



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 10:37 am BST

PDB ID : 4CFH  
Title : Structure of an active form of mammalian AMPK  
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Deposited on : 2013-11-18  
Resolution : 3.24 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

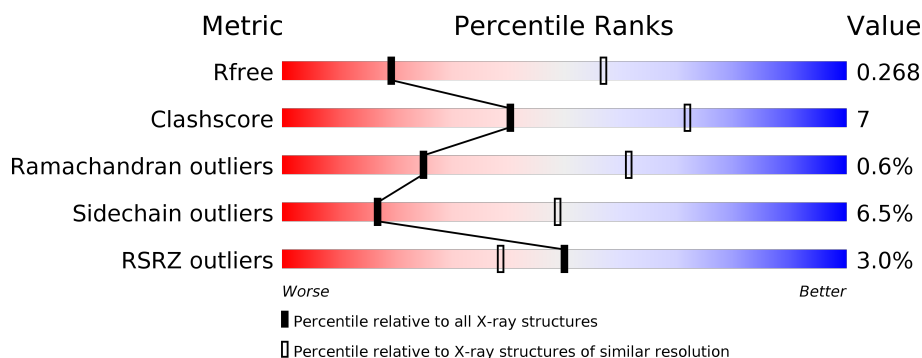
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.24 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	1619 (3.28-3.20)
Clashscore	141614	1755 (3.28-3.20)
Ramachandran outliers	138981	1728 (3.28-3.20)
Sidechain outliers	138945	1727 (3.28-3.20)
RSRZ outliers	127900	1567 (3.28-3.20)

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 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	493	<div> <div>3%</div> <div> <div></div> <div>62%</div> <div>17%</div> <div>•</div> <div>18%</div> </div> </div>
2	B	87	<div> <div>61%</div> <div>17%</div> <div>•</div> <div>21%</div> </div>
3	C	27	<div> <div>70%</div> <div>7%</div> <div>22%</div> </div>
4	E	330	<div> <div>2%</div> <div> <div></div> <div>67%</div> <div>22%</div> <div>•</div> <div>9%</div> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6429 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 SUB-UNIT ALPHA-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	P	S	0	0	0
			3217	2059	565	576	1	16			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	expression tag	UNP P54645
A	-17	SER	-	expression tag	UNP P54645
A	-16	HIS	-	expression tag	UNP P54645
A	-15	HIS	-	expression tag	UNP P54645
A	-14	HIS	-	expression tag	UNP P54645
A	-13	HIS	-	expression tag	UNP P54645
A	-12	HIS	-	expression tag	UNP P54645
A	-11	HIS	-	expression tag	UNP P54645
A	-10	SER	-	expression tag	UNP P54645
A	-9	SER	-	expression tag	UNP P54645
A	-8	GLY	-	expression tag	UNP P54645
A	-7	LEU	-	expression tag	UNP P54645
A	-6	GLU	-	expression tag	UNP P54645
A	-5	VAL	-	expression tag	UNP P54645
A	-4	LEU	-	expression tag	UNP P54645
A	-3	PHE	-	expression tag	UNP P54645
A	-2	GLN	-	expression tag	UNP P54645
A	-1	GLY	-	expression tag	UNP P54645
A	0	PRO	-	expression tag	UNP P54645
A	1	MET	-	expression tag	UNP P54645
A	471	LEU	-	SEE REMARK 999	UNP P54645
A	472	GLU	-	SEE REMARK 999	UNP P54645
A	473	VAL	-	SEE REMARK 999	UNP P54645
A	474	LEU	-	SEE REMARK 999	UNP P54645

- Molecule 2 is a protein called 5'-AMP-ACTIVATED PROTEIN KINASE SUBUNIT BETA-

2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	69	Total	C	N	O	S	0	0	0
			550	360	92	95	3			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	186	MET	-	expression tag	UNP O43741

- Molecule 3 is a protein called 5'-AMP-ACTIVATED PROTEIN KINASE CATALYTIC SUB-UNIT ALPHA-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	21	Total	C	N	O	S	0	0	0
			164	107	26	29	2			

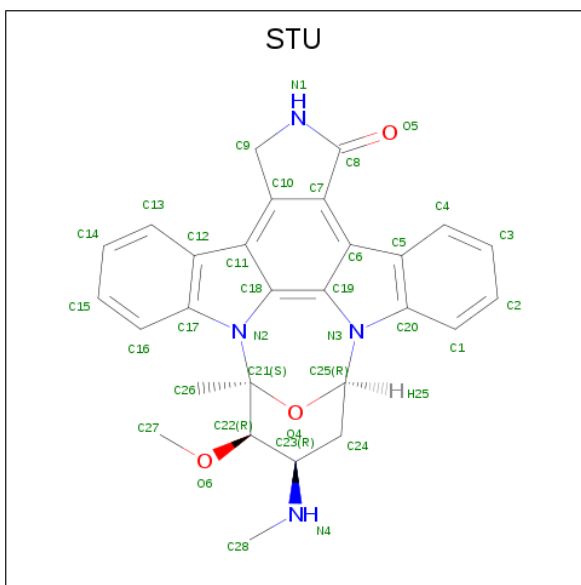
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	522	PHE	-	SEE REMARK 999	UNP P54645
C	523	GLN	-	SEE REMARK 999	UNP P54645

- Molecule 4 is a protein called 5'-AMP-ACTIVATED PROTEIN KINASE SUBUNIT GAMMA-1.

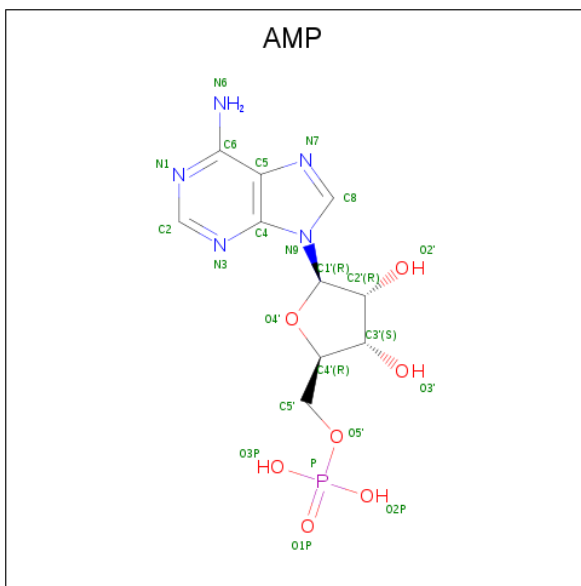
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	301	Total	C	N	O	S	0	0	0
			2417	1568	403	439	7			

- Molecule 5 is STAUROSPORINE (three-letter code: STU) (formula: C<sub>28</sub>H<sub>26</sub>N<sub>4</sub>O<sub>3</sub>).



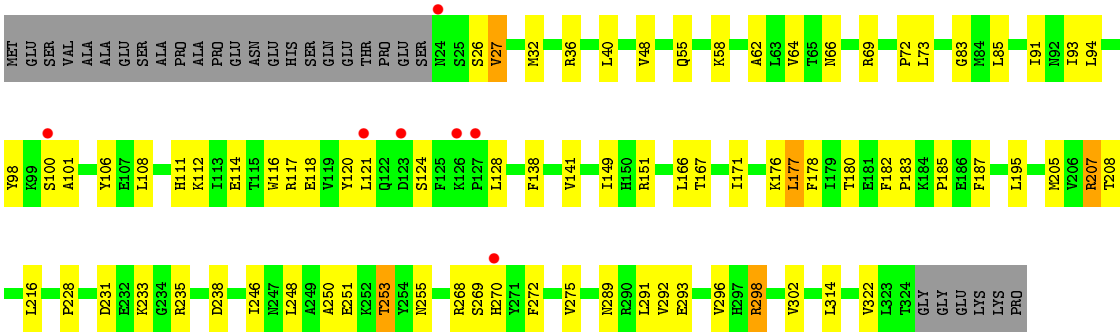
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			35	28	4	3		

- Molecule 6 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula:  $\text{C}_{10}\text{H}_{14}\text{N}_5\text{O}_7\text{P}$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	E	1	Total 23	C 10	N 5	O 7	P 1	0	0
6	E	1	Total 23	C 10	N 5	O 7	P 1	0	0





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	133.92Å 133.92Å 141.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.53 – 3.24 29.53 – 3.24	Depositor EDS
% Data completeness (in resolution range)	93.0 (29.53-3.24) 86.9 (29.53-3.24)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.46 (at 3.24Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.233 , 0.268 0.233 , 0.268	Depositor DCC
$R_{free}$ test set	989 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	108.5	Xtriage
Anisotropy	0.239	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 56.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6429	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.38% 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: TPO, STU, 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	0.22	0/3282	0.42	0/4440
2	B	0.22	0/565	0.42	0/772
3	C	0.22	0/167	0.36	0/223
4	E	0.23	0/2468	0.42	0/3353
All	All	0.23	0/6482	0.42	0/8788

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3217	0	3182	53	0
2	B	550	0	585	10	0
3	C	164	0	165	0	0
4	E	2417	0	2477	36	0
5	A	35	0	26	4	0
6	E	46	0	24	2	0
All	All	6429	0	6459	93	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 7.

All (93) 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:A:23:GLY:HA3	1:A:30:VAL:HB	1.70	0.72
1:A:264:GLU:HA	1:A:269:LYS:HE3	1.75	0.68
1:A:363:ARG:HB3	2:B:227:PRO:HG3	1.76	0.67
1:A:142:PRO:HD3	1:A:203:ILE:HG12	1.76	0.66
1:A:417:ILE:HG23	1:A:422:TYR:HB2	1.77	0.65
4:E:93:ILE:HG21	4:E:108:LEU:HD13	1.77	0.65
2:B:226:ASP:HB3	2:B:229:LEU:HG	1.76	0.65
4:E:73:LEU:HD21	4:E:85:LEU:HB2	1.79	0.64
1:A:139:ASP:HB2	1:A:160:LEU:HD12	1.80	0.64
1:A:348:ASP:N	1:A:348:ASP:OD1	2.29	0.63
1:A:143:GLU:HB3	5:A:1550:STU:H281	1.79	0.63
1:A:179:TYR:HA	1:A:202:VAL:HG21	1.80	0.62
5:A:1550:STU:H261	5:A:1550:STU:H16	1.81	0.62
1:A:45:LYS:HB3	1:A:91:MET:HB2	1.81	0.61
1:A:185:ILE:HD13	1:A:223:PHE:HB3	1.83	0.60
4:E:69:ARG:NH2	6:E:1325:AMP:O2P	2.33	0.58
1:A:139:ASP:O	1:A:144:ASN:ND2	2.34	0.58
1:A:463:PHE:HB2	2:B:241:LEU:HB3	1.86	0.57
1:A:233:THR:HG21	1:A:242:ILE:HG23	1.86	0.57
4:E:40:LEU:HD12	4:E:166:LEU:HD11	1.86	0.57
1:A:358:ARG:HD2	1:A:359:PRO:HD2	1.87	0.57
1:A:138:ARG:NH1	1:A:190:TYR:OH	2.37	0.56
1:A:277:PHE:HB3	1:A:278:PRO:HD2	1.87	0.56
1:A:363:ARG:NH1	2:B:222:ASN:O	2.39	0.56
1:A:436:ARG:NH1	1:A:445:SER:OG	2.40	0.55
4:E:289:ASN:O	4:E:293:GLU:HG2	2.07	0.54
4:E:32:MET:O	4:E:138:PHE:N	2.42	0.53
1:A:35:HIS:CE1	1:A:37:LEU:HB2	2.43	0.53
1:A:105:ILE:O	1:A:107:LYS:N	2.42	0.53
1:A:360:HIS:HB3	1:A:363:ARG:HG2	1.90	0.53
1:A:117:ARG:NH1	1:A:267:TRP:O	2.37	0.53
1:A:166:ASP:N	1:A:166:ASP:OD1	2.41	0.52
1:A:166:ASP:OD2	2:B:258:ARG:NH1	2.42	0.52
1:A:105:ILE:HG22	1:A:106:CYS:H	1.74	0.52
1:A:428:ASN:HB2	1:A:431:TYR:HB3	1.90	0.52
4:E:91:ILE:HG23	4:E:216:LEU:HD22	1.91	0.52
1:A:173:SER:HA	1:A:184:VAL:HG13	1.92	0.51
4:E:205:MET:HB3	4:E:228:PRO:HG2	1.92	0.51
4:E:94:LEU:O	4:E:98:TYR:HB2	2.10	0.51
4:E:177:LEU:HB3	4:E:178:PHE:HD1	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:207:ARG:HH21	4:E:231:ASP:HA	1.76	0.51
4:E:187:PHE:HB2	4:E:195:LEU:HD21	1.93	0.49
1:A:216:ASP:HB3	1:A:222:LEU:HB2	1.94	0.48
1:A:396:TRP:HB2	2:B:215:VAL:HG11	1.94	0.48
1:A:219:VAL:HG23	1:A:220:PRO:HD3	1.95	0.48
1:A:25:GLY:H	5:A:1550:STU:C26	2.27	0.48
4:E:27:VAL:HG11	4:E:185:PRO:HG3	1.95	0.48
4:E:233:LYS:HB2	4:E:235:ARG:HG3	1.96	0.47
1:A:422:TYR:CE1	1:A:436:ARG:HG3	2.50	0.47
4:E:55:GLN:HA	4:E:112:LYS:HA	1.97	0.47
1:A:366:PHE:HD1	4:E:64:VAL:HG13	1.80	0.46
2:B:208:LEU:HD12	2:B:209:PRO:HD2	1.97	0.46
2:B:209:PRO:HA	2:B:210:PRO:HD3	1.78	0.46
4:E:149:ILE:HG12	4:E:151:ARG:H	1.81	0.46
4:E:58:LYS:HB2	4:E:58:LYS:HE3	1.81	0.46
4:E:100:SER:HA	4:E:101:ALA:HA	1.53	0.45
1:A:137:HIS:NE2	1:A:157:ASP:O	2.47	0.45
4:E:248:LEU:HA	4:E:253:THR:HG23	1.98	0.45
1:A:192:GLY:N	1:A:194:GLU:OE2	2.48	0.45
4:E:298:ARG:NH1	6:E:1325:AMP:O1P	2.32	0.45
4:E:176:LYS:HG2	4:E:292:VAL:HG21	1.98	0.45
1:A:250:GLN:HB2	1:A:256:ARG:HA	1.97	0.45
1:A:162:ASN:ND2	1:A:170:LEU:HD22	2.32	0.45
1:A:233:THR:HB	1:A:237:LEU:HD23	2.00	0.44
1:A:258:THR:O	1:A:262:ILE:HG13	2.18	0.44
4:E:180:THR:HG22	4:E:182:PHE:H	1.82	0.44
4:E:138:PHE:CE1	4:E:322:VAL:HG23	2.52	0.44
1:A:410:MET:HE3	1:A:451:LEU:HB2	1.99	0.44
1:A:185:ILE:HD11	1:A:227:CYS:SG	2.58	0.43
4:E:251:GLU:HB2	4:E:253:THR:HG22	2.00	0.43
1:A:465:SER:HB3	2:B:239:ASN:HB3	2.01	0.43
4:E:141:VAL:HG22	4:E:171:ILE:HD13	2.00	0.43
1:A:247:HIS:CG	1:A:257:ALA:HB2	2.54	0.43
4:E:62:ALA:O	4:E:66:ASN:ND2	2.43	0.43
4:E:117:ARG:O	4:E:121:LEU:HB3	2.18	0.43
2:B:242:TYR:O	2:B:254:SER:N	2.52	0.43
4:E:269:SER:OG	4:E:270:HIS:N	2.52	0.43
4:E:83:GLY:HA2	4:E:128:LEU:HD13	2.00	0.43
1:A:193:PRO:O	1:A:197:ILE:HG12	2.20	0.42
4:E:93:ILE:HG12	4:E:116:TRP:CZ2	2.55	0.42
4:E:48:VAL:O	4:E:72:PRO:HD2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:ARG:HB3	1:A:460:LEU:HD23	2.01	0.41
4:E:246:ILE:O	4:E:250:ALA:N	2.52	0.41
1:A:36:GLU:HG3	1:A:37:LEU:HD13	2.02	0.41
1:A:116:SER:HB2	1:A:207:LEU:HB3	2.01	0.41
1:A:146:LEU:HD21	5:A:1550:STU:H241	2.03	0.41
1:A:274:LYS:HG3	1:A:280:ASP:H	1.86	0.41
1:A:342:LEU:HD23	1:A:342:LEU:HA	1.95	0.41
1:A:373:ARG:HA	1:A:373:ARG:HD3	1.91	0.41
4:E:180:THR:O	4:E:183:PRO:HD2	2.21	0.41
4:E:238:ASP:OD1	4:E:268:ARG:NH2	2.54	0.41
1:A:468:ASP:OD1	1:A:468:ASP:N	2.54	0.41
4:E:114:GLU:O	4:E:118:GLU:HG3	2.21	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	395/493 (80%)	358 (91%)	35 (9%)	2 (0%)	29	64
2	B	67/87 (77%)	61 (91%)	6 (9%)	0	100	100
3	C	19/27 (70%)	19 (100%)	0	0	100	100
4	E	299/330 (91%)	280 (94%)	16 (5%)	3 (1%)	15	50
All	All	780/937 (83%)	718 (92%)	57 (7%)	5 (1%)	25	61

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	E	26	SER
1	A	27	PHE
1	A	106	CYS

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Mol	Chain	Res	Type
4	E	27	VAL
4	E	124	SER

### 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	346/446 (78%)	323 (93%)	23 (7%)	16	48
2	B	66/81 (82%)	62 (94%)	4 (6%)	18	51
3	C	18/23 (78%)	16 (89%)	2 (11%)	6	24
4	E	274/299 (92%)	257 (94%)	17 (6%)	18	51
All	All	704/849 (83%)	658 (94%)	46 (6%)	17	49

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	PHE
1	A	41	LYS
1	A	88	ASP
1	A	111	LEU
1	A	112	ASP
1	A	166	ASP
1	A	178	ASN
1	A	189	LEU
1	A	194	GLU
1	A	207	LEU
1	A	219	VAL
1	A	224	LYS
1	A	228	ASP
1	A	247	HIS
1	A	348	ASP
1	A	357	THR
1	A	366	PHE
1	A	371	THR
1	A	401	ARG

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Mol	Chain	Res	Type
1	A	405	ARG
1	A	407	ASN
1	A	445	SER
1	A	457	ARG
2	B	219	LYS
2	B	236	VAL
2	B	251	MET
2	B	258	ARG
3	C	537	GLU
3	C	542	LEU
4	E	36	ARG
4	E	106	TYR
4	E	111	HIS
4	E	120	TYR
4	E	167	THR
4	E	177	LEU
4	E	207	ARG
4	E	208	THR
4	E	253	THR
4	E	255	ASN
4	E	272	PHE
4	E	275	VAL
4	E	291	LEU
4	E	296	VAL
4	E	298	ARG
4	E	302	VAL
4	E	314	LEU

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

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	0.88	0	10,14,16	1.02	1 (10%)

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	-	5/9/11/13	-

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	172	TPO	O-C-CA	-2.16	119.12	124.78

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	172	TPO	C-CA-CB-CG2
1	A	172	TPO	CB-OG1-P-O1P
1	A	172	TPO	CB-OG1-P-O2P
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

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

3 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
5	STU	A	1550	-	30,42,42	0.97	2 (6%)	31,68,68	2.23	8 (25%)
6	AMP	E	1325	-	22,25,25	0.91	1 (4%)	25,38,38	1.18	2 (8%)
6	AMP	E	1326	-	22,25,25	0.88	1 (4%)	25,38,38	1.19	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
5	STU	A	1550	-	-	1/4/42/42	-
6	AMP	E	1325	-	-	3/6/26/26	0/3/3/3
6	AMP	E	1326	-	-	0/6/26/26	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	E	1325	AMP	C5-C4	2.54	1.47	1.40
6	E	1326	AMP	C5-C4	2.45	1.47	1.40
5	A	1550	STU	O4-C25	-2.27	1.40	1.43
5	A	1550	STU	C10-C11	-2.07	1.39	1.42

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1550	STU	C9-N1-C8	-5.56	108.50	113.85
5	A	1550	STU	C26-C21-C22	-4.65	103.58	112.64
5	A	1550	STU	C10-C9-N1	4.52	106.36	101.76
5	A	1550	STU	C27-O6-C22	-4.15	107.29	114.44
5	A	1550	STU	C7-C8-N1	3.85	110.27	106.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1550	STU	O5-C8-C7	-3.44	125.24	129.32
6	E	1326	AMP	N3-C2-N1	-3.05	123.91	128.68
6	E	1325	AMP	N3-C2-N1	-3.04	123.93	128.68
6	E	1325	AMP	C4-C5-N7	-2.64	106.64	109.40
6	E	1326	AMP	C4-C5-N7	-2.59	106.70	109.40
5	A	1550	STU	C3-C4-C5	-2.31	116.98	120.86
5	A	1550	STU	C13-C12-C17	2.21	122.19	119.39

There are no chirality outliers.

All (4) torsion outliers are listed below:

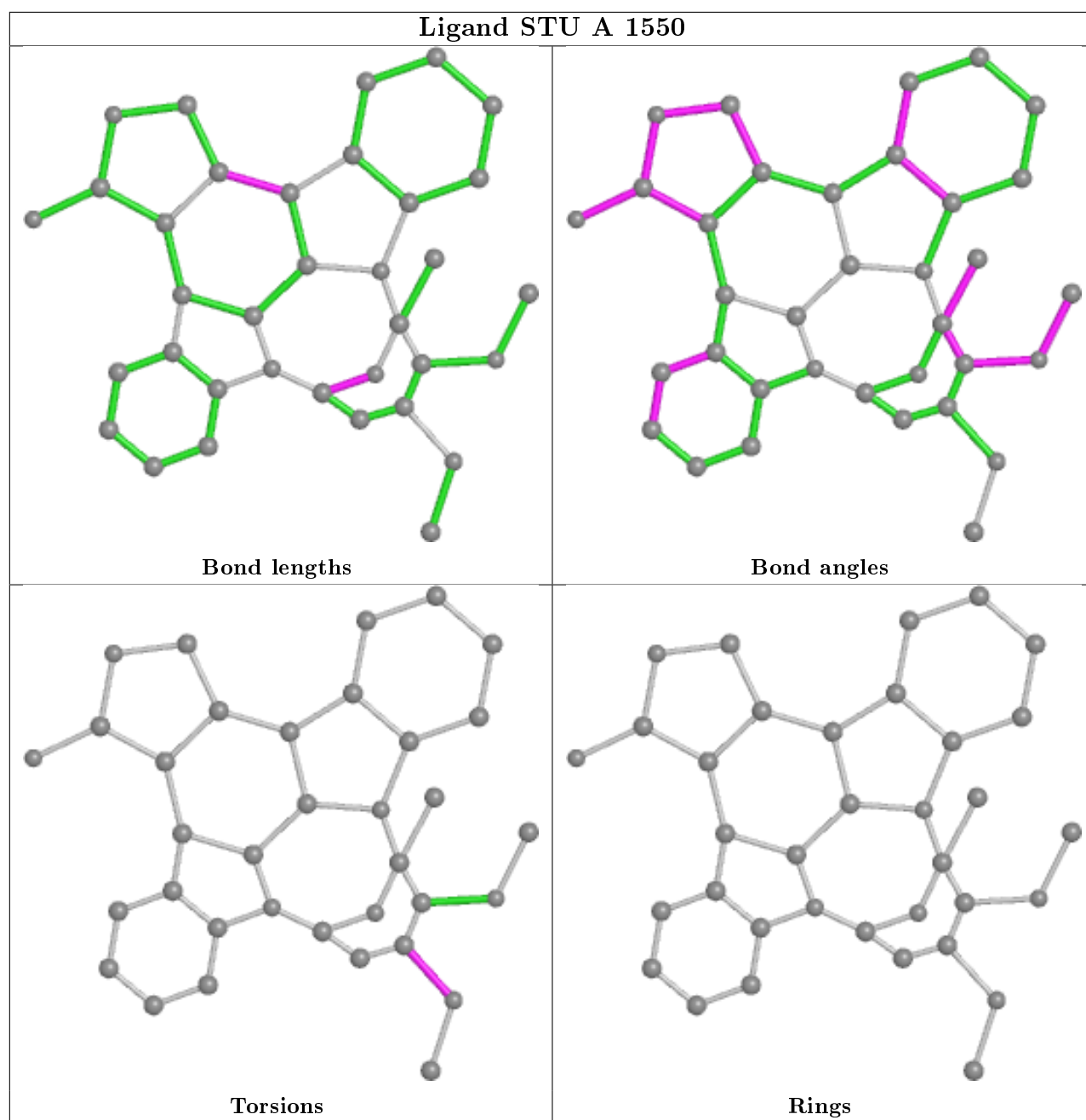
Mol	Chain	Res	Type	Atoms
6	E	1325	AMP	C5'-O5'-P-O2P
6	E	1325	AMP	C5'-O5'-P-O3P
6	E	1325	AMP	C5'-O5'-P-O1P
5	A	1550	STU	C24-C23-N4-C28

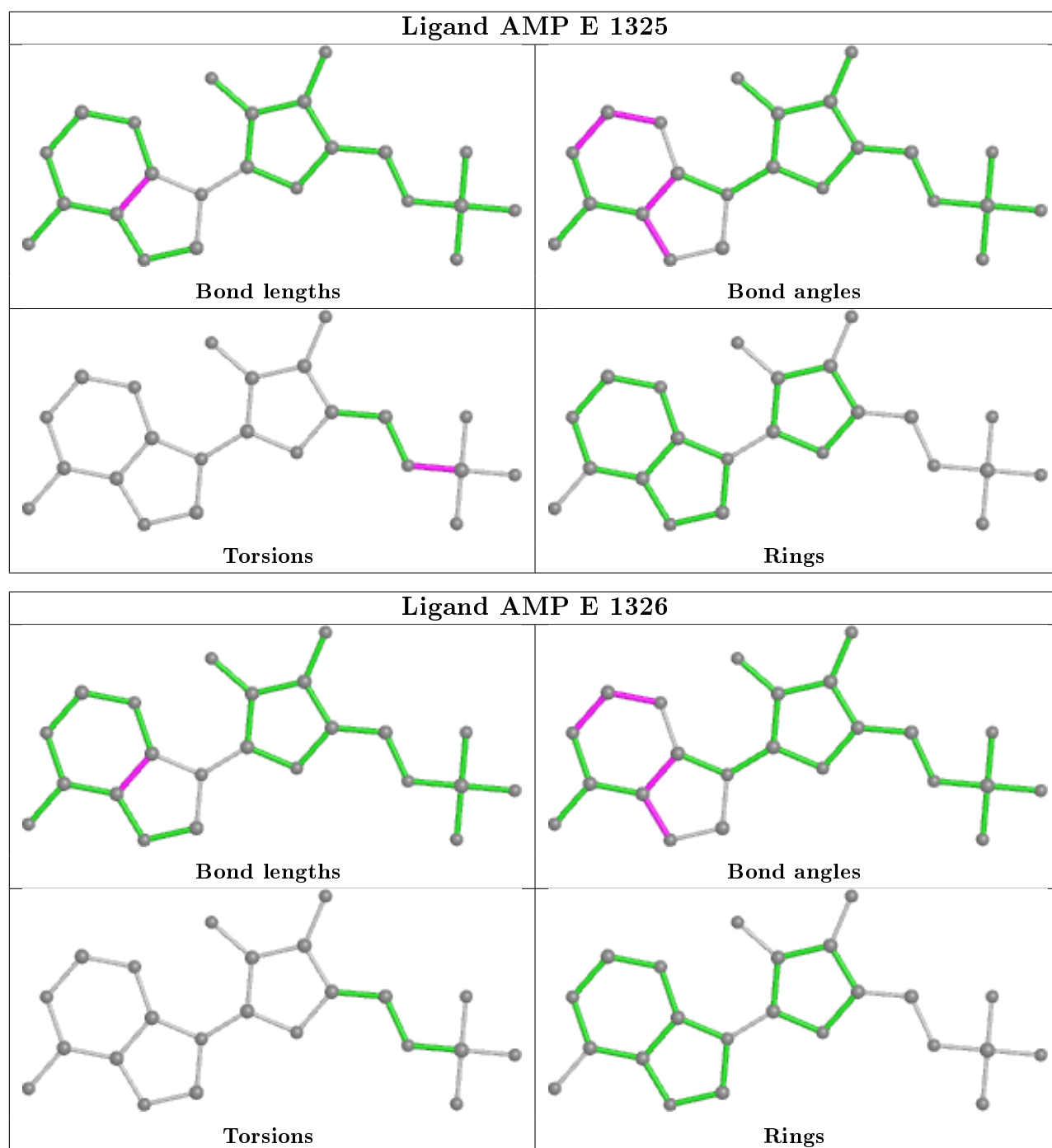
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1550	STU	4	0
6	E	1325	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	401/493 (81%)	0.14	17 (4%) 36 25	66, 108, 156, 196	0
2	B	69/87 (79%)	-0.08	0 100 100	62, 82, 134, 178	0
3	C	21/27 (77%)	-0.10	0 100 100	72, 96, 130, 166	0
4	E	301/330 (91%)	-0.05	7 (2%) 60 50	61, 89, 157, 199	0
All	All	792/937 (84%)	0.04	24 (3%) 50 39	61, 97, 157, 199	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	470	ILE	4.7
4	E	123	ASP	4.6
1	A	174	CYS	3.8
1	A	374	ALA	3.7
1	A	109	GLY	3.4
1	A	469	GLU	3.2
4	E	24	ASN	2.9
4	E	100	SER	2.7
1	A	437	LYS	2.6
1	A	443	THR	2.6
4	E	121	LEU	2.5
4	E	270	HIS	2.5
1	A	235	GLN	2.4
1	A	373	ARG	2.3
1	A	370	GLU	2.3
4	E	126	LYS	2.2
1	A	108	ASN	2.2
1	A	324	TYR	2.1
4	E	127	PRO	2.1
1	A	28	GLY	2.1
1	A	468	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	106	CYS	2.1
1	A	457	ARG	2.1
1	A	321	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	0.94	0.14	77,107,118,121	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

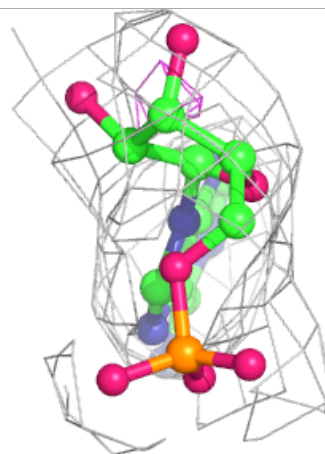
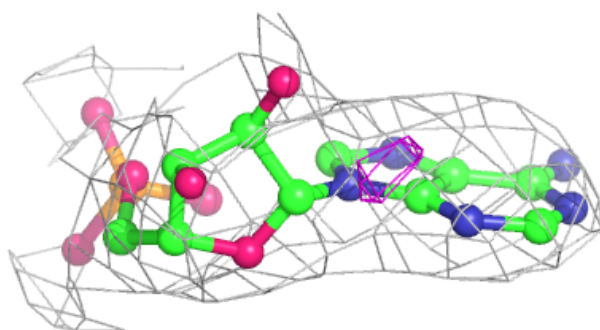
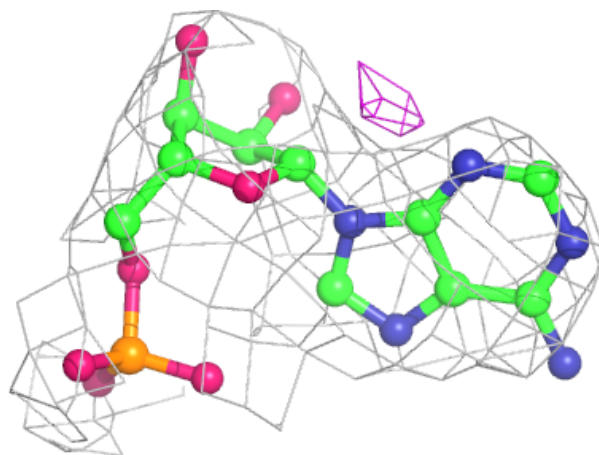
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
6	AMP	E	1325	23/23	0.94	0.23	67,74,79,89	0
5	STU	A	1550	35/35	0.95	0.21	84,110,121,126	0
6	AMP	E	1326	23/23	0.95	0.20	75,80,84,84	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.

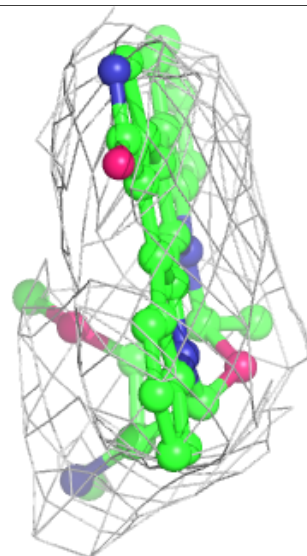
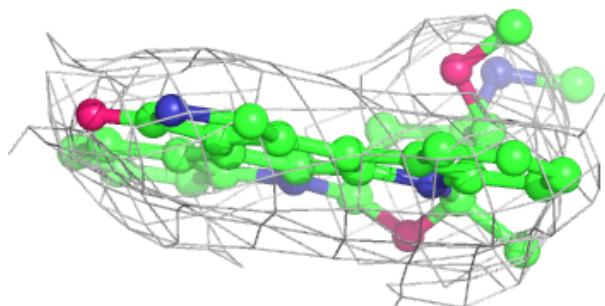
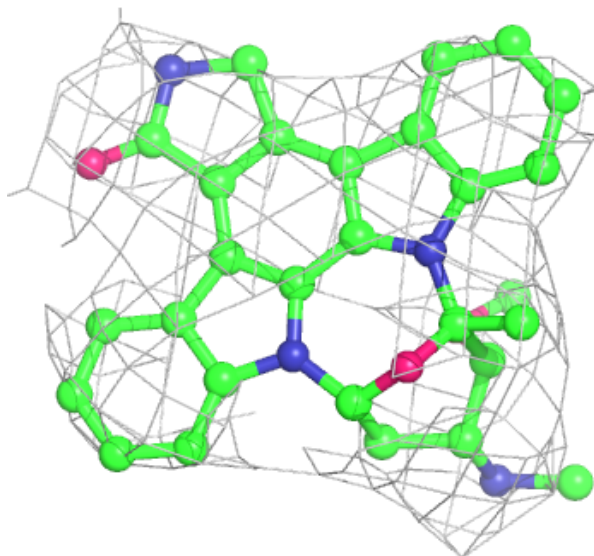
**Electron density around AMP E 1325:**

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



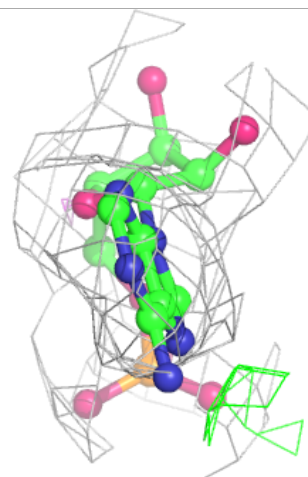
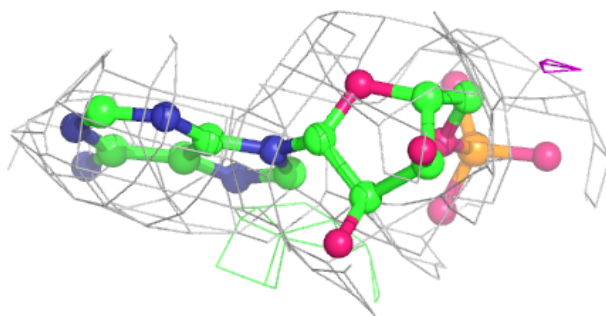
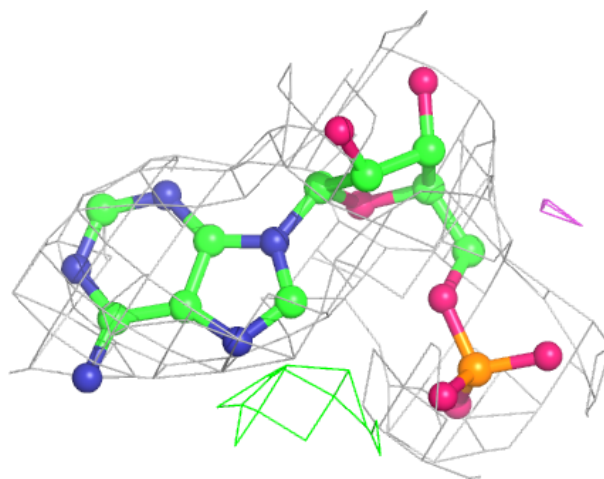
**Electron density around STU A 1550:**

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



**Electron density around AMP E 1326:**

$2mF_o-DF_c$  (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.