



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 06:05 AM UTC

PDB ID : 6E4T / pdb_00006e4t
Title : Structure of AMPK bound to activator
Authors : Calabrese, M.F.; Kurumbail, R.G.
Deposited on : 2018-07-18
Resolution : 3.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

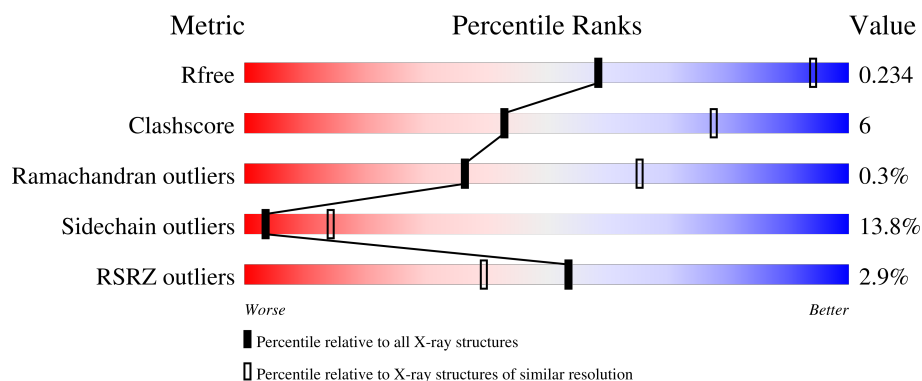
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	503	0% 56% 14% • 28%
2	B	204	3% 53% 23% • 22%
3	C	330	4% 65% 17% • 15%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 6369 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
1	A	362	Total	C	N	O	P	S	0	0	0
			2880	1848	492	521	1	18			

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	-	linker	UNP P54645
A	518	SER	-	linker	UNP P54645
A	519	GLY	-	linker	UNP P54645
A	520	GLY	-	linker	UNP P54645
A	521	PRO	-	linker	UNP P54645
A	522	GLY	-	linker	UNP P54645
A	523	GLY	-	linker	UNP P54645
A	524	SER	-	linker	UNP P54645

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	159	Total	C	N	O	P	S	0	0	0
			1208	783	201	221	1	2			

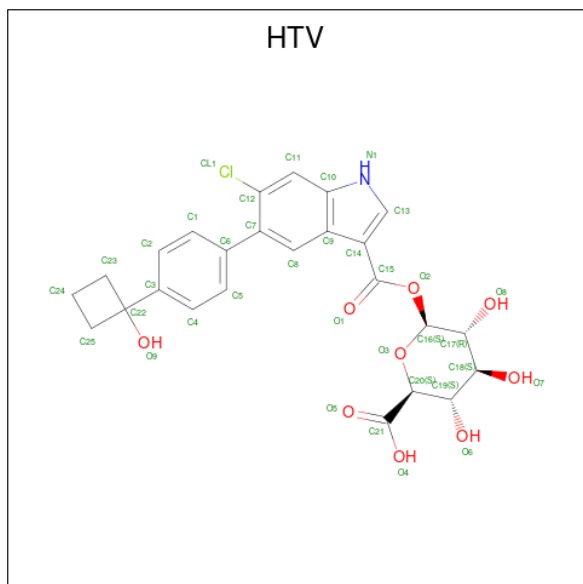
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	67	MET	-	initiating methionine	UNP P80386

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

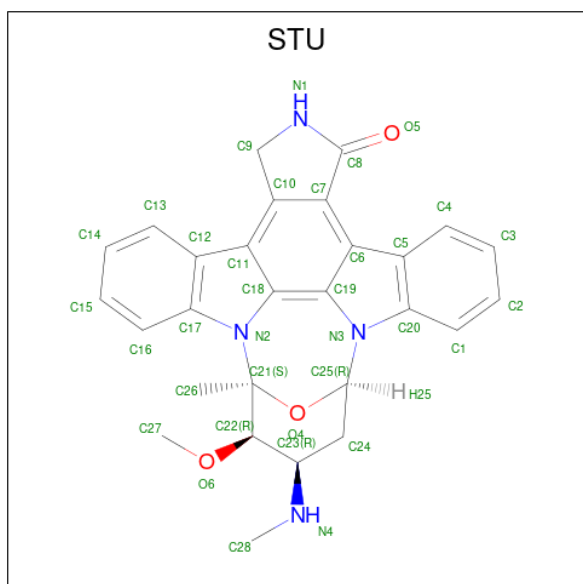
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	281	Total	C	N	O	S	0	0	0
			2127	1379	350	392	6			

- Molecule 4 is 1-O-{6-chloro-5-[4-(1-hydroxycyclobutyl)phenyl]-1H-indole-3-carbonyl}-beta-D-glucopyranuronic acid (CCD ID: HTV) (formula: $C_{25}H_{24}ClNO_9$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	Cl	N	O	0
			36	25	1	1	9	

- Molecule 5 is STAUROSPORINE (CCD ID: STU) (formula: $C_{28}H_{26}N_4O_3$).

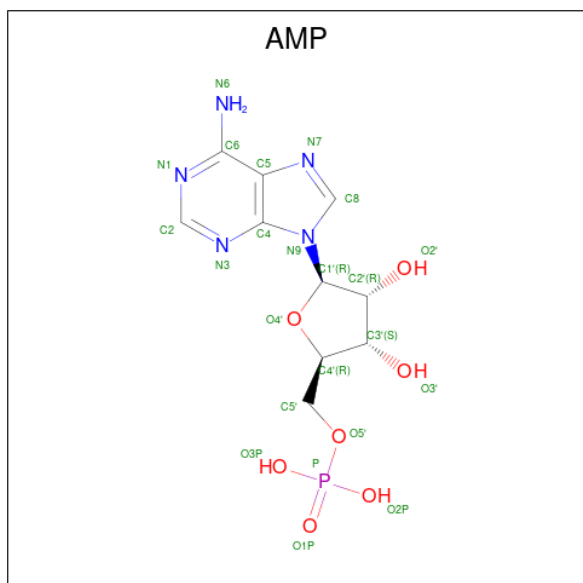


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			35	28	4	3		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

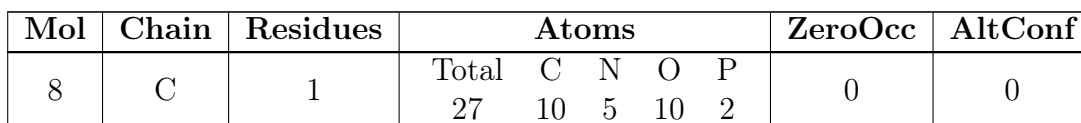
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	4	Total	Cl	0	0
			4	4		
6	B	1	Total	Cl	0	0
			1	1		

- Molecule 7 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: C₁₀H₁₄N₅O₇P).



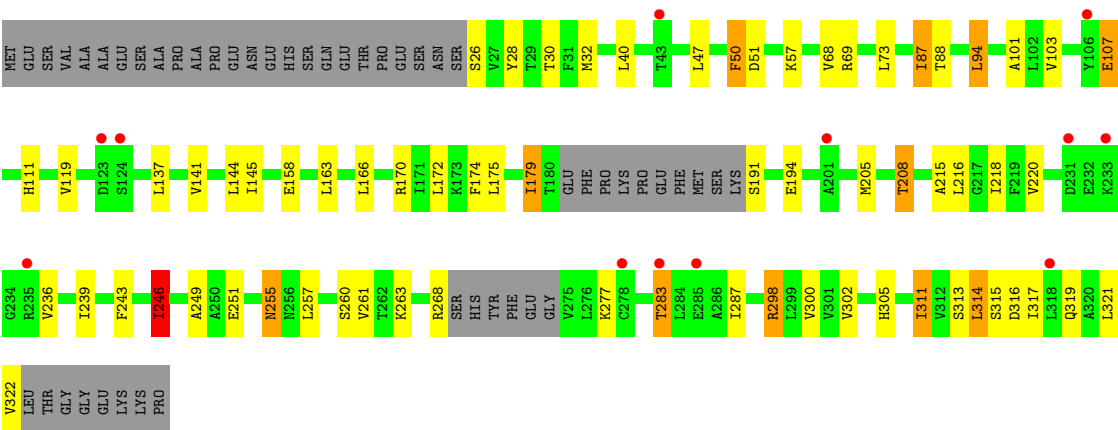
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
7	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 8 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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- Diagram illustrating the Lewis structure of the sulfate ion (SO_4^{2-}). The central sulfur atom (S) is bonded to four oxygen atoms (O). Two oxygen atoms are double-bonded to sulfur (top and bottom), and two are single-bonded (left and right). Each single-bonded oxygen has a negative charge (O^-). The sulfur atom has a formal charge of +6, indicated by a green 'S' with a '+' sign. The oxygen atoms are labeled O1, O2, O3, and O4 in green. The overall charge is -2, indicated by a green '2-'.

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



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	123.61Å 123.61Å 402.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.01 – 3.40 30.01 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (30.01-3.40) 99.7 (30.01-3.40)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 3.39Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.216 , 0.245 (Not available) , 0.234	Depositor DCC
R_{free} test set	1318 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	88.4	Xtriage
Anisotropy	0.574	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 117.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6369	wwPDB-VP
Average B, all atoms (Å ²)	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, SO4, CL, TPO, SEP, STU, AMP, HTV

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	0.90	0/2933	1.46	28/3968 (0.7%)
2	B	0.92	1/1231 (0.1%)	1.37	12/1689 (0.7%)
3	C	0.89	0/2167	1.48	22/2967 (0.7%)
All	All	0.90	1/6331 (0.0%)	1.45	62/8624 (0.7%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	258	LYS	CA-C	5.34	1.57	1.53

The worst 5 of 62 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	55	LEU	N-CA-C	-7.59	99.04	110.28
1	A	139	ASP	CA-CB-CG	7.11	119.70	112.60
1	A	456	SER	N-CA-C	-7.08	104.61	113.18
3	C	208	THR	CA-C-N	6.86	130.74	120.31
3	C	208	THR	C-N-CA	6.86	130.74	120.31

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	2880	0	2842	36	0
2	B	1208	0	1143	20	0
3	C	2127	0	2080	20	0
4	A	36	0	0	0	0
5	A	35	0	26	7	0
6	A	4	0	0	0	0
6	B	1	0	0	0	0
7	C	46	0	24	1	0
8	C	27	0	12	0	0
9	C	5	0	0	0	0
All	All	6369	0	6127	77	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

The worst 5 of 77 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)
5:A:602:STU:H283	5:A:602:STU:H272	1.27	1.08
5:A:602:STU:H16	5:A:602:STU:H261	1.51	0.90
5:A:602:STU:H272	5:A:602:STU:C28	2.02	0.90
5:A:602:STU:H283	5:A:602:STU:C27	2.05	0.85
1:A:218:HIS:HD2	1:A:221:THR:H	1.27	0.82

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	355/503 (71%)	326 (92%)	28 (8%)	1 (0%)	36 65
2	B	152/204 (74%)	138 (91%)	13 (9%)	1 (1%)	18 47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	275/330 (83%)	258 (94%)	17 (6%)	0	100	100
All	All	782/1037 (75%)	722 (92%)	58 (7%)	2 (0%)	36	65

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	120	GLU
1	A	427	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	101/448 (22%)	88 (87%)	13 (13%)	4	16
2	B	121/184 (66%)	107 (88%)	14 (12%)	5	20
3	C	226/299 (76%)	191 (84%)	35 (16%)	2	10
All	All	448/931 (48%)	386 (86%)	62 (14%)	3	14

5 of 62 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	47	LEU
3	C	302	VAL
3	C	111	HIS
3	C	298	ARG
3	C	316	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	66	ASN
3	C	92	ASN
2	B	110	ASN

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Mol	Chain	Res	Type
2	B	124	GLN
2	B	150	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
2	SEP	B	108	2	8,9,10	1.45	2 (25%)	7,12,14	4.25	3 (42%)
1	TPO	A	172	1	8,10,11	1.71	1 (12%)	10,14,16	2.16	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
2	SEP	B	108	2	-	6/6/8/10	-
1	TPO	A	172	1	-	1/9/11/13	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	172	TPO	P-OG1	-4.54	1.51	1.59
2	B	108	SEP	CB-CA	2.22	1.58	1.52
2	B	108	SEP	P-O1P	2.07	1.56	1.50

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	108	SEP	OG-CB-CA	9.39	117.28	108.14
2	B	108	SEP	O3P-P-OG	4.78	119.13	106.67
1	A	172	TPO	O3P-P-O2P	3.73	121.78	107.80
1	A	172	TPO	P-OG1-CB	-3.64	113.42	123.33
2	B	108	SEP	O3P-P-O2P	-3.05	96.38	107.80

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	172	TPO	O-C-CA-CB
2	B	108	SEP	N-CA-CB-OG
2	B	108	SEP	C-CA-CB-OG
2	B	108	SEP	CB-OG-P-O2P
2	B	108	SEP	CB-OG-P-O3P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 5 are monoatomic - leaving 6 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	$\# Z > 2$
4	HTV	A	601	-	38,40,40	1.56	7 (18%)	54,61,61	1.70	11 (20%)
5	STU	A	602	-	39,42,42	2.56	15 (38%)	50,68,68	2.87	19 (38%)
9	SO4	C	404	-	4,4,4	0.27	0	6,6,6	0.18	0
8	ADP	C	403	-	28,29,29	0.45	0	43,45,45	0.68	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	AMP	C	402	-	25,25,25	0.19	0	37,38,38	0.38	0
7	AMP	C	401	-	25,25,25	0.28	0	37,38,38	0.89	1 (2%)

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	HTV	A	601	-	-	2/22/50/50	0/5/5/5
5	STU	A	602	-	-	1/4/42/42	-
8	ADP	C	403	-	-	3/16/32/32	0/3/3/3
7	AMP	C	402	-	-	3/10/26/26	0/3/3/3
7	AMP	C	401	-	-	1/10/26/26	0/3/3/3

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	602	STU	C9-C10	-5.28	1.45	1.50
5	A	602	STU	C5-C20	5.20	1.49	1.41
4	A	601	HTV	O5-C21	5.13	1.37	1.22
5	A	602	STU	C12-C17	5.08	1.48	1.41
5	A	602	STU	C11-C10	4.91	1.49	1.40

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	602	STU	C11-C10-C7	-7.06	116.60	121.25
5	A	602	STU	C7-C8-N1	5.76	111.88	106.38
4	A	601	HTV	C11-C12-C7	5.76	126.44	121.97
5	A	602	STU	C6-C7-C10	5.73	124.75	120.76
5	A	602	STU	C9-C10-C11	5.51	132.91	128.73

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	601	HTV	C19-C20-C21-O4
4	A	601	HTV	C19-C20-C21-O5
7	C	402	AMP	C5'-O5'-P-O2P
7	C	402	AMP	C5'-O5'-P-O3P

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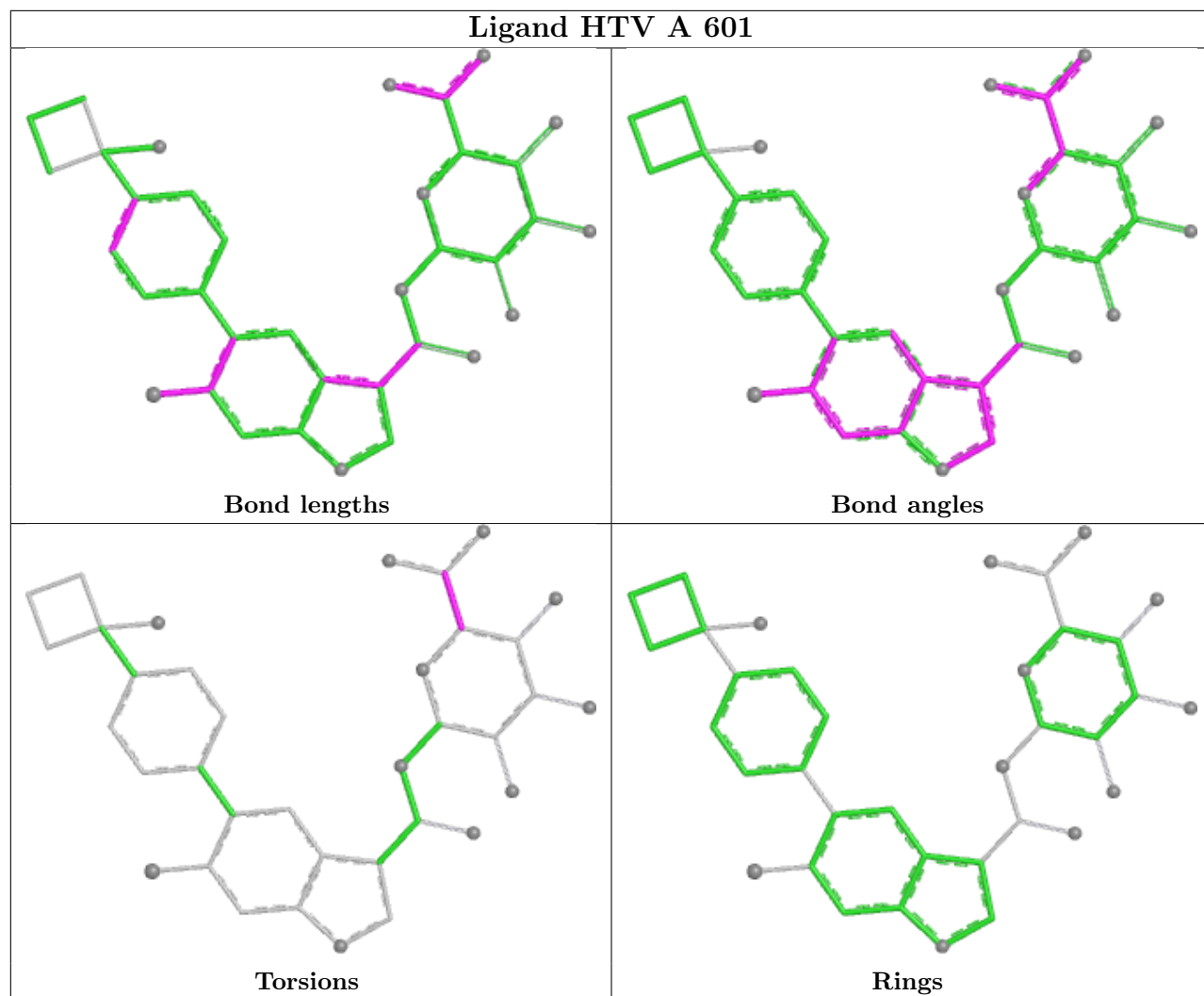
Mol	Chain	Res	Type	Atoms
8	C	403	ADP	C5'-O5'-PA-O1A

There are no ring outliers.

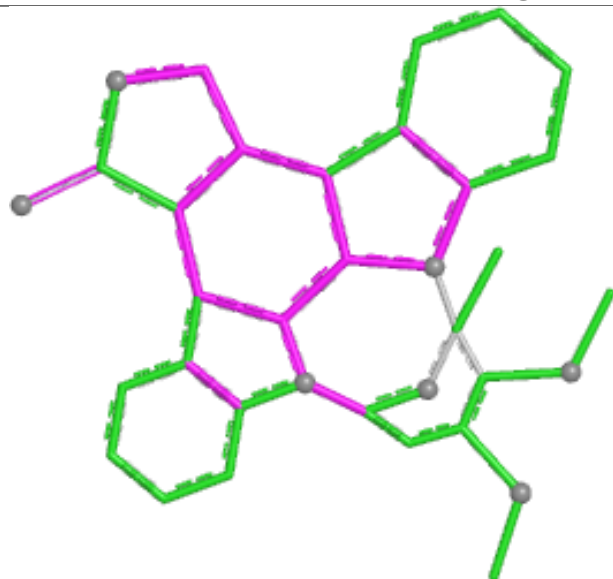
2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	602	STU	7	0
7	C	401	AMP	1	0

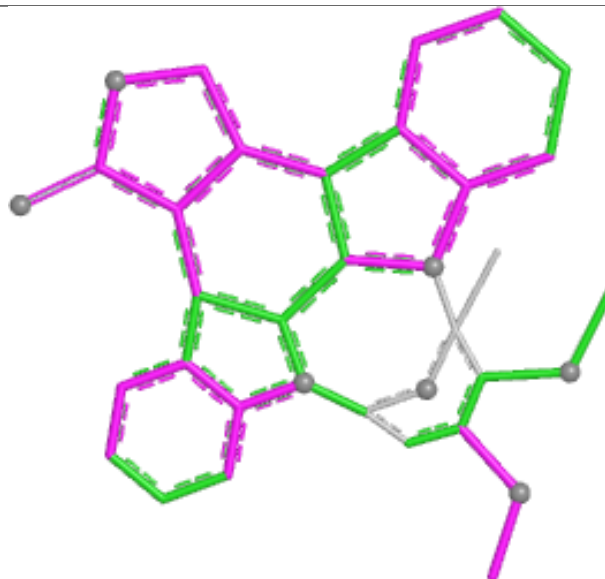
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



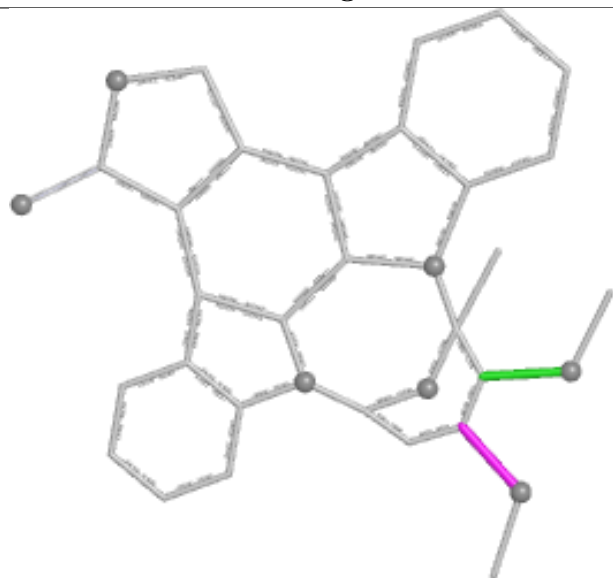
Ligand STU A 602



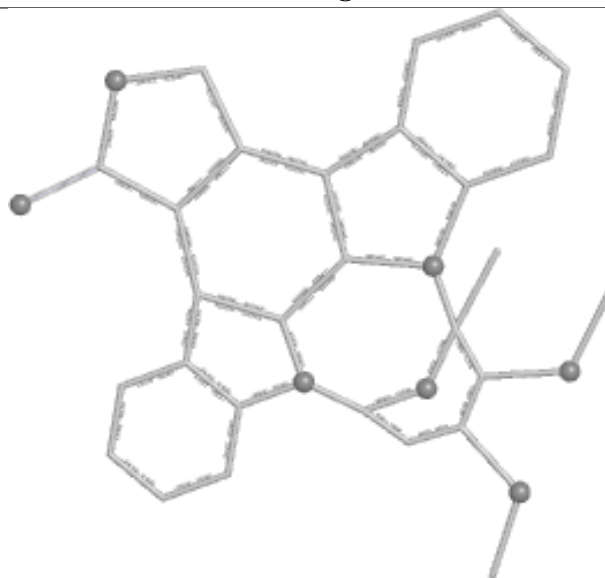
Bond lengths



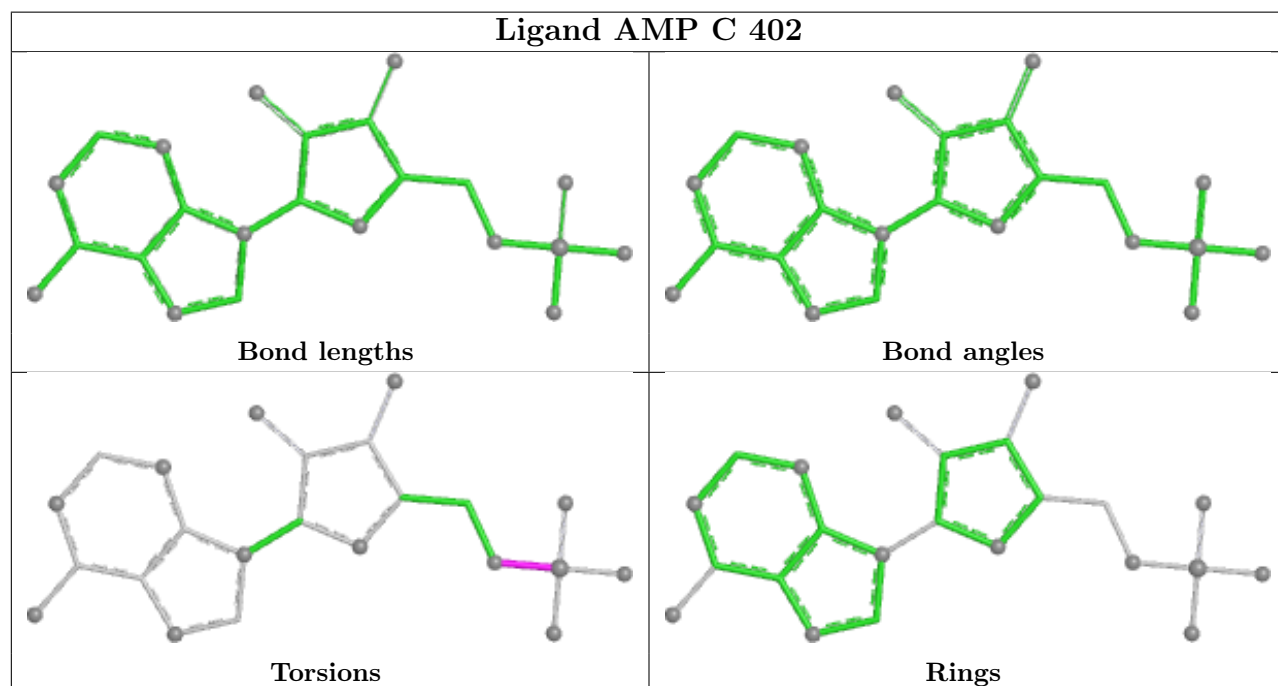
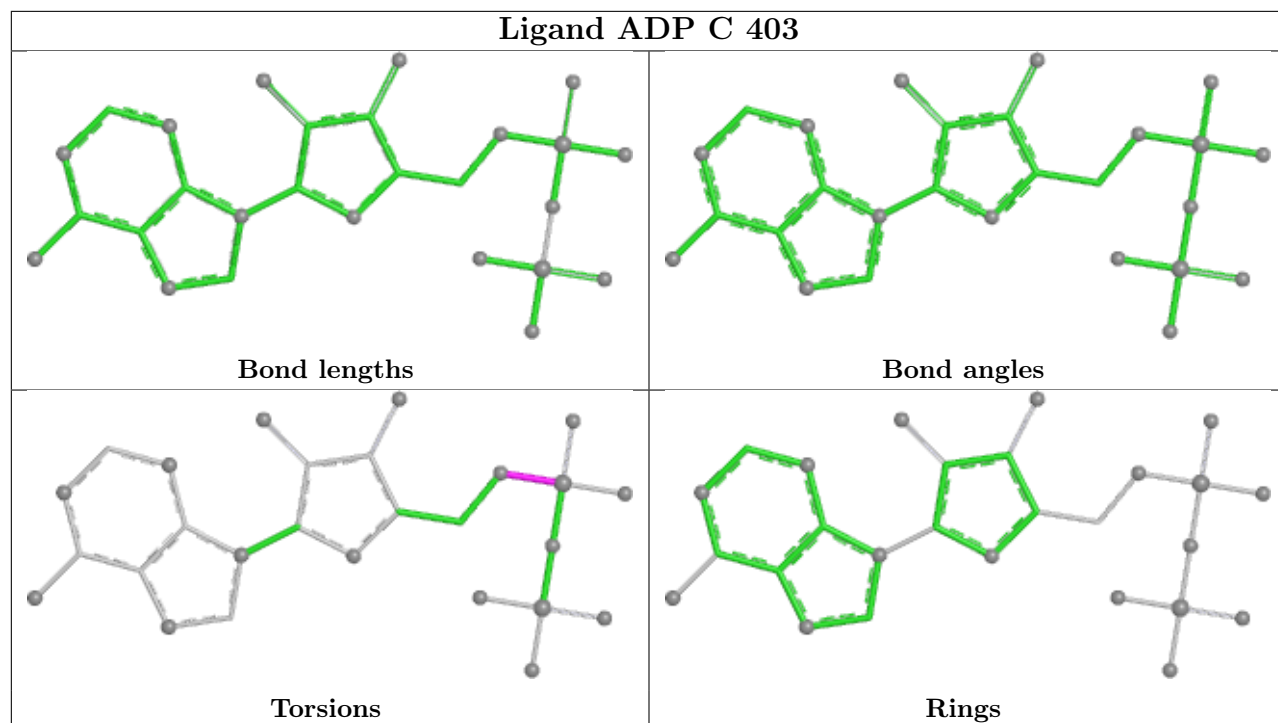
Bond angles

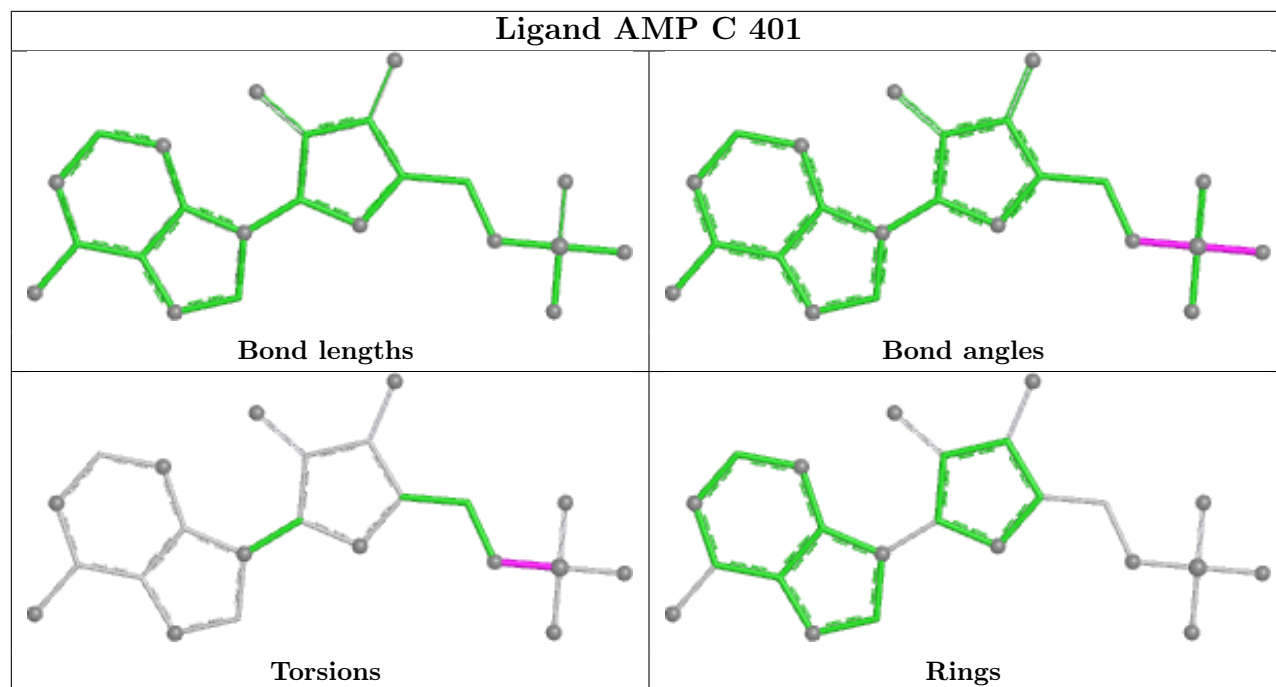


Torsions



Rings





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	361/503 (71%)	-0.16	4 (1%) 78 65	56, 89, 174, 196	0
2	B	158/204 (77%)	0.11	7 (4%) 39 28	71, 103, 140, 162	0
3	C	281/330 (85%)	0.26	12 (4%) 40 29	83, 139, 188, 205	0
All	All	800/1037 (77%)	0.04	23 (2%) 53 39	56, 107, 181, 205	0

The worst 5 of 23 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	231	ASP	4.4
1	A	399	GLY	3.5
3	C	124	SER	3.5
1	A	547	ALA	3.3
1	A	70	LEU	2.9

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SEP	B	108	10/11	0.93	0.08	98,105,107,109	0
1	TPO	A	172	11/12	0.96	0.07	91,92,98,101	0

6.3 Carbohydrates [i](#)

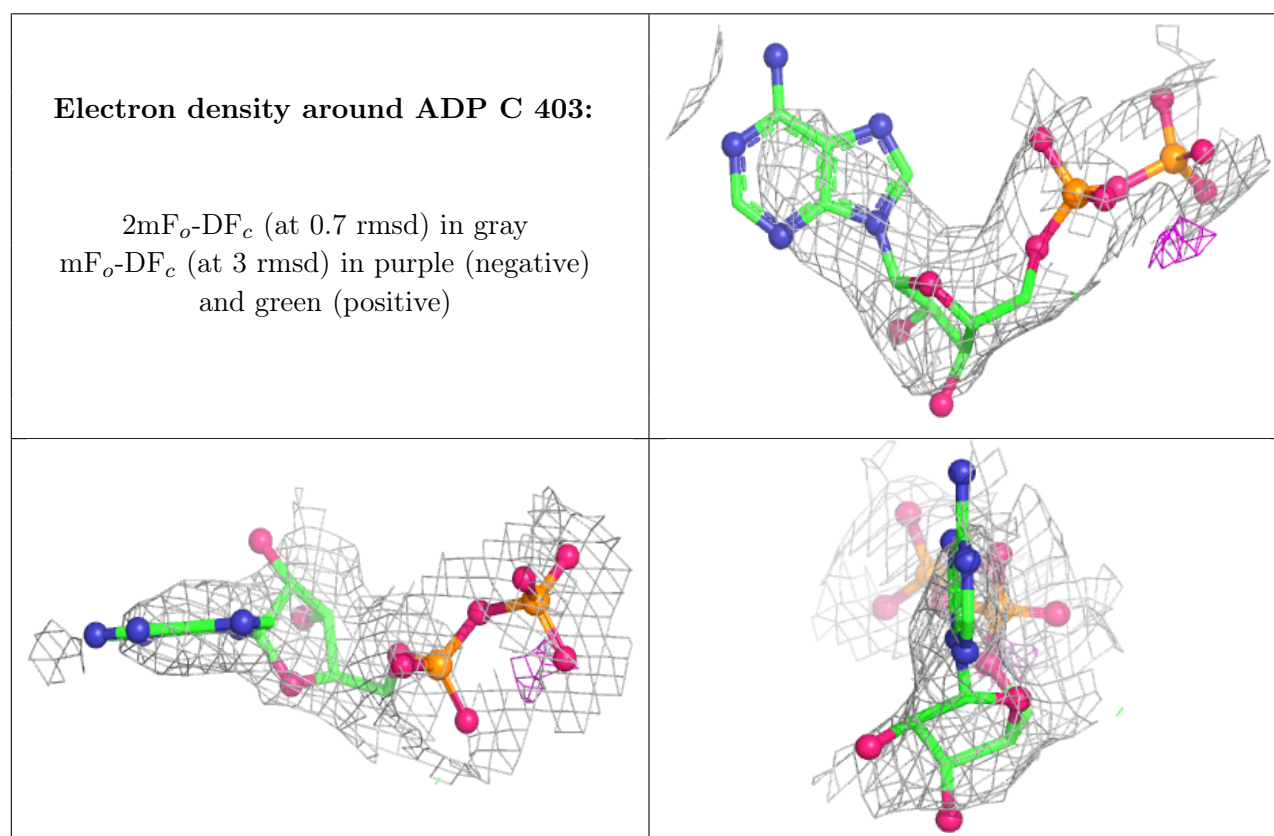
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

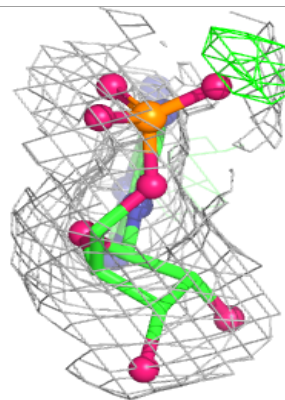
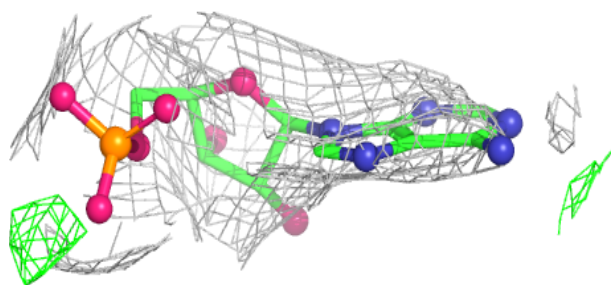
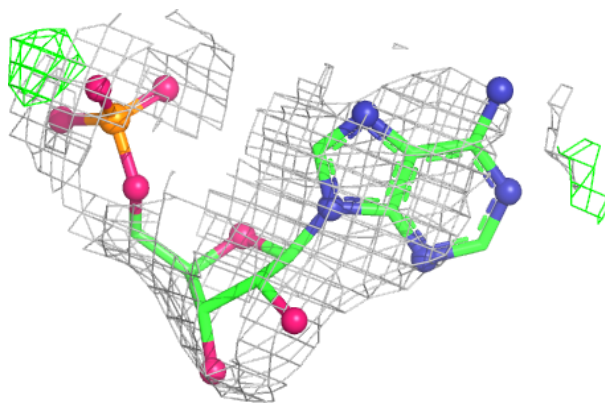
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
9	SO4	C	404	5/5	0.82	0.12	154,155,156,156	0
6	CL	A	606	1/1	0.84	0.12	117,117,117,117	0
8	ADP	C	403	27/27	0.88	0.12	183,193,200,202	0
7	AMP	C	401	23/23	0.92	0.10	159,172,185,187	0
7	AMP	C	402	23/23	0.92	0.11	148,165,171,176	0
4	HTV	A	601	36/36	0.94	0.10	82,96,144,145	0
6	CL	B	301	1/1	0.94	0.10	97,97,97,97	0
5	STU	A	602	35/35	0.97	0.09	63,68,72,75	0
6	CL	A	603	1/1	0.98	0.08	65,65,65,65	0
6	CL	A	604	1/1	0.98	0.06	57,57,57,57	0
6	CL	A	605	1/1	0.98	0.06	85,85,85,85	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.



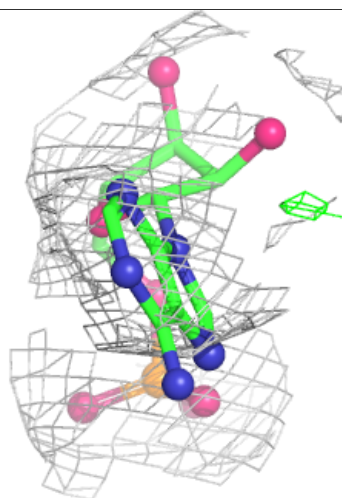
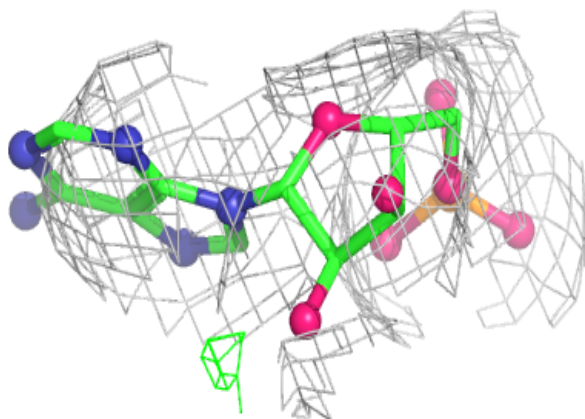
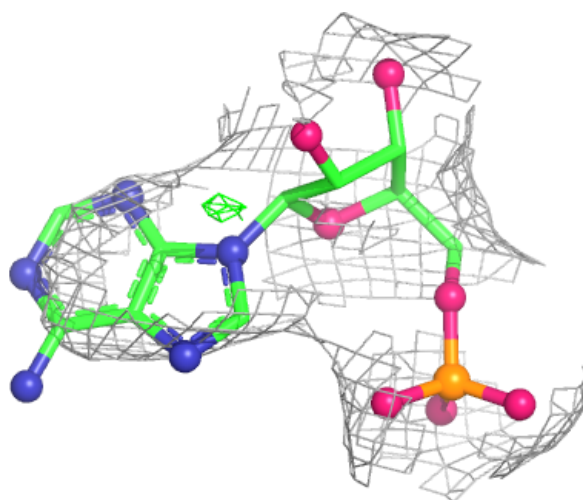
Electron density around AMP C 401:

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



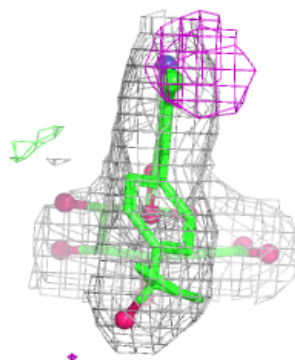
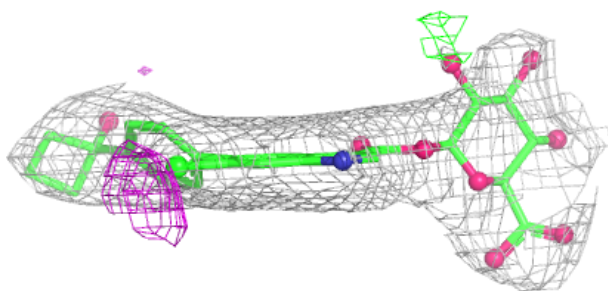
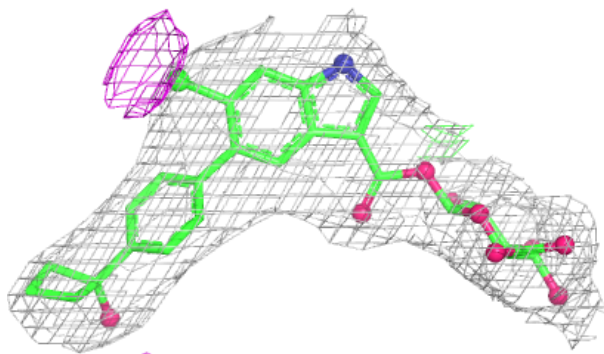
Electron density around AMP C 402:

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



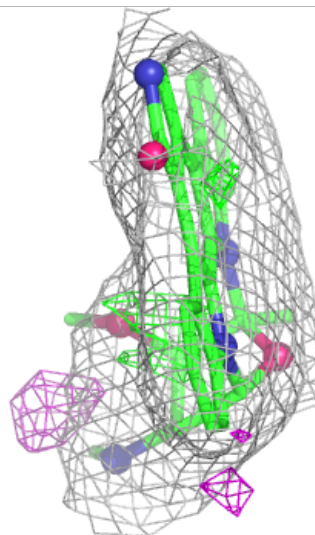
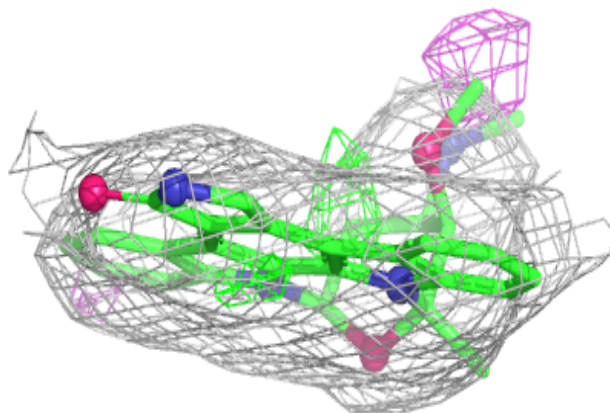
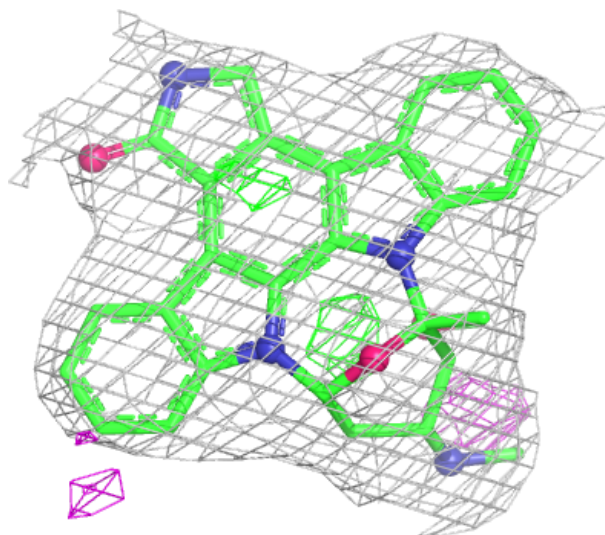
Electron density around HTV A 601:

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



Electron density around STU A 602:

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.