Full wwPDB X-ray Structure Validation Report (i)

Feb 28, 2023 – 03:28 pm GMT

PDB ID	:	6YMI
Title	:	Crystal structure of the SAM-SAH riboswitch with AMP.
Authors	:	Huang, L.; Lilley, D.M.J.
Deposited on	:	2020-04-08
Resolution	:	2.50 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 (i) 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 (i)) were used in the production of this report:

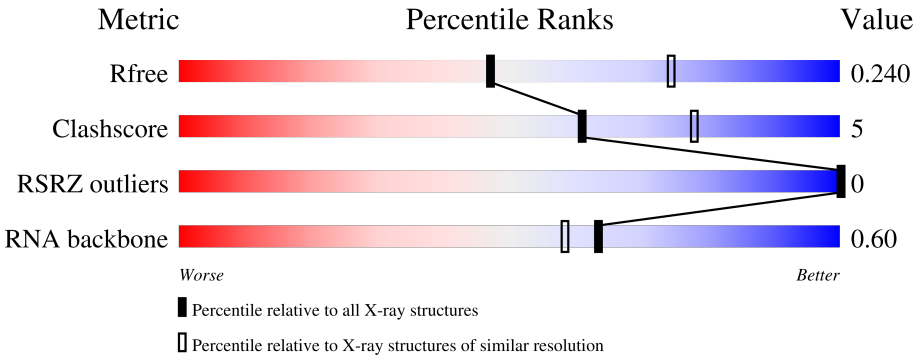
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.32.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.32.1

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 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)
RNA backbone	3102	1008 (2.84-2.16)

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 <=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	26	85% 12% .
1	C	26	73% 23% .
1	F	26	65% 31% .
1	I	26	77% 19% .
1	M	26	77% 19% .
1	O	26	85% 15%

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Mol	Chain	Length	Quality of chain
2	B	9	56% 33% 11%
2	D	9	56% 44%
2	G	9	67% 33%
2	J	9	67% 33%
2	N	9	56% 44%
2	P	9	44% 44% 11%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 4679 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.

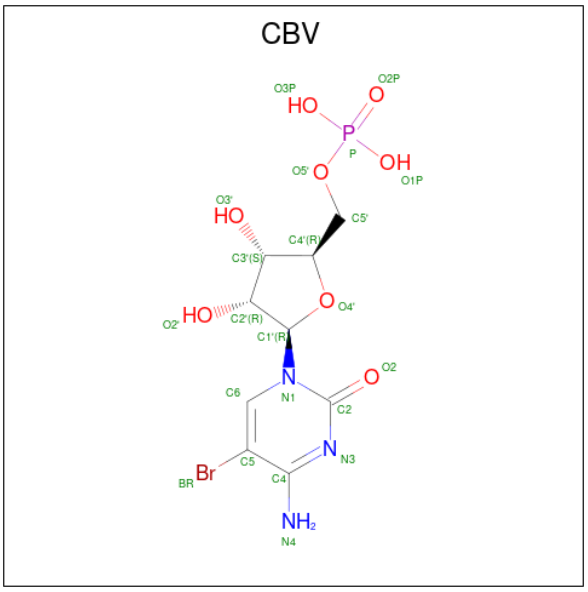
- Molecule 1 is a RNA chain called Chains: A,C,F,I,M,O.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	25	Total	C	N	O	P	0	0	0
			529	237	94	174	24			
1	C	25	Total	C	N	O	P	0	0	0
			529	237	94	174	24			
1	F	25	Total	C	N	O	P	0	0	0
			529	237	94	174	24			
1	I	25	Total	C	N	O	P	0	0	0
			529	237	94	174	24			
1	M	25	Total	C	N	O	P	0	0	0
			529	237	94	174	24			
1	O	26	Total	Br	C	N	O	P	0	0	0
			550	1	246	97	181	25			

- Molecule 2 is a RNA chain called Chains: B,D,G,J,N,P.

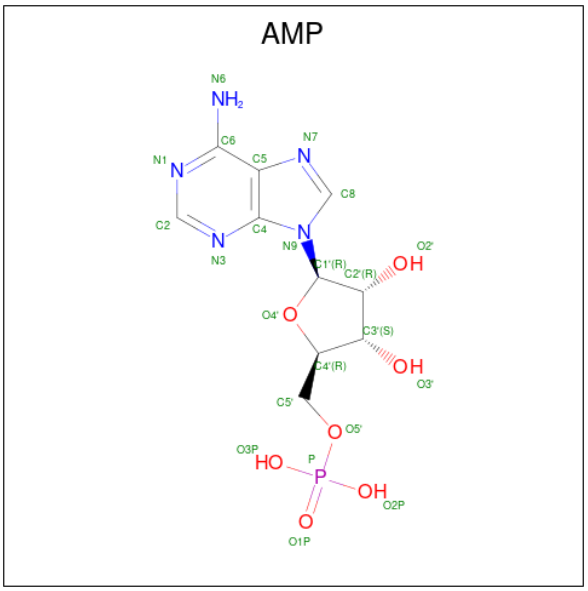
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	9	Total	C	N	O	P	0	0	0
			192	87	37	60	8			
2	D	9	Total	C	N	O	P	0	0	0
			192	87	37	60	8			
2	G	9	Total	C	N	O	P	0	0	0
			192	87	37	60	8			
2	J	9	Total	C	N	O	P	0	0	0
			192	87	37	60	8			
2	N	9	Total	C	N	O	P	0	0	0
			192	87	37	60	8			
2	P	9	Total	C	N	O	P	0	0	0
			192	87	37	60	8			

- Molecule 3 is 5-BROMOCYTIDINE 5'-(DIHYDROGEN PHOSPHATE) (three-letter code: CBV) (formula: $C_9H_{13}BrN_3O_8P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	Br	C	N	O	P	0	0
			21	1	9	3	7	1		
3	C	1	Total	Br	C	N	O	P	0	0
			21	1	9	3	7	1		
3	F	1	Total	Br	C	N	O	P	0	0
			21	1	9	3	7	1		
3	I	1	Total	Br	C	N	O	P	0	0
			21	1	9	3	7	1		
3	M	1	Total	Br	C	N	O	P	0	0
			21	1	9	3	7	1		

- Molecule 4 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: $C_{10}H_{14}N_5O_7P$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	F	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	G	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	I	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	M	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	N	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	O	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	O	0	0
			1	1		
5	F	4	Total	O	0	0
			4	4		
5	G	3	Total	O	0	0
			3	3		

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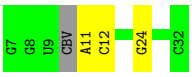
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	I	1	Total O 1 1	0	0
5	M	8	Total O 8 8	0	0
5	O	3	Total O 3 3	0	0

3 Residue-property plots (i)

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

Chain A:



- Molecule 1: Chains: A,C,F,I,M,O

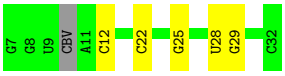
Chain C:

- Molecule 1: Chains: A,C,F,I,M,O

Chain F:

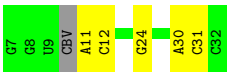
- Molecule 1: Chains: A,C,F,I,M,O

Chain I:



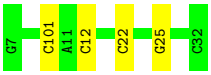
- Molecule 1: Chains: A,C,F,I,M,O

Chain M:



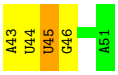
- Molecule 1: Chains: A,C,F,I,M,O

Chain O:



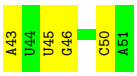
• Molecule 2: Chains: B,D,G,J,N,P

Chain B:



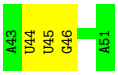
• Molecule 2: Chains: B,D,G,J,N,P

Chain D:



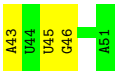
• Molecule 2: Chains: B,D,G,J,N,P

Chain G:



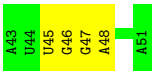
• Molecule 2: Chains: B,D,G,J,N,P

Chain J:



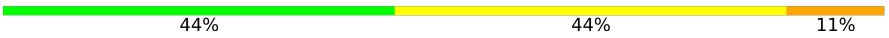
• Molecule 2: Chains: B,D,G,J,N,P

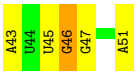
Chain N:



• Molecule 2: Chains: B,D,G,J,N,P

Chain P:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	86.51Å 147.36Å 74.84Å 90.00° 91.36° 90.00°	Depositor
Resolution (Å)	74.59 – 2.50 74.82 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.1 (74.59-2.50) 99.0 (74.82-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.03 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660, PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.214 , 0.242 0.216 , 0.240	Depositor DCC
R_{free} test set	1621 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	60.3	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.23 , 15.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.030 for -1/2*h+1/2*k,3/2*h+1/2*k,-l 0.033 for -1/2*h-1/2*k,-3/2*h+1/2*k,-l 0.167 for 1/2*h+1/2*k,3/2*h-1/2*k,-l 0.269 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.034 for -h,-k,l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	4679	wwPDB-VP
Average B, all atoms (Å^2)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: AMP, CBV

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	0.21	0/589	0.76	0/914
1	C	0.23	0/589	0.80	0/914
1	F	0.22	0/589	0.78	0/914
1	I	0.21	0/589	0.77	0/914
1	M	0.21	0/589	0.79	0/914
1	O	0.21	0/589	0.76	0/914
2	B	0.35	0/215	0.90	0/334
2	D	0.29	0/215	0.87	0/334
2	G	0.32	0/215	0.87	0/334
2	J	0.28	0/215	0.86	0/334
2	N	0.31	0/215	0.88	0/334
2	P	0.26	0/215	0.76	0/334
All	All	0.24	0/4824	0.80	0/7488

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

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	529	0	273	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	529	0	273	3	0
1	F	529	0	273	5	0
1	I	529	0	273	3	0
1	M	529	0	273	6	0
1	O	550	0	282	2	0
2	B	192	0	99	2	0
2	D	192	0	99	1	1
2	G	192	0	99	1	0
2	J	192	0	99	1	0
2	N	192	0	99	2	0
2	P	192	0	99	2	1
3	A	21	0	10	2	0
3	C	21	0	11	0	0
3	F	21	0	10	1	0
3	I	21	0	11	0	0
3	M	21	0	10	2	0
4	A	46	0	24	3	0
4	C	23	0	12	0	0
4	F	23	0	12	0	0
4	G	23	0	12	3	0
4	I	23	0	12	0	0
4	M	23	0	12	0	0
4	N	23	0	12	1	0
4	O	23	0	12	0	0
5	A	1	0	0	0	0
5	F	4	0	0	1	0
5	G	3	0	0	0	0
5	I	1	0	0	0	0
5	M	8	0	0	1	0
5	O	3	0	0	0	0
All	All	4679	0	2401	30	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (30) 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:M:31:C:OP2	5:M:201:HOH:O	1.95	0.83
1:F:25:G:OP1	5:F:201:HOH:O	2.12	0.67
1:A:11:A:H2'	1:A:12:C:C6	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:11:A:H2'	1:M:12:C:C6	2.35	0.62
1:F:11:A:H2'	1:F:12:C:C6	2.35	0.61
1:A:11:A:H8	3:A:101:CBV:H2'	1.69	0.57
1:F:24:G:N3	4:G:101:AMP:H2	2.04	0.55
2:G:44:U:OP2	4:G:101:AMP:O2'	2.24	0.54
1:A:11:A:C8	3:A:101:CBV:H2'	2.44	0.53
1:M:24:G:N3	4:N:101:AMP:H2	2.06	0.53
1:A:24:G:N3	4:A:102:AMP:H2	2.08	0.52
1:F:30:A:H2'	1:F:31:C:C6	2.47	0.49
1:I:25:G:H4'	2:J:43:A:C4	2.48	0.49
4:G:101:AMP:C8	1:O:22:C:H4'	2.47	0.48
4:A:102:AMP:C8	1:I:22:C:H4'	2.49	0.48
1:M:11:A:H8	3:M:101:CBV:H2'	1.78	0.47
1:O:25:G:H4'	2:P:43:A:C4	2.49	0.47
4:A:102:AMP:O2'	2:B:44:U:OP2	2.34	0.45
3:F:101:CBV:H6	3:F:101:CBV:O5'	2.17	0.45
1:I:28:U:H2'	1:I:29:G:H8	1.81	0.45
1:C:25:G:H4'	2:D:43:A:C4	2.52	0.45
1:M:30:A:H2'	1:M:31:C:C6	2.52	0.44
2:P:46:G:C2	2:P:47:G:C8	3.06	0.44
2:B:43:A:H5"	2:B:45:U:OP2	2.18	0.43
1:M:11:A:C8	3:M:101:CBV:H2'	2.54	0.42
2:N:47:G:H2'	2:N:48:A:C8	2.54	0.42
1:C:28:U:H2'	1:C:29:G:H8	1.85	0.42
1:F:14:A:H2'	1:F:15:C:O4'	2.20	0.41
1:C:30:A:H2'	1:C:31:C:C6	2.56	0.41
2:N:47:G:H2'	2:N:48:A:H8	1.84	0.41

All (1) 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)
2:D:50:C:O2	2:P:51:A:O2'[4_445]	2.16	0.04

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

There are no protein molecules in this entry.

5.3.2 Protein sidechains (i)

There are no protein molecules in this entry.

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	23/26 (88%)	0	0
1	C	23/26 (88%)	1 (4%)	0
1	F	23/26 (88%)	0	0
1	I	23/26 (88%)	1 (4%)	0
1	M	23/26 (88%)	0	0
1	O	24/26 (92%)	1 (4%)	0
2	B	8/9 (88%)	2 (25%)	0
2	D	8/9 (88%)	2 (25%)	0
2	G	8/9 (88%)	2 (25%)	0
2	J	8/9 (88%)	2 (25%)	0
2	N	8/9 (88%)	2 (25%)	0
2	P	8/9 (88%)	2 (25%)	0
All	All	187/210 (89%)	15 (8%)	0

All (15) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	45	U
2	B	46	G
1	C	12	C
2	D	45	U
2	D	46	G
2	G	45	U
2	G	46	G
1	I	12	C
2	J	45	U
2	J	46	G
2	N	45	U
2	N	46	G
1	O	12	C
2	P	45	U
2	P	46	G

There are no RNA pucker outliers to report.

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

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
1	CBV	O	101	1	19,22,23	1.08	3 (15%)	27,32,35	0.70	1 (3%)

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
1	CBV	O	101	1	-	0/7/25/26	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	O	101	CBV	C4-C5	2.56	1.48	1.42
1	O	101	CBV	C6-C5	2.55	1.39	1.34
1	O	101	CBV	C6-N1	-2.18	1.34	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	O	101	CBV	O2-C2-N3	-2.07	118.96	122.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

14 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	$\#\|Z\| > 2$
3	CBV	C	101	-	19,22,23	1.09	3 (15%)	27,32,35	0.69	1 (3%)
4	AMP	N	101	-	22,25,25	0.85	1 (4%)	25,38,38	1.18	2 (8%)
4	AMP	G	101	-	22,25,25	0.86	1 (4%)	25,38,38	1.16	2 (8%)
3	CBV	F	101	1	19,22,23	1.06	3 (15%)	27,32,35	0.69	0
4	AMP	A	103	-	22,25,25	0.89	1 (4%)	25,38,38	1.19	2 (8%)
4	AMP	A	102	-	22,25,25	0.85	1 (4%)	25,38,38	1.17	2 (8%)
4	AMP	F	102	-	22,25,25	0.88	1 (4%)	25,38,38	1.25	2 (8%)
4	AMP	I	102	-	22,25,25	0.87	1 (4%)	25,38,38	1.19	2 (8%)
3	CBV	A	101	1	19,22,23	1.08	3 (15%)	27,32,35	0.68	1 (3%)
3	CBV	M	101	1	19,22,23	1.06	3 (15%)	27,32,35	0.70	0
3	CBV	I	101	-	19,22,23	1.06	3 (15%)	27,32,35	0.69	1 (3%)
4	AMP	M	102	-	22,25,25	0.89	1 (4%)	25,38,38	1.24	2 (8%)
4	AMP	C	102	-	22,25,25	0.89	1 (4%)	25,38,38	1.18	2 (8%)
4	AMP	O	201	-	22,25,25	0.88	1 (4%)	25,38,38	1.18	2 (8%)

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
3	CBV	C	101	-	-	0/7/25/26	0/2/2/2
4	AMP	N	101	-	-	2/6/26/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	AMP	G	101	-	-	0/6/26/26	0/3/3/3
3	CBV	F	101	1	-	0/7/25/26	0/2/2/2
4	AMP	A	103	-	-	4/6/26/26	0/3/3/3
4	AMP	A	102	-	-	1/6/26/26	0/3/3/3
4	AMP	F	102	-	-	0/6/26/26	0/3/3/3
4	AMP	I	102	-	-	2/6/26/26	0/3/3/3
3	CBV	A	101	1	-	0/7/25/26	0/2/2/2
3	CBV	M	101	1	-	0/7/25/26	0/2/2/2
3	CBV	I	101	-	-	0/7/25/26	0/2/2/2
4	AMP	M	102	-	-	0/6/26/26	0/3/3/3
4	AMP	C	102	-	-	1/6/26/26	0/3/3/3
4	AMP	O	201	-	-	5/6/26/26	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	101	CBV	C6-C5	2.59	1.39	1.34
3	M	101	CBV	C4-C5	2.58	1.48	1.42
3	C	101	CBV	C4-C5	2.58	1.48	1.42
3	F	101	CBV	C4-C5	2.56	1.48	1.42
3	A	101	CBV	C4-C5	2.56	1.48	1.42
3	I	101	CBV	C6-C5	2.55	1.39	1.34
4	C	102	AMP	C5-C4	2.52	1.47	1.40
4	A	103	AMP	C5-C4	2.51	1.47	1.40
4	G	101	AMP	C5-C4	2.51	1.47	1.40
4	O	201	AMP	C5-C4	2.50	1.47	1.40
4	N	101	AMP	C5-C4	2.49	1.47	1.40
3	A	101	CBV	C6-C5	2.49	1.39	1.34
4	I	102	AMP	C5-C4	2.48	1.47	1.40
3	I	101	CBV	C4-C5	2.48	1.48	1.42
4	M	102	AMP	C5-C4	2.47	1.47	1.40
4	A	102	AMP	C5-C4	2.46	1.47	1.40
4	F	102	AMP	C5-C4	2.44	1.47	1.40
3	M	101	CBV	C6-C5	2.43	1.39	1.34
3	F	101	CBV	C6-C5	2.39	1.39	1.34
3	M	101	CBV	C6-N1	-2.24	1.34	1.38
3	A	101	CBV	C6-N1	-2.22	1.34	1.38
3	F	101	CBV	C6-N1	-2.21	1.34	1.38
3	I	101	CBV	C6-N1	-2.12	1.34	1.38
3	C	101	CBV	C6-N1	-2.09	1.34	1.38

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	F	102	AMP	N3-C2-N1	-3.17	123.73	128.68
4	M	102	AMP	N3-C2-N1	-3.16	123.74	128.68
4	A	103	AMP	N3-C2-N1	-3.14	123.78	128.68
4	C	102	AMP	N3-C2-N1	-3.12	123.80	128.68
4	I	102	AMP	N3-C2-N1	-3.11	123.83	128.68
4	O	201	AMP	N3-C2-N1	-3.05	123.91	128.68
4	N	101	AMP	C4-C5-N7	-2.92	106.36	109.40
4	M	102	AMP	C4-C5-N7	-2.86	106.42	109.40
4	A	102	AMP	C4-C5-N7	-2.86	106.42	109.40
4	F	102	AMP	C4-C5-N7	-2.84	106.44	109.40
4	G	101	AMP	C4-C5-N7	-2.80	106.48	109.40
4	A	102	AMP	N3-C2-N1	-2.61	124.61	128.68
4	O	201	AMP	C4-C5-N7	-2.56	106.73	109.40
4	A	103	AMP	C4-C5-N7	-2.54	106.75	109.40
4	N	101	AMP	N3-C2-N1	-2.50	124.77	128.68
4	I	102	AMP	C4-C5-N7	-2.48	106.81	109.40
4	C	102	AMP	C4-C5-N7	-2.47	106.82	109.40
4	G	101	AMP	N3-C2-N1	-2.35	125.01	128.68
3	C	101	CBV	O2-C2-N3	-2.18	118.79	122.33
3	I	101	CBV	O2-C2-N3	-2.05	119.00	122.33
3	A	101	CBV	O2-C2-N3	-2.02	119.04	122.33

There are no chirality outliers.

All (15) torsion outliers are listed below:

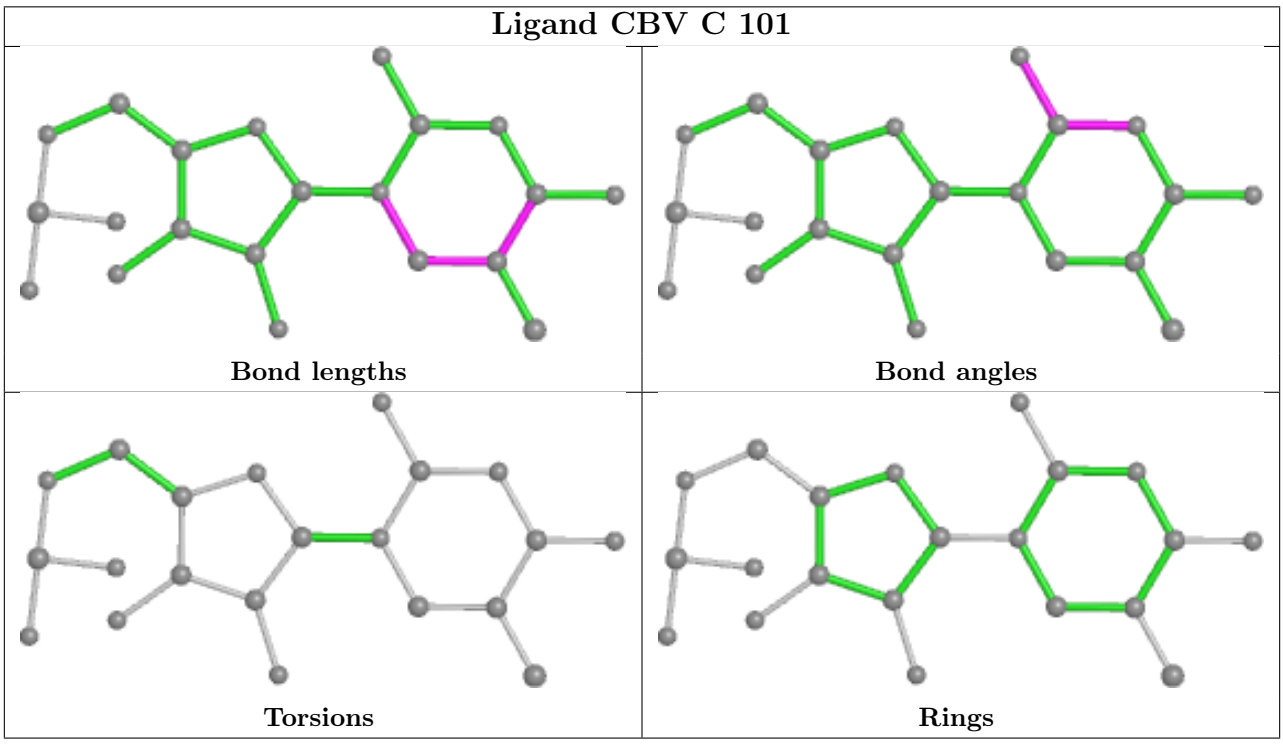
Mol	Chain	Res	Type	Atoms
4	A	103	AMP	C5'-O5'-P-O2P
4	A	103	AMP	C5'-O5'-P-O3P
4	I	102	AMP	O4'-C4'-C5'-O5'
4	N	101	AMP	O4'-C4'-C5'-O5'
4	N	101	AMP	C3'-C4'-C5'-O5'
4	O	201	AMP	C5'-O5'-P-O1P
4	O	201	AMP	C5'-O5'-P-O2P
4	O	201	AMP	C5'-O5'-P-O3P
4	O	201	AMP	O4'-C4'-C5'-O5'
4	O	201	AMP	C3'-C4'-C5'-O5'
4	I	102	AMP	C3'-C4'-C5'-O5'
4	A	103	AMP	C5'-O5'-P-O1P
4	C	102	AMP	O4'-C4'-C5'-O5'
4	A	102	AMP	C5'-O5'-P-O2P
4	A	103	AMP	O4'-C4'-C5'-O5'

There are no ring outliers.

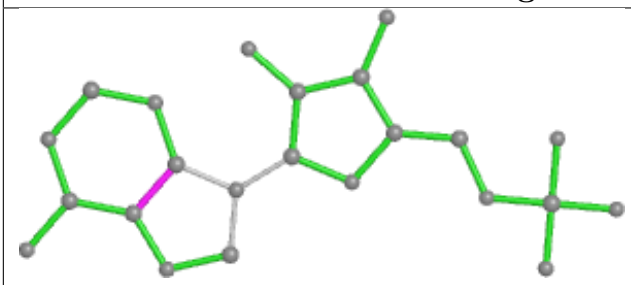
6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	N	101	AMP	1	0
4	G	101	AMP	3	0
3	F	101	CBV	1	0
4	A	102	AMP	3	0
3	A	101	CBV	2	0
3	M	101	CBV	2	0

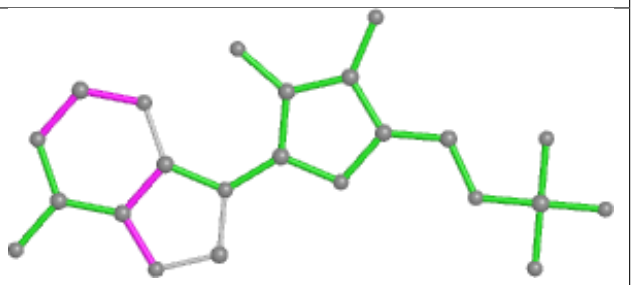
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.



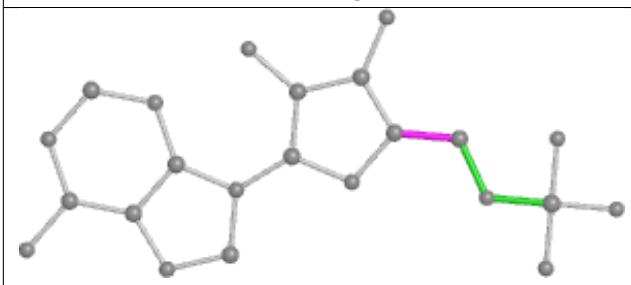
Ligand AMP N 101



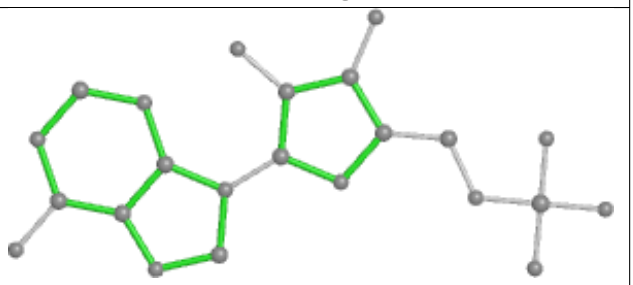
Bond lengths



Bond angles

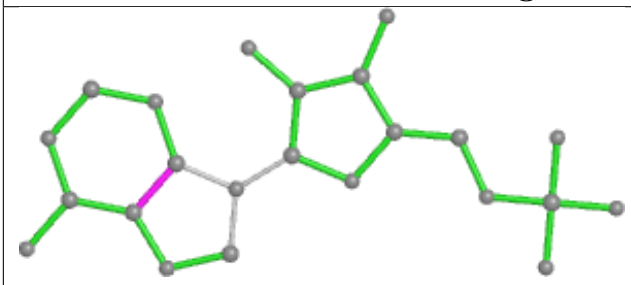


Torsions



Rings

Ligand AMP G 101



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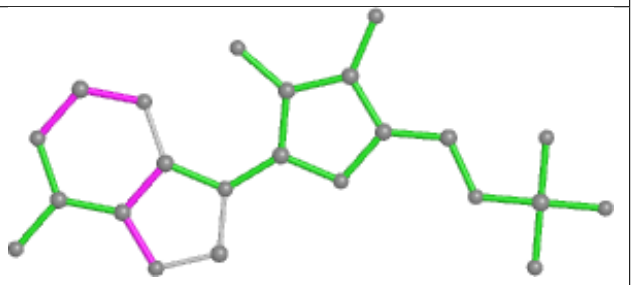

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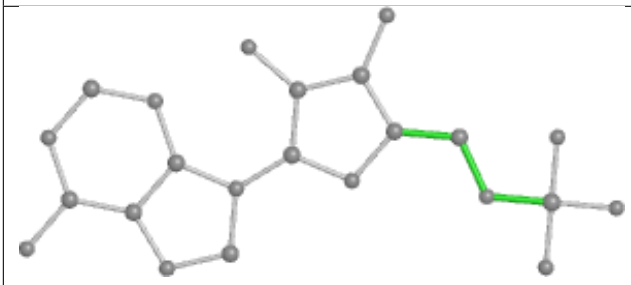

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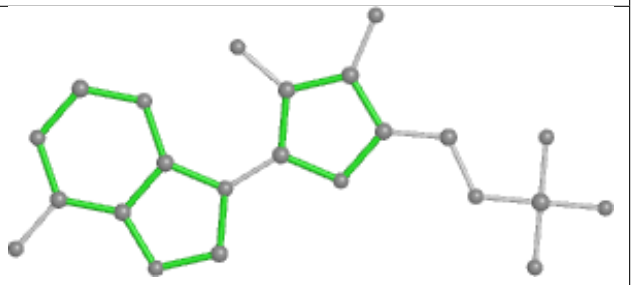
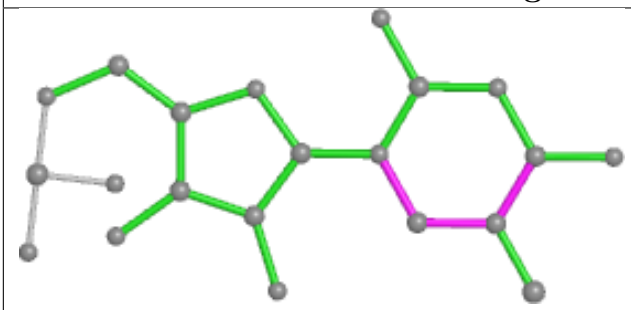


Rings

Ligand CBV F 101



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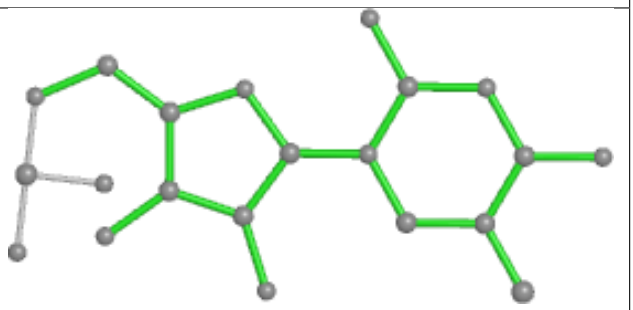
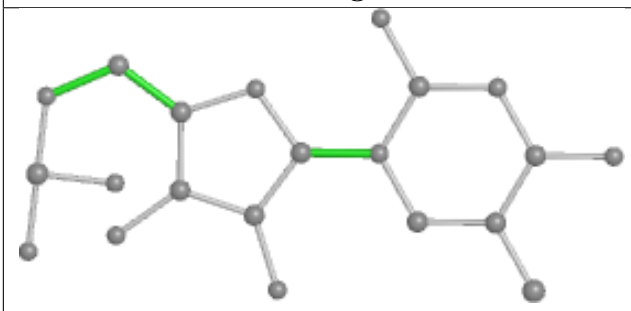


Bond angles



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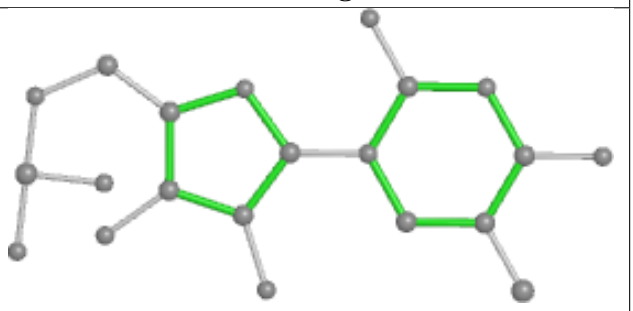
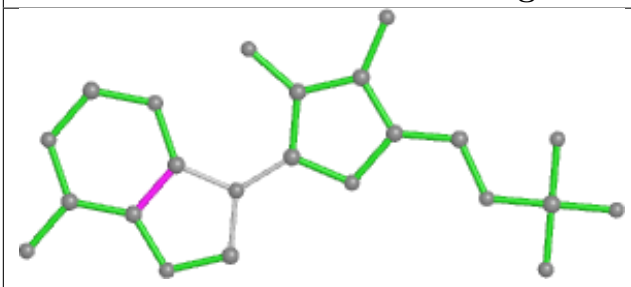


Rings

Ligand AMP A 103



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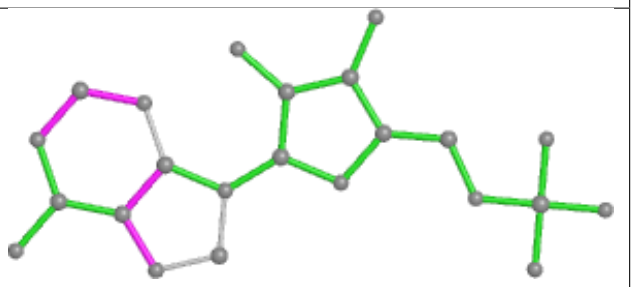

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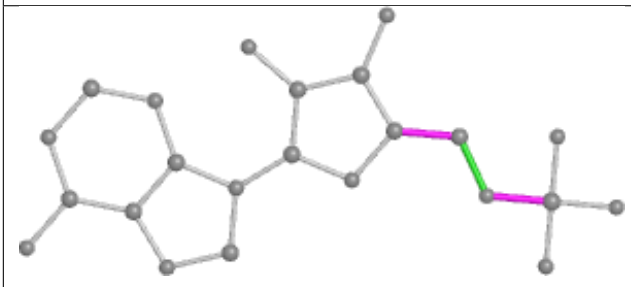

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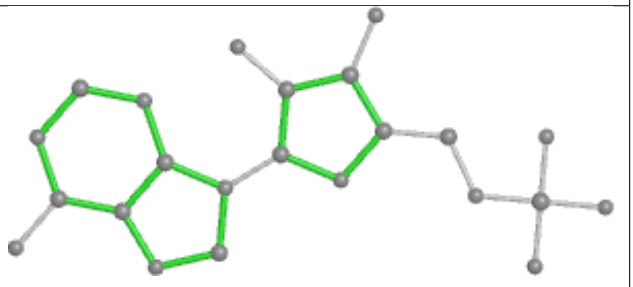
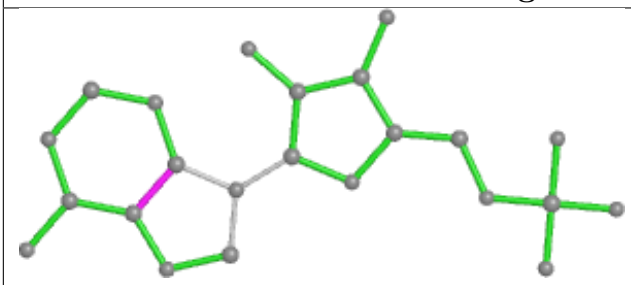


Rings

Ligand AMP A 102



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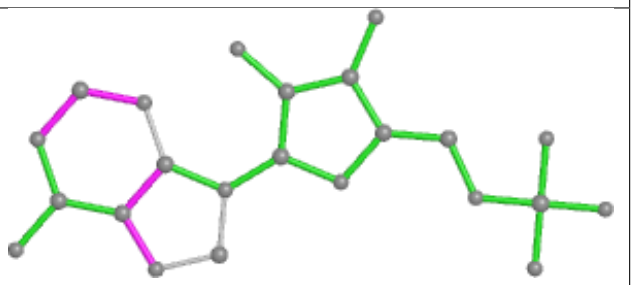

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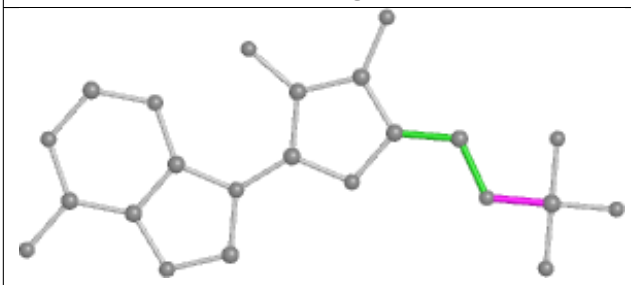

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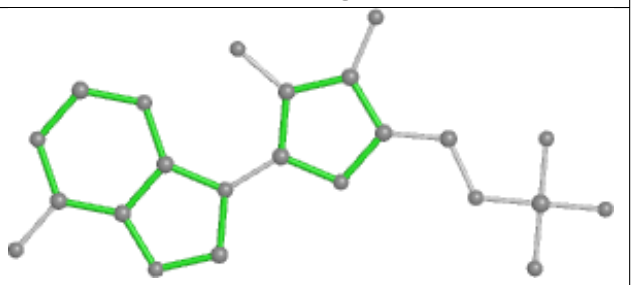

Rings

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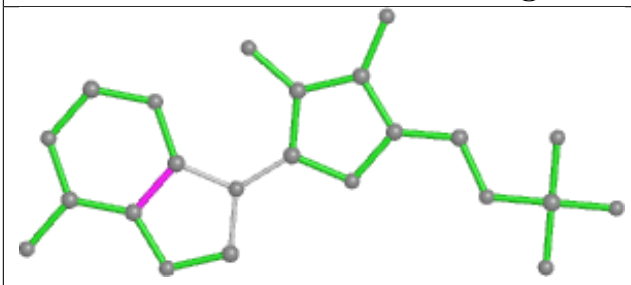

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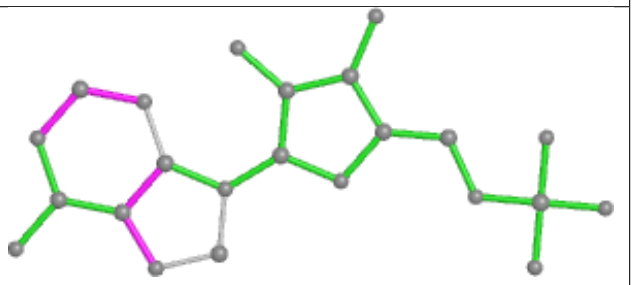
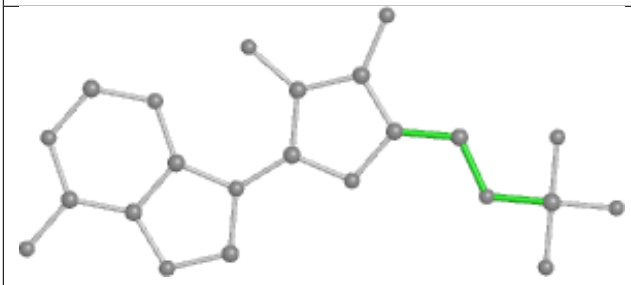


Bond angles



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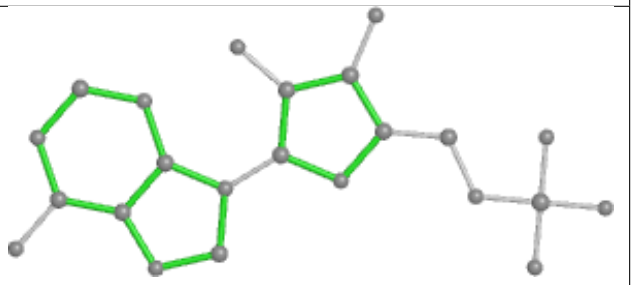
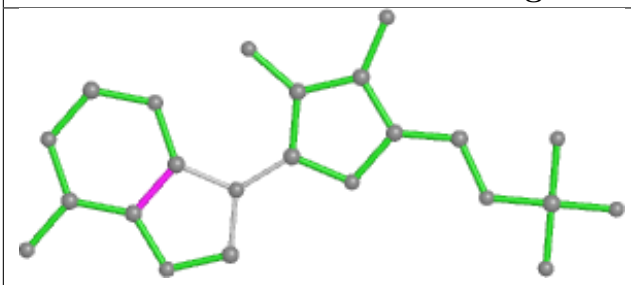
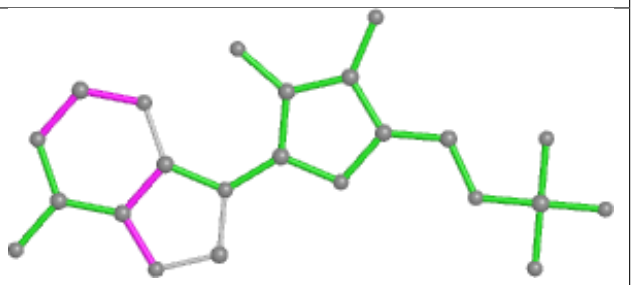


Rings

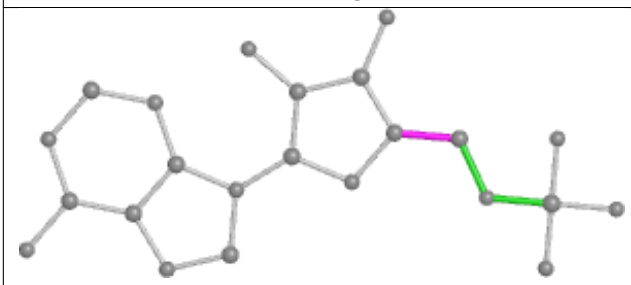
Ligand AMP I 102



Bond lengths



Bond angles



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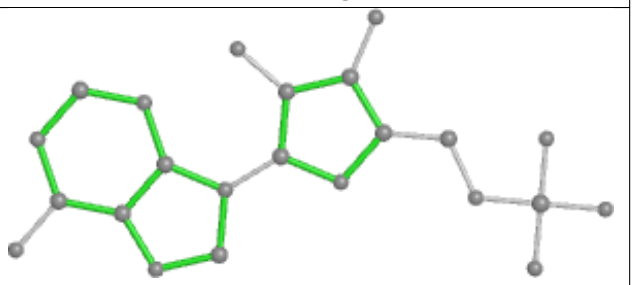
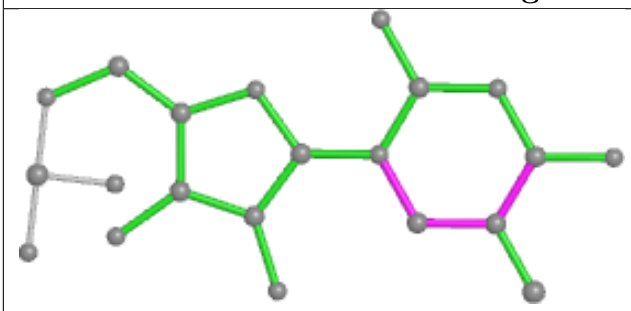
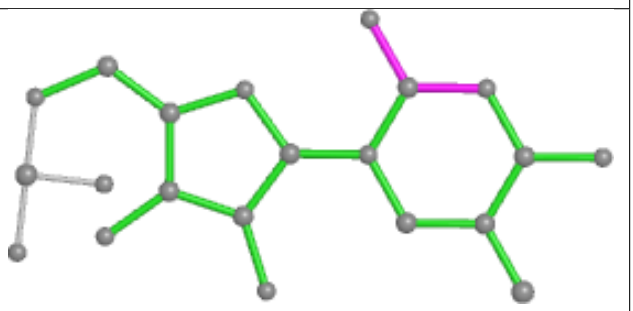


Rings

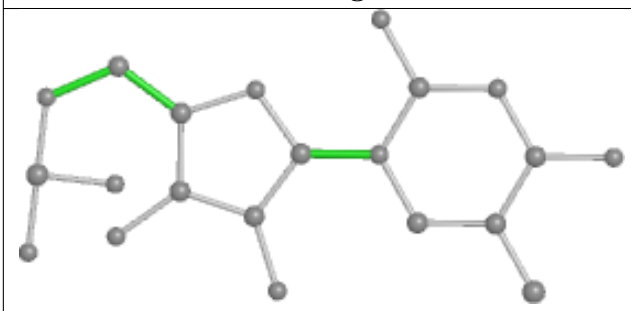
Ligand CBV A 101



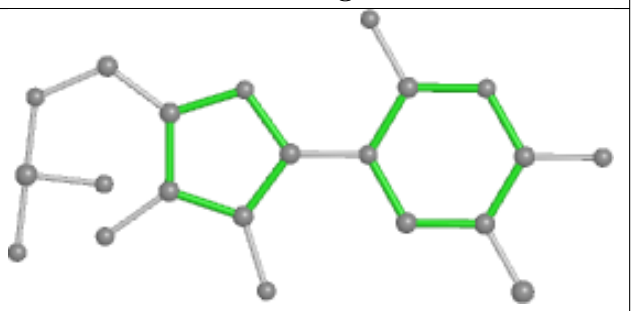
Bond lengths



Bond angles

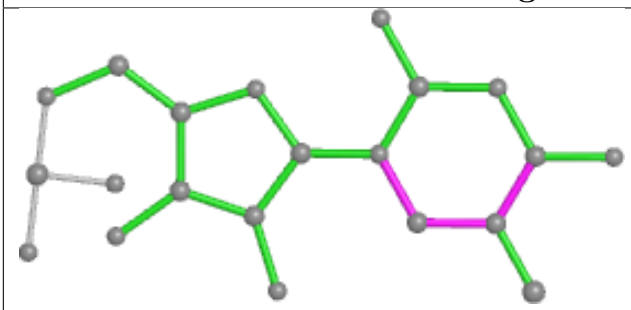


Torsions

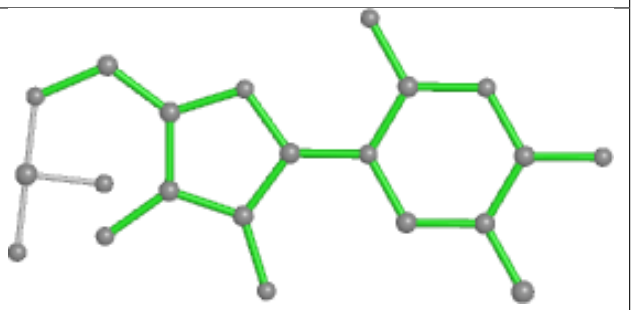


Rings

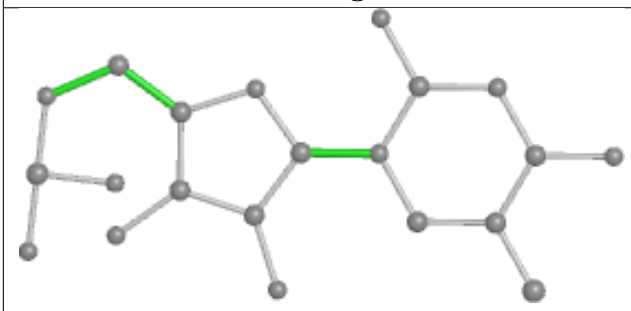
Ligand CBV M 101



Bond lengths



Bond angles



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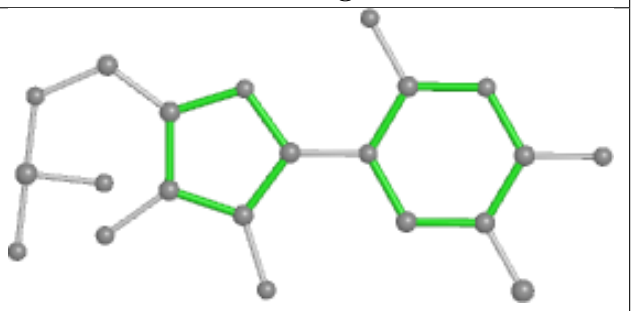
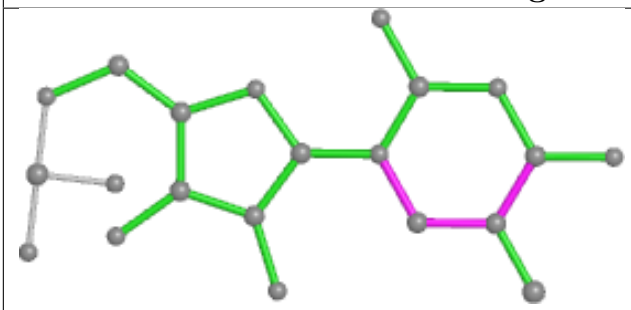
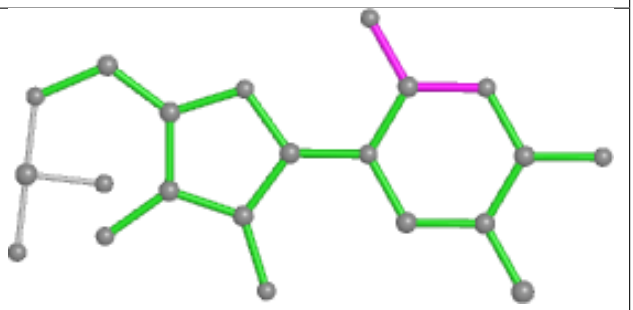


Rings

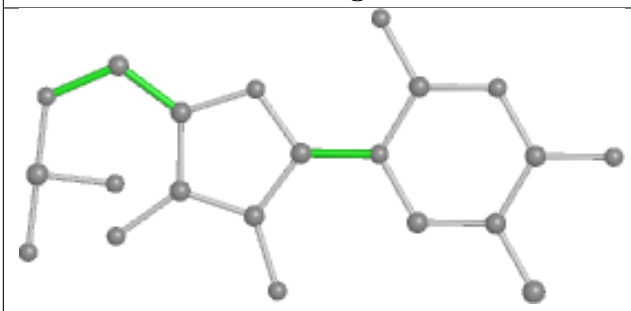
Ligand CBV I 101



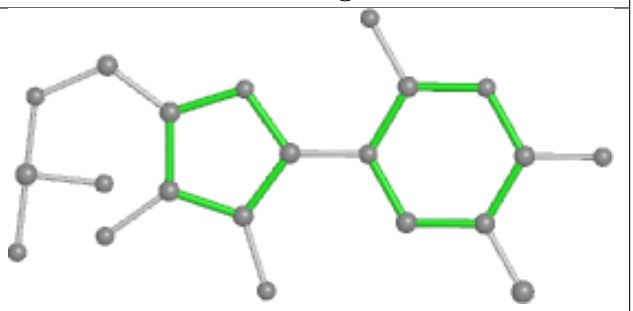
Bond lengths



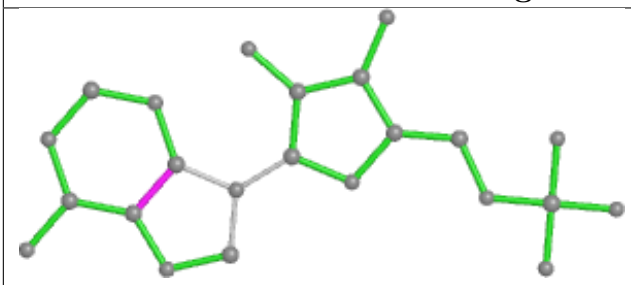
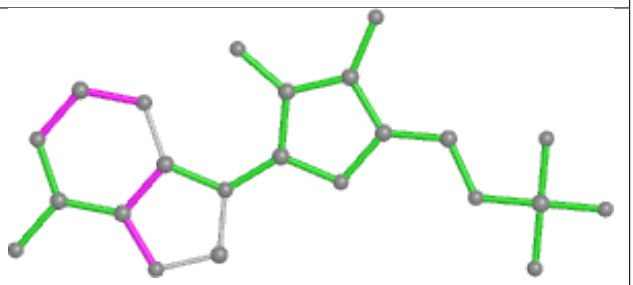
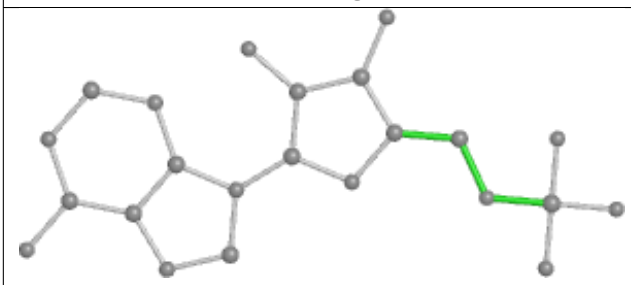
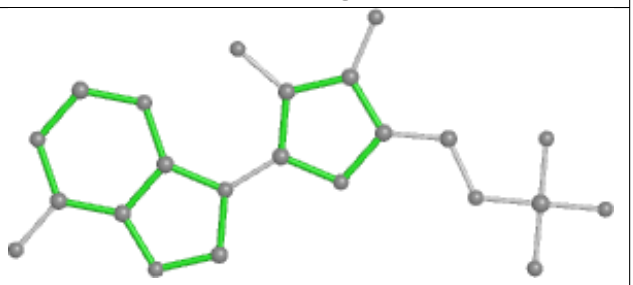
Bond angles

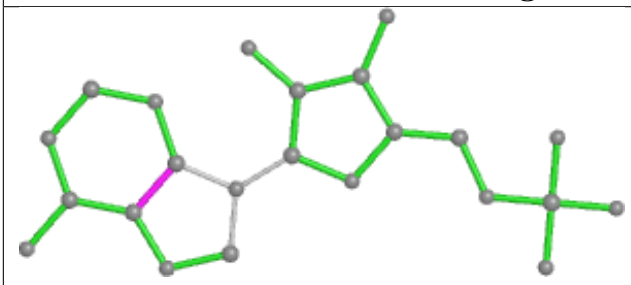
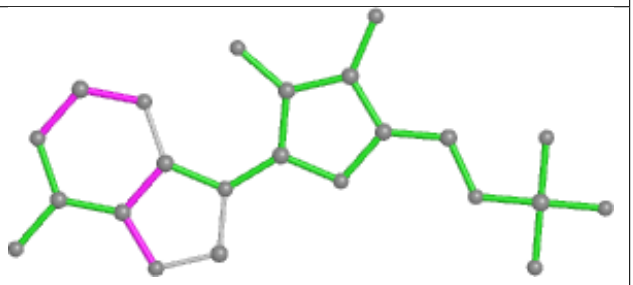
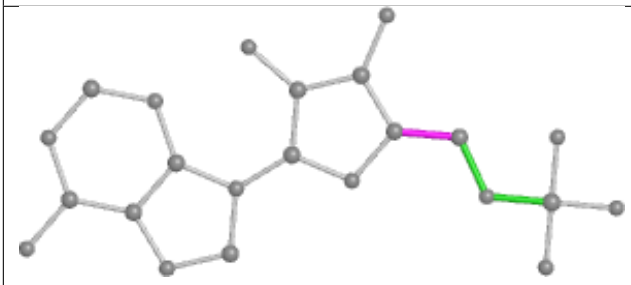
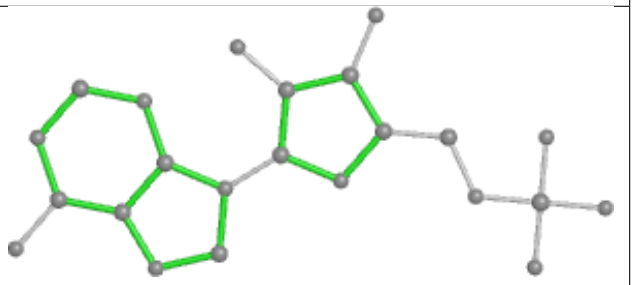


Torsions

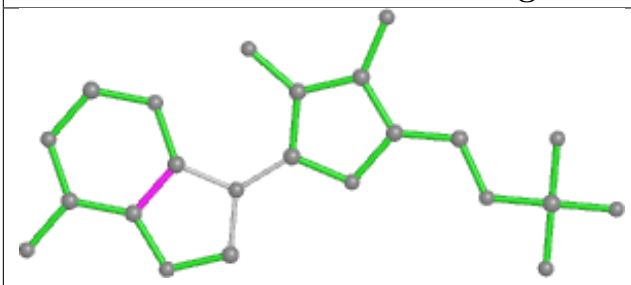


Rings

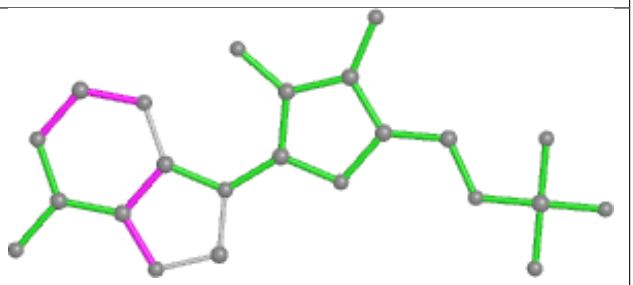
Ligand AMP M 102	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand AMP C 102	
 Bond lengths	 Bond angles
 Torsions	 Rings

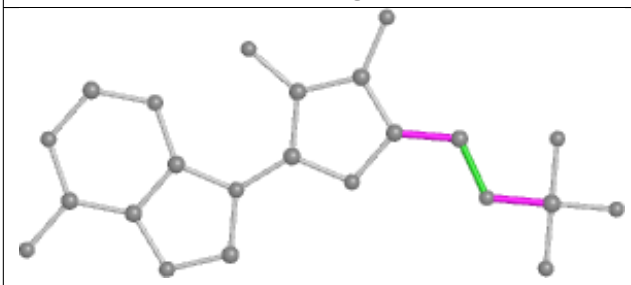
Ligand AMP O 201



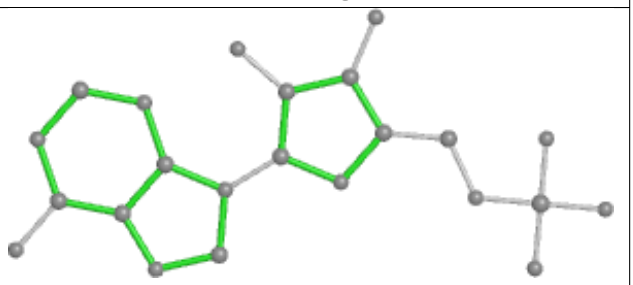
Bond lengths



Bond angles



Torsions



Rings

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^{th} 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	25/26 (96%)	-0.81	0	100	100	46, 53, 74, 88	0
1	C	25/26 (96%)	-0.76	0	100	100	58, 83, 104, 107	0
1	F	25/26 (96%)	-0.77	0	100	100	46, 51, 71, 82	0
1	I	25/26 (96%)	-0.52	0	100	100	66, 86, 101, 103	0
1	M	25/26 (96%)	-0.85	0	100	100	45, 52, 71, 83	0
1	O	25/26 (96%)	-0.62	0	100	100	60, 86, 103, 106	0
2	B	9/9 (100%)	-0.95	0	100	100	52, 56, 60, 67	0
2	D	9/9 (100%)	-0.70	0	100	100	59, 66, 93, 103	0
2	G	9/9 (100%)	-0.84	0	100	100	49, 51, 62, 69	0
2	J	9/9 (100%)	-0.46	0	100	100	60, 66, 106, 112	0
2	N	9/9 (100%)	-0.83	0	100	100	46, 53, 56, 64	0
2	P	9/9 (100%)	-0.49	0	100	100	58, 64, 102, 104	0
All	All	204/210 (97%)	-0.72	0	100	100	45, 64, 102, 112	0

There are no RSRZ outliers to report.

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

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^{th} 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
1	CBV	O	101	21/22	0.93	0.10	58,77,89,113	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

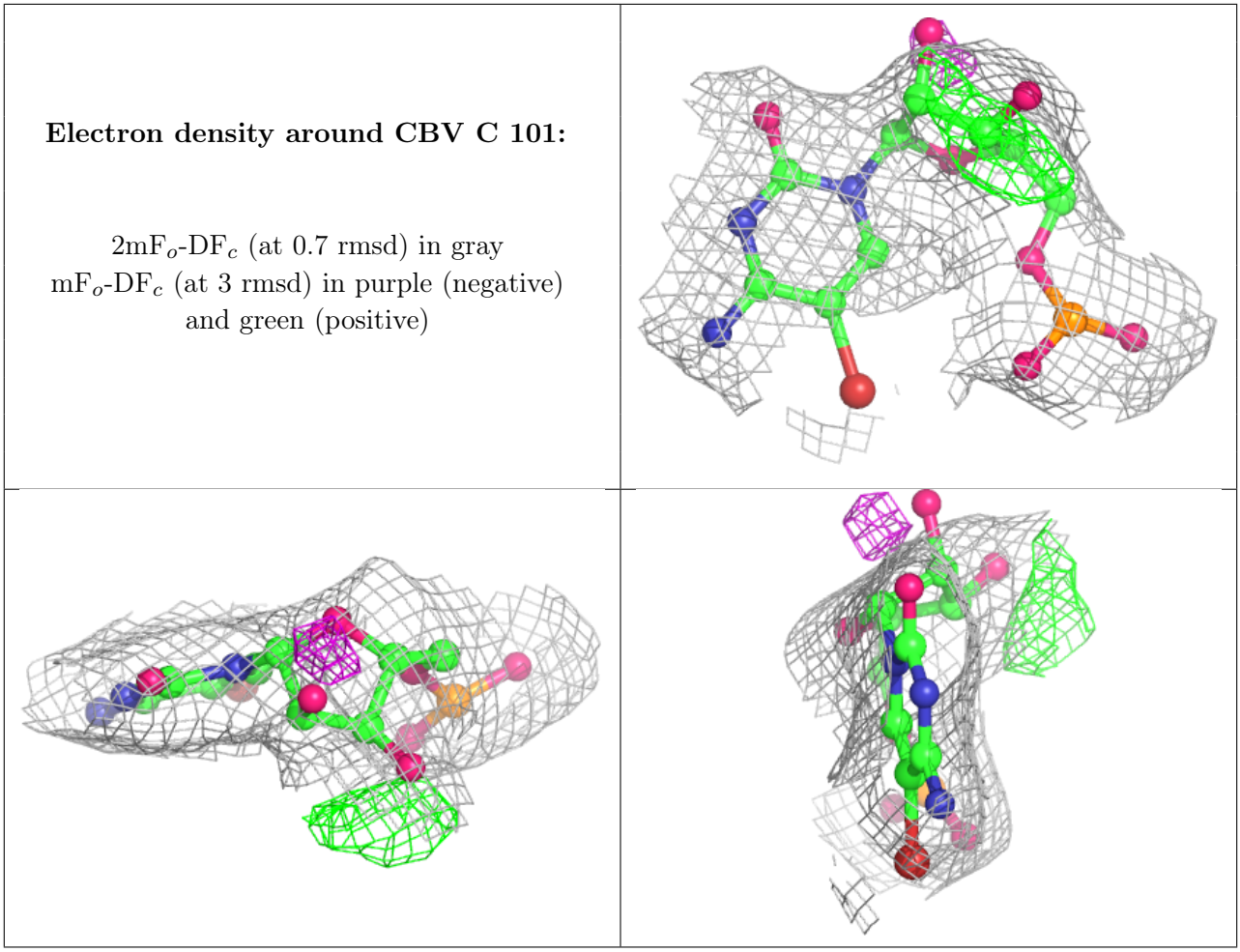
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
3	CBV	C	101	21/22	0.86	0.15	64,87,98,127	0
4	AMP	A	102	23/23	0.87	0.23	74,97,121,123	0
3	CBV	I	101	21/22	0.88	0.15	71,84,93,113	0
4	AMP	G	101	23/23	0.88	0.21	72,90,115,118	0
4	AMP	C	102	23/23	0.89	0.18	84,88,99,101	0
4	AMP	I	102	23/23	0.89	0.17	79,84,94,95	0
4	AMP	N	101	23/23	0.89	0.17	76,81,114,122	0
4	AMP	O	201	23/23	0.89	0.19	77,81,104,107	0
3	CBV	A	101	21/22	0.93	0.12	43,52,63,72	0
3	CBV	M	101	21/22	0.93	0.14	41,48,70,77	0
3	CBV	F	101	21/22	0.94	0.11	40,52,66,82	0
4	AMP	F	102	23/23	0.94	0.17	59,61,70,79	0
4	AMP	A	103	23/23	0.94	0.14	64,68,79,84	0
4	AMP	M	102	23/23	0.96	0.15	65,68,74,82	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.

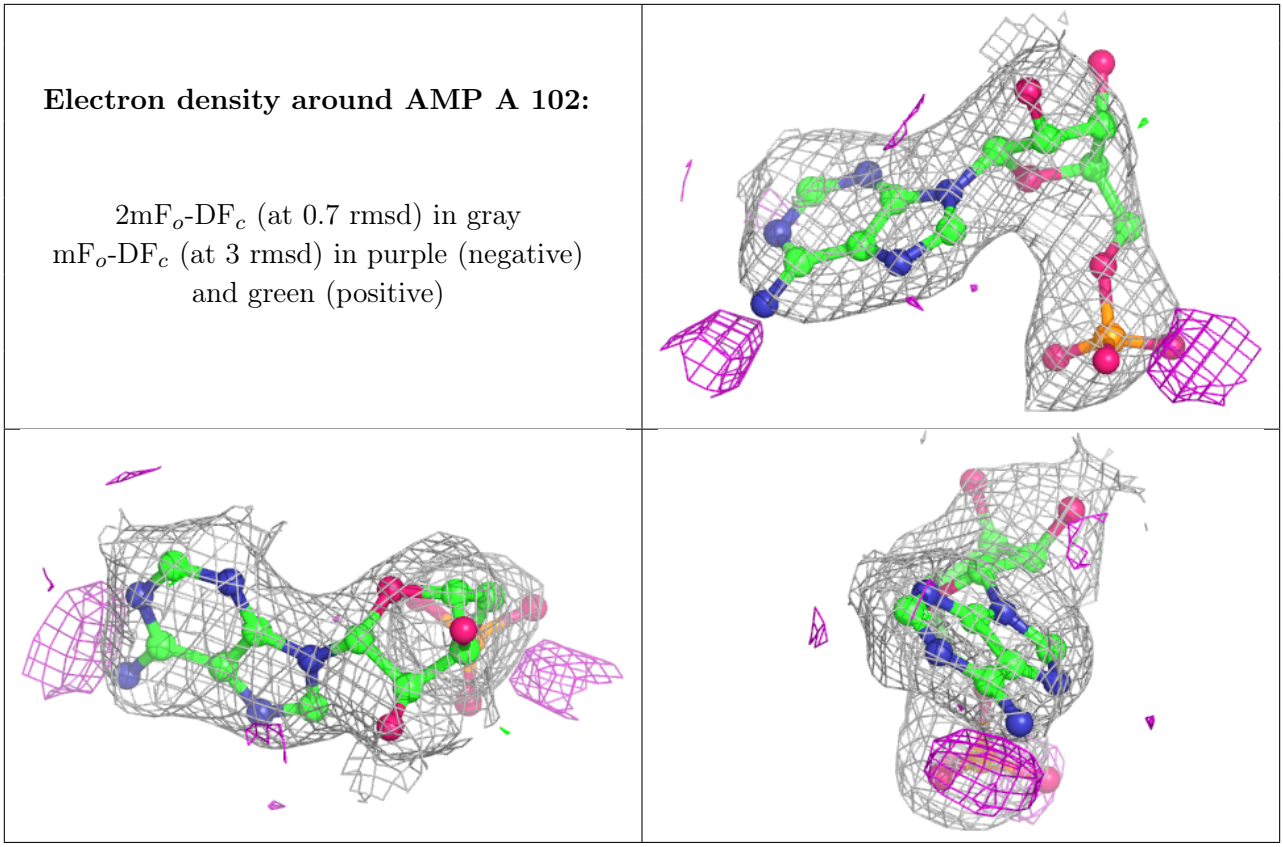
Electron density around CBV C 101:

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



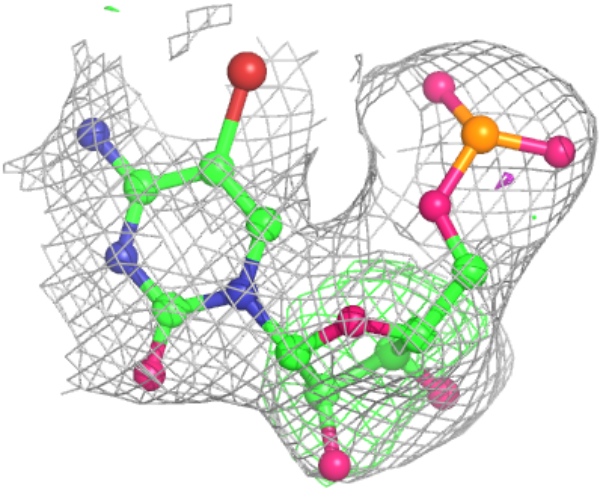
Electron density around AMP A 102:

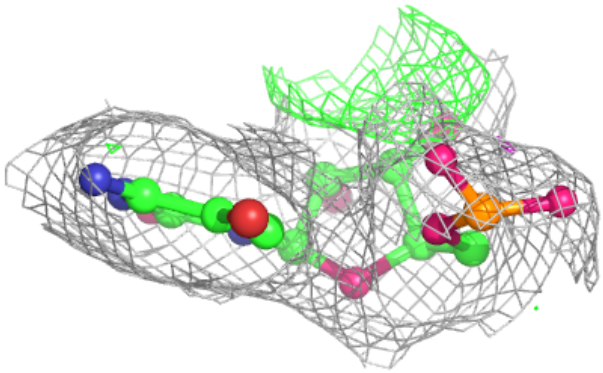
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

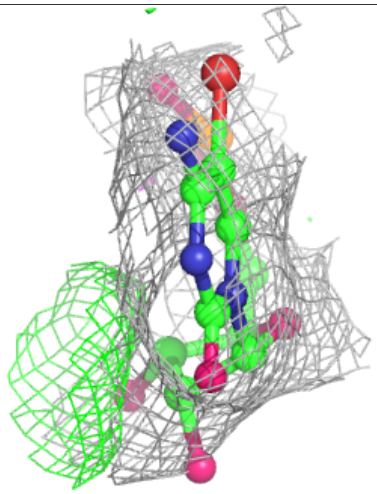


Electron density around CBV I 101:

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







Electron density around AMP G 101:

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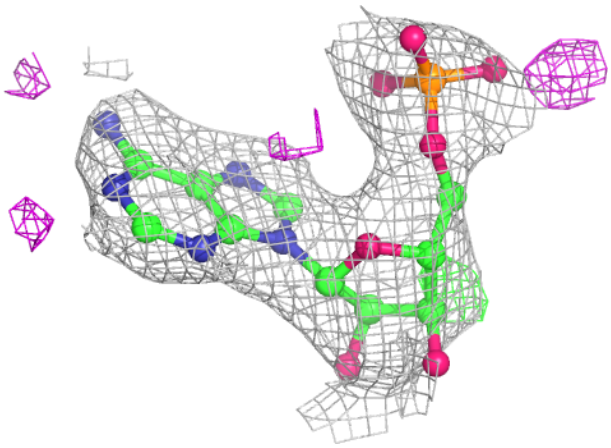
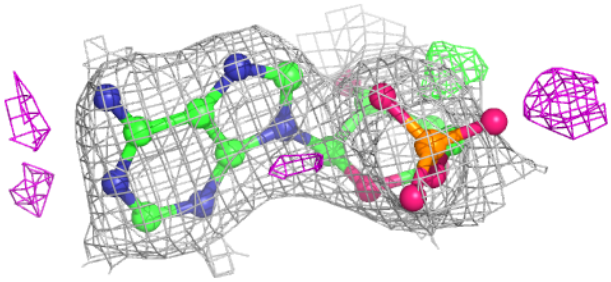
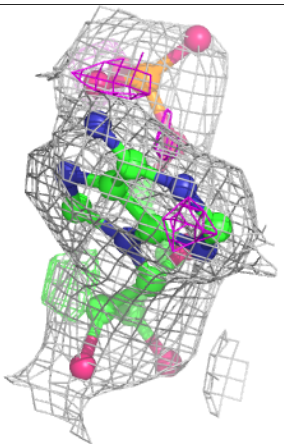
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

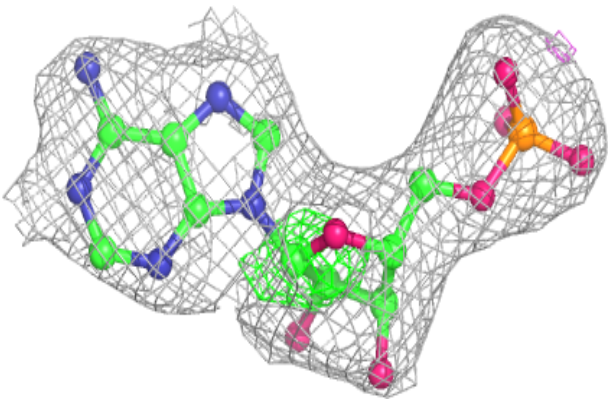


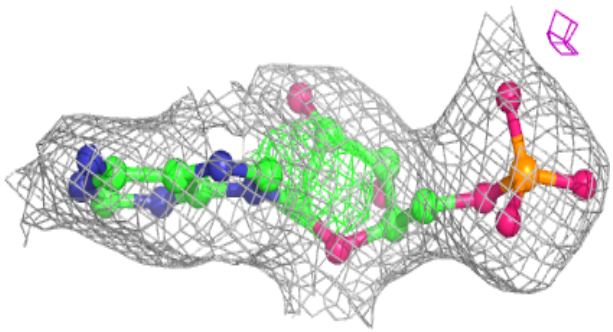


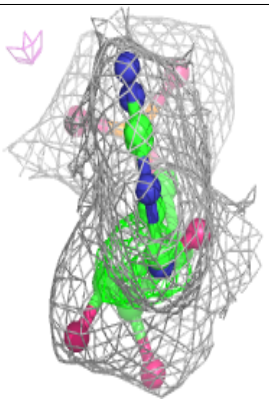


Electron density around AMP C 102:

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

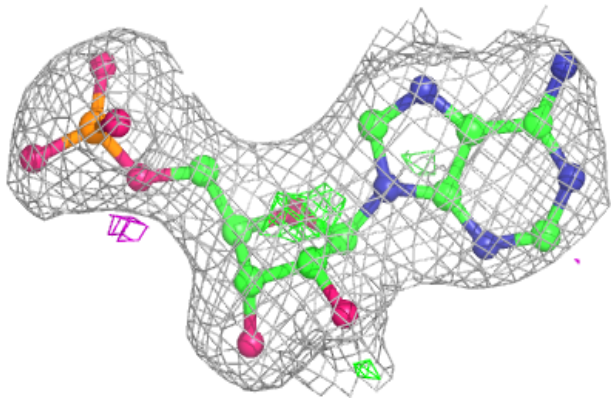


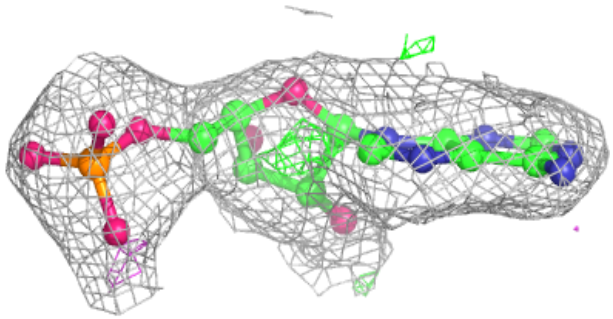


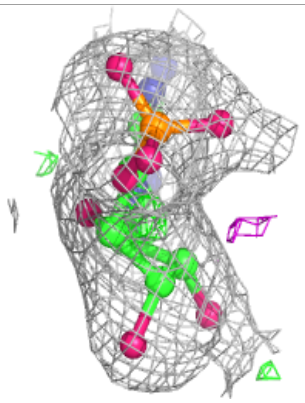


Electron density around AMP I 102:

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

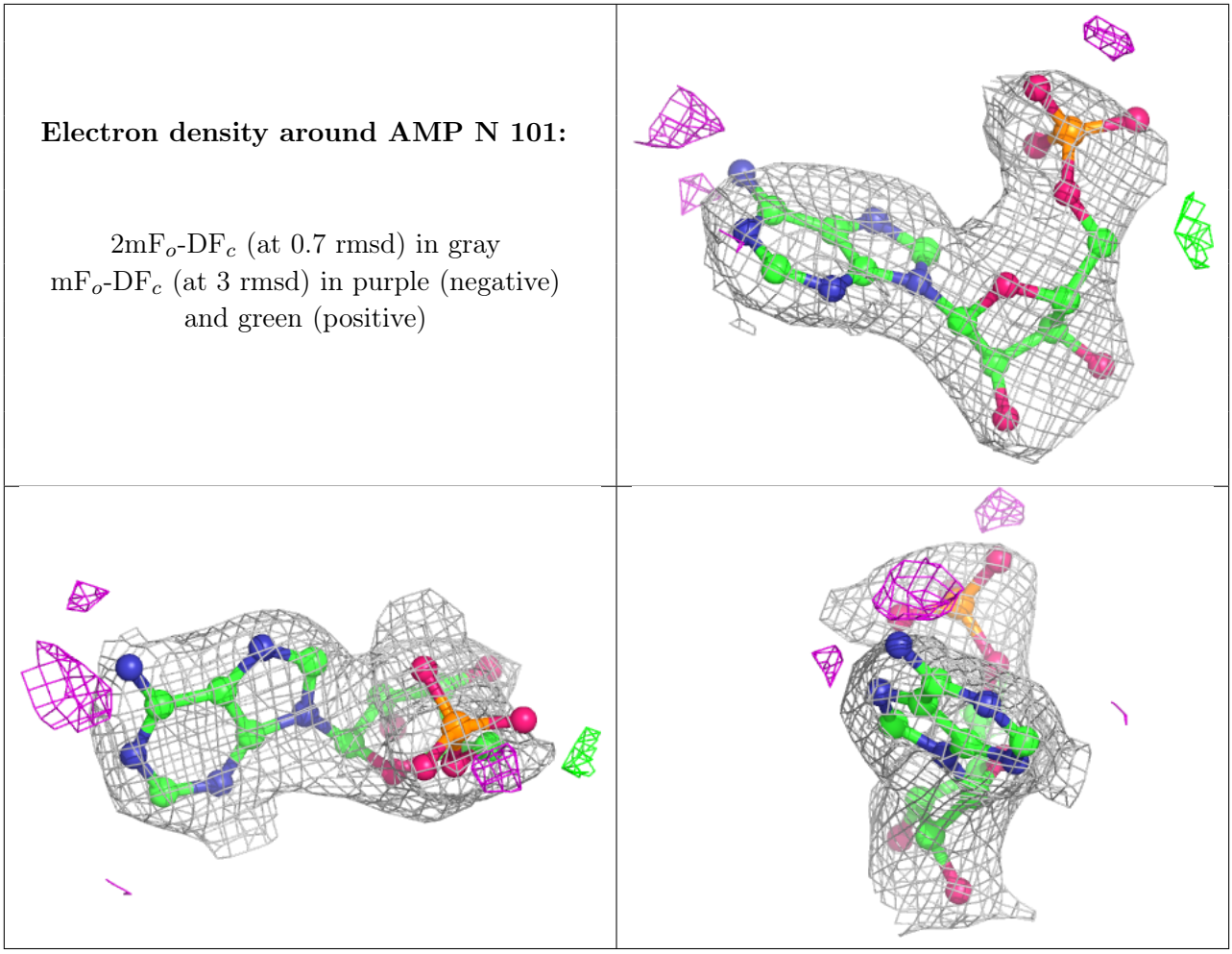






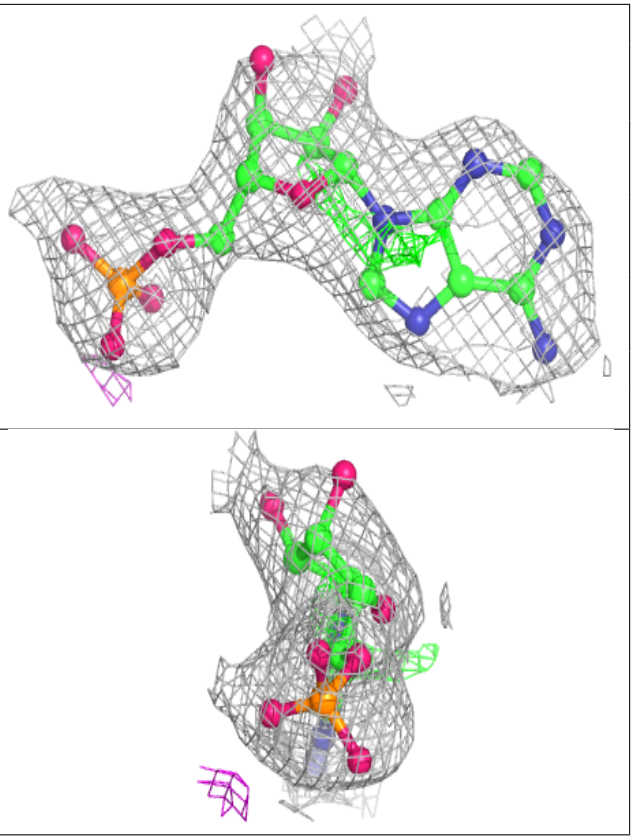
Electron density around AMP N 101:

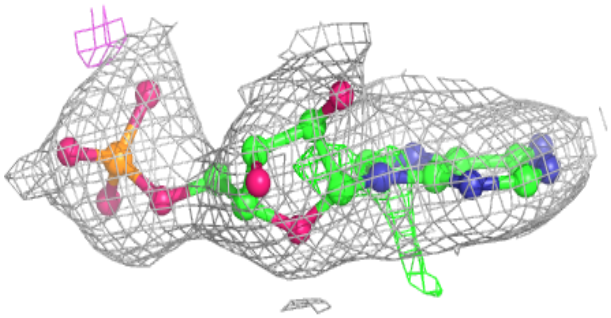
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around AMP O 201:

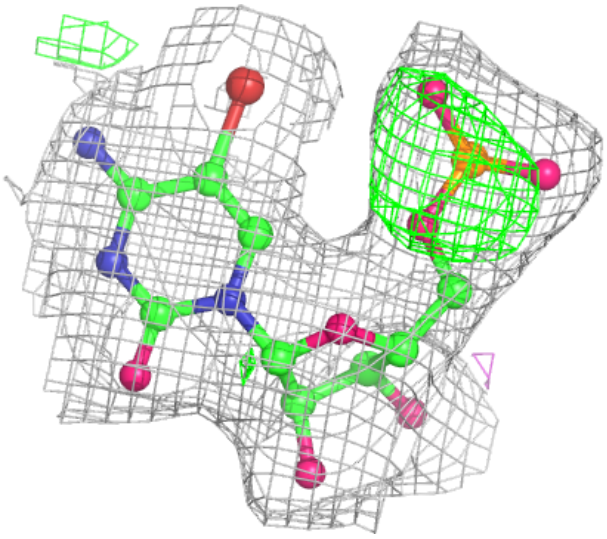
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

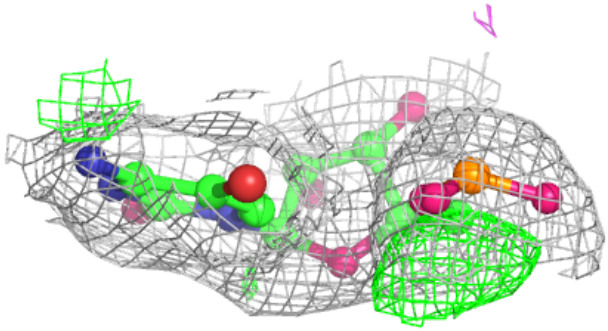


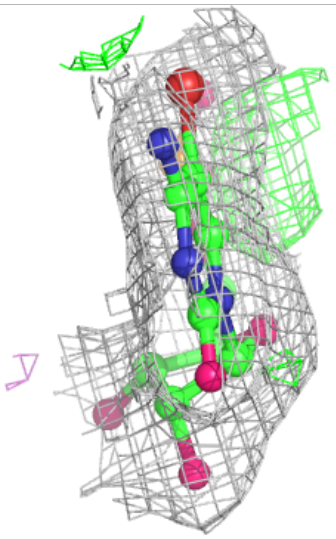


Electron density around CBV A 101:

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

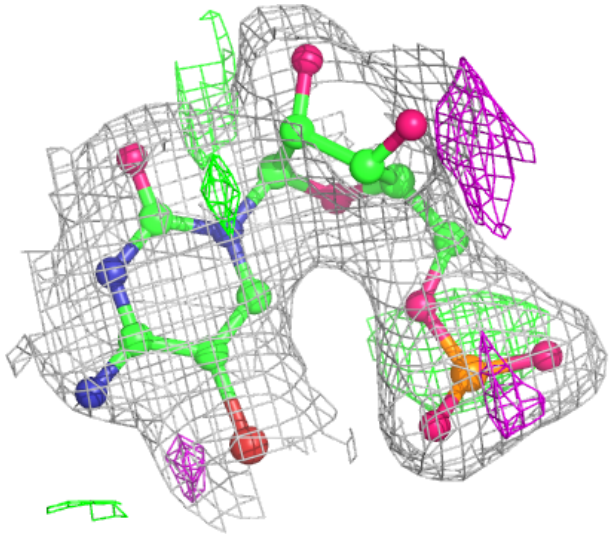


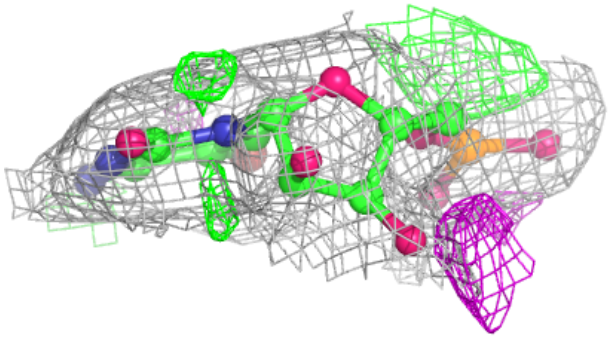


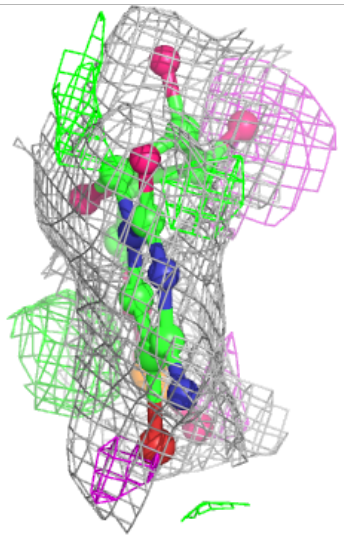


Electron density around CBV M 101:

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

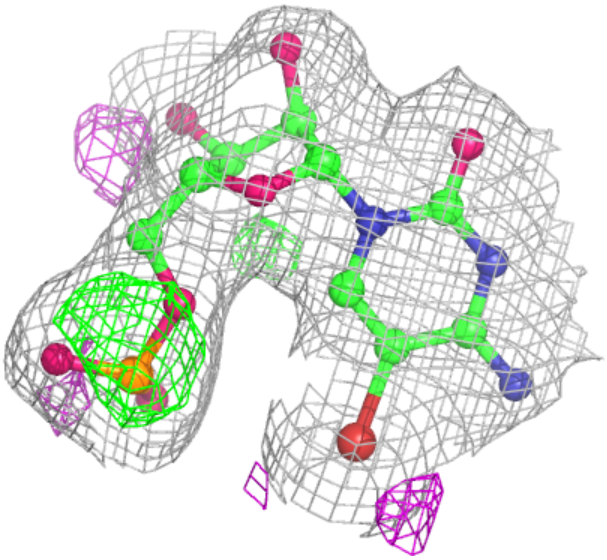


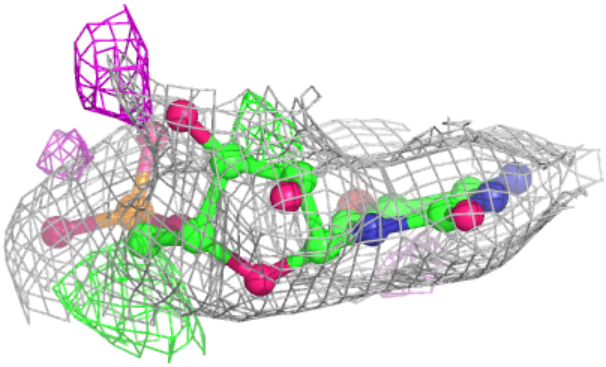




Electron density around CBV F 101:

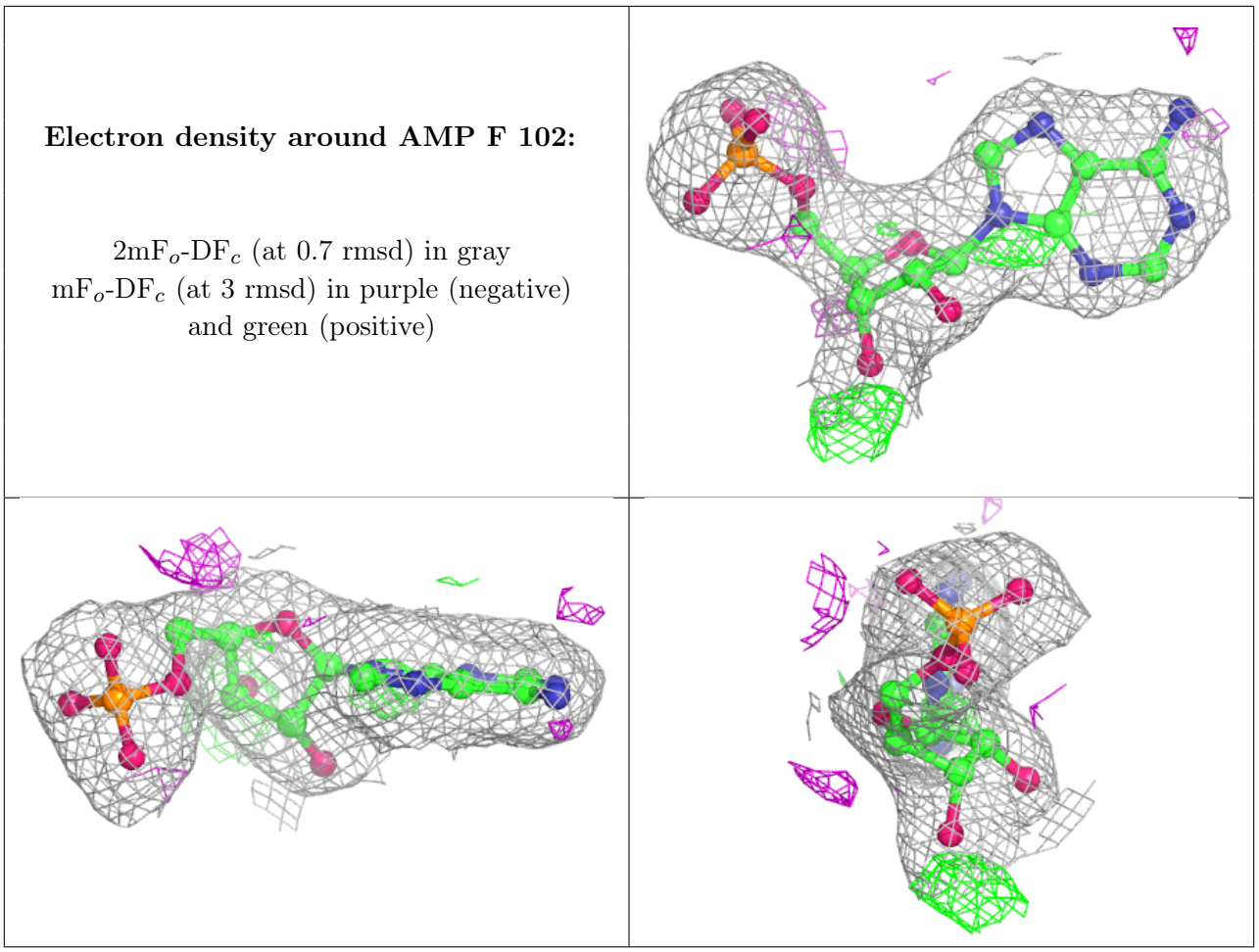
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)





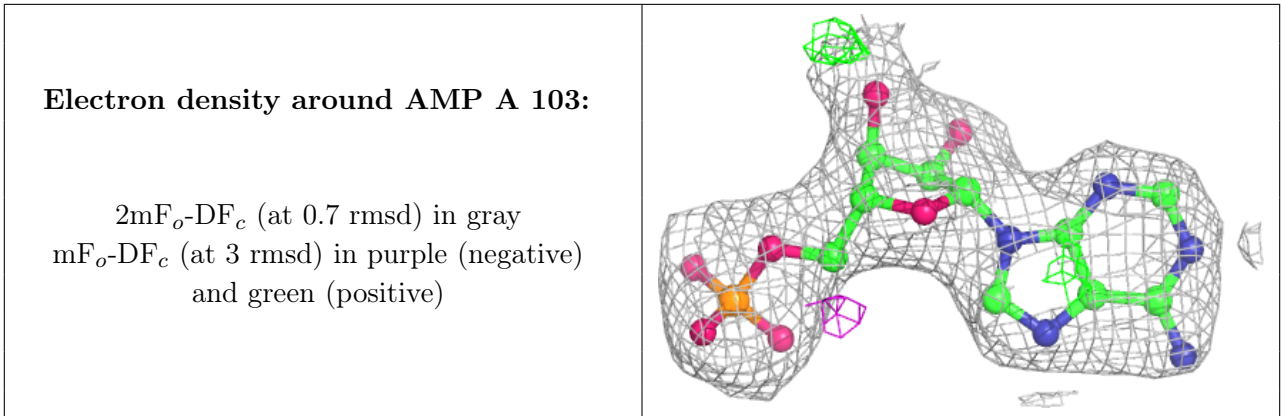
Electron density around AMP F 102:

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



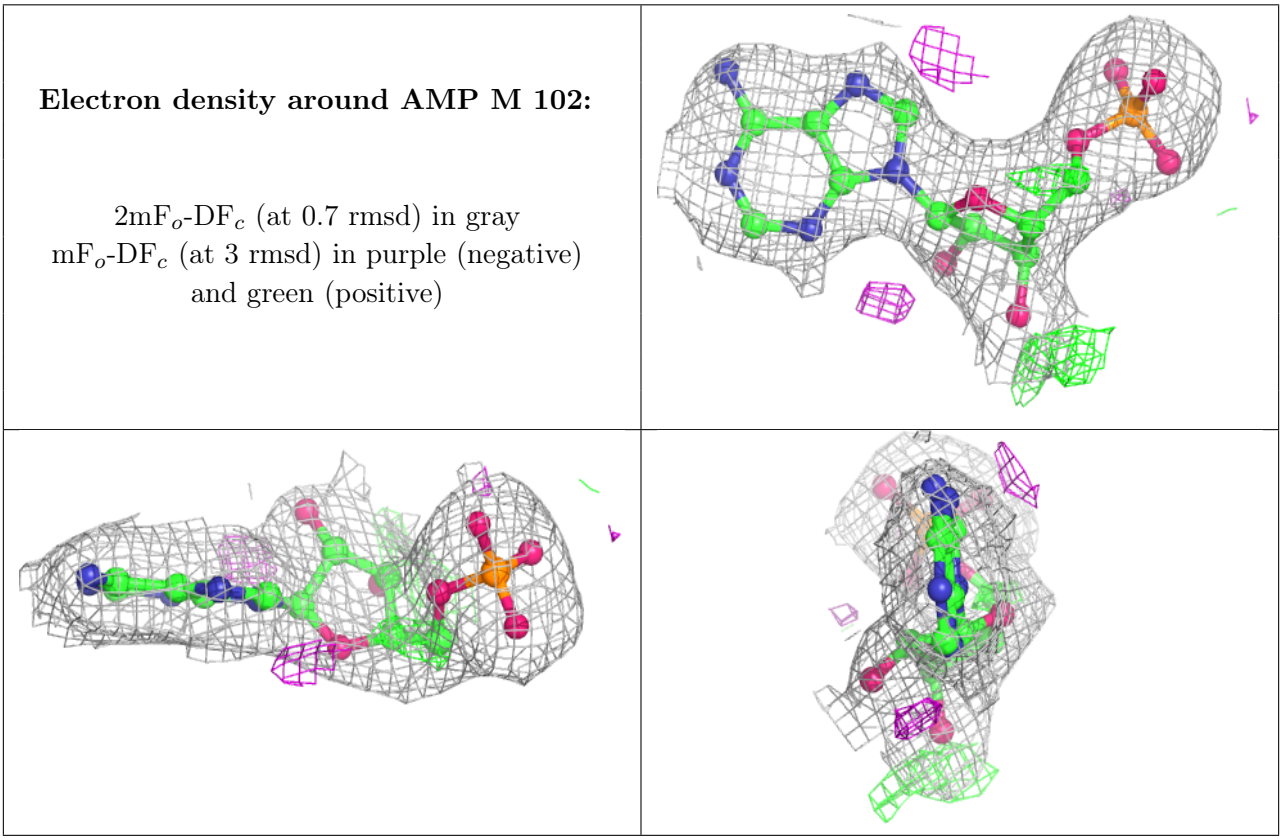
Electron density around AMP A 103:

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



Electron density around AMP M 102:

$2mF_o$-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.