



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 10:25 pm BST

PDB ID : 5UFU  
Title : Structure of AMPK bound to activator  
Authors : Calabrese, M.F.; Kurumbail, R.G.  
Deposited on : 2017-01-05  
Resolution : 3.45 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](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.

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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 : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

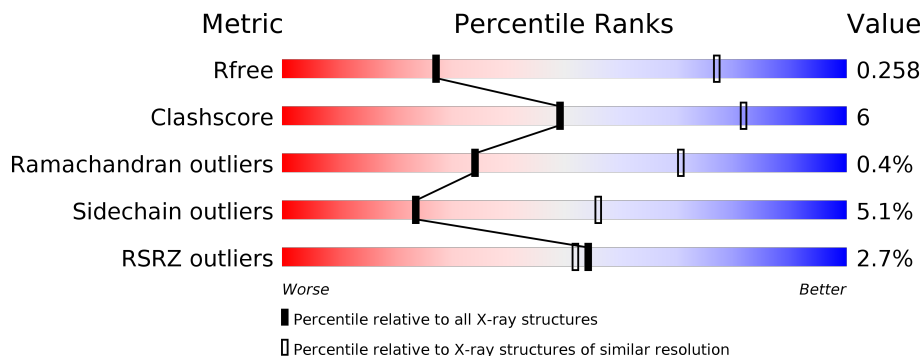
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.45 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	1291 (3.52-3.40)
Clashscore	141614	1372 (3.52-3.40)
Ramachandran outliers	138981	1337 (3.52-3.40)
Sidechain outliers	138945	1338 (3.52-3.40)
RSRZ outliers	127900	1205 (3.52-3.40)

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

Mol	Chain	Length	Quality of chain
1	A	503	 4% 60% 13% 23%
2	B	204	 65% 13% 22%
3	C	330	 4% 78% 8% 12%

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 6125 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 5'-AMP-activated protein kinase catalytic subunit alpha-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	369	2802	1793	481	510	1	17	0	0	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P54645
A	517	ALA	-	see remark 999	UNP P54645
A	518	SER	-	see remark 999	UNP P54645
A	519	GLY	-	see remark 999	UNP P54645
A	520	GLY	-	see remark 999	UNP P54645
A	521	PRO	-	see remark 999	UNP P54645
A	522	GLY	-	see remark 999	UNP P54645
A	523	GLY	-	see remark 999	UNP P54645
A	524	SER	-	see remark 999	UNP P54645

- Molecule 2 is a protein called 5'-AMP-activated protein kinase subunit beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	159	1172	762	198	210	2	0	0	0

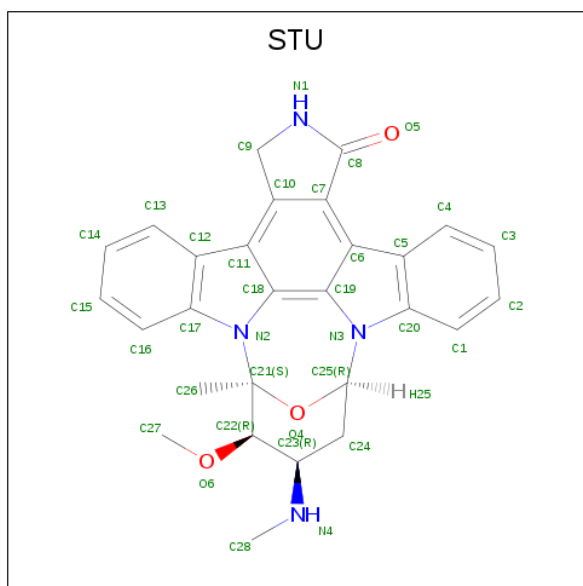
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	67	MET	-	initiating methionine	UNP P80386
B	108	ASP	SER	engineered mutation	UNP P80386

- Molecule 3 is a protein called 5'-AMP-activated protein kinase subunit gamma-1.

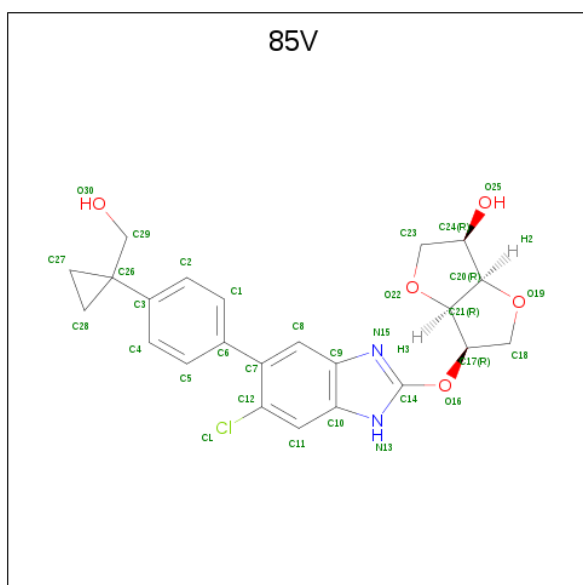
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	289	1999	1276	344	373	6	0	0	0

- Molecule 4 is STAUROSPORINE (three-letter code: STU) (formula:  $C_{28}H_{26}N_4O_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	35	28	4	3	0	0

- Molecule 5 is 1,4:3,6-dianhydro-2-O-(6-chloro-5-{4-[1-(hydroxymethyl)cyclopropyl]phenyl}-1H-benzimidazol-2-yl)-D-mannitol (three-letter code: 85V) (formula:  $C_{23}H_{23}ClN_2O_5$ ).

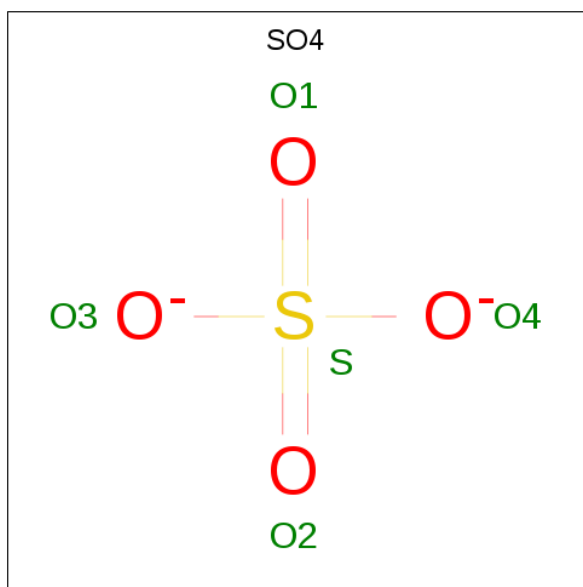


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Cl	N			O
5	A	1	31	23	1	2	5	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

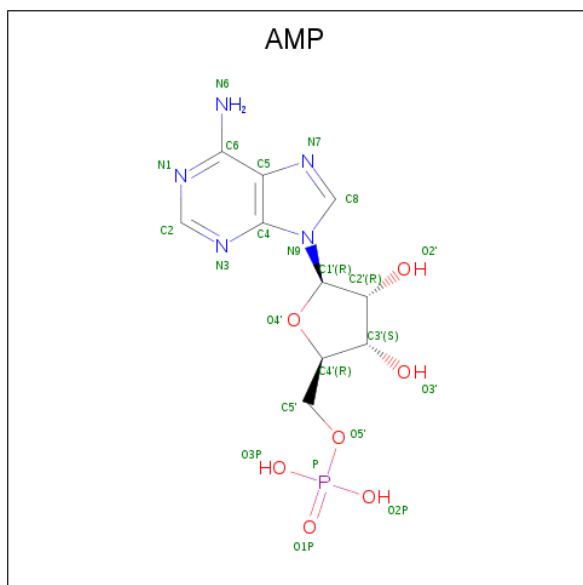
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	3	Total Cl 3 3	0	0

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



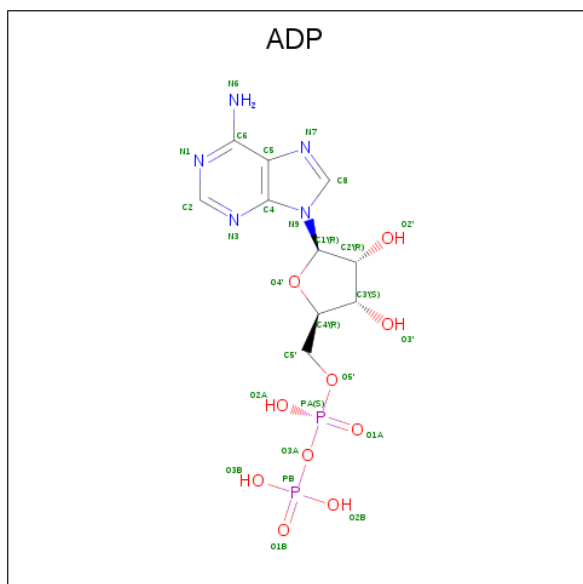
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total O S 5 4 1	0	0
7	C	1	Total O S 5 4 1	0	0

- Molecule 8 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
8	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
8	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 9 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).

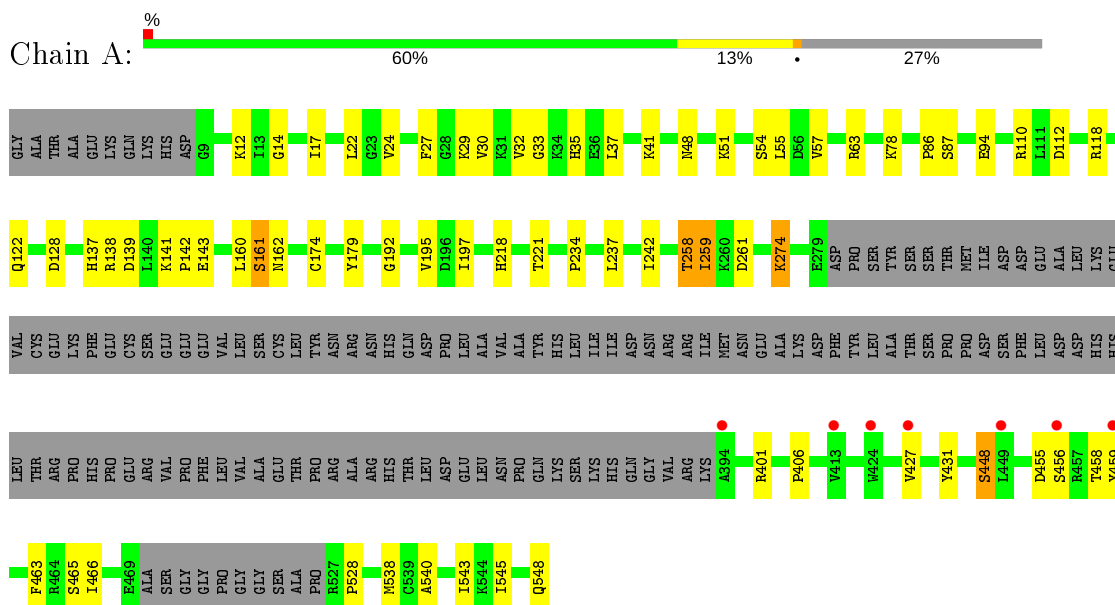


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

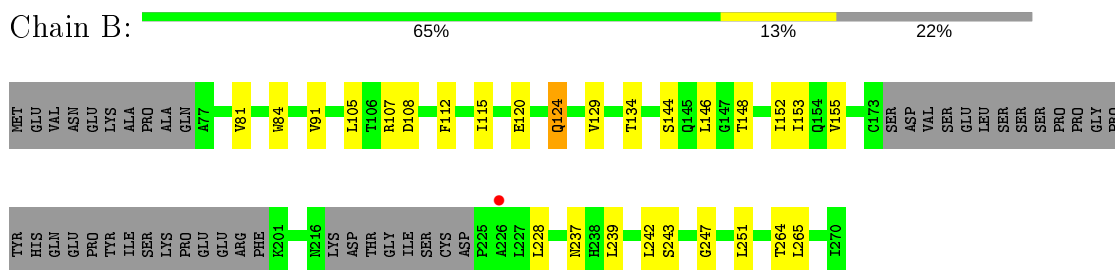
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

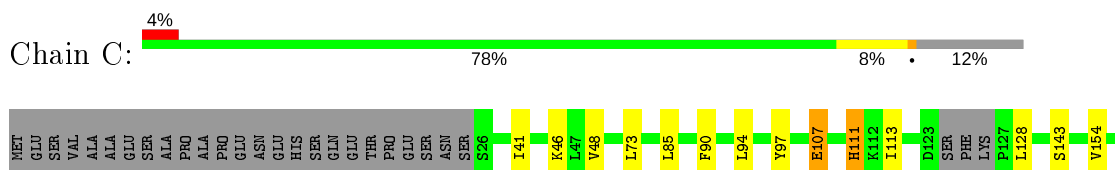
- Molecule 1: 5'-AMP-activated protein kinase catalytic subunit alpha-1

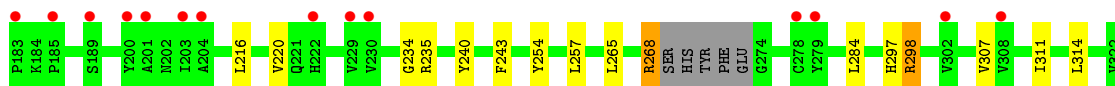


- Molecule 2: 5'-AMP-activated protein kinase subunit beta-1



- Molecule 3: 5'-AMP-activated protein kinase subunit gamma-1





LEU  
THR  
GLY  
GLY  
GLU  
LYS  
LYS  
PRO



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	124.26Å 124.26Å 401.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.43 – 3.45 47.43 – 3.45	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.43-3.45) 99.8 (47.43-3.45)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.57 (at 3.48Å)	Xtrriage
Refinement program	BUSTER 2.11.7	Depositor
R, $R_{free}$	0.193 , 0.238 0.216 , 0.258	Depositor DCC
$R_{free}$ test set	1242 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	83.1	Xtrriage
Anisotropy	0.513	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 101.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	6125	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	90.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% 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: 85V, ADP, CL, TPO, STU, 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	0.49	0/2856	0.72	1/3886 (0.0%)
2	B	0.49	0/1206	0.72	0/1663
3	C	0.47	0/2039	0.66	0/2809
All	All	0.48	0/6101	0.70	1/8358 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	14	GLY	N-CA-C	-5.08	100.39	113.10

There are no chirality outliers.

There are no planarity outliers.

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

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2802	0	2617	36	0
2	B	1172	0	1070	17	0
3	C	1999	0	1720	20	0
4	A	35	0	26	1	0
5	A	31	0	0	0	0
6	A	3	0	0	0	0
7	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	C	5	0	0	0	0
8	C	46	0	24	2	0
9	C	27	0	12	0	0
All	All	6125	0	5469	68	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.

All (68) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:107:GLU:O	3:C:111:HIS:HB2	1.92	0.70
1:A:427:VAL:HB	1:A:431:TYR:HD2	1.60	0.67
1:A:141:LYS:HG3	1:A:143:GLU:HB2	1.78	0.64
1:A:160:LEU:HD13	1:A:174:CYS:HB2	1.81	0.62
3:C:97:TYR:CD1	3:C:107:GLU:HG3	2.35	0.62
1:A:78:LYS:H	1:A:94:GLU:HG2	1.65	0.61
4:A:601:STU:H16	4:A:601:STU:H261	1.81	0.61
3:C:90:PHE:O	3:C:94:LEU:HB2	2.01	0.60
1:A:128:ASP:HB2	1:A:259:ILE:HG21	1.84	0.59
3:C:97:TYR:CG	3:C:107:GLU:HG3	2.38	0.58
1:A:274:LYS:HD3	1:A:274:LYS:H	1.69	0.57
2:B:107:ARG:HD2	2:B:112:PHE:CZ	2.40	0.57
1:A:48:ASN:HB3	1:A:51:LYS:HB3	1.87	0.56
1:A:218:HIS:HD2	1:A:221:THR:HG23	1.70	0.56
3:C:243:PHE:HB3	8:C:402:AMP:H5'1	1.86	0.56
1:A:455:ASP:HB3	1:A:458:THR:HG22	1.87	0.55
2:B:239:LEU:HD11	2:B:251:LEU:HD22	1.88	0.55
3:C:235:ARG:HA	3:C:307:VAL:HA	1.88	0.55
1:A:218:HIS:CD2	1:A:221:THR:HG23	2.41	0.54
1:A:24:VAL:HG12	1:A:29:LYS:HA	1.90	0.53
1:A:427:VAL:HB	1:A:431:TYR:CD2	2.45	0.52
2:B:107:ARG:HB3	2:B:112:PHE:CD2	2.44	0.52
1:A:401:ARG:O	1:A:548:GLN:HB3	2.10	0.52
1:A:465:SER:HB3	2:B:237:ASN:HB3	1.93	0.51
2:B:144:SER:HB3	2:B:148:THR:H	1.75	0.51
2:B:124:GLN:HG3	2:B:152:ILE:HD12	1.94	0.50
3:C:297:HIS:HA	3:C:314:LEU:HD22	1.94	0.50
3:C:94:LEU:HD11	3:C:254:TYR:HB2	1.93	0.49
2:B:265:LEU:HD12	3:C:48:VAL:HG22	1.93	0.49
1:A:137:HIS:HA	1:A:161:SER:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:LEU:HD12	1:A:30:VAL:HG12	1.96	0.48
2:B:84:TRP:CE2	2:B:129:VAL:HG21	2.49	0.47
3:C:94:LEU:HG	3:C:257:LEU:HD11	1.97	0.47
2:B:91:VAL:HG22	2:B:129:VAL:HG13	1.96	0.47
1:A:448:SER:HB3	1:A:466:ILE:HD11	1.97	0.47
1:A:137:HIS:HE1	1:A:139:ASP:O	1.97	0.46
1:A:540:ALA:HB2	2:B:251:LEU:HD11	1.97	0.46
3:C:265:LEU:HD22	3:C:268:ARG:NH1	2.30	0.46
3:C:265:LEU:HD22	3:C:268:ARG:HH12	1.80	0.46
1:A:234:PRO:HD2	1:A:237:LEU:HD22	1.97	0.46
2:B:228:LEU:HD11	3:C:46:LYS:HB2	1.98	0.45
1:A:35:HIS:HE1	1:A:37:LEU:HD12	1.82	0.45
1:A:258:THR:H	1:A:261:ASP:HB2	1.81	0.44
1:A:463:PHE:HB2	2:B:239:LEU:HB3	1.98	0.44
1:A:33:GLY:O	1:A:41:LYS:HA	2.17	0.44
1:A:12:LYS:HG2	1:A:17:ILE:HG22	1.98	0.44
3:C:298:ARG:HH11	8:C:402:AMP:H8	1.66	0.44
1:A:27:PHE:CE1	1:A:57:VAL:HG11	2.53	0.43
1:A:86:PRO:HG3	2:B:155:VAL:HG23	2.00	0.43
3:C:216:LEU:O	3:C:220:VAL:HG23	2.18	0.43
2:B:91:VAL:HG12	2:B:105:LEU:HD12	2.00	0.42
3:C:73:LEU:HD21	3:C:85:LEU:HB2	2.01	0.42
1:A:118:ARG:O	1:A:122:GLN:HG3	2.19	0.42
3:C:97:TYR:HB3	3:C:107:GLU:HG3	2.01	0.42
1:A:543:ILE:C	1:A:545:ILE:H	2.22	0.42
3:C:73:LEU:HD22	3:C:113:ILE:HG21	2.01	0.42
1:A:138:ARG:NH2	1:A:162:ASN:HB2	2.35	0.42
1:A:237:LEU:HB3	1:A:242:ILE:HD11	2.02	0.41
3:C:240:TYR:CD1	3:C:265:LEU:HD21	2.54	0.41
1:A:406:PRO:HG3	1:A:459:TYR:CZ	2.55	0.41
3:C:41:ILE:HD11	3:C:154:VAL:HG11	2.01	0.41
1:A:455:ASP:CG	1:A:456:SER:H	2.23	0.41
1:A:197:ILE:HD11	1:A:259:ILE:HG13	2.03	0.41
1:A:142:PRO:HD2	1:A:179:TYR:CE1	2.55	0.41
1:A:192:GLY:O	1:A:195:VAL:HG22	2.21	0.41
2:B:120:GLU:HA	2:B:155:VAL:HG13	2.02	0.40
2:B:107:ARG:HD2	2:B:112:PHE:CE2	2.56	0.40
2:B:81:VAL:HG22	2:B:115:ILE:HG12	2.03	0.40

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	362/503 (72%)	333 (92%)	28 (8%)	1 (0%)	41	75
2	B	153/204 (75%)	141 (92%)	11 (7%)	1 (1%)	22	60
3	C	283/330 (86%)	270 (95%)	12 (4%)	1 (0%)	34	70
All	All	798/1037 (77%)	744 (93%)	51 (6%)	3 (0%)	34	70

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	528	PRO
2	B	247	GLY
3	C	234	GLY

### 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	281/448 (63%)	268 (95%)	13 (5%)	27	59
2	B	115/185 (62%)	107 (93%)	8 (7%)	15	46
3	C	174/299 (58%)	166 (95%)	8 (5%)	27	59
All	All	570/932 (61%)	541 (95%)	29 (5%)	24	56

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	VAL
1	A	54	SER
1	A	55	LEU
1	A	63	ARG
1	A	87	SER
1	A	110	ARG
1	A	112	ASP
1	A	161	SER
1	A	258	THR
1	A	259	ILE
1	A	274	LYS
1	A	448	SER
1	A	538	MET
2	B	108	ASP
2	B	124	GLN
2	B	134	THR
2	B	146	LEU
2	B	153	ILE
2	B	242	LEU
2	B	243	SER
2	B	264	THR
3	C	107	GLU
3	C	111	HIS
3	C	128	LEU
3	C	143	SER
3	C	268	ARG
3	C	284	LEU
3	C	298	ARG
3	C	311	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	218	HIS
1	A	548	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](#)

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	TPO	A	172	1	8,10,11	1.49	1 (12%)	10,14,16	1.69	3 (30%)

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	TPO	A	172	1	-	3/9/11/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	172	TPO	P-OG1	-3.46	1.52	1.59

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	172	TPO	P-OG1-CB	-2.60	115.36	123.21
1	A	172	TPO	O2P-P-OG1	2.11	115.46	105.99
1	A	172	TPO	O3P-P-O1P	2.05	118.69	110.68

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	172	TPO	CB-OG1-P-O1P
1	A	172	TPO	CB-OG1-P-O3P
1	A	172	TPO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 3 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
9	ADP	C	403	-	24,29,29	0.66	0	29,45,45	0.96	1 (3%)
4	STU	A	601	-	30,42,42	1.13	2 (6%)	31,68,68	0.83	2 (6%)
7	SO4	A	606	-	4,4,4	0.16	0	6,6,6	0.09	0
7	SO4	C	404	-	4,4,4	0.14	0	6,6,6	0.22	0
8	AMP	C	402	-	22,25,25	0.66	0	25,38,38	1.26	3 (12%)
5	85V	A	602	-	35,36,36	0.75	1 (2%)	40,55,55	0.59	0
8	AMP	C	401	-	22,25,25	0.62	0	25,38,38	1.25	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
4	STU	A	601	-	-	1/4/42/42	-
9	ADP	C	403	-	-	2/12/32/32	0/3/3/3
5	85V	A	602	-	-	0/15/43/43	0/6/6/6
8	AMP	C	401	-	-	0/6/26/26	0/3/3/3
8	AMP	C	402	-	-	2/6/26/26	0/3/3/3

All (3) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	601	STU	C8-N1	4.60	1.38	1.35
5	A	602	85V	C14-N13	-2.60	1.31	1.34
4	A	601	STU	C9-C10	-2.34	1.48	1.50

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	401	AMP	P-O5'-C5'	4.84	131.62	118.30
8	C	402	AMP	P-O5'-C5'	3.94	129.15	118.30
8	C	402	AMP	O2P-P-O5'	2.81	114.22	106.73
4	A	601	STU	C16-C17-N2	-2.34	129.46	132.29
8	C	402	AMP	C5-C6-N6	2.27	123.80	120.35
8	C	401	AMP	C5-C6-N6	2.27	123.80	120.35
9	C	403	ADP	C5-C6-N6	2.26	123.79	120.35
4	A	601	STU	C1-C20-N3	-2.16	129.66	132.25

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	C	402	AMP	C3'-C4'-C5'-O5'
9	C	403	ADP	C5'-O5'-PA-O3A
9	C	403	ADP	C5'-O5'-PA-O1A
4	A	601	STU	C24-C23-N4-C28
8	C	402	AMP	O4'-C4'-C5'-O5'

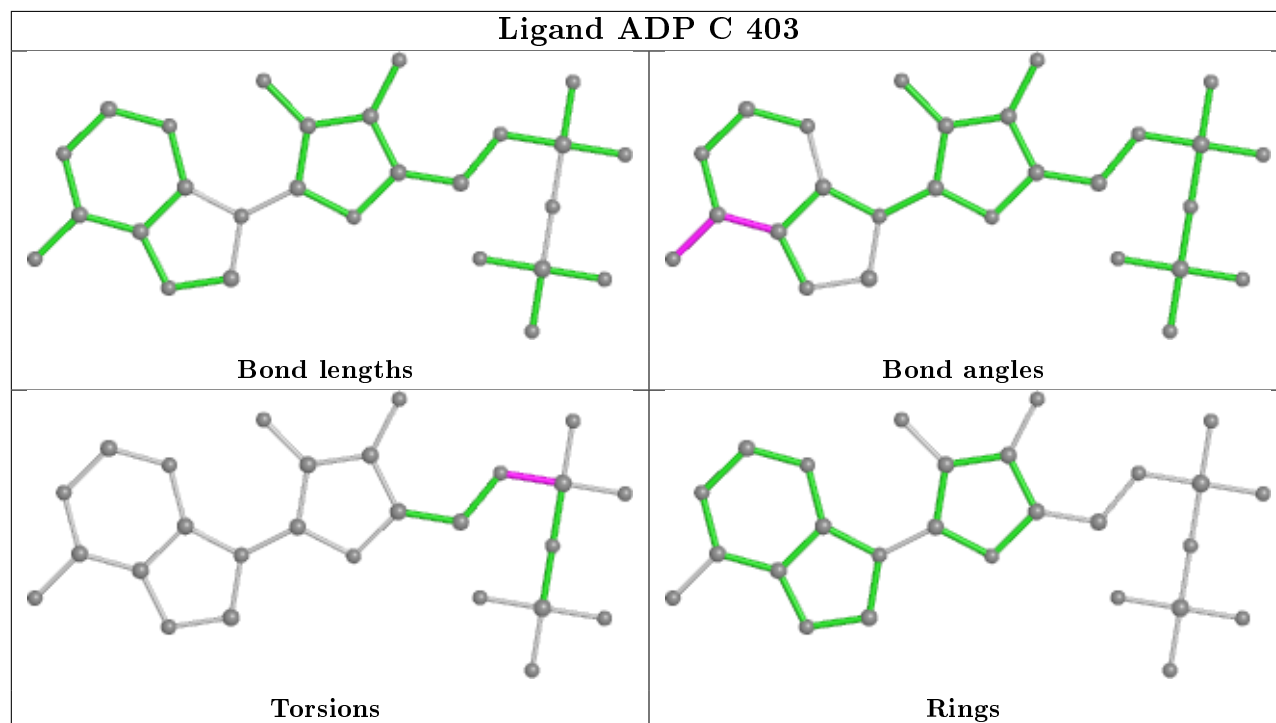
There are no ring outliers.

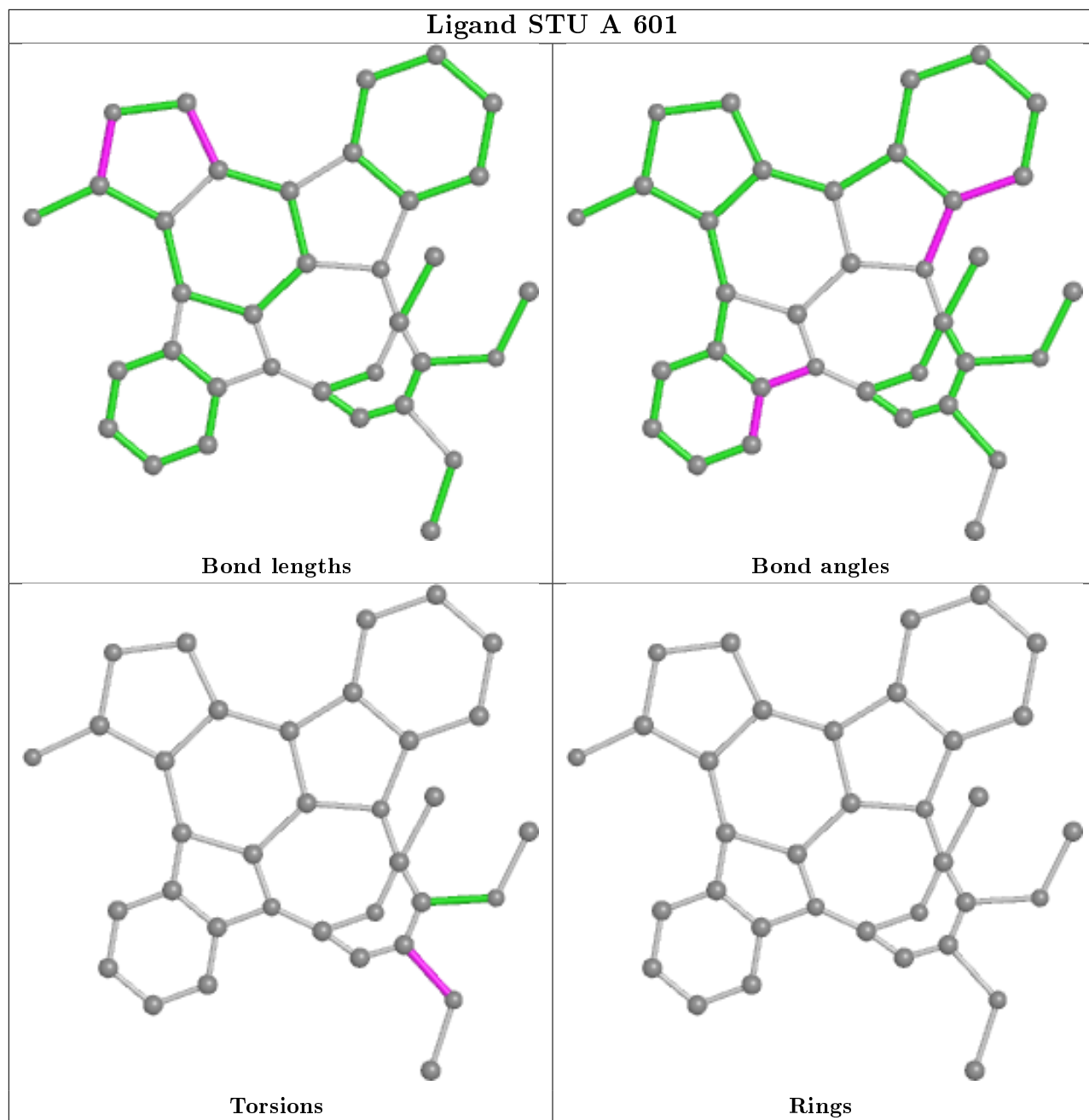
2 monomers are involved in 3 short contacts:

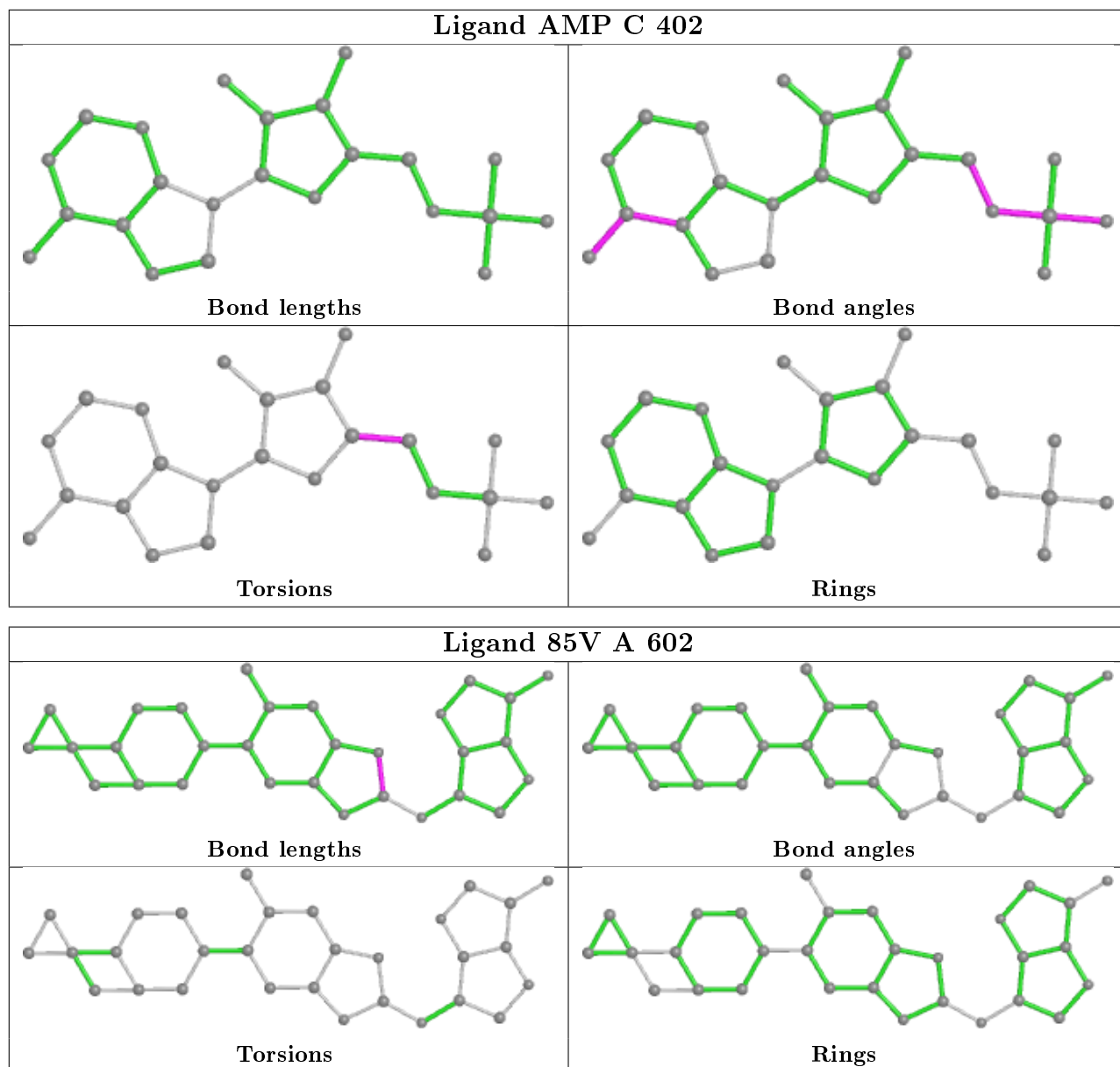
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	601	STU	1	0
8	C	402	AMP	2	0

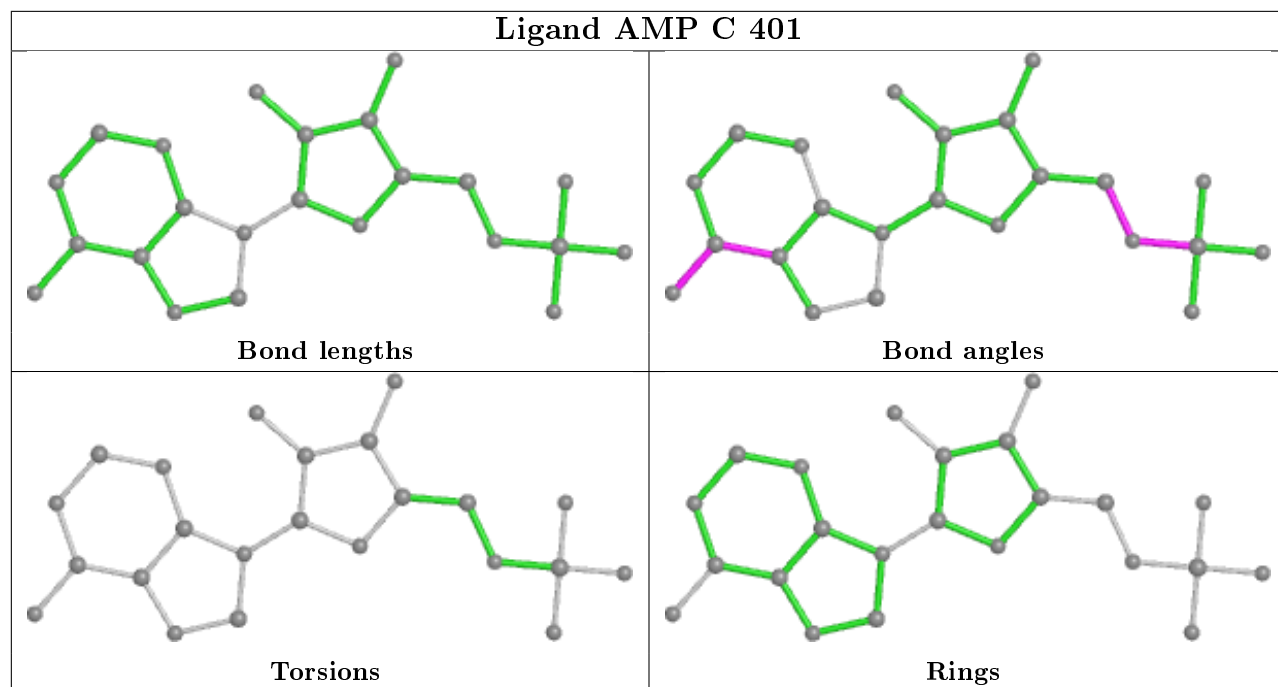
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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









## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	368/503 (73%)	0.06	7 (1%) 66 64	40, 69, 131, 152	0
2	B	159/204 (77%)	0.06	1 (0%) 89 87	49, 81, 112, 136	0
3	C	289/330 (87%)	0.29	14 (4%) 30 30	69, 110, 171, 180	0
All	All	816/1037 (78%)	0.14	22 (2%) 54 52	40, 87, 159, 180	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	230	VAL	4.3
1	A	394	ALA	3.9
1	A	459	TYR	3.8
3	C	201	ALA	3.4
3	C	308	VAL	3.4
3	C	278	CYS	3.4
3	C	302	VAL	3.0
3	C	183	PRO	2.9
3	C	229	VAL	2.9
1	A	424	TRP	2.9
3	C	189	SER	2.8
3	C	279	TYR	2.7
3	C	204	ALA	2.4
1	A	456	SER	2.3
1	A	449	LEU	2.3
1	A	427	VAL	2.3
3	C	200	TYR	2.2
3	C	203	ILE	2.2
1	A	413	VAL	2.2
3	C	185	PRO	2.2
3	C	222	HIS	2.1
2	B	226	ALA	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [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(Å <sup>2</sup> )	Q<0.9
1	TPO	A	172	11/12	0.96	0.18	70,74,79,80	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

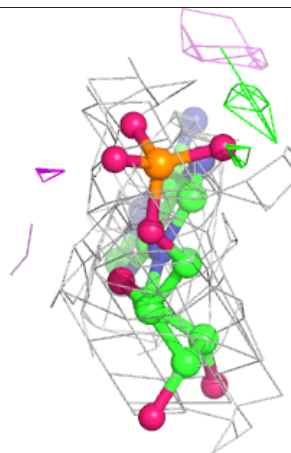
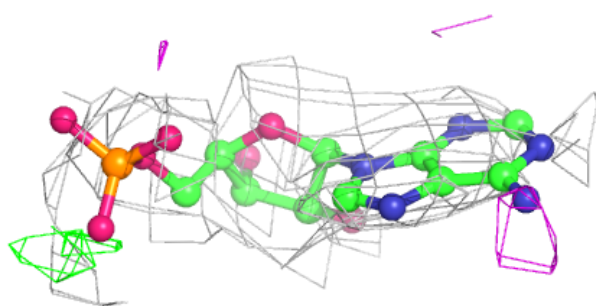
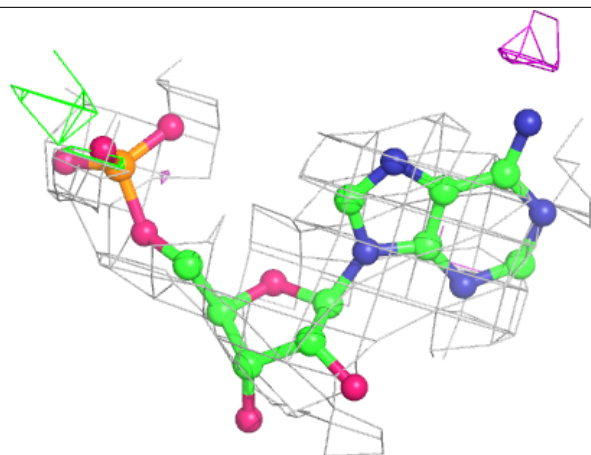
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
8	AMP	C	402	23/23	0.77	0.33	181,193,209,212	0
6	CL	A	604	1/1	0.80	0.40	75,75,75,75	0
7	SO4	A	606	5/5	0.82	0.19	174,174,174,174	0
7	SO4	C	404	5/5	0.83	0.20	141,141,141,142	0
8	AMP	C	401	23/23	0.88	0.27	128,143,148,153	0
9	ADP	C	403	27/27	0.89	0.25	157,160,179,181	0
6	CL	A	603	1/1	0.89	0.28	51,51,51,51	0
5	85V	A	602	31/31	0.97	0.27	46,52,56,58	0
6	CL	A	605	1/1	0.98	0.23	53,53,53,53	0
4	STU	A	601	35/35	0.98	0.26	48,53,60,64	0

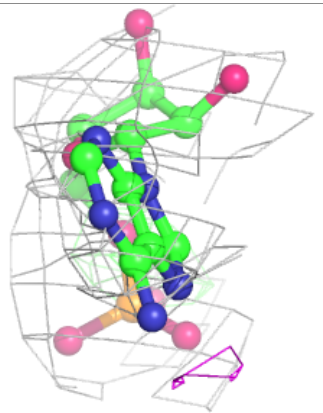
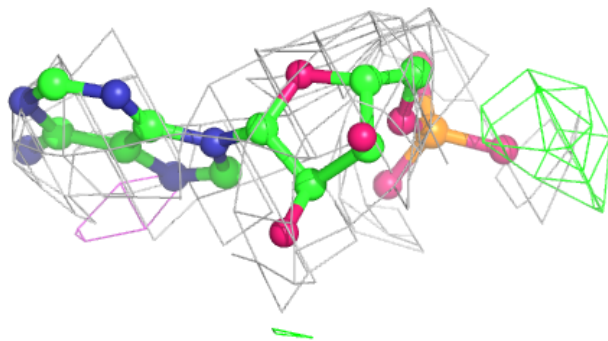
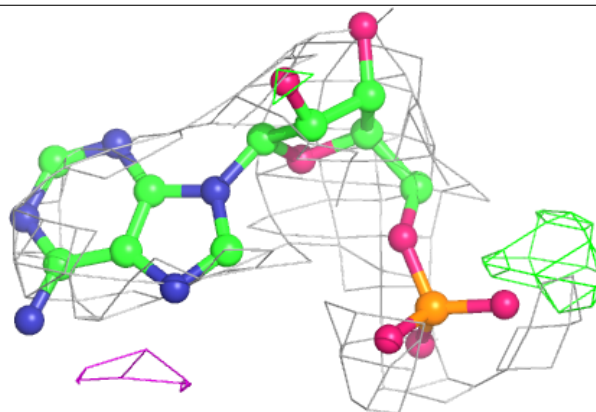
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around AMP C 402:**

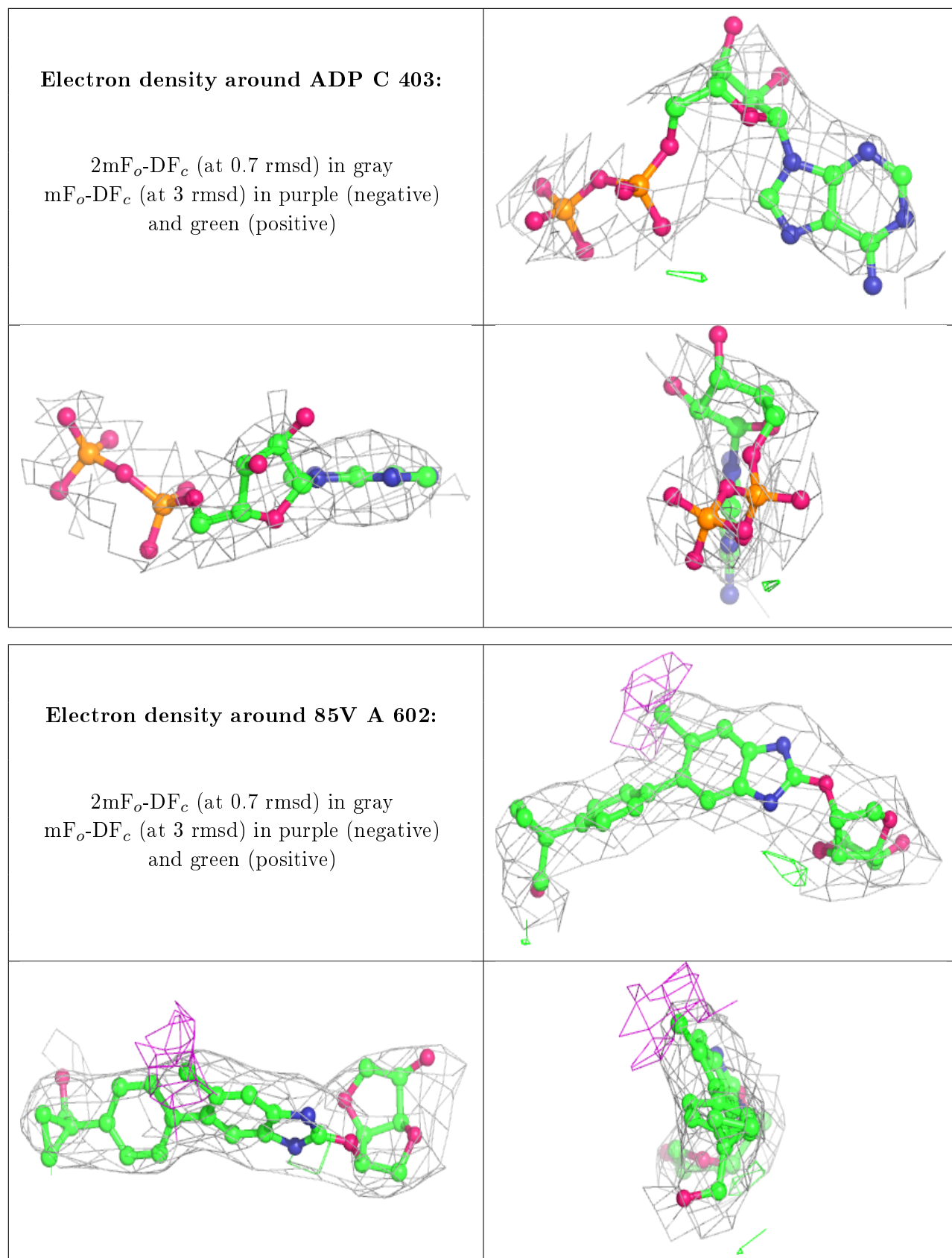
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

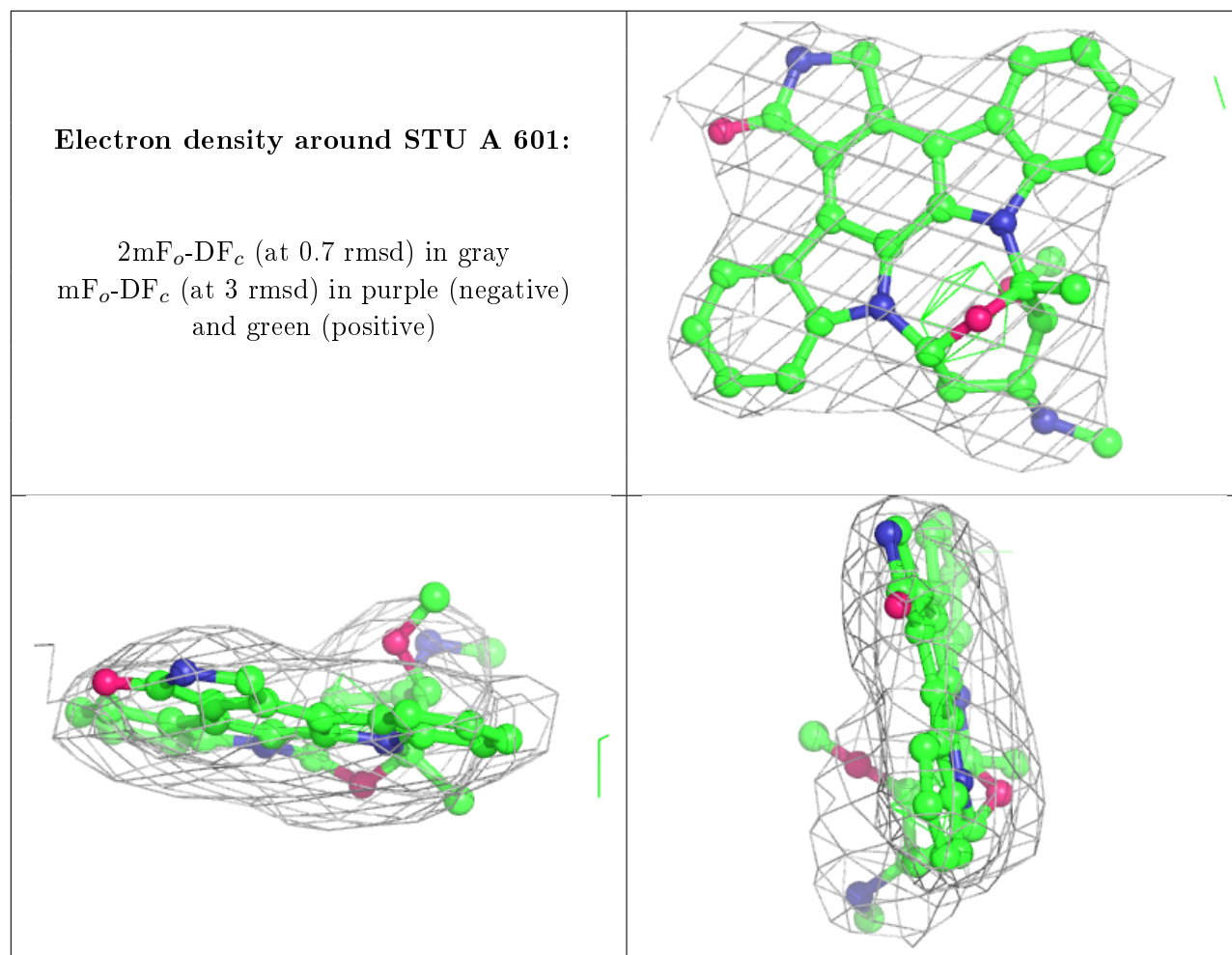
**Electron density around AMP C 401:**

$2mF_o-DF_c$  (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.