



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 26, 2024 – 04:11 PM EDT

PDB ID : 9C6M
Title : Yasminevirus c12orf29, a 5' to 3' RNA ligase, K73M mutant
Authors : Hu, Y.; Lopez, V.A.; Tagliabracci, V.S.; Tomchick, D.R.
Deposited on : 2024-06-07
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure 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/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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.002 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.38.3

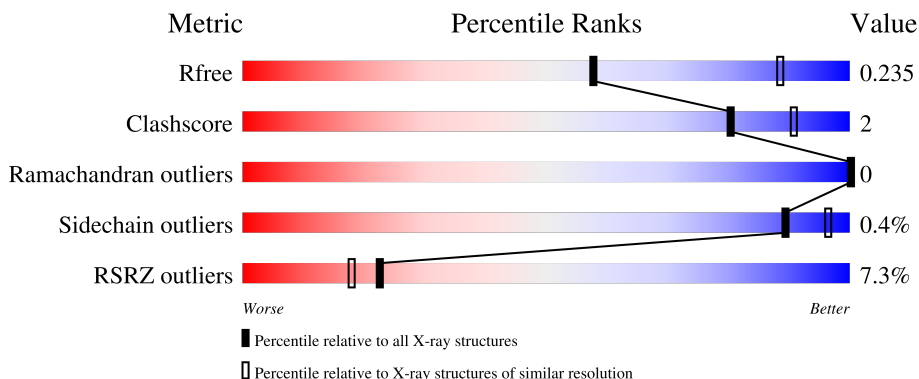
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	164625	3775 (2.60-2.60)
Clashscore	180529	4181 (2.60-2.60)
Ramachandran outliers	177936	4129 (2.60-2.60)
Sidechain outliers	177891	4129 (2.60-2.60)
RSRZ outliers	164620	3775 (2.60-2.60)

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	277	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 84%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">5% 84% 6% 10%</p>
1	B	277	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 83%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">7% 83% 6% 11%</p>
1	C	277	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 83%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">8% 83% 5% 12%</p>
1	D	277	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 79%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 14%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">6% 79% 8% 14%</p>
1	E	277	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 82%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 13%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">6% 82% 5% 13%</p>

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Mol	Chain	Length	Quality of chain
1	F	277	 <p>5% 81% 5% 14%</p>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 22888 atoms, of which 11302 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 RNA ligase1.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S	Se			
1	A	248	3846	1213	1914	347	358	8	6	0	0	0
1	B	246	3803	1200	1890	344	355	8	6	0	0	0
1	C	243	3779	1193	1883	342	347	8	6	0	0	0
1	D	239	3706	1173	1840	333	346	8	6	0	0	0
1	E	241	3736	1182	1858	336	346	8	6	0	0	0
1	F	237	3703	1170	1845	335	339	8	6	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP A0A5K0UB63
A	73	MSE	LYS	engineered mutation	UNP A0A5K0UB63
B	0	SER	-	expression tag	UNP A0A5K0UB63
B	73	MSE	LYS	engineered mutation	UNP A0A5K0UB63
C	0	SER	-	expression tag	UNP A0A5K0UB63
C	73	MSE	LYS	engineered mutation	UNP A0A5K0UB63
D	0	SER	-	expression tag	UNP A0A5K0UB63
D	73	MSE	LYS	engineered mutation	UNP A0A5K0UB63
E	0	SER	-	expression tag	UNP A0A5K0UB63
E	73	MSE	LYS	engineered mutation	UNP A0A5K0UB63
F	0	SER	-	expression tag	UNP A0A5K0UB63
F	73	MSE	LYS	engineered mutation	UNP A0A5K0UB63

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
2	A	1	43	10	12	5	13	3	0	0
2	B	1	43	10	12	5	13	3	0	0
2	C	1	43	10	12	5	13	3	0	0
2	D	1	43	10	12	5	13	3	0	0
2	E	1	43	10	12	5	13	3	0	0
2	F	1	43	10	12	5	13	3	0	0

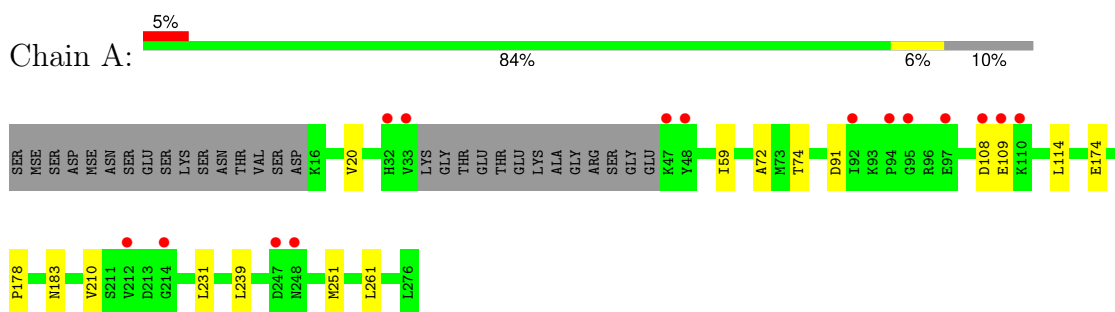
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	12	Total	O	0	0
			12	12		
3	C	6	Total	O	0	0
			6	6		
3	D	9	Total	O	0	0
			9	9		
3	E	9	Total	O	0	0
			9	9		
3	F	10	Total	O	0	0
			10	10		

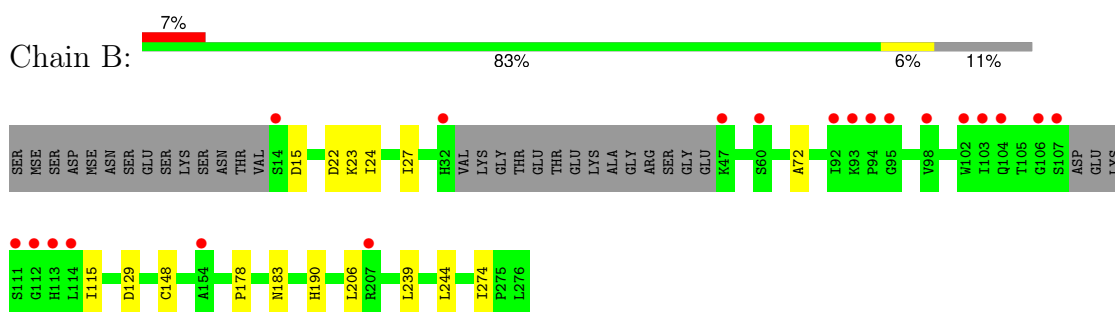
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and 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.

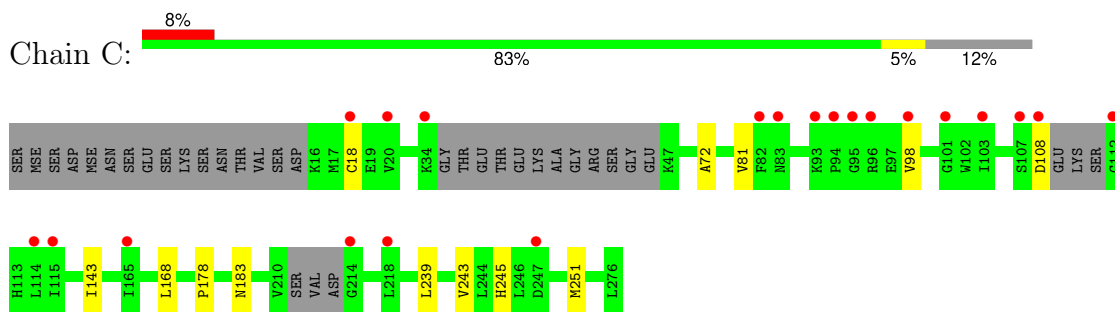
- Molecule 1: RNA ligase1



- Molecule 1: RNA ligase1

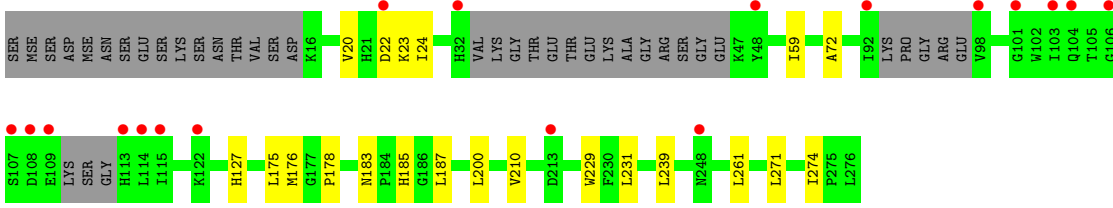


- Molecule 1: RNA ligase1

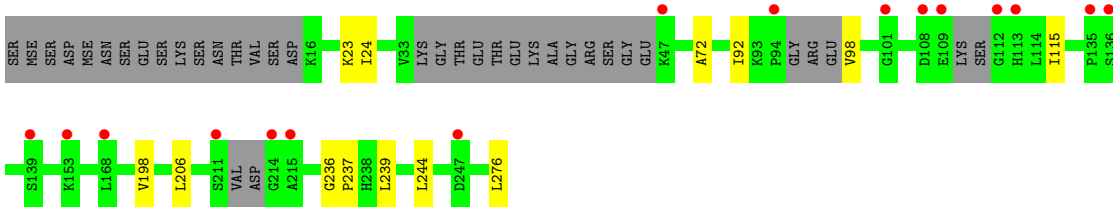
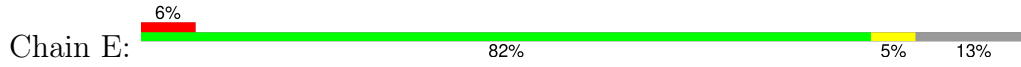


- Molecule 1: RNA ligase1

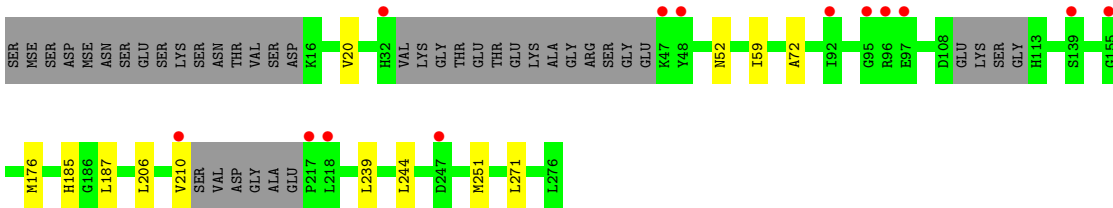
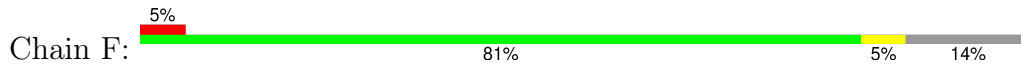




• Molecule 1: RNA ligase1



• Molecule 1: RNA ligase1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	103.40Å 86.05Å 114.31Å 90.00° 105.19° 90.00°	Depositor
Resolution (Å)	43.70 – 2.60 43.70 – 2.60	Depositor EDS
% Data completeness (in resolution range)	75.3 (43.70-2.60) 84.1 (43.70-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.54Å)	Xtrriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.195 , 0.235 0.194 , 0.235	Depositor DCC
R_{free} test set	57842 reflections (3.87%)	wwPDB-VP
Wilson B-factor (Å ²)	39.8	Xtrriage
Anisotropy	0.129	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	22888	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.47% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.25	0/1969	0.48	0/2653
1	B	0.26	0/1949	0.48	0/2625
1	C	0.25	0/1931	0.48	0/2598
1	D	0.25	0/1900	0.47	0/2559
1	E	0.25	0/1912	0.48	0/2573
1	F	0.25	0/1893	0.48	0/2547
All	All	0.25	0/11554	0.48	0/15555

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1932	1914	1914	8	0
1	B	1913	1890	1890	8	0
1	C	1896	1883	1883	8	0
1	D	1866	1840	1840	15	0
1	E	1878	1858	1858	8	0
1	F	1858	1845	1845	9	0
2	A	31	12	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	31	12	12	1	0
2	C	31	12	12	1	0
2	D	31	12	12	2	0
2	E	31	12	12	1	0
2	F	31	12	12	0	0
3	A	11	0	0	0	0
3	B	12	0	0	0	0
3	C	6	0	0	0	0
3	D	9	0	0	0	0
3	E	9	0	0	0	0
3	F	10	0	0	0	0
All	All	11586	11302	11302	56	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (56) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:176:MSE:HE2	1:D:187:LEU:HD11	1.41	1.02
1:F:176:MSE:HE2	1:F:187:LEU:HD11	1.47	0.97
1:C:72:ALA:HB1	1:C:239:LEU:HD22	1.76	0.67
1:D:72:ALA:HB1	1:D:239:LEU:HD22	1.79	0.65
1:D:176:MSE:HE1	1:D:271:LEU:CD2	2.27	0.65
1:D:59:ILE:HD13	1:D:210:VAL:HG22	1.80	0.64
1:F:176:MSE:HE1	1:F:271:LEU:CD2	2.28	0.63
1:D:20:VAL:HG11	2:D:301:ATP:H3'	1.81	0.63
1:F:72:ALA:HB1	1:F:239:LEU:HD22	1.81	0.61
1:B:72:ALA:HB1	1:B:239:LEU:HD22	1.83	0.61
1:E:72:ALA:HB1	1:E:239:LEU:HD22	1.83	0.61
1:D:176:MSE:HE3	1:D:185:HIS:CD2	2.37	0.59
1:F:176:MSE:HE3	1:F:185:HIS:CD2	2.39	0.58
1:B:239:LEU:HD21	1:B:274:ILE:HD12	1.84	0.58
1:A:59:ILE:HD13	1:A:210:VAL:HG22	1.84	0.57
1:E:24:ILE:HD12	2:E:301:ATP:H5'1	1.86	0.57
1:D:231:LEU:HD21	1:D:261:LEU:HD22	1.86	0.56
1:B:23:LYS:HG2	1:B:115:ILE:HD11	1.88	0.56
1:A:74:THR:HG23	1:A:174:GLU:HG2	1.88	0.55
1:A:72:ALA:HB1	1:A:239:LEU:HD22	1.89	0.54
1:C:245:HIS:NE2	1:C:251:MSE:HE2	2.22	0.54
1:B:178:PRO:HA	1:B:183:ASN:HB2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:176:MSE:HE1	1:D:271:LEU:HD22	1.90	0.54
1:F:20:VAL:HA	1:F:251:MSE:HE1	1.89	0.54
1:B:206:LEU:HD12	1:B:244:LEU:HD11	1.91	0.52
1:C:18:CYS:HB2	1:C:251:MSE:HE3	1.92	0.52
1:F:176:MSE:HE1	1:F:271:LEU:HD23	1.93	0.51
1:D:200:LEU:HD13	1:D:229:TRP:CH2	2.46	0.51
1:D:178:PRO:HA	1:D:183:ASN:HB2	1.94	0.50
1:F:206:LEU:HD13	1:F:244:LEU:HD11	1.96	0.48
1:E:206:LEU:CD1	1:E:244:LEU:HD11	2.43	0.48
1:C:178:PRO:HA	1:C:183:ASN:HB2	1.96	0.48
1:D:271:LEU:O	1:D:274:ILE:HG22	2.14	0.47
1:B:24:ILE:HD13	2:B:301:ATP:H5'1	1.95	0.47
1:E:198:VAL:HG13	1:E:276:LEU:CD2	2.45	0.47
1:D:176:MSE:HE1	1:D:271:LEU:HD23	1.97	0.47
1:A:178:PRO:HA	1:A:183:ASN:HB2	1.97	0.46
1:E:92:ILE:HD13	1:E:98:VAL:HG12	1.98	0.46
1:A:231:LEU:HD21	1:A:261:LEU:HD22	1.97	0.46
1:D:24:ILE:HD12	2:D:301:ATP:H5'2	1.97	0.45
1:F:176:MSE:HE1	1:F:271:LEU:HD22	1.96	0.45
1:B:129:ASP:OD2	1:B:190:HIS:ND1	2.42	0.45
1:C:143:ILE:HG22	1:C:143:ILE:O	2.17	0.45
1:C:98:VAL:O	1:C:98:VAL:HG13	2.16	0.45
1:C:243:VAL:HG11	2:C:301:ATP:C2	2.52	0.44
1:B:27:ILE:HD11	1:B:244:LEU:HG	1.99	0.44
1:D:127:HIS:HA	1:D:175:LEU:HD21	2.00	0.44
1:D:22:ASP:OD1	1:D:23:LYS:N	2.47	0.44
1:A:91:ASP:HA	1:A:114:LEU:O	2.18	0.44
1:A:20:VAL:HA	1:A:251:MSE:HE1	1.99	0.44
1:F:59:ILE:HG23	1:F:210:VAL:HG13	2.00	0.44
1:E:23:LYS:HG2	1:E:115:ILE:HD11	1.99	0.43
1:E:198:VAL:HG13	1:E:276:LEU:HD21	2.01	0.43
1:C:81:VAL:HG21	1:C:168:LEU:HB2	2.01	0.42
1:E:236:GLY:N	1:E:237:PRO:HD2	2.35	0.41
1:A:108:ASP:OD1	1:A:109:GLU:N	2.54	0.41

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 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	244/277 (88%)	240 (98%)	4 (2%)	0	100	100
1	B	240/277 (87%)	231 (96%)	9 (4%)	0	100	100
1	C	235/277 (85%)	226 (96%)	9 (4%)	0	100	100
1	D	231/277 (83%)	226 (98%)	5 (2%)	0	100	100
1	E	231/277 (83%)	224 (97%)	7 (3%)	0	100	100
1	F	229/277 (83%)	225 (98%)	4 (2%)	0	100	100
All	All	1410/1662 (85%)	1372 (97%)	38 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	212/229 (93%)	212 (100%)	0	100	100
1	B	210/229 (92%)	207 (99%)	3 (1%)	62	82
1	C	207/229 (90%)	206 (100%)	1 (0%)	86	95
1	D	205/229 (90%)	205 (100%)	0	100	100
1	E	206/229 (90%)	206 (100%)	0	100	100
1	F	204/229 (89%)	203 (100%)	1 (0%)	86	95
All	All	1244/1374 (90%)	1239 (100%)	5 (0%)	89	96

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	15	ASP
1	B	22	ASP
1	B	148	CYS
1	C	108	ASP
1	F	52	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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	$\# Z > 2$
2	ATP	B	301	-	28,33,33	0.62	0	34,52,52	1.02	3 (8%)
2	ATP	E	301	-	28,33,33	0.60	0	34,52,52	1.05	2 (5%)
2	ATP	C	301	-	28,33,33	0.69	0	34,52,52	0.96	1 (2%)
2	ATP	A	301	-	28,33,33	0.62	0	34,52,52	1.08	4 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ATP	D	301	-	28,33,33	0.61	0	34,52,52	1.12	3 (8%)
2	ATP	F	301	-	28,33,33	0.62	0	34,52,52	1.06	2 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ATP	B	301	-	-	7/18/38/38	0/3/3/3
2	ATP	E	301	-	-	8/18/38/38	0/3/3/3
2	ATP	C	301	-	-	7/18/38/38	0/3/3/3
2	ATP	A	301	-	-	5/18/38/38	0/3/3/3
2	ATP	D	301	-	-	7/18/38/38	0/3/3/3
2	ATP	F	301	-	-	8/18/38/38	0/3/3/3

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	ATP	C4'-O4'-C1'	-3.82	106.42	109.92
2	F	301	ATP	C4'-O4'-C1'	-2.88	107.29	109.92
2	A	301	ATP	C4'-O4'-C1'	-2.82	107.34	109.92
2	E	301	ATP	C4'-O4'-C1'	-2.49	107.65	109.92
2	A	301	ATP	O4'-C1'-N9	-2.43	105.52	108.75
2	A	301	ATP	C5-C6-N6	2.36	123.91	120.31
2	B	301	ATP	C5-C6-N6	2.35	123.90	120.31
2	E	301	ATP	C5-C6-N6	2.35	123.89	120.31
2	F	301	ATP	C5-C6-N6	2.35	123.89	120.31
2	D	301	ATP	C5-C6-N6	2.35	123.89	120.31
2	C	301	ATP	C5-C6-N6	2.27	123.77	120.31
2	B	301	ATP	O4'-C1'-N9	-2.15	105.89	108.75
2	D	301	ATP	O3'-C3'-C4'	-2.06	105.16	111.08
2	B	301	ATP	O3'-C3'-C4'	-2.02	105.28	111.08
2	A	301	ATP	O3'-C3'-C4'	-2.01	105.32	111.08

There are no chirality outliers.

All (42) torsion outliers are listed below:

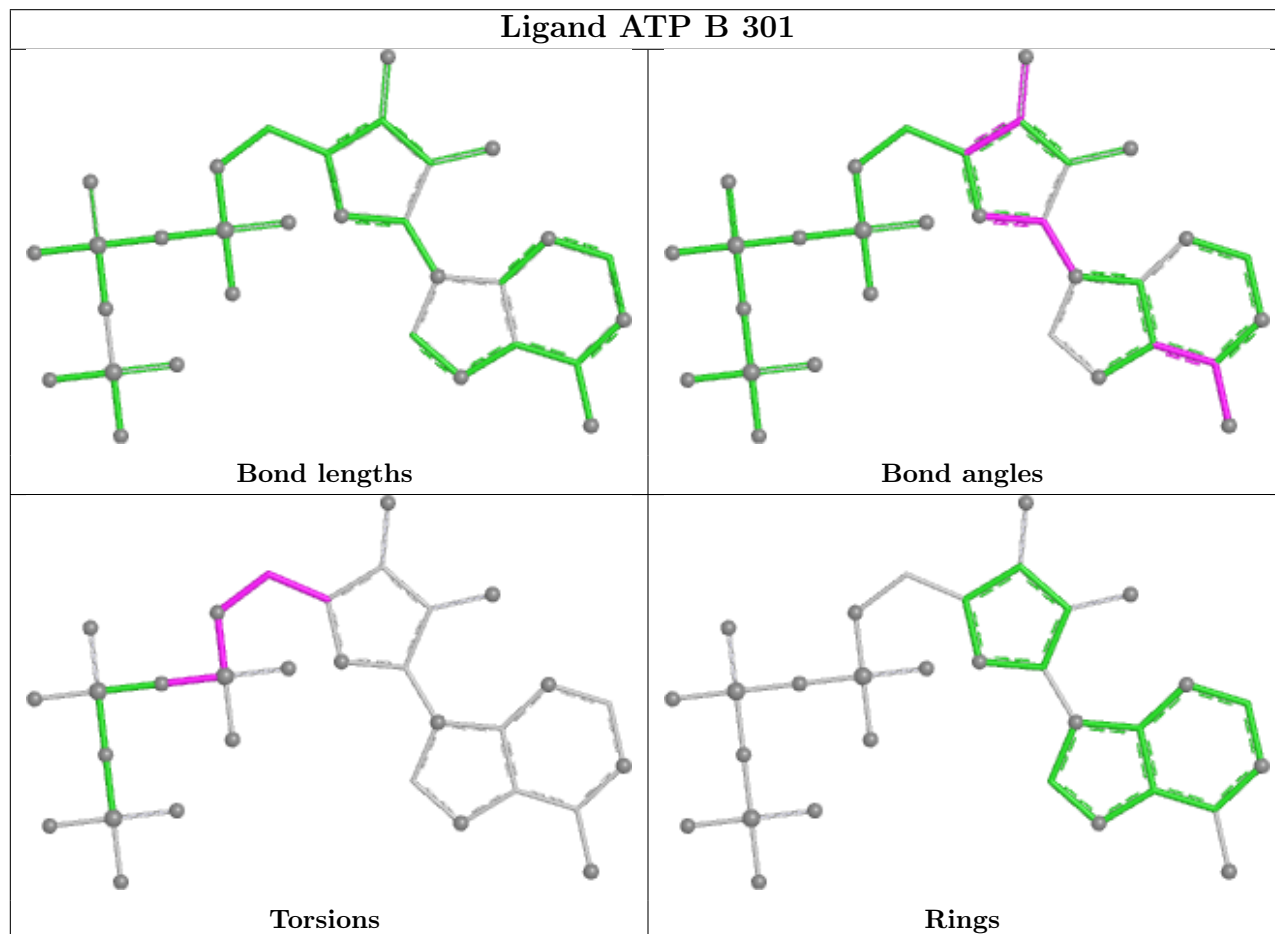
Mol	Chain	Res	Type	Atoms
2	A	301	ATP	PB-O3A-PA-O5'
2	B	301	ATP	C5'-O5'-PA-O2A
2	B	301	ATP	C5'-O5'-PA-O3A
2	C	301	ATP	C5'-O5'-PA-O2A
2	C	301	ATP	C5'-O5'-PA-O3A
2	D	301	ATP	C5'-O5'-PA-O2A
2	D	301	ATP	C5'-O5'-PA-O3A
2	E	301	ATP	PB-O3A-PA-O5'
2	E	301	ATP	C5'-O5'-PA-O2A
2	E	301	ATP	C5'-O5'-PA-O3A
2	F	301	ATP	C5'-O5'-PA-O2A
2	F	301	ATP	C5'-O5'-PA-O3A
2	F	301	ATP	O4'-C4'-C5'-O5'
2	F	301	ATP	C3'-C4'-C5'-O5'
2	A	301	ATP	O4'-C4'-C5'-O5'
2	A	301	ATP	C3'-C4'-C5'-O5'
2	B	301	ATP	C3'-C4'-C5'-O5'
2	D	301	ATP	C3'-C4'-C5'-O5'
2	E	301	ATP	O4'-C4'-C5'-O5'
2	E	301	ATP	C3'-C4'-C5'-O5'
2	B	301	ATP	O4'-C4'-C5'-O5'
2	C	301	ATP	O4'-C4'-C5'-O5'
2	C	301	ATP	C3'-C4'-C5'-O5'
2	D	301	ATP	O4'-C4'-C5'-O5'
2	B	301	ATP	PB-O3A-PA-O5'
2	C	301	ATP	PB-O3A-PA-O5'
2	F	301	ATP	PB-O3A-PA-O5'
2	C	301	ATP	C5'-O5'-PA-O1A
2	B	301	ATP	C4'-C5'-O5'-PA
2	F	301	ATP	C4'-C5'-O5'-PA
2	D	301	ATP	PB-O3A-PA-O5'
2	C	301	ATP	C4'-C5'-O5'-PA
2	E	301	ATP	C4'-C5'-O5'-PA
2	B	301	ATP	PB-O3A-PA-O2A
2	E	301	ATP	PB-O3A-PA-O1A
2	E	301	ATP	PB-O3A-PA-O2A
2	F	301	ATP	PB-O3A-PA-O1A
2	D	301	ATP	C4'-C5'-O5'-PA
2	A	301	ATP	PB-O3A-PA-O1A
2	A	301	ATP	PB-O3A-PA-O2A
2	D	301	ATP	PB-O3A-PA-O2A
2	F	301	ATP	PB-O3A-PA-O2A

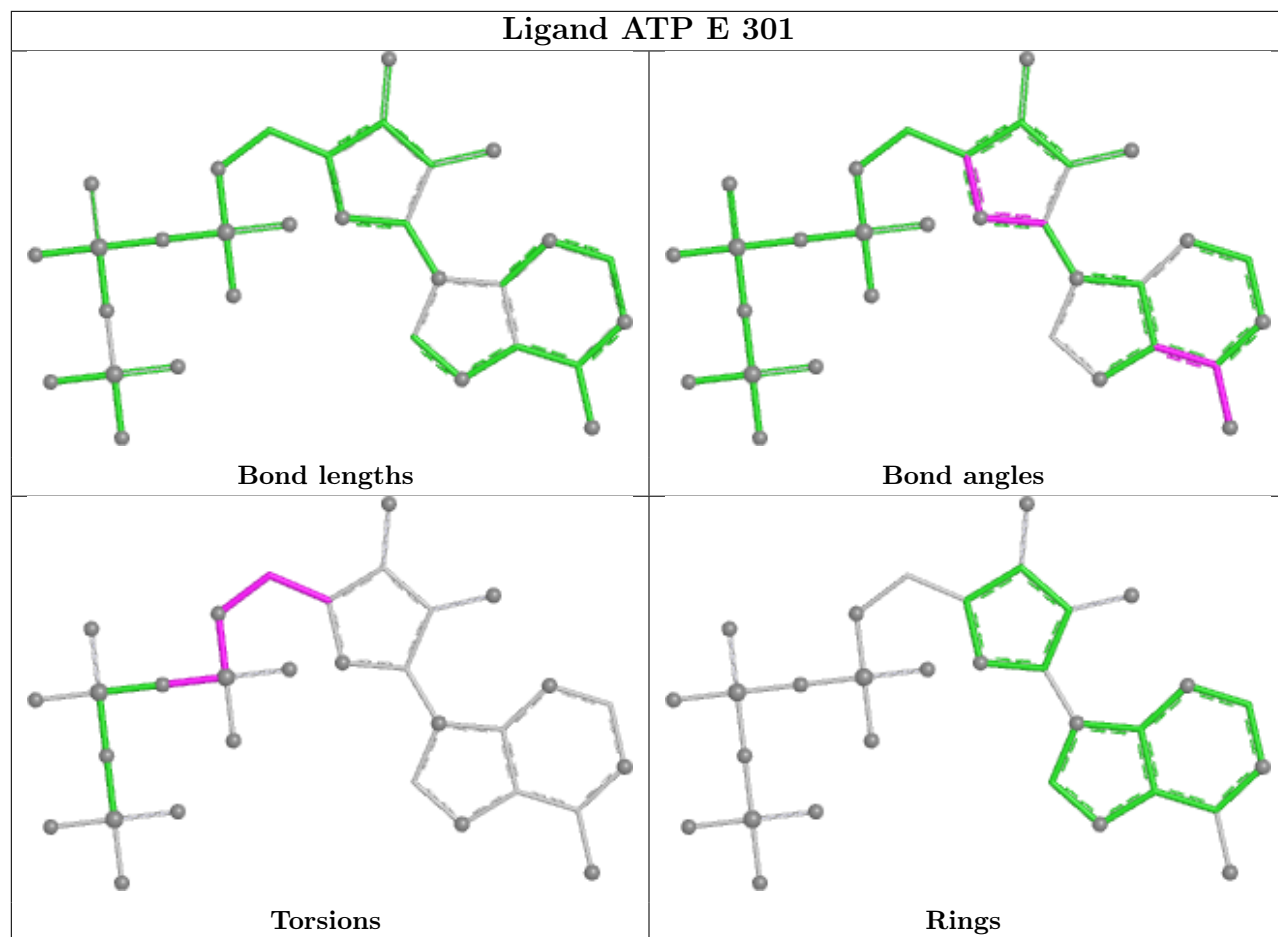
There are no ring outliers.

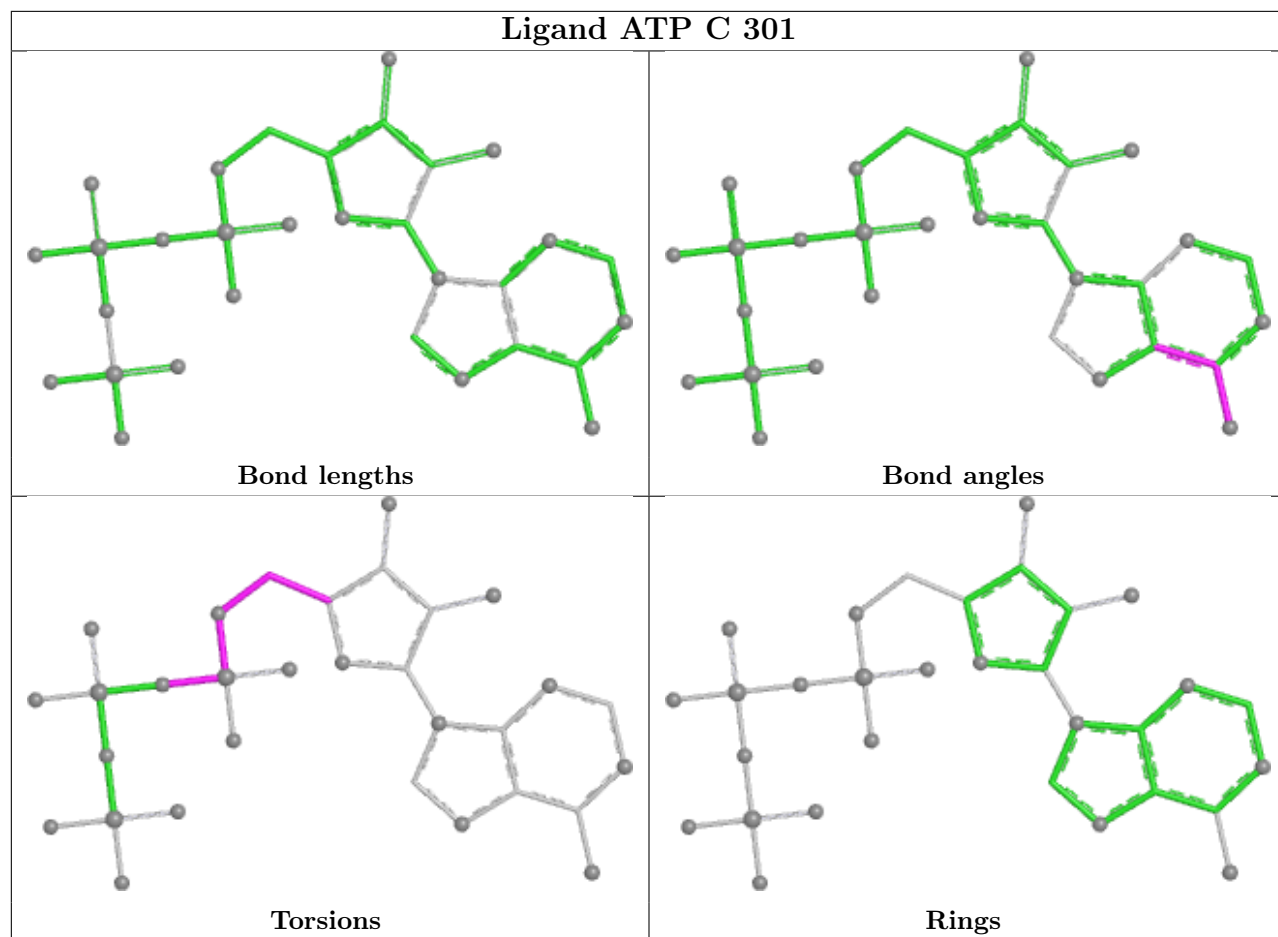
4 monomers are involved in 5 short contacts:

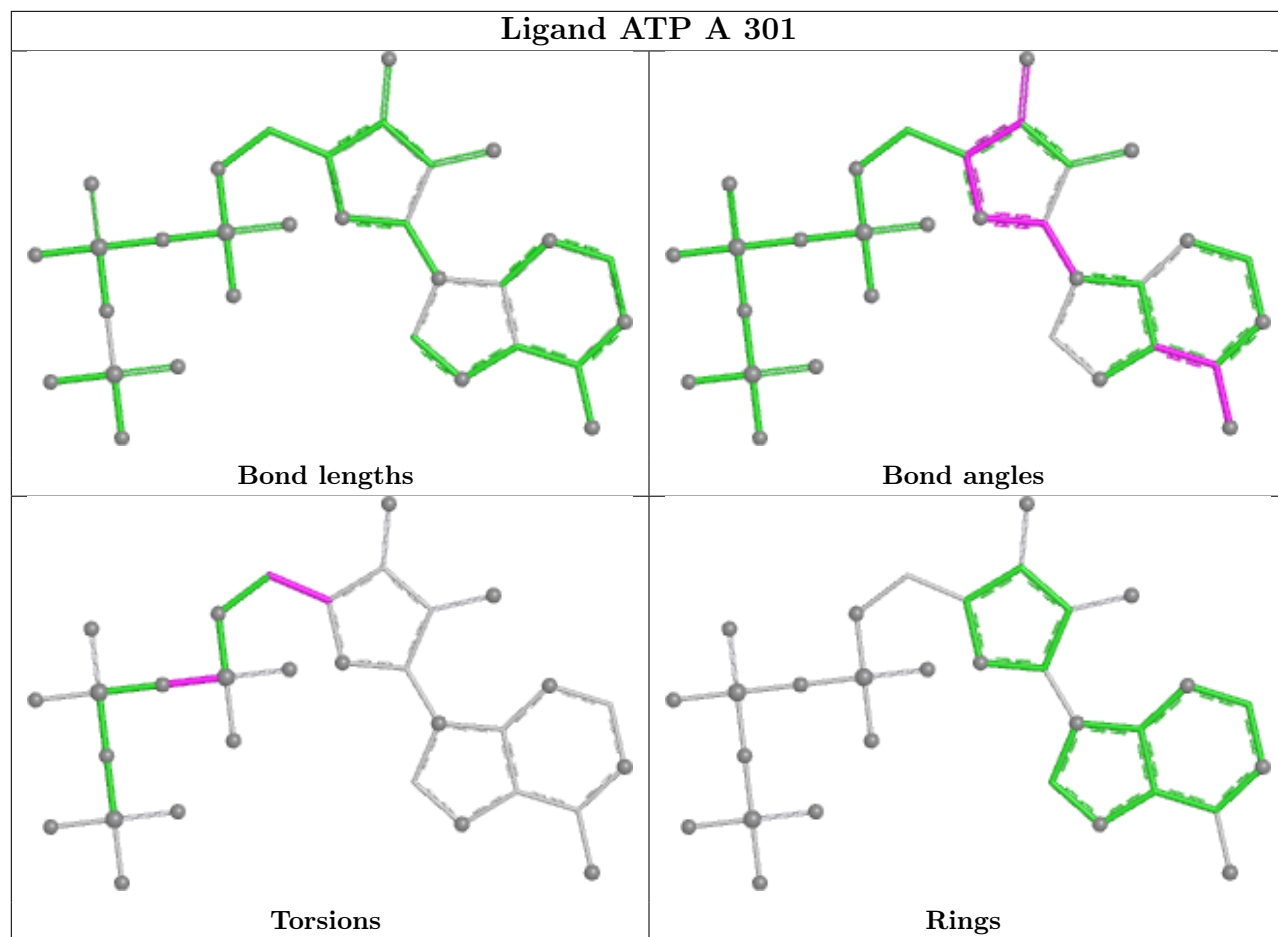
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	ATP	1	0
2	E	301	ATP	1	0
2	C	301	ATP	1	0
2	D	301	ATP	2	0

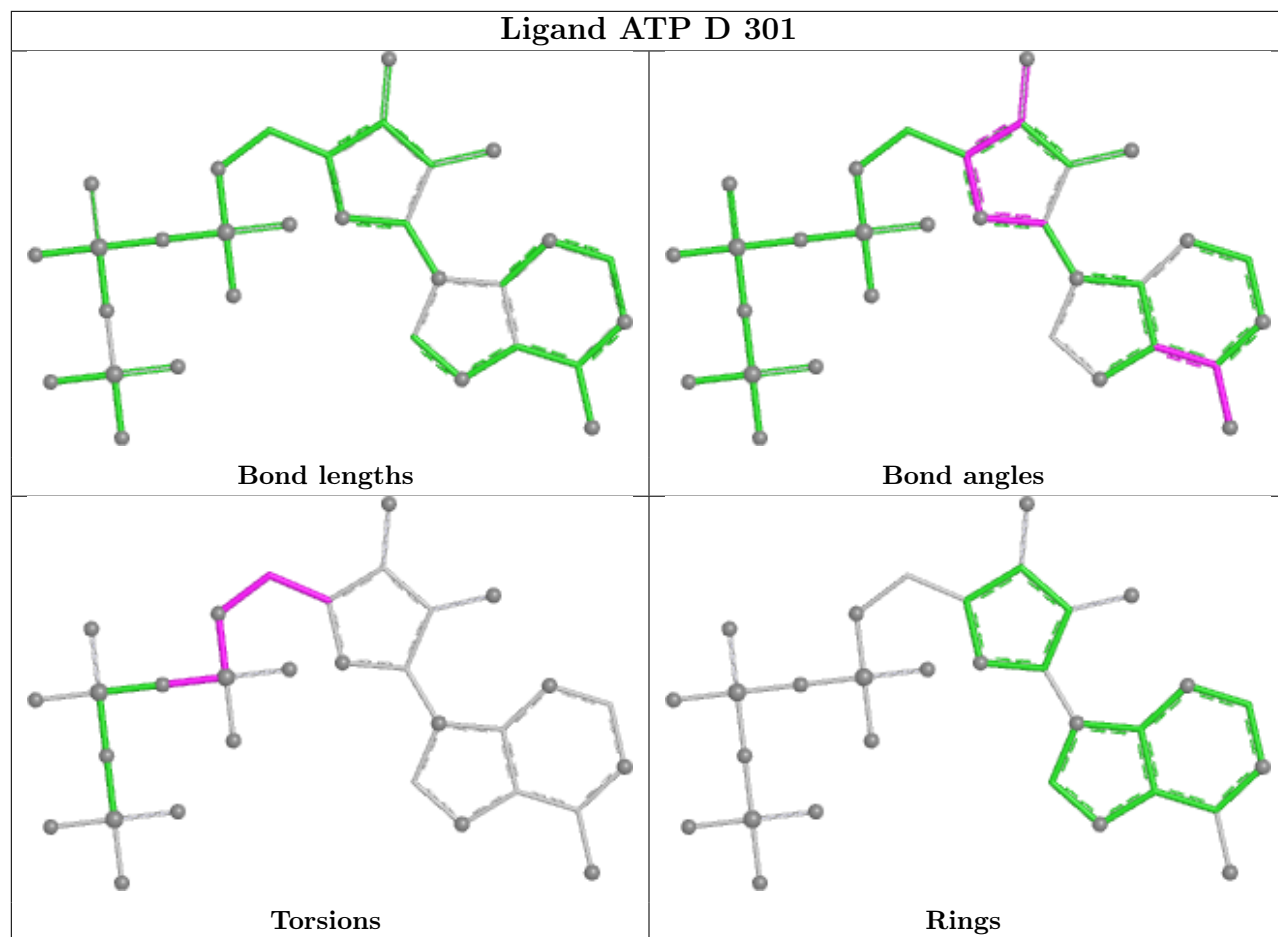
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

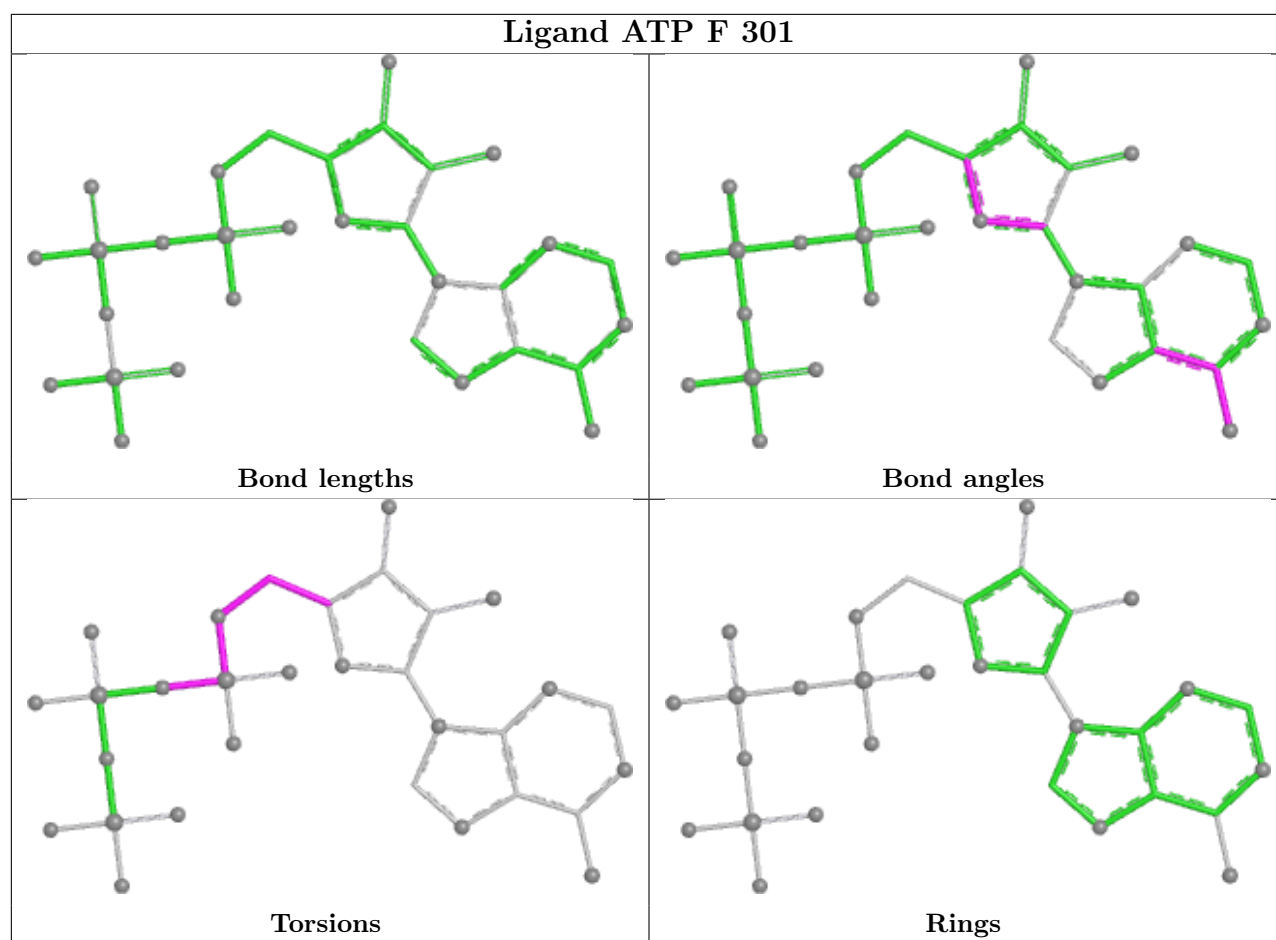












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 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, 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	242/277 (87%)	0.13	15 (6%) 28 22	21, 43, 89, 115	0
1	B	240/277 (86%)	0.39	20 (8%) 19 15	21, 47, 92, 116	0
1	C	237/277 (85%)	0.65	21 (8%) 17 14	26, 57, 95, 110	0
1	D	233/277 (84%)	0.25	18 (7%) 21 17	23, 46, 79, 112	0
1	E	235/277 (84%)	0.29	16 (6%) 25 20	22, 48, 86, 119	0
1	F	231/277 (83%)	0.35	13 (5%) 31 25	22, 48, 77, 107	0
All	All	1418/1662 (85%)	0.34	103 (7%) 22 18	21, 48, 88, 119	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	214	GLY	5.1
1	E	112	GLY	4.8
1	E	215	ALA	4.8
1	C	112	GLY	4.7
1	D	113	HIS	4.5
1	B	113	HIS	4.3
1	D	114	LEU	4.0
1	C	107	SER	3.9
1	E	94	PRO	3.9
1	D	32	HIS	3.9
1	A	212	VAL	3.8
1	D	107	SER	3.7
1	B	47	LYS	3.6
1	A	33	VAL	3.6
1	B	32	HIS	3.4
1	D	109	GLU	3.3
1	B	111	SER	3.3
1	C	165	ILE	3.3
1	D	106	GLY	3.3

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Mol	Chain	Res	Type	RSRZ
1	E	211	SER	3.2
1	F	32	HIS	3.2
1	E	247	ASP	3.2
1	B	107	SER	3.2
1	E	139	SER	3.1
1	F	155	GLY	3.1
1	A	94	PRO	3.1
1	F	48	TYR	3.1
1	F	47	LYS	3.1
1	C	94	PRO	3.0
1	B	114	LEU	3.0
1	D	108	ASP	3.0
1	C	95	GLY	3.0
1	C	218	LEU	2.9
1	B	103	ILE	2.9
1	B	93	LYS	2.9
1	B	112	GLY	2.9
1	D	104	GLN	2.9
1	D	115	ILE	2.8
1	E	109	GLU	2.8
1	F	210	VAL	2.8
1	C	34	LYS	2.8
1	C	214	GLY	2.8
1	B	104	GLN	2.7
1	C	96	ARG	2.7
1	E	153	LYS	2.6
1	B	94	PRO	2.6
1	C	98	VAL	2.6
1	F	217	PRO	2.6
1	C	82	PHE	2.6
1	C	115	ILE	2.6
1	F	139	SER	2.5
1	C	83	ASN	2.5
1	D	98	VAL	2.5
1	E	135	PRO	2.5
1	A	247	ASP	2.5
1	F	96	ARG	2.5
1	D	101	GLY	2.5
1	C	108	ASP	2.5
1	F	97	GLU	2.5
1	B	60	SER	2.4
1	C	247	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	103	ILE	2.4
1	A	214	GLY	2.4
1	B	14	SER	2.4
1	B	106	GLY	2.4
1	E	108	ASP	2.4
1	A	110	LYS	2.3
1	B	207	ARG	2.3
1	C	20	VAL	2.3
1	E	168	LEU	2.3
1	B	92	ILE	2.3
1	E	101	GLY	2.3
1	A	47	LYS	2.3
1	A	248	ASN	2.2
1	B	102	TRP	2.2
1	C	101	GLY	2.2
1	A	92	ILE	2.2
1	A	97	GLU	2.2
1	A	108	ASP	2.2
1	A	109	GLU	2.2
1	A	48	TYR	2.2
1	C	114	LEU	2.2
1	C	93	LYS	2.1
1	D	122	LYS	2.1
1	D	92	ILE	2.1
1	D	213	ASP	2.1
1	F	92	ILE	2.1
1	E	47	LYS	2.1
1	D	103	ILE	2.1
1	D	248	ASN	2.1
1	B	154	ALA	2.1
1	B	95	GLY	2.1
1	F	95	GLY	2.1
1	A	32	HIS	2.1
1	F	218	LEU	2.0
1	F	247	ASP	2.0
1	A	95	GLY	2.0
1	E	113	HIS	2.0
1	C	18	CYS	2.0
1	B	98	VAL	2.0
1	E	136	SER	2.0
1	D	22	ASP	2.0
1	D	48	TYR	2.0

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 [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

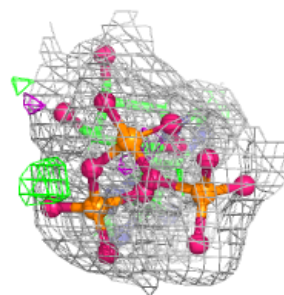
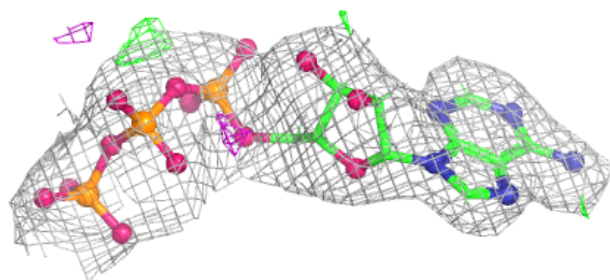
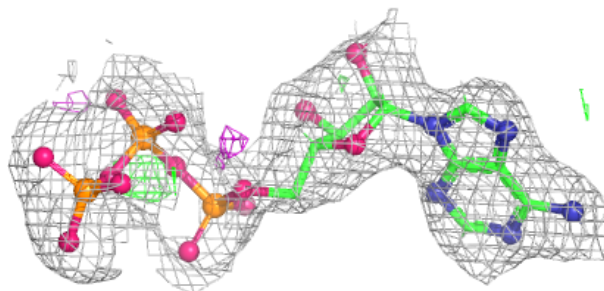
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ATP	C	301	31/31	0.85	0.12	41,56,86,90	0
2	ATP	E	301	31/31	0.85	0.13	33,52,98,110	0
2	ATP	D	301	31/31	0.88	0.09	31,43,87,94	0
2	ATP	B	301	31/31	0.89	0.11	34,50,89,104	0
2	ATP	F	301	31/31	0.90	0.10	32,49,85,87	0
2	ATP	A	301	31/31	0.92	0.09	22,39,63,69	0

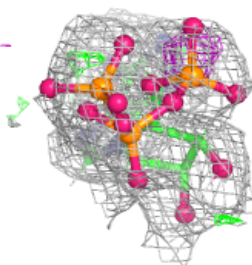
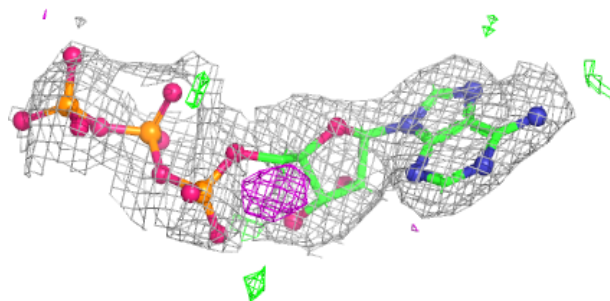
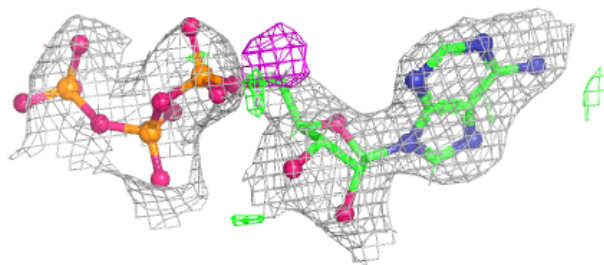
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP C 301:

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

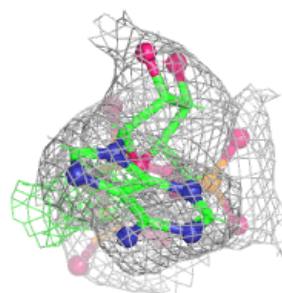
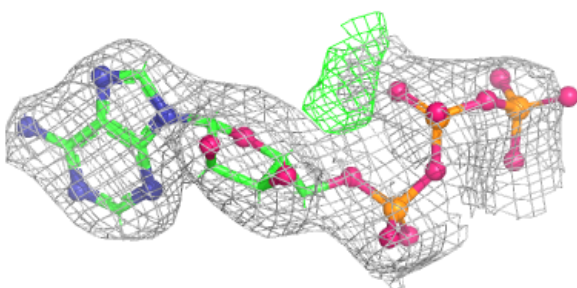
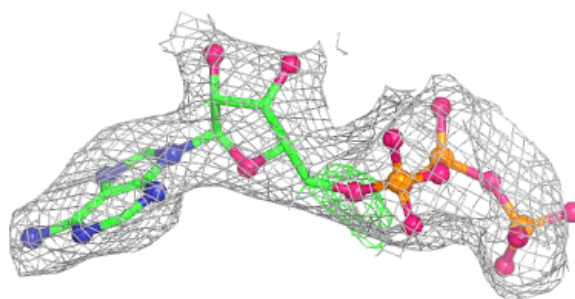
**Electron density around ATP E 301:**

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

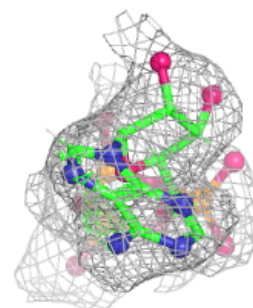
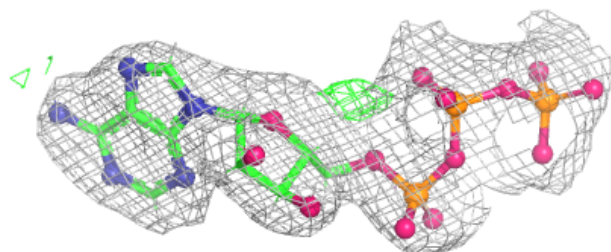
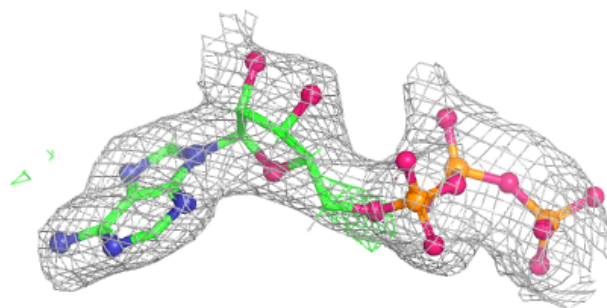


Electron density around ATP D 301:

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

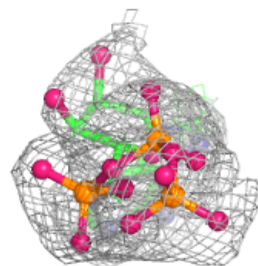
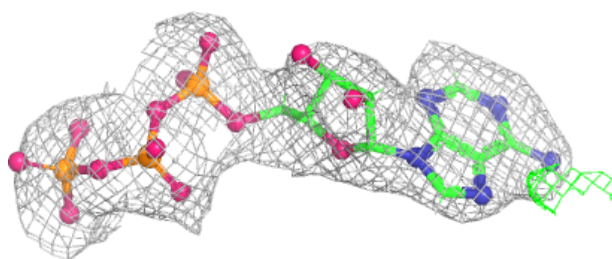
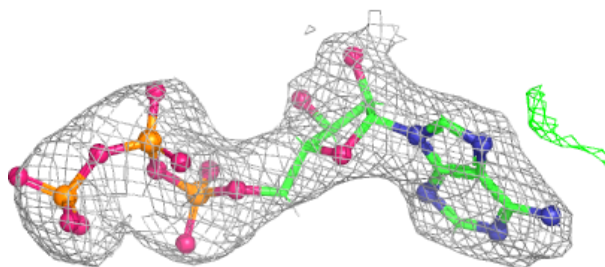
**Electron density around ATP B 301:**

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

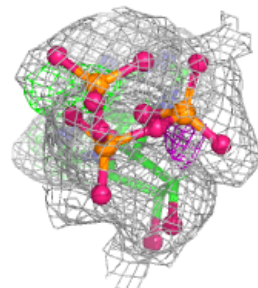
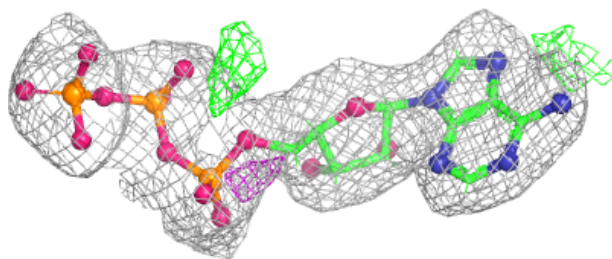
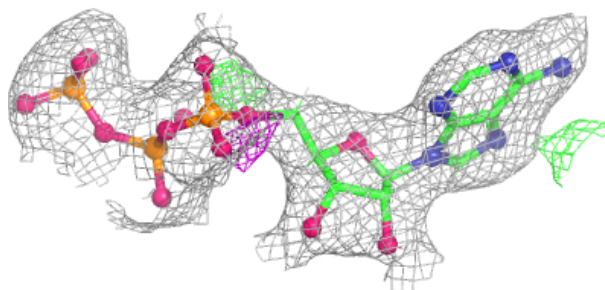


Electron density around ATP F 301:

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

**Electron density around ATP A 301:**

$2mF_o-DF_c$ (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.