



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 03:27 am BST

PDB ID : 2P2F
Title : Acetyl-CoA Synthetase, wild-type with acetate, AMP, and CoA bound
Authors : Reger, A.S.; Gulick, A.M.
Deposited on : 2007-03-07
Resolution : 2.58 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 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

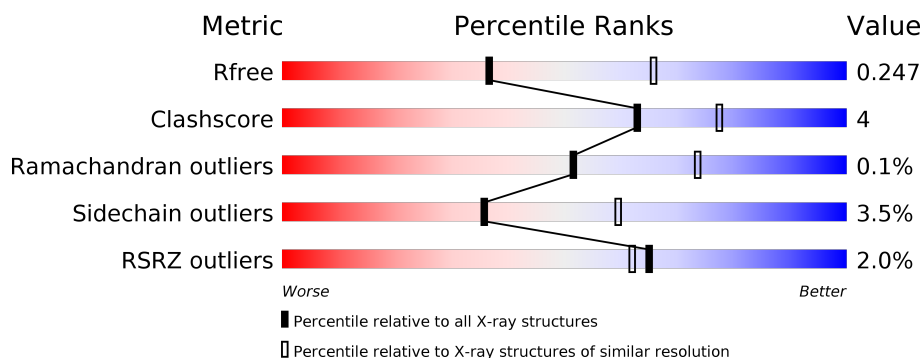
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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 ≥ 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	652	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>10%</div> <div>••</div> </div> </div>
1	B	652	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>9%</div> <div>••</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	653	-	-	X	-
2	ACT	B	653	-	-	X	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10137 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 Acetyl-coenzyme A synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	637	Total	C	N	O	S	0	0	0
			4926	3137	839	932	18			
1	B	632	Total	C	N	O	S	0	0	0
			4902	3120	842	922	18			

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



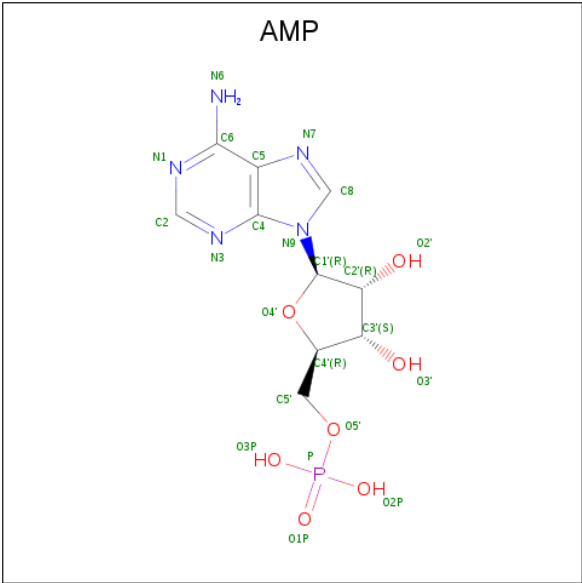
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is COENZYME A (three-letter code: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
3	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 4 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C₁₀H₁₄N₅O₇P).



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

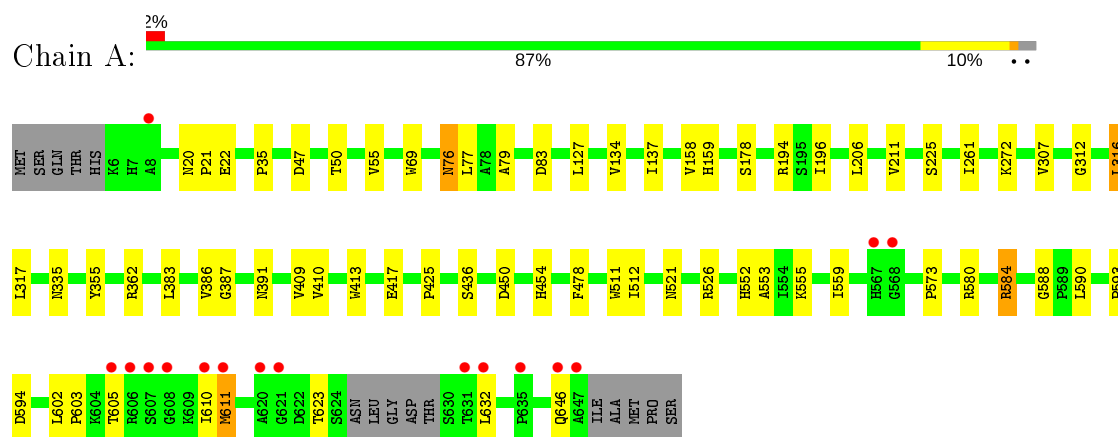
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	86	Total 86	O 86	0	0
5	B	73	Total 73	O 73	0	0

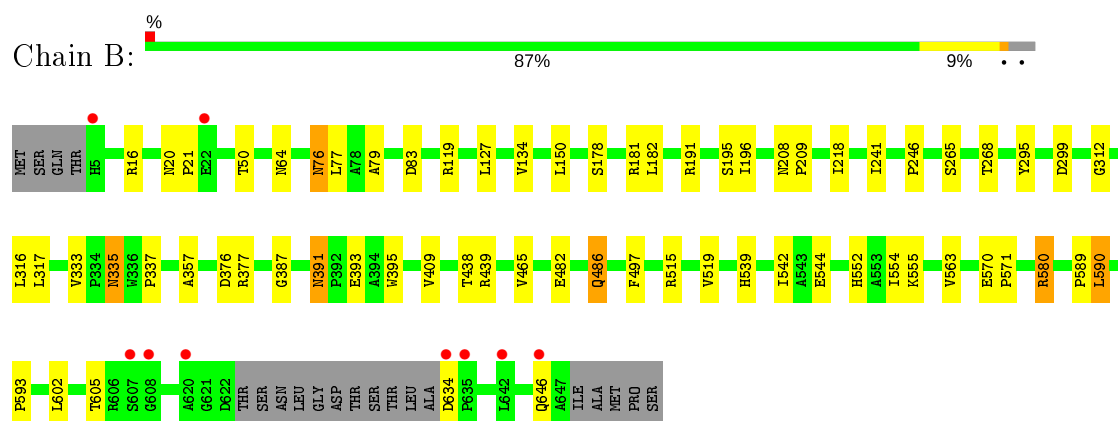
3 Residue-property plots [i](#)

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: Acetyl-coenzyme A synthetase



• Molecule 1: Acetyl-coenzyme A synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.91Å 94.93Å 161.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	81.65 – 2.58 24.73 – 2.58	Depositor EDS
% Data completeness (in resolution range)	99.0 (81.65-2.58) 99.1 (24.73-2.58)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.44 (at 2.57Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.183 , 0.247 0.185 , 0.247	Depositor DCC
R_{free} test set	2261 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.020 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10137	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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: COA, AMP, ACT

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	0.69	1/5057 (0.0%)	0.74	2/6902 (0.0%)
1	B	0.67	0/5034	0.72	0/6866
All	All	0.68	1/10091 (0.0%)	0.73	2/13768 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	22	GLU	CG-CD	5.52	1.60	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	ASP	CB-CG-OD2	5.38	123.14	118.30
1	A	590	LEU	CA-CB-CG	5.00	126.80	115.30

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	4926	0	4731	38	0
1	B	4902	0	4727	38	0
2	A	4	0	3	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	4	0	3	5	0
3	A	48	0	32	4	0
3	B	48	0	32	5	0
4	A	23	0	12	1	0
4	B	23	0	12	1	0
5	A	86	0	0	1	0
5	B	73	0	0	5	0
All	All	10137	0	9552	82	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 4.

All (82) 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:B:387:GLY:HA2	2:B:653:ACT:H3	1.45	0.97
1:B:387:GLY:CA	2:B:653:ACT:H3	2.10	0.80
2:B:653:ACT:H2	5:B:1064:HOH:O	1.86	0.75
2:A:653:ACT:H2	3:A:990:COA:S1P	2.27	0.74
1:B:191:ARG:NH1	3:B:991:COA:O4A	2.27	0.68
1:A:611:MET:CE	1:A:611:MET:HA	2.24	0.68
1:A:611:MET:HA	1:A:611:MET:HE3	1.79	0.64
3:B:991:COA:O2A	3:B:991:COA:OAP	2.10	0.64
1:B:552:HIS:HD2	1:B:555:LYS:H	1.48	0.61
1:B:357:ALA:HB2	3:B:991:COA:H32	1.83	0.60
1:A:76:ASN:HD22	1:A:79:ALA:H	1.51	0.59
1:B:580:ARG:HD2	1:B:593:PRO:O	2.03	0.58
1:B:552:HIS:CD2	1:B:555:LYS:H	2.22	0.57
1:A:410:VAL:HA	1:A:425:PRO:HG2	1.85	0.57
1:B:134:VAL:HG11	1:B:178:SER:HB2	1.88	0.55
1:A:134:VAL:HG11	1:A:178:SER:HB2	1.89	0.55
1:B:119:ARG:HG2	1:B:246:PRO:HA	1.89	0.54
1:A:20:ASN:HB2	1:A:21:PRO:CD	2.38	0.53
1:A:552:HIS:HD2	1:A:555:LYS:H	1.55	0.53
1:A:387:GLY:HA2	2:A:653:ACT:H1	1.92	0.52
1:B:539:HIS:HB3	1:B:542:ILE:HD12	1.91	0.51
1:A:194:ARG:HD3	1:B:337:PRO:O	2.11	0.50
1:B:465:VAL:HG12	1:B:497:PHE:HA	1.94	0.50
1:A:20:ASN:HB2	1:A:21:PRO:HD2	1.93	0.49
2:B:653:ACT:C	4:B:998:AMP:O3P	2.60	0.49
1:B:552:HIS:HD2	1:B:554:ILE:H	1.59	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:990:COA:O2A	3:A:990:COA:O6A	2.31	0.49
1:B:265:SER:HA	5:B:1020:HOH:O	2.12	0.49
1:A:552:HIS:CD2	1:A:555:LYS:H	2.30	0.48
1:B:552:HIS:CD2	1:B:554:ILE:H	2.31	0.48
1:B:515:ARG:HG3	1:B:519:VAL:HG12	1.93	0.48
1:B:20:ASN:HB2	1:B:21:PRO:CD	2.43	0.48
1:A:137:ILE:HD12	1:A:159:HIS:HB2	1.94	0.48
3:A:990:COA:OAP	3:A:990:COA:O2A	2.20	0.48
1:B:312:GLY:O	1:B:316:LEU:HB3	2.14	0.47
1:A:584:ARG:O	1:A:588:GLY:HA2	2.15	0.47
1:A:386:VAL:HG22	1:A:387:GLY:N	2.30	0.47
1:A:355:TYR:CE1	1:A:386:VAL:HB	2.50	0.47
1:A:611:MET:CE	1:A:611:MET:CA	2.94	0.45
1:B:76:ASN:HD22	1:B:79:ALA:H	1.63	0.45
1:B:191:ARG:CZ	1:B:333:VAL:HG12	2.46	0.45
2:B:653:ACT:CH3	5:B:1064:HOH:O	2.57	0.45
1:A:450:ASP:OD2	1:A:454:HIS:HB2	2.17	0.45
1:A:196:ILE:HG21	3:A:990:COA:H2A	1.97	0.45
1:B:376:ASP:O	1:B:377:ARG:HB2	2.17	0.45
1:A:312:GLY:O	1:A:316:LEU:HB3	2.16	0.45
2:A:653:ACT:C	4:A:999:AMP:O3P	2.65	0.44
1:B:150:LEU:HD23	1:B:150:LEU:HA	1.85	0.44
1:B:182:LEU:HD11	1:B:218:ILE:HG13	2.00	0.44
1:A:383:LEU:HB2	1:A:409:VAL:HG22	1.99	0.44
1:A:362:ARG:HD2	1:A:553:ALA:O	2.18	0.44
1:B:391:ASN:HB3	5:B:1065:HOH:O	2.17	0.44
1:B:395:TRP:CZ2	1:B:409:VAL:HB	2.52	0.43
1:A:134:VAL:HA	1:A:158:VAL:O	2.18	0.43
1:A:316:LEU:HD12	1:A:317:LEU:HD12	2.00	0.43
1:B:335:ASN:H	1:B:335:ASN:HD22	1.67	0.43
1:A:206:LEU:HD22	1:A:211:VAL:HG11	2.00	0.43
1:A:261:ILE:HD12	1:A:478:PHE:HB2	2.00	0.43
1:B:208:ASN:HA	1:B:209:PRO:HD2	1.89	0.43
1:A:413:TRP:CD1	1:A:512:ILE:HG12	2.53	0.43
1:A:573:PRO:HG2	1:B:393:GLU:HB3	2.01	0.43
1:A:580:ARG:HD2	1:A:593:PRO:O	2.19	0.43
1:B:439:ARG:HG2	1:B:439:ARG:HH11	1.84	0.42
1:B:589:PRO:HB2	3:B:991:COA:H4B	2.00	0.42
1:A:555:LYS:NZ	1:A:594:ASP:OD1	2.52	0.42
1:A:602:LEU:HA	1:A:603:PRO:HD3	1.92	0.42
1:A:272:LYS:HA	1:A:272:LYS:HD2	1.78	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:417:GLU:OE2	1:A:526:ARG:NH2	2.47	0.42
1:A:436:SER:HB2	1:A:511:TRP:CE2	2.55	0.41
1:B:196:ILE:HG21	3:B:991:COA:H2A	2.02	0.41
1:B:482:GLU:O	1:B:486:GLN:HB2	2.20	0.41
1:B:570:GLU:HA	1:B:571:PRO:HD3	1.97	0.41
1:B:590:LEU:HD22	5:B:1052:HOH:O	2.20	0.41
1:A:559:ILE:HB	1:A:593:PRO:HA	2.02	0.41
1:A:77:LEU:C	1:A:77:LEU:HD23	2.41	0.41
1:A:35:PRO:HD2	5:A:1004:HOH:O	2.21	0.41
1:B:542:ILE:HG12	1:B:563:VAL:CG1	2.51	0.41
1:A:355:TYR:CZ	1:A:386:VAL:HB	2.56	0.40
1:A:55:VAL:O	1:A:69:TRP:HA	2.21	0.40
1:B:295:TYR:CE1	1:B:299:ASP:HB2	2.56	0.40
1:B:77:LEU:HD23	1:B:77:LEU:C	2.41	0.40
1:B:544:GLU:HG3	1:B:602:LEU:HD12	2.04	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	633/652 (97%)	614 (97%)	19 (3%)	0	100	100
1	B	628/652 (96%)	607 (97%)	20 (3%)	1 (0%)	47	69
All	All	1261/1304 (97%)	1221 (97%)	39 (3%)	1 (0%)	51	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	438	THR

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	507/538 (94%)	490 (97%)	17 (3%)	37	60
1	B	508/538 (94%)	489 (96%)	19 (4%)	34	57
All	All	1015/1076 (94%)	979 (96%)	36 (4%)	36	59

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	THR
1	A	76	ASN
1	A	83	ASP
1	A	127	LEU
1	A	225	SER
1	A	307	VAL
1	A	316	LEU
1	A	335	ASN
1	A	391	ASN
1	A	521	ASN
1	A	584	ARG
1	A	605	THR
1	A	610	ILE
1	A	611	MET
1	A	623	THR
1	A	632	LEU
1	A	646	GLN
1	B	16	ARG
1	B	50	THR
1	B	64	ASN
1	B	76	ASN
1	B	83	ASP
1	B	127	LEU
1	B	181	ARG
1	B	195	SER
1	B	241	ILE
1	B	268	THR

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Mol	Chain	Res	Type
1	B	317	LEU
1	B	335	ASN
1	B	391	ASN
1	B	486	GLN
1	B	580	ARG
1	B	590	LEU
1	B	605	THR
1	B	634	ASP
1	B	646	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	ASN
1	A	335	ASN
1	A	445	GLN
1	A	481	HIS
1	A	525	HIS
1	A	552	HIS
1	A	646	GLN
1	B	64	ASN
1	B	76	ASN
1	B	89	ASN
1	B	335	ASN
1	B	344	GLN
1	B	486	GLN
1	B	552	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 carbohydrates in this entry.

5.6 Ligand geometry

6 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
3	COA	B	991	-	41,50,50	1.82	3 (7%)	52,75,75	1.34	4 (7%)
4	AMP	B	998	-	22,25,25	1.26	2 (9%)	25,38,38	1.33	2 (8%)
2	ACT	A	653	4	1,3,3	1.94	0	0,3,3	0.00	-
4	AMP	A	999	2	22,25,25	1.19	3 (13%)	25,38,38	1.26	3 (12%)
3	COA	A	990	-	41,50,50	1.64	3 (7%)	52,75,75	1.50	4 (7%)
2	ACT	B	653	-	1,3,3	0.72	0	0,3,3	0.00	-

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
3	COA	B	991	-	-	13/44/64/64	0/3/3/3
4	AMP	A	999	2	-	0/6/26/26	0/3/3/3
3	COA	A	990	-	-	18/44/64/64	0/3/3/3
4	AMP	B	998	-	-	0/6/26/26	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	991	COA	O9P-C9P	9.43	1.42	1.23
3	A	990	COA	O9P-C9P	8.59	1.40	1.23
3	B	991	COA	C2A-N3A	4.16	1.38	1.32
3	A	990	COA	C2A-N3A	3.96	1.38	1.32
4	B	998	AMP	O4'-C1'	3.30	1.45	1.41
3	B	991	COA	C2A-N1A	2.99	1.39	1.33
4	A	999	AMP	C2-N3	2.79	1.36	1.32
4	B	998	AMP	C5-C4	2.52	1.47	1.40
3	A	990	COA	C2A-N1A	2.44	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	999	AMP	C5-C4	2.16	1.46	1.40
4	A	999	AMP	C2'-C1'	-2.08	1.50	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	990	COA	N3A-C2A-N1A	-6.27	118.88	128.68
3	B	991	COA	P2A-O3A-P1A	-4.85	116.19	132.83
3	B	991	COA	N3A-C2A-N1A	-4.77	121.23	128.68
3	A	990	COA	CEP-CBP-CAP	4.03	115.81	108.82
3	A	990	COA	P2A-O3A-P1A	-3.71	120.10	132.83
3	A	990	COA	CEP-CBP-CCP	-3.63	102.31	108.23
4	B	998	AMP	C4-C5-N7	-3.60	105.65	109.40
4	A	999	AMP	N3-C2-N1	-3.04	123.92	128.68
4	A	999	AMP	N6-C6-N1	2.93	124.67	118.57
3	B	991	COA	CEP-CBP-CAP	2.74	113.58	108.82
3	B	991	COA	CDP-CBP-CAP	2.70	113.50	108.82
4	A	999	AMP	C5-C6-N6	-2.37	116.75	120.35
4	B	998	AMP	N3-C2-N1	-2.19	125.26	128.68

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	991	COA	O4B-C4B-C5B-O5B
3	B	991	COA	CDP-CBP-CCP-O6A
3	B	991	COA	CEP-CBP-CCP-O6A
3	B	991	COA	CAP-CBP-CCP-O6A
3	B	991	COA	C9P-CAP-CBP-CCP
3	B	991	COA	C9P-CAP-CBP-CDP
3	B	991	COA	C9P-CAP-CBP-CEP
3	A	990	COA	O4B-C4B-C5B-O5B
3	A	990	COA	OAP-CAP-CBP-CCP
3	A	990	COA	OAP-CAP-CBP-CDP
3	A	990	COA	OAP-CAP-CBP-CEP
3	A	990	COA	O9P-C9P-CAP-CBP
3	A	990	COA	N8P-C9P-CAP-CBP
3	A	990	COA	C5P-C6P-C7P-N8P
3	A	990	COA	S1P-C2P-C3P-N4P
3	B	991	COA	C3B-C4B-C5B-O5B
3	A	990	COA	C3B-C4B-C5B-O5B
3	A	990	COA	C6P-C5P-N4P-C3P

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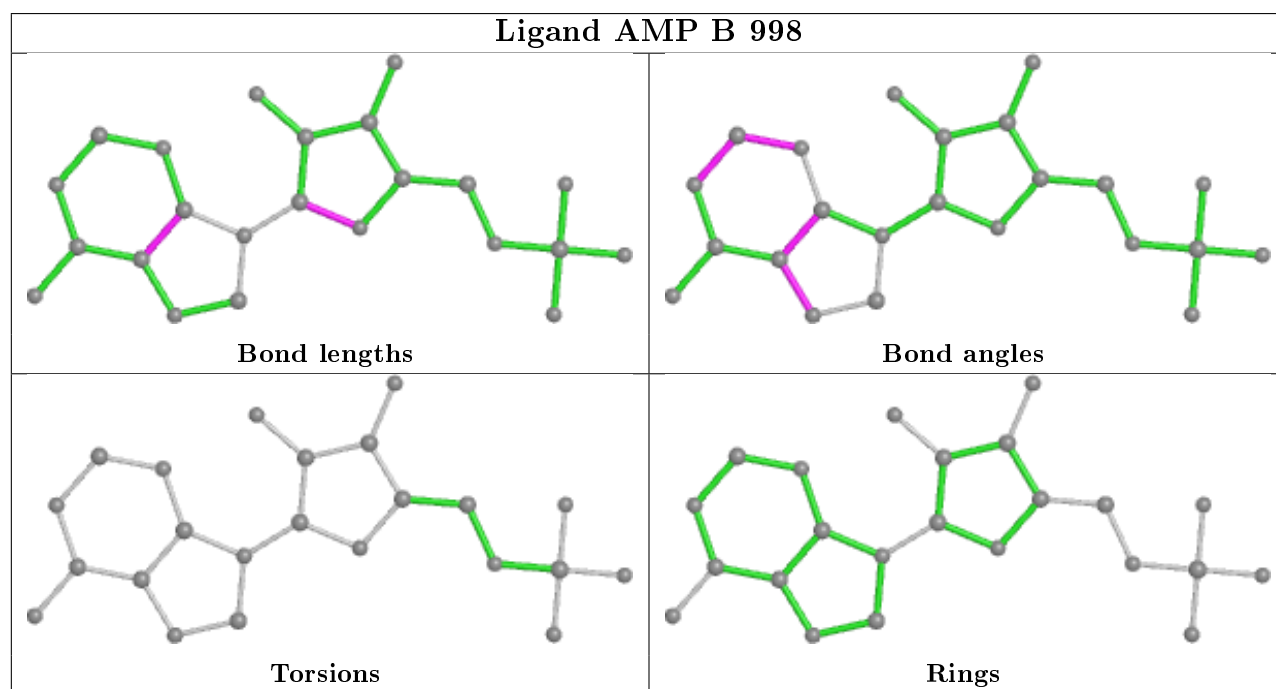
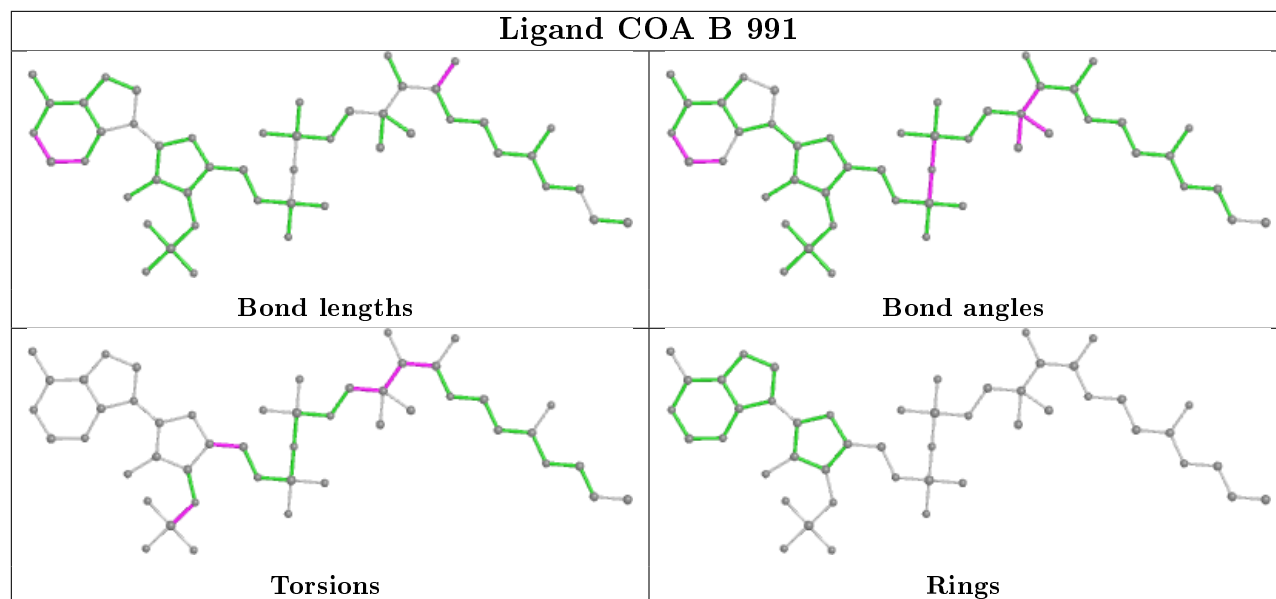
Mol	Chain	Res	Type	Atoms
3	A	990	COA	O5P-C5P-N4P-C3P
3	A	990	COA	P1A-O3A-P2A-O6A
3	A	990	COA	P2A-O3A-P1A-O1A
3	B	991	COA	O9P-C9P-CAP-CBP
3	B	991	COA	N8P-C9P-CAP-CBP
3	B	991	COA	C3B-O3B-P3B-O7A
3	A	990	COA	C2B-C3B-O3B-P3B
3	B	991	COA	C3B-O3B-P3B-O8A
3	B	991	COA	C3B-O3B-P3B-O9A
3	A	990	COA	CCP-O6A-P2A-O3A
3	A	990	COA	P2A-O3A-P1A-O2A
3	A	990	COA	C5B-O5B-P1A-O1A
3	A	990	COA	CCP-O6A-P2A-O4A

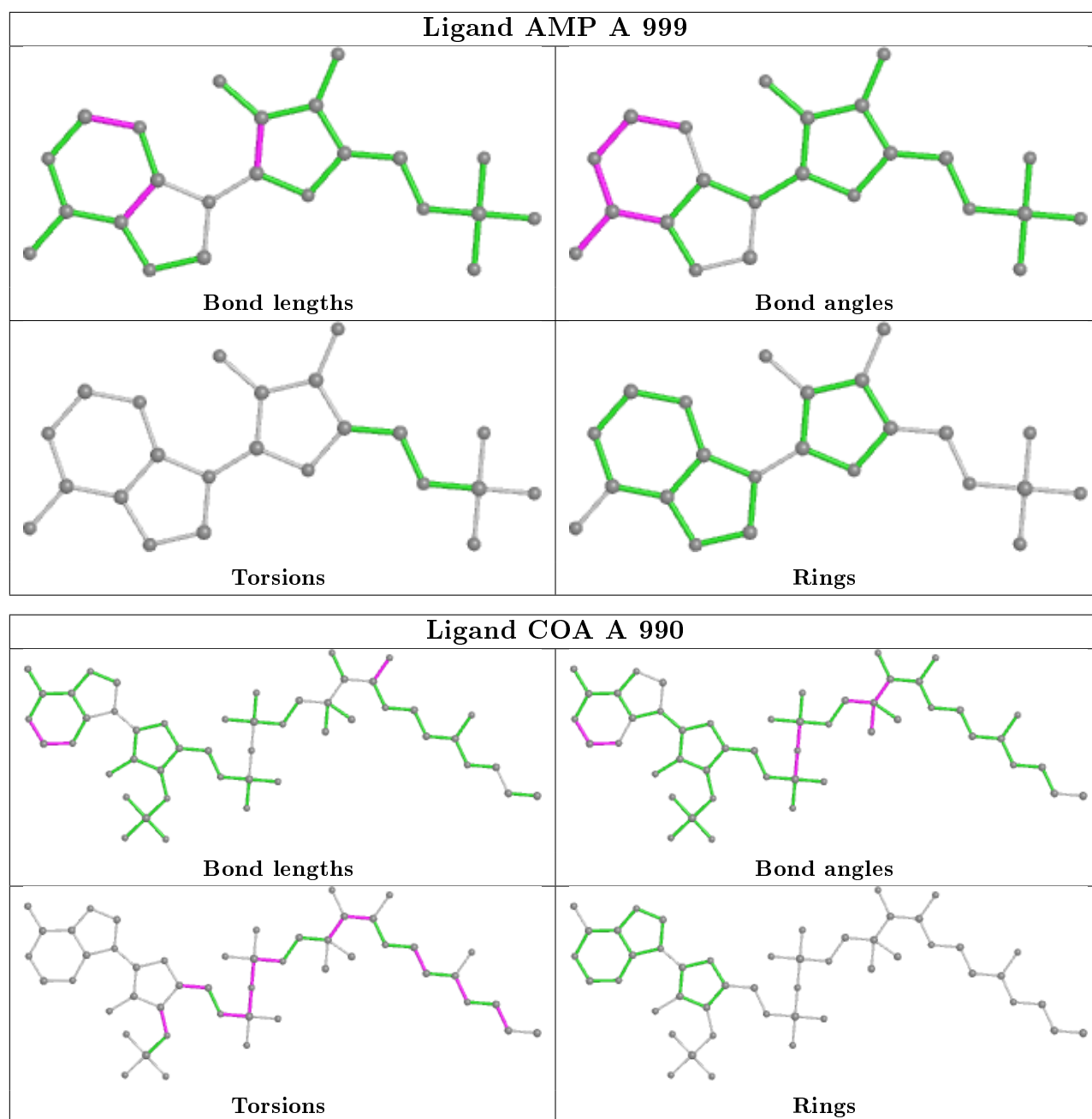
There are no ring outliers.

6 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	991	COA	5	0
4	B	998	AMP	1	0
2	A	653	ACT	3	0
4	A	999	AMP	1	0
3	A	990	COA	4	0
2	B	653	ACT	5	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, 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	637/652 (97%)	-0.26	16 (2%) 57 53	5, 16, 34, 53	0
1	B	632/652 (96%)	-0.29	9 (1%) 75 73	9, 17, 35, 53	0
All	All	1269/1304 (97%)	-0.27	25 (1%) 65 62	5, 16, 34, 53	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	635	PRO	4.9
1	B	5	HIS	4.4
1	A	608	GLY	4.1
1	A	607	SER	3.9
1	A	646	GLN	3.6
1	B	607	SER	3.5
1	B	646	GLN	3.4
1	A	567	HIS	3.4
1	A	568	GLY	3.0
1	A	632	LEU	3.0
1	B	620	ALA	2.9
1	A	606	ARG	2.8
1	A	631	THR	2.7
1	B	608	GLY	2.7
1	A	605	THR	2.6
1	B	22	GLU	2.6
1	A	611	MET	2.5
1	B	642	LEU	2.3
1	A	610	ILE	2.2
1	A	647	ALA	2.2
1	A	621	GLY	2.1
1	A	620	ALA	2.1
1	A	635	PRO	2.1
1	B	634	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	8	ALA	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 carbohydrates in this entry.

6.4 Ligands [i](#)

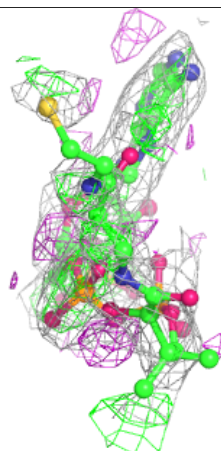
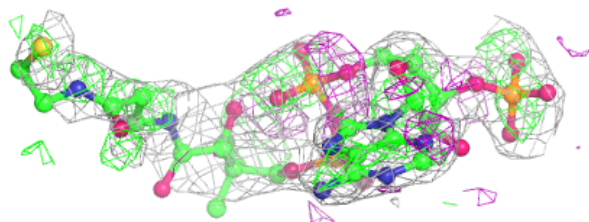
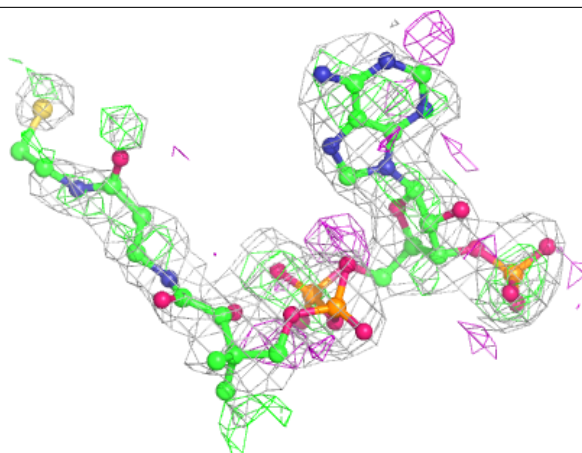
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	ACT	B	653	4/4	0.63	0.46	58,59,59,59	0
3	COA	B	991	48/48	0.77	0.28	22,39,58,58	0
3	COA	A	990	48/48	0.82	0.25	25,47,53,55	0
2	ACT	A	653	4/4	0.95	0.20	48,48,48,49	0
4	AMP	B	998	23/23	0.98	0.09	10,13,20,20	0
4	AMP	A	999	23/23	0.98	0.08	8,10,22,24	0

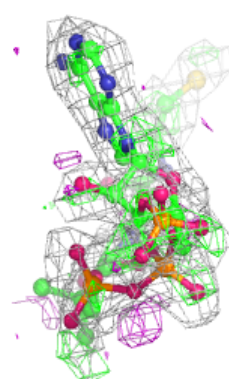
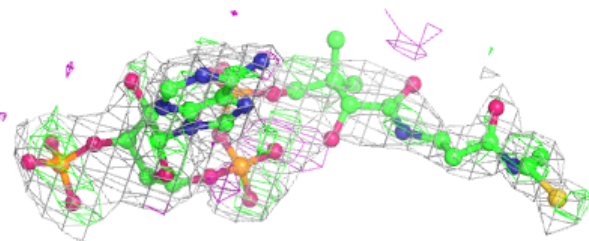
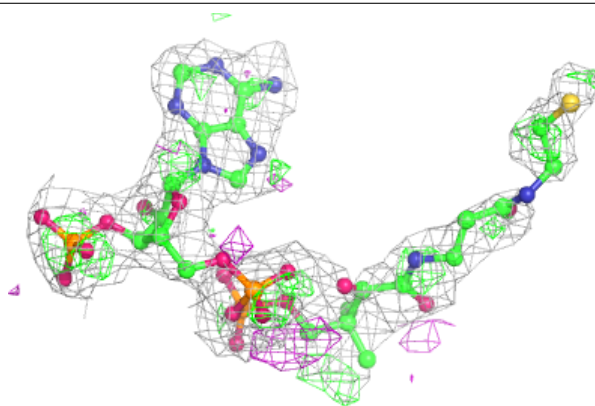
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around COA B 991:

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

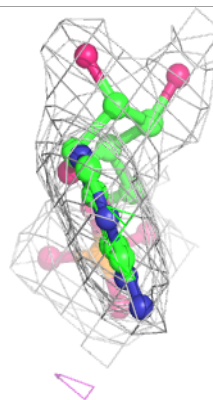
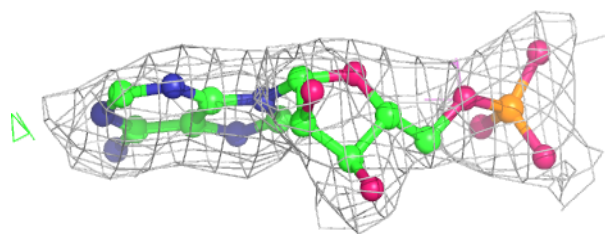
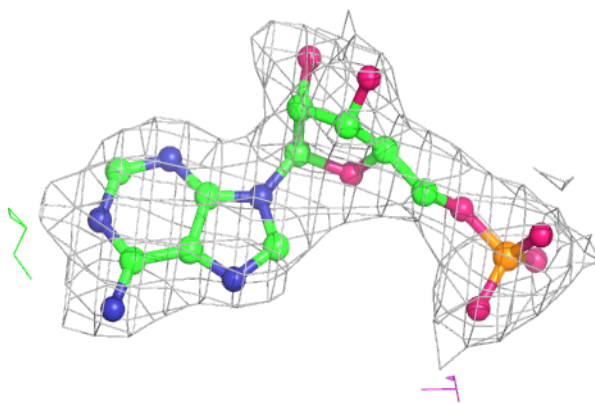
**Electron density around COA A 990:**

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

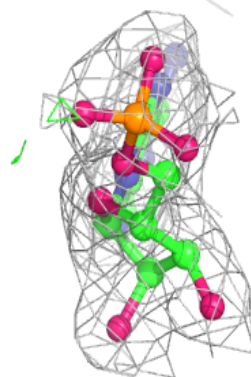
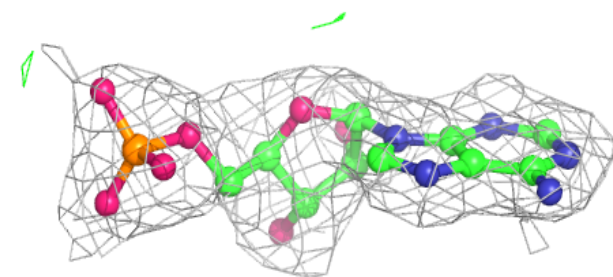
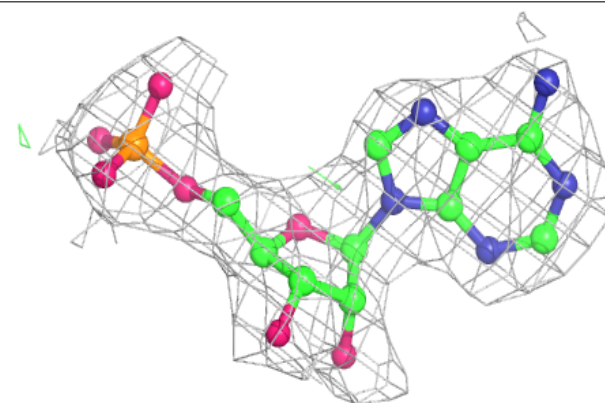


Electron density around AMP B 998:

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

**Electron density around AMP A 999:**

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



6.5 Other polymers

There are no such residues in this entry.