



# wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 24, 2023 – 12:38 AM EDT

PDB ID : 5KOD  
Title : Crystal Structure of GH3.5 Acyl Acid Amido Synthetase from *Arabidopsis thaliana*  
Authors : Jez, J.M.; Westfall, C.S.; Zubieta, C.  
Deposited on : 2016-06-30  
Resolution : 2.20 Å (reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

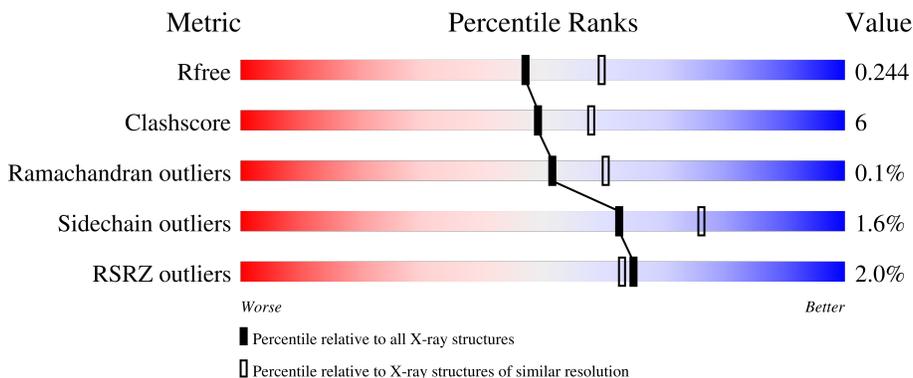
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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

Mol	Chain	Length	Quality of chain
1	A	612	 2% 83% 10% • 6%
1	B	612	 % 82% 11% • 6%
1	C	612	 3% 79% 13% • 8%
1	D	612	 % 82% 10% 8%

## 2 Entry composition [i](#)

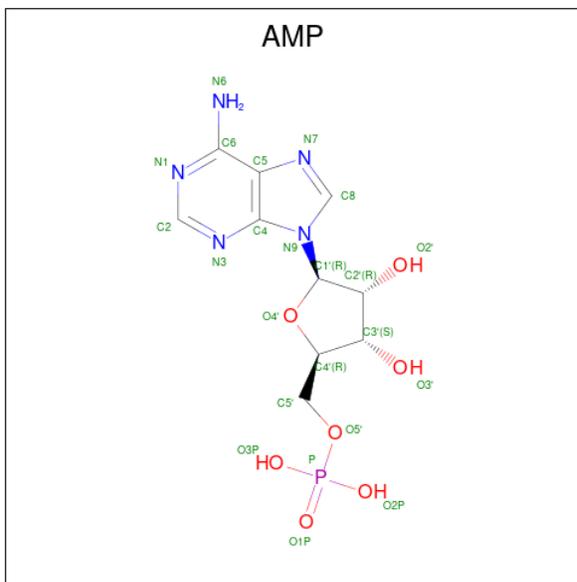
There are 5 unique types of molecules in this entry. The entry contains 19306 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Indole-3-acetic acid-amido synthetase GH3.5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	573	Total	C	N	O	S	0	3	0
			4565	2920	750	870	25			
1	B	573	Total	C	N	O	S	0	0	0
			4570	2924	749	872	25			
1	C	566	Total	C	N	O	S	0	2	0
			4507	2882	740	860	25			
1	D	566	Total	C	N	O	S	0	1	0
			4509	2888	739	856	26			

- Molecule 2 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>5</sub>O<sub>7</sub>P).



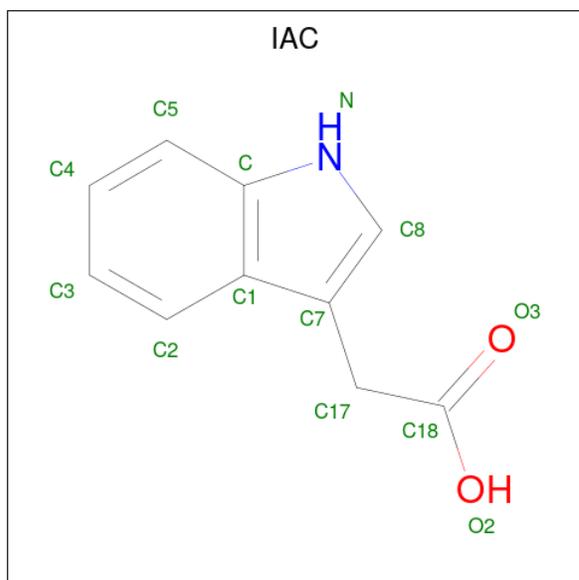
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
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		

- Molecule 3 is 1H-INDOL-3-YLACETIC ACID (three-letter code: IAC) (formula: C<sub>10</sub>H<sub>9</sub>NO<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	10	1	2		
3	B	1	Total	C	N	O	0	0
			13	10	1	2		
3	C	1	Total	C	N	O	0	0
			13	10	1	2		
3	D	1	Total	C	N	O	0	0
			13	10	1	2		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).

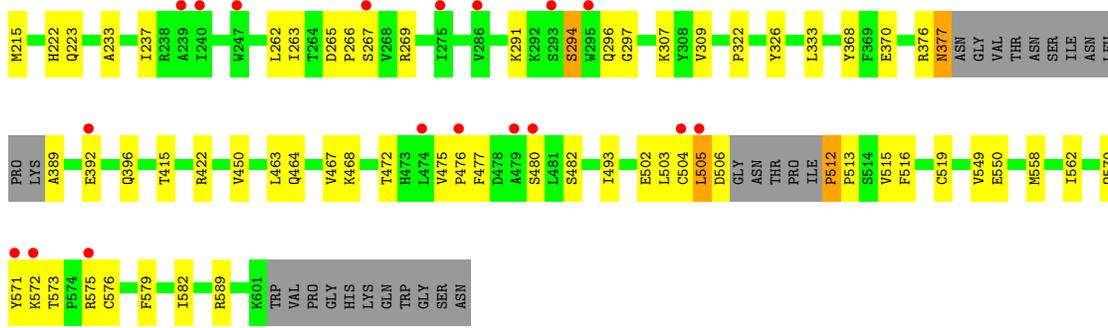


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

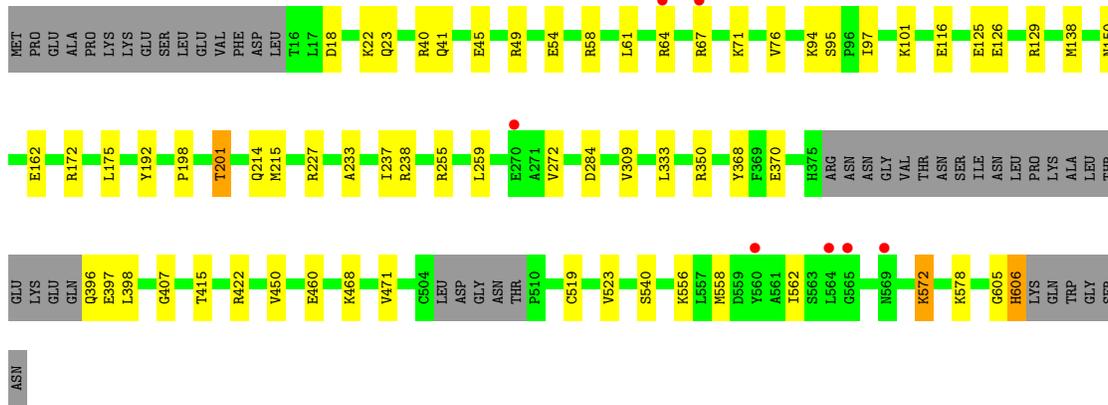
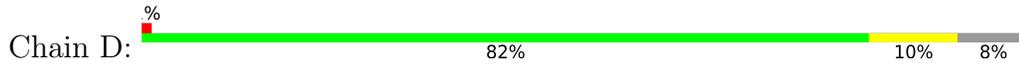
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	282	Total	O	0	0
			282	282		
5	B	260	Total	O	0	0
			260	260		
5	C	215	Total	O	0	0
			215	215		
5	D	244	Total	O	0	0
			244	244		





● Molecule 1: Indole-3-acetic acid-amido synthetase GH3.5



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.61Å 143.52Å 102.32Å 90.00° 114.72° 90.00°	Depositor
Resolution (Å)	42.16 – 2.20 42.16 – 2.20	Depositor EDS
% Data completeness (in resolution range)	89.5 (42.16-2.20) 89.5 (42.16-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.80 (at 2.20Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.205 , 0.245 0.207 , 0.244	Depositor DCC
$R_{free}$ test set	5422 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.7	Xtrriage
Anisotropy	0.125	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 42.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.033 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	19306	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.65% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: IAC, SO4, AMP

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	0.45	0/4679	0.58	0/6352
1	B	0.53	2/4673 (0.0%)	0.62	5/6342 (0.1%)
1	C	0.57	2/4609 (0.0%)	0.69	8/6250 (0.1%)
1	D	0.45	0/4616	0.59	3/6266 (0.0%)
All	All	0.50	4/18577 (0.0%)	0.62	16/25210 (0.1%)

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	C	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	571	TYR	C-N	6.33	1.48	1.34
1	B	501	TRP	NE1-CE2	-5.60	1.30	1.37
1	C	513	PRO	N-CD	5.24	1.55	1.47
1	C	476	PRO	N-CD	5.18	1.55	1.47

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	184	PHE	O-C-N	-16.38	96.49	122.70
1	C	184	PHE	C-N-CA	10.63	148.26	121.70
1	C	184	PHE	CA-C-N	8.44	135.77	117.20
1	C	505	LEU	CA-CB-CG	-6.89	99.46	115.30
1	C	512	PRO	C-N-CD	6.48	142.01	128.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	184	PHE	Mainchain

## 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	4565	0	4547	57	0
1	B	4570	0	4563	38	0
1	C	4507	0	4510	68	3
1	D	4509	0	4495	52	3
2	A	23	0	12	0	0
2	B	23	0	12	0	0
2	C	23	0	12	0	0
2	D	23	0	12	0	0
3	A	13	0	8	2	0
3	B	13	0	8	3	0
3	C	13	0	8	2	0
3	D	13	0	8	2	0
4	B	10	0	0	0	0
5	A	282	0	0	9	1
5	B	260	0	0	5	0
5	C	215	0	0	5	1
5	D	244	0	0	6	0
All	All	19306	0	18195	215	4

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 215 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:C:480:SER:OG	1:C:504:CYS:CA	1.65	1.43
1:C:480:SER:OG	1:C:504:CYS:CB	1.77	1.31
1:C:480:SER:OG	1:C:504:CYS:C	1.70	1.29

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:GLN:HB2	5:A:1029:HOH:O	1.14	1.23
1:C:480:SER:OG	1:C:504:CYS:O	1.57	1.18

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:377:ASN:CB	1:D:572:LYS:CE[1_556]	2.02	0.18
1:C:377:ASN:CB	1:D:572:LYS:CD[1_556]	2.10	0.10
1:C:377:ASN:CB	1:D:572:LYS:NZ[1_556]	2.11	0.09
5:A:901:HOH:O	5:C:861:HOH:O[1_655]	2.16	0.04

## 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	572/612 (94%)	561 (98%)	10 (2%)	1 (0%)	47	55
1	B	567/612 (93%)	554 (98%)	11 (2%)	2 (0%)	34	37
1	C	560/612 (92%)	547 (98%)	13 (2%)	0	100	100
1	D	561/612 (92%)	551 (98%)	10 (2%)	0	100	100
All	All	2260/2448 (92%)	2213 (98%)	44 (2%)	3 (0%)	51	60

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	505	LEU
1	A	565	GLY
1	B	506	ASP

### 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	522/557 (94%)	514 (98%)	8 (2%)	65	78
1	B	523/557 (94%)	512 (98%)	11 (2%)	53	67
1	C	516/557 (93%)	503 (98%)	13 (2%)	47	60
1	D	515/557 (92%)	513 (100%)	2 (0%)	91	96
All	All	2076/2228 (93%)	2042 (98%)	34 (2%)	62	76

5 of 34 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	505	LEU
1	C	506	ASP
1	D	201	THR
1	B	277	LYS
1	B	245	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	377	ASN
1	C	464	GLN
1	D	606	HIS
1	D	23	GLN
1	B	23	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

10 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
4	SO4	B	703	-	4,4,4	0.09	0	6,6,6	0.49	0
2	AMP	A	701	-	22,25,25	0.82	1 (4%)	25,38,38	1.39	2 (8%)
2	AMP	C	701	-	22,25,25	0.88	1 (4%)	25,38,38	1.27	3 (12%)
3	IAC	C	702	-	13,14,14	2.37	6 (46%)	14,19,19	1.92	6 (42%)
3	IAC	B	702	-	13,14,14	2.45	7 (53%)	14,19,19	2.35	6 (42%)
4	SO4	B	704	-	4,4,4	0.14	0	6,6,6	0.18	0
2	AMP	D	701	-	22,25,25	0.82	1 (4%)	25,38,38	1.43	5 (20%)
2	AMP	B	701	-	22,25,25	0.88	1 (4%)	25,38,38	1.35	3 (12%)
3	IAC	D	702	-	13,14,14	2.38	7 (53%)	14,19,19	2.45	7 (50%)
3	IAC	A	702	-	13,14,14	2.41	7 (53%)	14,19,19	2.27	5 (35%)

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	701	-	-	0/6/26/26	0/3/3/3
2	AMP	C	701	-	-	3/6/26/26	0/3/3/3
3	IAC	C	702	-	-	1/4/4/4	0/2/2/2
3	IAC	B	702	-	-	0/4/4/4	0/2/2/2
2	AMP	D	701	-	-	0/6/26/26	0/3/3/3
2	AMP	B	701	-	-	0/6/26/26	0/3/3/3
3	IAC	D	702	-	-	0/4/4/4	0/2/2/2
3	IAC	A	702	-	-	0/4/4/4	0/2/2/2

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	702	IAC	C3-C2	4.31	1.46	1.36
3	D	702	IAC	C3-C2	4.29	1.46	1.36
3	A	702	IAC	C3-C2	4.28	1.46	1.36
3	C	702	IAC	C3-C2	4.16	1.46	1.36
3	A	702	IAC	C-N	3.70	1.50	1.38

The worst 5 of 37 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	702	IAC	C2-C1-C	-4.16	112.64	118.17
2	A	701	AMP	N3-C2-N1	-3.90	122.59	128.68
3	B	702	IAC	C2-C1-C	-3.82	113.10	118.17
2	D	701	AMP	N3-C2-N1	-3.77	122.79	128.68
3	B	702	IAC	C1-C-N	-3.76	99.55	107.92

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	701	AMP	C5'-O5'-P-O2P
2	C	701	AMP	C5'-O5'-P-O3P
3	C	702	IAC	C18-C17-C7-C1
2	C	701	AMP	C5'-O5'-P-O1P

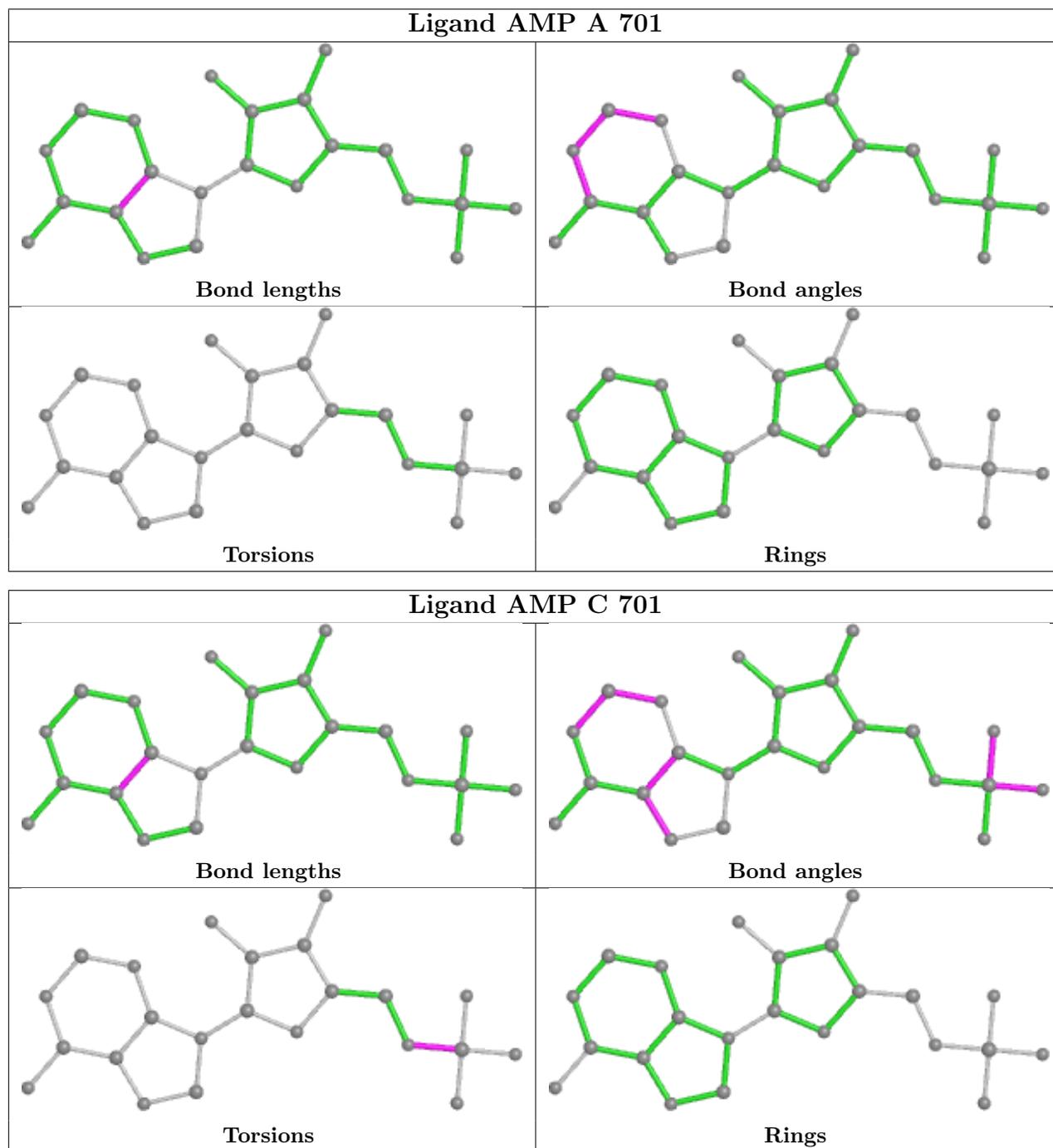
There are no ring outliers.

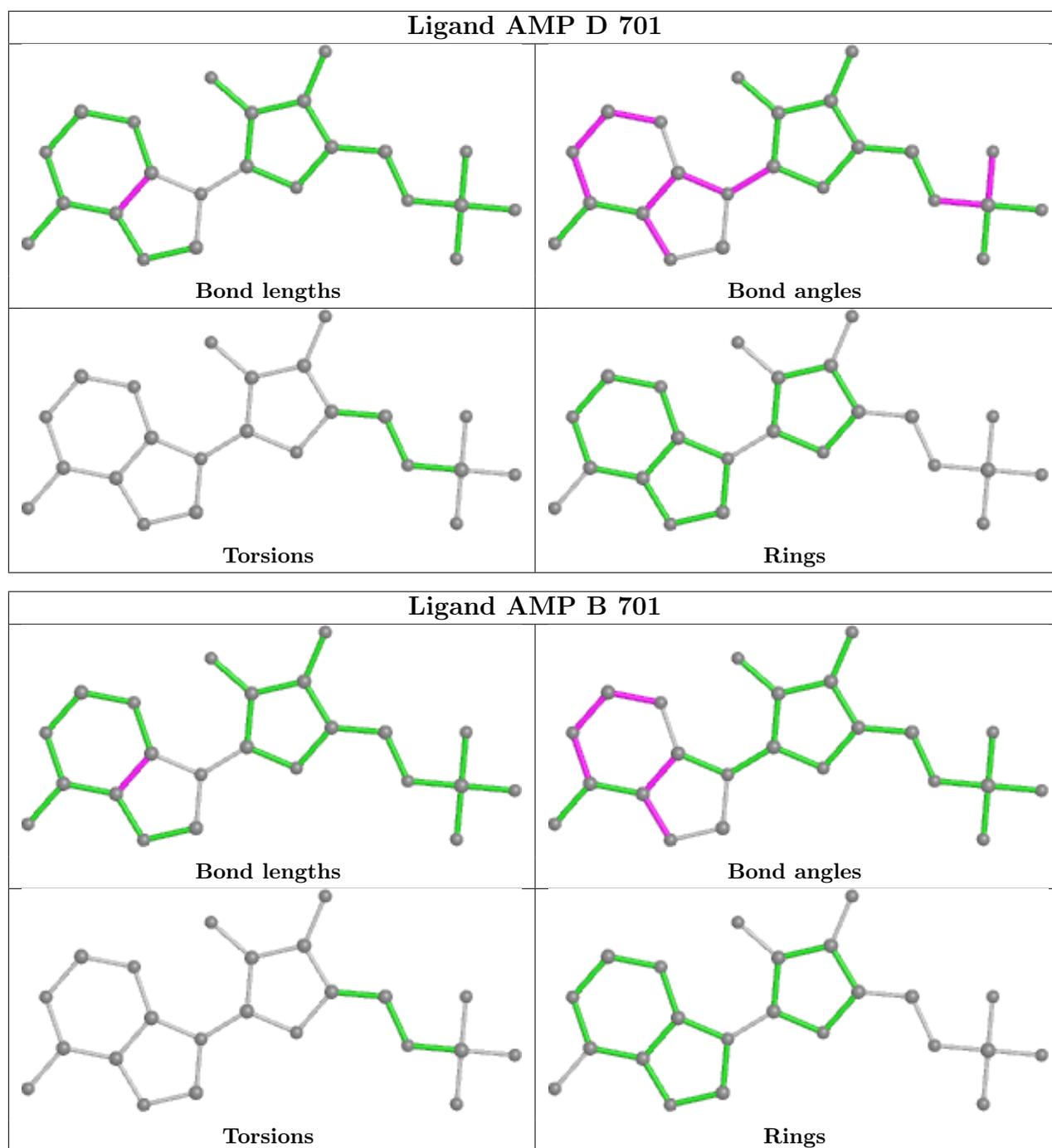
4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	702	IAC	2	0
3	B	702	IAC	3	0
3	D	702	IAC	2	0
3	A	702	IAC	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 [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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	573/612 (93%)	-0.23	11 (1%) 66 65	16, 27, 48, 70	0
1	B	573/612 (93%)	-0.24	7 (1%) 79 77	13, 27, 48, 68	0
1	C	566/612 (92%)	-0.12	21 (3%) 41 39	18, 31, 59, 79	0
1	D	566/612 (92%)	-0.29	7 (1%) 79 77	16, 27, 50, 68	0
All	All	2278/2448 (93%)	-0.22	46 (2%) 65 63	13, 28, 51, 79	0

The worst 5 of 46 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	571	TYR	6.5
1	C	480	SER	5.3
1	D	564	LEU	4.6
1	A	508	ASN	4.2
1	A	564	LEU	4.0

### 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](#)

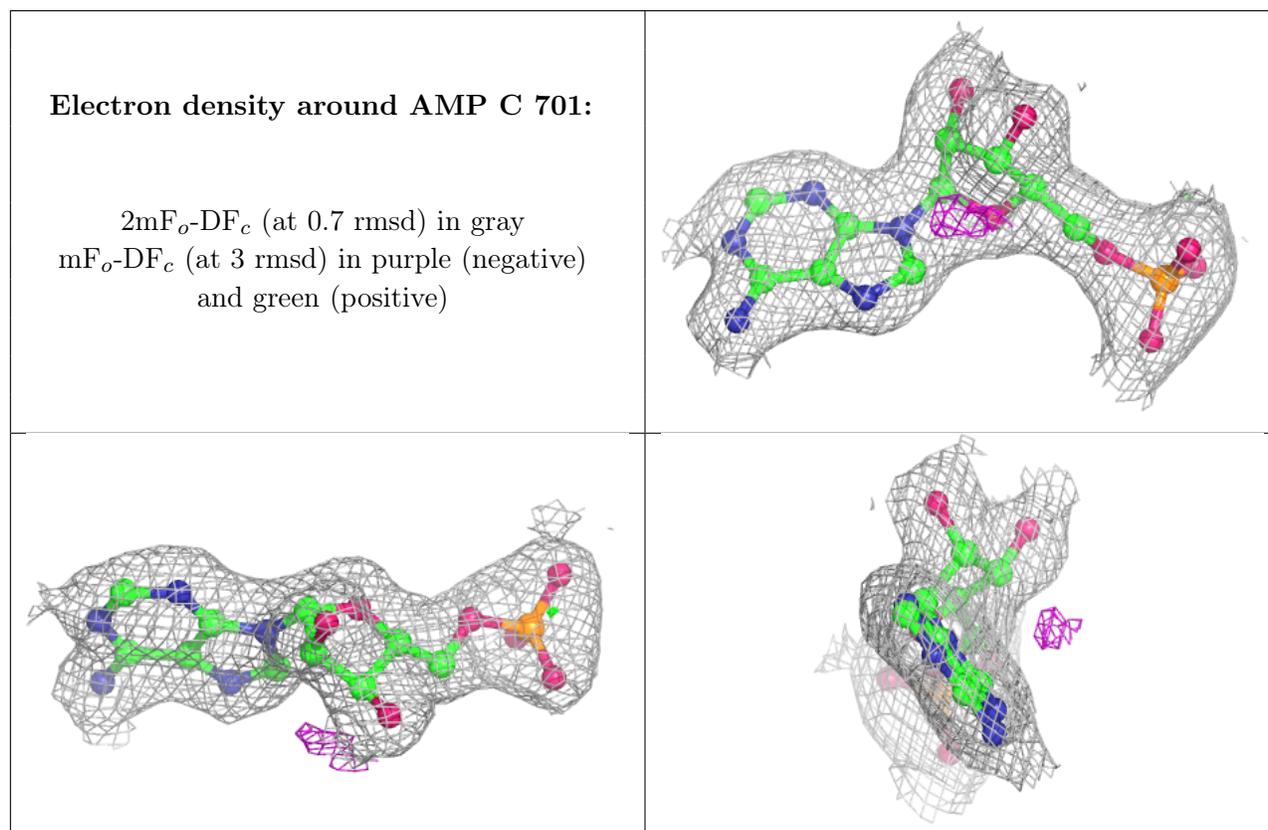
There are no monosaccharides in this entry.

### 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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

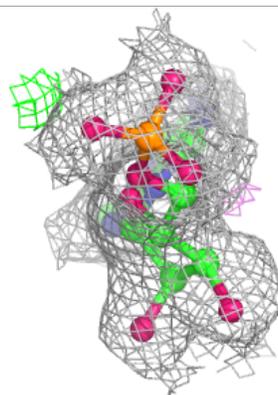
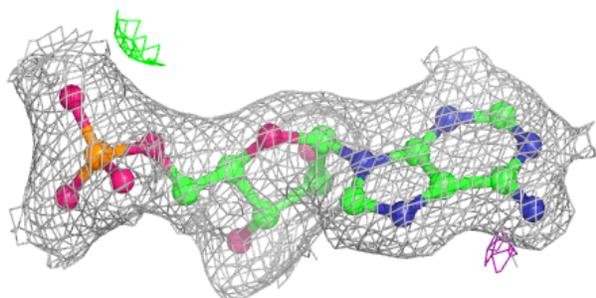
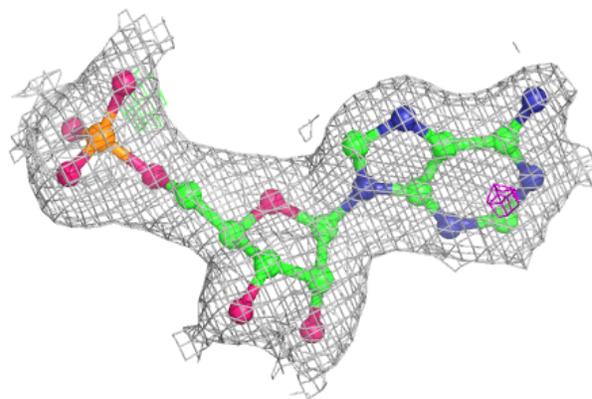
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	SO4	B	703	5/5	0.77	0.24	47,48,57,62	0
3	IAC	C	702	13/13	0.84	0.25	34,44,52,53	0
4	SO4	B	704	5/5	0.84	0.22	72,76,77,78	0
3	IAC	A	702	13/13	0.88	0.30	9,35,41,42	0
3	IAC	D	702	13/13	0.88	0.24	10,33,41,43	0
3	IAC	B	702	13/13	0.89	0.23	5,29,35,38	0
2	AMP	C	701	23/23	0.97	0.12	12,17,25,29	0
2	AMP	D	701	23/23	0.97	0.11	10,16,22,25	0
2	AMP	A	701	23/23	0.97	0.12	12,18,23,27	0
2	AMP	B	701	23/23	0.97	0.13	12,16,19,27	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.

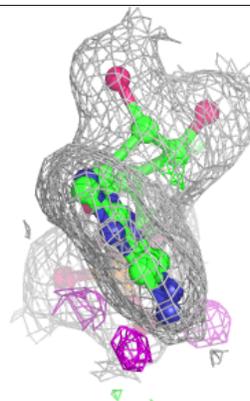
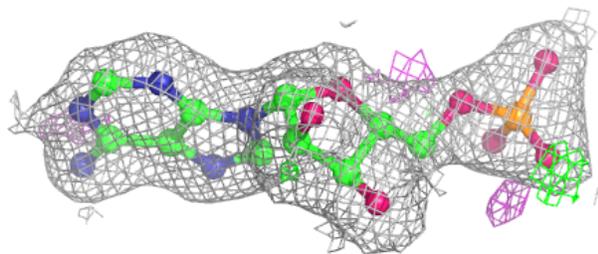
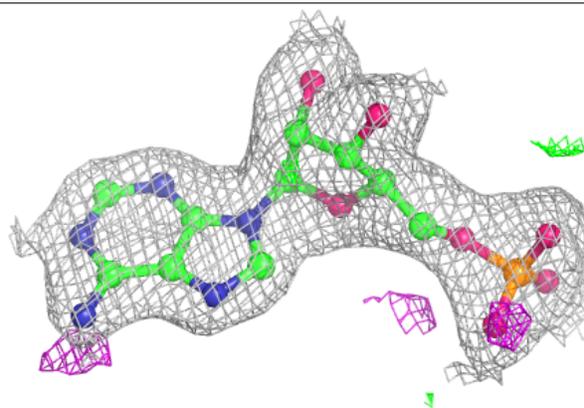


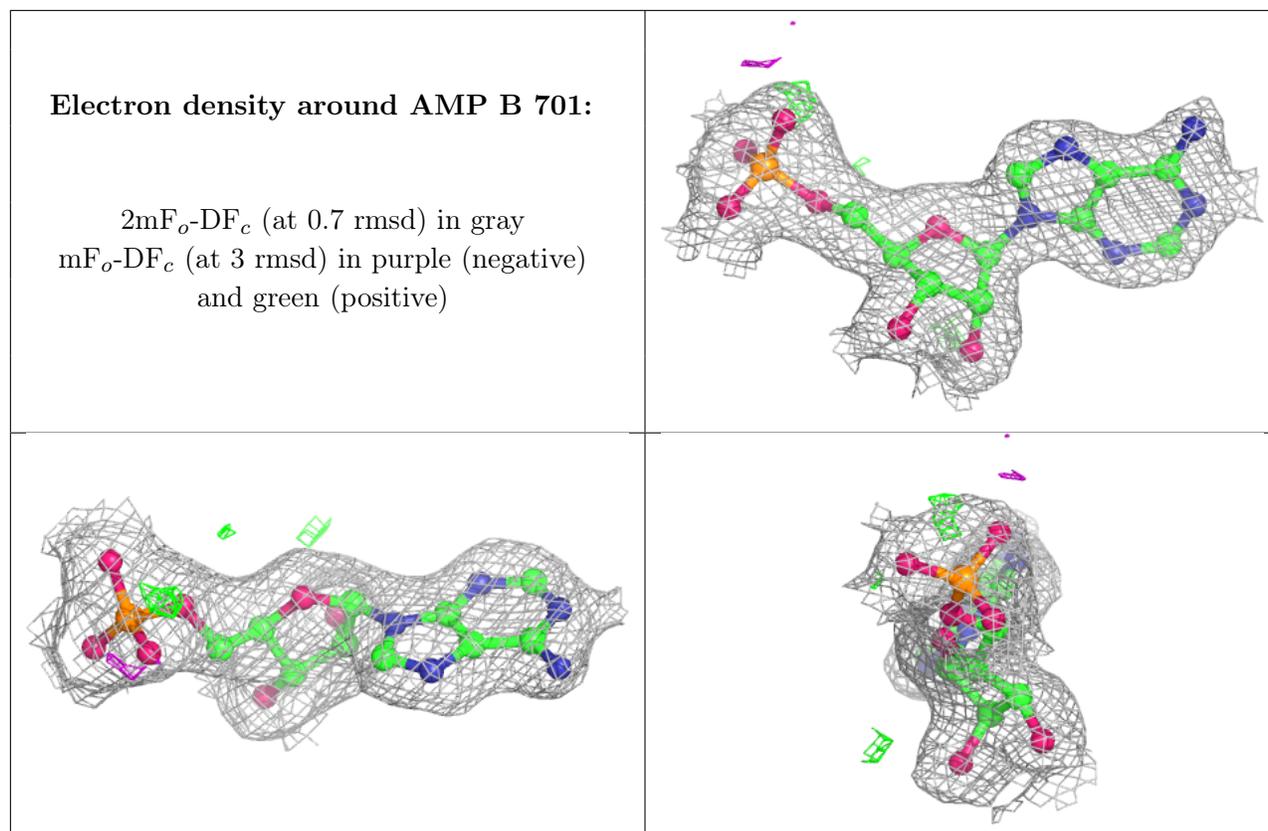
**Electron density around AMP D 701:**

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

**Electron density around AMP A 701:**

$2mF_o-DF_c$  (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.