



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:12 AM UTC

PDB ID : 7WJV / pdb_00007wvjv
Title : Crystal structure of human liver FBPase complexed with an covalent inhibitor
Authors : Cao, H.; Huang, Y.; Ren, Y.; Wan, J.
Deposited on : 2022-01-08
Resolution : 1.72 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 ⓘ](#)) 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

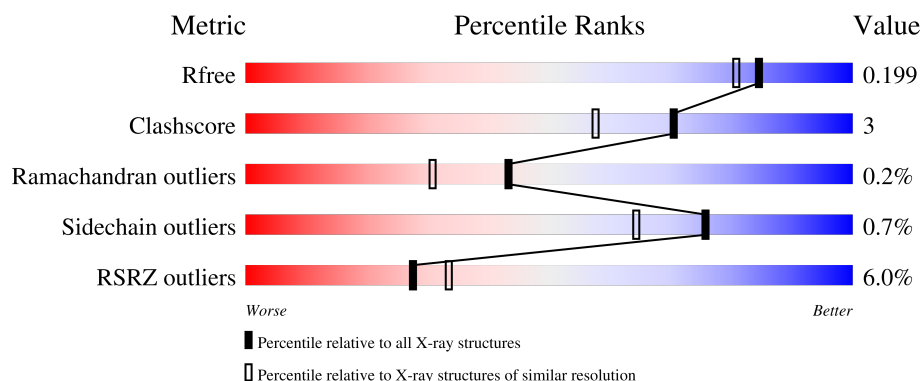
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.72 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	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	348	6% 84% 7% 7%
1	B	348	5% 89% 5% 6%
1	C	348	6% 86% 7% 6%
1	D	348	6% 84% 7% 8%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11179 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 Fructose-1,6-bisphosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	1	0
			2471	1572	416	466	17			
1	B	327	Total	C	N	O	S	0	0	0
			2485	1581	417	470	17			
1	C	326	Total	C	N	O	S	0	1	0
			2488	1583	419	469	17			
1	D	320	Total	C	N	O	S	0	0	0
			2443	1556	409	461	17			

There are 40 discrepancies between the modelled and reference sequences:

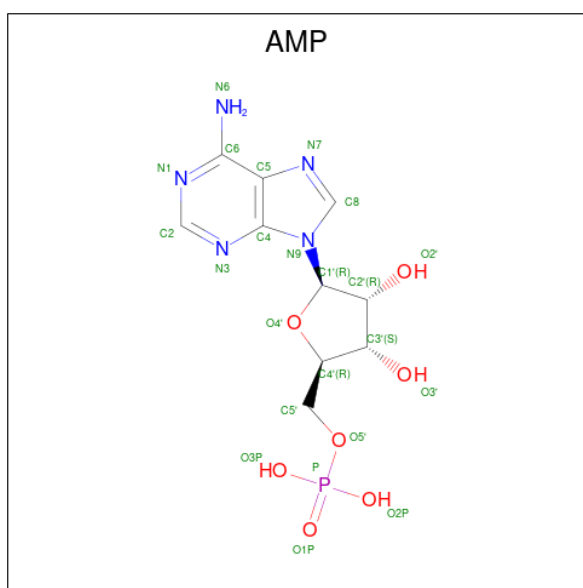
Chain	Residue	Modelled	Actual	Comment	Reference
A	338	LEU	-	expression tag	UNP Q2TU34
A	339	GLU	-	expression tag	UNP Q2TU34
A	340	HIS	-	expression tag	UNP Q2TU34
A	341	HIS	-	expression tag	UNP Q2TU34
A	342	HIS	-	expression tag	UNP Q2TU34
A	343	HIS	-	expression tag	UNP Q2TU34
A	344	HIS	-	expression tag	UNP Q2TU34
A	345	HIS	-	expression tag	UNP Q2TU34
A	346	HIS	-	expression tag	UNP Q2TU34
A	347	HIS	-	expression tag	UNP Q2TU34
B	338	LEU	-	expression tag	UNP Q2TU34
B	339	GLU	-	expression tag	UNP Q2TU34
B	340	HIS	-	expression tag	UNP Q2TU34
B	341	HIS	-	expression tag	UNP Q2TU34
B	342	HIS	-	expression tag	UNP Q2TU34
B	343	HIS	-	expression tag	UNP Q2TU34
B	344	HIS	-	expression tag	UNP Q2TU34
B	345	HIS	-	expression tag	UNP Q2TU34
B	346	HIS	-	expression tag	UNP Q2TU34
B	347	HIS	-	expression tag	UNP Q2TU34
C	338	LEU	-	expression tag	UNP Q2TU34

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Chain	Residue	Modelled	Actual	Comment	Reference
C	339	GLU	-	expression tag	UNP Q2TU34
C	340	HIS	-	expression tag	UNP Q2TU34
C	341	HIS	-	expression tag	UNP Q2TU34
C	342	HIS	-	expression tag	UNP Q2TU34
C	343	HIS	-	expression tag	UNP Q2TU34
C	344	HIS	-	expression tag	UNP Q2TU34
C	345	HIS	-	expression tag	UNP Q2TU34
C	346	HIS	-	expression tag	UNP Q2TU34
C	347	HIS	-	expression tag	UNP Q2TU34
D	338	LEU	-	expression tag	UNP Q2TU34
D	339	GLU	-	expression tag	UNP Q2TU34
D	340	HIS	-	expression tag	UNP Q2TU34
D	341	HIS	-	expression tag	UNP Q2TU34
D	342	HIS	-	expression tag	UNP Q2TU34
D	343	HIS	-	expression tag	UNP Q2TU34
D	344	HIS	-	expression tag	UNP Q2TU34
D	345	HIS	-	expression tag	UNP Q2TU34
D	346	HIS	-	expression tag	UNP Q2TU34
D	347	HIS	-	expression tag	UNP Q2TU34

- Molecule 2 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



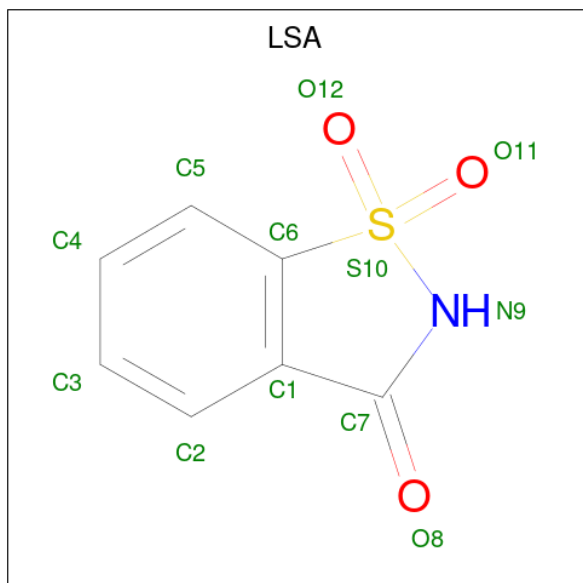
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

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

- Molecule 3 is 1,2-BENZISOTHAZOL-3(2H)-ONE 1,1-DIOXIDE (CCD ID: LSA) (formula: $C_7H_5NO_3S$).

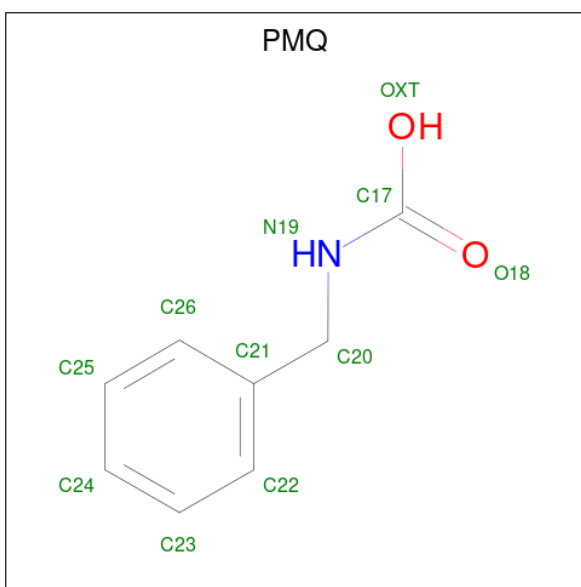


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	7	1	3	1		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

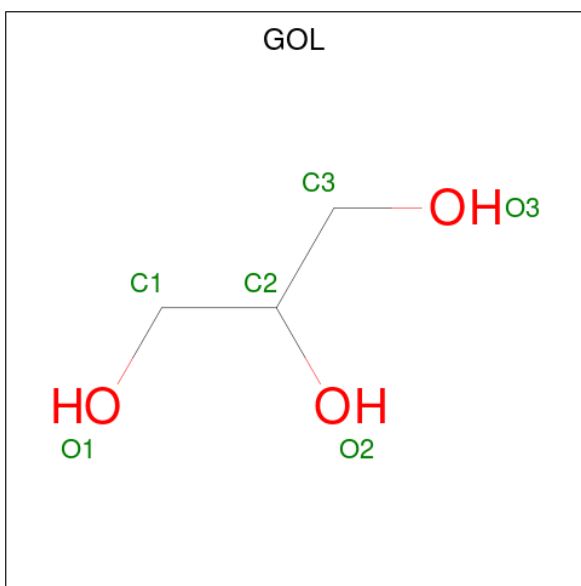
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		

- Molecule 5 is benzylcarbamic acid (CCD ID: PMQ) (formula: $C_8H_9NO_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			10	8	1	1		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	D	1	Total	C	O	0	0
			6	3	3		

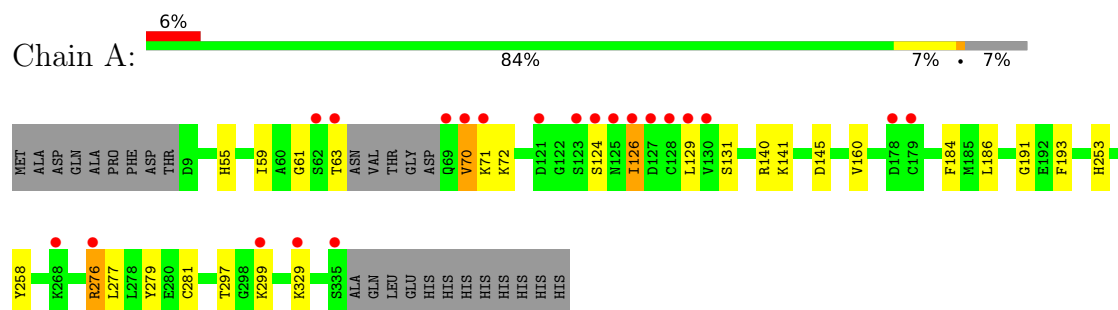
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	291	Total 291	O 291	0	0
7	B	328	Total 328	O 328	0	0
7	C	270	Total 270	O 270	0	0
7	D	269	Total 269	O 269	0	0

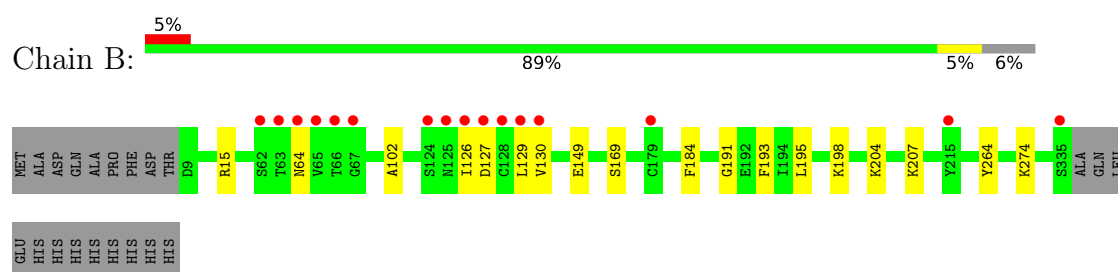
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

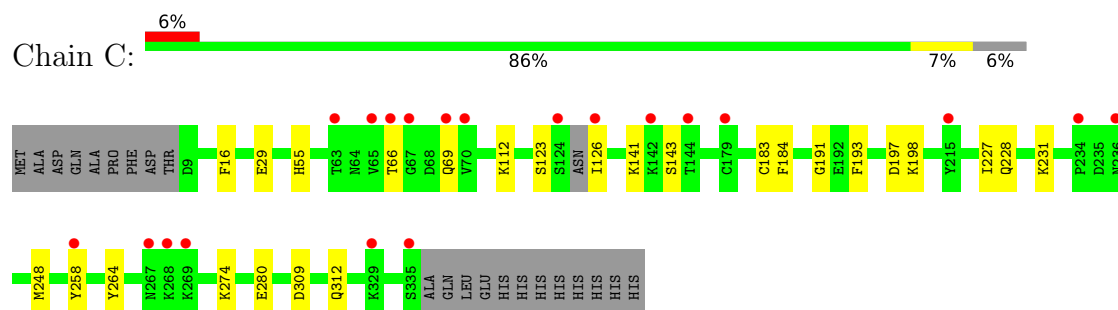
• Molecule 1: Fructose-1,6-bisphosphatase 1



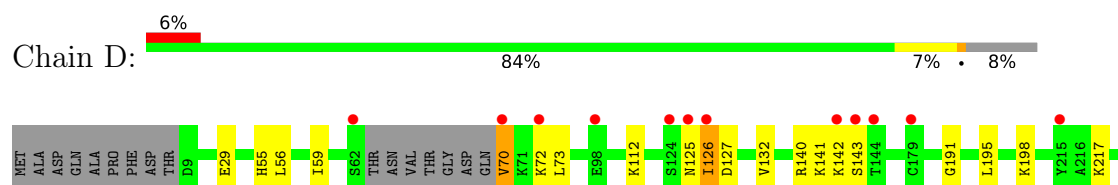
• Molecule 1: Fructose-1,6-bisphosphatase 1

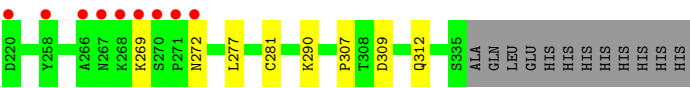


• Molecule 1: Fructose-1,6-bisphosphatase 1



• Molecule 1: Fructose-1,6-bisphosphatase 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.64Å 83.62Å 278.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.45 – 1.72 46.45 – 1.72	Depositor EDS
% Data completeness (in resolution range)	99.8 (46.45-1.72) 99.8 (46.45-1.72)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 1.73Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.174 , 0.198 0.175 , 0.199	Depositor DCC
R_{free} test set	2000 reflections (1.20%)	wwPDB-VP
Wilson B-factor (Å ²)	20.2	Xtriage
Anisotropy	0.153	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11179	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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: LSA, PMQ, MG, GOL, 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	0.36	1/2514 (0.0%)	0.55	2/3395 (0.1%)
1	B	0.30	0/2529	0.50	0/3416
1	C	0.29	0/2534	0.51	0/3423
1	D	0.38	0/2486	0.55	0/3358
All	All	0.33	1/10063 (0.0%)	0.53	2/13592 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	70	VAL	C-N	6.62	1.42	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	ILE	N-CA-C	-5.90	104.99	113.07
1	A	61	GLY	N-CA-C	5.73	118.73	111.85

There are no chirality outliers.

There are no planarity outliers.

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	2471	0	2514	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2485	0	2527	20	0
1	C	2488	0	2529	16	0
1	D	2443	0	2485	16	0
2	A	23	0	12	0	0
2	B	23	0	12	0	0
2	C	23	0	12	0	0
2	D	23	0	12	0	0
3	A	12	0	5	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	B	10	0	8	5	0
6	B	6	0	8	0	0
6	C	6	0	8	0	0
6	D	6	0	8	0	0
7	A	291	0	0	5	0
7	B	328	0	0	2	0
7	C	270	0	0	3	0
7	D	269	0	0	4	0
All	All	11179	0	10140	69	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 3.

All (69) 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:127:ASP:HA	5:B:501:PMQ:H23	1.52	0.90
1:B:102:ALA:HB3	1:B:149:GLU:HG2	1.71	0.72
1:C:198:LYS:NZ	7:C:502:HOH:O	2.28	0.64
1:B:102:ALA:CB	1:B:149:GLU:HG2	2.29	0.63
1:C:264:TYR:CZ	1:C:274:LYS:HD3	2.34	0.62
1:A:140:ARG:NH2	7:A:502:HOH:O	2.31	0.62
1:A:140:ARG:HG3	1:A:160:VAL:HG21	1.81	0.61
1:B:129:LEU:HB2	5:B:501:PMQ:HN19	1.66	0.60
1:D:56:LEU:HD21	1:D:70:VAL:HB	1.83	0.59
1:A:191:GLY:HA3	1:C:191:GLY:HA3	1.87	0.55
1:D:72:LYS:HZ2	1:D:73:LEU:H	1.53	0.55
1:B:191:GLY:HA3	1:D:191:GLY:HA3	1.88	0.54
1:B:129:LEU:H	5:B:501:PMQ:C22	2.18	0.54
1:A:70:VAL:HG22	1:A:126:ILE:CG2	2.39	0.53
1:A:71:LYS:NZ	7:A:501:HOH:O	2.30	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:HIS:NE2	7:C:503:HOH:O	2.33	0.51
1:A:297:THR:OG1	1:A:299:LYS:HG2	2.10	0.51
1:D:195:LEU:HD21	1:D:198:LYS:HG2	1.92	0.51
1:C:248:MET:HE1	1:C:280:GLU:HB3	1.93	0.50
1:A:141:LYS:NZ	7:A:509:HOH:O	2.45	0.49
1:C:16:PHE:HE2	1:C:198:LYS:HE2	1.76	0.49
1:A:276[A]:ARG:HG2	1:A:279:TYR:OH	2.12	0.49
1:C:183:CYS:HB2	1:C:197:ASP:HB2	1.95	0.49
1:A:70:VAL:HG22	1:A:126:ILE:HG22	1.95	0.48
1:D:126:ILE:HD12	1:D:132:VAL:HG11	1.94	0.48
1:C:29:GLU:OE2	1:C:112:LYS:HG2	2.13	0.48
1:C:258:TYR:OH	1:D:127:ASP:HB2	2.14	0.48
1:D:277:LEU:HA	1:D:281:CYS:HB2	1.97	0.46
1:D:277:LEU:HD23	1:D:307:PRO:HB3	1.98	0.46
1:C:66:THR:CB	1:C:69:GLN:HG3	2.46	0.46
1:A:253:HIS:CD2	5:B:501:PMQ:H25	2.50	0.46
1:B:184:PHE:HB3	1:B:193:PHE:HB3	1.98	0.46
1:C:123:SER:HA	1:C:126:ILE:HB	1.97	0.46
1:A:129:LEU:O	1:B:169:SER:N	2.37	0.45
1:A:253:HIS:CD2	1:B:129:LEU:HD23	2.51	0.45
1:B:264:TYR:CZ	1:B:274:LYS:HD3	2.52	0.45
1:D:290:LYS:NZ	7:D:504:HOH:O	2.48	0.45
1:A:258:TYR:CD2	1:B:126:ILE:HD11	2.52	0.45
1:C:228:GLN:NE2	7:C:501:HOH:O	2.20	0.44
1:B:207:LYS:NZ	7:B:608:HOH:O	2.50	0.44
1:D:309:ASP:HB3	1:D:312:GLN:HB3	2.00	0.44
1:B:204:LYS:CA	1:B:204:LYS:HE2	2.47	0.44
1:A:72:LYS:HD3	7:A:739:HOH:O	2.18	0.43
1:A:258:TYR:CE1	1:B:126:ILE:HG13	2.52	0.43
1:B:129:LEU:HB2	5:B:501:PMQ:N19	2.32	0.43
1:D:29:GLU:OE1	1:D:112:LYS:HG2	2.19	0.43
1:B:204:LYS:HE2	1:B:204:LYS:HA	1.99	0.43
1:C:141:LYS:HE2	1:C:143:SER:OG	2.19	0.43
1:A:131:SER:OG	1:B:130:VAL:HG13	2.18	0.43
1:A:277:LEU:HA	1:A:281:CYS:HB2	2.01	0.43
1:C:227:ILE:HG22	1:C:231:LYS:HD2	2.01	0.43
1:B:195:LEU:HD21	1:B:198:LYS:HG2	2.01	0.42
1:C:184:PHE:HB3	1:C:193:PHE:HB3	2.01	0.42
1:A:184:PHE:HB3	1:A:193:PHE:HB3	2.00	0.42
1:D:55:HIS:HA	1:D:59:ILE:HG22	2.02	0.42
1:A:71:LYS:HA	1:A:71:LYS:HD3	1.81	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:LYS:NZ	1:A:145:ASP:O	2.39	0.41
1:A:186:LEU:HB2	1:A:193:PHE:CE1	2.56	0.41
1:D:272:ASN:HA	7:D:507:HOH:O	2.20	0.41
1:B:15:ARG:HD3	7:D:503:HOH:O	2.20	0.41
1:C:309:ASP:HB3	1:C:312:GLN:HB3	2.02	0.41
1:B:198:LYS:NZ	7:B:616:HOH:O	2.54	0.41
1:D:140:ARG:O	1:D:142:LYS:HD3	2.21	0.41
1:A:329:LYS:HA	1:A:329:LYS:HD2	1.88	0.41
1:A:55:HIS:HA	1:A:59:ILE:HG22	2.03	0.40
1:D:126:ILE:HA	7:D:506:HOH:O	2.20	0.40
1:D:141:LYS:NZ	1:D:143:SER:HB2	2.36	0.40
1:A:126:ILE:HG21	1:A:126:ILE:HD13	1.81	0.40
1:A:299:LYS:HE2	7:A:738:HOH:O	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	319/348 (92%)	312 (98%)	6 (2%)	1 (0%)	36	24
1	B	325/348 (93%)	319 (98%)	5 (2%)	1 (0%)	36	24
1	C	323/348 (93%)	317 (98%)	6 (2%)	0	100	100
1	D	316/348 (91%)	310 (98%)	6 (2%)	0	100	100
All	All	1283/1392 (92%)	1258 (98%)	23 (2%)	2 (0%)	43	31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	SER
1	B	64	ASN

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	269/291 (92%)	266 (99%)	3 (1%)	65	50
1	B	270/291 (93%)	270 (100%)	0	100	100
1	C	270/291 (93%)	270 (100%)	0	100	100
1	D	266/291 (91%)	261 (98%)	5 (2%)	50	27
All	All	1075/1164 (92%)	1067 (99%)	8 (1%)	76	64

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	THR
1	A	276[A]	ARG
1	A	276[B]	ARG
1	D	70	VAL
1	D	125	ASN
1	D	126	ILE
1	D	217	LYS
1	D	269	LYS

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

Mol	Chain	Res	Type
1	A	272	ASN
1	C	236	ASN
1	C	272	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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$
2	AMP	B	502	-	25,25,25	1.35	5 (20%)	37,38,38	1.80	8 (21%)
6	GOL	B	503	-	5,5,5	0.69	0	5,5,5	1.00	0
5	PMQ	B	501	1	10,10,11	1.43	1 (10%)	10,11,13	1.75	2 (20%)
6	GOL	C	402	-	5,5,5	0.82	0	5,5,5	1.03	0
2	AMP	A	401	-	25,25,25	1.39	4 (16%)	37,38,38	1.90	8 (21%)
3	LSA	A	402	-	13,13,13	4.90	5 (38%)	20,20,20	3.52	12 (60%)
2	AMP	C	401	-	25,25,25	1.43	5 (20%)	37,38,38	1.80	8 (21%)
6	GOL	D	402	-	5,5,5	0.76	0	5,5,5	1.00	0
2	AMP	D	401	-	25,25,25	1.43	5 (20%)	37,38,38	1.78	8 (21%)

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	AMP	B	502	-	-	0/10/26/26	0/3/3/3
6	GOL	B	503	-	-	0/4/4/4	-
5	PMQ	B	501	1	-	1/4/4/5	0/1/1/1
6	GOL	C	402	-	-	0/4/4/4	-
2	AMP	A	401	-	-	0/10/26/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	LSA	A	402	-	-	-	0/2/2/2
2	AMP	C	401	-	-	0/10/26/26	0/3/3/3
6	GOL	D	402	-	-	2/4/4/4	-
2	AMP	D	401	-	-	0/10/26/26	0/3/3/3

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	LSA	C6-S10	-15.83	1.56	1.75
3	A	402	LSA	C1-C6	5.91	1.46	1.38
2	C	401	AMP	C5-C4	4.63	1.47	1.39
2	A	401	AMP	C5-C4	4.63	1.47	1.39
2	D	401	AMP	C5-C4	4.54	1.47	1.39
2	B	502	AMP	C5-C4	4.41	1.47	1.39
5	B	501	PMQ	C17-N19	4.25	1.43	1.33
3	A	402	LSA	S10-N9	-3.02	1.59	1.65
2	D	401	AMP	C5-C6	2.57	1.48	1.41
2	B	502	AMP	C5-C6	2.53	1.48	1.41
2	D	401	AMP	C8-N7	2.50	1.36	1.31
2	A	401	AMP	C5-N7	-2.49	1.34	1.39
2	C	401	AMP	C5-C6	2.47	1.47	1.41
2	D	401	AMP	C5-N7	-2.44	1.34	1.39
3	A	402	LSA	C2-C1	-2.42	1.36	1.39
2	B	502	AMP	C8-N7	2.38	1.36	1.31
2	A	401	AMP	C8-N7	2.36	1.36	1.31
2	C	401	AMP	C5-N7	-2.35	1.34	1.39
2	C	401	AMP	C8-N7	2.33	1.36	1.31
2	A	401	AMP	C5-C6	2.32	1.47	1.41
2	B	502	AMP	C5-N7	-2.16	1.35	1.39
2	B	502	AMP	C4-N9	-2.09	1.33	1.37
3	A	402	LSA	C7-N9	-2.08	1.34	1.37
2	C	401	AMP	C4-N9	-2.03	1.33	1.37
2	D	401	AMP	C4-N9	-2.03	1.33	1.37

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	LSA	C6-C1-C7	-9.57	103.72	112.23
2	A	401	AMP	C5-C4-N3	-5.80	118.73	126.72
2	D	401	AMP	C5-C4-N3	-5.58	119.03	126.72
2	C	401	AMP	C5-C4-N3	-5.36	119.33	126.72
2	B	502	AMP	C5-C4-N3	-5.36	119.33	126.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	LSA	C6-S10-N9	5.32	101.34	93.19
3	A	402	LSA	C7-N9-S10	-5.06	107.07	114.96
2	A	401	AMP	N3-C4-N9	4.98	135.64	127.17
3	A	402	LSA	O11-S10-O12	-4.92	110.21	116.32
5	B	501	PMQ	O18-C17-N19	-4.83	119.17	125.03
2	B	502	AMP	N3-C4-N9	4.71	135.17	127.17
2	C	401	AMP	N3-C4-N9	4.70	135.17	127.17
2	D	401	AMP	N3-C4-N9	4.40	134.65	127.17
3	A	402	LSA	C1-C7-N9	4.04	116.89	109.76
2	A	401	AMP	C2-N3-C4	3.58	120.58	111.83
2	D	401	AMP	C2-N3-C4	3.57	120.54	111.83
2	B	502	AMP	C2-N3-C4	3.48	120.33	111.83
2	D	401	AMP	N3-C2-N1	-3.43	123.39	128.58
2	A	401	AMP	N3-C2-N1	-3.31	123.57	128.58
2	B	502	AMP	N3-C2-N1	-3.31	123.57	128.58
2	B	502	AMP	C4-N9-C8	3.30	109.20	105.74
2	C	401	AMP	C2-N3-C4	3.27	119.82	111.83
2	C	401	AMP	C4-N9-C8	3.27	109.17	105.74
3	A	402	LSA	O8-C7-N9	-3.24	122.17	124.25
2	C	401	AMP	N3-C2-N1	-3.11	123.87	128.58
2	D	401	AMP	C4-C5-N7	-3.03	107.12	110.58
2	A	401	AMP	C4-N9-C8	2.84	108.72	105.74
2	A	401	AMP	C4-C5-N7	-2.80	107.38	110.58
3	A	402	LSA	O8-C7-C1	-2.78	120.93	127.68
3	A	402	LSA	C3-C2-C1	-2.78	114.77	119.80
3	A	402	LSA	C4-C3-C2	2.77	123.66	120.24
2	C	401	AMP	C4-C5-N7	-2.72	107.48	110.58
3	A	402	LSA	C2-C1-C7	2.64	134.43	129.32
2	D	401	AMP	C4-N9-C8	2.46	108.32	105.74
2	B	502	AMP	C4-C5-N7	-2.41	107.82	110.58
2	A	401	AMP	C5-N7-C8	2.31	107.09	103.45
5	B	501	PMQ	C21-C20-N19	-2.30	108.47	112.61
3	A	402	LSA	C4-C5-C6	-2.28	114.68	118.76
2	C	401	AMP	C5-N7-C8	2.21	106.92	103.45
2	B	502	AMP	N9-C8-N7	-2.20	110.82	113.94
2	A	401	AMP	N6-C6-N1	2.17	123.22	118.38
2	D	401	AMP	C5-N7-C8	2.13	106.80	103.45
2	C	401	AMP	N9-C8-N7	-2.12	110.93	113.94
2	B	502	AMP	C5-N7-C8	2.04	106.66	103.45
2	D	401	AMP	C2-N1-C6	2.02	122.05	118.73
3	A	402	LSA	C3-C4-C5	2.01	122.72	120.24

There are no chirality outliers.

All (3) torsion outliers are listed below:

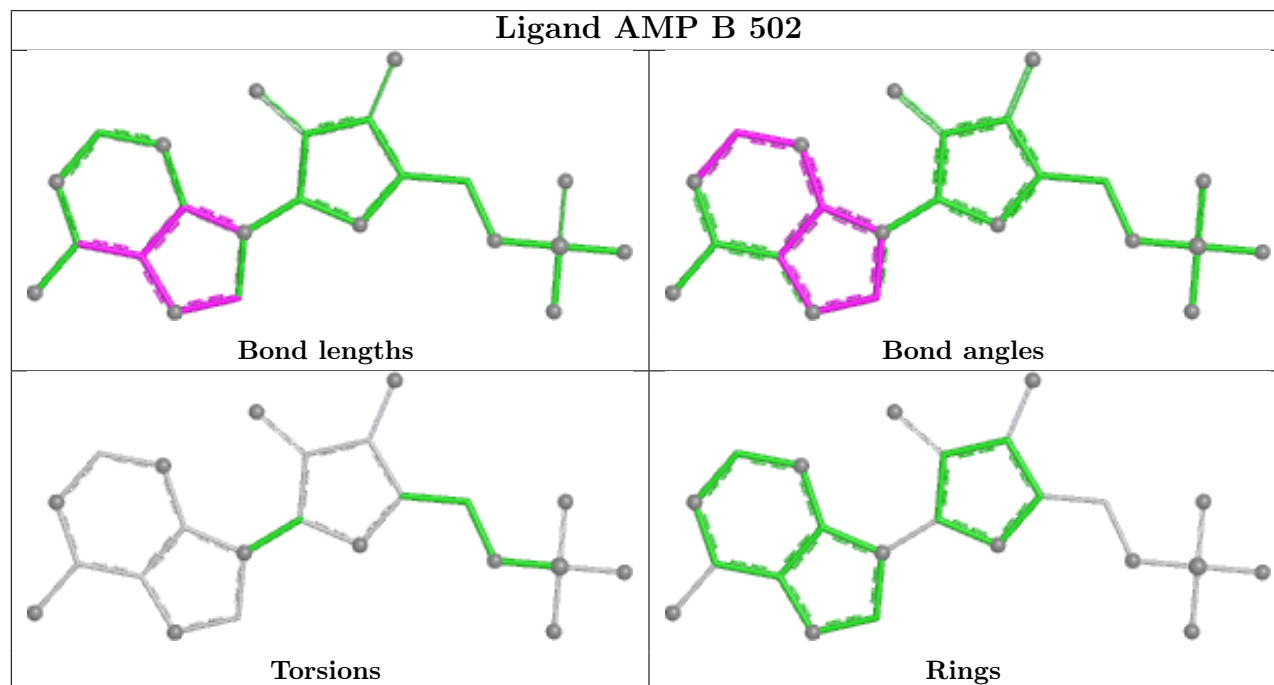
Mol	Chain	Res	Type	Atoms
6	D	402	GOL	C1-C2-C3-O3
6	D	402	GOL	O2-C2-C3-O3
5	B	501	PMQ	O18-C17-N19-C20

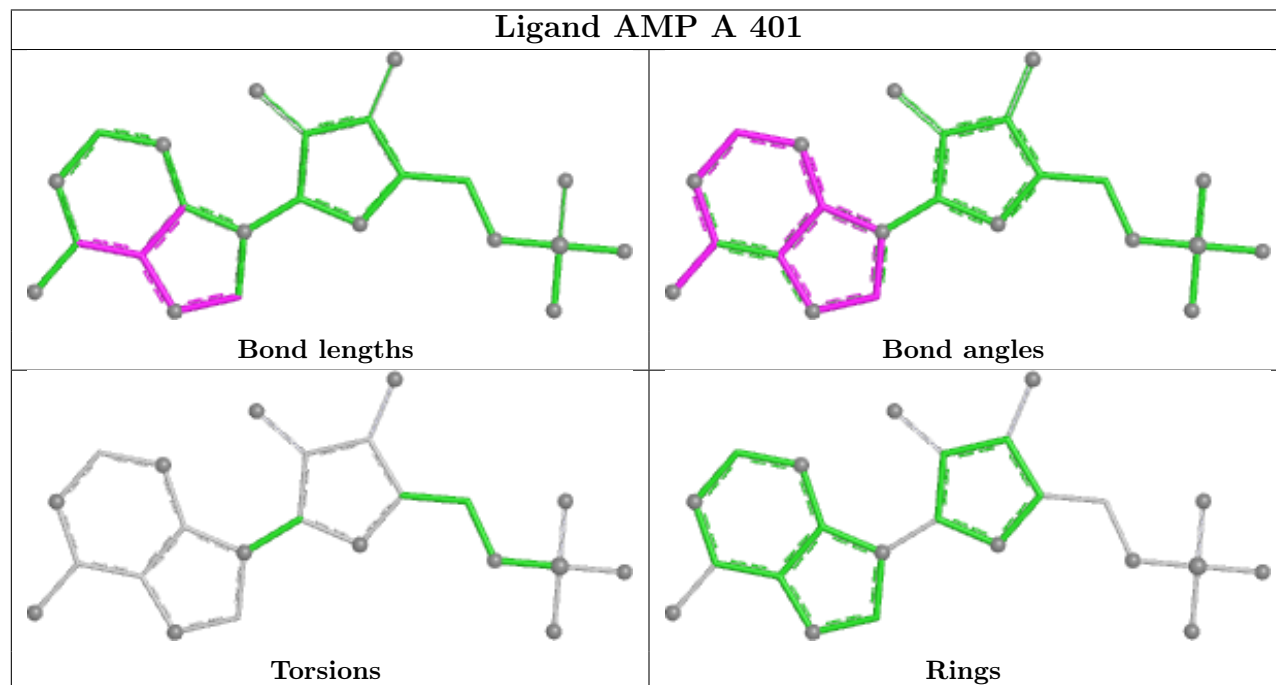
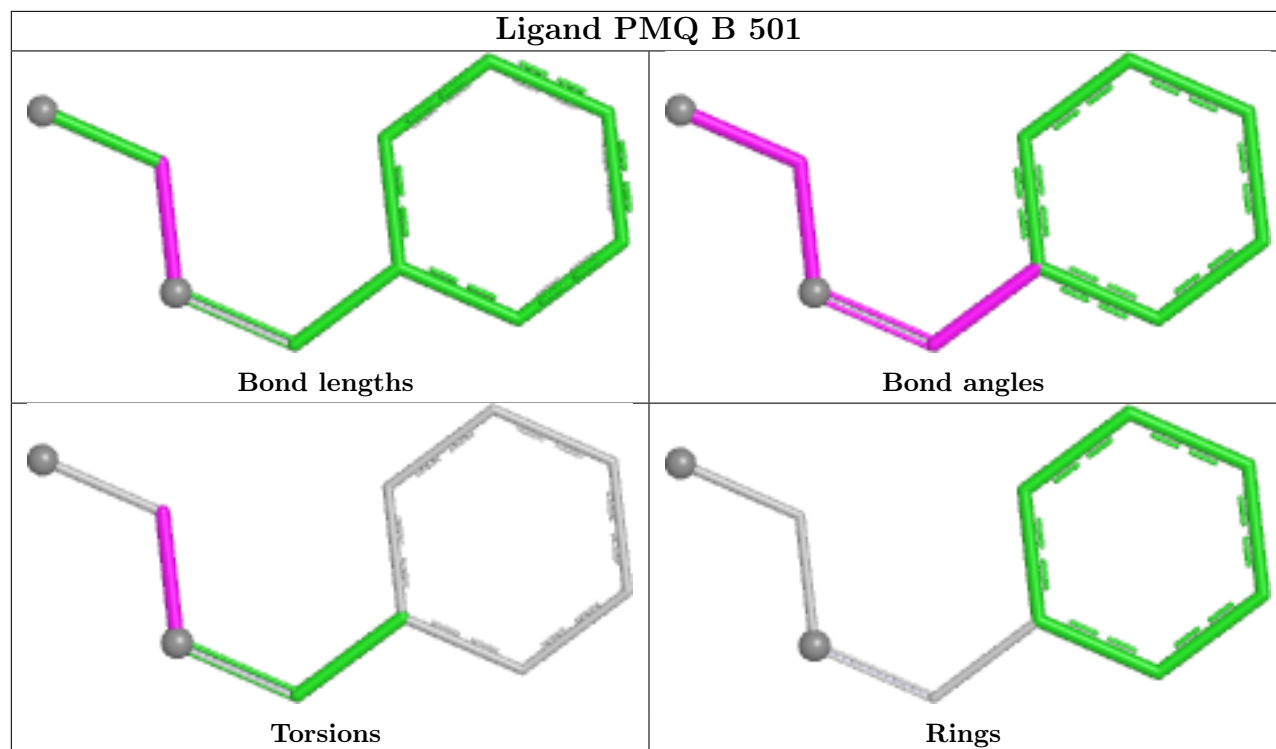
There are no ring outliers.

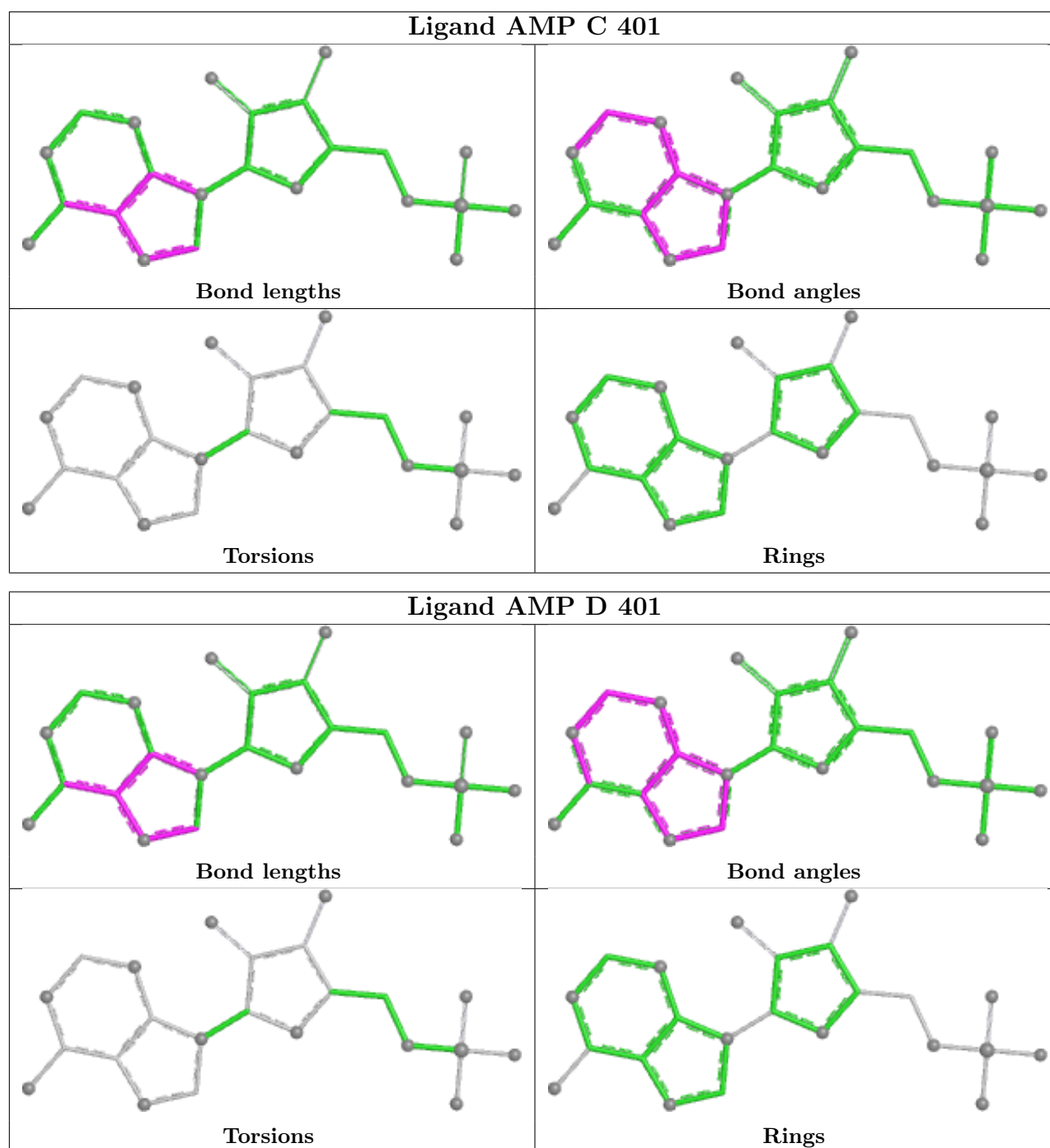
1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	501	PMQ	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	322/348 (92%)	0.07	21 (6%) 25 30	11, 20, 38, 57	1 (0%)
1	B	327/348 (93%)	-0.12	16 (4%) 35 41	12, 18, 36, 57	0
1	C	326/348 (93%)	0.18	20 (6%) 27 33	13, 22, 45, 57	1 (0%)
1	D	320/348 (91%)	0.05	21 (6%) 24 30	13, 19, 41, 55	0
All	All	1295/1392 (93%)	0.04	78 (6%) 27 33	11, 20, 41, 57	2 (0%)

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	70	VAL	7.1
1	A	63	THR	7.1
1	D	126	ILE	6.6
1	C	66	THR	6.5
1	A	128	CYS	6.2
1	D	125	ASN	5.8
1	A	62	SER	5.7
1	A	125	ASN	5.7
1	A	126	ILE	5.6
1	C	126	ILE	5.5
1	A	70	VAL	5.0
1	D	268	LYS	5.0
1	B	67	GLY	4.8
1	D	269	LYS	4.8
1	A	124	SER	4.7
1	B	129	LEU	4.6
1	B	66	THR	4.6
1	D	124	SER	4.6
1	D	271	PRO	4.4
1	C	67	GLY	4.4
1	B	65	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	127	ASP	4.2
1	B	130	VAL	4.0
1	A	130	VAL	3.9
1	B	126	ILE	3.9
1	D	266	ALA	3.8
1	C	258	TYR	3.7
1	B	62	SER	3.7
1	B	64	ASN	3.7
1	C	124	SER	3.6
1	A	335	SER	3.5
1	C	144	THR	3.5
1	D	267	ASN	3.4
1	A	71	LYS	3.3
1	B	125	ASN	3.3
1	C	65	VAL	3.3
1	C	179	CYS	3.3
1	B	335	SER	3.2
1	D	144	THR	3.2
1	C	236	ASN	3.1
1	C	215	TYR	3.1
1	C	268	LYS	3.1
1	A	127	ASP	3.1
1	B	63	THR	3.0
1	A	69	GLN	2.9
1	C	335	SER	2.9
1	D	62	SER	2.9
1	A	179	CYS	2.9
1	A	129	LEU	2.9
1	C	269	LYS	2.9
1	C	234	PRO	2.8
1	B	128	CYS	2.8
1	D	143	SER	2.8
1	A	276[A]	ARG	2.7
1	A	299	LYS	2.6
1	A	329	LYS	2.6
1	D	98	GLU	2.6
1	D	179	CYS	2.6
1	C	70	VAL	2.5
1	B	215	TYR	2.5
1	D	258	TYR	2.5
1	D	270	SER	2.4
1	D	215	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	267	ASN	2.3
1	B	179	CYS	2.2
1	A	123	SER	2.2
1	A	268	LYS	2.2
1	C	69	GLN	2.2
1	C	63	THR	2.2
1	D	220	ASP	2.2
1	C	142	LYS	2.2
1	B	124	SER	2.2
1	D	72	LYS	2.1
1	A	121	ASP	2.1
1	A	178	ASP	2.1
1	D	272	ASN	2.0
1	C	329	LYS	2.0
1	D	142	LYS	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.

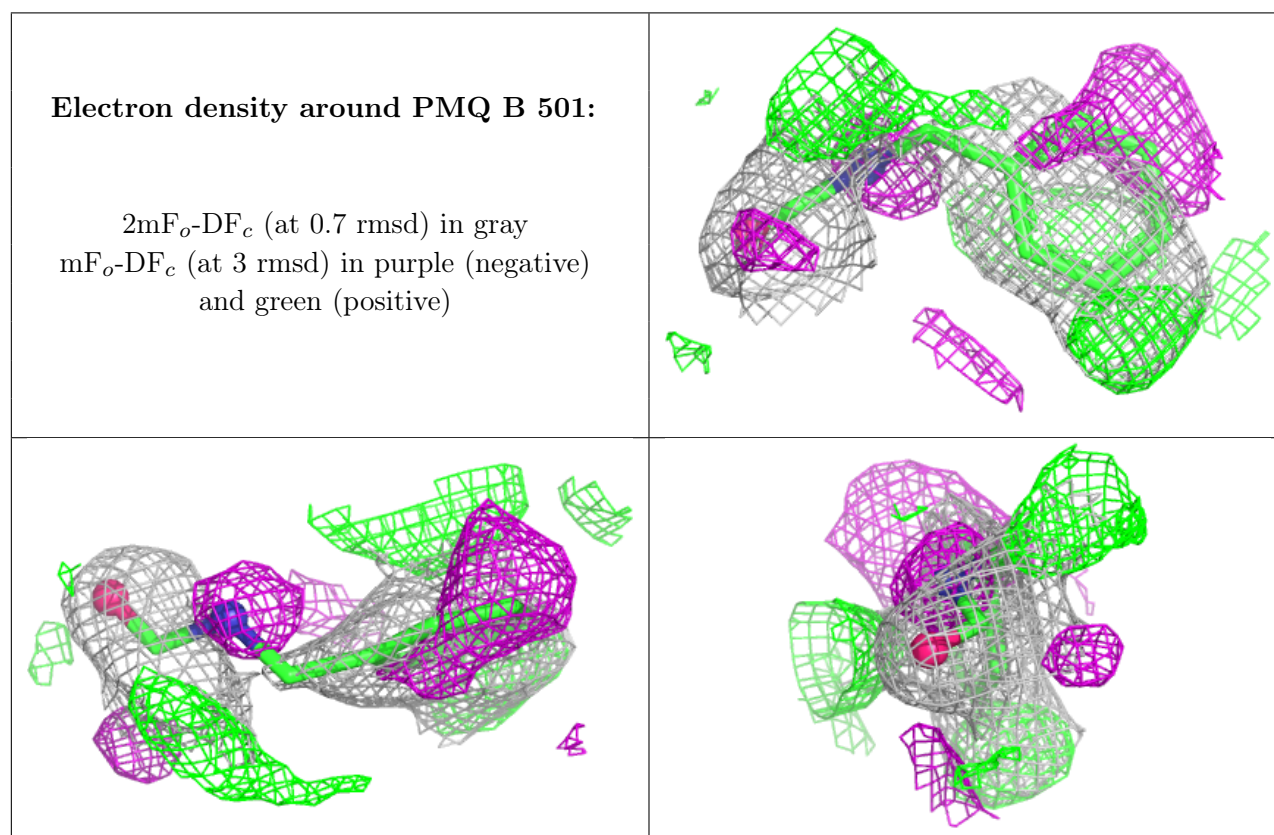
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	LSA	A	402	12/12	0.58	0.21	43,53,63,76	0
5	PMQ	B	501	10/11	0.61	0.22	37,41,44,53	0
6	GOL	D	402	6/6	0.85	0.16	23,39,46,46	0
6	GOL	C	402	6/6	0.88	0.13	26,36,40,43	0
6	GOL	B	503	6/6	0.94	0.12	21,26,28,31	0
2	AMP	C	401	23/23	0.97	0.07	17,21,23,25	0
4	MG	A	403	1/1	0.97	0.07	21,21,21,21	0
2	AMP	B	502	23/23	0.98	0.04	14,17,19,19	0

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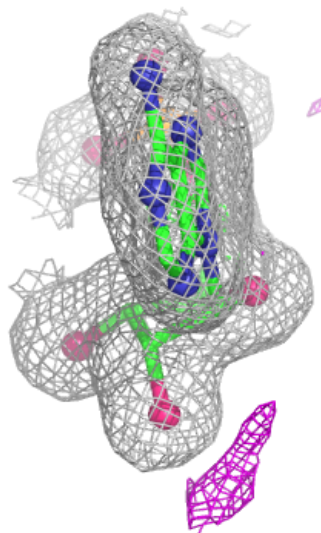
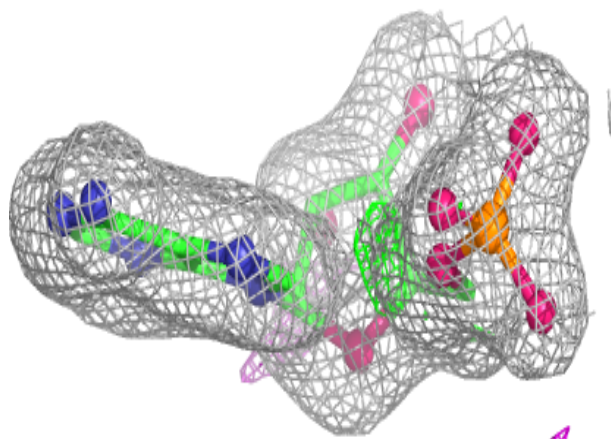
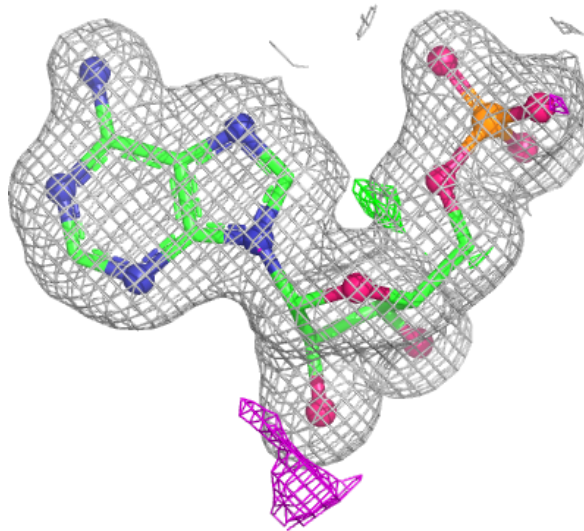
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	AMP	A	401	23/23	0.98	0.05	16,18,21,23	0
2	AMP	D	401	23/23	0.98	0.04	15,18,20,24	0
4	MG	B	504	1/1	0.99	0.02	15,15,15,15	0

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.



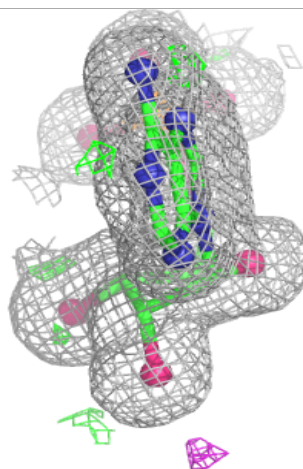
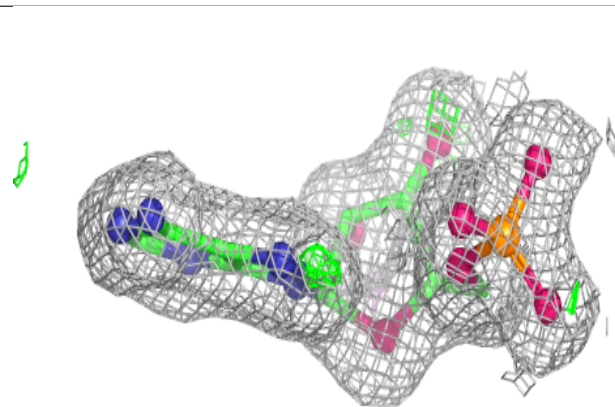
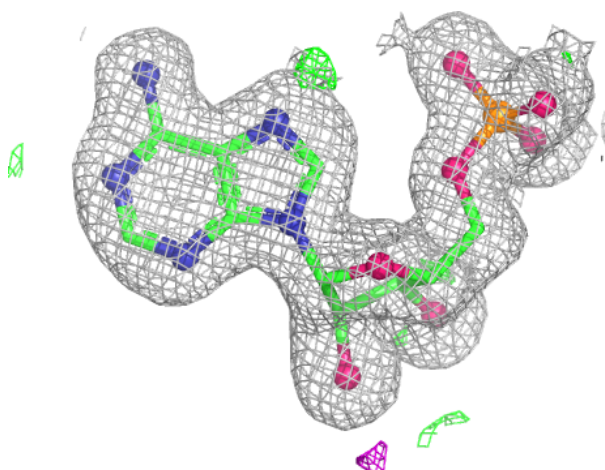
Electron density around AMP C 401:

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



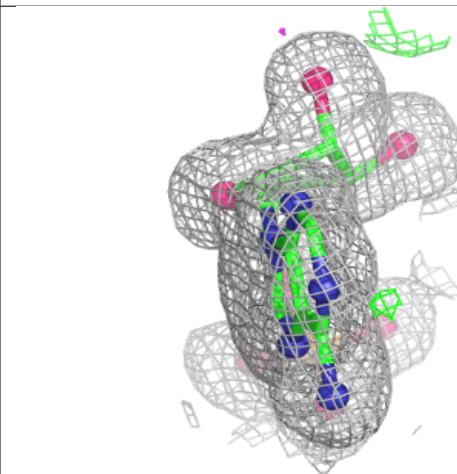
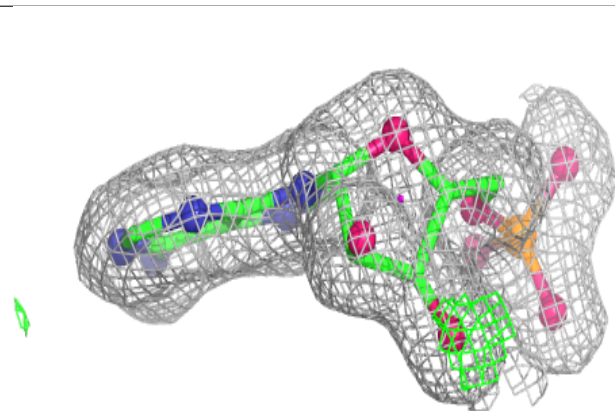
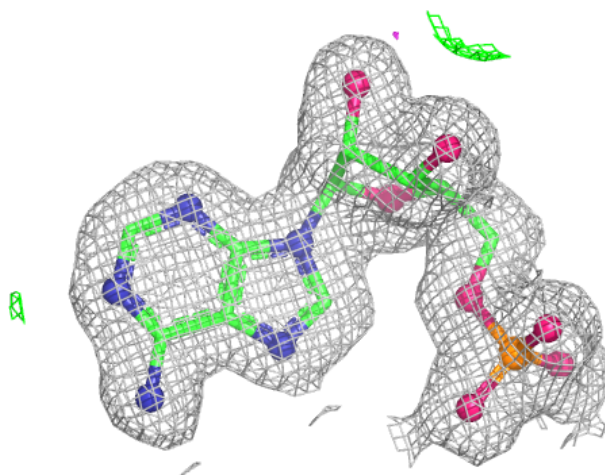
Electron density around AMP B 502:

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



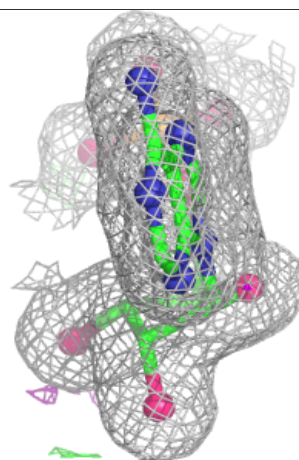
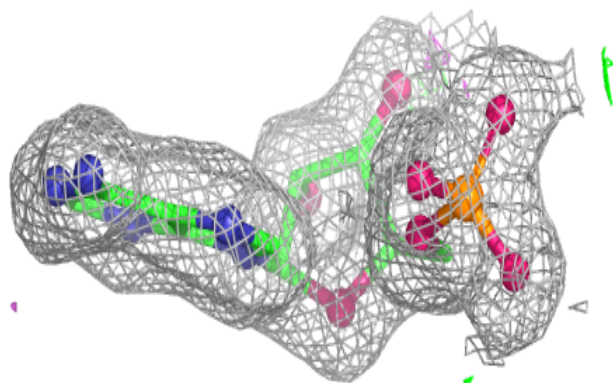
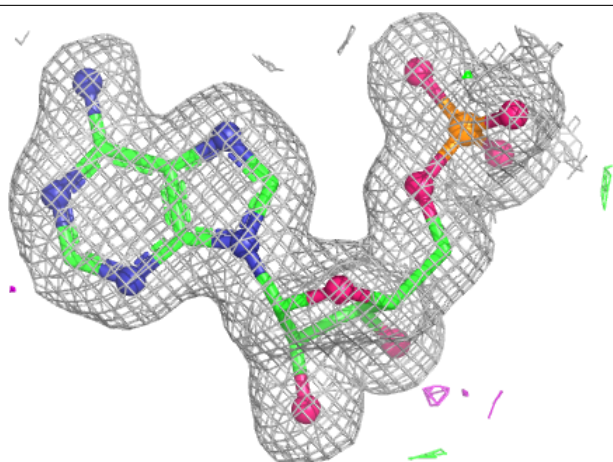
Electron density around AMP A 401:

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



Electron density around AMP D 401:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.