



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:14 PM UTC

PDB ID : 3X09 / pdb_00003x09
Title : Crystal structure of PIP4KIIBETA F205L complex with AMP
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Deposited on : 2014-10-09
Resolution : 2.70 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

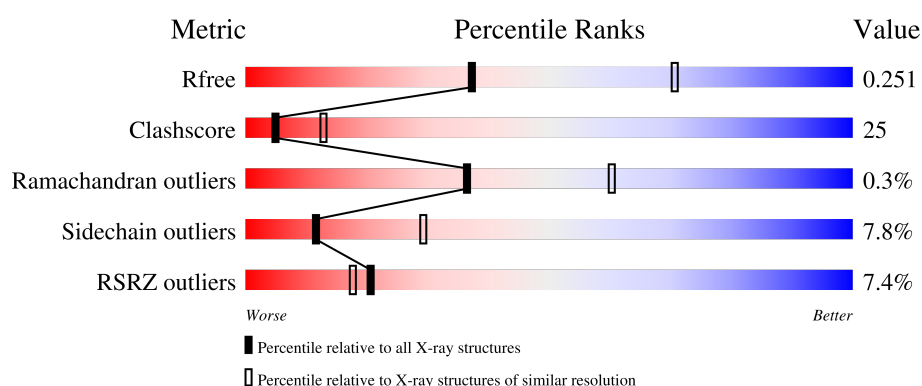
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	393	4% 50% 26% • 20%
1	B	393	7% 40% 34% • 22%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5274 atoms, of which 46 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.

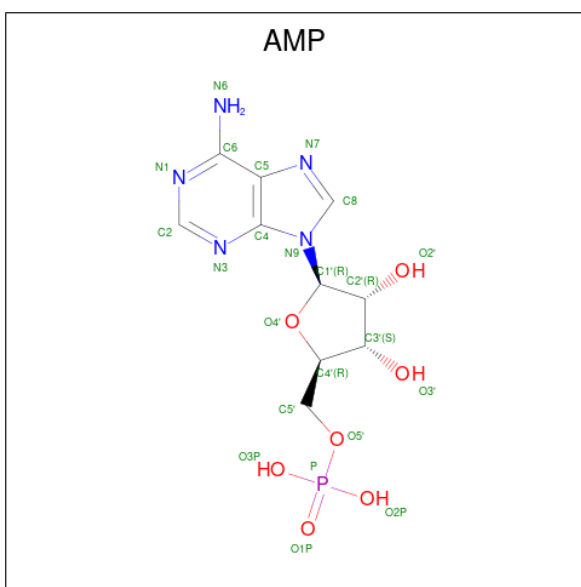
- Molecule 1 is a protein called Phosphatidylinositol 5-phosphate 4-kinase type-2 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	315	Total	C	N	O	S	0	0	0
			2592	1648	442	488	14			
1	B	305	Total	C	N	O	S	0	0	0
			2508	1603	431	461	13			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	GLY	-	expression tag	UNP P78356
A	25	PRO	-	expression tag	UNP P78356
A	26	ASN	-	expression tag	UNP P78356
A	27	CYS	-	expression tag	UNP P78356
A	28	ALA	-	expression tag	UNP P78356
A	29	PRO	-	expression tag	UNP P78356
A	30	GLY	-	expression tag	UNP P78356
A	205	LEU	PHE	engineered mutation	UNP P78356
B	24	GLY	-	expression tag	UNP P78356
B	25	PRO	-	expression tag	UNP P78356
B	26	ASN	-	expression tag	UNP P78356
B	27	CYS	-	expression tag	UNP P78356
B	28	ALA	-	expression tag	UNP P78356
B	29	PRO	-	expression tag	UNP P78356
B	30	GLY	-	expression tag	UNP P78356
B	205	LEU	PHE	engineered mutation	UNP P78356

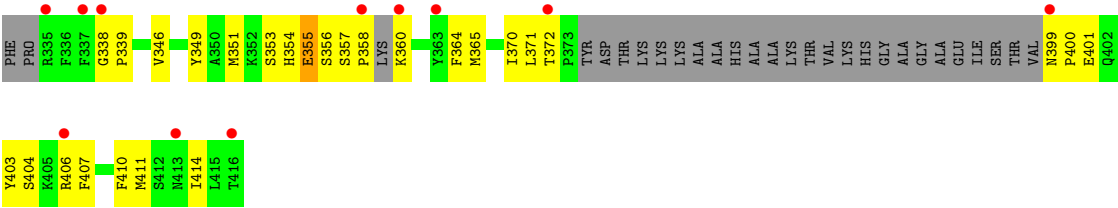
- Molecule 2 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
2	A	1	Total 33	C 10	H 10	N 5	O 7	P 1	0	0
2	A	1	Total 35	C 10	H 12	N 5	O 7	P 1	0	0
2	A	1	Total 35	C 10	H 12	N 5	O 7	P 1	0	0
2	B	1	Total 35	C 10	H 12	N 5	O 7	P 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	23	Total	O	0	0
			23	23		
3	B	13	Total	O	0	0
			13	13		



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	108.27Å 183.42Å 107.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	53.64 – 2.70 53.64 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.5 (53.64-2.70) 99.5 (53.64-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.42 (at 2.69Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.7.3_928)	Depositor
R, R_{free}	0.203 , 0.260 0.198 , 0.251	Depositor DCC
R_{free} test set	1479 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	65.2	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 66.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.020 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5274	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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

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	0.56	0/2643	0.92	5/3554 (0.1%)
1	B	0.57	1/2559 (0.0%)	0.89	1/3439 (0.0%)
All	All	0.57	1/5202 (0.0%)	0.90	6/6993 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	99	CYS	CA-C	-6.19	1.49	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	GLU	N-CA-C	5.96	118.94	108.75
1	A	204	VAL	N-CA-C	-5.84	104.63	111.00
1	B	213	ARG	N-CA-CB	-5.49	101.79	111.39
1	A	235	LEU	CA-C-N	5.31	125.07	119.76
1	A	235	LEU	C-N-CA	5.31	125.07	119.76
1	A	254	GLU	N-CA-C	5.05	116.47	111.07

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	2592	0	2563	107	0
1	B	2508	0	2490	157	0
2	A	69	34	36	3	0
2	B	23	12	12	3	0
3	A	23	0	0	3	0
3	B	13	0	0	1	0
All	All	5228	46	5101	259	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 25.

All (259) 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:218:LYS:HE2	1:B:403:TYR:OH	1.43	1.18
1:B:225:GLU:HG2	1:B:241:ASN:HB2	1.25	1.14
1:A:77:ILE:HD11	1:B:77:ILE:HD11	1.31	1.11
1:A:42:ILE:HD12	1:A:42:ILE:H	1.22	1.01
1:B:222:VAL:HG12	1:B:223:ALA:H	1.28	0.97
1:A:77:ILE:HD11	1:B:77:ILE:CD1	1.95	0.97
1:B:94:LYS:HB2	1:B:190:THR:HB	1.45	0.96
1:B:222:VAL:HG12	1:B:223:ALA:N	1.80	0.96
1:A:138:ARG:HG2	1:A:138:ARG:HH11	1.30	0.93
1:B:110:PHE:CE2	1:B:182:GLN:HG2	2.04	0.92
1:B:218:LYS:CE	1:B:278:ASP:C	2.44	0.90
1:A:155:GLU:OE2	1:A:155:GLU:N	2.05	0.90
1:B:179:LEU:HG	1:B:263:LYS:HG2	1.53	0.89
1:B:218:LYS:HZ3	1:B:278:ASP:HA	1.36	0.87
1:A:415:LEU:O	1:A:416:THR:HG23	1.75	0.87
1:A:77:ILE:CD1	1:B:77:ILE:HD11	2.04	0.87
1:B:222:VAL:CG1	1:B:223:ALA:H	1.88	0.86
1:B:218:LYS:NZ	1:B:278:ASP:HA	1.91	0.84
1:B:139:PHE:O	1:B:139:PHE:CD2	2.29	0.84
1:B:258:LYS:O	1:B:262:GLU:HG2	1.77	0.83
1:A:204:VAL:HG23	2:A:501:AMP:C6	2.14	0.82
1:B:218:LYS:HE2	1:B:403:TYR:HH	1.38	0.82
1:B:139:PHE:O	1:B:139:PHE:HD2	1.61	0.82
1:A:166:LYS:HD2	1:A:274:LEU:HD21	1.61	0.81
1:B:215:TYR:HB3	1:B:217:LEU:HD21	1.61	0.80
1:B:225:GLU:CG	1:B:241:ASN:HB2	2.08	0.80
1:B:243:PHE:HE2	1:B:414:ILE:HG22	1.45	0.80
1:B:78:LYS:HD2	1:B:94:LYS:HE2	1.63	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:MET:HE2	1:A:67:MET:HA	1.61	0.80
1:A:138:ARG:HH11	1:A:138:ARG:CG	1.94	0.80
1:B:268:VAL:HG12	1:B:404:SER:HB2	1.63	0.80
1:A:42:ILE:H	1:A:42:ILE:CD1	1.95	0.79
1:B:252:VAL:CG1	1:B:256:SER:HB2	2.10	0.79
1:B:118:GLN:O	1:B:122:THR:HB	1.82	0.79
1:B:218:LYS:HG2	1:B:403:TYR:OH	1.82	0.79
1:B:218:LYS:CE	1:B:403:TYR:OH	2.29	0.78
1:A:42:ILE:HD12	1:A:42:ILE:N	1.98	0.78
1:B:219:GLY:O	1:B:406:ARG:HD3	1.83	0.77
1:B:222:VAL:CG1	1:B:223:ALA:N	2.47	0.77
1:B:218:LYS:HE3	1:B:278:ASP:O	1.83	0.77
1:B:218:LYS:HE3	1:B:278:ASP:C	2.10	0.76
1:B:217:LEU:HB2	1:B:411:MET:HE1	1.67	0.76
1:B:236:PRO:HG2	1:B:238:PHE:CE1	2.20	0.76
1:A:256:SER:HB3	1:A:351:MET:CE	2.16	0.76
1:B:283:VAL:HG22	1:B:365:MET:HG2	1.66	0.76
1:B:139:PHE:CD2	1:B:139:PHE:C	2.64	0.75
1:B:218:LYS:NZ	1:B:278:ASP:CA	2.50	0.74
1:B:277:MET:O	1:B:278:ASP:HB2	1.87	0.73
1:B:252:VAL:HG13	1:B:256:SER:HB2	1.69	0.73
1:B:256:SER:HB3	1:B:351:MET:HE2	1.70	0.73
1:B:34:LYS:NZ	1:B:122:THR:HG21	2.04	0.72
1:B:236:PRO:HG2	1:B:238:PHE:HE1	1.52	0.72
1:B:288:VAL:CG1	1:B:339:PRO:HB2	2.18	0.72
1:A:295:GLU:O	1:A:298:VAL:HG12	1.90	0.71
1:A:256:SER:HB3	1:A:351:MET:HE2	1.73	0.71
1:A:260:PHE:CD2	1:A:415:LEU:HD11	2.26	0.70
1:A:141:THR:HG22	1:A:142:THR:O	1.91	0.70
1:A:204:VAL:HG13	1:A:366:ALA:HB3	1.74	0.70
1:B:243:PHE:HE2	1:B:414:ILE:CG2	2.04	0.70
1:A:204:VAL:HG23	2:A:501:AMP:N6	2.07	0.70
1:A:67:MET:HA	1:A:67:MET:CE	2.22	0.69
1:B:46:LEU:HD21	1:B:149:ILE:HD13	1.72	0.69
1:B:240:ASP:OD1	1:B:240:ASP:N	2.19	0.68
1:A:140:LEU:HB2	1:A:149:ILE:HB	1.76	0.67
1:B:109:ARG:CZ	1:B:172:VAL:HG22	2.25	0.67
1:B:224:ARG:C	1:B:225:GLU:HG3	2.19	0.67
1:A:208:ARG:HH22	1:A:342:PHE:HA	1.61	0.66
1:A:399:ASN:HB3	1:A:400:PRO:CD	2.24	0.66
1:A:141:THR:CG2	1:A:142:THR:O	2.44	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:LYS:HB2	1:A:190:THR:HB	1.78	0.65
1:B:407:PHE:O	1:B:411:MET:HG2	1.97	0.65
1:A:218:LYS:HD2	1:A:403:TYR:OH	1.97	0.65
1:A:219:GLY:O	1:A:406:ARG:HD2	1.96	0.65
1:B:78:LYS:HD2	1:B:94:LYS:CE	2.26	0.65
1:A:349:TYR:CD1	1:A:366:ALA:HB2	2.32	0.65
1:B:268:VAL:CG1	1:B:404:SER:HB2	2.27	0.65
1:B:246:GLU:O	1:B:246:GLU:HG3	1.97	0.64
1:B:191:VAL:O	1:B:192:ASP:HB2	1.97	0.64
1:B:218:LYS:CE	1:B:278:ASP:O	2.44	0.63
1:B:243:PHE:CE2	1:B:414:ILE:CG2	2.81	0.63
1:B:37:ARG:HG2	1:B:88:ASN:OD1	1.98	0.63
1:B:254:GLU:HG2	1:B:258:LYS:HE3	1.79	0.63
1:B:120:SER:O	1:B:142:THR:HB	1.98	0.63
1:A:399:ASN:HB3	1:A:400:PRO:HD2	1.82	0.62
1:A:399:ASN:CB	1:A:400:PRO:CD	2.77	0.62
1:A:218:LYS:HG2	1:A:279:TYR:O	2.00	0.61
1:B:240:ASP:HB3	1:B:410:PHE:CZ	2.36	0.61
1:A:261:LEU:HD21	1:A:412:SER:HA	1.83	0.61
1:B:205:LEU:HG	1:B:211:VAL:HG21	1.81	0.61
1:B:238:PHE:HB3	1:B:242:ASP:HB2	1.82	0.60
1:A:242:ASP:O	1:A:246:GLU:HG2	2.01	0.60
1:B:218:LYS:CE	1:B:278:ASP:HA	2.31	0.60
1:B:217:LEU:CB	1:B:411:MET:HE1	2.30	0.60
1:B:213:ARG:NH2	1:B:246:GLU:OE1	2.33	0.60
1:A:206:SER:HB2	1:A:364:PHE:CE2	2.37	0.60
1:A:415:LEU:O	1:A:416:THR:CG2	2.47	0.59
1:B:244:LEU:HD11	1:B:410:PHE:HE1	1.67	0.59
1:B:230:GLU:O	1:B:233:LYS:HG2	2.02	0.59
1:B:206:SER:HB2	1:B:364:PHE:CE2	2.38	0.59
1:A:43:LEU:HD21	1:A:140:LEU:HD11	1.85	0.58
1:A:277:MET:O	1:A:278:ASP:HB2	2.03	0.58
1:B:166:LYS:HD2	1:B:274:LEU:HD21	1.85	0.58
1:A:138:ARG:CG	1:A:138:ARG:NH1	2.60	0.58
1:A:224:ARG:O	1:A:241:ASN:OD1	2.22	0.57
1:B:204:VAL:HG23	2:B:501:AMP:C6	2.39	0.57
1:B:218:LYS:CE	1:B:278:ASP:CA	2.83	0.57
1:B:243:PHE:CE2	1:B:414:ILE:HG22	2.34	0.56
1:A:51:ASN:HB2	1:A:122:THR:HG21	1.87	0.56
1:A:67:MET:HE2	1:A:68:PRO:HD3	1.87	0.56
1:B:236:PRO:CG	1:B:238:PHE:CE1	2.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:277:MET:HE3	1:B:403:TYR:CD2	2.41	0.55
1:B:204:VAL:HG23	2:B:501:AMP:N1	2.21	0.55
1:B:224:ARG:CZ	1:B:239:LYS:HD3	2.37	0.55
1:A:217:LEU:HD13	1:A:414:ILE:HD11	1.87	0.55
1:B:399:ASN:N	1:B:400:PRO:CD	2.69	0.54
1:A:127:ILE:HG12	1:A:127:ILE:O	2.08	0.54
1:B:34:LYS:HD3	1:B:122:THR:HG23	1.90	0.53
1:B:218:LYS:NZ	1:B:278:ASP:C	2.66	0.53
1:A:77:ILE:CG1	1:B:77:ILE:HD11	2.38	0.53
1:B:34:LYS:HZ3	1:B:122:THR:HG21	1.73	0.53
1:A:157:VAL:HG21	1:A:197:TYR:CD2	2.44	0.53
1:B:355:GLU:N	1:B:355:GLU:OE2	2.42	0.53
1:B:204:VAL:HA	1:B:349:TYR:CD2	2.43	0.53
1:B:218:LYS:HZ3	1:B:278:ASP:CA	2.13	0.53
1:A:128:ASN:C	1:A:128:ASN:HD22	2.17	0.53
1:A:241:ASN:O	1:A:245:ASN:HB2	2.08	0.53
1:B:60:VAL:O	1:B:104:ARG:NH2	2.37	0.52
1:A:128:ASN:HB3	1:A:140:LEU:HD23	1.90	0.52
1:B:34:LYS:CE	1:B:122:THR:HG21	2.40	0.52
1:B:218:LYS:HE2	1:B:278:ASP:C	2.30	0.52
1:B:48:TRP:HH2	1:B:77:ILE:HD13	1.74	0.52
1:B:140:LEU:HB2	1:B:149:ILE:HB	1.91	0.52
1:B:189:LEU:HD12	1:B:189:LEU:N	2.24	0.52
1:B:81:ASN:O	1:B:91:SER:HB3	2.10	0.52
1:A:268:VAL:CG1	1:A:404:SER:HA	2.40	0.51
1:B:227:SER:OG	1:B:230:GLU:HB2	2.10	0.51
1:B:43:LEU:HD21	1:B:140:LEU:HD11	1.91	0.51
1:A:234:ASP:O	1:A:236:PRO:HD3	2.11	0.51
1:A:371:LEU:O	1:A:372:THR:CB	2.58	0.51
1:B:166:LYS:HD2	1:B:274:LEU:CD2	2.41	0.50
1:B:233:LYS:HB3	1:B:233:LYS:HZ3	1.75	0.50
1:B:218:LYS:CG	1:B:403:TYR:OH	2.56	0.50
1:A:204:VAL:HG23	2:A:501:AMP:N1	2.27	0.49
1:B:215:TYR:CB	1:B:217:LEU:HD21	2.36	0.49
1:A:204:VAL:CG1	1:A:366:ALA:HB3	2.40	0.49
1:B:34:LYS:CE	1:B:122:THR:CG2	2.90	0.49
1:B:281:LEU:HD12	1:B:282:LEU:N	2.27	0.49
1:A:66:LEU:O	1:A:67:MET:HE3	2.12	0.49
1:A:274:LEU:O	1:A:275:LYS:HB2	2.12	0.49
1:B:63:PRO:HD3	1:B:101:MET:HE1	1.95	0.49
1:B:189:LEU:N	1:B:189:LEU:CD1	2.76	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:ARG:NH1	1:A:172:VAL:HG22	2.29	0.48
1:A:185:GLY:HA2	3:A:603:HOH:O	2.13	0.48
1:B:216:ASP:C	1:B:217:LEU:HD23	2.38	0.48
1:A:74:TYR:CD1	1:A:74:TYR:C	2.92	0.48
1:A:77:ILE:HG22	1:A:95:PHE:HB3	1.95	0.48
1:B:99:CYS:N	1:B:100:PRO:CD	2.77	0.48
1:B:224:ARG:O	1:B:225:GLU:HG3	2.12	0.48
1:A:250:LEU:HD22	1:A:363:TYR:CE2	2.49	0.47
1:B:354:HIS:ND1	1:B:356:SER:HB2	2.29	0.47
1:B:254:GLU:CG	1:B:258:LYS:HE3	2.42	0.47
1:B:43:LEU:HD21	1:B:140:LEU:CD1	2.44	0.47
1:A:357:SER:OG	1:A:358:PRO:HD2	2.14	0.47
1:A:141:THR:HG23	1:A:145:ARG:HA	1.97	0.47
1:A:204:VAL:HG13	1:A:366:ALA:CB	2.43	0.47
1:A:217:LEU:CD1	1:A:414:ILE:HD11	2.44	0.47
1:B:218:LYS:HE2	1:B:278:ASP:HA	1.94	0.47
1:B:225:GLU:HA	1:B:239:LYS:HB2	1.97	0.47
1:B:256:SER:HB3	1:B:351:MET:CE	2.42	0.47
1:B:289:ASP:O	1:B:292:GLU:N	2.46	0.47
1:B:224:ARG:NH2	1:B:239:LYS:HD3	2.31	0.46
1:B:181:PRO:HG2	1:B:370:ILE:HG12	1.96	0.46
1:B:411:MET:O	1:B:414:ILE:HG12	2.15	0.46
1:B:113:ASP:HB3	1:B:116:ASP:OD1	2.16	0.46
1:B:256:SER:O	1:B:260:PHE:HB3	2.16	0.46
1:A:219:GLY:O	1:A:406:ARG:CD	2.62	0.46
1:B:230:GLU:HA	1:B:233:LYS:HD3	1.98	0.46
1:B:79:VAL:O	1:B:92:ARG:HA	2.15	0.46
1:A:67:MET:HE2	1:A:68:PRO:CD	2.46	0.45
1:A:141:THR:CG2	1:A:142:THR:N	2.76	0.45
1:B:243:PHE:CE1	1:B:248:GLN:HB3	2.51	0.45
1:B:233:LYS:CB	1:B:233:LYS:NZ	2.79	0.45
1:A:56:GLU:HG2	3:A:621:HOH:O	2.16	0.45
1:B:57:LEU:O	1:B:104:ARG:NH2	2.50	0.45
1:B:110:PHE:CZ	1:B:182:GLN:HG2	2.47	0.45
1:A:166:LYS:CD	1:A:274:LEU:HD21	2.40	0.44
1:B:215:TYR:HB3	1:B:217:LEU:CD2	2.41	0.44
1:A:160:MET:O	1:A:164:LEU:HB2	2.16	0.44
1:B:37:ARG:HD3	1:B:88:ASN:OD1	2.17	0.44
1:B:96:LYS:HB3	1:B:188:ARG:HB3	1.98	0.44
1:B:281:LEU:HD12	1:B:281:LEU:C	2.42	0.44
1:A:43:LEU:O	1:A:47:MET:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:LEU:HD23	1:A:106:LEU:HA	1.84	0.44
1:B:104:ARG:HA	1:B:107:ARG:NH1	2.33	0.44
1:A:35:LEU:HB3	1:A:88:ASN:ND2	2.33	0.44
1:A:138:ARG:NH1	1:A:138:ARG:CB	2.81	0.44
1:A:160:MET:CE	1:A:164:LEU:HD13	2.47	0.44
1:A:155:GLU:N	1:A:155:GLU:CD	2.74	0.44
1:B:218:LYS:NZ	1:B:278:ASP:O	2.49	0.44
1:B:218:LYS:HG3	1:B:279:TYR:O	2.17	0.44
1:A:84:PHE:O	1:A:85:ASN:C	2.61	0.43
1:A:218:LYS:HE2	1:A:224:ARG:CZ	2.48	0.43
1:B:159:GLU:O	1:B:162:ASN:HB3	2.18	0.43
1:A:103:PHE:HA	1:A:106:LEU:HB2	2.00	0.43
1:A:251:HIS:O	1:A:353:SER:HA	2.17	0.43
1:B:251:HIS:HB2	1:B:357:SER:HB2	1.99	0.43
1:A:190:THR:HA	1:A:194:VAL:O	2.19	0.43
1:A:343:ASP:OD2	1:A:345:SER:OG	2.26	0.43
1:B:219:GLY:N	1:B:240:ASP:OD2	2.52	0.43
1:B:371:LEU:O	1:B:372:THR:HG23	2.18	0.43
1:A:63:PRO:HB2	1:A:65:MET:O	2.19	0.43
1:A:186:MET:HG3	1:A:199:VAL:HG22	2.00	0.43
1:B:74:TYR:CD1	1:B:74:TYR:C	2.96	0.43
1:B:225:GLU:HB3	1:B:242:ASP:OD1	2.19	0.43
1:A:77:ILE:HD12	1:B:79:VAL:HG22	2.01	0.43
1:A:179:LEU:HG	1:A:263:LYS:HB3	2.00	0.43
1:B:157:VAL:O	1:B:161:HIS:HD2	2.01	0.43
1:B:204:VAL:HA	1:B:349:TYR:HD2	1.83	0.43
1:A:218:LYS:HB3	1:A:403:TYR:OH	2.19	0.43
1:B:207:HIS:CE1	1:B:346:VAL:HG11	2.53	0.42
1:B:36:PHE:HD1	1:B:47:MET:HB3	1.84	0.42
1:B:34:LYS:HE2	1:B:122:THR:CG2	2.48	0.42
1:B:64:VAL:HG23	1:B:65:MET:HG2	2.01	0.42
1:A:43:LEU:CD2	1:A:140:LEU:HD11	2.50	0.42
1:A:48:TRP:HB2	1:A:89:LEU:HD21	2.01	0.42
1:A:268:VAL:HG13	1:A:279:TYR:OH	2.20	0.42
2:B:501:AMP:N6	3:B:611:HOH:O	2.52	0.42
1:B:233:LYS:HB3	1:B:233:LYS:NZ	2.34	0.42
1:B:224:ARG:O	1:B:225:GLU:CG	2.67	0.42
1:A:206:SER:HB2	1:A:364:PHE:CZ	2.55	0.42
1:B:218:LYS:HE2	1:B:278:ASP:CA	2.50	0.42
1:A:243:PHE:HZ	1:A:414:ILE:HG23	1.85	0.41
1:A:141:THR:HG23	1:A:142:THR:N	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ARG:CG	1:B:88:ASN:OD1	2.66	0.41
1:B:233:LYS:HG2	1:B:233:LYS:H	1.53	0.41
1:B:121:VAL:O	1:B:121:VAL:HG12	2.19	0.41
1:A:110:PHE:CE1	1:A:182:GLN:HB3	2.55	0.41
1:A:153:SER:O	1:A:157:VAL:HG23	2.20	0.41
1:B:213:ARG:CZ	1:B:246:GLU:OE1	2.68	0.41
1:B:244:LEU:HD11	1:B:410:PHE:CE1	2.52	0.41
1:B:224:ARG:C	1:B:225:GLU:CG	2.91	0.41
1:A:160:MET:HE1	1:A:164:LEU:HD13	2.03	0.41
1:B:267:ASP:O	1:B:270:PHE:HB3	2.21	0.41
1:A:283:VAL:HG22	1:A:365:MET:HG2	2.02	0.41
1:B:288:VAL:HG12	1:B:339:PRO:HB2	1.98	0.41
1:A:60:VAL:HA	1:A:61:PRO:HD2	1.88	0.41
1:A:109:ARG:HG2	3:A:610:HOH:O	2.21	0.41
1:A:109:ARG:CZ	1:A:172:VAL:HG22	2.51	0.41
1:A:209:LEU:N	1:A:209:LEU:HD23	2.36	0.41
1:A:82:HIS:HB2	1:B:74:TYR:CZ	2.56	0.40
1:B:257:LYS:O	1:B:261:LEU:HG	2.20	0.40
1:A:219:GLY:O	1:A:406:ARG:NE	2.54	0.40
1:B:150:LYS:HE3	1:B:150:LYS:HB2	1.88	0.40
1:B:358:PRO:O	1:B:360:LYS:N	2.54	0.40
1:A:50:VAL:HG21	1:A:121:VAL:HG11	2.03	0.40
1:B:124:SER:HB3	1:B:143:TYR:CD2	2.57	0.40
1:A:205:LEU:HD22	1:A:211:VAL:HG21	2.04	0.40
1:A:276:ILE:HG22	1:A:277:MET:N	2.35	0.40
1:B:276:ILE:HG22	1:B:277:MET:N	2.36	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	305/393 (78%)	285 (93%)	19 (6%)	1 (0%)	36	60
1	B	291/393 (74%)	274 (94%)	16 (6%)	1 (0%)	36	60
All	All	596/786 (76%)	559 (94%)	35 (6%)	2 (0%)	36	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	338	GLY
1	A	61	PRO

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	293/352 (83%)	275 (94%)	18 (6%)	17	40
1	B	282/352 (80%)	255 (90%)	27 (10%)	8	20
All	All	575/704 (82%)	530 (92%)	45 (8%)	11	29

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ARG
1	A	42	ILE
1	A	67	MET
1	A	76	LYS
1	A	77	ILE
1	A	127	ILE
1	A	128	ASN
1	A	138	ARG
1	A	141	THR
1	A	149	ILE
1	A	214	LYS
1	A	224	ARG
1	A	295	GLU
1	A	298	VAL

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Mol	Chain	Res	Type
1	A	348	VAL
1	A	356	SER
1	A	360	LYS
1	A	406	ARG
1	B	37	ARG
1	B	42	ILE
1	B	104	ARG
1	B	122	THR
1	B	123	ARG
1	B	128	ASN
1	B	138	ARG
1	B	139	PHE
1	B	155	GLU
1	B	165	LYS
1	B	178	THR
1	B	189	LEU
1	B	205	LEU
1	B	208	ARG
1	B	210	THR
1	B	224	ARG
1	B	230	GLU
1	B	233	LYS
1	B	235	LEU
1	B	240	ASP
1	B	252	VAL
1	B	266	ARG
1	B	280	SER
1	B	285	ILE
1	B	353	SER
1	B	355	GLU
1	B	401	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	115	GLN
1	A	118	GLN
1	A	175	HIS
1	A	241	ASN
1	A	273	GLN
1	A	408	ASN
1	B	115	GLN

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Mol	Chain	Res	Type
1	B	128	ASN
1	B	161	HIS
1	B	175	HIS
1	B	207	HIS
1	B	251	HIS

5.3.3 RNA [i](#)

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](#)

4 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	$\# Z > 2$
2	AMP	A	501	-	25,25,25	1.51	4 (16%)	37,38,38	1.95	9 (24%)
2	AMP	A	502	-	25,25,25	1.54	4 (16%)	37,38,38	1.93	7 (18%)
2	AMP	A	503	-	25,25,25	1.42	4 (16%)	37,38,38	2.01	8 (21%)
2	AMP	B	501	-	25,25,25	1.43	5 (20%)	37,38,38	1.97	11 (29%)

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	A	501	-	-	6/10/26/26	0/3/3/3
2	AMP	A	502	-	-	1/10/26/26	0/3/3/3
2	AMP	A	503	-	-	3/10/26/26	0/3/3/3
2	AMP	B	501	-	-	2/10/26/26	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	AMP	C5-C4	5.26	1.48	1.39
2	A	502	AMP	C5-C4	5.05	1.48	1.39
2	B	501	AMP	C5-C4	4.62	1.47	1.39
2	A	503	AMP	C5-C4	4.45	1.47	1.39
2	A	502	AMP	C5-N7	-3.06	1.33	1.39
2	A	501	AMP	C5-C6	2.97	1.49	1.41
2	A	502	AMP	C5-C6	2.77	1.48	1.41
2	B	501	AMP	C5-C6	2.72	1.48	1.41
2	A	503	AMP	C5-C6	2.62	1.48	1.41
2	A	501	AMP	C8-N7	2.58	1.36	1.31
2	A	502	AMP	C8-N7	2.51	1.36	1.31
2	A	503	AMP	C5-N7	-2.40	1.34	1.39
2	B	501	AMP	C8-N7	2.39	1.36	1.31
2	A	503	AMP	C8-N7	2.23	1.36	1.31
2	B	501	AMP	C4-N9	-2.15	1.33	1.37
2	B	501	AMP	C5-N7	-2.14	1.35	1.39
2	A	501	AMP	C5-N7	-2.01	1.35	1.39

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	503	AMP	C5-C4-N3	-6.53	117.72	126.72
2	A	502	AMP	C5-C4-N3	-6.35	117.98	126.72
2	A	501	AMP	C5-C4-N3	-5.79	118.75	126.72
2	B	501	AMP	C5-C4-N3	-5.47	119.18	126.72
2	A	503	AMP	N3-C4-N9	5.10	135.83	127.17
2	A	502	AMP	N3-C4-N9	4.59	134.98	127.17
2	B	501	AMP	N3-C4-N9	4.51	134.84	127.17
2	A	501	AMP	N3-C4-N9	4.23	134.37	127.17
2	A	502	AMP	C4-C5-N7	-4.19	105.79	110.58
2	A	503	AMP	C4-C5-N7	-3.83	106.21	110.58
2	A	503	AMP	C2-N3-C4	3.69	120.85	111.83
2	A	501	AMP	C4-C5-N7	-3.67	106.39	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	AMP	C2-N3-C4	3.54	120.49	111.83
2	B	501	AMP	C4-C5-N7	-3.54	106.53	110.58
2	B	501	AMP	C4-N9-C8	3.51	109.43	105.74
2	A	501	AMP	C2-N3-C4	3.49	120.35	111.83
2	A	502	AMP	C2-N3-C4	3.36	120.03	111.83
2	B	501	AMP	N3-C2-N1	-3.26	123.64	128.58
2	A	502	AMP	C5-N7-C8	3.09	108.31	103.45
2	A	501	AMP	N3-C2-N1	-3.00	124.04	128.58
2	A	503	AMP	C5-N7-C8	2.97	108.12	103.45
2	A	503	AMP	C4-N9-C8	2.88	108.76	105.74
2	B	501	AMP	O3P-P-O5'	-2.79	99.40	106.67
2	A	501	AMP	O3P-P-O5'	-2.66	99.73	106.67
2	B	501	AMP	C5-N7-C8	2.66	107.62	103.45
2	A	503	AMP	N3-C2-N1	-2.63	124.60	128.58
2	A	503	AMP	N9-C8-N7	-2.48	110.42	113.94
2	B	501	AMP	N9-C8-N7	-2.48	110.42	113.94
2	A	501	AMP	C5-N7-C8	2.43	107.27	103.45
2	A	501	AMP	C2-N1-C6	2.40	122.67	118.73
2	A	502	AMP	N3-C2-N1	-2.30	125.11	128.58
2	B	501	AMP	C2-N1-C6	2.19	122.33	118.73
2	A	502	AMP	O3P-P-O5'	-2.18	101.00	106.67
2	B	501	AMP	C6-C5-N7	2.08	136.11	132.09
2	A	501	AMP	C4-N9-C8	2.08	107.92	105.74

There are no chirality outliers.

All (12) torsion outliers are listed below:

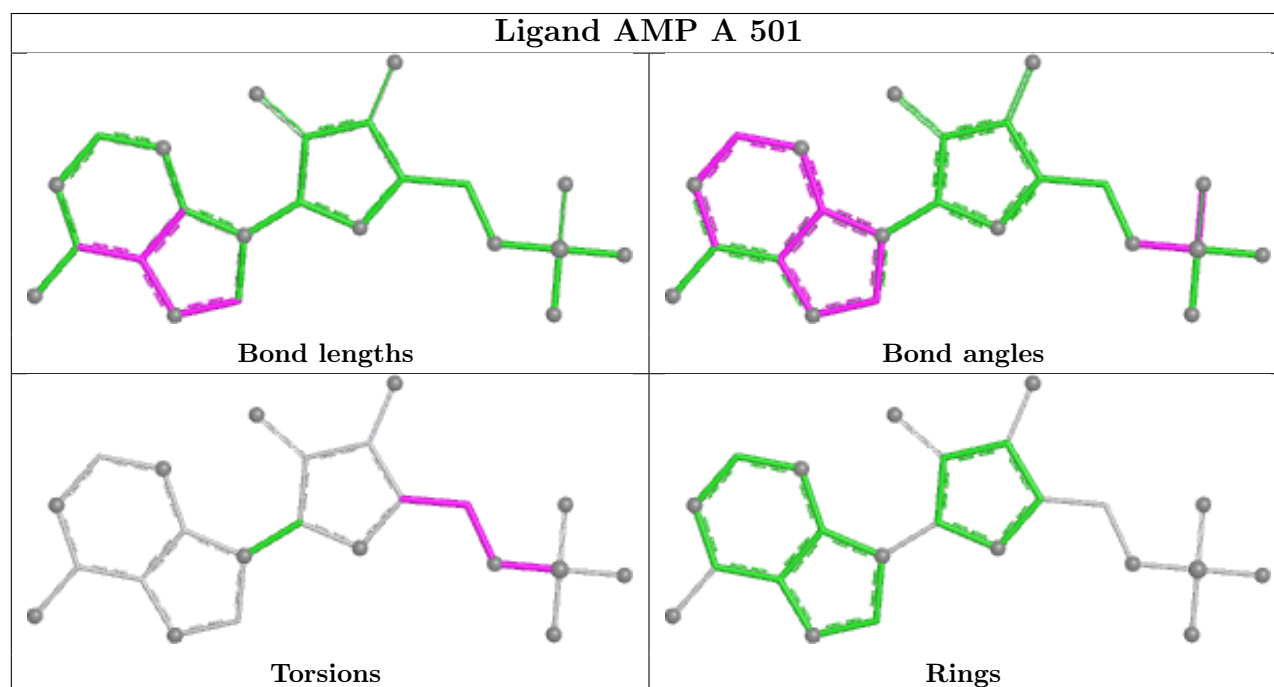
Mol	Chain	Res	Type	Atoms
2	A	501	AMP	C5'-O5'-P-O2P
2	A	501	AMP	C5'-O5'-P-O3P
2	A	501	AMP	O4'-C4'-C5'-O5'
2	A	503	AMP	C5'-O5'-P-O1P
2	A	503	AMP	C5'-O5'-P-O2P
2	A	503	AMP	C5'-O5'-P-O3P
2	A	501	AMP	C3'-C4'-C5'-O5'
2	B	501	AMP	O4'-C4'-C5'-O5'
2	B	501	AMP	C3'-C4'-C5'-O5'
2	A	501	AMP	C5'-O5'-P-O1P
2	A	502	AMP	C5'-O5'-P-O1P
2	A	501	AMP	C4'-C5'-O5'-P

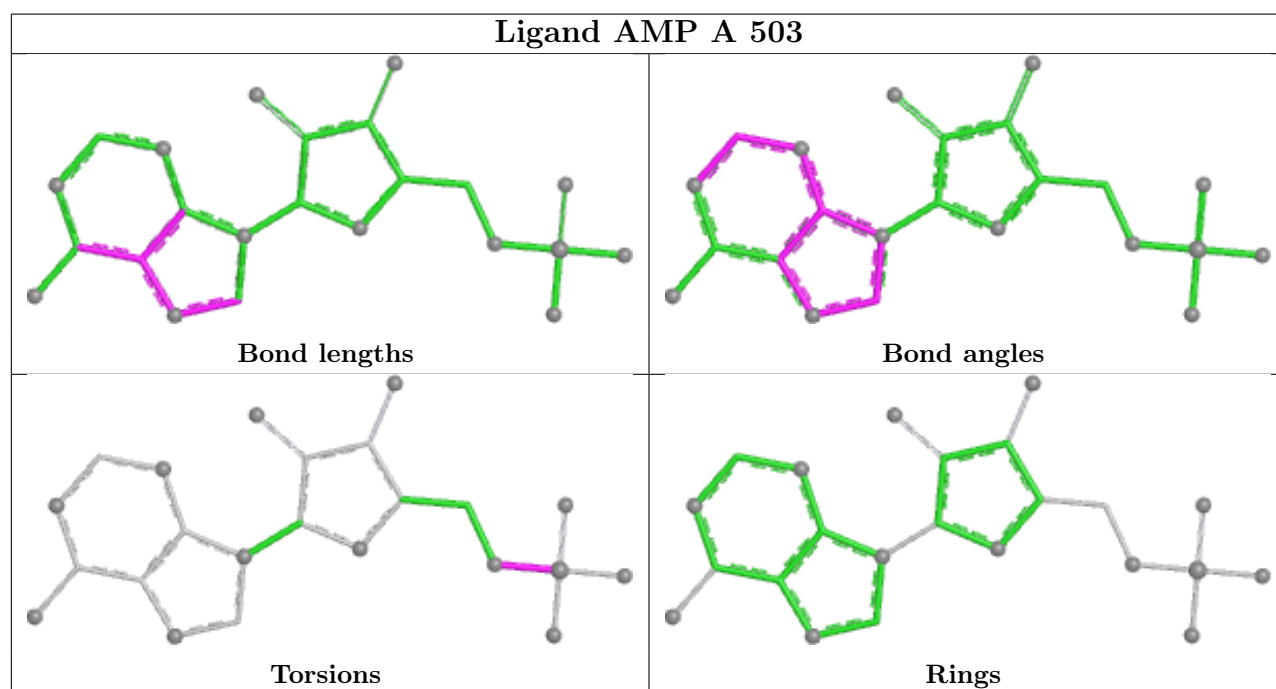
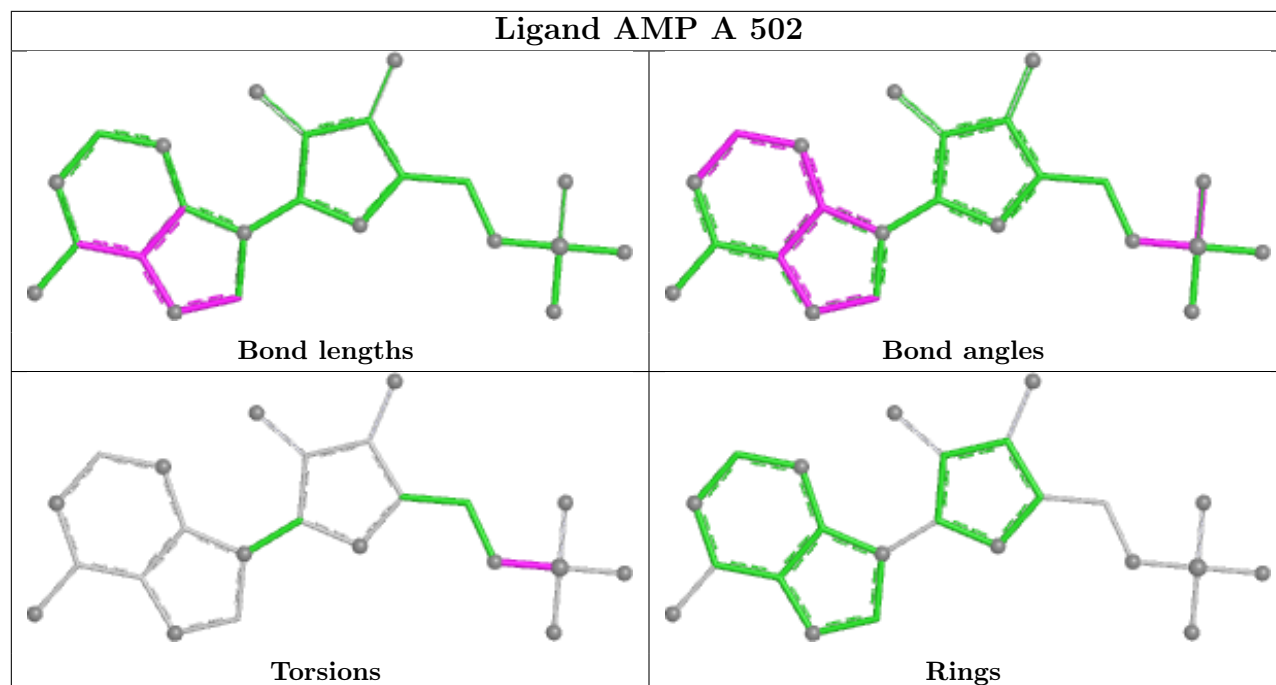
There are no ring outliers.

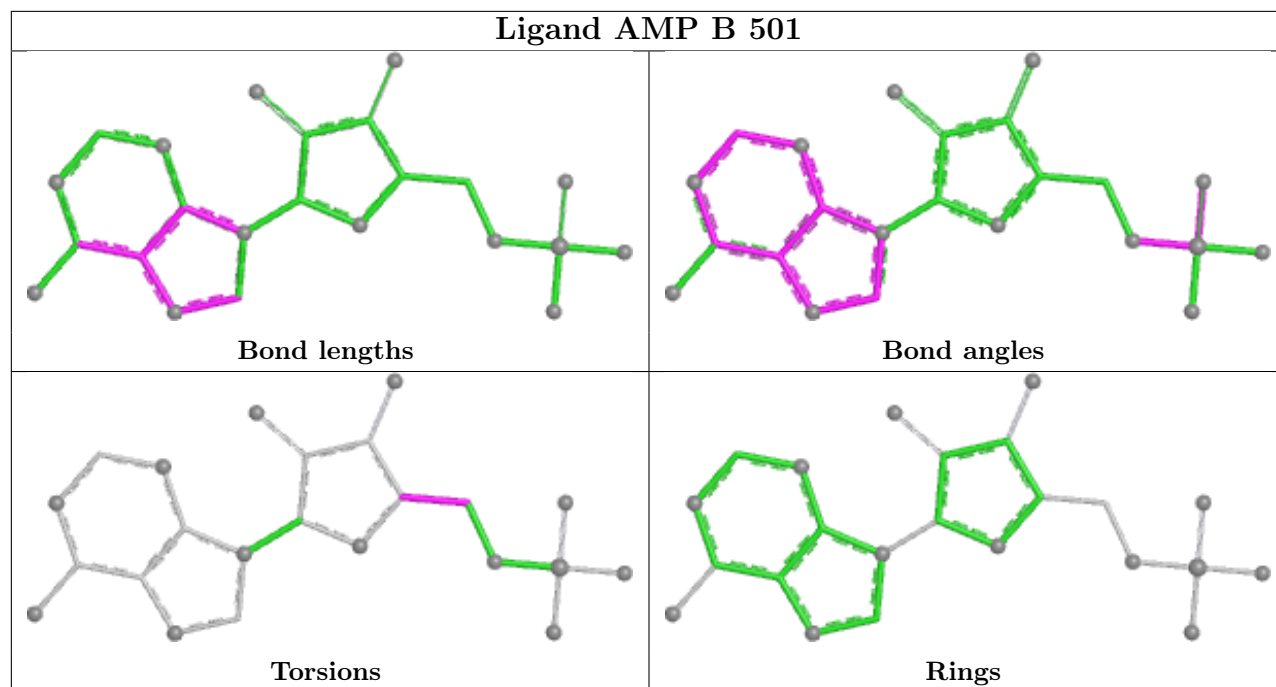
2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	AMP	3	0
2	B	501	AMP	3	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	315/393 (80%)	0.32	17 (5%) 31 27	37, 69, 122, 144	0
1	B	305/393 (77%)	0.43	29 (9%) 14 11	39, 74, 137, 161	0
All	All	620/786 (78%)	0.38	46 (7%) 20 18	37, 71, 134, 161	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	219	GLY	10.5
1	A	221	THR	6.8
1	A	398	VAL	4.3
1	A	416	THR	4.2
1	B	337	PHE	3.8
1	B	253	GLY	3.7
1	A	405	LYS	3.6
1	B	338	GLY	3.5
1	B	218	LYS	3.1
1	A	139	PHE	3.1
1	B	222	VAL	2.8
1	A	402	GLN	2.8
1	A	354	HIS	2.8
1	B	291	ALA	2.8
1	B	335	ARG	2.8
1	B	254	GLU	2.8
1	A	304	ASP	2.7
1	B	236	PRO	2.7
1	B	358	PRO	2.7
1	B	413	ASN	2.6
1	B	251	HIS	2.6
1	B	292	GLU	2.6
1	B	252	VAL	2.6
1	B	372	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	44	SER	2.5
1	A	406	ARG	2.4
1	A	399	ASN	2.4
1	A	302	ALA	2.4
1	B	249	LYS	2.4
1	B	139	PHE	2.3
1	B	247	GLY	2.3
1	B	258	LYS	2.2
1	B	33	VAL	2.2
1	B	363	TYR	2.2
1	A	232	ALA	2.2
1	A	228	ASP	2.2
1	B	360	LYS	2.2
1	B	248	GLN	2.2
1	A	401	GLU	2.1
1	B	406	ARG	2.1
1	B	399	ASN	2.1
1	A	252	VAL	2.1
1	B	235	LEU	2.1
1	B	416	THR	2.0
1	A	372	THR	2.0
1	A	176	GLY	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
2	AMP	A	501	23/23	0.79	0.13	64,140,183,194	0

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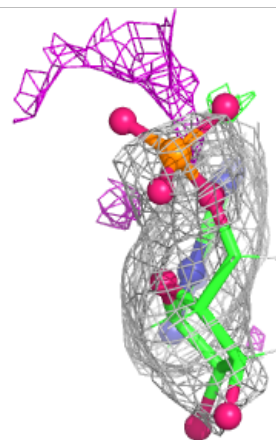
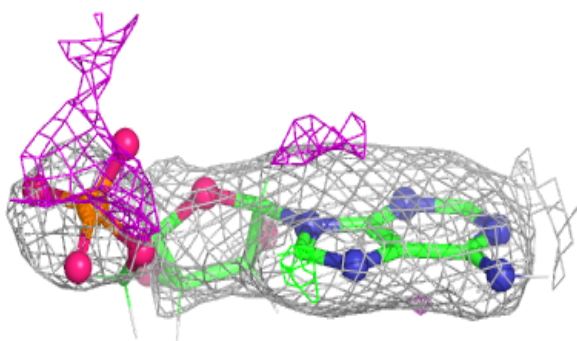
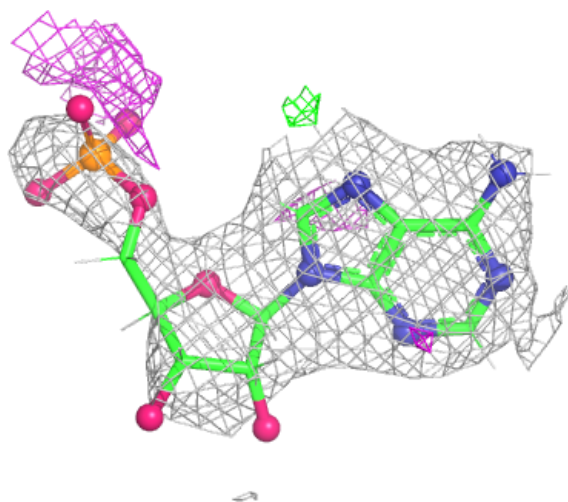
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	AMP	B	501	23/23	0.81	0.12	118,147,189,194	0
2	AMP	A	502	23/23	0.94	0.09	53,69,83,91	0
2	AMP	A	503	23/23	0.95	0.07	52,71,87,89	0

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.

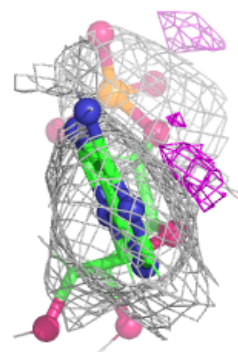
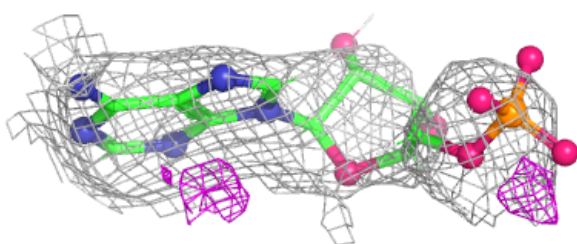
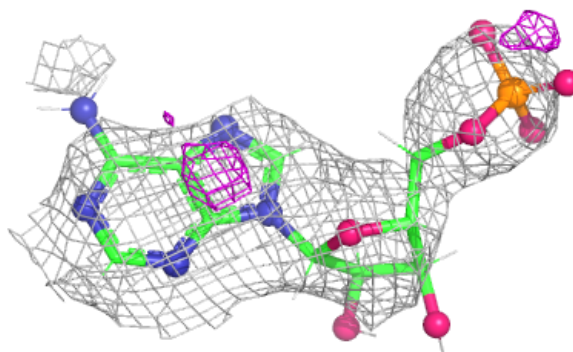
Electron density around AMP A 501:

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



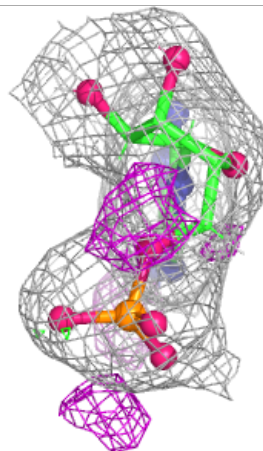
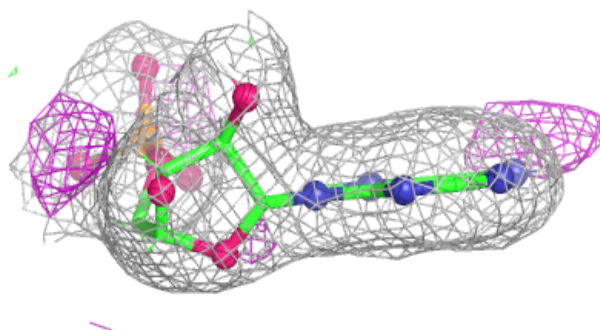
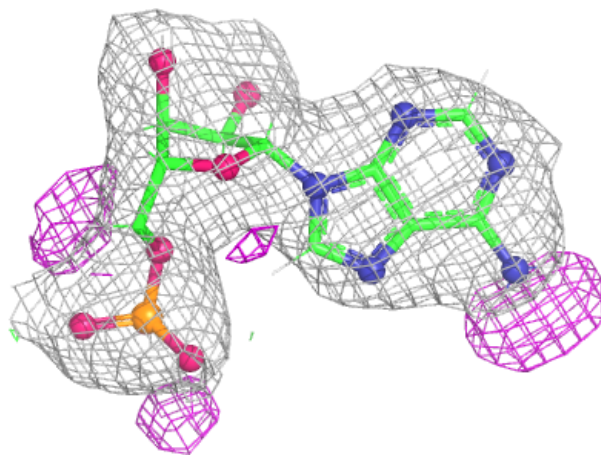
Electron density around AMP B 501:

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



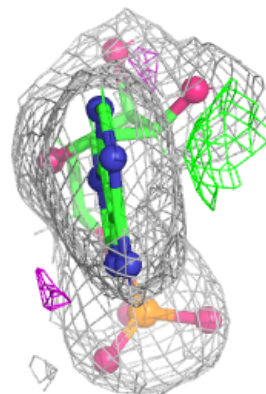
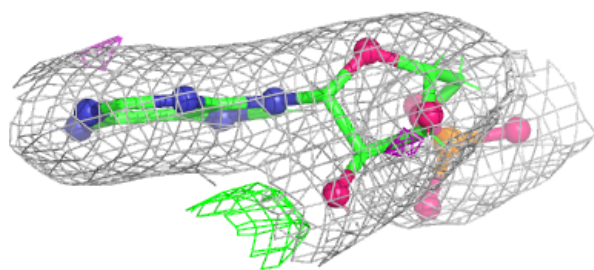
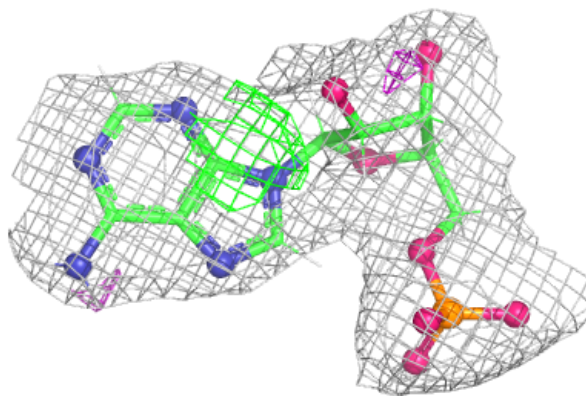
Electron density around AMP A 502:

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



Electron density around AMP A 503:

$2mF_o-DF_c$ (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.