



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 22, 2020 – 04:19 pm BST

PDB ID : 6Q45  
Title : F1-ATPase from *Fusobacterium nucleatum*  
Authors : Petri, J.; Nakatani, Y.; Montgomery, M.G.; Ferguson, S.A.; Aragao, D.; Leslie, A.G.W.; Heikal, A.; Walker, J.E.; Cook, G.M.  
Deposited on : 2018-12-05  
Resolution : 3.60 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](mailto: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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

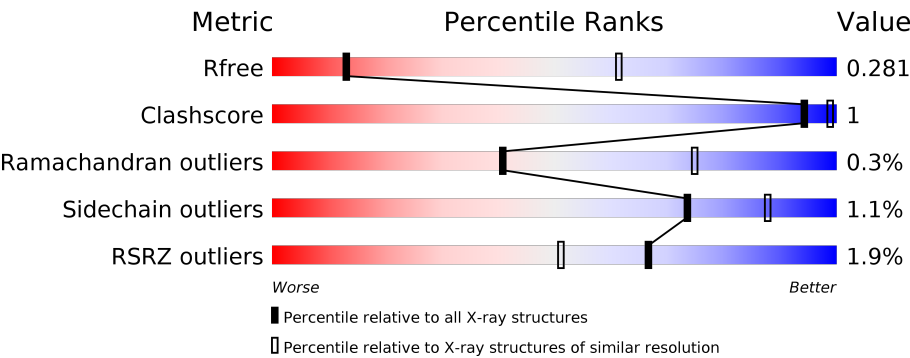
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.60 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))
$R_{free}$	130704	1257 (3.70-3.50)
Clashscore	141614	1353 (3.70-3.50)
Ramachandran outliers	138981	1307 (3.70-3.50)
Sidechain outliers	138945	1307 (3.70-3.50)
RSRZ outliers	127900	1161 (3.70-3.50)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	500	<div><div></div><div>93%<div>5%</div></div></div>
1	B	500	<div><div>2%</div><div>91%<div>5%</div></div></div>
1	C	500	<div><div></div><div>88%<div>6%6%</div></div></div>
1	I	500	<div><div>2%</div><div>93%<div>5%</div></div></div>
1	J	500	<div><div>3%</div><div>91%<div>6%</div></div></div>
1	K	500	<div><div></div><div>86%<div>7%7%</div></div></div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	D	462	<div><div></div><div>95%</div><div></div><div>.</div></div>
2	E	462	<div><div>%</div><div></div><div>95%</div><div>.</div></div>
2	F	462	<div><div></div><div>94%</div><div>5%</div><div></div></div>
2	L	462	<div><div></div><div>95%</div><div>5%</div><div>.</div></div>
2	M	462	<div><div>11%</div><div></div><div>96%</div><div>.</div><div>.</div></div>
2	N	462	<div><div></div><div>94%</div><div>6%</div><div></div></div>
3	G	282	<div><div>5%</div><div></div><div>90%</div><div>9%</div></div>
3	O	282	<div><div>2%</div><div></div><div>90%</div><div>10%</div></div>
4	H	134	<div><div></div><div>96%</div><div></div><div>.</div><div>.</div></div>
4	P	134	<div><div>7%</div><div></div><div>99%</div><div>.</div></div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 49762 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 synthase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	476	Total	C	N	O	S	0	0	0
			3650	2316	625	696	13			
1	B	474	Total	C	N	O	S	0	0	0
			3633	2304	622	694	13			
1	C	471	Total	C	N	O	S	0	0	0
			3613	2292	620	688	13			
1	I	474	Total	C	N	O	S	0	0	0
			3633	2304	622	694	13			
1	J	469	Total	C	N	O	S	0	0	0
			3597	2282	618	684	13			
1	K	466	Total	C	N	O	S	0	0	0
			3566	2262	611	680	13			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	460	Total	C	N	O	S	0	0	0
			3519	2227	591	692	9			
2	E	460	Total	C	N	O	S	0	0	0
			3519	2227	591	692	9			
2	F	462	Total	C	N	O	S	0	0	0
			3533	2236	594	694	9			
2	L	459	Total	C	N	O	S	0	0	0
			3511	2222	590	691	8			
2	M	459	Total	C	N	O	S	0	0	0
			3511	2222	590	691	8			
2	N	460	Total	C	N	O	S	0	0	0
			3519	2227	591	692	9			

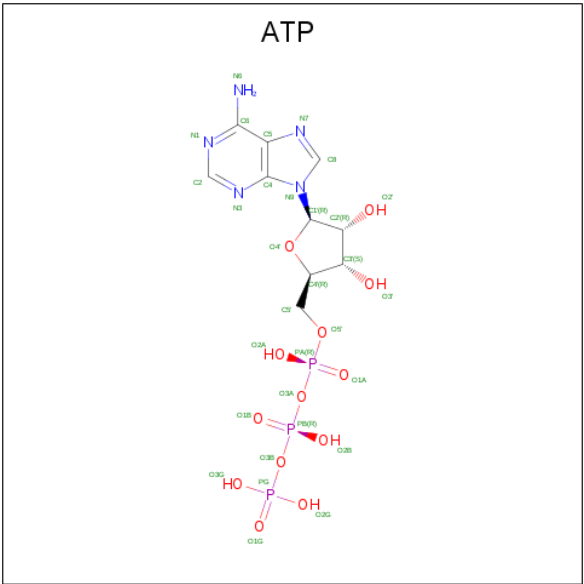
- Molecule 3 is a protein called ATP synthase gamma chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	281	Total	C	N	O	S	0	0	0
			2248	1416	386	438	8			
3	O	281	Total	C	N	O	S	0	0	0
			2248	1416	386	438	8			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	134	Total	C	N	O	S	0	0	0
			1062	671	177	211	3			
4	P	134	Total	C	N	O	S	0	0	0
			1062	671	177	211	3			

- Molecule 5 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>).



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

Continued on next page...

*Continued from previous page...*

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

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	J	1	Total	Mg	0	0
			1	1		
6	D	1	Total	Mg	0	0
			1	1		
6	K	1	Total	Mg	0	0
			1	1		
6	B	1	Total	Mg	0	0
			1	1		
6	I	1	Total	Mg	0	0
			1	1		
6	C	1	Total	Mg	0	0
			1	1		
6	A	1	Total	Mg	0	0
			1	1		
6	N	1	Total	Mg	0	0
			1	1		
6	L	1	Total	Mg	0	0
			1	1		
6	F	1	Total	Mg	0	0
			1	1		

- Molecule 7 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	3	Total	O	0	0
			3	3		
8	B	3	Total	O	0	0
			3	3		
8	C	3	Total	O	0	0
			3	3		
8	D	4	Total	O	0	0
			4	4		
8	F	4	Total	O	0	0
			4	4		
8	I	3	Total	O	0	0
			3	3		
8	J	3	Total	O	0	0
			3	3		
8	K	3	Total	O	0	0
			3	3		

Continued on next page...

*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	4	Total	O	0	0
			4	4		
8	N	4	Total	O	0	0
			4	4		

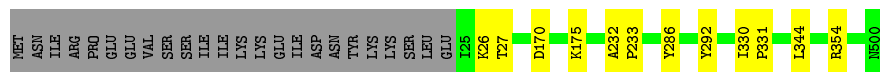


### 3 Residue-property plots [i](#)


These plots are drawn for all protein, RNA and DNA 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 electron 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 dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). 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.

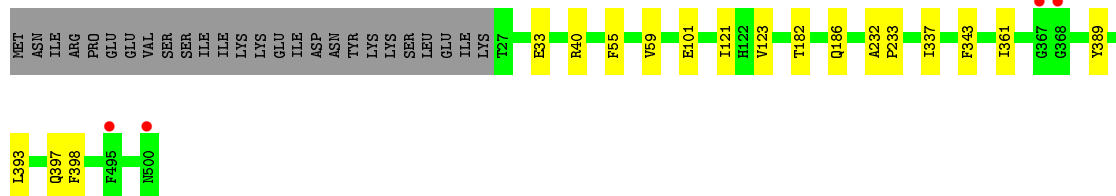
- Molecule 1: ATP synthase subunit alpha

Chain A: 




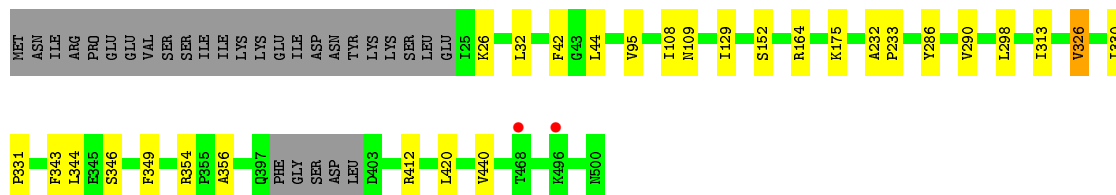
- Molecule 1: ATP synthase subunit alpha

Chain B: 




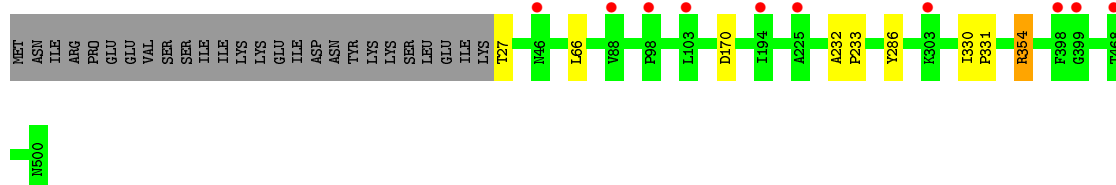
- Molecule 1: ATP synthase subunit alpha

Chain C: 

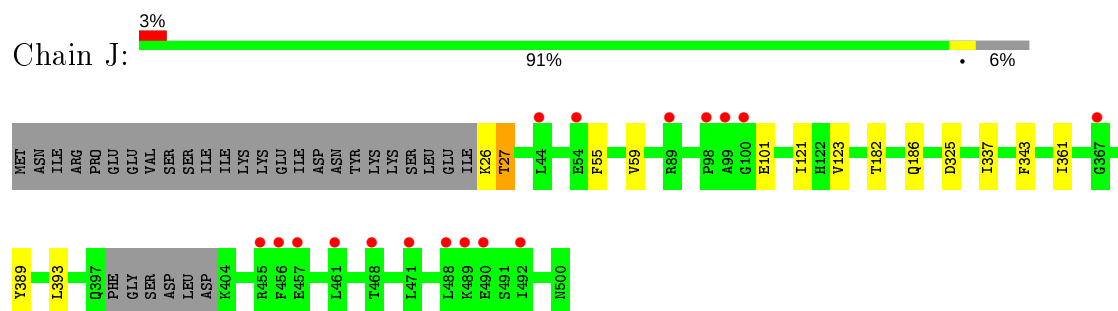


- Molecule 1: ATP synthase subunit alpha

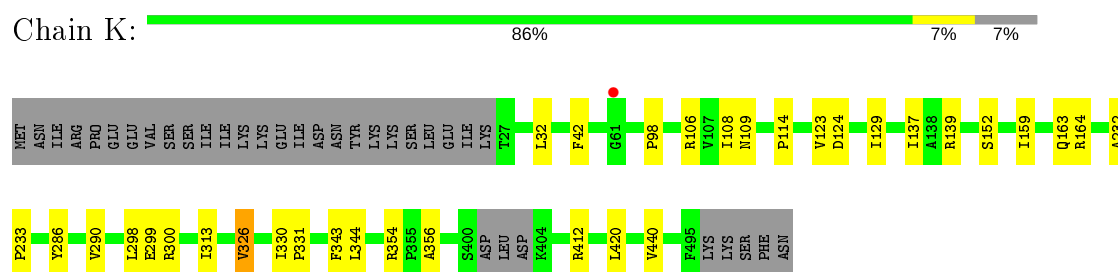
Chain I: 



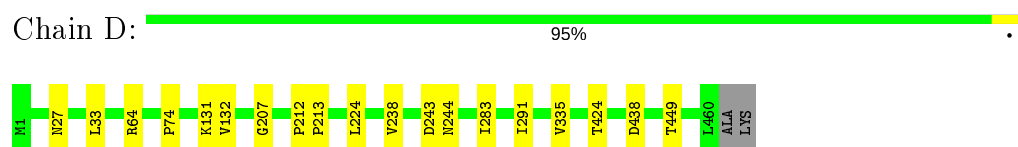
- Molecule 1: ATP synthase subunit alpha



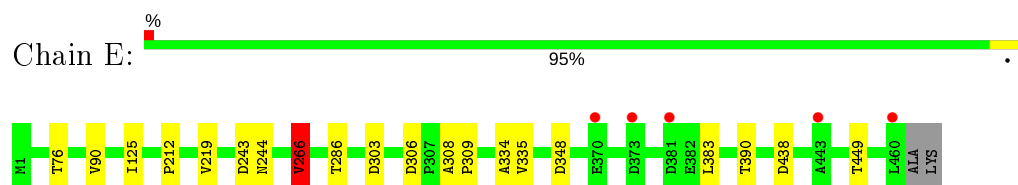
- Molecule 1: ATP synthase subunit alpha



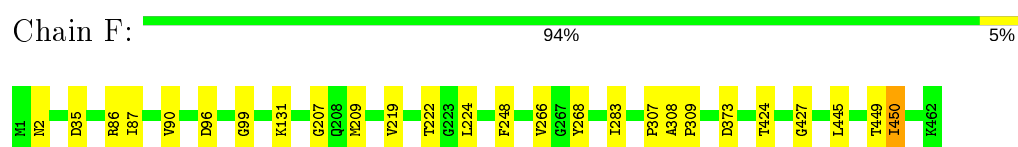
- Molecule 2: ATP synthase subunit beta



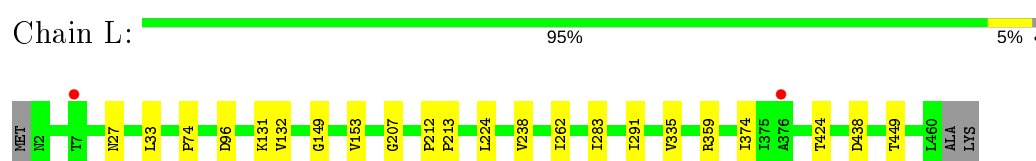
- Molecule 2: ATP synthase subunit beta



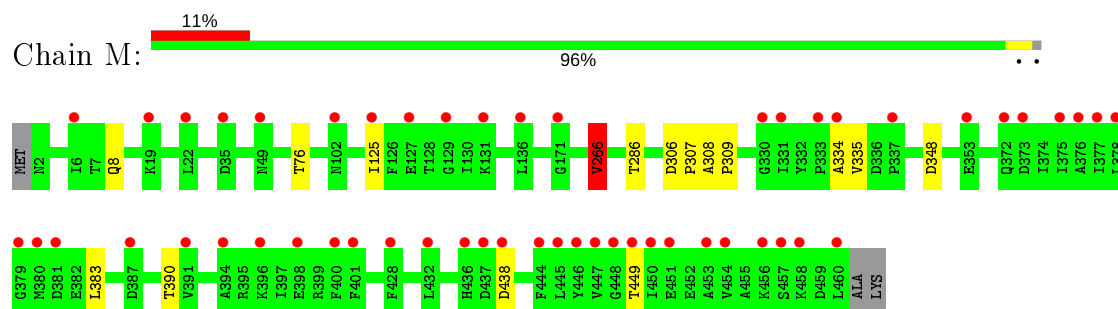
- Molecule 2: ATP synthase subunit beta



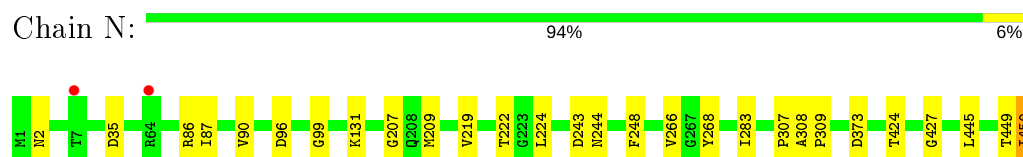
- Molecule 2: ATP synthase subunit beta



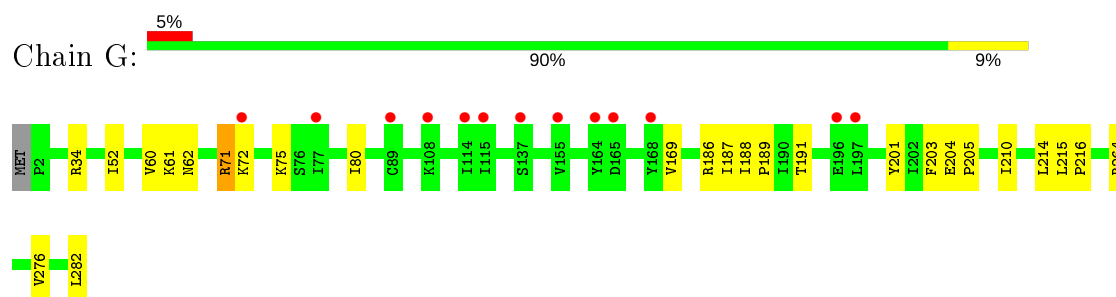
- Molecule 2: ATP synthase subunit beta



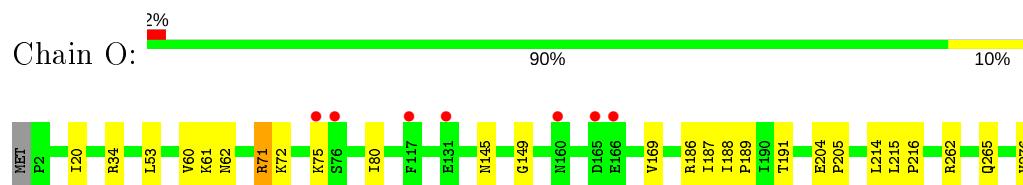
- Molecule 2: ATP synthase subunit beta



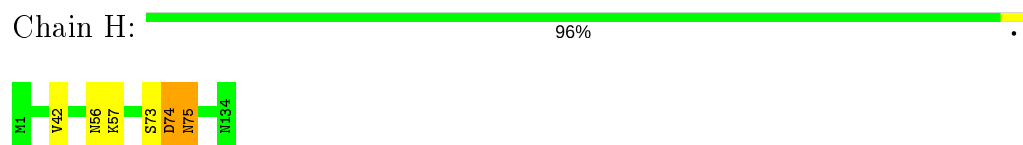
- Molecule 3: ATP synthase gamma chain



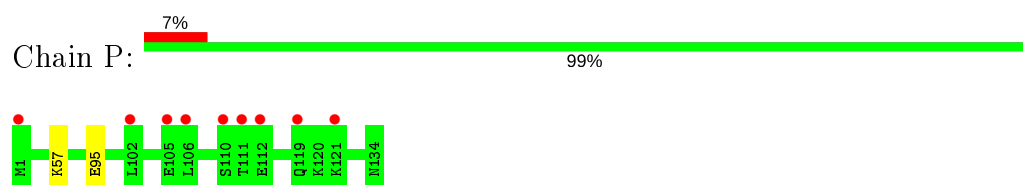
- Molecule 3: ATP synthase gamma chain



- Molecule 4: ATP synthase epsilon chain



- Molecule 4: ATP synthase epsilon chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.94Å 200.21Å 201.72Å 90.00° 102.20° 90.00°	Depositor
Resolution (Å)	49.34 – 3.60 49.29 – 3.60	Depositor EDS
% Data completeness (in resolution range)	96.9 (49.34-3.60) 96.9 (49.29-3.60)	Depositor EDS
$R_{merge}$	0.21	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.94 (at 3.57Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.237 , 0.280 0.238 , 0.281	Depositor DCC
$R_{free}$ test set	4872 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	96.3	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 48.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	49762	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.23% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MG, ATP, ADP

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.29	0/3704	0.46	0/4997
1	B	0.29	0/3687	0.46	0/4975
1	C	0.29	0/3665	0.47	0/4943
1	I	0.29	0/3687	0.46	0/4975
1	J	0.29	0/3649	0.46	0/4921
1	K	0.29	0/3618	0.47	0/4882
2	D	0.28	0/3572	0.47	0/4837
2	E	0.29	0/3572	0.47	0/4837
2	F	0.29	0/3586	0.48	0/4855
2	L	0.29	0/3564	0.47	0/4827
2	M	0.29	0/3564	0.47	0/4827
2	N	0.29	0/3572	0.48	0/4837
3	G	0.30	0/2281	0.47	0/3069
3	O	0.30	0/2281	0.47	0/3069
4	H	0.30	0/1074	0.48	0/1441
4	P	0.29	0/1074	0.48	0/1441
All	All	0.29	0/50150	0.47	0/67733

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	A	0	1
1	C	0	1
1	I	0	1
1	K	0	1
3	G	0	3
3	O	0	3
All	All	0	10

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 10 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	354	ARG	Sidechain
1	C	354	ARG	Sidechain
3	G	186	ARG	Sidechain
3	G	34	ARG	Sidechain
3	G	71	ARG	Sidechain

## 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3650	0	3741	5	0
1	B	3633	0	3717	9	0
1	C	3613	0	3708	15	0
1	I	3633	0	3717	5	0
1	J	3597	0	3693	8	0
1	K	3566	0	3651	17	0
2	D	3519	0	3567	8	0
2	E	3519	0	3568	7	0
2	F	3533	0	3585	10	0
2	L	3511	0	3555	11	0
2	M	3511	0	3556	8	0
2	N	3519	0	3567	12	0
3	G	2248	0	2290	15	0
3	O	2248	0	2290	19	0
4	H	1062	0	1094	4	0
4	P	1062	0	1094	0	0
5	A	31	0	12	0	0
5	B	31	0	12	0	0
5	C	31	0	12	0	0
5	I	31	0	12	0	0
5	J	31	0	12	0	0
5	K	31	0	12	0	0
6	A	1	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
6	F	1	0	0	0	0
6	I	1	0	0	0	0
6	J	1	0	0	0	0
6	K	1	0	0	0	0
6	L	1	0	0	0	0
6	N	1	0	0	0	0
7	D	27	0	12	0	0
7	F	27	0	12	0	0
7	L	27	0	12	0	0
7	N	27	0	12	0	0
8	A	3	0	0	0	0
8	B	3	0	0	0	0
8	C	3	0	0	0	0
8	D	4	0	0	0	0
8	F	4	0	0	0	0
8	I	3	0	0	0	0
8	J	3	0	0	0	0
8	K	3	0	0	0	0
8	L	4	0	0	0	0
8	N	4	0	0	0	0
All	All	49762	0	50513	134	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

The worst 5 of 134 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)
3:O:53:LEU:HD11	3:O:214:LEU:HD21	1.64	0.80
2:F:308:ALA:HB3	2:F:309:PRO:HD3	1.75	0.69
2:N:308:ALA:HB3	2:N:309:PRO:HD3	1.76	0.66
2:M:306:ASP:HA	3:O:265:GLN:OE1	1.96	0.65
3:O:188:ILE:HG22	3:O:189:PRO:HD3	1.80	0.64

There are no symmetry-related clashes.

## 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	474/500 (95%)	453 (96%)	21 (4%)	0	100	100
1	B	472/500 (94%)	446 (94%)	26 (6%)	0	100	100
1	C	467/500 (93%)	443 (95%)	23 (5%)	1 (0%)	47	79
1	I	472/500 (94%)	452 (96%)	20 (4%)	0	100	100
1	J	465/500 (93%)	440 (95%)	24 (5%)	1 (0%)	47	79
1	K	462/500 (92%)	439 (95%)	22 (5%)	1 (0%)	47	79
2	D	458/462 (99%)	421 (92%)	36 (8%)	1 (0%)	47	79
2	E	458/462 (99%)	422 (92%)	33 (7%)	3 (1%)	22	61
2	F	460/462 (100%)	423 (92%)	34 (7%)	3 (1%)	22	61
2	L	457/462 (99%)	418 (92%)	38 (8%)	1 (0%)	47	79
2	M	457/462 (99%)	421 (92%)	33 (7%)	3 (1%)	22	61
2	N	458/462 (99%)	420 (92%)	35 (8%)	3 (1%)	22	61
3	G	279/282 (99%)	267 (96%)	12 (4%)	0	100	100
3	O	279/282 (99%)	267 (96%)	12 (4%)	0	100	100
4	H	132/134 (98%)	122 (92%)	7 (5%)	3 (2%)	6	38
4	P	132/134 (98%)	128 (97%)	3 (2%)	1 (1%)	19	59
All	All	6382/6604 (97%)	5982 (94%)	379 (6%)	21 (0%)	41	75

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	H	74	ASP
4	H	75	ASN
2	M	266	VAL
2	E	266	VAL
4	H	57	LYS



### 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	390/414 (94%)	387 (99%)	3 (1%)	81	91
1	B	388/414 (94%)	387 (100%)	1 (0%)	92	97
1	C	386/414 (93%)	384 (100%)	2 (0%)	88	95
1	I	388/414 (94%)	386 (100%)	2 (0%)	88	95
1	J	384/414 (93%)	383 (100%)	1 (0%)	92	97
1	K	380/414 (92%)	378 (100%)	2 (0%)	88	95
2	D	382/383 (100%)	379 (99%)	3 (1%)	81	91
2	E	382/383 (100%)	373 (98%)	9 (2%)	49	75
2	F	383/383 (100%)	378 (99%)	5 (1%)	69	86
2	L	381/383 (100%)	378 (99%)	3 (1%)	81	91
2	M	381/383 (100%)	372 (98%)	9 (2%)	49	75
2	N	382/383 (100%)	377 (99%)	5 (1%)	69	86
3	G	255/256 (100%)	249 (98%)	6 (2%)	49	75
3	O	255/256 (100%)	249 (98%)	6 (2%)	49	75
4	H	120/120 (100%)	119 (99%)	1 (1%)	81	91
4	P	120/120 (100%)	119 (99%)	1 (1%)	81	91
All	All	5357/5534 (97%)	5298 (99%)	59 (1%)	73	88

5 of 59 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	G	191	THR
1	K	124	ASP
3	O	75	LYS
3	G	276	VAL
4	H	56	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	J	78	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 10 are monoatomic - leaving 10 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$
7	ADP	D	600	6	24,29,29	1.02	2 (8%)	29,45,45	1.51	4 (13%)
5	ATP	B	600	6	26,33,33	0.96	2 (7%)	31,52,52	1.61	5 (16%)
5	ATP	A	600	6	26,33,33	0.99	2 (7%)	31,52,52	1.47	5 (16%)
7	ADP	L	600	6	24,29,29	1.03	2 (8%)	29,45,45	1.43	4 (13%)
5	ATP	C	600	6	26,33,33	0.98	2 (7%)	31,52,52	1.52	5 (16%)
5	ATP	J	600	6	26,33,33	1.01	2 (7%)	31,52,52	1.55	5 (16%)
5	ATP	I	600	6	26,33,33	0.95	1 (3%)	31,52,52	1.56	6 (19%)
5	ATP	K	600	6	26,33,33	0.98	1 (3%)	31,52,52	1.62	5 (16%)
7	ADP	F	600	6	24,29,29	1.02	2 (8%)	29,45,45	1.37	4 (13%)
7	ADP	N	600	6	24,29,29	1.01	2 (8%)	29,45,45	1.47	4 (13%)

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
7	ADP	D	600	6	-	0/12/32/32	0/3/3/3
5	ATP	B	600	6	-	0/18/38/38	0/3/3/3
5	ATP	A	600	6	-	0/18/38/38	0/3/3/3
7	ADP	L	600	6	-	1/12/32/32	0/3/3/3
5	ATP	C	600	6	-	0/18/38/38	0/3/3/3
5	ATP	J	600	6	-	0/18/38/38	0/3/3/3
5	ATP	I	600	6	-	0/18/38/38	0/3/3/3
5	ATP	K	600	6	-	0/18/38/38	0/3/3/3
7	ADP	F	600	6	-	3/12/32/32	0/3/3/3
7	ADP	N	600	6	-	0/12/32/32	0/3/3/3

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	K	600	ATP	C5-C4	2.63	1.47	1.40
5	J	600	ATP	C5-C4	2.61	1.47	1.40
5	C	600	ATP	C5-C4	2.56	1.47	1.40
7	L	600	ADP	C5-C4	2.55	1.47	1.40
7	D	600	ADP	C5-C4	2.55	1.47	1.40

The worst 5 of 47 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	600	ATP	PA-O3A-PB	-4.22	118.34	132.83
5	K	600	ATP	PA-O3A-PB	-3.98	119.15	132.83
5	I	600	ATP	N3-C2-N1	-3.77	122.79	128.68
7	N	600	ADP	PA-O3A-PB	-3.68	120.20	132.83
7	L	600	ADP	N3-C2-N1	-3.68	122.93	128.68

There are no chirality outliers.

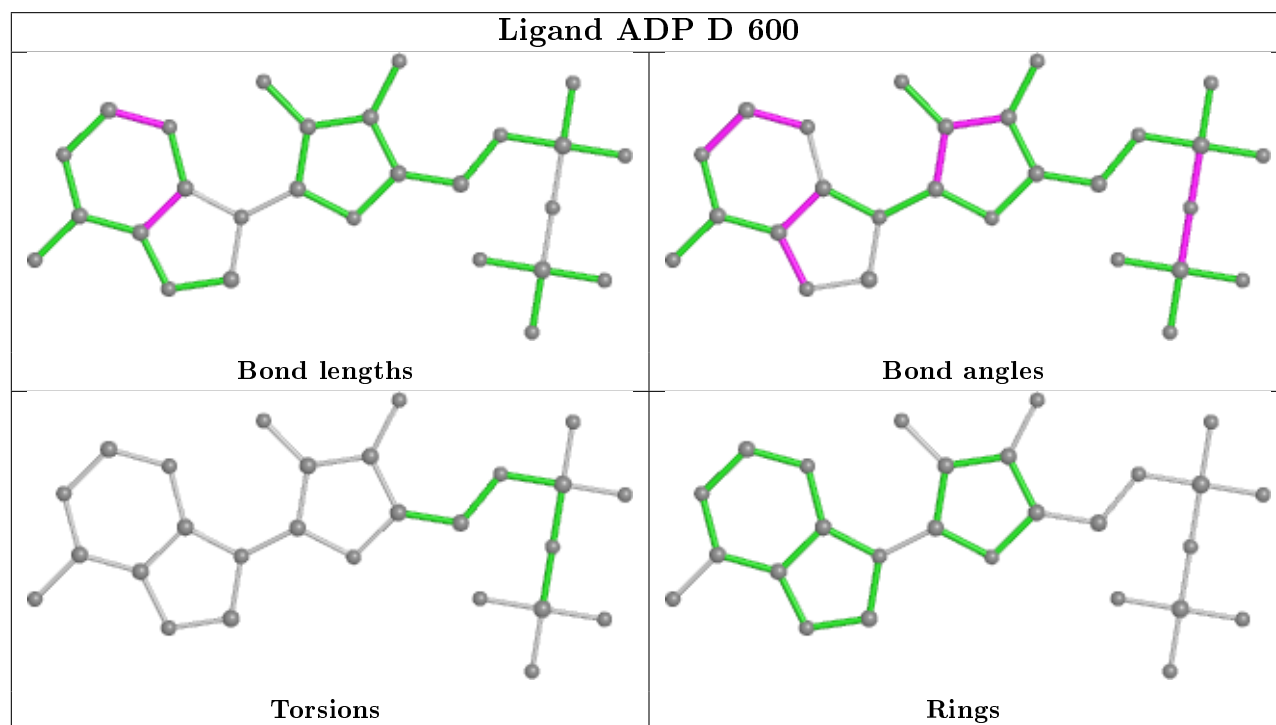
All (4) torsion outliers are listed below:

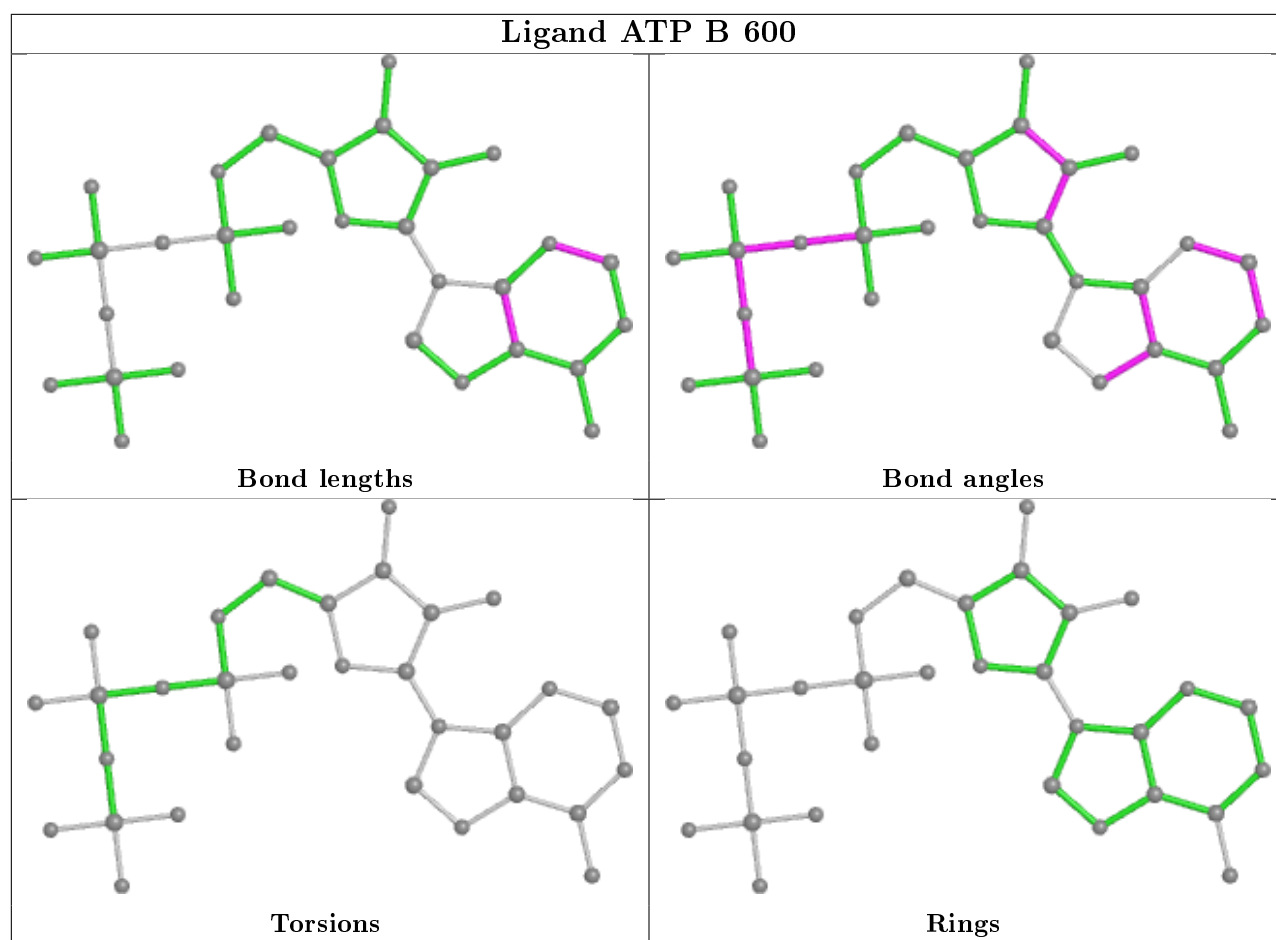
Mol	Chain	Res	Type	Atoms
7	F	600	ADP	O4'-C4'-C5'-O5'
7	F	600	ADP	C3'-C4'-C5'-O5'
7	F	600	ADP	C5'-O5'-PA-O1A
7	L	600	ADP	C5'-O5'-PA-O1A

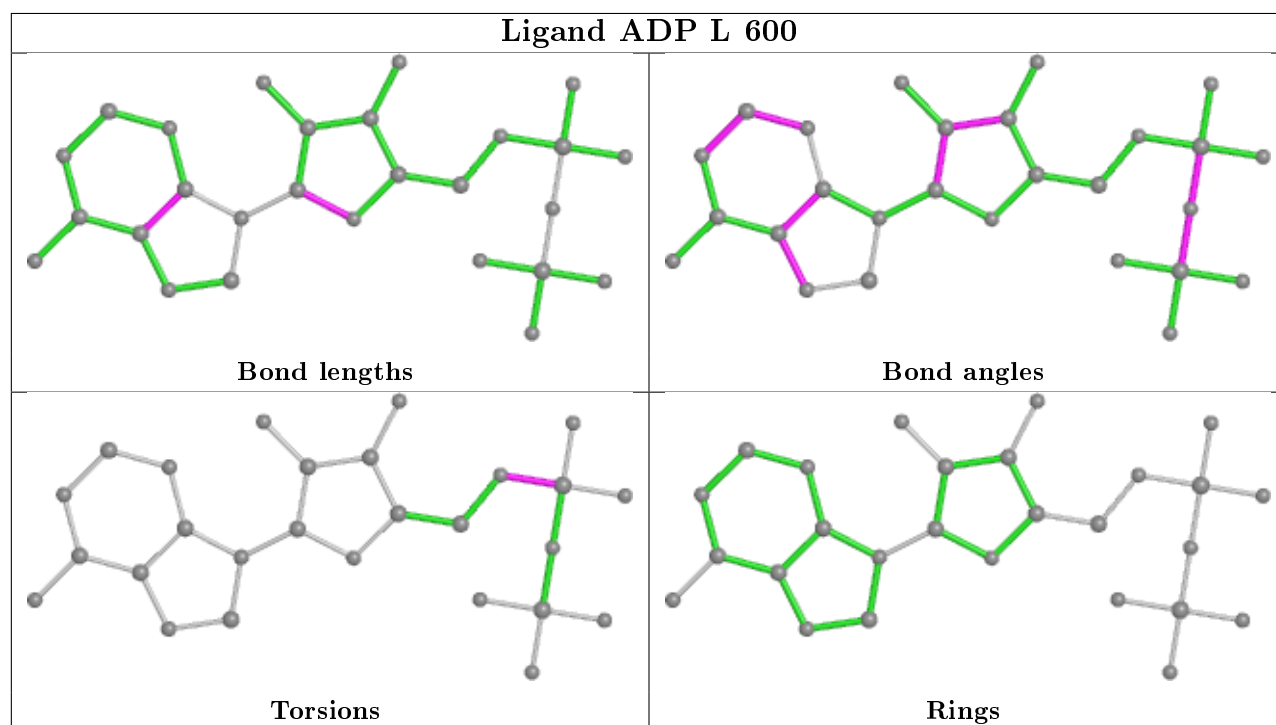
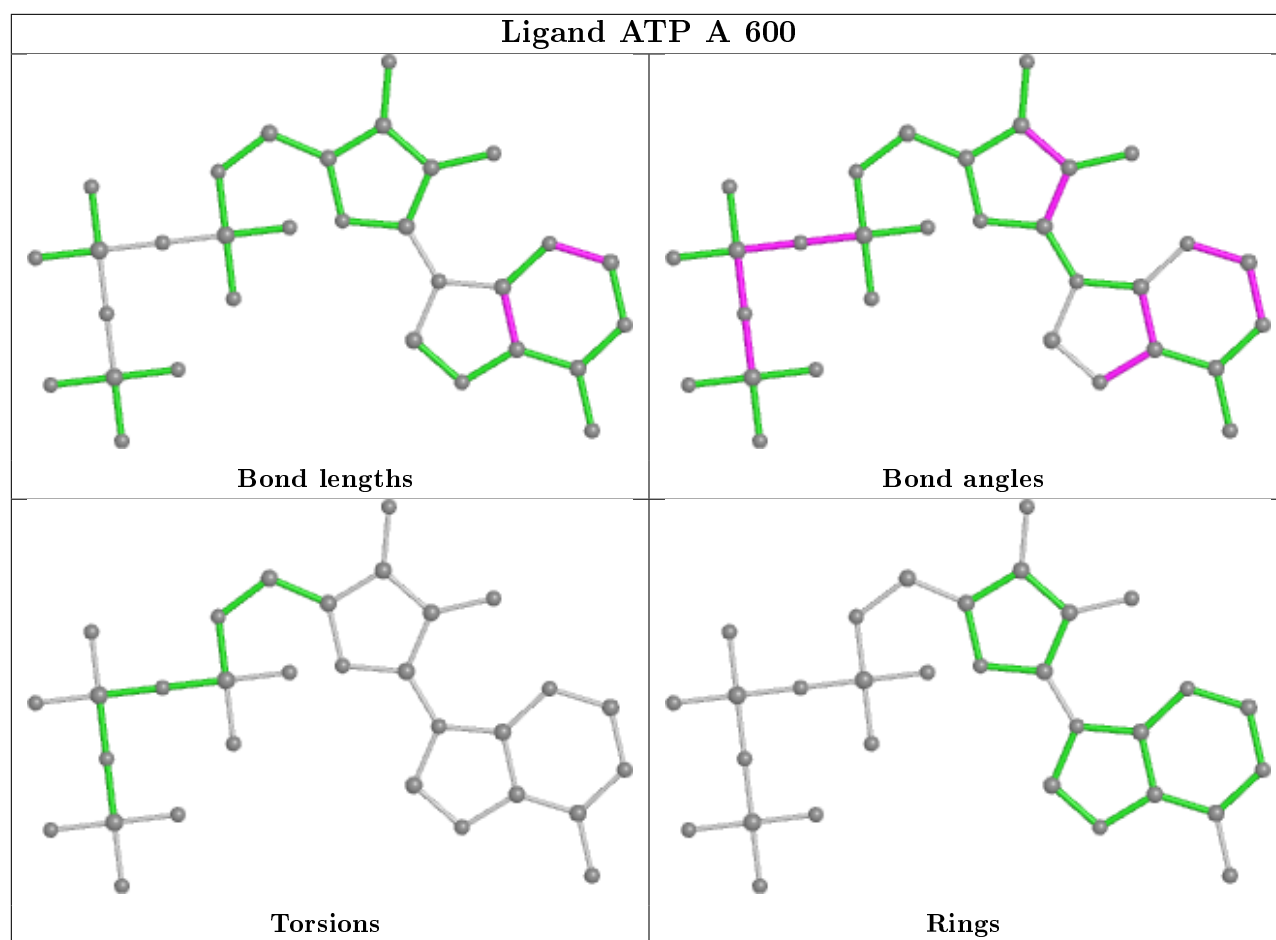
There are no ring outliers.

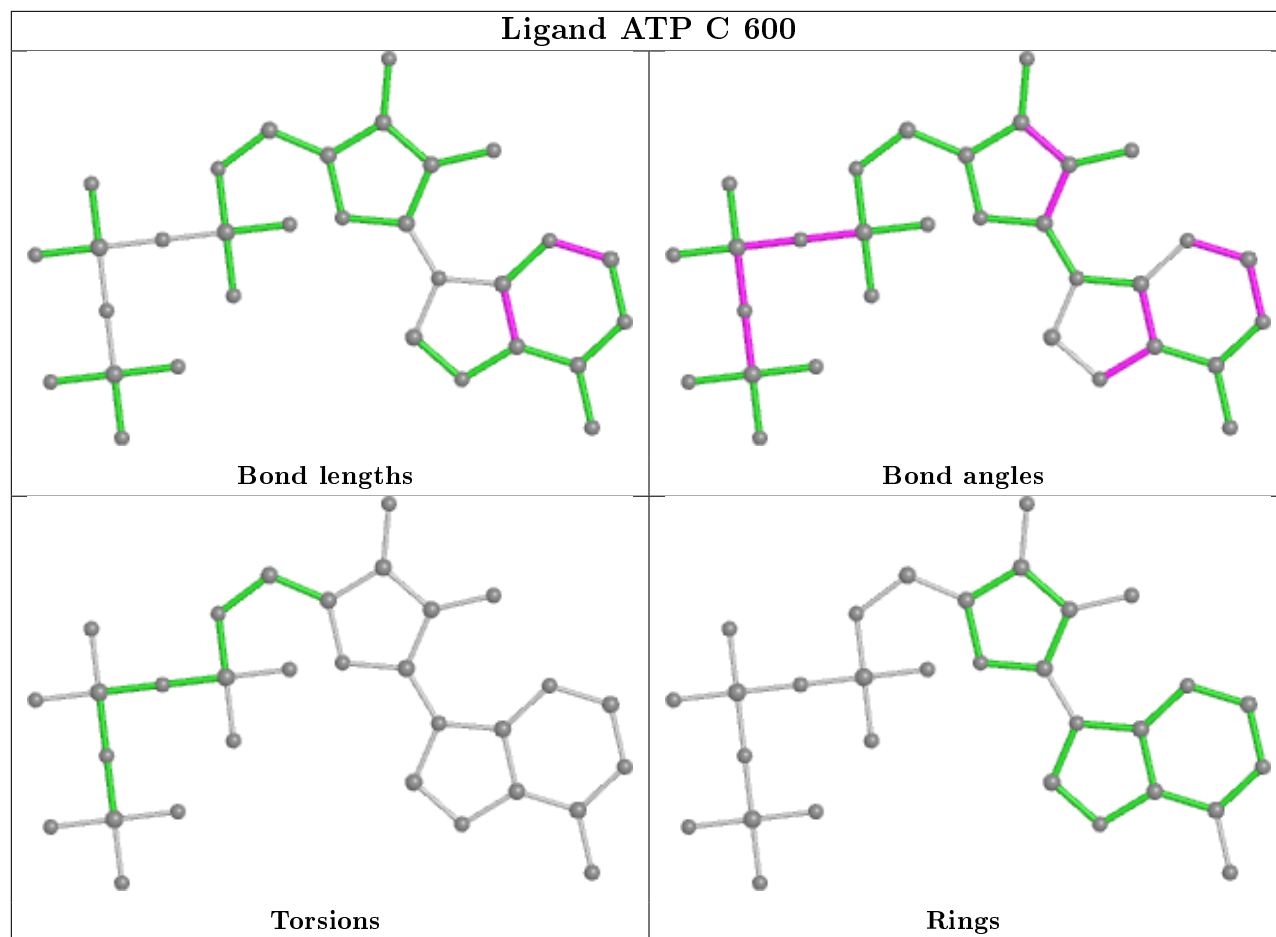
No monomer is involved in short contacts.

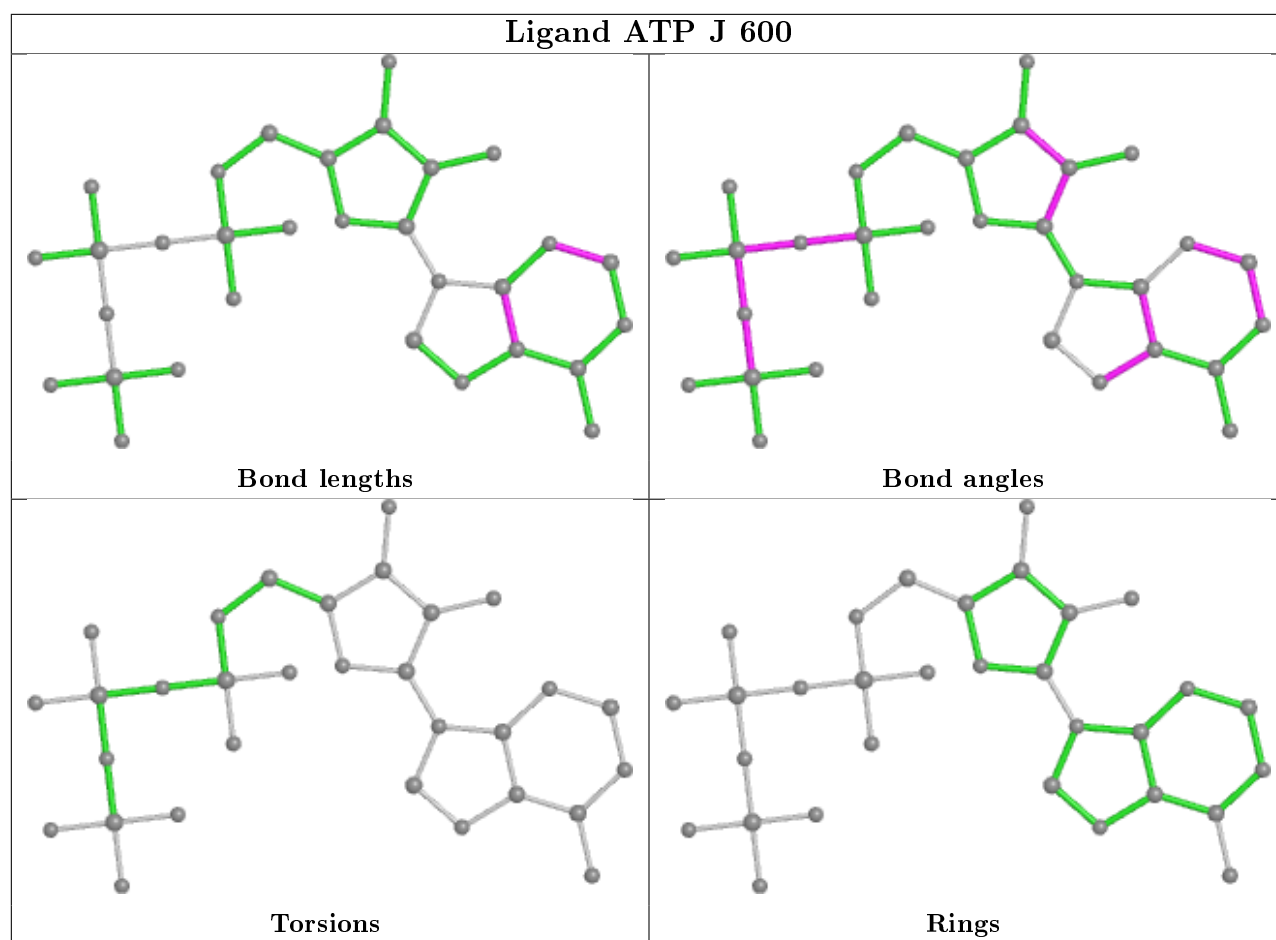
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



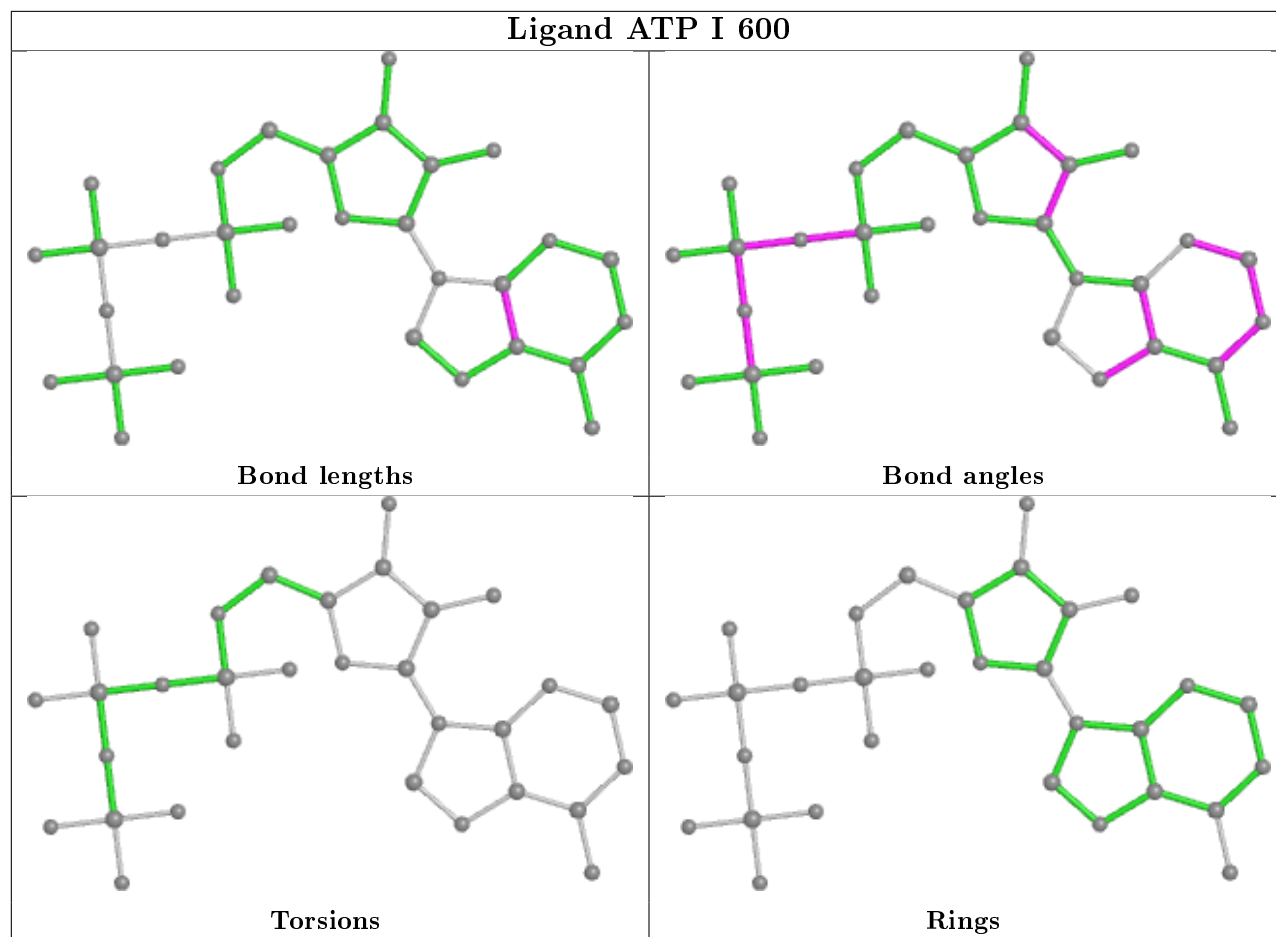


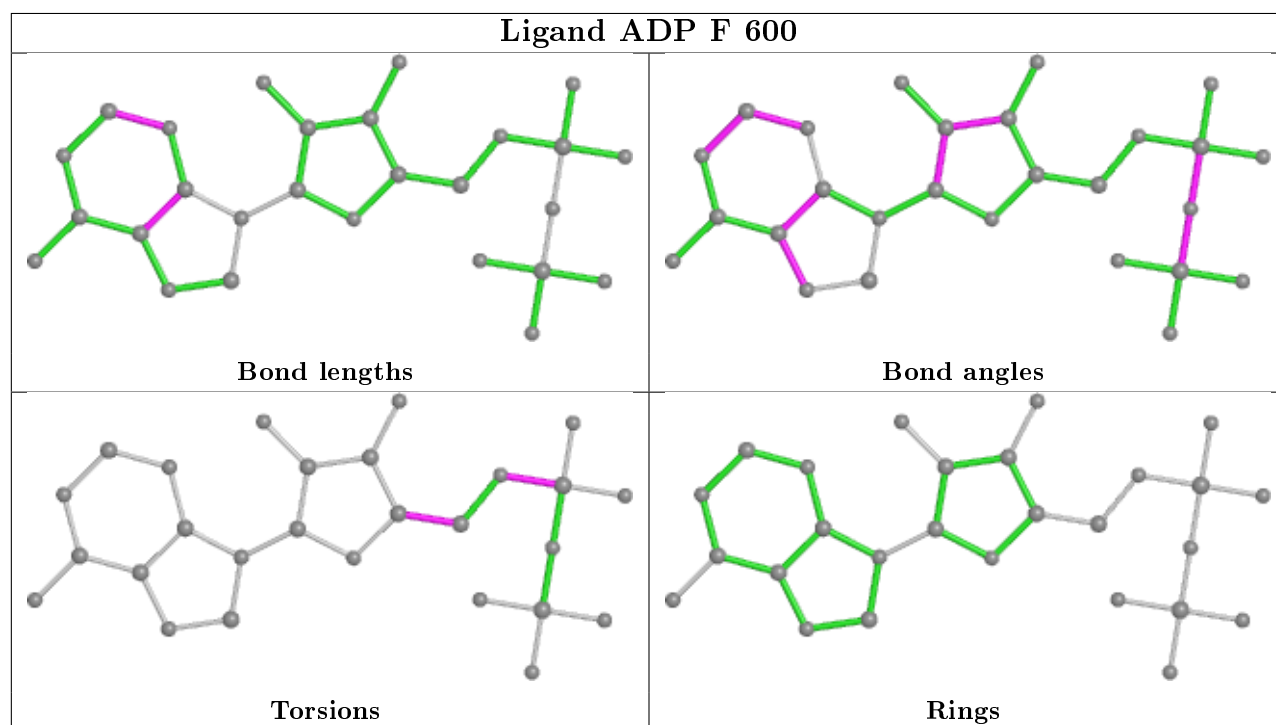
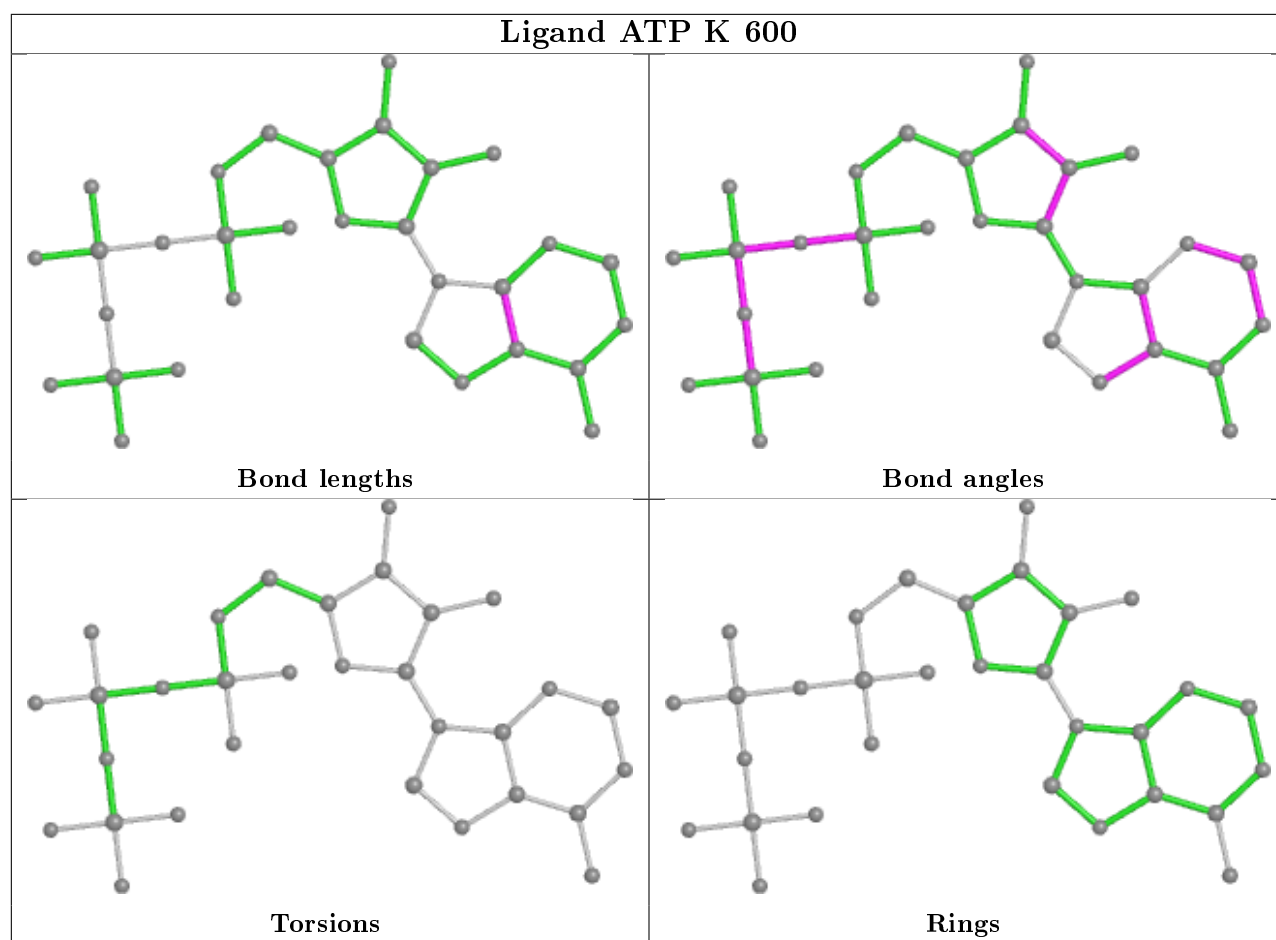


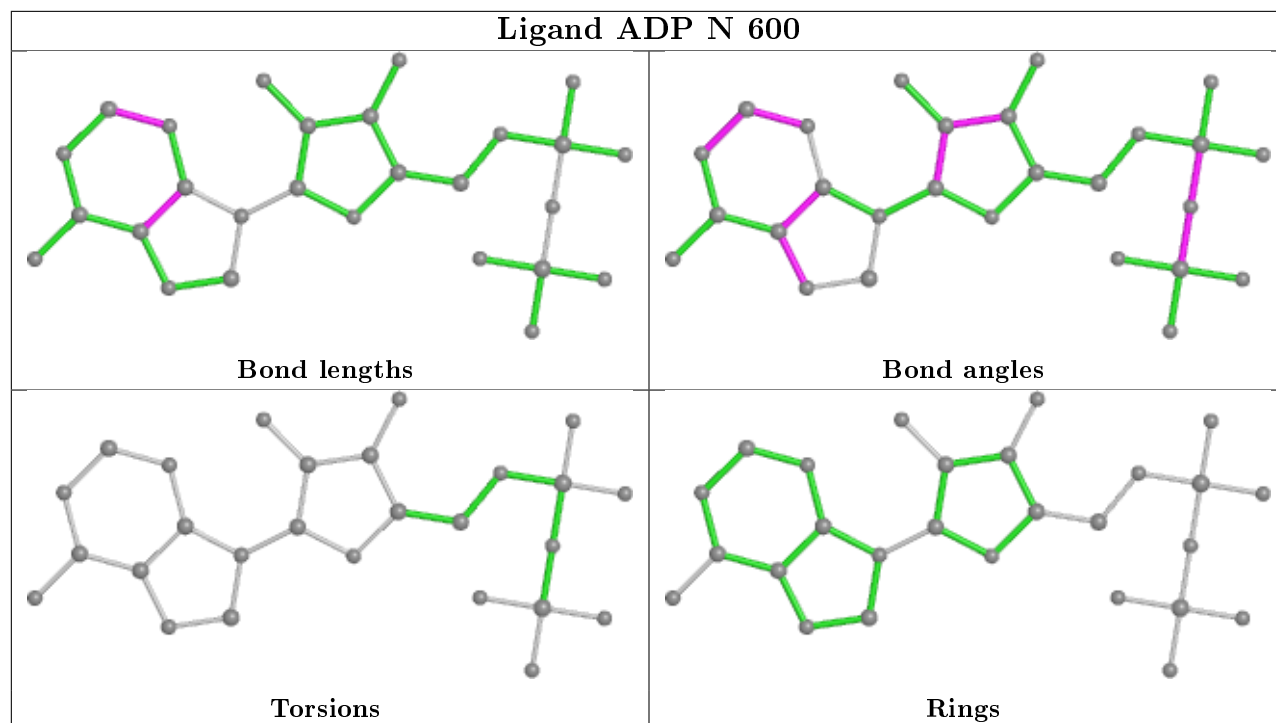












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	476/500 (95%)	-0.28	0 100 100	54, 85, 121, 147	0
1	B	474/500 (94%)	-0.25	4 (0%) 86 75	52, 94, 154, 190	0
1	C	471/500 (94%)	-0.26	2 (0%) 92 86	53, 87, 134, 158	0
1	I	474/500 (94%)	0.16	10 (2%) 63 48	94, 136, 168, 196	0
1	J	469/500 (93%)	0.22	17 (3%) 42 28	92, 145, 182, 210	0
1	K	466/500 (93%)	-0.14	1 (0%) 95 91	64, 101, 143, 166	0
2	D	460/462 (99%)	-0.32	0 100 100	50, 79, 124, 153	0
2	E	460/462 (99%)	-0.05	5 (1%) 80 68	55, 96, 151, 200	0
2	F	462/462 (100%)	-0.28	0 100 100	53, 90, 124, 162	0
2	L	459/462 (99%)	-0.08	2 (0%) 92 86	66, 111, 148, 181	0
2	M	459/462 (99%)	0.62	53 (11%) 4 3	98, 172, 211, 247	0
2	N	460/462 (99%)	-0.17	2 (0%) 92 86	76, 107, 146, 168	0
3	G	281/282 (99%)	0.31	13 (4%) 32 20	54, 135, 185, 210	0
3	O	281/282 (99%)	0.10	7 (2%) 57 41	91, 130, 165, 188	0
4	H	134/134 (100%)	0.09	0 100 100	102, 131, 186, 214	0
4	P	134/134 (100%)	0.22	9 (6%) 17 10	77, 108, 194, 218	0
All	All	6420/6604 (97%)	-0.04	125 (1%) 66 51	50, 108, 176, 247	0

The worst 5 of 125 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	M	136	LEU	4.3
2	M	373	ASP	4.3
4	P	111	THR	4.1
2	M	432	LEU	3.9
1	J	100	GLY	3.7

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

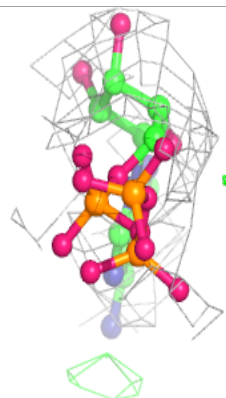
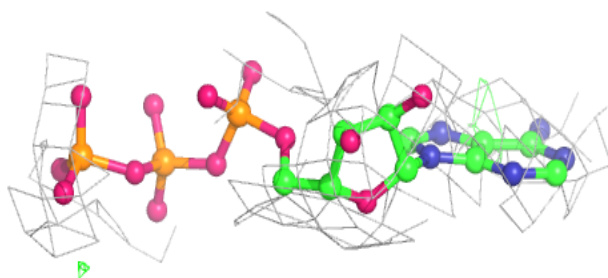
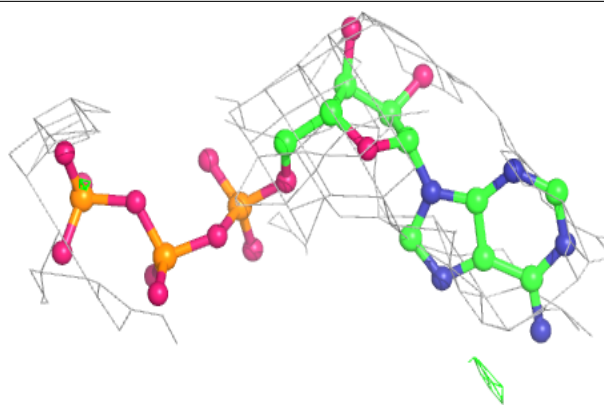
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	ATP	J	600	31/31	0.91	0.24	118,131,144,145	0
6	MG	D	601	1/1	0.93	0.17	66,66,66,66	0
6	MG	I	601	1/1	0.93	0.12	101,101,101,101	0
5	ATP	B	600	31/31	0.94	0.23	77,84,90,90	0
6	MG	N	601	1/1	0.94	0.15	75,75,75,75	0
7	ADP	N	600	27/27	0.94	0.21	76,87,92,93	0
5	ATP	K	600	31/31	0.95	0.24	63,66,69,71	0
6	MG	F	601	1/1	0.95	0.19	70,70,70,70	0
5	ATP	C	600	31/31	0.95	0.22	67,75,83,84	0
7	ADP	L	600	27/27	0.95	0.24	71,77,82,83	0
5	ATP	I	600	31/31	0.96	0.17	90,97,104,104	0
6	MG	A	601	1/1	0.96	0.18	56,56,56,56	0
7	ADP	D	600	27/27	0.96	0.21	69,71,75,77	0
6	MG	C	601	1/1	0.96	0.25	67,67,67,67	0
5	ATP	A	600	31/31	0.96	0.19	56,62,67,68	0
7	ADP	F	600	27/27	0.96	0.19	75,83,87,88	0
6	MG	B	601	1/1	0.97	0.17	82,82,82,82	0
6	MG	L	601	1/1	0.97	0.18	76,76,76,76	0
6	MG	K	601	1/1	0.97	0.23	68,68,68,68	0
6	MG	J	601	1/1	0.99	0.30	114,114,114,114	0

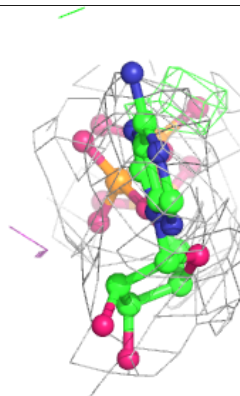
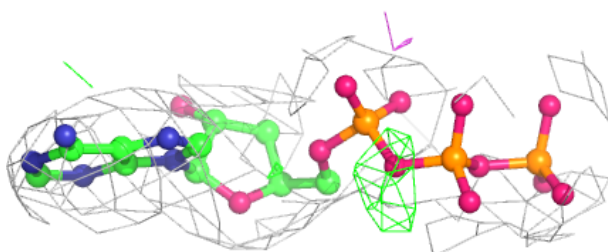
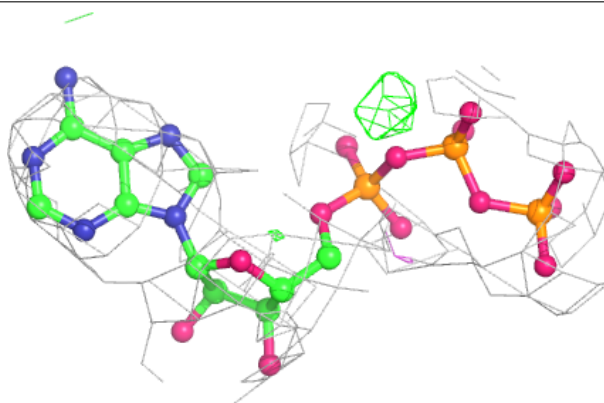
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around ATP J 600:**

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

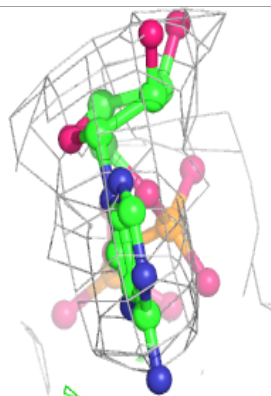
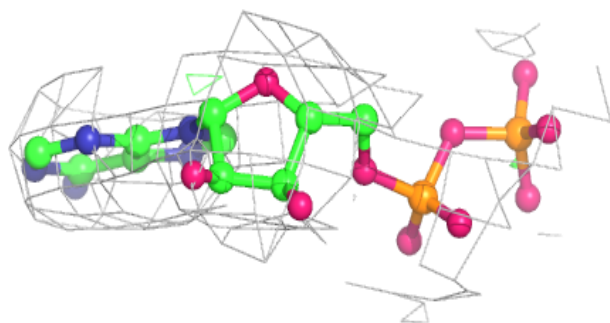
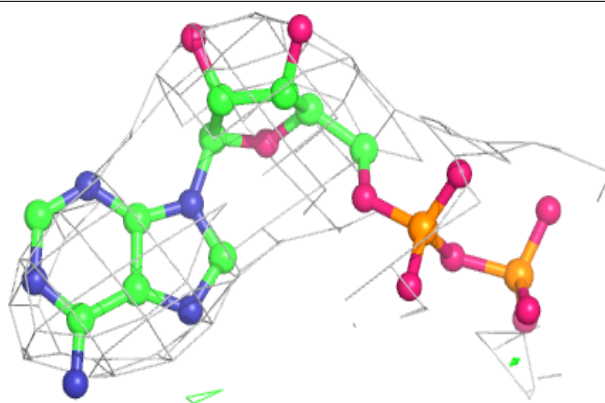
**Electron density around ATP B 600:**

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

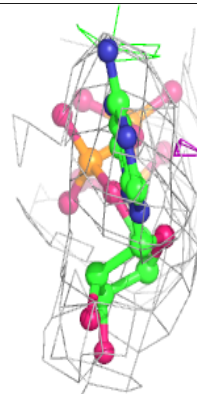
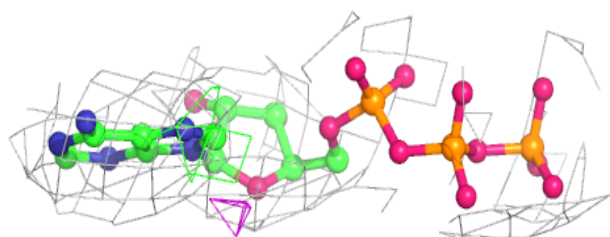
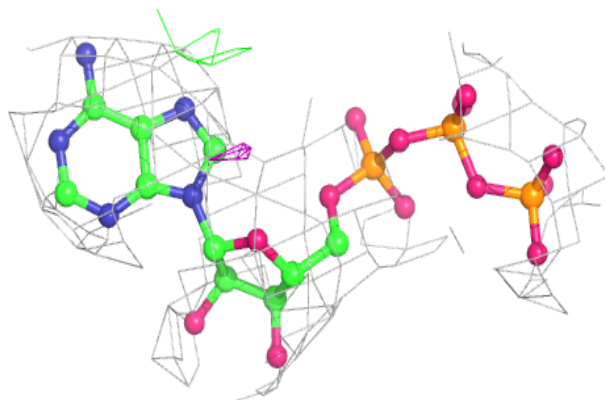


**Electron density around ADP N 600:**

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

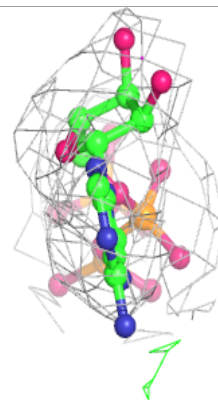
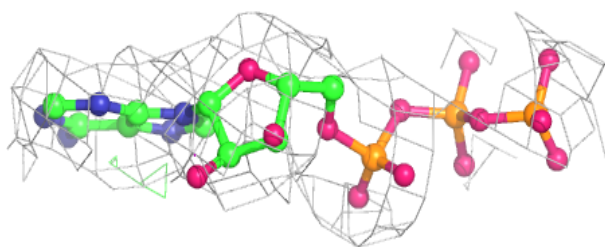
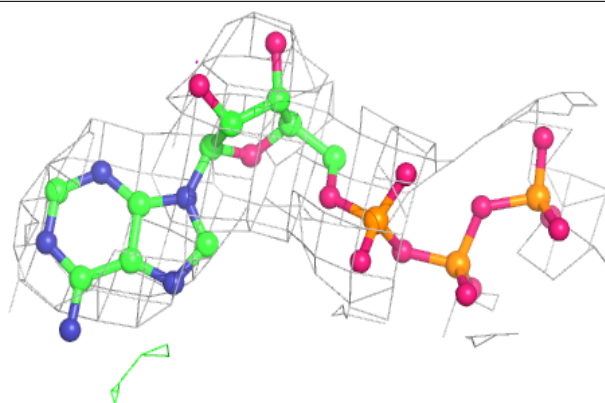
**Electron density around ATP K 600:**

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

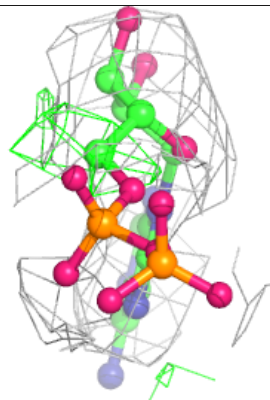
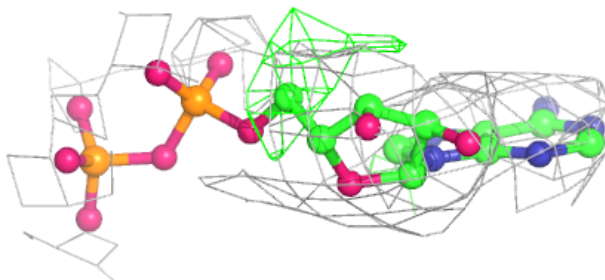
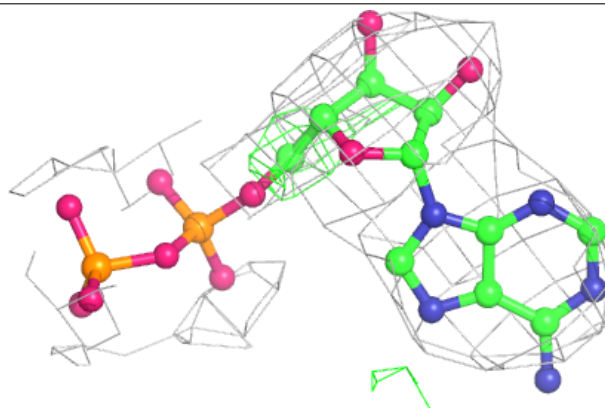


**Electron density around ATP C 600:**

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

**Electron density around ADP L 600:**

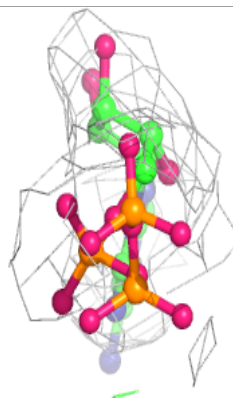
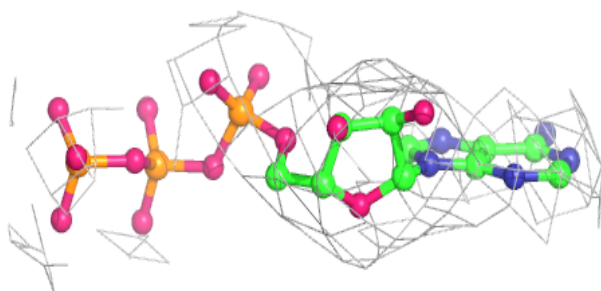
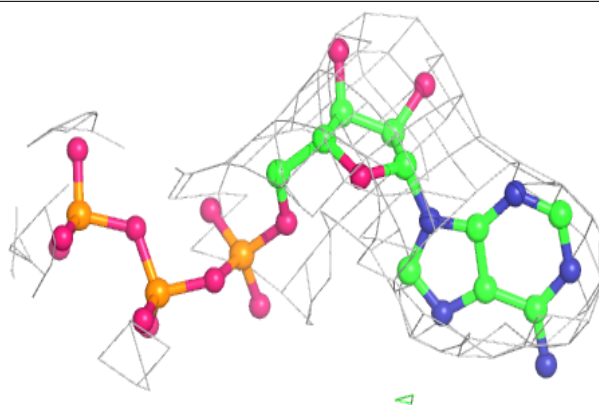
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



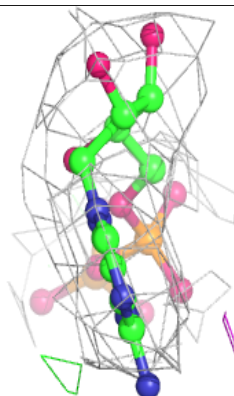
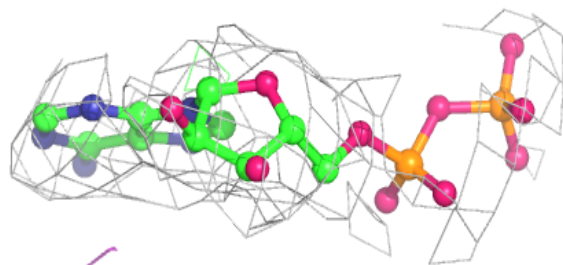
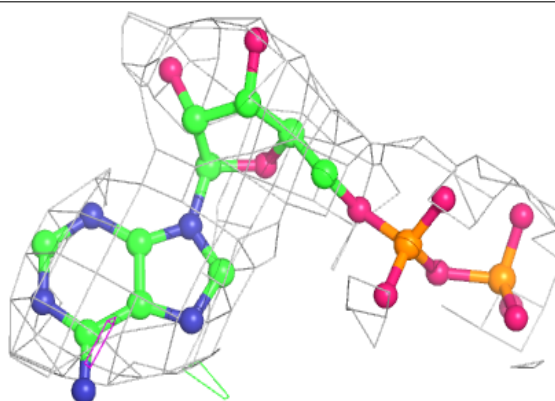


**Electron density around ATP I 600:**

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

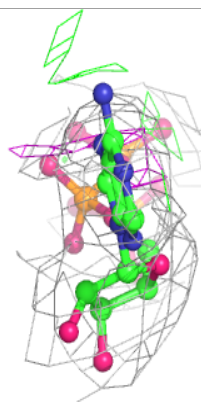
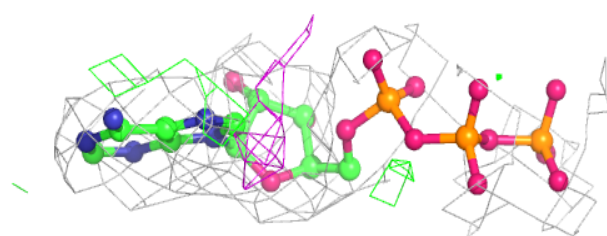
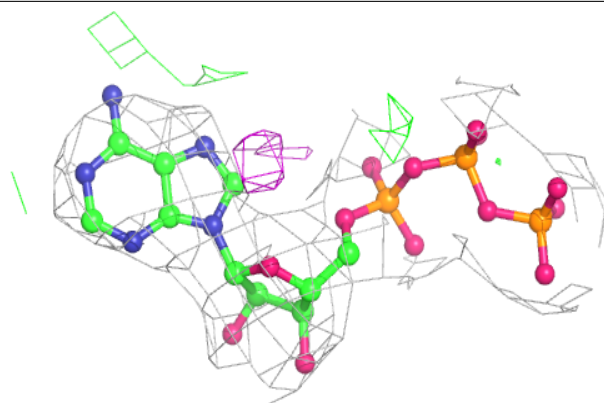
**Electron density around ADP D 600:**

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

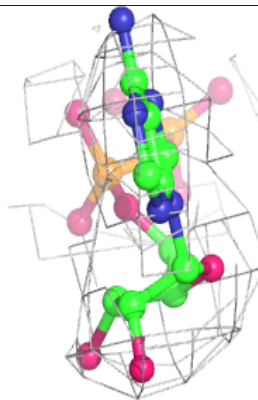
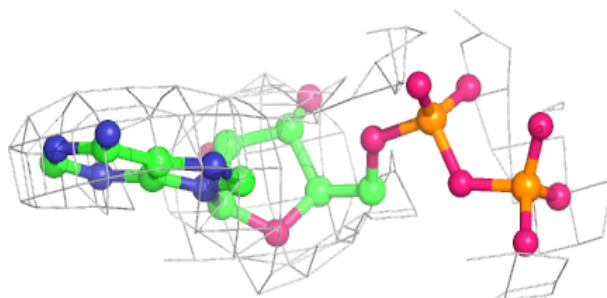
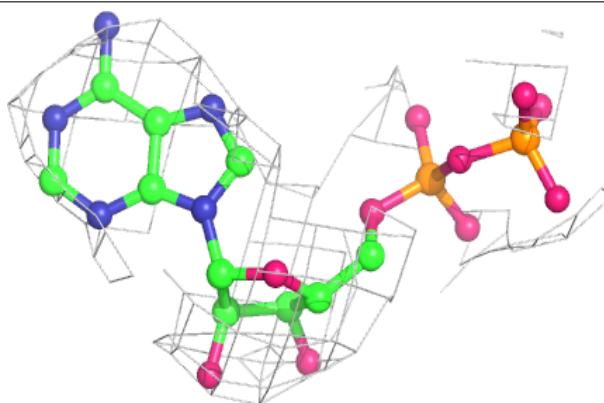


**Electron density around ATP A 600:**

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

**Electron density around ADP F 600:**

$2mF_o-DF_c$  (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.