Full wwPDB X-ray Structure Validation Report (i)

Jun 11, 2024 – 11:21 PM EDT

PDB ID	:	1TUU
Title	:	Acetate Kinase crystallized with ATPgS
Authors	:	Gorrell, A.; Lawrence, S.H.; Ferry, J.G.
Deposited on	:	2004-06-25
Resolution	:	2.50 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 (i) 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 (i)) were used in the production of this report:

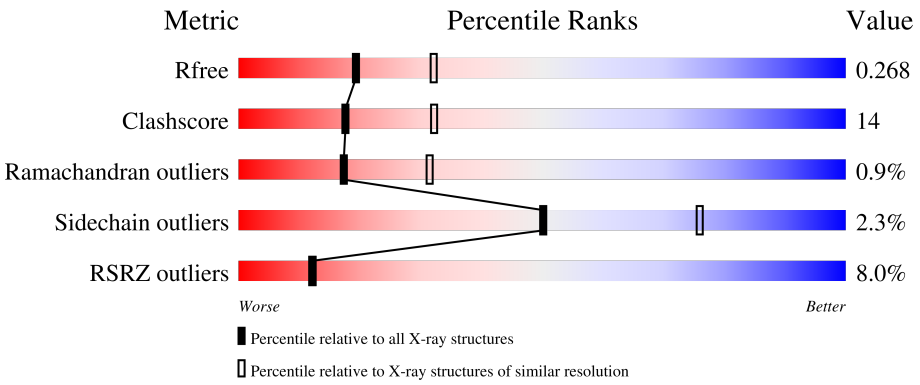
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 <=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	399	2% 73% 26% •
1	B	399	14% 67% 31% •

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
6	PIS	B	2854	-	-	X	-

2 Entry composition (i)

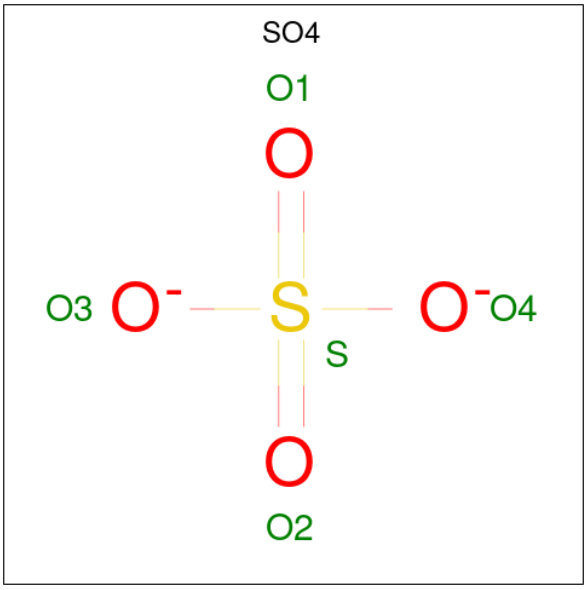
There are 7 unique types of molecules in this entry. The entry contains 6189 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Acetate kinase.

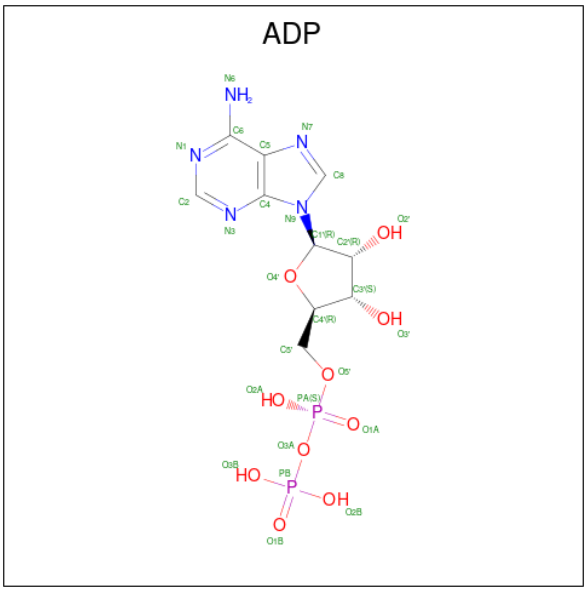
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	399	3038	1920	507	592	19	0	0	0
1	B	398	3028	1915	506	588	19	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



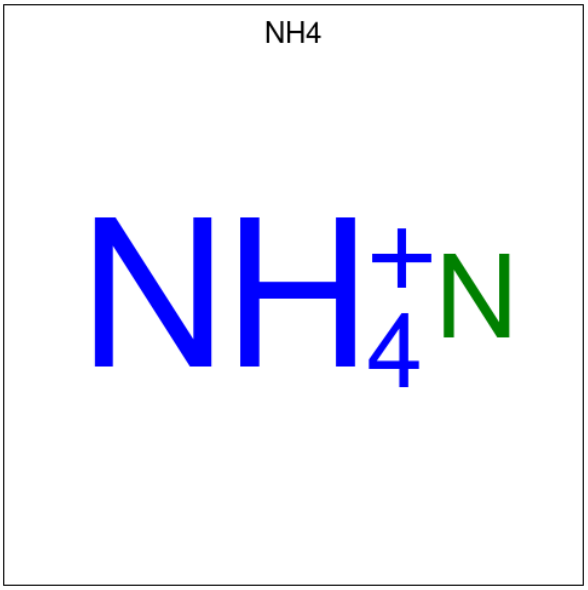
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



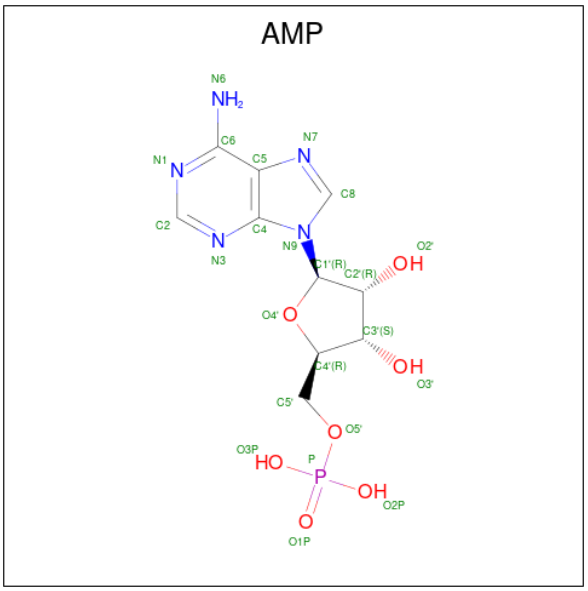
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	27	10	5	10	2	0	0

- Molecule 4 is AMMONIUM ION (three-letter code: NH4) (formula: H_4N).



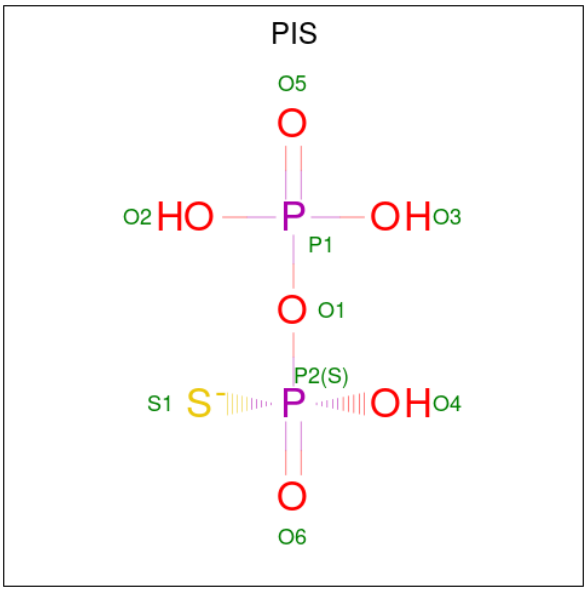
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	N		
4	B	1	1	1	0	0

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



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

- Molecule 6 is TRIHYDROGEN THIODIPHOSPHATE (three-letter code: PIS) (formula: $H_3O_6P_2S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	O	P	S		
6	B	1	9	6	2	1	0	0

- Molecule 7 is water.

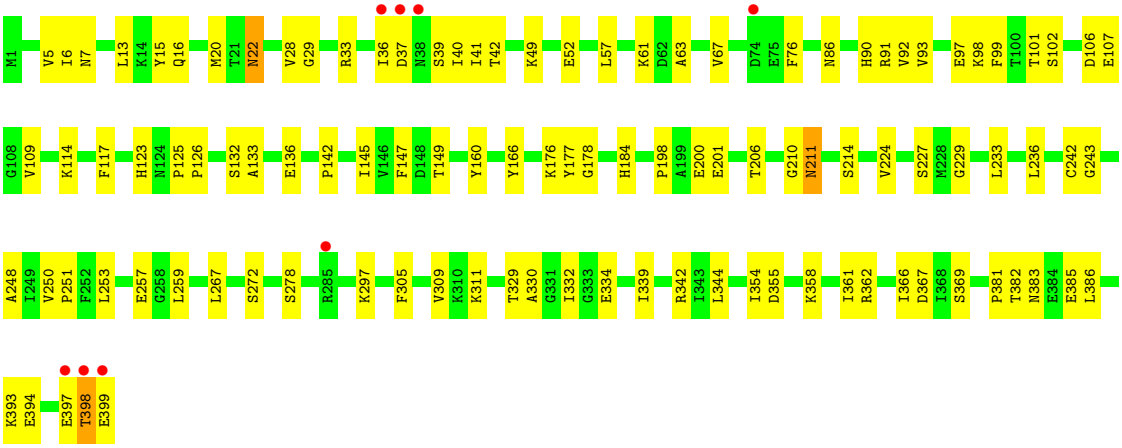
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	35	35	35	0	0
			Total	O		
7	B	23	23	23	0	0

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.

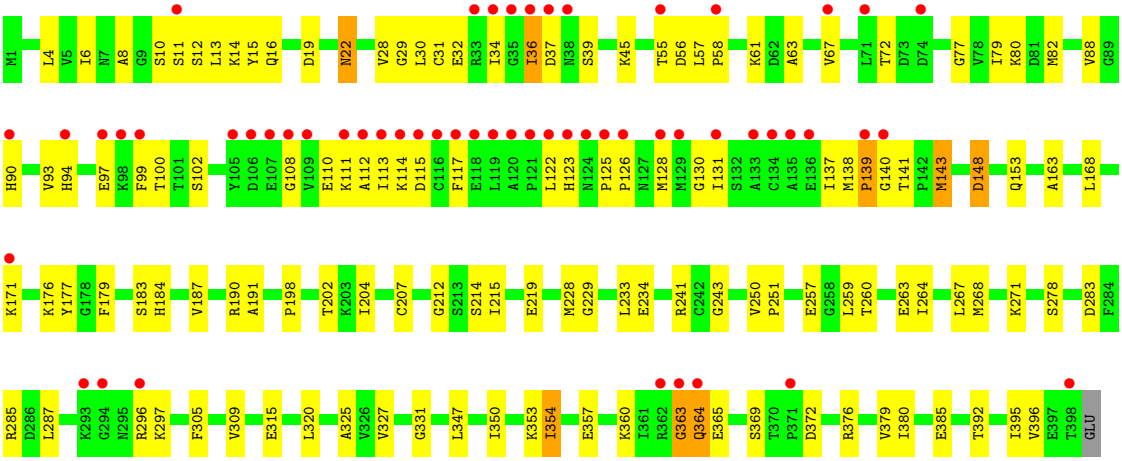
- Molecule 1: Acetate kinase

Chain A:



- Molecule 1: Acetate kinase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	179.08Å 66.11Å 81.67Å 90.00° 103.31° 90.00°	Depositor
Resolution (Å)	29.63 – 2.50 29.63 – 2.49	Depositor EDS
% Data completeness (in resolution range)	92.3 (29.63-2.50) 91.5 (29.63-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.17 (at 2.48Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.213 , 0.276 0.208 , 0.268	Depositor DCC
R_{free} test set	2084 reflections (6.91%)	wwPDB-VP
Wilson B-factor (Å^2)	34.8	Xtriage
Anisotropy	0.785	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 47.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6189	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: SO4, ADP, AMP, NH4, PIS

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	0.46	1/3086 (0.0%)	0.63	0/4159
1	B	0.41	0/3076	0.60	0/4147
All	All	0.44	1/6162 (0.0%)	0.62	0/8306

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	242	CYS	CB-SG	-10.77	1.64	1.82

There are no bond angle outliers.

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	3038	0	3077	79	0
1	B	3028	0	3071	89	0
2	A	5	0	0	0	0
3	A	27	0	12	2	0
4	B	1	0	0	0	0
5	B	23	0	12	2	0
6	B	9	0	0	6	0
7	A	35	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	23	0	0	0	0
All	All	6189	0	6172	168	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 14.

All (168) 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)
6:B:2854:PIS:S1	6:B:2854:PIS:P2	1.52	1.51
1:A:176:LYS:HG3	1:A:229:GLY:HA2	1.57	0.84
1:A:22:ASN:H	1:A:22:ASN:HD22	1.25	0.82
1:B:39:SER:HB3	1:B:56:ASP:HA	1.63	0.80
1:B:97:GLU:HG3	1:B:177:TYR:HE1	1.46	0.80
1:A:6:ILE:HG23	1:A:13:LEU:HD11	1.63	0.80
1:B:16:GLN:HG2	1:B:28:VAL:HG22	1.64	0.79
1:B:36:ILE:HG22	1:B:37:ASP:H	1.49	0.78
1:B:331:GLY:HA3	5:B:2853:AMP:H5'2	1.66	0.77
1:B:176:LYS:HG3	1:B:229:GLY:HA2	1.67	0.76
1:A:332:ILE:HG12	3:A:401:ADP:H1'	1.70	0.72
1:B:100:THR:HA	1:B:153:GLN:HG3	1.69	0.72
1:A:257:GLU:HB2	1:A:259:LEU:HG	1.70	0.72
1:A:253:LEU:HD11	1:B:168:LEU:HD22	1.71	0.71
1:A:97:GLU:HG3	1:A:177:TYR:OH	1.90	0.71
1:A:123:HIS:O	1:A:126:PRO:HD2	1.91	0.69
6:B:2854:PIS:S1	6:B:2854:PIS:O6	2.48	0.69
1:B:16:GLN:NE2	1:B:385:GLU:HG2	2.06	0.69
1:A:125:PRO:HB2	1:A:126:PRO:HD3	1.74	0.69
1:B:19:ASP:HB3	1:B:22:ASN:HD21	1.58	0.68
1:B:14:LYS:HG2	1:B:30:LEU:HD13	1.75	0.67
1:B:63:ALA:O	1:B:67:VAL:HG23	1.94	0.67
1:B:97:GLU:HG3	1:B:177:TYR:CE1	2.28	0.67
1:B:191:ALA:HB3	1:B:204:ILE:HD13	1.76	0.66
1:A:16:GLN:HG3	1:A:28:VAL:HG22	1.75	0.66
6:B:2854:PIS:S1	6:B:2854:PIS:O1	2.54	0.65
1:A:394:GLU:O	1:A:398:THR:HG22	1.97	0.64
1:A:16:GLN:NE2	1:A:385:GLU:HG2	2.13	0.64
1:B:45:LYS:HD2	1:B:77:GLY:O	1.98	0.63
1:A:132:SER:O	1:A:136:GLU:HG3	1.99	0.62
1:A:123:HIS:C	1:A:126:PRO:HD2	2.19	0.62
1:B:212:GLY:HA3	6:B:2854:PIS:S1	2.40	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:278:SER:HA	1:B:297:LYS:HB3	1.82	0.61
1:B:241:ARG:NH1	6:B:2854:PIS:S1	2.69	0.61
1:A:358:LYS:HB2	1:A:358:LYS:NZ	2.15	0.61
1:B:94:HIS:HB2	1:B:122:LEU:CD2	2.30	0.61
1:A:339:ILE:HA	1:A:342:ARG:HD2	1.82	0.61
1:A:36:ILE:HG22	1:A:37:ASP:H	1.65	0.60
1:B:357:GLU:O	1:B:360:LYS:HG3	2.01	0.59
6:B:2854:PIS:S1	6:B:2854:PIS:O4	2.50	0.59
1:A:114:LYS:HA	1:A:117:PHE:CE1	2.38	0.58
1:B:183:SER:O	1:B:187:VAL:HG22	2.03	0.58
1:B:347:LEU:HD22	1:B:350:ILE:HD11	1.84	0.58
1:B:283:ASP:O	1:B:287:LEU:HG	2.04	0.58
1:A:184:HIS:CD2	1:A:206:THR:OG1	2.56	0.57
1:A:39:SER:CB	1:A:57:LEU:H	2.17	0.57
1:A:184:HIS:HD2	1:A:206:THR:OG1	1.87	0.57
1:B:8:ALA:HB2	1:B:130:GLY:HA3	1.87	0.57
1:B:117:PHE:HE1	1:B:126:PRO:HD2	1.69	0.57
1:B:57:LEU:HD22	1:B:63:ALA:HA	1.86	0.57
1:B:260:THR:OG1	1:B:263:GLU:HG3	2.05	0.57
1:B:22:ASN:ND2	1:B:22:ASN:H	2.02	0.56
1:B:184:HIS:HE1	1:B:214:SER:OG	1.88	0.56
1:A:22:ASN:H	1:A:22:ASN:ND2	1.99	0.56
1:A:61:LYS:HD2	1:A:136:GLU:OE1	2.05	0.56
1:B:392:THR:O	1:B:396:VAL:HG23	2.05	0.56
1:B:354:ILE:HA	1:B:369:SER:HA	1.87	0.56
1:B:117:PHE:CE1	1:B:125:PRO:HB3	2.41	0.55
1:A:383:ASN:ND2	1:A:386:LEU:HB2	2.20	0.55
1:B:365:GLU:HA	1:B:379:VAL:O	2.07	0.55
1:A:114:LYS:HA	1:A:117:PHE:CD1	2.41	0.54
1:A:329:THR:OG1	1:A:382:THR:HB	2.07	0.54
1:B:97:GLU:O	1:B:99:PHE:N	2.40	0.54
1:B:93:VAL:HA	1:B:148:ASP:HB2	1.90	0.53
1:A:166:TYR:CD2	1:B:297:LYS:HE2	2.44	0.53
1:B:22:ASN:H	1:B:22:ASN:HD22	1.55	0.52
1:A:311:LYS:HB3	1:B:315:GLU:HG3	1.90	0.52
1:B:39:SER:OG	1:B:57:LEU:HG	2.10	0.52
1:A:63:ALA:O	1:A:67:VAL:HG23	2.10	0.52
1:B:139:PRO:HG2	1:B:140:GLY:H	1.74	0.52
1:B:207:CYS:SG	1:B:215:ILE:HD12	2.51	0.51
1:A:267:LEU:C	1:A:267:LEU:HD23	2.31	0.51
1:B:61:LYS:HZ2	1:B:137:ILE:CD1	2.22	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:ALA:CB	1:B:204:ILE:HD13	2.39	0.51
1:A:98:LYS:HG3	1:A:99:PHE:CD2	2.46	0.51
1:A:358:LYS:HB2	1:A:358:LYS:HZ2	1.75	0.51
1:B:15:TYR:CE1	1:B:29:GLY:HA3	2.46	0.50
1:A:7:ASN:OD1	1:A:91:ARG:HD2	2.12	0.50
1:B:99:PHE:CZ	1:B:112:ALA:HB3	2.46	0.50
1:B:102:SER:HB2	1:B:395:ILE:HG13	1.91	0.50
1:B:128:MET:SD	1:B:131:ILE:HD12	2.52	0.50
1:A:6:ILE:CG2	1:A:13:LEU:HD11	2.39	0.50
1:A:22:ASN:HD22	1:A:22:ASN:N	1.93	0.50
1:A:86:ASN:O	1:A:142:PRO:HD2	2.12	0.50
1:A:354:ILE:HA	1:A:369:SER:HA	1.92	0.50
1:A:15:TYR:CZ	1:A:29:GLY:HA3	2.46	0.50
1:A:397:GLU:C	1:A:399:GLU:H	2.15	0.50
1:B:331:GLY:HA3	5:B:2853:AMP:C5'	2.40	0.50
1:A:106:ASP:OD1	1:A:109:VAL:HG23	2.12	0.49
1:B:61:LYS:HZ2	1:B:137:ILE:HD13	1.76	0.49
1:B:36:ILE:HG22	1:B:37:ASP:N	2.24	0.49
1:B:110:GLU:O	1:B:114:LYS:HG3	2.11	0.49
1:A:278:SER:HA	1:A:297:LYS:HB3	1.94	0.49
1:A:125:PRO:CB	1:A:126:PRO:HD3	2.43	0.49
1:A:233:LEU:O	1:A:243:GLY:HA2	2.13	0.49
1:B:187:VAL:HB	1:B:327:VAL:HG11	1.95	0.48
1:B:257:GLU:HB2	1:B:259:LEU:HG	1.94	0.48
1:A:250:VAL:HB	1:A:251:PRO:CD	2.43	0.48
1:A:33:ARG:HH11	1:A:33:ARG:HG2	1.78	0.48
1:A:149:THR:HG22	1:A:178:GLY:O	2.13	0.48
1:B:285:ARG:HH11	1:B:285:ARG:HG2	1.78	0.48
1:B:11:SER:HB2	1:B:34:ILE:O	2.14	0.48
1:B:6:ILE:HB	1:B:90:HIS:CD2	2.48	0.47
1:B:267:LEU:HA	1:B:271:LYS:HE3	1.94	0.47
1:B:228:MET:HA	1:B:234:GLU:O	2.14	0.47
1:B:233:LEU:O	1:B:243:GLY:HA2	2.13	0.47
1:A:22:ASN:ND2	1:A:22:ASN:N	2.61	0.47
1:A:102:SER:HA	1:A:145:ILE:O	2.15	0.47
1:B:30:LEU:HG	1:B:32:GLU:HG3	1.96	0.47
1:A:334:GLU:HG2	1:A:381:PRO:HA	1.96	0.47
1:A:40:ILE:HG12	1:A:41:ILE:N	2.30	0.46
1:A:224:VAL:O	1:A:224:VAL:HG12	2.15	0.46
1:B:108:GLY:HA2	1:B:111:LYS:HE3	1.98	0.46
1:A:224:VAL:HG12	7:A:402:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:ILE:CG1	3:A:401:ADP:H1'	2.42	0.46
1:A:236:LEU:HD21	1:A:309:VAL:HA	1.97	0.46
1:A:101:THR:O	1:A:102:SER:C	2.54	0.46
1:B:365:GLU:OE2	1:B:380:ILE:HG23	2.16	0.46
1:B:82:MET:O	1:B:82:MET:HG2	2.16	0.45
1:B:72:THR:HA	1:B:79:ILE:O	2.16	0.45
1:A:5:VAL:HG11	1:A:385:GLU:HA	1.99	0.45
1:A:361:ILE:HD12	1:A:366:ILE:HD13	1.98	0.45
1:B:250:VAL:HB	1:B:251:PRO:CD	2.46	0.45
1:A:160:TYR:HA	1:A:176:LYS:HB3	1.98	0.45
1:A:90:HIS:HB2	1:A:145:ILE:HD13	1.99	0.44
1:A:198:PRO:HD2	1:A:201:GLU:OE1	2.16	0.44
1:A:366:ILE:HG12	1:A:367:ASP:N	2.32	0.44
1:B:363:GLY:O	1:B:364:GLN:HB2	2.16	0.44
1:A:20:MET:HB3	1:A:393:LYS:HE2	1.99	0.44
1:B:6:ILE:HG23	1:B:13:LEU:HD21	1.99	0.44
1:A:383:ASN:HD21	1:A:386:LEU:HB2	1.81	0.44
1:A:329:THR:O	1:A:330:ALA:HB3	2.17	0.44
1:B:22:ASN:HD22	1:B:22:ASN:C	2.21	0.44
1:B:88:VAL:HB	1:B:143:MET:HG2	2.00	0.44
1:B:94:HIS:HB2	1:B:122:LEU:HD22	2.00	0.44
1:B:325:ALA:HA	1:B:376:ARG:O	2.17	0.43
1:B:39:SER:CB	1:B:57:LEU:H	2.32	0.43
1:A:305:PHE:O	1:A:309:VAL:HG23	2.17	0.43
1:A:40:ILE:HD11	1:A:52:GLU:HG3	2.01	0.43
1:B:12:SER:HA	1:B:31:CYS:O	2.19	0.43
1:B:198:PRO:O	1:B:202:THR:HG23	2.19	0.43
1:A:210:GLY:O	1:A:211:ASN:C	2.57	0.43
1:B:278:SER:HA	1:B:297:LYS:CB	2.48	0.42
1:A:355:ASP:OD2	1:A:355:ASP:C	2.57	0.42
1:B:4:LEU:HG	1:B:6:ILE:HD11	2.01	0.42
1:B:264:ILE:O	1:B:268:MET:HG3	2.18	0.42
1:B:219:GLU:HB2	1:B:320:LEU:CD2	2.49	0.42
1:B:36:ILE:CG2	1:B:37:ASP:H	2.22	0.42
1:A:250:VAL:HB	1:A:251:PRO:HD3	2.01	0.42
1:A:33:ARG:HG2	1:A:33:ARG:NH1	2.35	0.41
1:B:171:LYS:O	1:B:171:LYS:HG2	2.20	0.41
1:B:190:ARG:HB2	1:B:380:ILE:HD13	2.03	0.41
1:A:16:GLN:CD	1:A:385:GLU:HG2	2.41	0.41
1:A:214:SER:HA	1:A:227:SER:OG	2.21	0.41
1:B:15:TYR:CZ	1:B:29:GLY:HA3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:PHE:O	1:B:309:VAL:HG23	2.20	0.41
1:A:92:VAL:O	1:A:147:PHE:HA	2.20	0.41
1:A:344:LEU:HD13	1:A:354:ILE:HG21	2.02	0.41
1:B:353:LYS:HE2	1:B:372:ASP:OD2	2.20	0.41
1:A:248:ALA:O	1:A:251:PRO:HG2	2.21	0.41
1:A:93:VAL:HG21	1:A:123:HIS:CB	2.51	0.40
1:A:49:LYS:HB3	1:A:76:PHE:CD2	2.56	0.40
1:A:272:SER:HB3	1:B:163:ALA:O	2.21	0.40
1:A:61:LYS:HA	1:A:133:ALA:HB1	2.03	0.40
1:B:113:ILE:HG21	1:B:128:MET:SD	2.61	0.40
1:B:285:ARG:HG2	1:B:285:ARG:NH1	2.36	0.40
1:B:138:MET:HB3	1:B:141:THR:OG1	2.22	0.40
1:B:179:PHE:CD1	1:B:179:PHE:N	2.88	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	397/399 (100%)	382 (96%)	13 (3%)	2 (0%)	29 48
1	B	396/399 (99%)	358 (90%)	33 (8%)	5 (1%)	12 21
All	All	793/798 (99%)	740 (93%)	46 (6%)	7 (1%)	17 31

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	362	ARG
1	B	58	PRO
1	B	139	PRO
1	B	363	GLY
1	A	398	THR

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Mol	Chain	Res	Type
1	B	364	GLN
1	B	36	ILE

5.3.2 Protein sidechains (i)

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	328/328 (100%)	323 (98%)	5 (2%)	65	85
1	B	327/328 (100%)	317 (97%)	10 (3%)	40	67
All	All	655/656 (100%)	640 (98%)	15 (2%)	50	76

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	42	THR
1	A	107	GLU
1	A	200	GLU
1	A	211	ASN
1	B	10	SER
1	B	22	ASN
1	B	55	THR
1	B	80	LYS
1	B	115	ASP
1	B	123	HIS
1	B	143	MET
1	B	148	ASP
1	B	296	ARG
1	B	354	ILE

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	16	GLN

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Mol	Chain	Res	Type
1	A	22	ASN
1	A	127	ASN
1	A	184	HIS
1	A	364	GLN
1	A	383	ASN
1	B	22	ASN
1	B	90	HIS
1	B	94	HIS
1	B	123	HIS
1	B	124	ASN
1	B	184	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 1 is modelled with single atom - leaving 4 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	SO4	A	400	-	4,4,4	0.19	0	6,6,6	0.17	0
3	ADP	A	401	-	24,29,29	1.60	5 (20%)	29,45,45	1.15	2 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
6	PIS	B	2854	-	5,8,8	2.54	3 (60%)	8,13,13	4.56	6 (75%)
5	AMP	B	2853	-	21,25,25	1.79	4 (19%)	23,38,38	1.66	4 (17%)

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
6	PIS	B	2854	-	-	0/4/6/6	-
3	ADP	A	401	-	-	4/12/32/32	0/3/3/3
5	AMP	B	2853	-	-	1/6/26/26	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	2853	AMP	C3'-C4'	5.74	1.67	1.53
6	B	2854	PIS	P2-O6	4.02	1.55	1.46
3	A	401	ADP	PA-O3A	4.00	1.63	1.59
3	A	401	ADP	O4'-C1'	3.32	1.45	1.40
6	B	2854	PIS	P2-O4	-2.98	1.49	1.56
5	B	2853	AMP	O3'-C3'	2.97	1.50	1.43
3	A	401	ADP	C2'-C3'	2.93	1.61	1.53
5	B	2853	AMP	C2-N3	2.61	1.36	1.32
3	A	401	ADP	C8-N7	-2.55	1.30	1.34
5	B	2853	AMP	C2'-C3'	2.48	1.60	1.53
6	B	2854	PIS	P1-O2	-2.41	1.45	1.54
3	A	401	ADP	PA-O2A	-2.08	1.45	1.55

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	B	2854	PIS	O2-P1-O5	-6.53	85.38	110.83
6	B	2854	PIS	O1-P1-O5	-6.27	78.06	111.04
6	B	2854	PIS	O3-P1-O5	-6.07	87.19	110.83
6	B	2854	PIS	O2-P1-O1	5.06	121.62	104.64
5	B	2853	AMP	C4-C5-N7	4.13	113.71	109.34
5	B	2853	AMP	C4'-O4'-C1'	3.40	113.04	109.92
6	B	2854	PIS	O3-P1-O1	3.19	115.32	104.64
6	B	2854	PIS	O3-P1-O2	3.06	119.29	107.80
5	B	2853	AMP	N6-C6-N1	2.97	124.67	118.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	2853	AMP	N3-C2-N1	-2.88	124.76	128.67
3	A	401	ADP	C4'-O4'-C1'	2.51	112.23	109.92
3	A	401	ADP	O4'-C4'-C3'	-2.01	101.16	105.15

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	ADP	PA-O3A-PB-O2B
3	A	401	ADP	O4'-C4'-C5'-O5'
3	A	401	ADP	C3'-C4'-C5'-O5'
3	A	401	ADP	PA-O3A-PB-O3B
5	B	2853	AMP	O4'-C4'-C5'-O5'

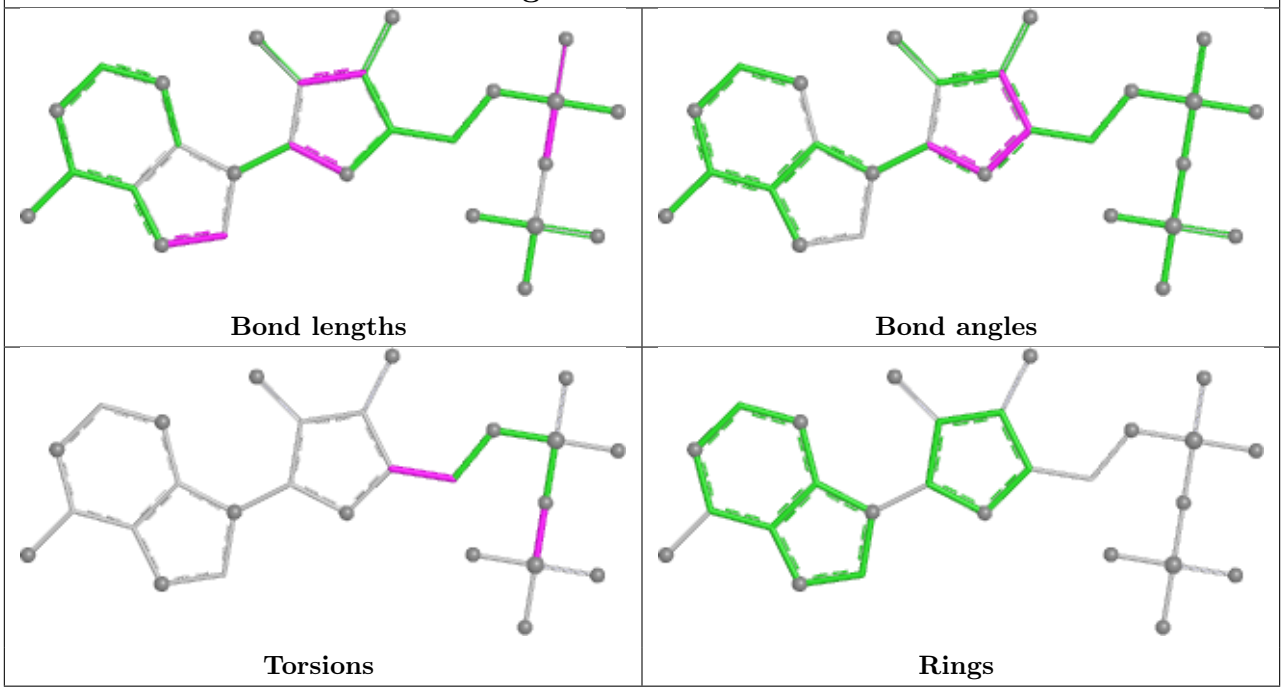
There are no ring outliers.

3 monomers are involved in 10 short contacts:

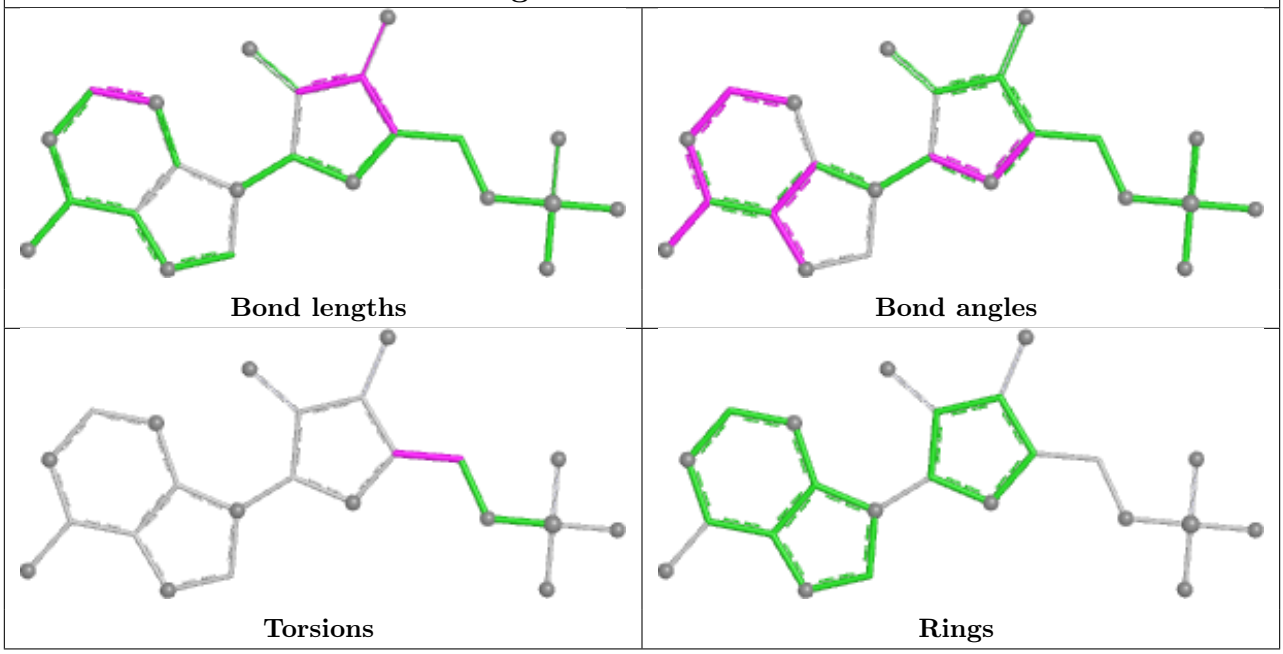
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	ADP	2	0
6	B	2854	PIS	6	0
5	B	2853	AMP	2	0

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 then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand ADP A 401



Ligand AMP B 2853



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	399/399 (100%)	-0.06	8 (2%)	65	68	20, 34, 49, 62	0
1	B	398/399 (99%)	0.57	56 (14%)	2	2	21, 40, 72, 78	0
All	All	797/798 (99%)	0.25	64 (8%)	12	12	20, 36, 65, 78	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	117	PHE	6.7
1	B	98	LYS	6.7
1	B	107	GLU	5.9
1	B	363	GLY	5.8
1	B	111	LYS	5.7
1	B	123	HIS	5.5
1	B	122	LEU	5.2
1	B	118	GLU	5.0
1	B	119	LEU	5.0
1	B	124	ASN	5.0
1	B	36	ILE	4.7
1	B	125	PRO	4.6
1	B	58	PRO	4.5
1	B	109	VAL	4.4
1	B	128	MET	4.3
1	B	126	PRO	4.3
1	B	136	GLU	4.0
1	B	34	ILE	4.0
1	A	37	ASP	3.9
1	B	115	ASP	3.9
1	A	74	ASP	3.7
1	B	134	CYS	3.7
1	B	37	ASP	3.7
1	A	398	THR	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	36	ILE	3.7
1	B	108	GLY	3.6
1	B	99	PHE	3.6
1	B	112	ALA	3.6
1	B	90	HIS	3.5
1	B	131	ILE	3.5
1	B	135	ALA	3.4
1	B	116	CYS	3.3
1	B	371	PRO	3.2
1	B	114	LYS	3.1
1	B	105	TYR	3.1
1	B	362	ARG	3.0
1	B	140	GLY	2.9
1	B	106	ASP	2.8
1	A	38	ASN	2.8
1	B	129	MET	2.7
1	B	294	GLY	2.6
1	B	139	PRO	2.6
1	A	399	GLU	2.5
1	B	293	LYS	2.5
1	B	296	ARG	2.5
1	B	33	ARG	2.4
1	B	113	ILE	2.4
1	B	94	HIS	2.4
1	A	397	GLU	2.4
1	B	11	SER	2.4
1	B	121	PRO	2.3
1	A	285	ARG	2.3
1	B	67	VAL	2.3
1	B	35	GLY	2.3
1	B	171	LYS	2.3
1	B	398	THR	2.3
1	B	38	ASN	2.2
1	B	97	GLU	2.2
1	B	71	LEU	2.2
1	B	74	ASP	2.1
1	B	120	ALA	2.1
1	B	364	GLN	2.1
1	B	55	THR	2.1
1	B	133	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 monosaccharides in this entry.

6.4 Ligands (i)

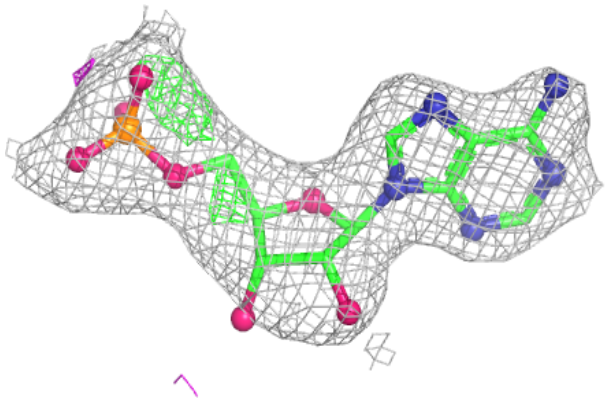
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

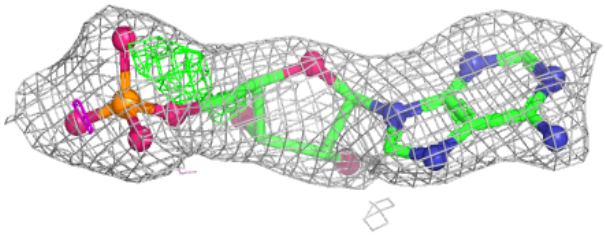
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	NH4	B	2852	1/1	0.88	0.60	36,36,36,36	0
5	AMP	B	2853	23/23	0.89	0.20	50,53,57,58	0
6	PIS	B	2854	9/9	0.90	0.19	83,84,85,86	0
3	ADP	A	401	27/27	0.93	0.18	39,49,51,55	0
2	SO4	A	400	5/5	0.99	0.17	34,36,37,38	0

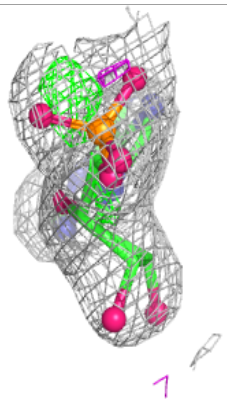
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around AMP B 2853:

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

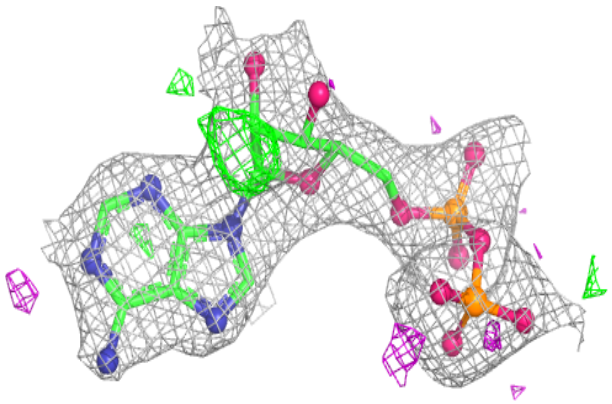


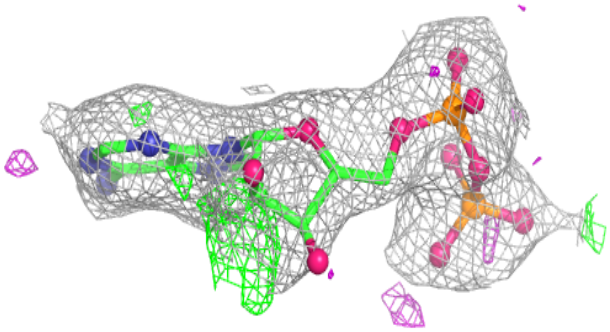


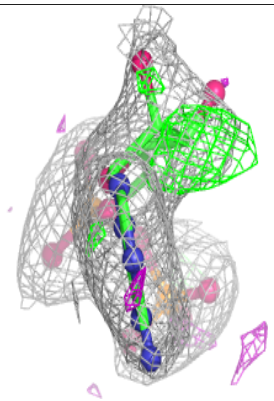


Electron density around ADP A 401:

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







6.5 Other polymers (i)

There are no such residues in this entry.