



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 10:04 AM UTC

PDB ID : 4YB7 / pdb_00004yb7
Title : Adenosine triphosphate phosphoribosyltransferase from *Campylobacter jejuni*
in complex with ATP
Authors : Mittelstaedt, G.; Moggre, G.-J.; Parker, E.J.
Deposited on : 2015-02-18
Resolution : 2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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/XrayValidationReportHelp>

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:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

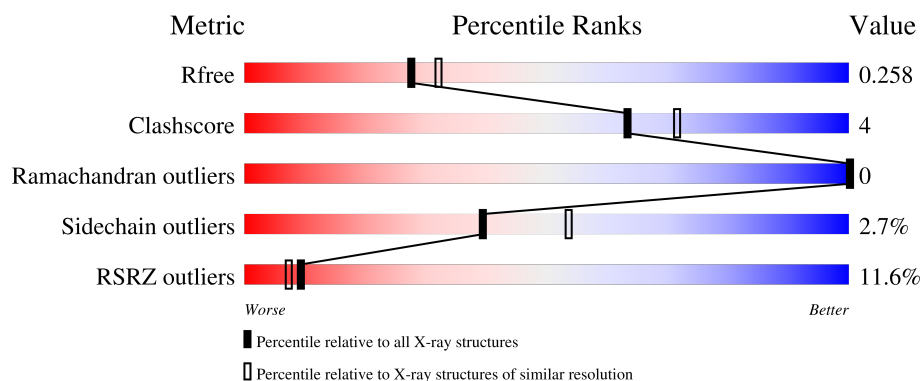
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	300	8% 91% 7% ..
1	B	300	10% 89% 8% ..
1	C	300	12% 86% 11% ..
1	D	300	11% 91% 6% ..
1	E	300	8% 91% 6% ..

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Mol	Chain	Length	Quality of chain
1	F	300	6% 90% 7% ..
1	G	300	13% 89% 7% ..
1	H	300	17% 89% 9% .
1	I	300	15% 88% 9% ..
1	J	300	6% 91% 7% ..
1	K	300	13% 88% 8% ..
1	L	300	18% 88% 8% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 26987 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	296	Total	C	N	O	S	0	0	0
			2228	1418	381	416	13			
1	C	294	Total	C	N	O	S	0	1	0
			2198	1399	376	409	14			
1	D	294	Total	C	N	O	S	0	0	0
			2168	1381	370	404	13			
1	E	295	Total	C	N	O	S	0	0	0
			2203	1395	377	418	13			
1	I	292	Total	C	N	O	S	0	0	0
			2169	1377	376	403	13			
1	K	294	Total	C	N	O	S	0	1	0
			2168	1380	369	405	14			
1	A	296	Total	C	N	O	S	0	0	0
			2202	1404	379	406	13			
1	F	293	Total	C	N	O	S	0	1	0
			2191	1389	375	413	14			
1	G	292	Total	C	N	O	S	0	0	0
			2149	1364	365	407	13			
1	H	296	Total	C	N	O	S	5	0	0
			2162	1372	372	405	13			
1	J	296	Total	C	N	O	S	0	0	0
			2218	1409	379	417	13			
1	L	293	Total	C	N	O	S	0	1	0
			2148	1360	368	406	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP Q5HSJ4
C	0	GLY	-	expression tag	UNP Q5HSJ4
D	0	GLY	-	expression tag	UNP Q5HSJ4
E	0	GLY	-	expression tag	UNP Q5HSJ4
I	0	GLY	-	expression tag	UNP Q5HSJ4

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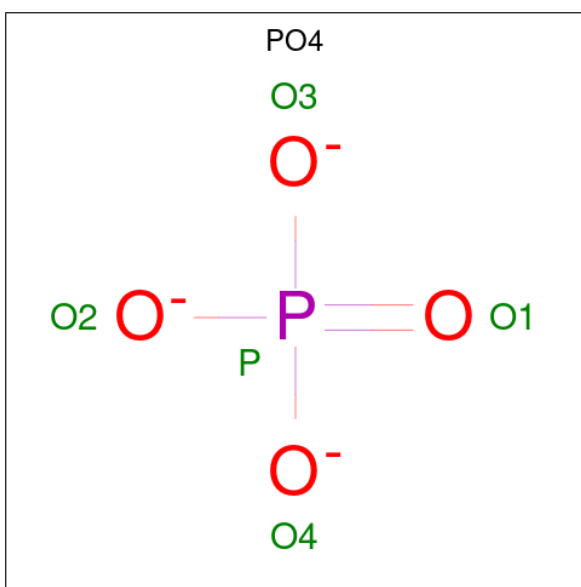
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Chain	Residue	Modelled	Actual	Comment	Reference
K	0	GLY	-	expression tag	UNP Q5HSJ4
A	0	GLY	-	expression tag	UNP Q5HSJ4
F	0	GLY	-	expression tag	UNP Q5HSJ4
G	0	GLY	-	expression tag	UNP Q5HSJ4
H	0	GLY	-	expression tag	UNP Q5HSJ4
J	0	GLY	-	expression tag	UNP Q5HSJ4
L	0	GLY	-	expression tag	UNP Q5HSJ4

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

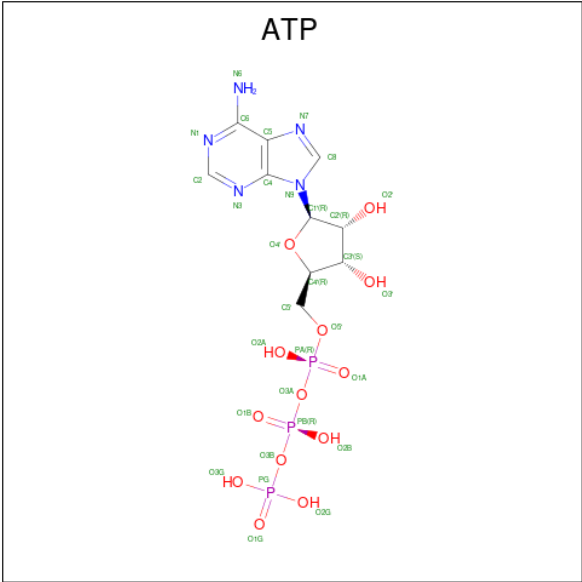
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		
2	D	1	Total	Mg	0	0
			1	1		
2	E	1	Total	Mg	0	0
			1	1		
2	I	1	Total	Mg	0	0
			1	1		
2	K	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		
2	F	1	Total	Mg	0	0
			1	1		
2	G	1	Total	Mg	0	0
			1	1		
2	H	1	Total	Mg	0	0
			1	1		
2	J	1	Total	Mg	0	0
			1	1		
2	L	1	Total	Mg	0	0
			1	1		

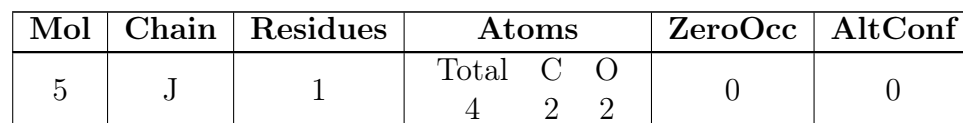
- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	E	1	Total	O	P	0	0
			5	4	1		
3	I	1	Total	O	P	0	0
			5	4	1		
3	J	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



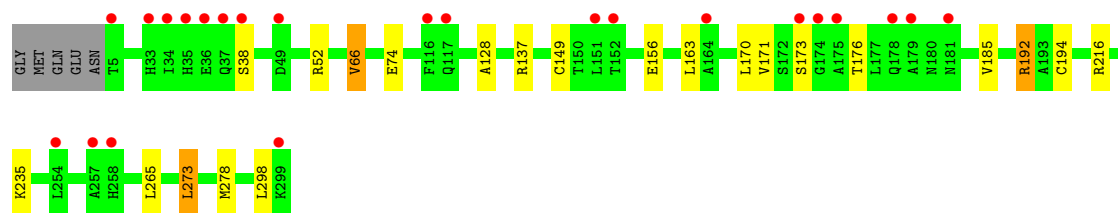


- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|------------------|---------|---------|
| 6 | B | 52 | Total O
52 52 | 0 | 0 |
| 6 | C | 44 | Total O
44 44 | 0 | 0 |
| 6 | D | 31 | Total O
31 31 | 0 | 0 |
| 6 | E | 19 | Total O
19 19 | 0 | 0 |
| 6 | I | 22 | Total O
22 22 | 0 | 0 |
| 6 | K | 15 | Total O
15 15 | 0 | 0 |
| 6 | A | 68 | Total O
68 68 | 0 | 0 |
| 6 | F | 29 | Total O
29 29 | 0 | 0 |
| 6 | G | 34 | Total O
34 34 | 0 | 0 |
| 6 | H | 12 | Total O
12 12 | 0 | 0 |
| 6 | J | 26 | Total O
26 26 | 0 | 0 |

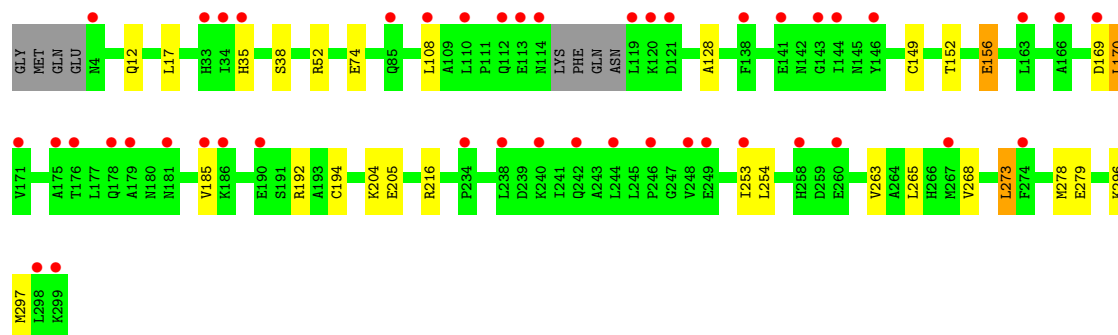
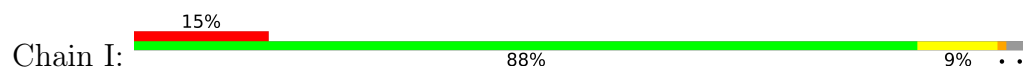


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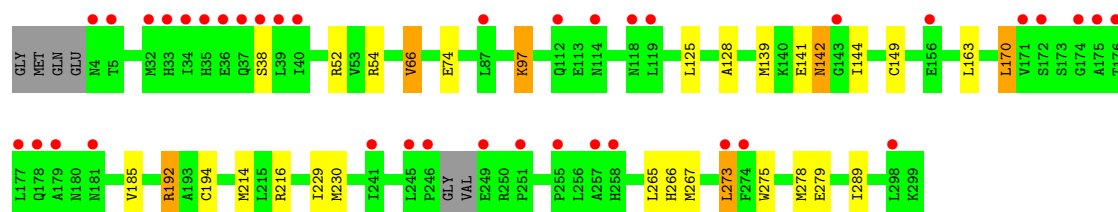
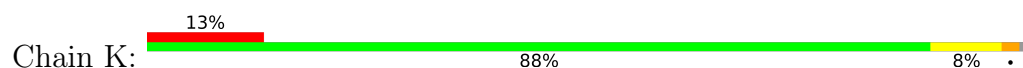
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	L	13	Total	O	0	0
			13	13		



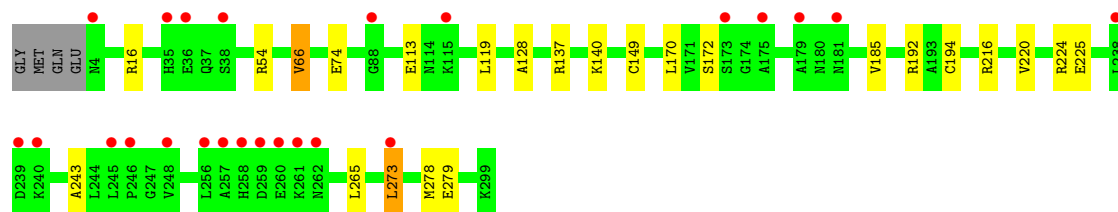
- Molecule 1: ATP phosphoribosyltransferase



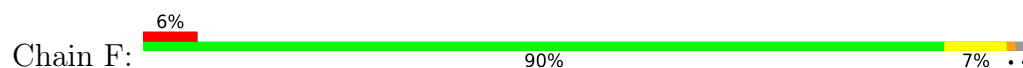
- Molecule 1: ATP phosphoribosyltransferase



- Molecule 1: ATP phosphoribosyltransferase

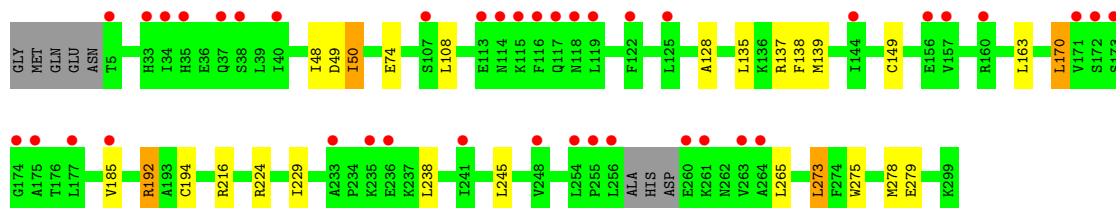
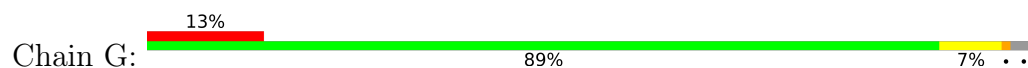


- Molecule 1: ATP phosphoribosyltransferase

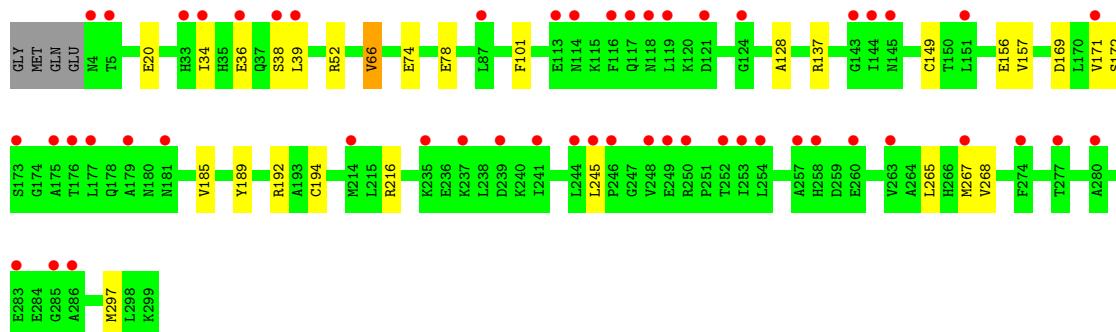
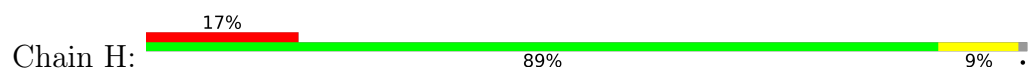




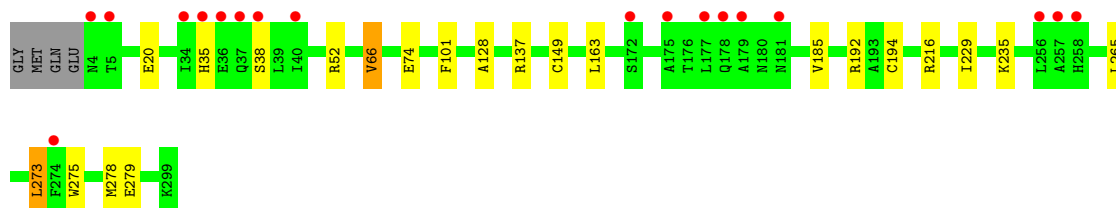
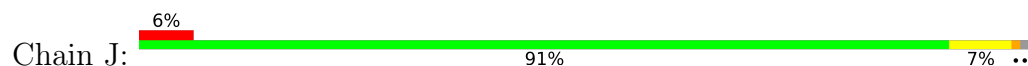
- Molecule 1: ATP phosphoribosyltransferase



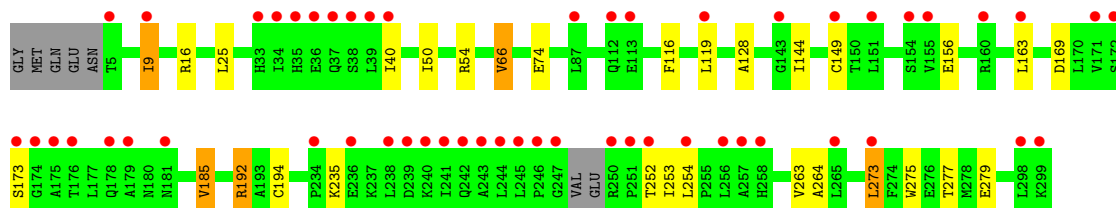
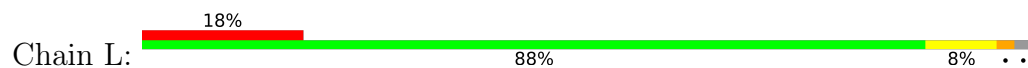
- Molecule 1: ATP phosphoribosyltransferase



- Molecule 1: ATP phosphoribosyltransferase



- Molecule 1: ATP phosphoribosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	91.67Å 91.83Å 154.90Å 101.11° 95.21° 118.14°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.1 (50.00-2.20) 98.1 (50.00-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.231 , 0.254 0.235 , 0.258	Depositor DCC
R_{free} test set	10624 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.053 for k,h,-h-k-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	26987	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ATP, PO4, ACY, 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	0.68	0/2229	0.85	3/3020 (0.1%)
1	B	0.66	0/2257	0.82	4/3055 (0.1%)
1	C	0.64	0/2225	0.86	5/3011 (0.2%)
1	D	0.62	0/2195	0.81	2/2977 (0.1%)
1	E	0.58	0/2231	0.86	7/3026 (0.2%)
1	F	0.61	0/2216	0.80	3/2997 (0.1%)
1	G	0.60	0/2174	0.85	6/2950 (0.2%)
1	H	0.57	1/2190 (0.0%)	0.84	8/2974 (0.3%)
1	I	0.62	0/2197	0.80	2/2976 (0.1%)
1	J	0.60	0/2247	0.81	4/3042 (0.1%)
1	K	0.60	0/2194	0.83	5/2975 (0.2%)
1	L	0.60	0/2174	0.83	3/2952 (0.1%)
All	All	0.61	1/26529 (0.0%)	0.83	52/35955 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	34	ILE	C-N	-5.56	1.26	1.33

The worst 5 of 52 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	192	ARG	NE-CZ-NH2	8.94	127.25	119.20
1	F	192	ARG	NE-CZ-NH2	8.93	127.23	119.20
1	C	192	ARG	NE-CZ-NH2	8.90	127.21	119.20
1	L	192	ARG	NE-CZ-NH2	8.73	127.06	119.20
1	E	192	ARG	NE-CZ-NH2	8.71	127.04	119.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2202	0	2219	16	0
1	B	2228	0	2247	24	0
1	C	2198	0	2200	29	0
1	D	2168	0	2142	27	0
1	E	2203	0	2185	12	1
1	F	2191	0	2190	15	1
1	G	2149	0	2110	25	0
1	H	2162	0	2109	18	0
1	I	2169	0	2159	22	0
1	J	2218	0	2210	22	1
1	K	2168	0	2131	16	1
1	L	2148	0	2087	34	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
3	B	5	0	0	0	0
3	C	10	0	0	0	0
3	E	5	0	0	0	0
3	I	5	0	0	0	0
3	J	5	0	0	0	0
4	A	31	0	12	2	0
4	B	31	0	12	2	0
4	C	31	0	12	0	0
4	D	31	0	12	2	0
4	E	31	0	12	0	0
4	F	31	0	12	0	0
4	G	31	0	12	1	0
4	H	31	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	31	0	12	1	0
4	J	31	0	12	2	0
4	K	31	0	12	1	0
4	L	31	0	12	2	0
5	J	4	0	3	0	0
6	A	68	0	0	2	0
6	B	52	0	0	1	0
6	C	44	0	0	2	0
6	D	31	0	0	3	0
6	E	19	0	0	0	0
6	F	29	0	0	1	0
6	G	34	0	0	1	0
6	H	12	0	0	0	0
6	I	22	0	0	0	0
6	J	26	0	0	2	0
6	K	15	0	0	0	0
6	L	13	0	0	0	0
All	All	26987	0	26136	212	2

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 4.

The worst 5 of 212 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:139:MET:O	1:K:142:ASN:O	1.68	1.12
1:E:156:GLU:OE1	1:E:176:THR:HG21	1.50	1.11
1:B:245:LEU:HD21	1:B:273:LEU:HD11	1.28	1.10
1:D:138:PHE:CD2	1:D:139:MET:CE	2.38	1.07
1:D:138:PHE:HD2	1:D:139:MET:HE3	1.19	1.04

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:118:ASN:ND2	1:J:235:LYS:O[1_445]	1.72	0.48
1:E:235:LYS:NZ	1:K:141:GLU:OE2[1_655]	2.09	0.11

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	294/300 (98%)	291 (99%)	3 (1%)	0	100	100
1	B	294/300 (98%)	291 (99%)	3 (1%)	0	100	100
1	C	291/300 (97%)	288 (99%)	3 (1%)	0	100	100
1	D	290/300 (97%)	287 (99%)	3 (1%)	0	100	100
1	E	293/300 (98%)	290 (99%)	3 (1%)	0	100	100
1	F	290/300 (97%)	287 (99%)	3 (1%)	0	100	100
1	G	288/300 (96%)	285 (99%)	3 (1%)	0	100	100
1	H	294/300 (98%)	291 (99%)	3 (1%)	0	100	100
1	I	288/300 (96%)	284 (99%)	4 (1%)	0	100	100
1	J	294/300 (98%)	291 (99%)	3 (1%)	0	100	100
1	K	291/300 (97%)	288 (99%)	3 (1%)	0	100	100
1	L	290/300 (97%)	287 (99%)	3 (1%)	0	100	100
All	All	3497/3600 (97%)	3460 (99%)	37 (1%)	0	100	100

There are no Ramachandran outliers to report.

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 X-ray entries followed by that with respect to entries of similar resolution.

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	227/263 (86%)	220 (97%)	7 (3%)	35	48
1	B	234/263 (89%)	228 (97%)	6 (3%)	40	55

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	228/263 (87%)	221 (97%)	7 (3%)	35	48
1	D	219/263 (83%)	216 (99%)	3 (1%)	59	75
1	E	229/263 (87%)	222 (97%)	7 (3%)	35	48
1	F	228/263 (87%)	222 (97%)	6 (3%)	40	55
1	G	218/263 (83%)	212 (97%)	6 (3%)	38	52
1	H	215/263 (82%)	212 (99%)	3 (1%)	59	75
1	I	223/263 (85%)	216 (97%)	7 (3%)	35	48
1	J	230/263 (88%)	225 (98%)	5 (2%)	45	61
1	K	218/263 (83%)	209 (96%)	9 (4%)	27	37
1	L	215/263 (82%)	209 (97%)	6 (3%)	38	52
All	All	2684/3156 (85%)	2612 (97%)	72 (3%)	39	53

5 of 72 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	273	LEU
1	L	273	LEU
1	H	185	VAL
1	J	273	LEU
1	I	156	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 36 such sidechains are listed below:

Mol	Chain	Res	Type
1	J	114	ASN
1	L	242	GLN
1	J	133	GLN
1	L	33	HIS
1	I	12	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 31 ligands modelled in this entry, 12 are monoatomic - leaving 19 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$
4	ATP	A	302	2	32,33,33	1.49	5 (15%)	48,52,52	1.85	9 (18%)
3	PO4	C	303	-	4,4,4	0.91	0	6,6,6	0.38	0
4	ATP	L	302	2	32,33,33	1.54	5 (15%)	48,52,52	2.12	13 (27%)
3	PO4	J	302	-	4,4,4	0.87	0	6,6,6	0.52	0
4	ATP	D	302	2	32,33,33	1.49	4 (12%)	48,52,52	1.89	11 (22%)
4	ATP	I	303	2	32,33,33	1.40	6 (18%)	48,52,52	1.97	12 (25%)
4	ATP	G	302	2	32,33,33	1.51	4 (12%)	48,52,52	2.02	12 (25%)
4	ATP	J	303	2	32,33,33	1.29	4 (12%)	48,52,52	1.88	13 (27%)
4	ATP	E	303	2	32,33,33	1.45	5 (15%)	48,52,52	1.85	10 (20%)
3	PO4	E	302	-	4,4,4	0.80	0	6,6,6	0.62	0
4	ATP	B	303	2	32,33,33	1.48	6 (18%)	48,52,52	1.84	11 (22%)
5	ACY	J	304	-	3,3,3	0.85	0	3,3,3	0.71	0
3	PO4	B	302	-	4,4,4	0.87	0	6,6,6	0.57	0
4	ATP	C	304	2	32,33,33	1.49	4 (12%)	48,52,52	2.07	12 (25%)
4	ATP	K	302	2	32,33,33	1.46	4 (12%)	48,52,52	1.93	9 (18%)
3	PO4	C	302	-	4,4,4	1.01	0	6,6,6	0.52	0
4	ATP	H	302	2	32,33,33	1.55	7 (21%)	48,52,52	1.72	10 (20%)
4	ATP	F	302	2	32,33,33	1.53	6 (18%)	48,52,52	1.94	12 (25%)
3	PO4	I	302	-	4,4,4	0.90	0	6,6,6	0.47	0

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
4	ATP	C	304	2	-	5/22/38/38	0/3/3/3
4	ATP	A	302	2	-	6/22/38/38	0/3/3/3
4	ATP	D	302	2	-	3/22/38/38	0/3/3/3
4	ATP	J	303	2	-	2/22/38/38	0/3/3/3
4	ATP	L	302	2	-	8/22/38/38	0/3/3/3
4	ATP	E	303	2	-	9/22/38/38	0/3/3/3
4	ATP	K	302	2	-	6/22/38/38	0/3/3/3
4	ATP	I	303	2	-	5/22/38/38	0/3/3/3
4	ATP	G	302	2	-	2/22/38/38	0/3/3/3
4	ATP	B	303	2	-	4/22/38/38	0/3/3/3
4	ATP	H	302	2	-	5/22/38/38	0/3/3/3
4	ATP	F	302	2	-	4/22/38/38	0/3/3/3

The worst 5 of 60 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	K	302	ATP	C5-C4	5.02	1.48	1.39
4	E	303	ATP	C5-C4	4.88	1.47	1.39
4	L	302	ATP	C5-C4	4.86	1.47	1.39
4	H	302	ATP	C5-C4	4.83	1.47	1.39
4	D	302	ATP	C5-C4	4.81	1.47	1.39

The worst 5 of 134 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	302	ATP	C5-C4-N3	-7.37	116.57	126.72
4	C	304	ATP	C5-C4-N3	-6.95	117.14	126.72
4	G	302	ATP	C5-C4-N3	-6.84	117.29	126.72
4	K	302	ATP	C5-C4-N3	-6.48	117.80	126.72
4	I	303	ATP	C5-C4-N3	-6.43	117.87	126.72

There are no chirality outliers.

5 of 59 torsion outliers are listed below:

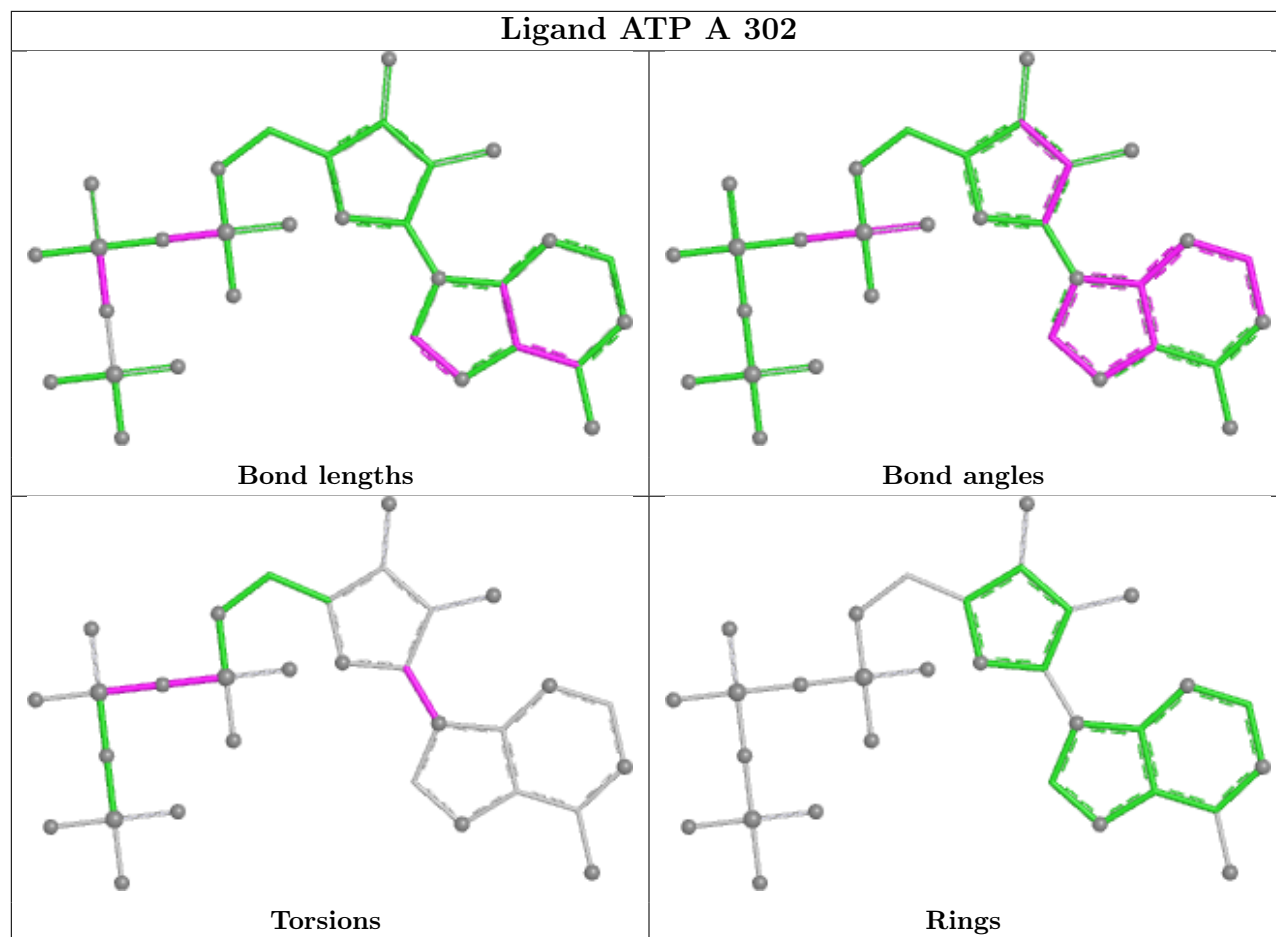
Mol	Chain	Res	Type	Atoms
4	E	303	ATP	C5'-O5'-PA-O1A
4	K	302	ATP	C5'-O5'-PA-O1A
4	K	302	ATP	O4'-C4'-C5'-O5'
4	K	302	ATP	C3'-C4'-C5'-O5'
4	H	302	ATP	C5'-O5'-PA-O3A

There are no ring outliers.

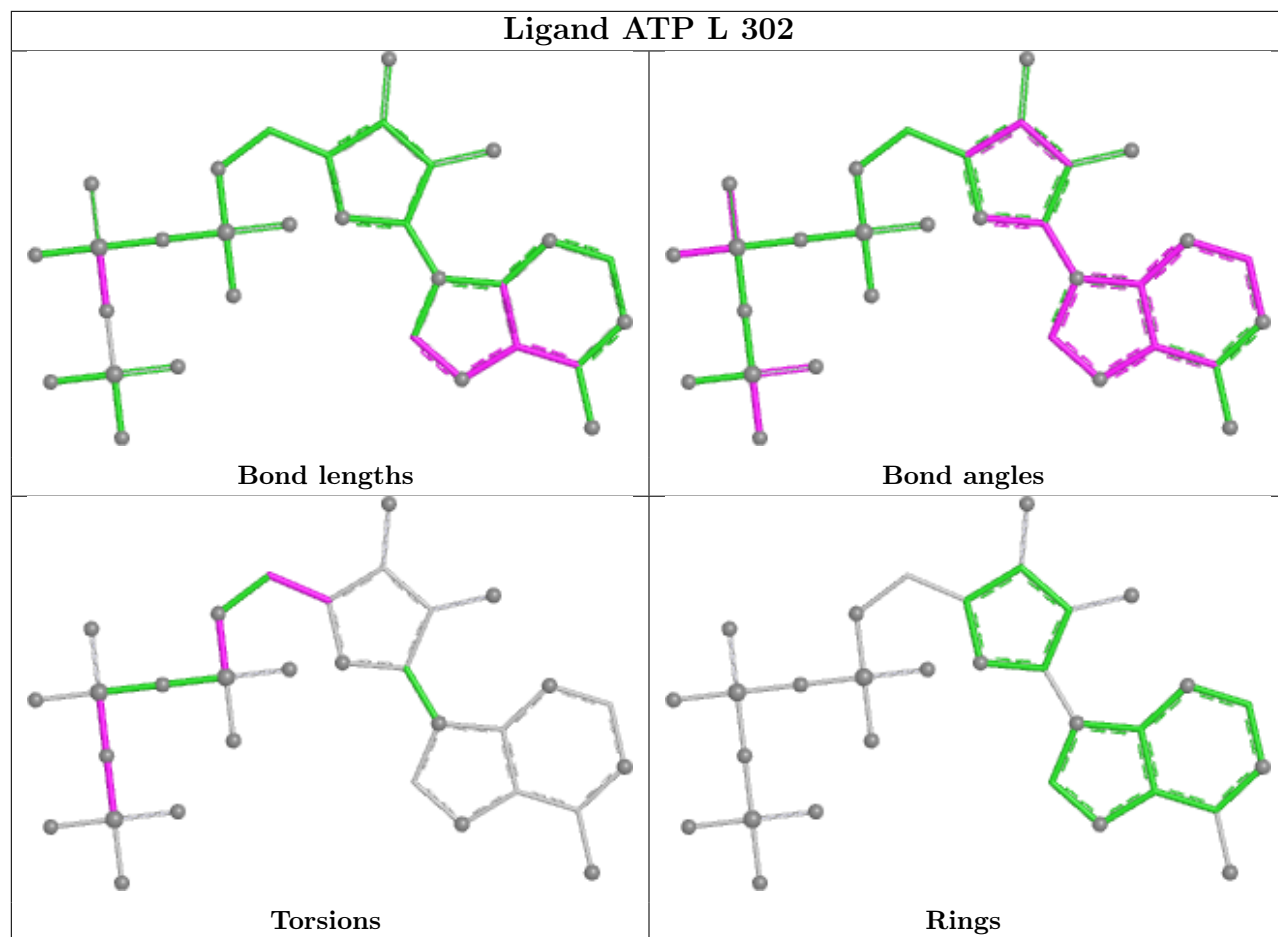
8 monomers are involved in 13 short contacts:

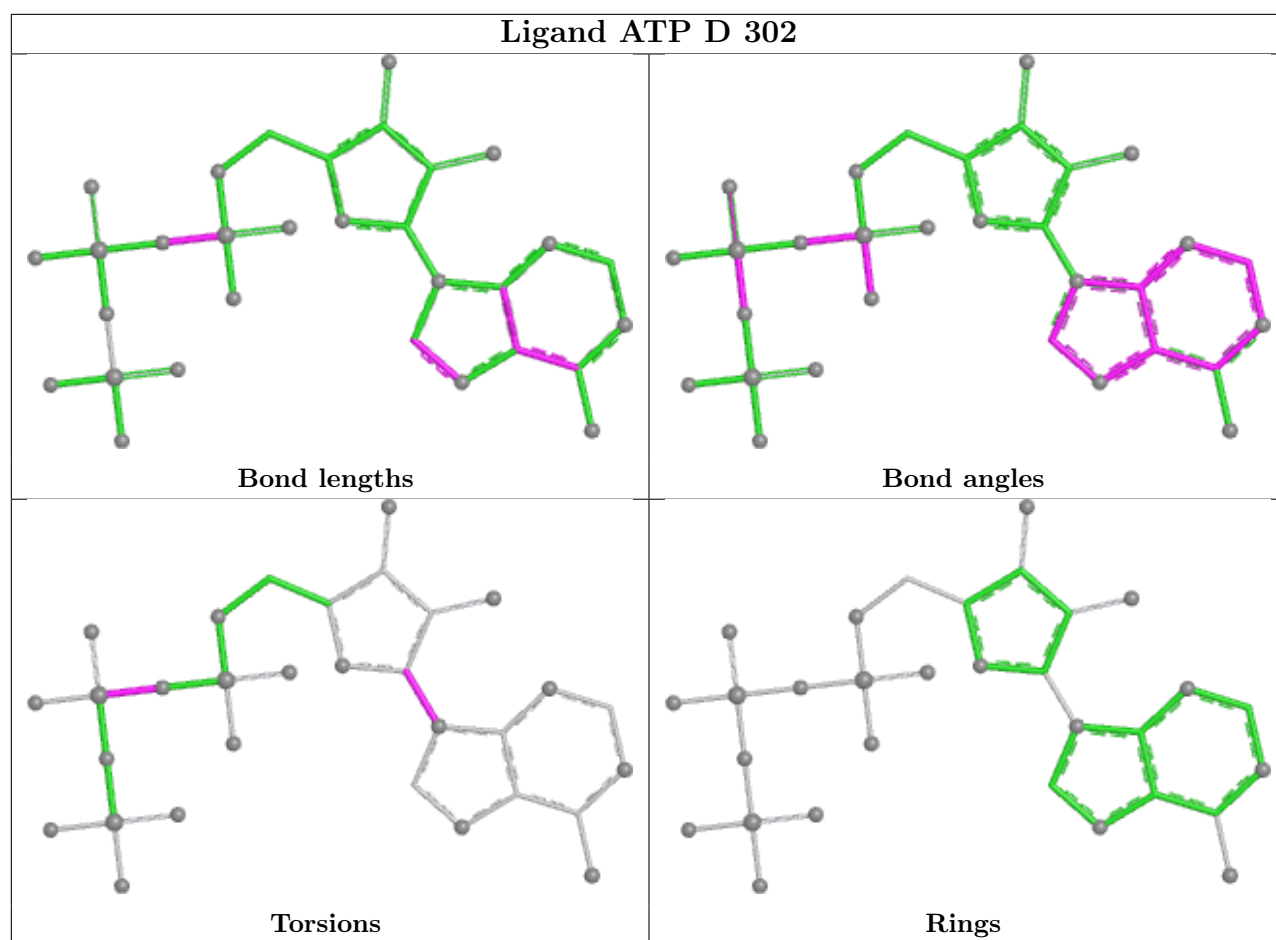
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	302	ATP	2	0
4	L	302	ATP	2	0
4	D	302	ATP	2	0
4	I	303	ATP	1	0
4	G	302	ATP	1	0
4	J	303	ATP	2	0
4	B	303	ATP	2	0
4	K	302	ATP	1	0

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.

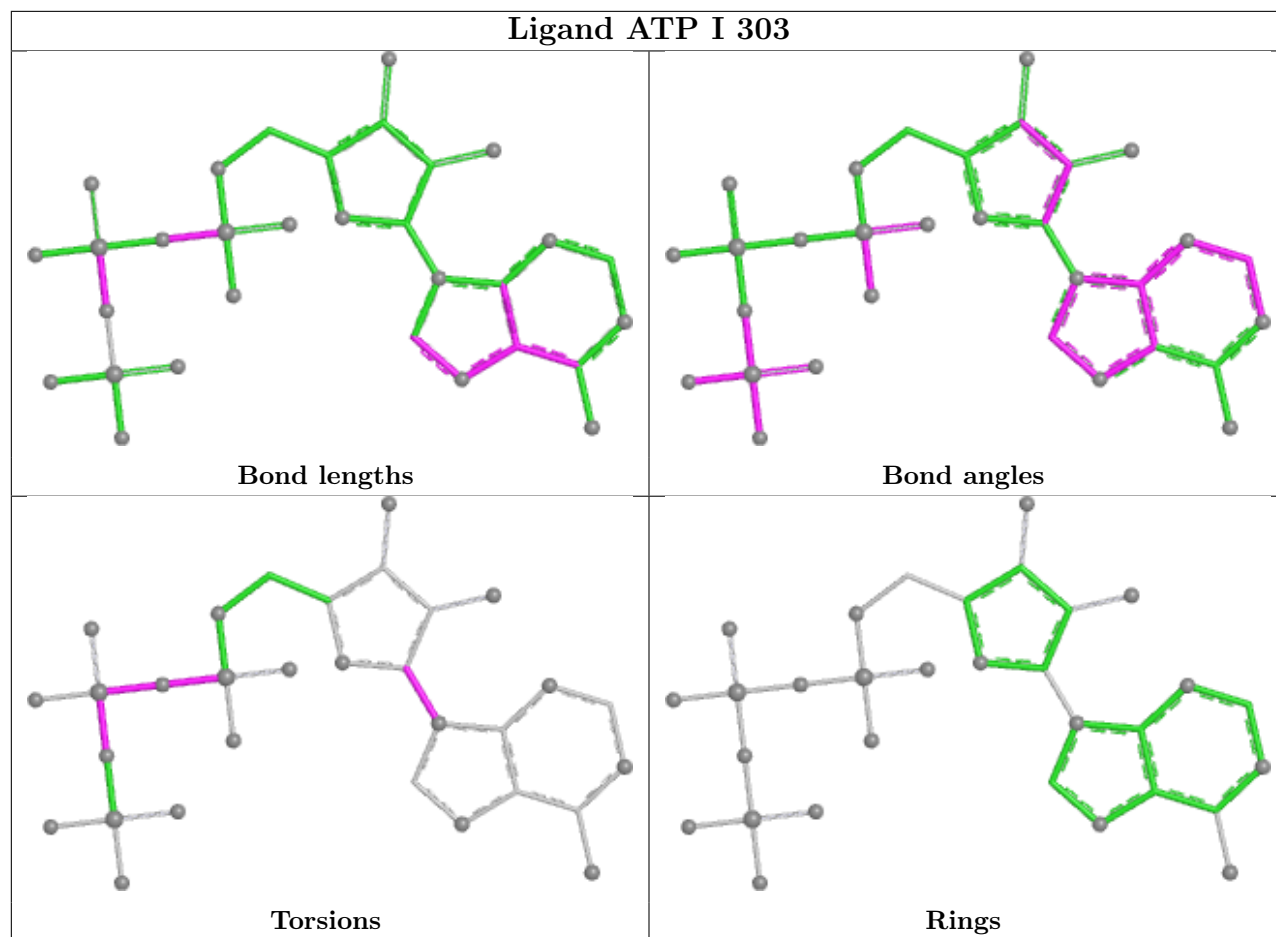


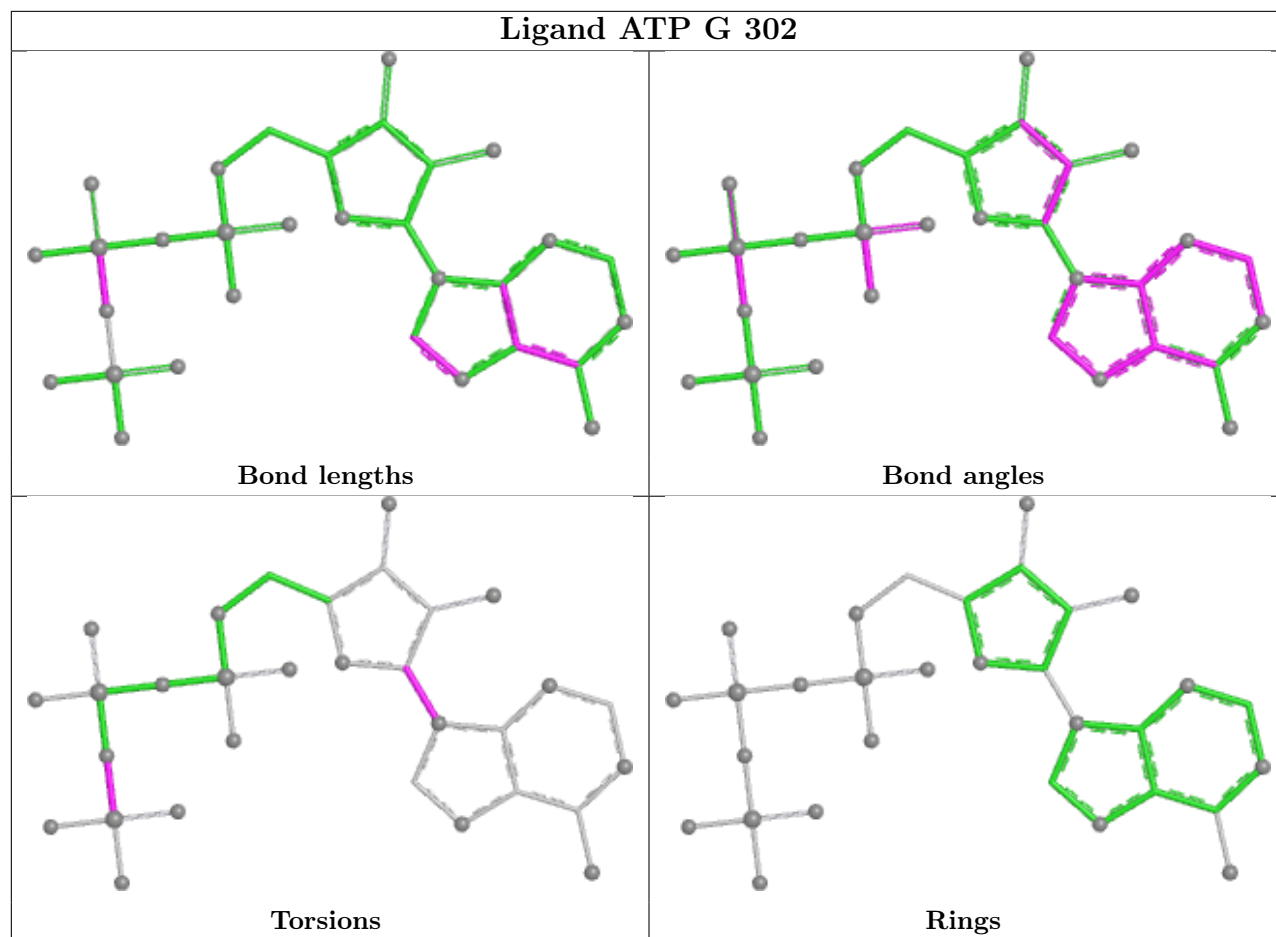
Ligand ATP L 302



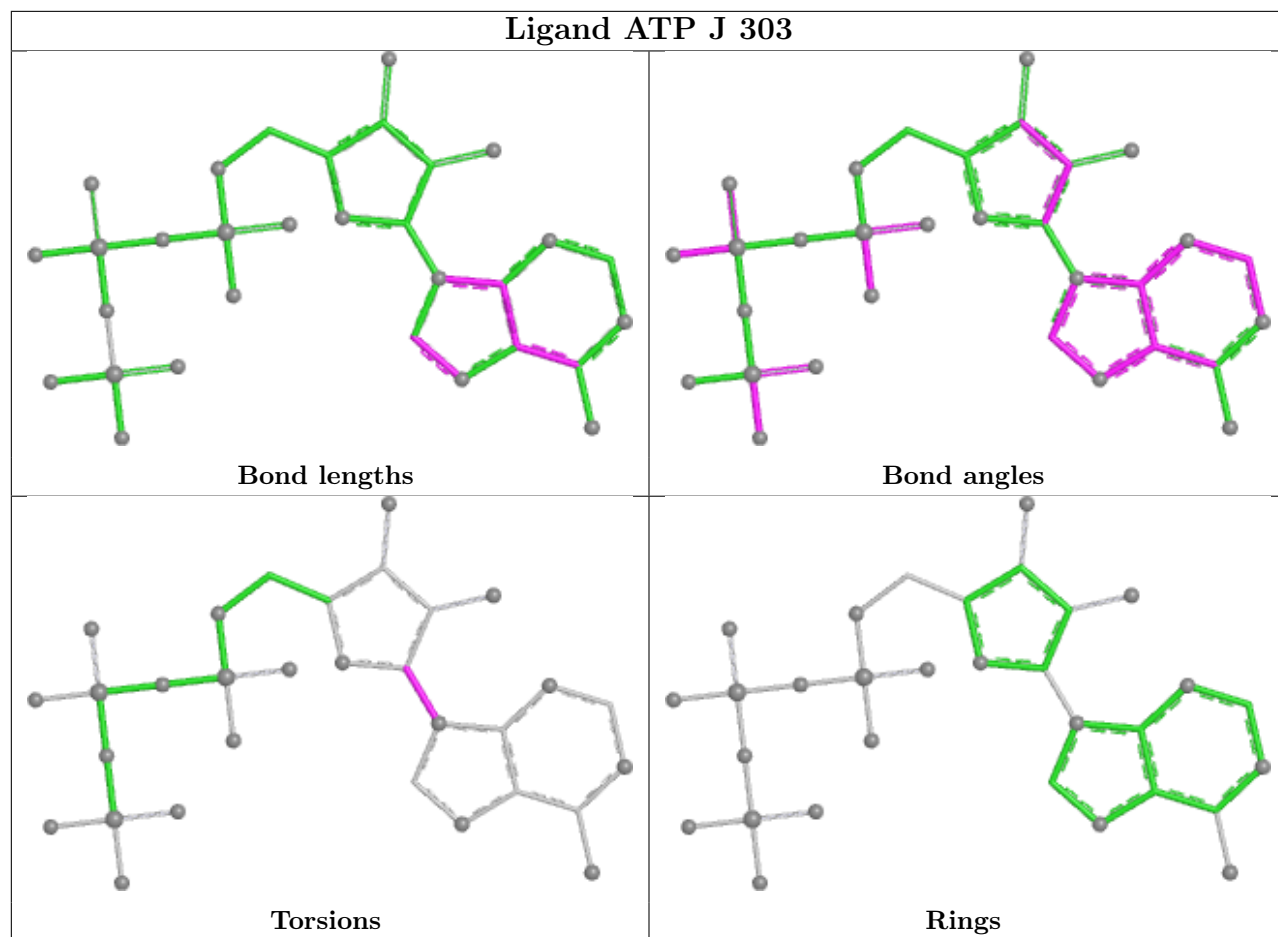


Ligand ATP I 303

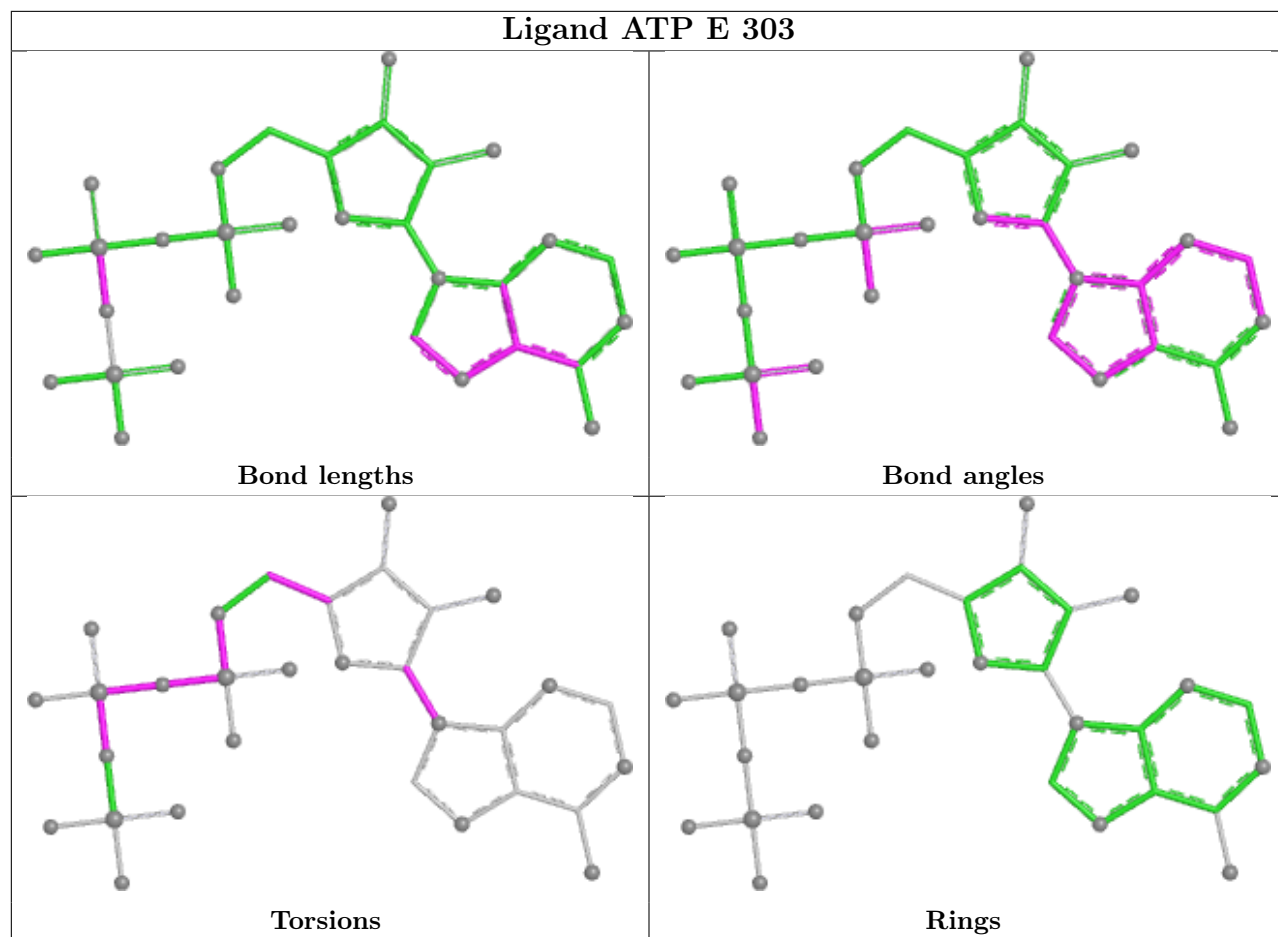


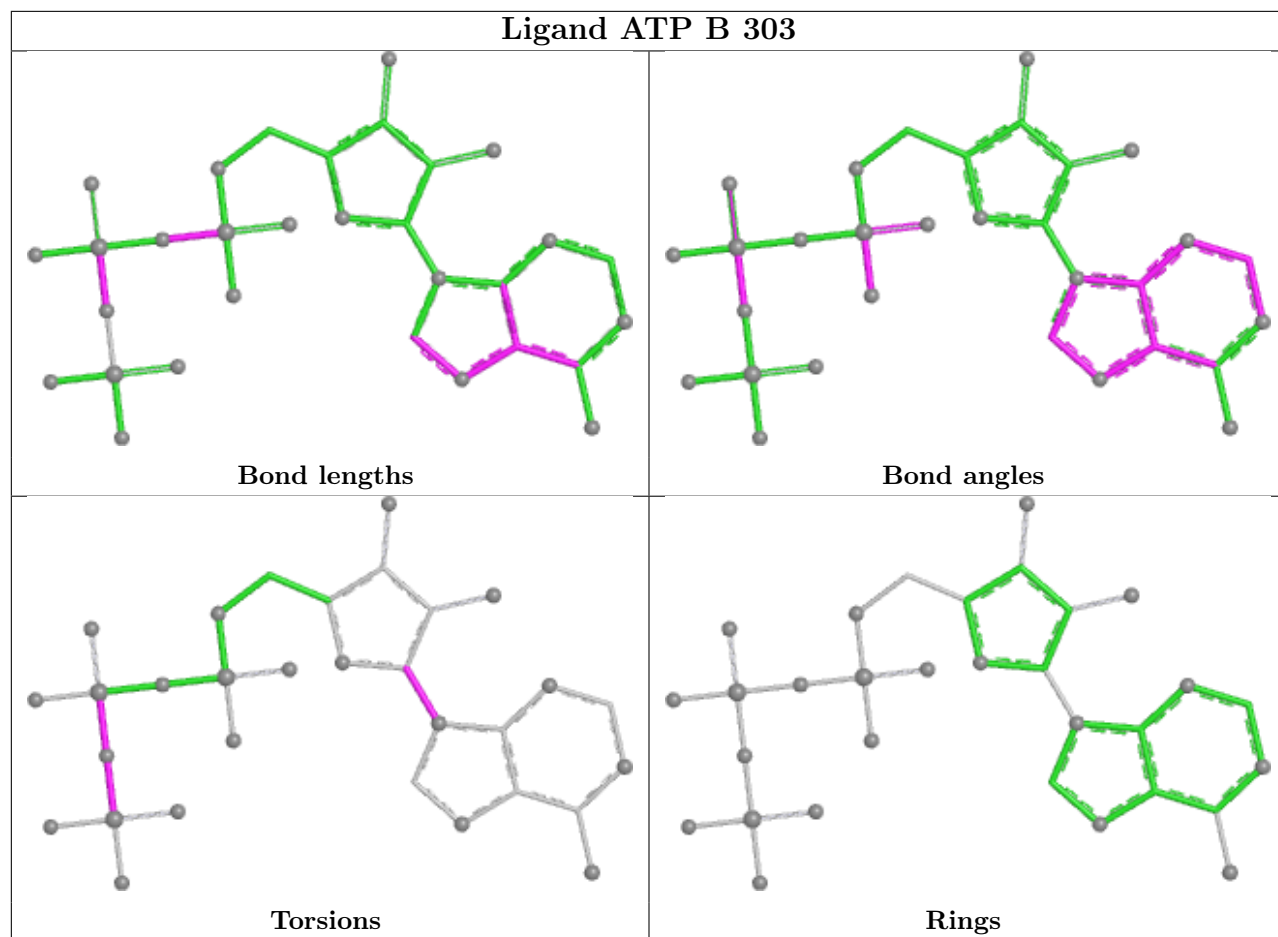


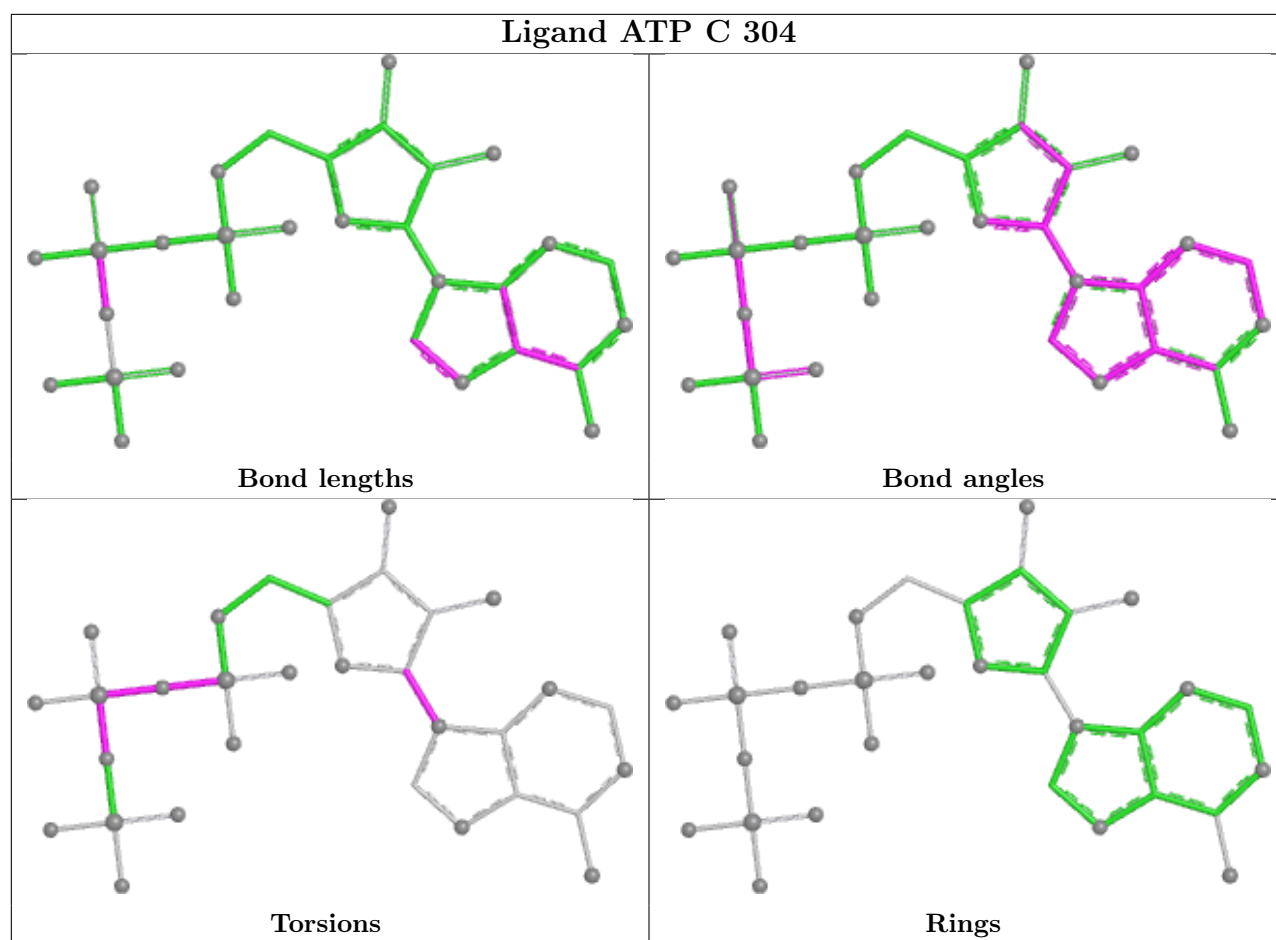
Ligand ATP J 303

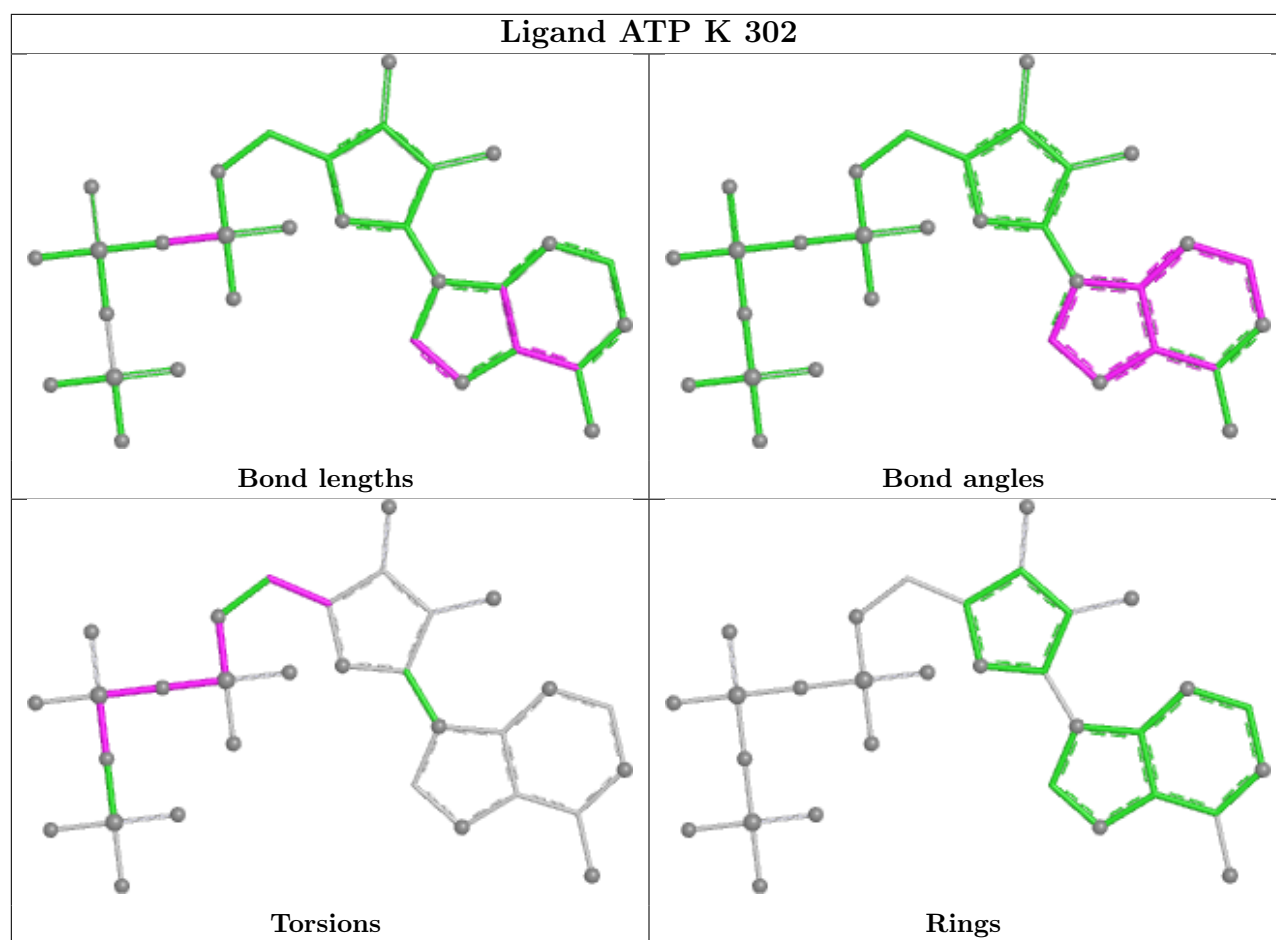


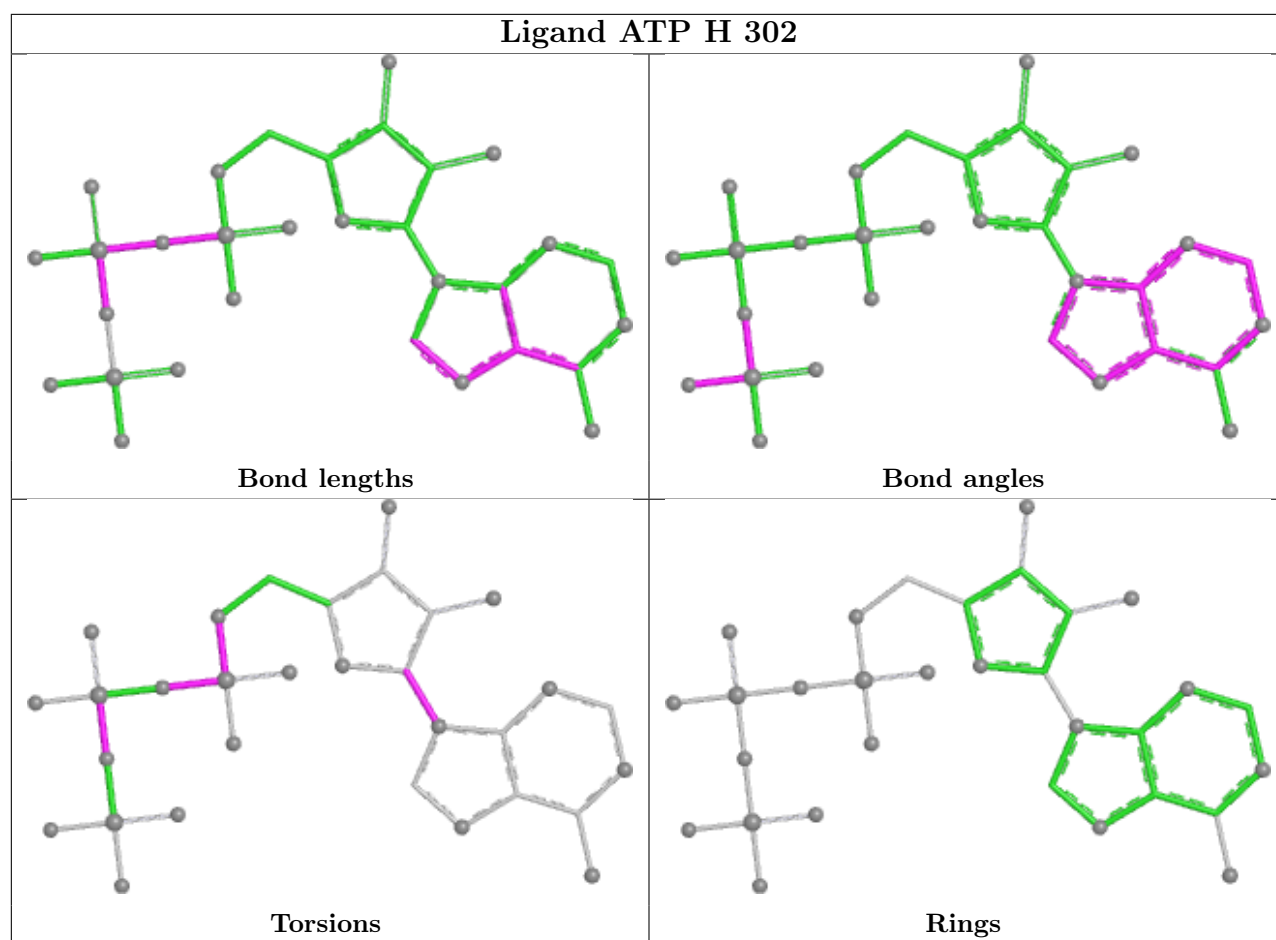
Ligand ATP E 303

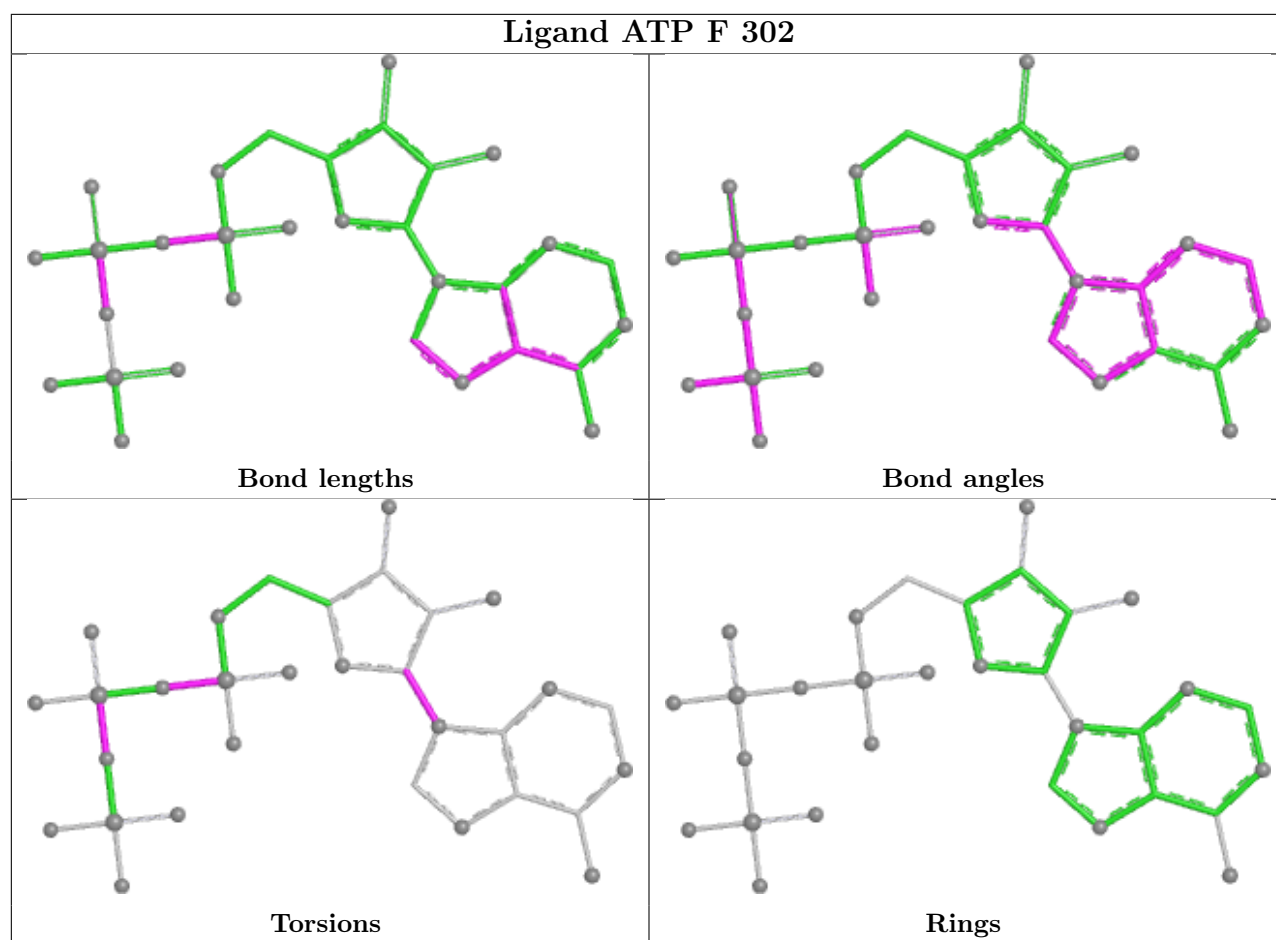












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	296/300 (98%)	0.45	24 (8%)	18 15	18, 35, 64, 92	0
1	B	296/300 (98%)	0.53	29 (9%)	13 10	18, 38, 72, 92	0
1	C	294/300 (98%)	0.64	36 (12%)	8 6	14, 39, 77, 94	1 (0%)
1	D	294/300 (98%)	0.78	34 (11%)	9 7	23, 43, 79, 99	0
1	E	295/300 (98%)	0.78	23 (7%)	19 16	23, 49, 78, 100	0
1	F	293/300 (97%)	0.56	19 (6%)	25 22	14, 42, 69, 85	1 (0%)
1	G	292/300 (97%)	0.82	40 (13%)	6 5	23, 44, 77, 99	0
1	H	295/300 (98%)	1.16	52 (17%)	4 3	33, 55, 85, 102	0
1	I	292/300 (97%)	0.91	45 (15%)	5 4	23, 48, 82, 92	0
1	J	296/300 (98%)	0.54	18 (6%)	27 24	23, 42, 69, 82	0
1	K	294/300 (98%)	1.06	38 (12%)	7 5	18, 52, 81, 102	1 (0%)
1	L	293/300 (97%)	1.15	53 (18%)	3 2	21, 52, 82, 91	1 (0%)
All	All	3530/3600 (98%)	0.78	411 (11%)	9 7	14, 45, 78, 102	4 (0%)

The worst 5 of 411 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	236	GLU	5.9
1	E	35	HIS	5.4
1	K	175	ALA	5.2
1	D	118	ASN	5.2
1	D	175	ALA	5.1

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

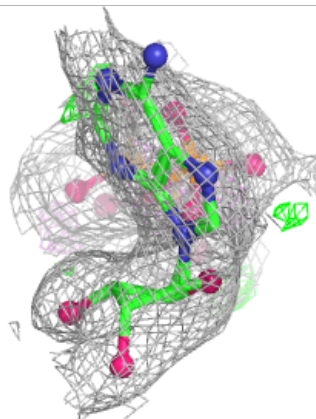
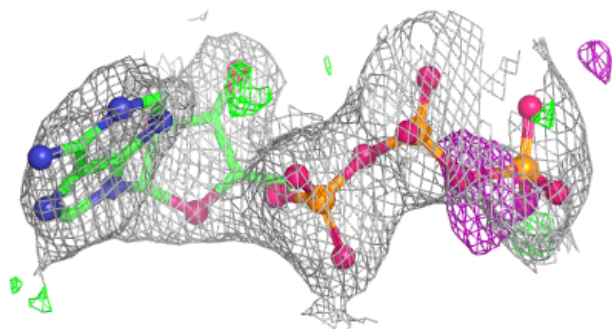
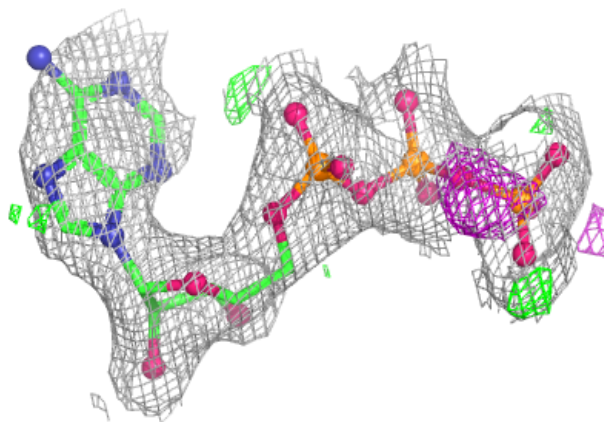
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	ACY	J	304	4/4	0.27	0.25	56,60,62,65	0
3	PO4	E	302	5/5	0.35	0.17	103,108,109,114	0
3	PO4	J	302	5/5	0.53	0.19	128,130,132,136	0
3	PO4	I	302	5/5	0.59	0.14	93,96,97,98	0
3	PO4	C	303	5/5	0.60	0.11	98,101,108,110	0
3	PO4	B	302	5/5	0.62	0.15	92,95,99,99	0
3	PO4	C	302	5/5	0.63	0.13	92,104,106,111	0
2	MG	K	301	1/1	0.73	0.16	70,70,70,70	0
2	MG	L	301	1/1	0.81	0.15	62,62,62,62	0
4	ATP	L	302	31/31	0.86	0.12	38,55,69,70	0
4	ATP	K	302	31/31	0.87	0.12	39,59,71,76	0
4	ATP	H	302	31/31	0.88	0.10	45,56,68,72	0
4	ATP	G	302	31/31	0.89	0.10	37,52,62,68	0
4	ATP	E	303	31/31	0.89	0.10	44,54,70,74	0
4	ATP	I	303	31/31	0.90	0.10	37,48,68,77	0
2	MG	E	301	1/1	0.90	0.07	59,59,59,59	0
4	ATP	D	302	31/31	0.92	0.10	42,56,70,71	0
4	ATP	A	302	31/31	0.92	0.10	27,45,61,62	0
4	ATP	F	302	31/31	0.93	0.09	28,42,58,59	0
2	MG	I	301	1/1	0.93	0.05	43,43,43,43	0
4	ATP	B	303	31/31	0.93	0.09	36,43,53,55	0
4	ATP	C	304	31/31	0.93	0.09	28,40,54,56	0
2	MG	G	301	1/1	0.93	0.06	51,51,51,51	0
4	ATP	J	303	31/31	0.94	0.09	30,44,60,66	0
2	MG	D	301	1/1	0.95	0.09	49,49,49,49	0
2	MG	H	301	1/1	0.95	0.05	55,55,55,55	0
2	MG	J	301	1/1	0.96	0.04	42,42,42,42	0
2	MG	B	301	1/1	0.96	0.08	39,39,39,39	0
2	MG	F	301	1/1	0.96	0.06	39,39,39,39	0
2	MG	C	301	1/1	0.97	0.06	45,45,45,45	0
2	MG	A	301	1/1	0.97	0.11	44,44,44,44	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

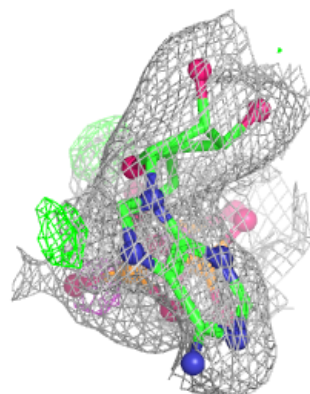
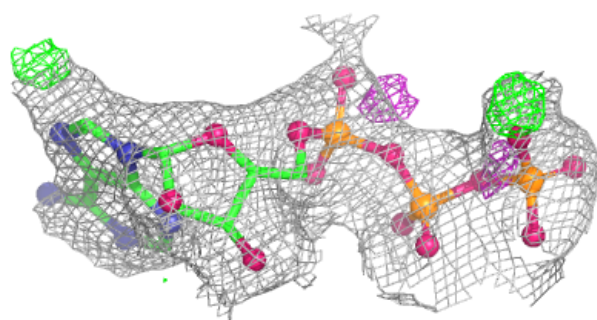
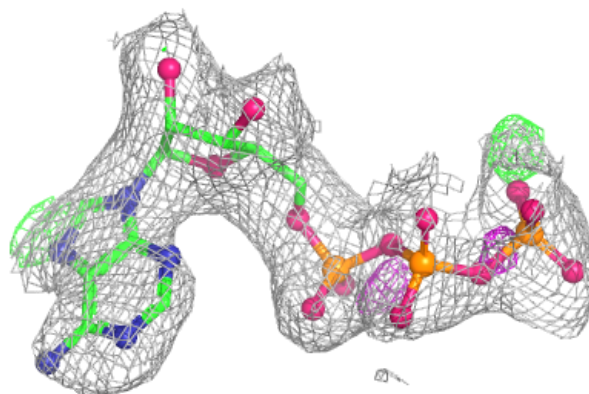
Electron density around ATP L 302:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

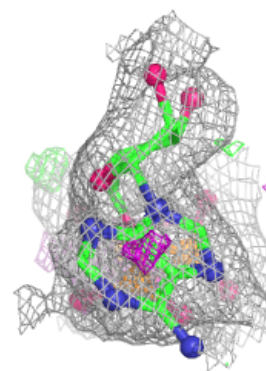
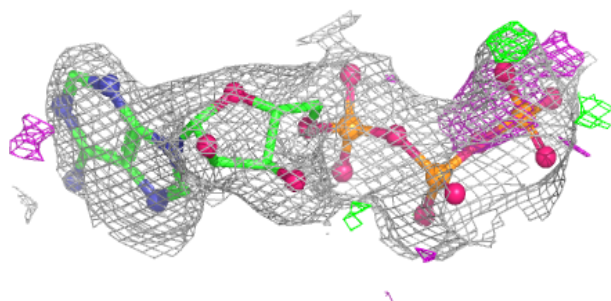
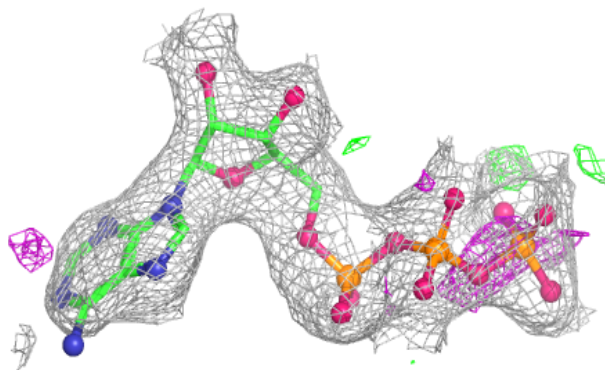


Electron density around ATP K 302:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

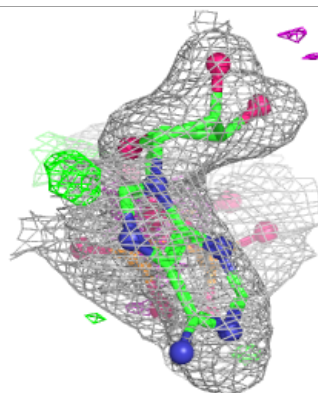
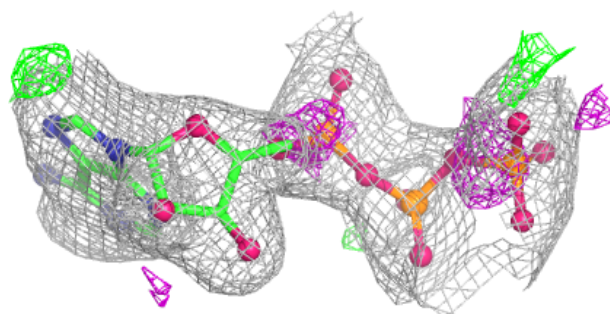
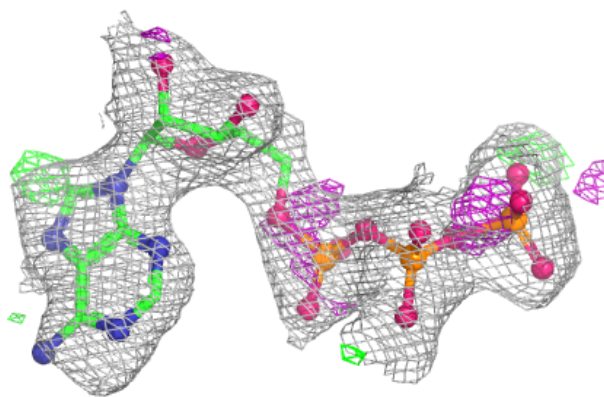
**Electron density around ATP H 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

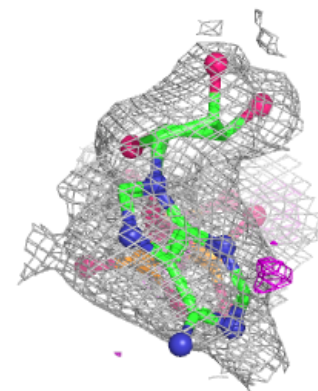
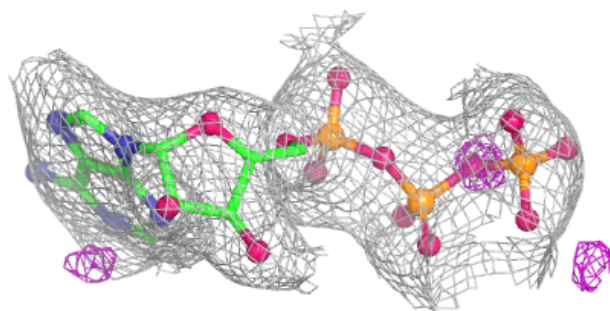
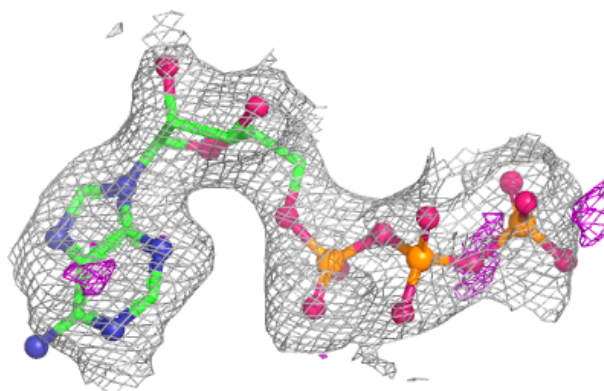


Electron density around ATP G 302:

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and green (positive)

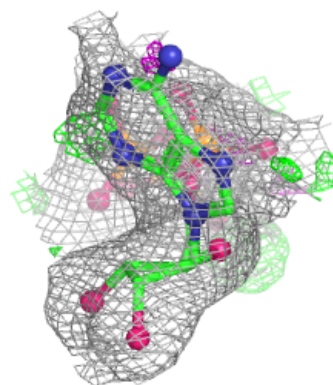
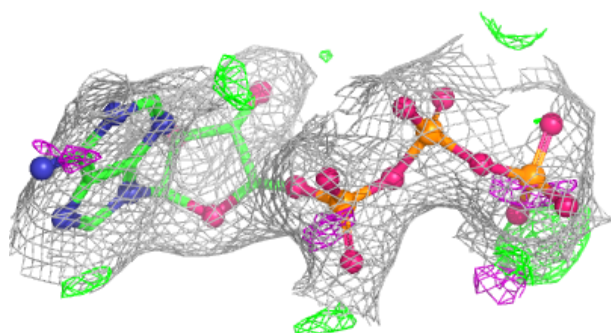
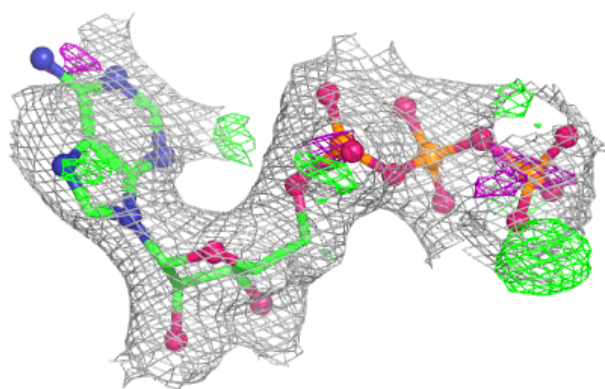
**Electron density around ATP E 303:**

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and green (positive)

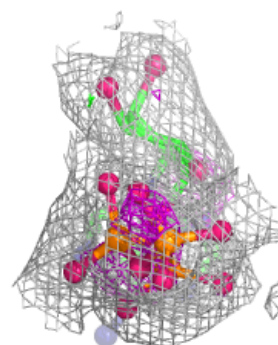
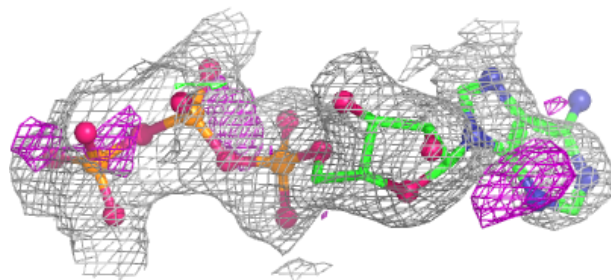
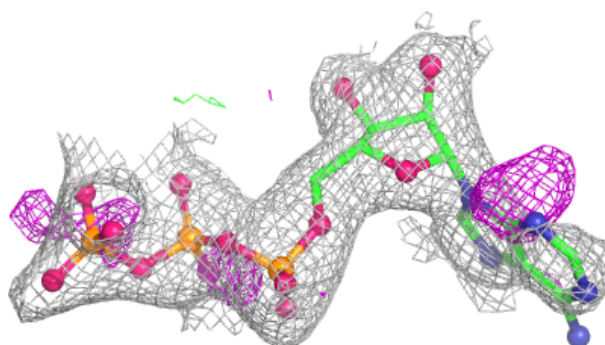


Electron density around ATP I 303:

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and green (positive)

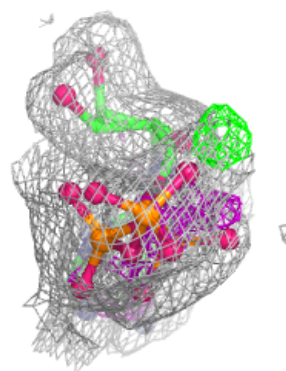
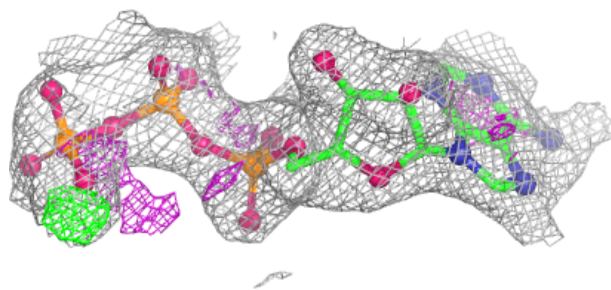
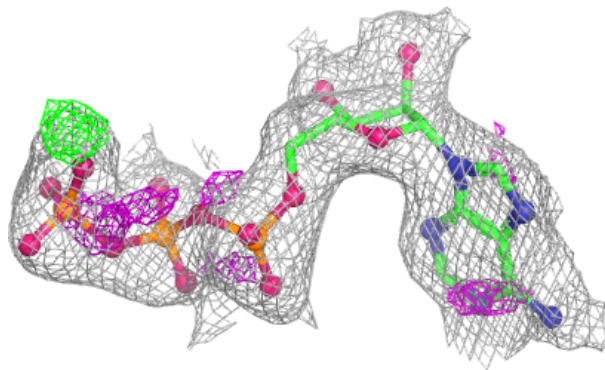
**Electron density around ATP D 302:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

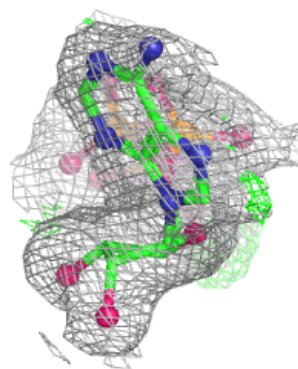
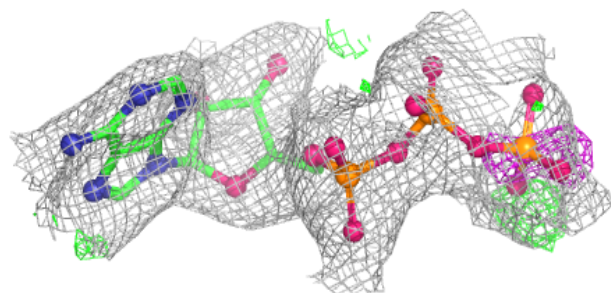
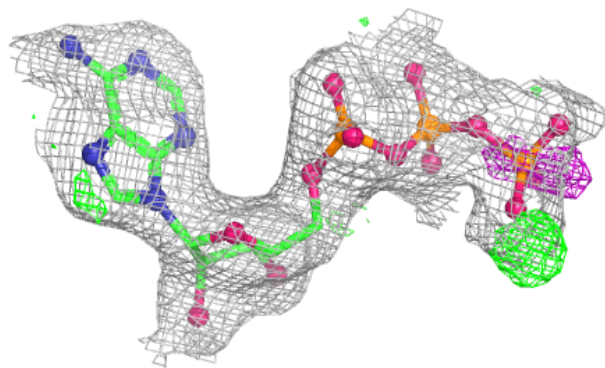


Electron density around ATP A 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

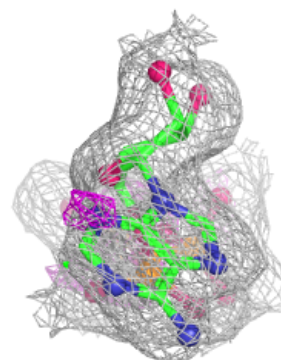
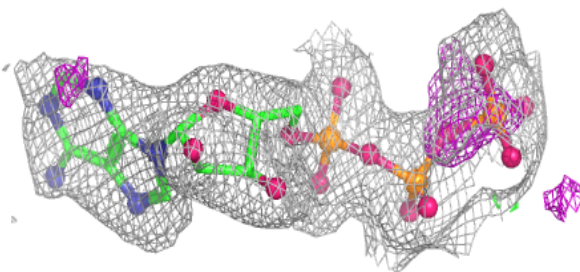
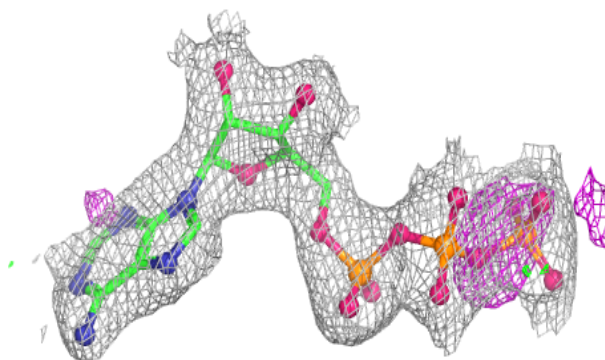
**Electron density around ATP F 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

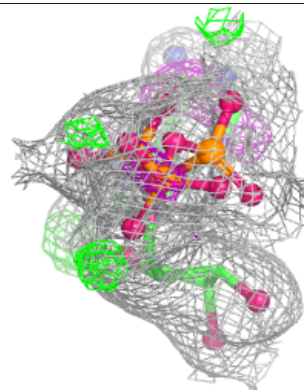
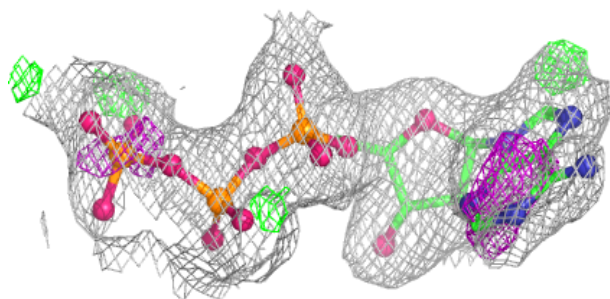
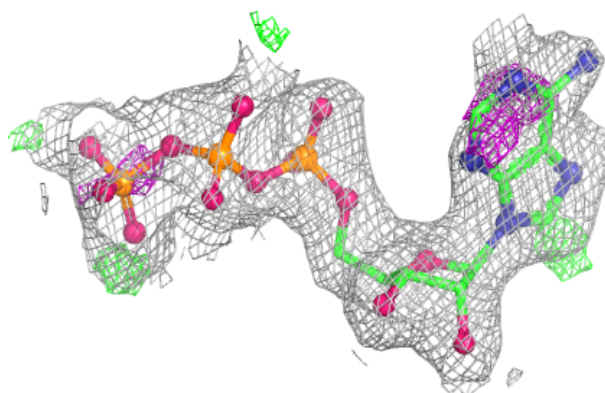


Electron density around ATP B 303:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

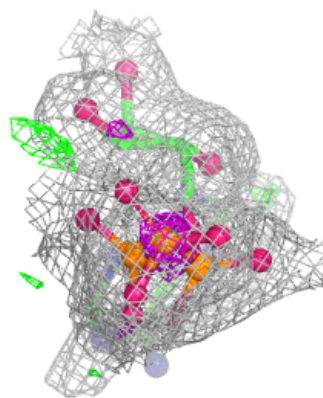
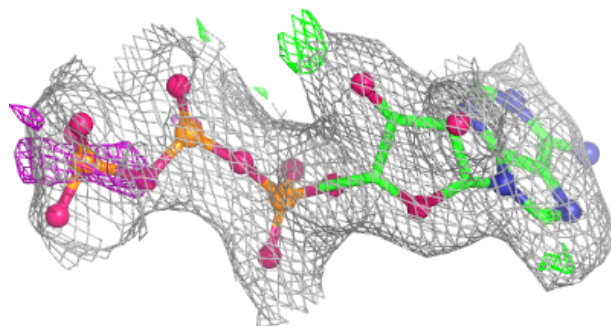
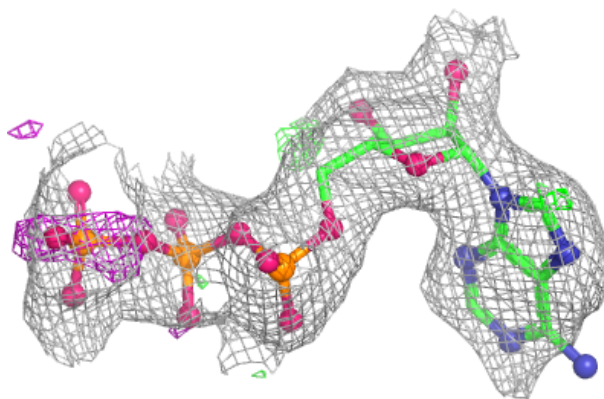
**Electron density around ATP C 304:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ATP J 303:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.