



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:04 PM UTC

PDB ID : 2Y8Q / pdb_00002y8q
Title : Structure of the regulatory fragment of mammalian AMPK in complex with one ADP
Authors : Xiao, B.; Sanders, M.J.; Underwood, E.; Heath, R.; Mayer, F.; Carmena, D.; Jing, C.; Walker, P.A.; Eccleston, J.F.; Haire, L.F.; Saiu, P.; Howell, S.A.; Aasland, R.; Martin, S.R.; Carling, D.; Gamblin, S.J.
Deposited on : 2011-02-09
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ 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](#) ①) 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

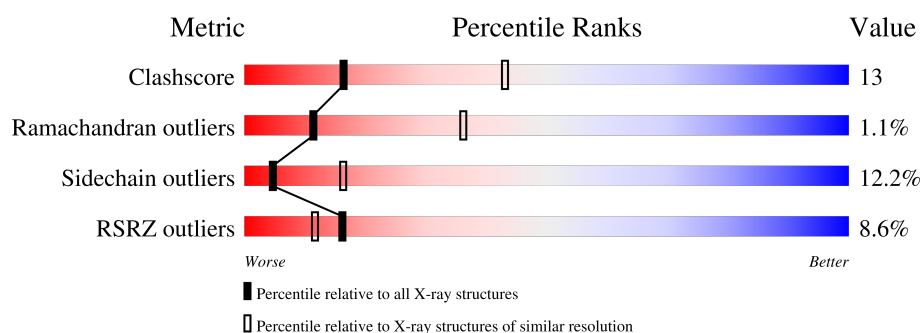
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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))
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	173	2% 42% 16% . 40%
2	B	87	7% 62% 15% . . 18%
3	E	330	9% 62% 24% 5% . 8%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3987 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	104	Total	C	N	O	S	0	0	0
			854	539	152	156	7			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	378	MET	-	expression tag	UNP P54645
A	379	SER	-	expression tag	UNP P54645
A	380	HIS	-	expression tag	UNP P54645
A	381	HIS	-	expression tag	UNP P54645
A	382	HIS	-	expression tag	UNP P54645
A	383	HIS	-	expression tag	UNP P54645
A	384	HIS	-	expression tag	UNP P54645
A	385	HIS	-	expression tag	UNP P54645
A	386	SER	-	expression tag	UNP P54645
A	387	GLY	-	expression tag	UNP P54645
A	388	LEU	-	expression tag	UNP P54645
A	389	VAL	-	expression tag	UNP P54645
A	390	PRO	-	expression tag	UNP P54645
A	391	ARG	-	expression tag	UNP P54645
A	392	GLY	-	expression tag	UNP P54645
A	393	SER	-	expression tag	UNP P54645
A	394	MET	-	expression tag	UNP P54645
A	395	ALA	-	expression tag	UNP P54645
A	545	ASN	-	expression tag	UNP P54645
A	546	SER	-	expression tag	UNP P54645
A	547	CYS	-	expression tag	UNP P54645
A	548	THR	-	expression tag	UNP P54645
A	549	VAL	-	expression tag	UNP P54645
A	550	ASN	-	expression tag	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	71	Total	C	N	O	S	0	0	0
			586	384	100	99	3			

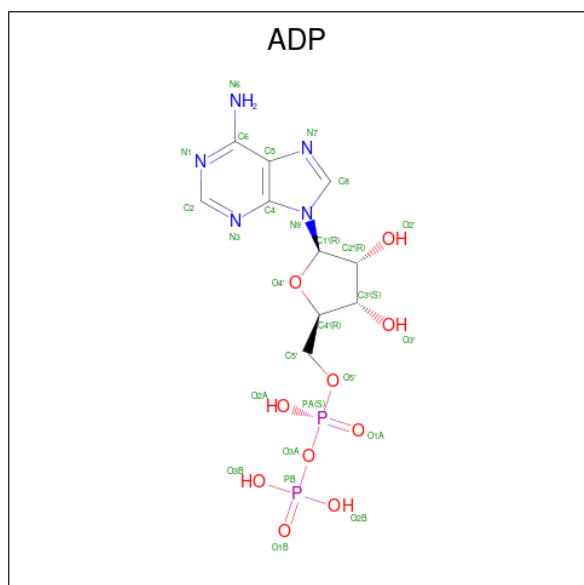
There is a discrepancy between the modelled and reference sequences:

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

- Molecule 3 is a protein called 5'-AMP-ACTIVATED PROTEIN KINASE SUBUNIT GAMMA-1.

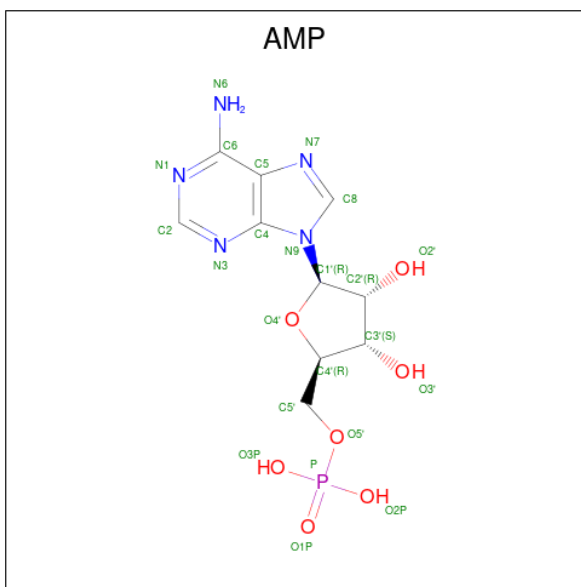
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	304	Total	C	N	O	S	0	0	0
			2441	1584	407	443	7			

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

- Molecule 5 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	6	Total	O	0	0
			6	6		
6	B	11	Total	O	0	0
			11	11		
6	E	39	Total	O	0	0
			39	39		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.03Å 119.92Å 130.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 20.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	93.8 (20.00-2.80) 93.6 (20.00-2.80)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.79Å)	Xtriage
Refinement program	REFMAC 5.5	Depositor
R, R_{free}	0.232 , 0.275 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	66.7	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3987	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% 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, ADP

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	0.55	0/872	0.83	0/1178
2	B	0.56	0/601	1.05	4/811 (0.5%)
3	E	0.75	3/2493 (0.1%)	1.16	14/3384 (0.4%)
All	All	0.68	3/3966 (0.1%)	1.08	18/5373 (0.3%)

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	A	0	1
2	B	0	1
3	E	0	4
All	All	0	6

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	273	GLU	C-N	20.07	1.62	1.33
3	E	274	GLY	N-CA	8.94	1.58	1.45
3	E	274	GLY	CA-C	8.23	1.63	1.51

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	274	GLY	O-C-N	21.57	150.74	122.70
3	E	272	PHE	CA-C-N	-17.44	94.45	122.53
3	E	272	PHE	C-N-CA	-17.44	94.45	122.53
3	E	273	GLU	CA-C-N	-12.56	96.80	121.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	273	GLU	C-N-CA	-12.56	96.80	121.41
3	E	272	PHE	O-C-N	10.14	136.08	122.59
3	E	274	GLY	CA-C-O	-8.74	105.36	120.57
2	B	235	HIS	N-CA-C	7.84	119.90	111.36
3	E	274	GLY	N-CA-C	-6.54	97.68	113.18
2	B	209	PRO	N-CA-C	6.40	118.50	110.70
3	E	270	HIS	N-CA-C	6.32	118.37	108.96
3	E	26	SER	N-CA-C	5.92	117.93	110.24
3	E	182	PHE	N-CA-C	5.89	122.84	109.81
3	E	276	LEU	N-CA-CB	-5.50	102.16	110.35
2	B	213	LEU	CA-C-N	5.24	127.25	120.44
2	B	213	LEU	C-N-CA	5.24	127.25	120.44
3	E	267	HIS	N-CA-C	-5.16	107.06	113.19
3	E	255	ASN	N-CA-C	5.01	116.75	111.28

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	546	SER	Peptide
2	B	208	LEU	Peptide
3	E	182	PHE	Peptide
3	E	272	PHE	Mainchain
3	E	273	GLU	Mainchain
3	E	274	GLY	Mainchain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	854	0	845	18	0
2	B	586	0	612	13	0
3	E	2441	0	2506	79	0
4	E	27	0	12	0	0
5	E	23	0	12	0	0
6	A	6	0	0	0	0
6	B	11	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	E	39	0	0	16	0
All	All	3987	0	3987	103	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 13.

All (103) 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:E:251:GLU:HG2	6:E:2024:HOH:O	1.68	0.94
3:E:121:LEU:C	3:E:122:GLN:HG3	2.01	0.84
2:B:208:LEU:O	2:B:210:PRO:CD	2.28	0.81
2:B:198:SER:HB2	2:B:201:ARG:HG3	1.62	0.81
2:B:208:LEU:O	2:B:210:PRO:HD3	1.83	0.78
1:A:526:PRO:HG2	3:E:128:LEU:HD23	1.67	0.77
3:E:276:LEU:HD22	3:E:276:LEU:N	2.04	0.73
3:E:323:LEU:O	3:E:324:THR:HG23	1.89	0.73
3:E:108:LEU:O	3:E:111:HIS:HB2	1.91	0.70
3:E:60:PHE:CE1	3:E:90:PHE:HB2	2.27	0.69
1:A:532:THR:H	3:E:161:ASN:HD21	1.39	0.68
3:E:248:LEU:O	6:E:2024:HOH:O	2.13	0.67
3:E:242:LYS:O	3:E:245:VAL:HB	1.96	0.66
2:B:208:LEU:O	2:B:210:PRO:HD2	1.95	0.65
3:E:121:LEU:O	3:E:122:GLN:HG3	1.97	0.65
1:A:543:ILE:O	1:A:547:CYS:HB3	1.98	0.64
3:E:112:LYS:HA	6:E:2009:HOH:O	1.96	0.64
3:E:251:GLU:CG	6:E:2024:HOH:O	2.37	0.63
3:E:91:ILE:HD13	3:E:242:LYS:HG2	1.81	0.62
3:E:274:GLY:H	3:E:276:LEU:H	1.48	0.62
3:E:273:GLU:HG3	3:E:296:VAL:HG22	1.83	0.61
3:E:275:VAL:C	3:E:276:LEU:HD22	2.25	0.60
3:E:323:LEU:C	3:E:324:THR:CG2	2.75	0.60
2:B:236:VAL:HG13	2:B:236:VAL:O	2.01	0.59
3:E:276:LEU:N	3:E:276:LEU:CD2	2.66	0.58
3:E:248:LEU:HD12	6:E:2023:HOH:O	2.03	0.58
3:E:239:ILE:HG21	3:E:273:GLU:HB3	1.84	0.58
3:E:291:LEU:HD13	3:E:299:LEU:HG	1.85	0.58
3:E:24:ASN:ND2	6:E:2001:HOH:O	2.36	0.58
2:B:210:PRO:HG2	2:B:212:LEU:HD12	1.86	0.57
3:E:272:PHE:CE2	3:E:276:LEU:HD21	2.40	0.57
3:E:282:GLU:HB2	3:E:287:ILE:CD1	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:273:GLU:HG3	3:E:296:VAL:CG2	2.35	0.56
3:E:194:GLU:HB2	6:E:2019:HOH:O	2.07	0.55
2:B:209:PRO:C	2:B:211:HIS:H	2.16	0.53
3:E:282:GLU:HB2	3:E:287:ILE:HD11	1.92	0.52
3:E:26:SER:HB3	3:E:29:THR:H	1.75	0.52
2:B:209:PRO:C	2:B:211:HIS:N	2.67	0.52
3:E:225:SER:HA	3:E:242:LYS:HD2	1.92	0.51
3:E:323:LEU:O	3:E:324:THR:CG2	2.58	0.51
3:E:56:VAL:HG23	6:E:2009:HOH:O	2.09	0.51
3:E:48:VAL:O	3:E:72:PRO:HD2	2.10	0.51
3:E:84:MET:HE3	3:E:153:PRO:HG3	1.93	0.50
1:A:445:SER:HB3	1:A:538:MET:HE1	1.93	0.50
3:E:274:GLY:N	3:E:276:LEU:H	2.09	0.50
3:E:257:LEU:HB2	6:E:2027:HOH:O	2.10	0.50
1:A:410:MET:HE3	1:A:451:LEU:HD22	1.93	0.49
3:E:266:GLN:C	3:E:268:ARG:H	2.18	0.49
3:E:245:VAL:HA	6:E:2022:HOH:O	2.12	0.49
3:E:218:ILE:HG23	3:E:222:HIS:HD2	1.79	0.48
3:E:91:ILE:HG13	3:E:245:VAL:HG11	1.95	0.48
1:A:447:MET:HE2	1:A:538:MET:HB3	1.96	0.48
1:A:528:PRO:HA	3:E:80:SER:CB	2.44	0.48
1:A:407:ASN:OD1	1:A:407:ASN:N	2.40	0.47
3:E:169:LYS:NZ	3:E:295:GLU:HB3	2.29	0.47
2:B:262:LYS:HE3	3:E:38:TYR:CZ	2.49	0.47
3:E:112:LYS:CA	6:E:2009:HOH:O	2.59	0.47
3:E:191:SER:HG	3:E:194:GLU:HG2	1.79	0.47
1:A:468:ASP:O	1:A:469:GLU:HB2	2.14	0.47
1:A:445:SER:HB3	1:A:538:MET:CE	2.45	0.46
3:E:111:HIS:O	6:E:2009:HOH:O	2.21	0.46
3:E:276:LEU:HD13	3:E:276:LEU:HA	1.57	0.46
3:E:40:LEU:HD21	3:E:137:LEU:HD22	1.97	0.46
3:E:243:PHE:C	3:E:245:VAL:H	2.24	0.46
3:E:121:LEU:N	6:E:2010:HOH:O	2.49	0.45
1:A:528:PRO:HA	3:E:80:SER:HB2	1.97	0.45
3:E:87:ILE:HG23	3:E:246:ILE:HG23	1.98	0.45
3:E:248:LEU:C	6:E:2024:HOH:O	2.56	0.45
3:E:56:VAL:N	6:E:2009:HOH:O	2.35	0.45
3:E:120:TYR:C	3:E:120:TYR:CD1	2.94	0.45
3:E:118:GLU:C	6:E:2010:HOH:O	2.60	0.45
1:A:416:ALA:CB	1:A:546:SER:HB3	2.47	0.45
3:E:305:HIS:O	3:E:306:ASP:HB2	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:191:SER:OG	3:E:194:GLU:HG2	2.16	0.44
3:E:241:SER:O	3:E:244:ASP:HB2	2.18	0.44
3:E:60:PHE:CD1	3:E:90:PHE:HB2	2.52	0.44
2:B:258:ARG:HG3	2:B:263:TYR:CE1	2.52	0.44
1:A:428:ASN:HD22	1:A:431:TYR:HB3	1.83	0.43
2:B:198:SER:CB	2:B:201:ARG:HG3	2.43	0.43
3:E:243:PHE:C	3:E:245:VAL:N	2.75	0.43
1:A:409:ILE:O	1:A:413:VAL:HG13	2.18	0.43
3:E:87:ILE:HD12	3:E:246:ILE:HG12	2.00	0.43
3:E:120:TYR:O	3:E:122:GLN:HG3	2.19	0.43
3:E:323:LEU:C	3:E:324:THR:HG22	2.44	0.43
3:E:188:MET:HE3	3:E:285:GLU:HB2	2.01	0.42
3:E:299:LEU:HB2	3:E:312:VAL:HG13	2.00	0.42
1:A:533:ILE:HG21	3:E:74:TRP:CD2	2.54	0.42
3:E:243:PHE:CE1	3:E:246:ILE:HD11	2.54	0.42
1:A:416:ALA:HB2	1:A:546:SER:HB3	2.01	0.42
3:E:120:TYR:O	3:E:122:GLN:CG	2.68	0.42
1:A:530:SER:O	1:A:534:GLU:HG3	2.20	0.41
1:A:451:LEU:HD23	2:B:205:PRO:HG3	2.02	0.41
3:E:184:LYS:HA	3:E:185:PRO:HD3	1.92	0.41
3:E:266:GLN:C	3:E:268:ARG:N	2.78	0.41
3:E:273:GLU:HG2	3:E:276:LEU:HB2	2.01	0.41
3:E:73:LEU:HD21	3:E:85:LEU:HB2	2.02	0.41
3:E:291:LEU:CD1	3:E:299:LEU:HG	2.50	0.41
3:E:23:SER:HB3	3:E:187:PHE:HD2	1.85	0.41
3:E:120:TYR:C	3:E:120:TYR:HD1	2.27	0.41
3:E:120:TYR:O	3:E:120:TYR:HD1	2.04	0.40
3:E:242:LYS:O	3:E:245:VAL:CB	2.68	0.40
3:E:55:GLN:HB2	3:E:58:LYS:HD2	2.02	0.40
2:B:247:LYS:O	2:B:250:VAL:HG12	2.22	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	100/173 (58%)	99 (99%)	1 (1%)	0	100	100
2	B	67/87 (77%)	58 (87%)	8 (12%)	1 (2%)	8	28
3	E	302/330 (92%)	276 (91%)	22 (7%)	4 (1%)	9	31
All	All	469/590 (80%)	433 (92%)	31 (7%)	5 (1%)	11	36

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	271	TYR
3	E	272	PHE
3	E	256	ASN
2	B	208	LEU
3	E	182	PHE

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	97/157 (62%)	85 (88%)	12 (12%)	4	16
2	B	67/81 (83%)	64 (96%)	3 (4%)	24	58
3	E	277/299 (93%)	238 (86%)	39 (14%)	3	12
All	All	441/537 (82%)	387 (88%)	54 (12%)	5	16

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	394	MET
1	A	407	ASN
1	A	413	VAL
1	A	420	LEU
1	A	433	ARG

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Mol	Chain	Res	Type
1	A	434	VAL
1	A	436	ARG
1	A	441	THR
1	A	445	SER
1	A	460	LEU
1	A	527	ARG
1	A	548	THR
2	B	211	HIS
2	B	236	VAL
2	B	272	ILE
3	E	24	ASN
3	E	40	LEU
3	E	54	LEU
3	E	65	THR
3	E	68	VAL
3	E	91	ILE
3	E	94	LEU
3	E	96	ARG
3	E	103	VAL
3	E	108	LEU
3	E	109	GLU
3	E	110	GLU
3	E	111	HIS
3	E	122	GLN
3	E	129	VAL
3	E	149	ILE
3	E	159	SER
3	E	169	LYS
3	E	180	THR
3	E	193	GLU
3	E	208	THR
3	E	224	VAL
3	E	239	ILE
3	E	241	SER
3	E	248	LEU
3	E	261	VAL
3	E	273	GLU
3	E	284	LEU
3	E	287	ILE
3	E	291	LEU
3	E	296	VAL
3	E	302	VAL

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Mol	Chain	Res	Type
3	E	308	VAL
3	E	309	LYS
3	E	311	ILE
3	E	315	SER
3	E	317	ILE
3	E	318	LEU
3	E	324	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	428	ASN
1	A	453	GLN
3	E	161	ASN
3	E	222	HIS
3	E	247	ASN
3	E	267	HIS
3	E	270	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	ADP	E	1327	-	28,29,29	1.47	5 (17%)	43,45,45	1.91	10 (23%)
5	AMP	E	1328	-	25,25,25	1.48	4 (16%)	37,38,38	1.94	10 (27%)

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	ADP	E	1327	-	-	3/16/32/32	0/3/3/3
5	AMP	E	1328	-	-	3/10/26/26	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1327	ADP	C5-C4	4.77	1.47	1.39
5	E	1328	AMP	C5-C4	4.74	1.47	1.39
4	E	1327	ADP	C5-C6	2.90	1.49	1.41
5	E	1328	AMP	C8-N7	2.65	1.36	1.31
5	E	1328	AMP	C5-C6	2.58	1.48	1.41
4	E	1327	ADP	C8-N7	2.48	1.36	1.31
5	E	1328	AMP	C5-N7	-2.48	1.34	1.39
4	E	1327	ADP	C5-N7	-2.14	1.35	1.39
4	E	1327	ADP	PA-O3A	2.11	1.61	1.59

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	1327	ADP	C5-C4-N3	-5.94	118.54	126.72
5	E	1328	AMP	C5-C4-N3	-5.28	119.44	126.72
4	E	1327	ADP	N3-C4-N9	4.72	135.20	127.17
5	E	1328	AMP	N3-C2-N1	-4.22	122.19	128.58
5	E	1328	AMP	N3-C4-N9	3.92	133.83	127.17
5	E	1328	AMP	C2-N3-C4	3.78	121.06	111.83
4	E	1327	ADP	C4-C5-N7	-3.66	106.40	110.58
4	E	1327	ADP	C2-N3-C4	3.65	120.74	111.83
5	E	1328	AMP	C4-C5-N7	-3.61	106.46	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	1327	ADP	C4-N9-C8	2.98	108.87	105.74
4	E	1327	ADP	N3-C2-N1	-2.94	124.14	128.58
4	E	1327	ADP	C5-N7-C8	2.80	107.84	103.45
5	E	1328	AMP	C2-N1-C6	2.74	123.23	118.73
5	E	1328	AMP	C5-N7-C8	2.66	107.64	103.45
5	E	1328	AMP	C6-C5-N7	2.52	136.95	132.09
4	E	1327	ADP	N9-C8-N7	-2.34	110.62	113.94
5	E	1328	AMP	C4-N9-C8	2.33	108.18	105.74
4	E	1327	ADP	O3B-PB-O2B	2.23	116.16	107.80
5	E	1328	AMP	N9-C8-N7	-2.07	111.00	113.94
4	E	1327	ADP	O3A-PA-O1A	-2.02	104.63	110.70

There are no chirality outliers.

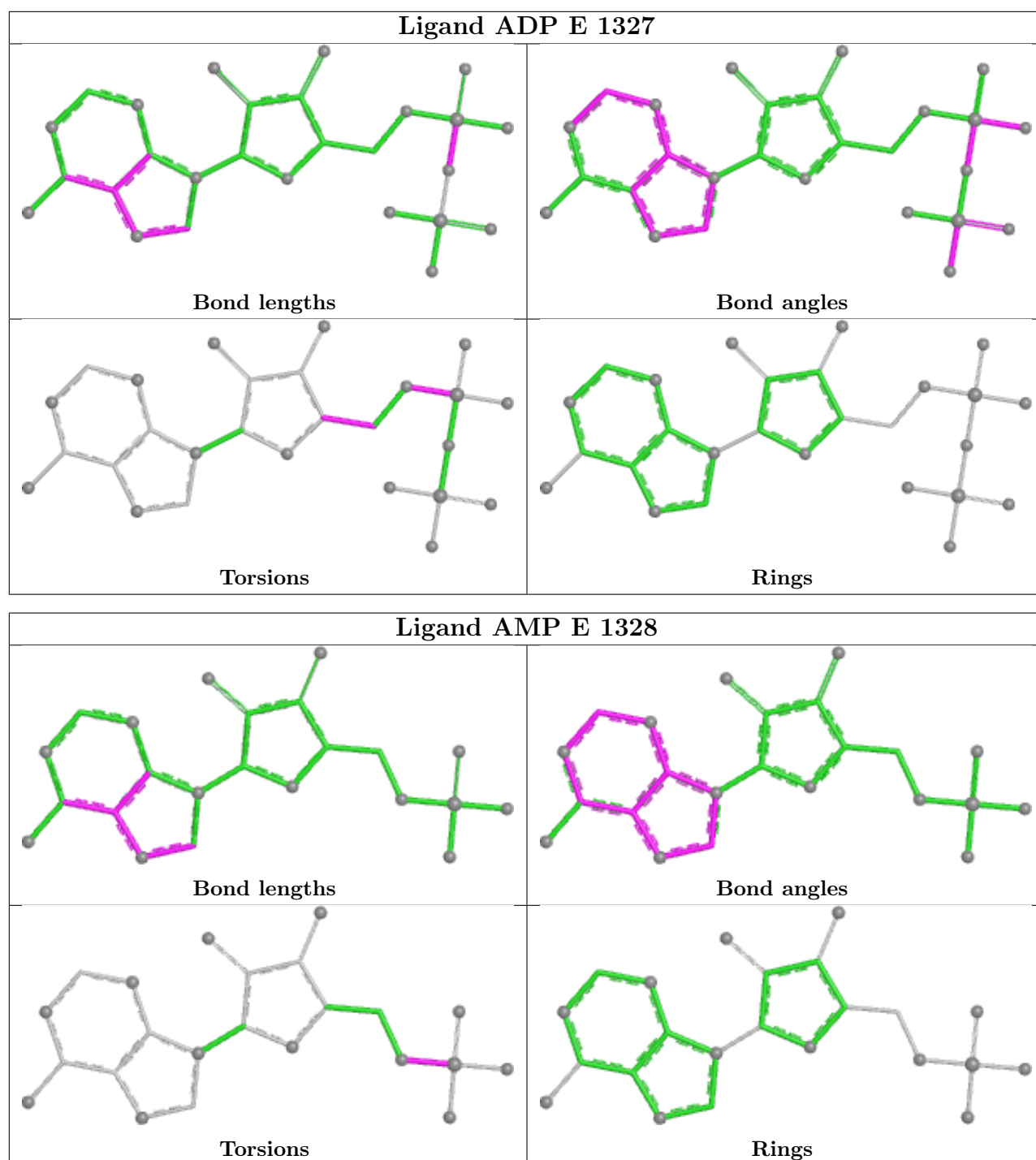
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	1327	ADP	C5'-O5'-PA-O1A
5	E	1328	AMP	C5'-O5'-P-O3P
4	E	1327	ADP	O4'-C4'-C5'-O5'
5	E	1328	AMP	C5'-O5'-P-O1P
4	E	1327	ADP	C3'-C4'-C5'-O5'
5	E	1328	AMP	C5'-O5'-P-O2P

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	E	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	273:GLU	C	274:GLY	N	1.62

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, 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	104/173 (60%)	0.12	4 (3%) 44 35	42, 64, 98, 116	0
2	B	71/87 (81%)	0.56	6 (8%) 16 12	43, 88, 117, 137	0
3	E	304/330 (92%)	0.37	31 (10%) 12 9	35, 60, 107, 128	6 (1%)
All	All	479/590 (81%)	0.34	41 (8%) 16 12	35, 62, 110, 137	6 (1%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	275	VAL	8.2
3	E	271	TYR	7.4
3	E	326	GLY	7.1
3	E	274	GLY	6.6
3	E	249	ALA	6.5
3	E	270	HIS	5.3
3	E	272	PHE	5.2
3	E	325	GLY	4.5
3	E	273	GLU	4.0
3	E	90	PHE	3.7
3	E	246	ILE	3.7
3	E	116	TRP	3.6
3	E	250	ALA	3.6
3	E	253	THR	3.5
1	A	524	VAL	3.5
3	E	125	PHE	3.4
3	E	232	GLU	3.2
2	B	213	LEU	3.1
3	E	23	SER	3.0
3	E	24	ASN	2.8
3	E	121	LEU	2.8
3	E	257	LEU	2.6
3	E	254	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	393	SER	2.5
3	E	108	LEU	2.4
3	E	252	LYS	2.4
3	E	324	THR	2.4
2	B	235	HIS	2.4
2	B	233	PRO	2.4
3	E	109	GLU	2.3
3	E	182	PHE	2.3
2	B	198	SER	2.2
3	E	25	SER	2.2
2	B	236	VAL	2.2
2	B	215	VAL	2.2
3	E	213	TYR	2.2
1	A	467	ASP	2.1
1	A	550	ASN	2.1
3	E	247	ASN	2.1
3	E	248	LEU	2.1
3	E	256	ASN	2.0

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 oligosaccharides in this entry.

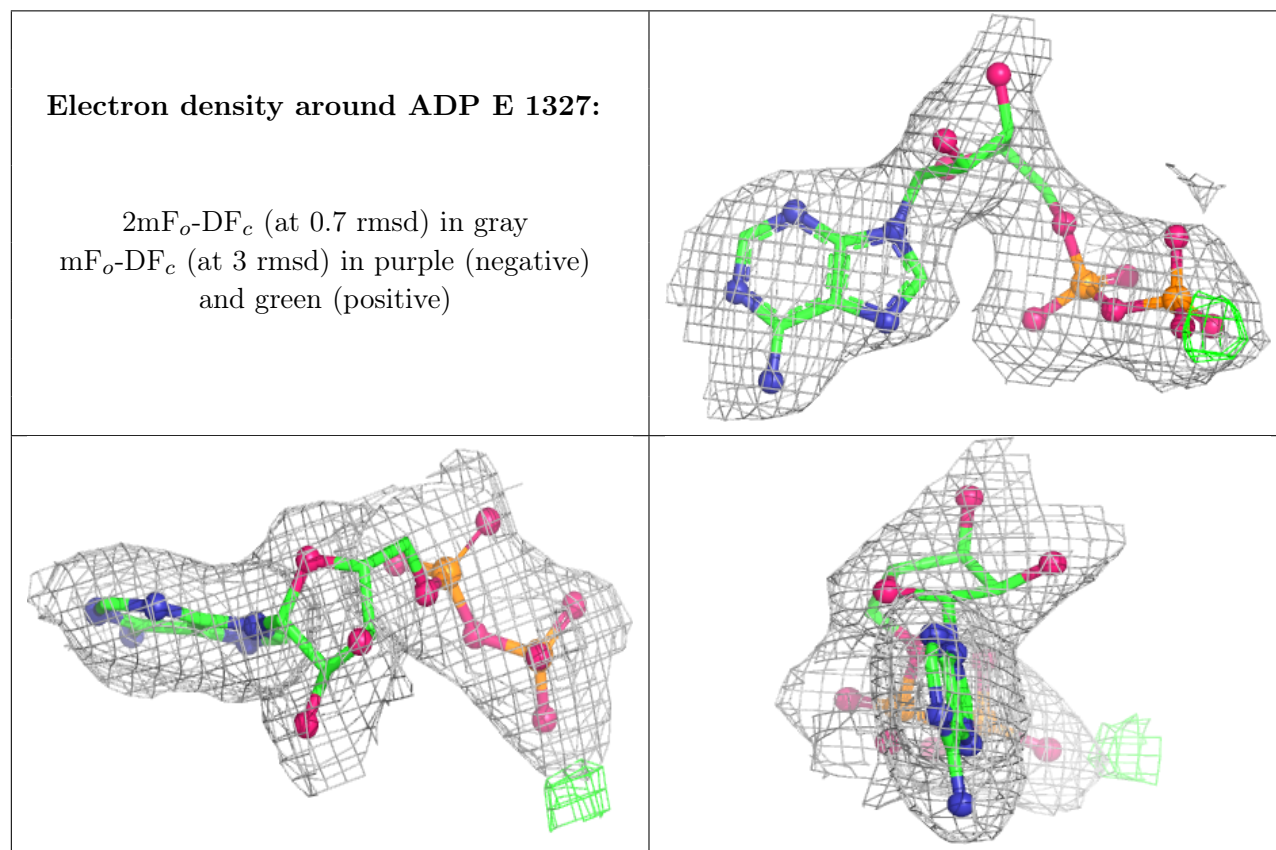
6.4 Ligands [i](#)

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, 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($\text{\AA}^2$)	Q<0.9
4	ADP	E	1327	27/27	0.96	0.07	44,48,51,53	0
5	AMP	E	1328	23/23	0.98	0.05	37,39,45,46	0

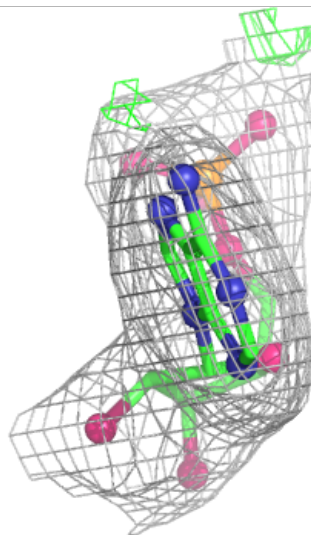
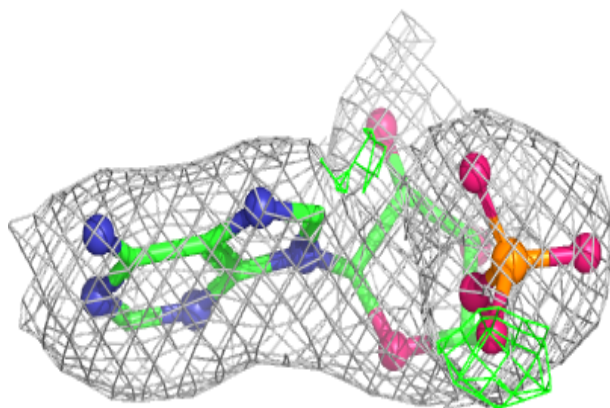
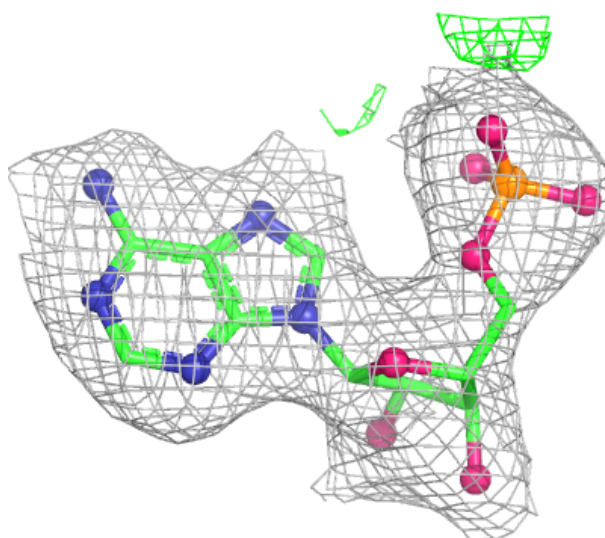
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

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 1328:

$2mF_o-DF_c$ (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.