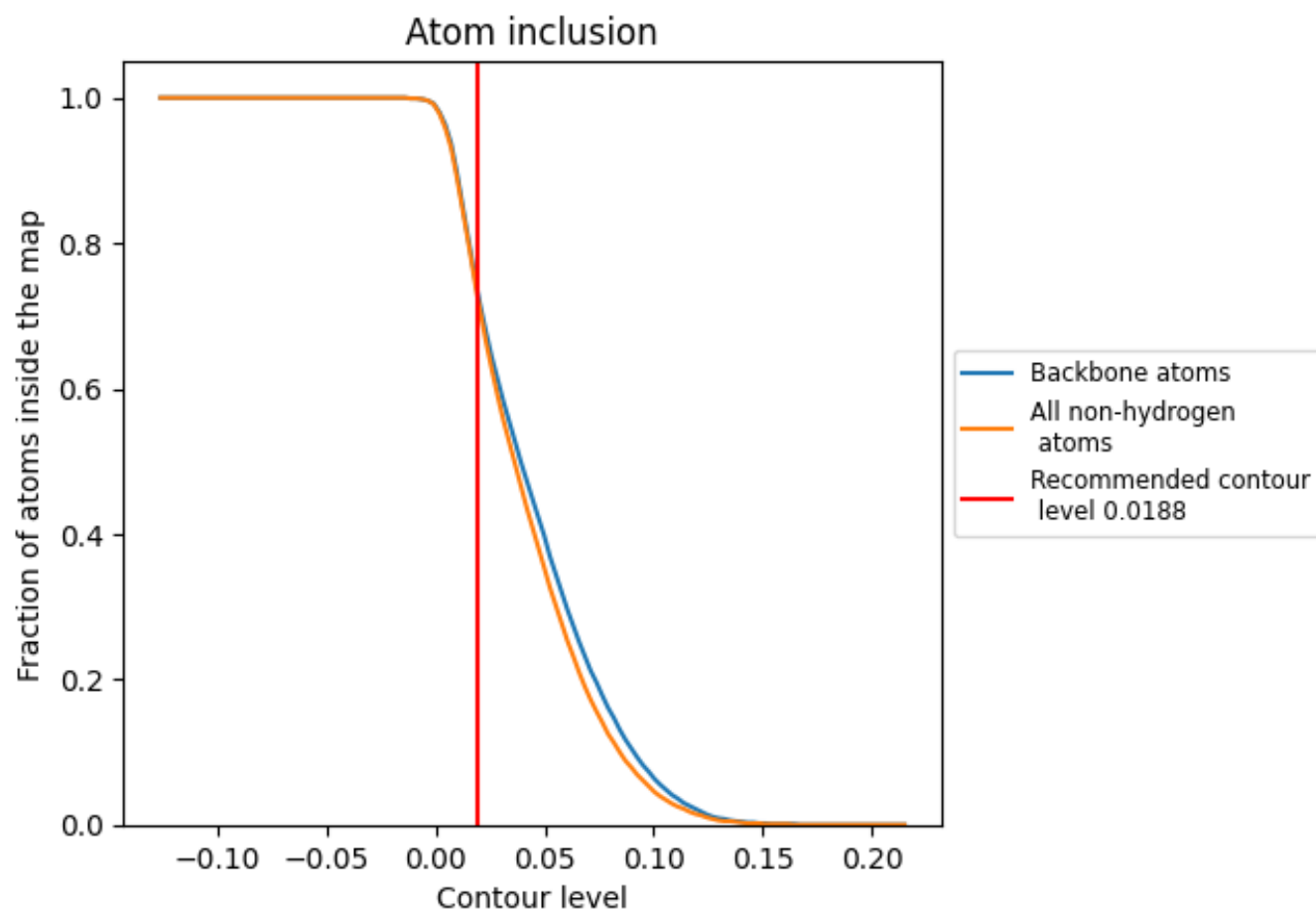
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0188).

9.4 Atom inclusion [i](#)



At the recommended contour level, 74% of all backbone atoms, 73% 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 (0.0188) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7320	 0.4080
8	 0.2180	 0.1430
A	 0.9370	 0.5380
B	 0.9380	 0.5220
C	 0.9370	 0.5230
D	 0.9530	 0.5380
E	 0.9300	 0.5150
F	 0.9590	 0.5470
G	 0.8610	 0.4620
H	 0.6720	 0.3810
I	 0.7440	 0.4000
J	 0.7690	 0.4680
K	 0.2380	 0.1050
L	 0.2690	 0.1310
M	 0.3010	 0.1290
N	 0.2690	 0.1800
O	 0.1940	 0.1060
P	 0.2040	 0.0940
Q	 0.2690	 0.1150
R	 0.2750	 0.1060
S	 0.8460	 0.4440
a	 0.1600	 0.0670
b	 0.4440	 0.2440
c	 0.8170	 0.3610
d	 0.4660	 0.2470
e	 0.0160	 0.1140
f	 0.0680	 0.0650
g	 0.0410	 0.0160
i	 0.1440	 0.0810
k	 0.0280	 0.0070
u	 0.0240	 0.0440

