



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 02:27 PM UTC

PDB ID : 4Q86 / pdb_00004q86
Title : YcaO with AMP Bound
Authors : Chekan, J.R.; Nair, S.K.
Deposited on : 2014-04-25
Resolution : 2.25 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

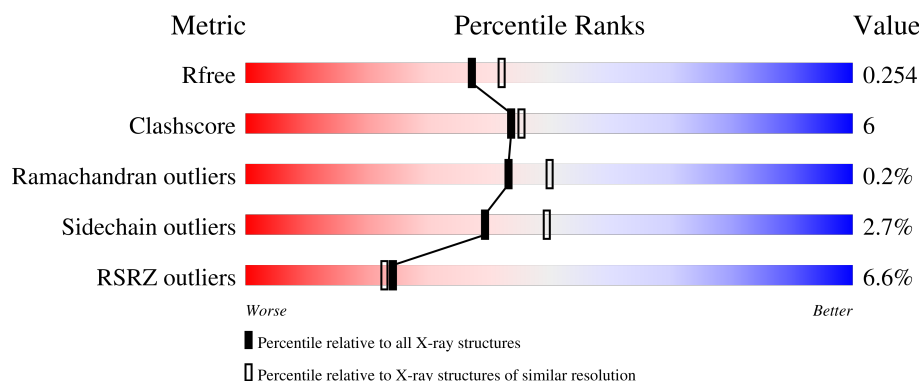
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	586	91% 7% ..
1	B	586	2% 90% 8% ...
1	C	586	2% 87% 9% ..
1	D	586	2% 90% 8% ..
1	E	586	3% 88% 10% ..

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Mol	Chain	Length	Quality of chain
1	F	586	40% 82% 13% ..
1	G	586	4% 85% 11% ..
1	H	586	87% 10% ...

2 Entry composition

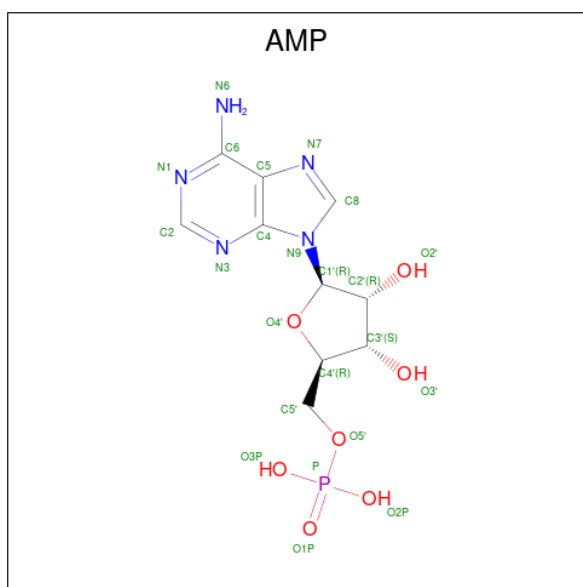
There are 4 unique types of molecules in this entry. The entry contains 38743 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ribosomal protein S12 methylthiotransferase accessory factor YcaO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	582	Total	C	N	O	S	0	0	0
			4599	2935	748	899	17			
1	B	583	Total	C	N	O	S	0	0	0
			4613	2946	750	900	17			
1	C	570	Total	C	N	O	S	0	0	0
			4508	2875	734	882	17			
1	D	582	Total	C	N	O	S	0	0	0
			4599	2935	748	899	17			
1	E	578	Total	C	N	O	S	0	0	0
			4567	2914	743	893	17			
1	F	574	Total	C	N	O	S	0	0	0
			4534	2890	739	888	17			
1	G	575	Total	C	N	O	S	0	0	0
			4541	2895	740	889	17			
1	H	583	Total	C	N	O	S	0	0	0
			4613	2946	750	900	17			

- Molecule 2 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: C₁₀H₁₄N₅O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	D	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	E	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	G	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	H	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Mg	0	0
			3	3		
3	B	3	Total	Mg	0	0
			3	3		
3	C	3	Total	Mg	0	0
			3	3		
3	D	3	Total	Mg	0	0
			3	3		
3	E	3	Total	Mg	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	3	Total 3	Mg 3	0	0
3	H	3	Total 3	Mg 3	0	0

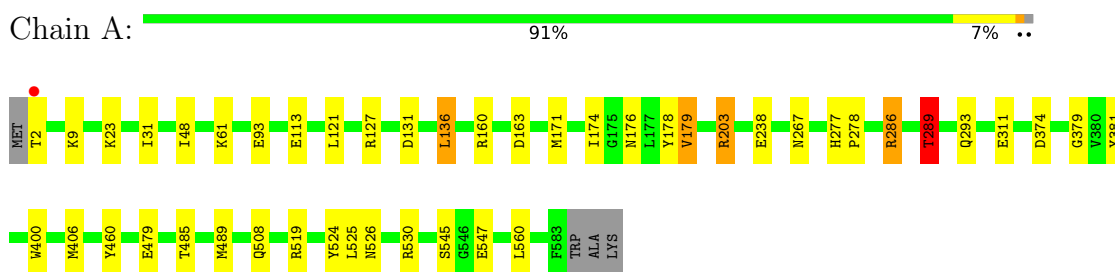
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	334	Total 334	O 334	0	0
4	B	336	Total 336	O 336	0	0
4	C	231	Total 231	O 231	0	0
4	D	255	Total 255	O 255	0	0
4	E	215	Total 215	O 215	0	0
4	F	111	Total 111	O 111	0	0
4	G	197	Total 197	O 197	0	0
4	H	308	Total 308	O 308	0	0

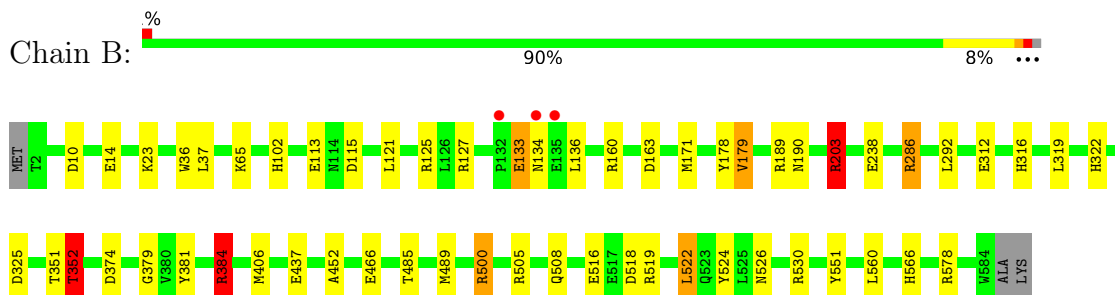
3 Residue-property plots [i](#)

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.

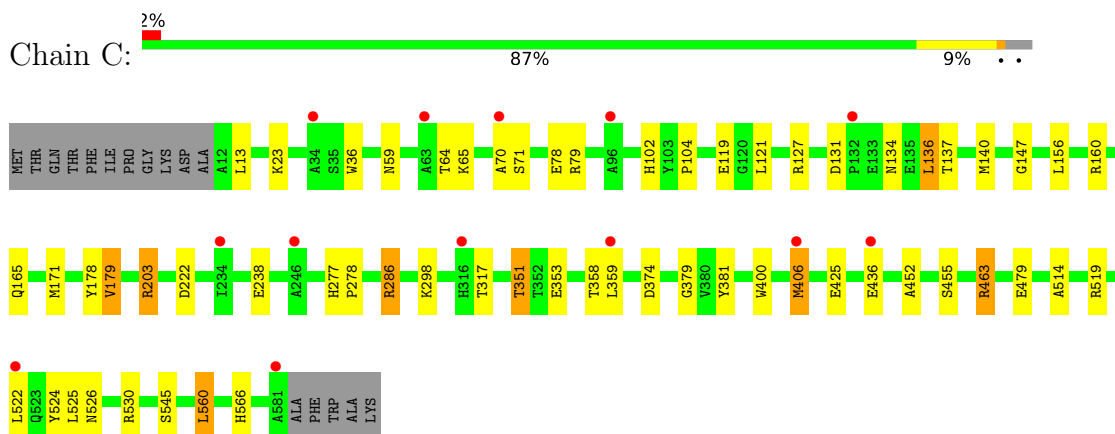
- Molecule 1: Ribosomal protein S12 methylthiotransferase accessory factor YcaO



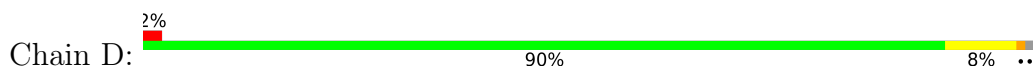
- Molecule 1: Ribosomal protein S12 methylthiotransferase accessory factor YcaO

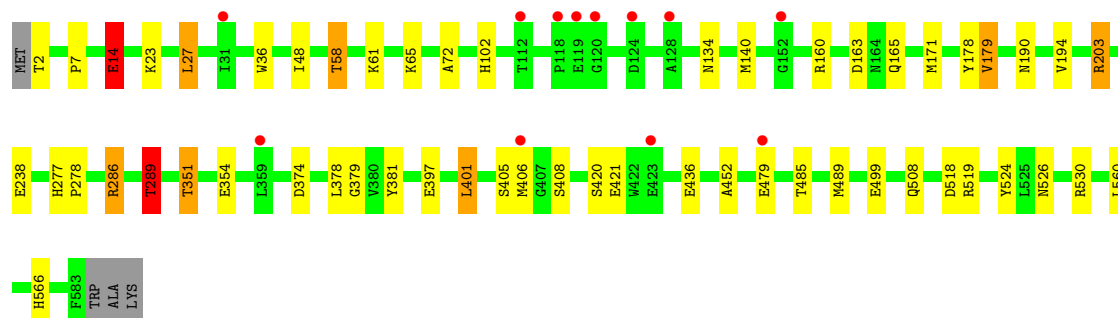


- Molecule 1: Ribosomal protein S12 methylthiotransferase accessory factor YcaO



- Molecule 1: Ribosomal protein S12 methylthiotransferase accessory factor YcaO



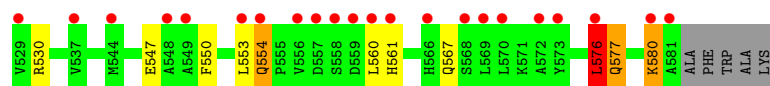


- Molecule 1: Ribosomal protein S12 methylthiotransferase accessory factor YcaO

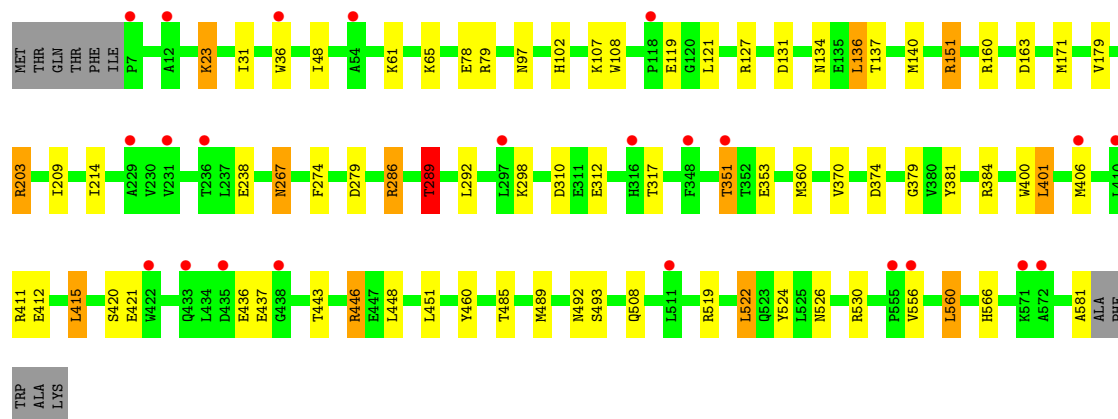
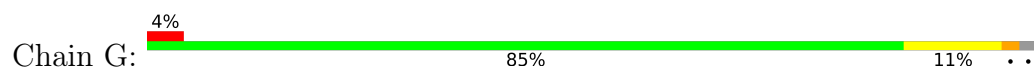
Chain E:

- Molecule 1: Ribosomal protein S12 methylthiotransferase accessory factor YcaO

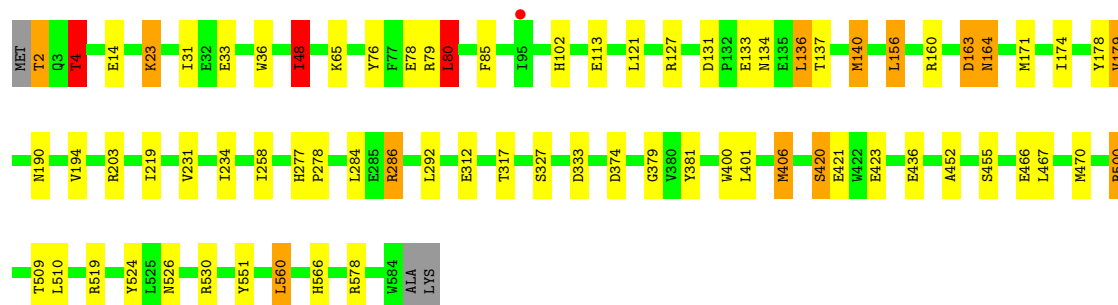
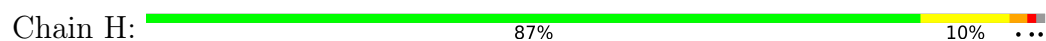
Chain F:



- Molecule 1: Ribosomal protein S12 methylthiotransferase accessory factor YcaO



- Molecule 1: Ribosomal protein S12 methylthiotransferase accessory factor YcaO



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	111.25Å 112.95Å 132.89Å 89.94° 73.51° 77.23°	Depositor
Resolution (Å)	127.14 – 2.25 127.14 – 2.25	Depositor EDS
% Data completeness (in resolution range)	93.1 (127.14-2.25) 93.1 (127.14-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.25Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.218 , 0.250 0.221 , 0.254	Depositor DCC
R_{free} test set	13386 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.223	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	38743	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.86	0/4713	0.86	6/6411 (0.1%)
1	B	0.90	0/4729	0.91	9/6434 (0.1%)
1	C	0.74	0/4619	0.87	7/6283 (0.1%)
1	D	0.81	0/4713	0.87	10/6411 (0.2%)
1	E	0.77	0/4680	0.86	8/6366 (0.1%)
1	F	0.68	0/4645	0.92	13/6317 (0.2%)
1	G	0.77	0/4653	0.89	12/6328 (0.2%)
1	H	0.90	2/4729 (0.0%)	0.95	14/6434 (0.2%)
All	All	0.81	2/37481 (0.0%)	0.89	79/50984 (0.2%)

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	F	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	164	ASN	N-CA	5.51	1.54	1.46
1	H	164	ASN	CA-C	-5.05	1.44	1.52

The worst 5 of 79 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	203	ARG	NE-CZ-NH2	16.54	134.09	119.20
1	H	203	ARG	NE-CZ-NH1	-11.88	109.62	121.50
1	F	119	GLU	N-CA-C	-10.11	94.24	110.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	384	ARG	CG-CD-NE	10.10	134.21	112.00
1	F	122	LEU	N-CA-C	-9.64	100.88	108.78

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	293	GLN	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4599	0	4372	36	0
1	B	4613	0	4382	51	0
1	C	4508	0	4284	49	0
1	D	4599	0	4372	46	0
1	E	4567	0	4343	51	0
1	F	4534	0	4309	69	0
1	G	4541	0	4317	66	0
1	H	4613	0	4382	77	0
2	A	23	0	12	1	0
2	B	23	0	12	0	0
2	C	23	0	12	1	0
2	D	23	0	12	2	0
2	E	23	0	12	2	0
2	G	23	0	12	0	0
2	H	23	0	12	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
3	D	3	0	0	0	0
3	E	3	0	0	0	0
3	G	3	0	0	0	0
3	H	3	0	0	0	0
4	A	334	0	0	7	0
4	B	336	0	0	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	231	0	0	16	0
4	D	255	0	0	16	0
4	E	215	0	0	19	0
4	F	111	0	0	24	0
4	G	197	0	0	16	0
4	H	308	0	0	25	0
All	All	38743	0	34845	434	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 6.

The worst 5 of 434 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:F:68:ALA:HB2	4:F:633:HOH:O	1.36	1.26
1:E:171:MET:SD	4:E:738:HOH:O	2.04	1.15
1:B:171:MET:SD	4:B:1021:HOH:O	2.02	1.15
1:G:78:GLU:OE2	1:G:79:ARG:NH1	1.86	1.08
1:F:550:PHE:HB3	1:F:553:LEU:HD12	1.33	1.06

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	580/586 (99%)	569 (98%)	10 (2%)	1 (0%)	43	50
1	B	581/586 (99%)	567 (98%)	13 (2%)	1 (0%)	43	50
1	C	568/586 (97%)	555 (98%)	12 (2%)	1 (0%)	43	50
1	D	580/586 (99%)	566 (98%)	13 (2%)	1 (0%)	43	50
1	E	576/586 (98%)	564 (98%)	11 (2%)	1 (0%)	43	50

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	572/586 (98%)	557 (97%)	14 (2%)	1 (0%)	43	50
1	G	573/586 (98%)	560 (98%)	12 (2%)	1 (0%)	43	50
1	H	581/586 (99%)	566 (97%)	14 (2%)	1 (0%)	43	50
All	All	4611/4688 (98%)	4504 (98%)	99 (2%)	8 (0%)	43	50

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	179	VAL
1	B	179	VAL
1	C	179	VAL
1	D	179	VAL
1	E	179	VAL

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	484/487 (99%)	477 (99%)	7 (1%)	59	70
1	B	485/487 (100%)	475 (98%)	10 (2%)	47	58
1	C	475/487 (98%)	463 (98%)	12 (2%)	42	52
1	D	484/487 (99%)	472 (98%)	12 (2%)	42	52
1	E	481/487 (99%)	470 (98%)	11 (2%)	44	55
1	F	477/487 (98%)	456 (96%)	21 (4%)	25	29
1	G	478/487 (98%)	461 (96%)	17 (4%)	31	39
1	H	485/487 (100%)	470 (97%)	15 (3%)	35	44
All	All	3849/3896 (99%)	3744 (97%)	105 (3%)	39	49

5 of 105 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	164	ASN

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Mol	Chain	Res	Type
1	F	577	GLN
1	H	292	LEU
1	F	286	ARG
1	F	437	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	561	HIS
1	G	567	GLN
1	C	364	ASN
1	C	59	ASN
1	H	195	GLN

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](#)

Of 28 ligands modelled in this entry, 21 are monoatomic - leaving 7 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
2	AMP	A	601	-	25,25,25	2.04	6 (24%)	37,38,38	2.48	13 (35%)
2	AMP	B	601	-	25,25,25	1.74	4 (16%)	37,38,38	1.94	8 (21%)
2	AMP	H	601	-	25,25,25	1.42	5 (20%)	37,38,38	2.31	9 (24%)
2	AMP	E	601	-	25,25,25	1.85	4 (16%)	37,38,38	2.03	10 (27%)
2	AMP	G	601	-	25,25,25	1.56	4 (16%)	37,38,38	2.20	15 (40%)
2	AMP	C	601	-	25,25,25	1.54	5 (20%)	37,38,38	2.16	13 (35%)
2	AMP	D	601	-	25,25,25	2.02	6 (24%)	37,38,38	2.12	12 (32%)

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	AMP	A	601	-	-	0/10/26/26	0/3/3/3
2	AMP	B	601	-	-	0/10/26/26	0/3/3/3
2	AMP	H	601	-	-	0/10/26/26	0/3/3/3
2	AMP	E	601	-	-	0/10/26/26	0/3/3/3
2	AMP	G	601	-	-	0/10/26/26	0/3/3/3
2	AMP	C	601	-	-	2/10/26/26	0/3/3/3
2	AMP	D	601	-	-	0/10/26/26	0/3/3/3

The worst 5 of 34 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	AMP	C5-C4	6.24	1.50	1.39
2	B	601	AMP	C5-C4	5.76	1.49	1.39
2	A	601	AMP	C4-N9	-5.61	1.26	1.37
2	E	601	AMP	C4-N9	-5.17	1.26	1.37
2	A	601	AMP	C5-C4	5.04	1.48	1.39

The worst 5 of 80 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	AMP	C2-N1-C6	7.32	130.75	118.73
2	H	601	AMP	C5-C4-N3	-6.81	117.34	126.72
2	E	601	AMP	C2-N1-C6	6.67	129.68	118.73
2	B	601	AMP	C5-C4-N3	-6.03	118.41	126.72
2	A	601	AMP	N3-C2-N1	-5.32	120.53	128.58

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	601	AMP	C5'-O5'-P-O2P
2	C	601	AMP	C5'-O5'-P-O3P

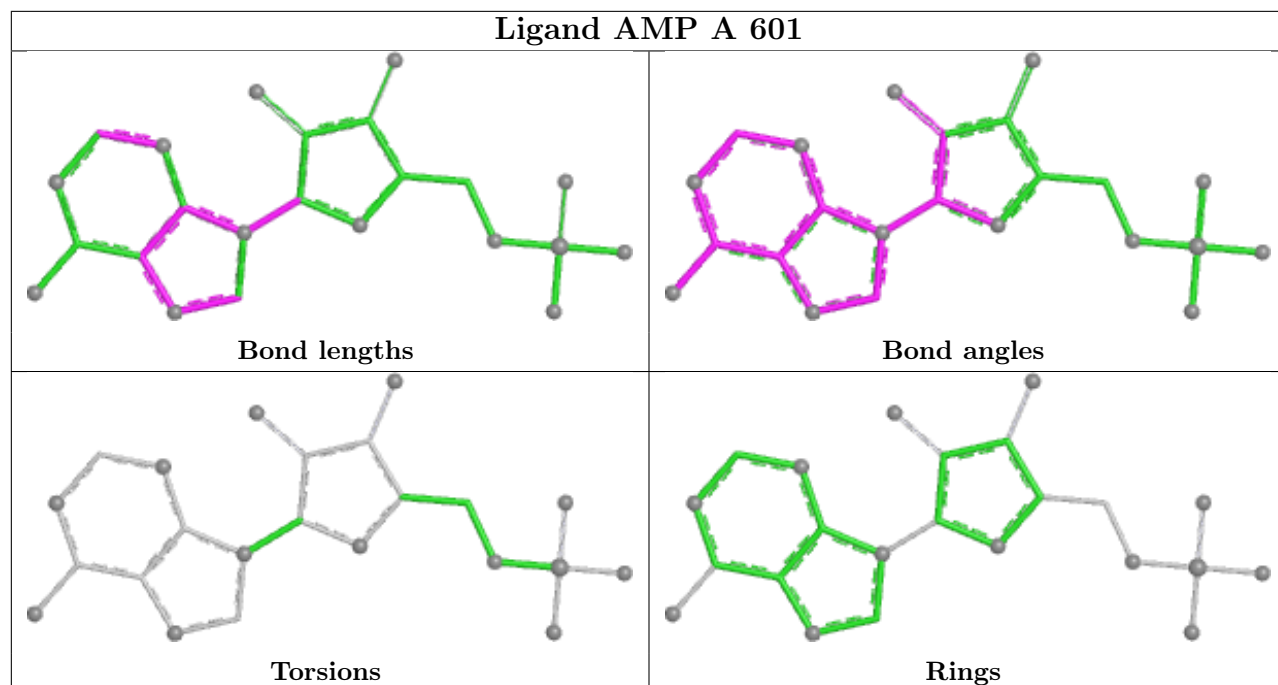
There are no ring outliers.

4 monomers are involved in 6 short contacts:

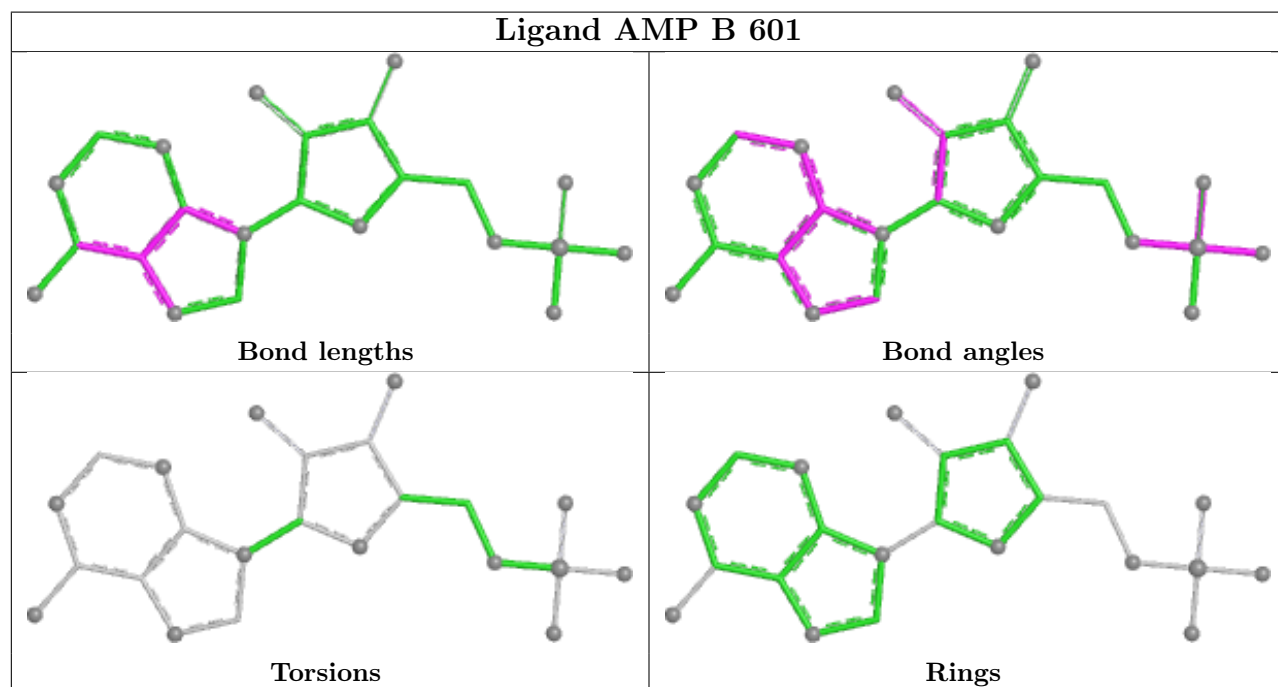
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	AMP	1	0
2	E	601	AMP	2	0
2	C	601	AMP	1	0
2	D	601	AMP	2	0

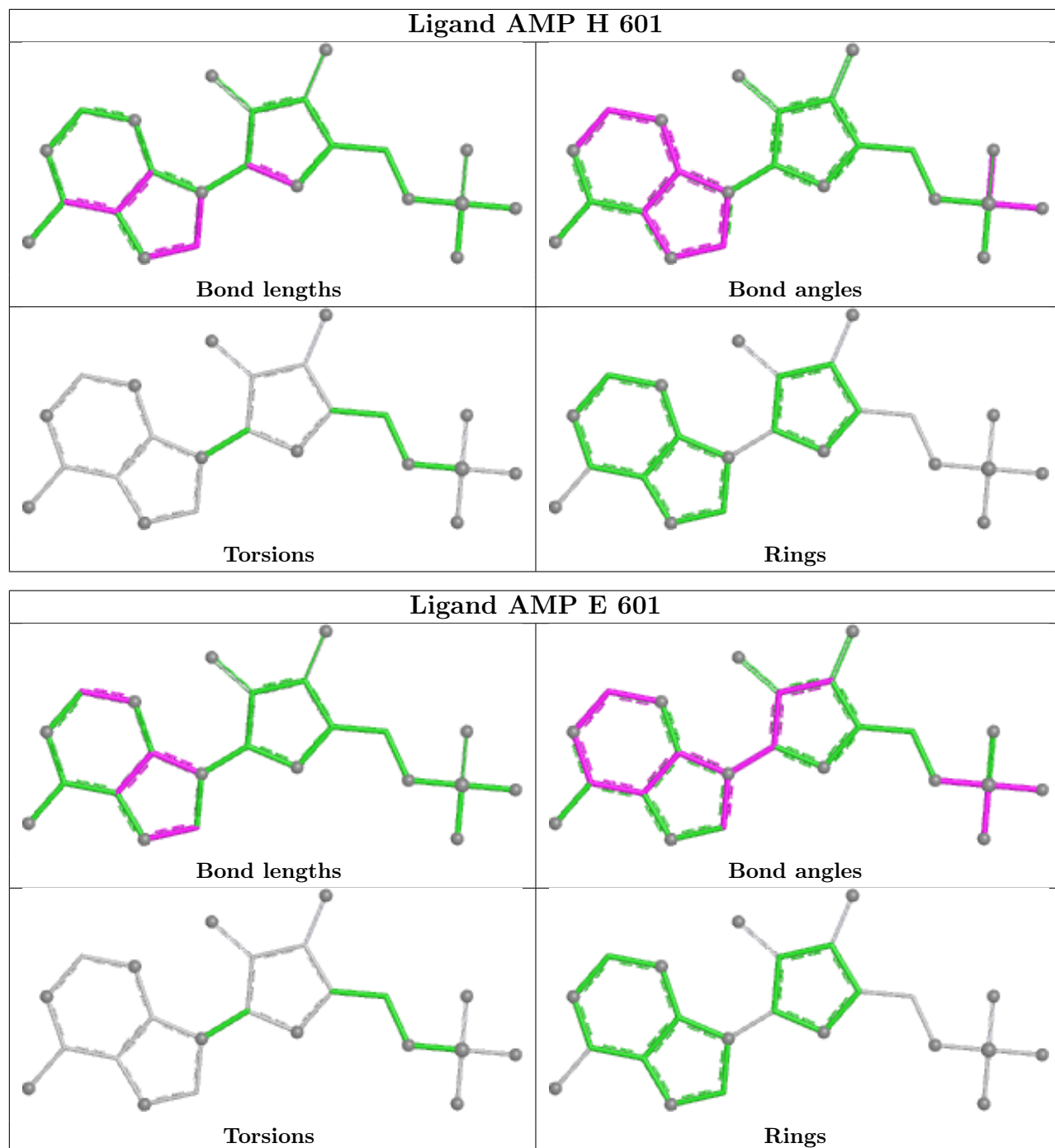
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand AMP A 601

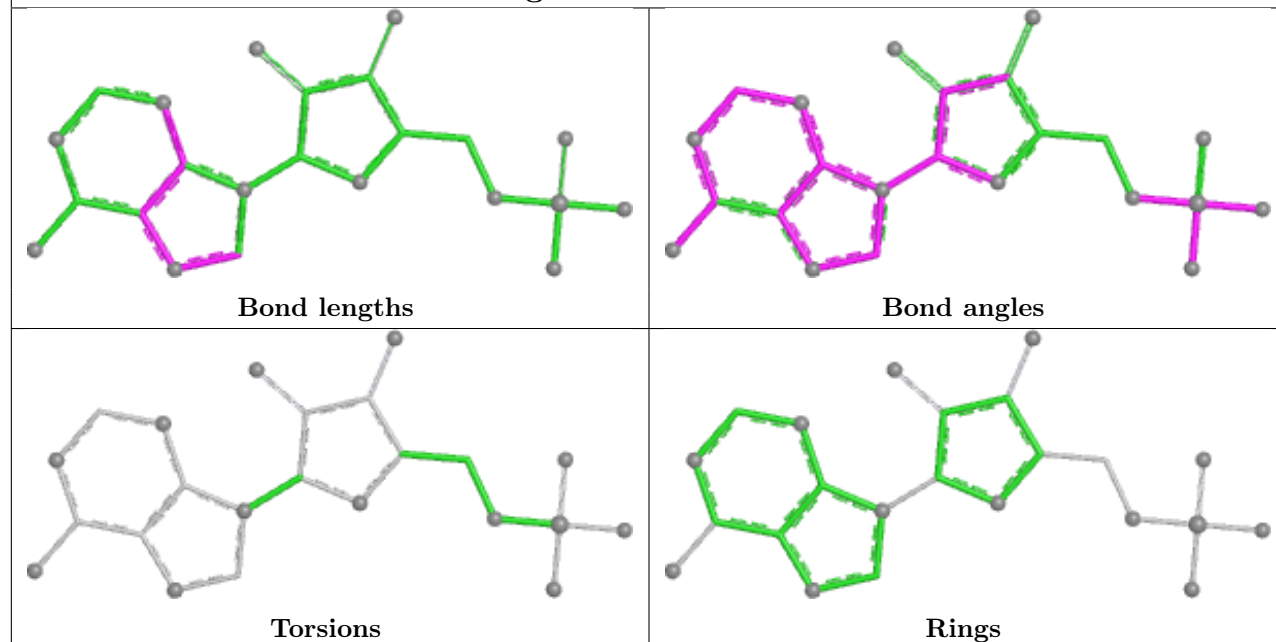


Ligand AMP B 601

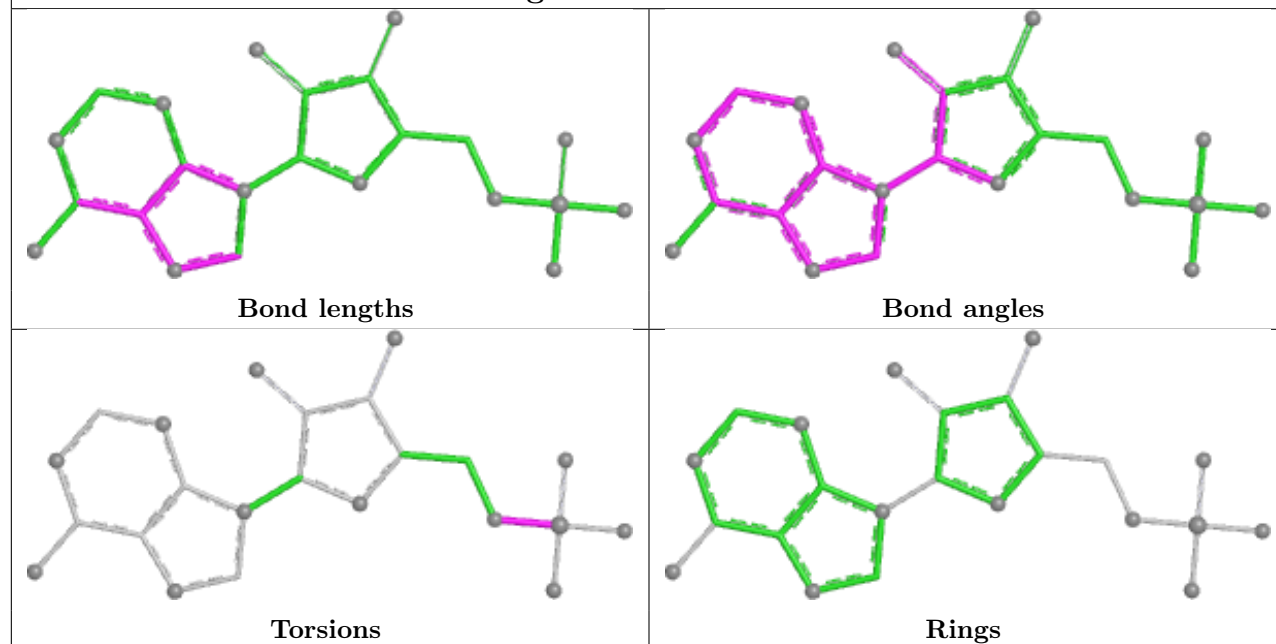


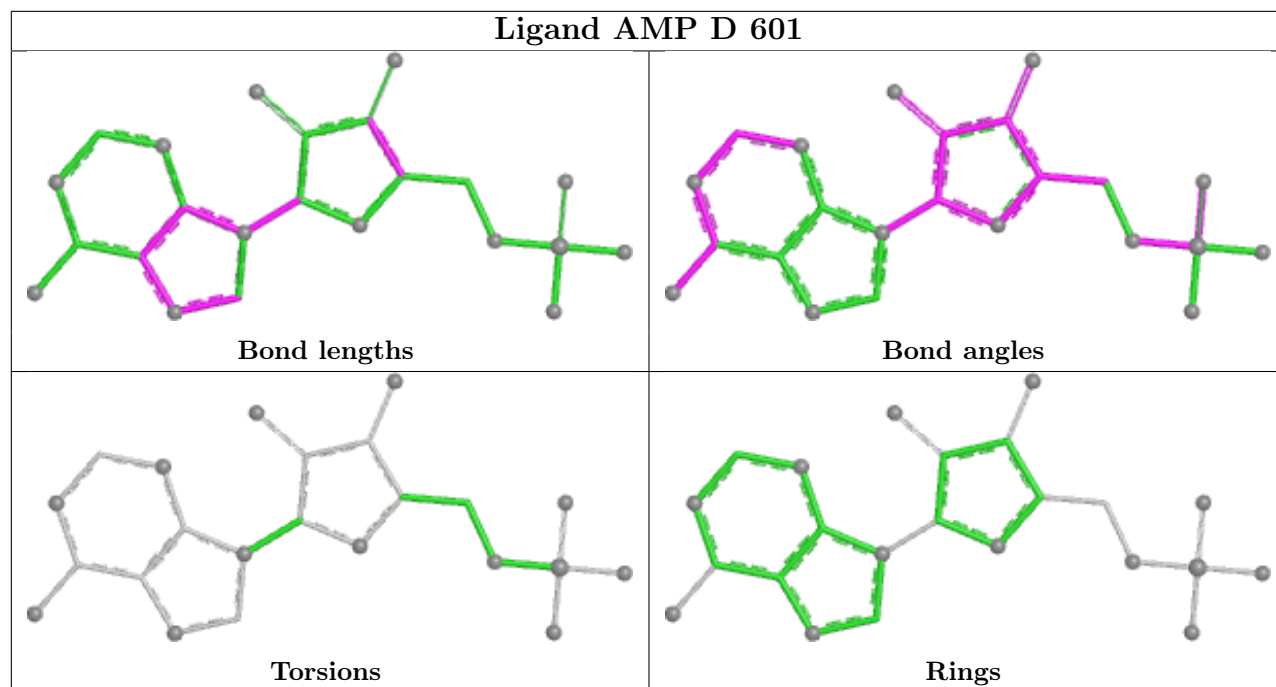


Ligand AMP G 601



Ligand AMP C 601





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	582/586 (99%)	-0.11	1 (0%) 91 92	17, 28, 45, 64	0
1	B	583/586 (99%)	-0.09	3 (0%) 87 88	15, 26, 47, 82	0
1	C	570/586 (97%)	0.43	13 (2%) 61 61	22, 39, 57, 93	0
1	D	582/586 (99%)	0.32	12 (2%) 63 64	21, 36, 57, 86	0
1	E	578/586 (98%)	0.59	19 (3%) 49 49	24, 40, 63, 87	0
1	F	574/586 (97%)	1.83	233 (40%) 0 0	35, 66, 93, 111	0
1	G	575/586 (98%)	0.56	23 (4%) 42 42	23, 40, 60, 92	0
1	H	583/586 (99%)	0.01	1 (0%) 91 92	14, 30, 49, 74	0
All	All	4627/4688 (98%)	0.44	305 (6%) 24 23	14, 36, 71, 111	0

The worst 5 of 305 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	185	ALA	5.0
1	F	297	LEU	4.8
1	F	186	GLY	4.5
1	F	341	TYR	4.5
1	F	359	LEU	4.3

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 oligosaccharides in this entry.

6.4 Ligands ⓘ

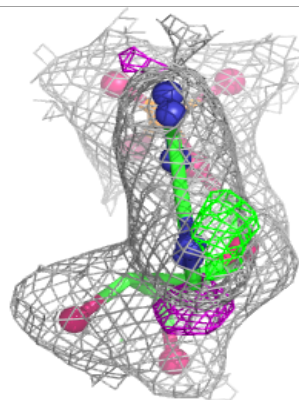
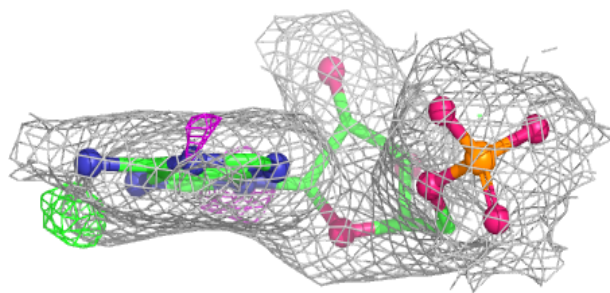
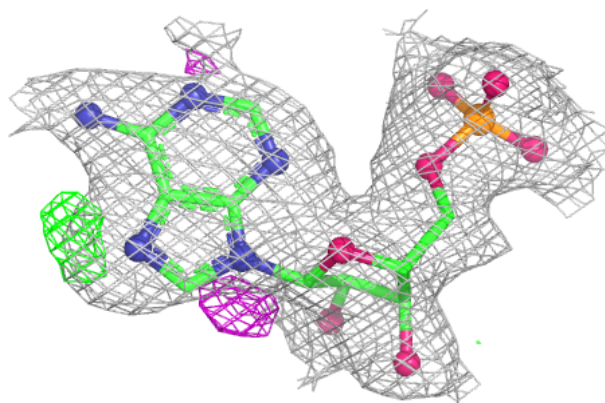
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	AMP	E	601	23/23	0.94	0.09	23,31,35,40	0
3	MG	E	602	1/1	0.94	0.05	27,27,27,27	0
3	MG	G	604	1/1	0.94	0.13	49,49,49,49	0
2	AMP	D	601	23/23	0.95	0.10	18,25,44,57	0
2	AMP	C	601	23/23	0.95	0.09	30,38,56,61	0
2	AMP	A	601	23/23	0.96	0.08	19,24,30,37	0
2	AMP	G	601	23/23	0.96	0.07	28,36,42,45	0
3	MG	H	604	1/1	0.96	0.04	29,29,29,29	0
3	MG	D	602	1/1	0.97	0.06	15,15,15,15	0
2	AMP	B	601	23/23	0.98	0.05	16,19,21,23	0
3	MG	D	603	1/1	0.98	0.07	28,28,28,28	0
3	MG	D	604	1/1	0.98	0.04	34,34,34,34	0
2	AMP	H	601	23/23	0.98	0.05	16,19,23,25	0
3	MG	C	603	1/1	0.98	0.03	26,26,26,26	0
3	MG	C	604	1/1	0.98	0.05	36,36,36,36	0
3	MG	A	602	1/1	0.99	0.03	14,14,14,14	0
3	MG	A	603	1/1	0.99	0.01	18,18,18,18	0
3	MG	E	604	1/1	0.99	0.03	39,39,39,39	0
3	MG	G	602	1/1	0.99	0.05	19,19,19,19	0
3	MG	G	603	1/1	0.99	0.03	34,34,34,34	0
3	MG	A	604	1/1	0.99	0.03	31,31,31,31	0
3	MG	H	602	1/1	0.99	0.02	15,15,15,15	0
3	MG	H	603	1/1	0.99	0.04	20,20,20,20	0
3	MG	C	602	1/1	0.99	0.05	17,17,17,17	0
3	MG	E	603	1/1	1.00	0.02	34,34,34,34	0
3	MG	B	603	1/1	1.00	0.03	17,17,17,17	0
3	MG	B	604	1/1	1.00	0.04	24,24,24,24	0
3	MG	B	602	1/1	1.00	0.06	7,7,7,7	0

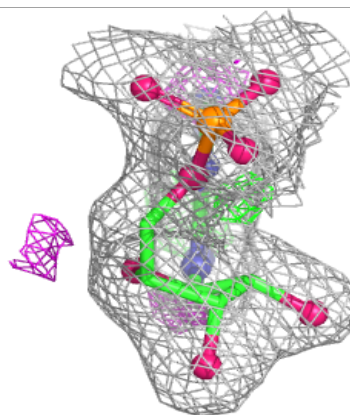
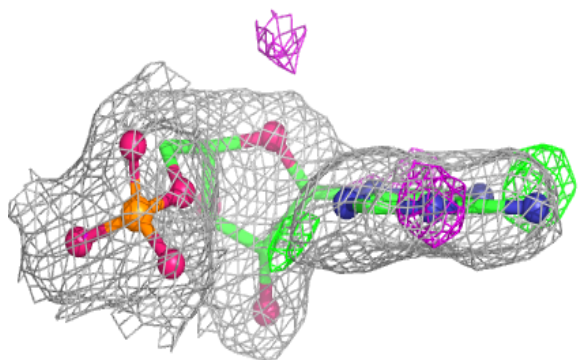
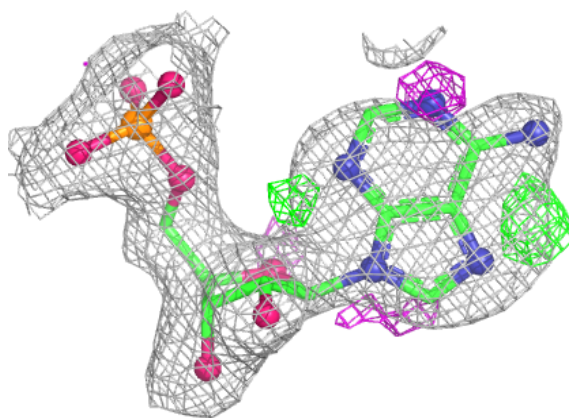
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around AMP E 601:

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

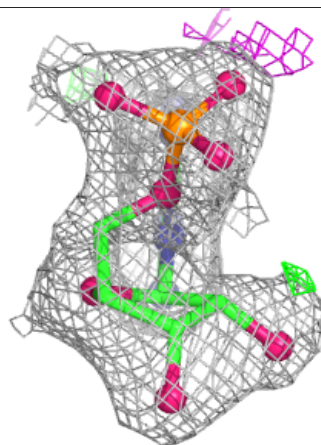
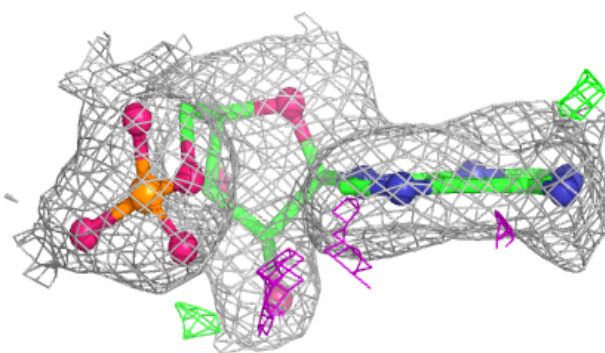
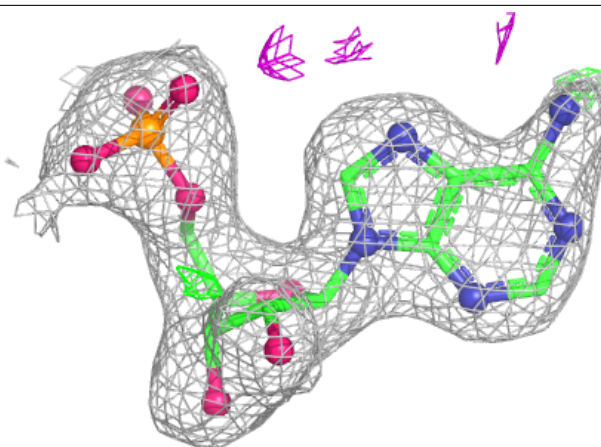
**Electron density around AMP D 601:**

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



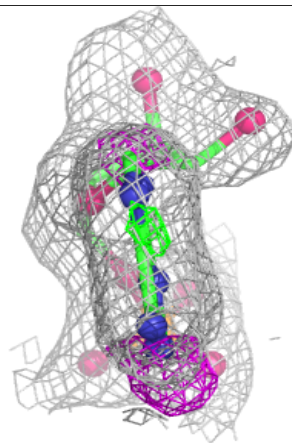
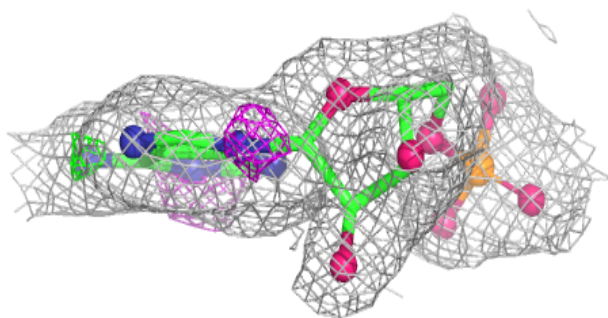
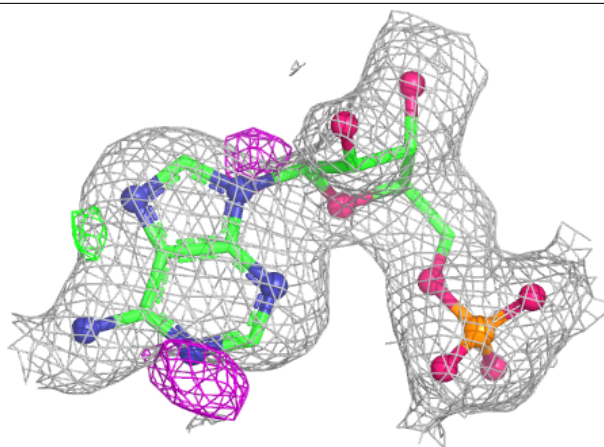
Electron density around AMP C 601:

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



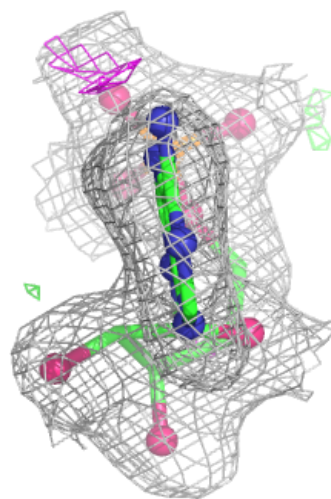
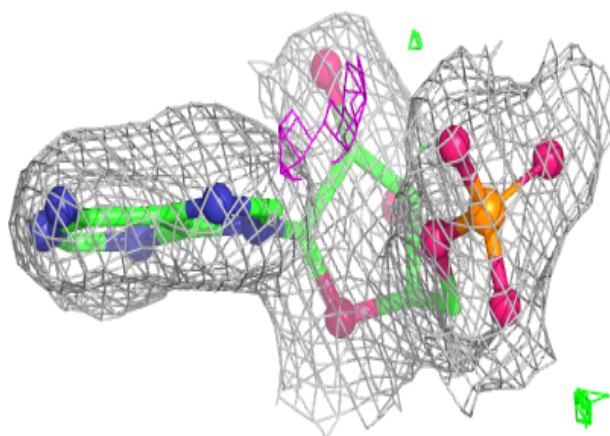
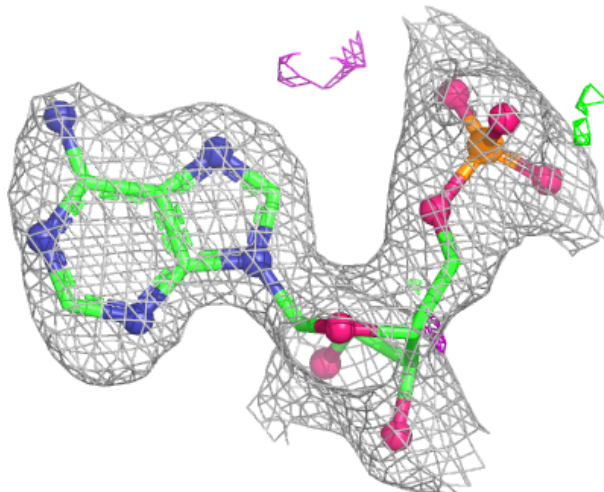
Electron density around AMP A 601:

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



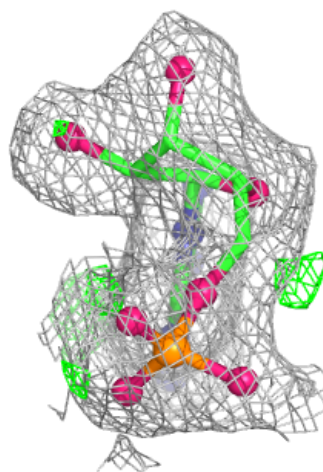
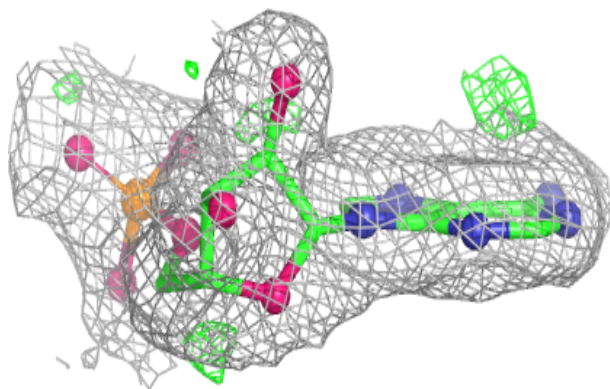
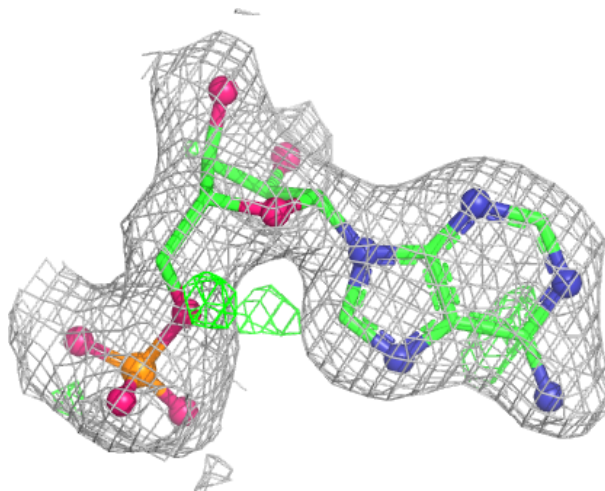
Electron density around AMP G 601:

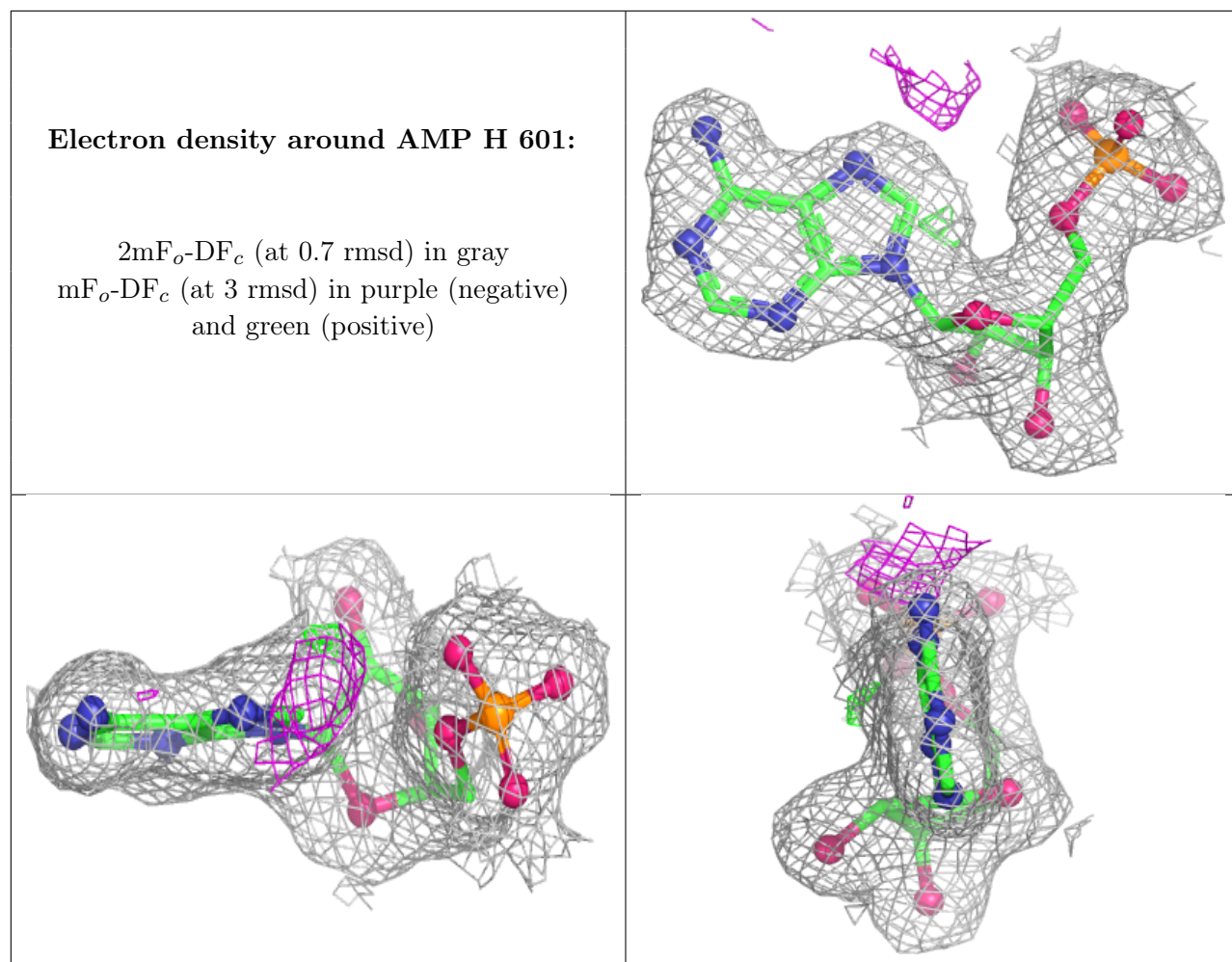
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around AMP B 601:

$2mF_o-DF_c$ (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.