



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2021 – 05:13 PM EDT

PDB ID : 3ETH
Title : Crystal structure of E. coli Purk in complex with MgATP
Authors : Holden, H.M.; Thoden, J.B.
Deposited on : 2008-10-08
Resolution : 1.60 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
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) : 2.23.2

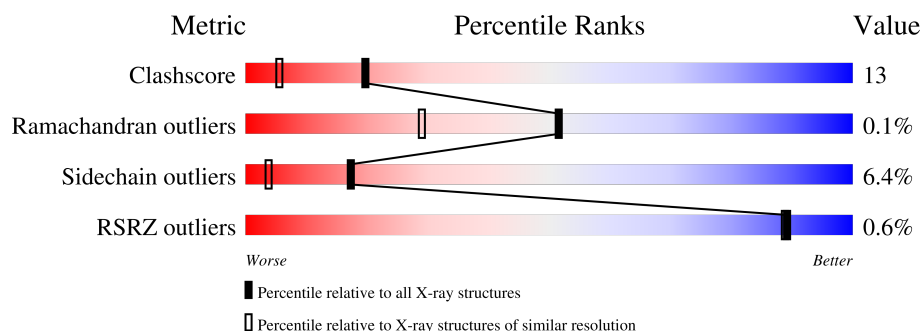
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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))
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	355	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>19%</div> <div>..</div> </div> </div>
1	B	355	<div> <div></div> <div>75%</div> <div>21%</div> <div>..</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6346 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.

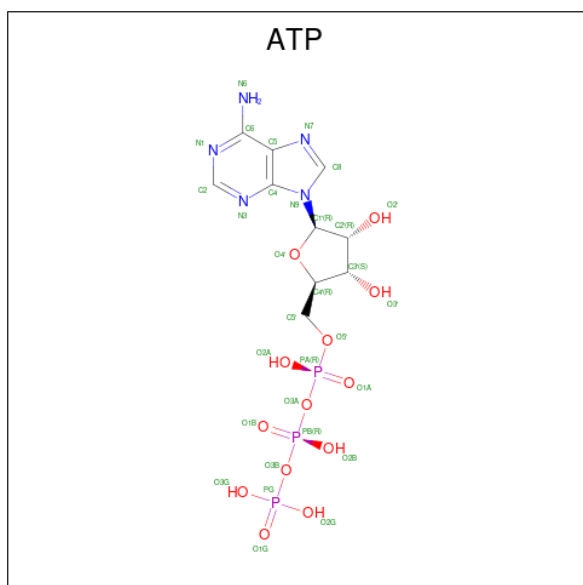
- Molecule 1 is a protein called Phosphoribosylaminoimidazole carboxylase ATPase subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	355	Total	C	N	O	S	0	4	0
			2802	1782	498	511	11			
1	B	355	Total	C	N	O	S	0	1	0
			2788	1774	495	508	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	GLN	GLU	engineered mutation	UNP P09029
A	205	ARG	GLN	engineered mutation	UNP P09029
B	61	GLN	GLU	engineered mutation	UNP P09029
B	205	ARG	GLN	engineered mutation	UNP P09029

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		
3	B	2	Total	Mg	0	0
			2	2		

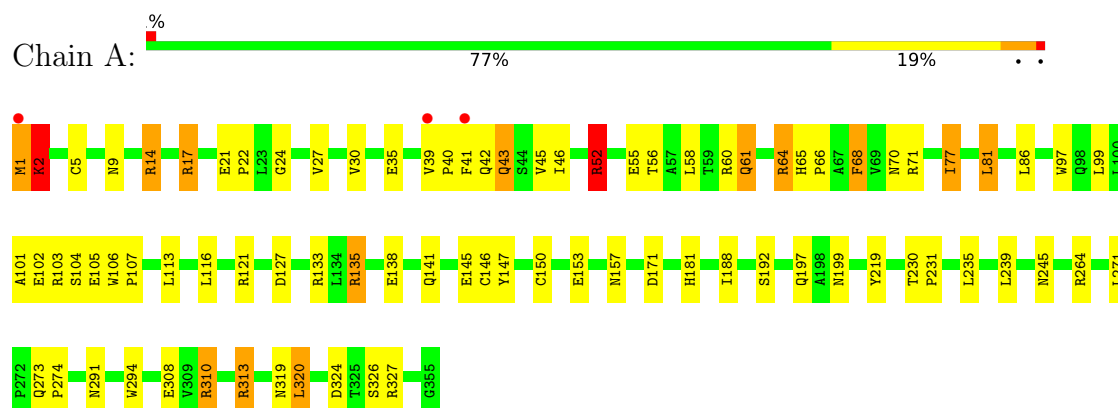
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	375	Total	O	0	0
			375	375		
4	B	315	Total	O	0	0
			315	315		

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: Phosphoribosylaminoimidazole carboxylase ATPase subunit



- Molecule 1: Phosphoribosylaminoimidazole carboxylase ATPase subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.10Å 57.11Å 59.22Å 77.46° 82.39° 77.26°	Depositor
Resolution (Å)	30.00 – 1.60 39.17 – 1.60	Depositor EDS
% Data completeness (in resolution range)	93.8 (30.00-1.60) 93.8 (39.17-1.60)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 1.60Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.175 , 0.228 0.179 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	48.9	Xtriage
Anisotropy	0.704	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 83.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.014 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6346	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.94% 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: ATP, MG

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	0.73	0/2888	1.33	19/3934 (0.5%)
1	B	0.72	0/2862	1.35	24/3900 (0.6%)
All	All	0.73	0/5750	1.34	43/7834 (0.5%)

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	327	ARG	NE-CZ-NH2	-13.09	113.75	120.30
1	B	313	ARG	NE-CZ-NH1	-12.42	114.09	120.30
1	A	14	ARG	NE-CZ-NH1	-11.34	114.63	120.30
1	B	17	ARG	NE-CZ-NH2	-9.71	115.45	120.30
1	A	14	ARG	NE-CZ-NH2	9.52	125.06	120.30
1	B	168	ARG	NE-CZ-NH2	7.98	124.29	120.30
1	B	320	LEU	CB-CG-CD2	-7.62	98.04	111.00
1	B	134	LEU	CB-CG-CD1	7.62	123.95	111.00
1	B	327	ARG	NE-CZ-NH1	7.27	123.94	120.30
1	B	92	LEU	CB-CG-CD2	-7.08	98.96	111.00
1	B	306	ASP	CB-CG-OD2	-7.07	111.94	118.30
1	A	17	ARG	NE-CZ-NH2	-7.05	116.78	120.30
1	A	327	ARG	NE-CZ-NH2	-6.78	116.91	120.30
1	B	320	LEU	CA-CB-CG	6.74	130.81	115.30
1	B	205	ARG	NE-CZ-NH2	6.64	123.62	120.30
1	B	168	ARG	NE-CZ-NH1	-6.42	117.09	120.30
1	B	112	ARG	NE-CZ-NH1	-6.39	117.10	120.30
1	A	271	LEU	CB-CG-CD2	-6.36	100.19	111.00
1	A	2	LYS	CB-CA-C	-6.33	97.74	110.40
1	A	135[A]	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	A	135[B]	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	B	210	LEU	CB-CG-CD1	6.23	121.59	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	153	GLU	CB-CA-C	-6.19	98.02	110.40
1	B	313	ARG	NE-CZ-NH2	6.18	123.39	120.30
1	A	1	MET	CG-SD-CE	6.02	109.83	100.20
1	B	229	VAL	CG1-CB-CG2	-5.98	101.33	110.90
1	B	33	ASP	CB-CG-OD2	5.94	123.65	118.30
1	B	122	ARG	NE-CZ-NH1	-5.88	117.36	120.30
1	B	236	ILE	CB-CA-C	-5.75	100.10	111.60
1	A	264	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	A	264	ARG	NE-CZ-NH2	-5.57	117.52	120.30
1	A	81	LEU	CB-CG-CD1	-5.46	101.71	111.00
1	A	135[A]	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	A	135[B]	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	B	327	ARG	CG-CD-NE	-5.28	100.71	111.80
1	B	306	ASP	CB-CG-OD1	5.27	123.05	118.30
1	B	44	SER	CB-CA-C	-5.25	100.13	110.10
1	A	113	LEU	CB-CG-CD2	5.18	119.80	111.00
1	A	45	VAL	CB-CA-C	-5.12	101.68	111.40
1	B	2	LYS	N-CA-C	5.11	124.79	111.00
1	A	52	ARG	NE-CZ-NH1	-5.08	117.76	120.30
1	B	339	LEU	CB-CG-CD1	-5.06	102.39	111.00
1	A	97	TRP	N-CA-CB	-5.06	101.49	110.60

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	2802	0	2779	87	0
1	B	2788	0	2767	56	0
2	A	31	0	12	0	0
2	B	31	0	12	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	375	0	0	10	0
4	B	315	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6346	0	5570	143	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 13.

All (143) 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:A:2:LYS:NZ	1:A:2:LYS:H	1.39	1.20
1:A:199:ASN:HB3	4:A:770:HOH:O	1.47	1.14
1:A:2:LYS:HZ3	1:A:2:LYS:N	1.61	0.98
1:A:310:ARG:HG3	1:A:313:ARG:HD3	1.47	0.93
1:A:2:LYS:H	1:A:2:LYS:HZ3	0.97	0.92
1:A:1:MET:HG3	1:A:24:GLY:O	1.71	0.90
1:A:146:CYS:HB3	1:A:150:CYS:SG	2.12	0.89
1:A:308:GLU:HB3	1:A:313:ARG:HH12	1.37	0.87
1:B:127:ASP:HB3	1:B:188:ILE:HG12	1.56	0.85
1:A:2:LYS:NZ	1:A:2:LYS:N	2.20	0.84
1:A:310:ARG:HD2	1:A:313:ARG:HD2	1.62	0.79
1:A:310:ARG:CG	1:A:313:ARG:HH11	2.00	0.74
1:A:310:ARG:CD	1:A:313:ARG:HD2	2.17	0.74
1:A:101:ALA:N	1:A:105:GLU:OE2	2.23	0.71
1:B:21:GLU:HB3	1:B:22:PRO:HD3	1.71	0.71
1:A:320:LEU:N	1:A:320:LEU:HD12	2.05	0.71
1:A:65:HIS:O	1:A:71:ARG:NH2	2.19	0.70
1:A:86:LEU:HA	4:A:524:HOH:O	1.93	0.69
1:A:273:GLN:NE2	4:A:467:HOH:O	2.28	0.67
1:B:52:ARG:HG3	4:B:445:HOH:O	1.93	0.67
1:A:135[A]:ARG:HD2	4:A:769:HOH:O	1.95	0.66
1:B:203:GLN:O	1:B:207:GLU:HG3	1.96	0.66
1:B:135:ARG:N	1:B:138:GLU:OE2	2.21	0.65
1:A:102:GLU:N	1:A:105:GLU:OE2	2.28	0.64
1:A:310:ARG:HG3	1:A:313:ARG:CD	2.26	0.63
1:A:52:ARG:N	1:A:52:ARG:HD2	2.14	0.62
1:A:116:LEU:HD11	1:A:133:ARG:HG2	1.81	0.61
1:B:52:ARG:HD2	1:B:53:TRP:N	2.15	0.61
1:A:77:ILE:CD1	1:A:239:LEU:HD12	2.31	0.60
1:A:230:THR:HG21	1:A:235:LEU:HD13	1.82	0.60
1:A:138:GLU:HB3	1:A:141:GLN:OE1	2.02	0.60
1:A:320:LEU:N	1:A:320:LEU:CD1	2.65	0.60
1:B:197:GLN:HG2	4:B:694:HOH:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:ASN:C	1:A:320:LEU:HD12	2.22	0.59
1:A:42:GLN:OE1	1:A:43:GLN:NE2	2.36	0.58
1:B:17:ARG:HG2	1:B:17:ARG:HH11	1.67	0.58
1:B:273:GLN:NE2	4:B:551:HOH:O	2.36	0.58
1:A:135[A]:ARG:NH1	4:A:769:HOH:O	2.35	0.57
1:A:46:ILE:HB	1:A:68:PHE:HA	1.86	0.57
1:A:310:ARG:HG2	1:A:313:ARG:HH11	1.69	0.57
1:A:308:GLU:HB3	1:A:313:ARG:NH1	2.14	0.57
1:A:1:MET:HG2	1:A:1:MET:O	2.03	0.57
1:A:1:MET:HE1	4:A:629:HOH:O	2.06	0.56
1:A:65:HIS:ND1	1:A:66:PRO:HD2	2.20	0.56
1:B:110:PHE:CE1	1:B:136:ALA:HA	2.40	0.56
1:B:52:ARG:HD2	1:B:52:ARG:C	2.26	0.56
1:B:2:LYS:CB	1:B:25:ILE:HG12	2.36	0.55
1:B:2:LYS:HB2	1:B:25:ILE:HA	1.89	0.54
1:B:84:LYS:HE3	1:B:97:TRP:CD1	2.43	0.54
1:B:40:PRO:HB3	1:B:43:GLN:HE21	1.73	0.54
1:A:65:HIS:CG	1:A:66:PRO:HD2	2.43	0.54
1:A:230:THR:CG2	1:A:235:LEU:HD13	2.39	0.53
1:A:2:LYS:NZ	1:A:2:LYS:CB	2.73	0.52
1:A:35:GLU:O	1:A:39:VAL:HG23	2.09	0.52
1:A:17:ARG:HA	1:A:27:VAL:HB	1.92	0.52
1:A:77:ILE:HD11	1:A:239:LEU:HD12	1.92	0.51
1:B:1:MET:SD	1:B:2:LYS:N	2.83	0.51
1:B:310:ARG:HE	1:B:313:ARG:HD2	1.76	0.51
1:A:101:ALA:CA	1:A:105:GLU:OE2	2.58	0.51
1:A:310:ARG:CG	1:A:313:ARG:HD3	2.32	0.50
1:A:310:ARG:CD	1:A:313:ARG:CD	2.88	0.50
1:A:1:MET:O	1:A:2:LYS:C	2.49	0.50
1:A:105:GLU:HA	4:A:530:HOH:O	2.11	0.50
1:A:2:LYS:H	1:A:2:LYS:HZ2	1.46	0.49
1:A:77:ILE:HD13	1:A:239:LEU:HD12	1.93	0.49
1:A:102:GLU:HA	1:A:147:TYR:HB3	1.94	0.49
1:A:52:ARG:HD2	1:A:52:ARG:H	1.77	0.49
1:A:106:TRP:N	1:A:107:PRO:CD	2.75	0.49
1:B:17:ARG:HA	1:B:27:VAL:HB	1.95	0.49
1:B:93:PRO:HD2	1:B:209:MET:SD	2.53	0.49
1:A:135[B]:ARG:HD2	4:A:769:HOH:O	2.13	0.48
1:A:68:PHE:HD2	1:A:71:ARG:HD3	1.78	0.48
1:B:121:ARG:HD2	1:B:150:CYS:SG	2.54	0.47
1:A:40:PRO:HA	1:A:43:GLN:HE22	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:LYS:HB2	1:B:25:ILE:HG12	1.97	0.47
1:A:70:ASN:ND2	1:A:219:TYR:CE1	2.83	0.47
1:B:126:TYR:CZ	1:B:129:ARG:NE	2.83	0.46
1:B:17:ARG:HG2	1:B:17:ARG:NH1	2.31	0.46
1:B:291:ASN:HB3	1:B:294:TRP:CE2	2.51	0.46
1:A:308:GLU:H	1:A:313:ARG:HH22	1.64	0.46
1:A:308:GLU:O	1:A:313:ARG:NH1	2.49	0.46
1:A:308:GLU:H	1:A:313:ARG:NH2	2.13	0.46
1:B:30:VAL:HG23	1:B:34:ALA:HB3	1.98	0.46
1:B:8:GLY:O	4:B:451:HOH:O	2.21	0.45
1:A:39:VAL:HB	1:A:41:PHE:HD2	1.81	0.45
1:B:87:PHE:CZ	1:B:236:ILE:HG21	2.51	0.45
1:B:310:ARG:HH21	1:B:313:ARG:HD2	1.81	0.45
1:B:2:LYS:HB2	1:B:25:ILE:HG23	1.98	0.45
1:B:2:LYS:CG	1:B:25:ILE:HG12	2.47	0.45
1:A:291:ASN:HB3	1:A:294:TRP:CE2	2.51	0.45
1:B:273:GLN:NE2	1:B:274:PRO:HD2	2.31	0.45
1:A:17:ARG:HG2	1:A:17:ARG:HH11	1.82	0.44
1:B:98:GLN:HG3	1:B:99:LEU:O	2.17	0.44
1:B:205:ARG:CZ	1:B:205:ARG:CB	2.95	0.44
1:A:171:ASP:OD1	1:A:171:ASP:C	2.55	0.44
1:A:230:THR:HB	1:A:231:PRO:HD2	1.99	0.44
1:B:121:ARG:HD2	1:B:121:ARG:HA	1.43	0.44
1:A:2:LYS:HZ3	1:A:2:LYS:CB	2.30	0.44
1:A:5:CYS:HB2	1:A:46:ILE:HD13	1.99	0.44
1:B:305:TYR:O	1:B:306:ASP:HB3	2.18	0.44
1:B:181:HIS:O	1:B:192:SER:HA	2.18	0.43
1:B:34:ALA:O	1:B:58:LEU:HB2	2.19	0.43
1:B:199:ASN:OD1	1:B:202:GLN:N	2.44	0.43
1:A:21:GLU:HB3	1:A:22:PRO:HD3	2.00	0.43
1:A:81:LEU:HD13	1:A:99:LEU:HD13	2.00	0.43
1:A:121:ARG:HD2	1:A:150:CYS:SG	2.58	0.43
1:A:77:ILE:HD13	1:A:77:ILE:HG23	1.81	0.43
1:B:73:VAL:O	1:B:74:PHE:C	2.56	0.43
1:B:273:GLN:OE1	4:B:550:HOH:O	2.21	0.43
1:A:65:HIS:CE1	1:A:66:PRO:HD2	2.54	0.42
1:B:66:PRO:HA	1:B:71:ARG:HH21	1.83	0.42
1:B:185:GLN:HE21	1:B:190:ARG:HG3	1.83	0.42
1:A:61:GLN:O	1:A:64:ARG:HG3	2.19	0.42
1:A:1:MET:O	1:A:2:LYS:O	2.37	0.42
1:A:199:ASN:OD1	4:A:681:HOH:O	2.21	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:ARG:CG	1:A:313:ARG:CD	2.94	0.42
1:B:350:ALA:O	1:B:353:LYS:HB2	2.19	0.42
1:A:30:VAL:HG23	1:A:30:VAL:O	2.19	0.42
1:A:39:VAL:HB	1:A:41:PHE:CD2	2.54	0.42
1:A:121:ARG:HD2	1:A:121:ARG:HA	1.60	0.41
1:A:39:VAL:HA	1:A:40:PRO:HD3	1.89	0.41
1:A:127:ASP:CG	1:A:188:ILE:HD12	2.41	0.41
1:B:56:THR:O	1:B:60:ARG:HG3	2.20	0.41
1:B:126:TYR:O	1:B:129:ARG:HD3	2.21	0.41
1:B:353:LYS:HD2	1:B:353:LYS:HA	1.63	0.41
1:A:56:THR:O	1:A:60:ARG:HG3	2.21	0.41
1:B:81:LEU:HD12	1:B:97:TRP:HZ3	1.85	0.41
1:B:40:PRO:HB3	1:B:43:GLN:NE2	2.36	0.41
1:B:319:ASN:C	1:B:320:LEU:HD12	2.41	0.41
1:A:273:GLN:NE2	1:A:274:PRO:HD2	2.35	0.41
1:A:308:GLU:O	1:A:313:ARG:CZ	2.69	0.41
1:A:324:ASP:OD2	1:A:326:SER:OG	2.38	0.41
1:B:171:ASP:OD1	1:B:171:ASP:C	2.59	0.41
1:A:52:ARG:HH11	1:A:52:ARG:HD3	1.74	0.40
1:A:68:PHE:CD2	1:A:71:ARG:HB2	2.55	0.40
1:B:46:ILE:HB	1:B:68:PHE:HA	2.02	0.40
1:B:65:HIS:CE1	1:B:67:ALA:HB3	2.56	0.40
1:B:89:LYS:O	1:B:89:LYS:CG	2.68	0.40
1:B:182:ASN:HD21	1:B:246:SER:HB2	1.86	0.40
1:B:9:ASN:HD22	1:B:9:ASN:HA	1.49	0.40
1:A:181:HIS:O	1:A:192:SER:HA	2.20	0.40
1:A:313:ARG:NH2	4:A:622:HOH:O	2.53	0.40
1:B:332:LEU:HB3	1:B:347:VAL:HG13	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	357/355 (101%)	352 (99%)	4 (1%)	1 (0%)	41	21
1	B	354/355 (100%)	351 (99%)	3 (1%)	0	100	100
All	All	711/710 (100%)	703 (99%)	7 (1%)	1 (0%)	51	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	LYS

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	300/296 (101%)	279 (93%)	21 (7%)	15	3
1	B	297/296 (100%)	279 (94%)	18 (6%)	18	4
All	All	597/592 (101%)	558 (94%)	39 (6%)	17	3

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	9	ASN
1	A	14	ARG
1	A	43	GLN
1	A	52	ARG
1	A	55[A]	GLU
1	A	55[B]	GLU
1	A	58	LEU
1	A	61	GLN
1	A	64	ARG
1	A	68	PHE
1	A	77	ILE
1	A	103	ARG
1	A	104	SER
1	A	145	GLU

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Mol	Chain	Res	Type
1	A	157	ASN
1	A	197	GLN
1	A	245	ASN
1	A	310	ARG
1	A	313	ARG
1	A	320	LEU
1	B	1	MET
1	B	2	LYS
1	B	9	ASN
1	B	30	VAL
1	B	52	ARG
1	B	103	ARG
1	B	121	ARG
1	B	129	ARG
1	B	140	GLU
1	B	145	GLU
1	B	202	GLN
1	B	205	ARG
1	B	210	LEU
1	B	245	ASN
1	B	313	ARG
1	B	320	LEU
1	B	340	PRO
1	B	353	LYS

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	43	GLN
1	A	154	GLN
1	A	182	ASN
1	A	185	GLN
1	A	202	GLN
1	A	273	GLN
1	B	9	ASN
1	B	61	GLN
1	B	182	ASN
1	B	185	GLN
1	B	201	GLN
1	B	273	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	ATP	B	400	3	26,33,33	0.93	0	31,52,52	2.01	9 (29%)
2	ATP	A	400	3	26,33,33	0.88	0	31,52,52	2.10	11 (35%)

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	ATP	B	400	3	-	3/18/38/38	0/3/3/3
2	ATP	A	400	3	-	3/18/38/38	0/3/3/3

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	ATP	C5-C6-N6	4.90	127.79	120.35
2	A	400	ATP	C5-C6-N1	-4.67	109.76	120.35
2	B	400	ATP	N3-C2-N1	-4.58	121.52	128.68
2	B	400	ATP	C2-N1-C6	4.40	126.28	118.75
2	A	400	ATP	C5-C6-N6	4.37	126.99	120.35
2	A	400	ATP	C2-N1-C6	4.30	126.11	118.75
2	A	400	ATP	O2G-PG-O3B	3.94	117.86	104.64
2	B	400	ATP	C5-C6-N1	-3.76	111.84	120.35
2	A	400	ATP	O4'-C4'-C3'	-3.06	99.06	105.11
2	A	400	ATP	PB-O3B-PG	-3.03	122.43	132.83
2	A	400	ATP	O2A-PA-O5'	2.89	121.17	107.75
2	B	400	ATP	C2'-C3'-C4'	-2.74	97.31	102.64
2	B	400	ATP	O5'-C5'-C4'	-2.57	100.14	108.99
2	A	400	ATP	O2'-C2'-C1'	-2.44	101.85	110.85
2	A	400	ATP	N3-C2-N1	-2.30	125.08	128.68
2	B	400	ATP	O4'-C4'-C5'	2.30	116.93	109.37
2	A	400	ATP	O3B-PG-O1G	-2.24	98.77	111.19
2	A	400	ATP	N6-C6-N1	2.18	123.09	118.57
2	B	400	ATP	PB-O3B-PG	-2.09	125.66	132.83
2	B	400	ATP	O4'-C1'-C2'	-2.07	103.90	106.93

There are no chirality outliers.

All (6) torsion outliers are listed below:

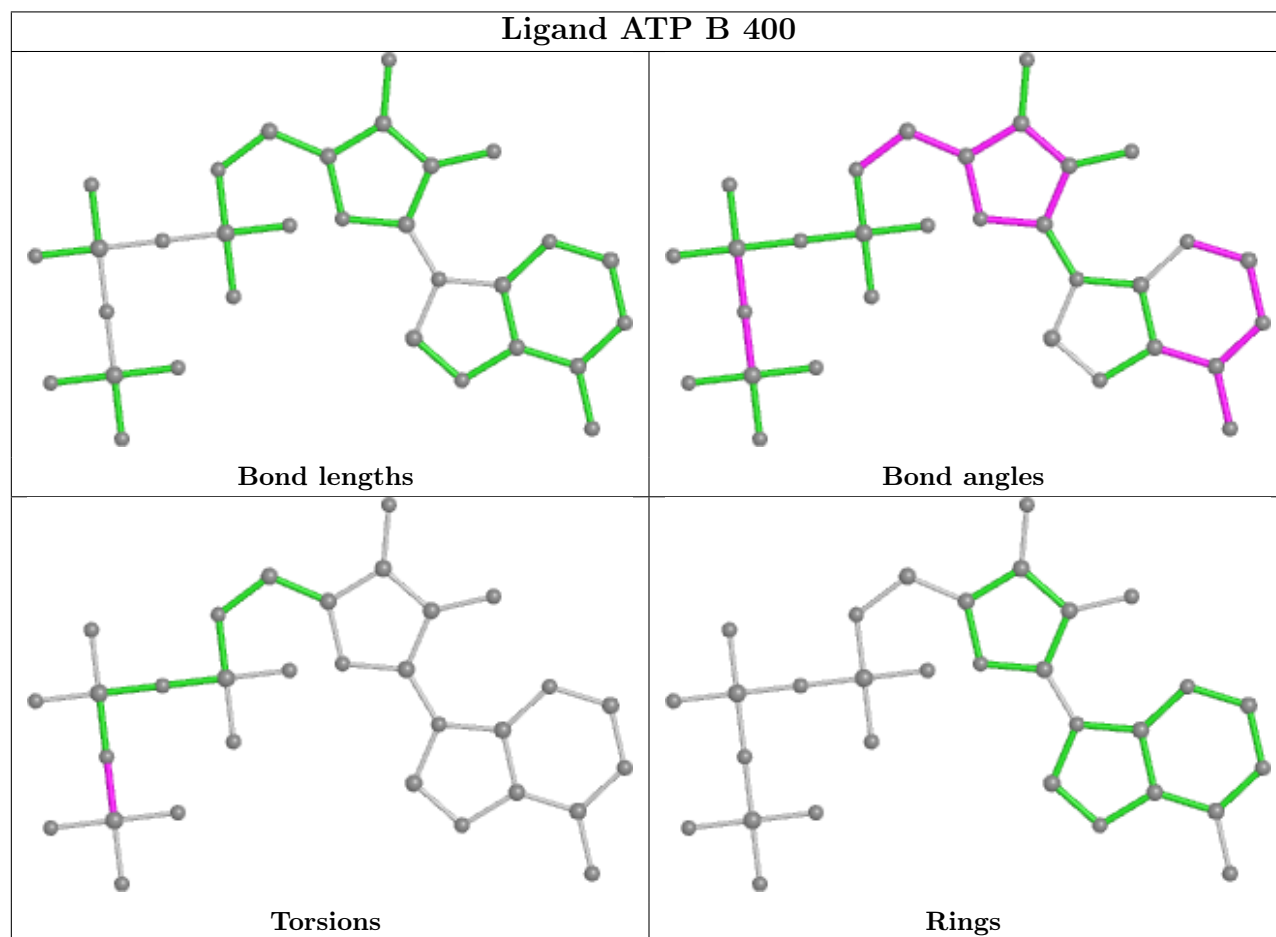
Mol	Chain	Res	Type	Atoms
2	A	400	ATP	PB-O3B-PG-O2G
2	B	400	ATP	PB-O3B-PG-O2G
2	A	400	ATP	PB-O3A-PA-O2A
2	A	400	ATP	PB-O3B-PG-O3G
2	B	400	ATP	PB-O3B-PG-O3G
2	B	400	ATP	PB-O3B-PG-O1G

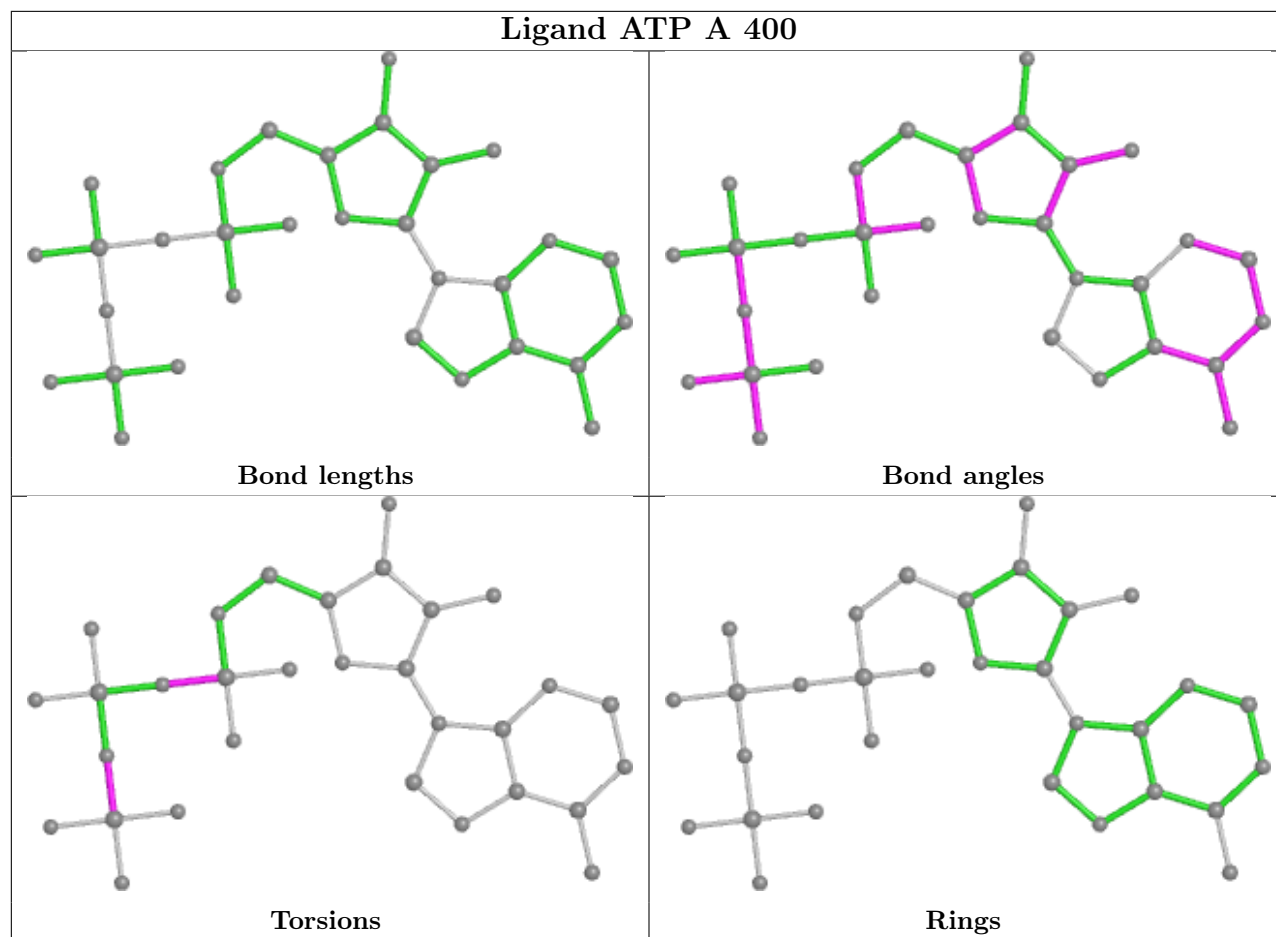
There are no ring outliers.

No monomer is involved in short contacts.

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 [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	355/355 (100%)	-0.47	3 (0%) 86 86	12, 21, 56, 84	0
1	B	355/355 (100%)	-0.56	1 (0%) 94 93	10, 22, 52, 86	0
All	All	710/710 (100%)	-0.52	4 (0%) 89 89	10, 21, 54, 86	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	41	PHE	3.1
1	A	1	MET	3.1
1	B	41	PHE	2.6
1	A	39	VAL	2.2

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
3	MG	A	402	1/1	0.97	0.04	16,16,16,16	0

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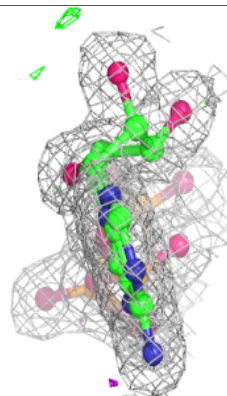
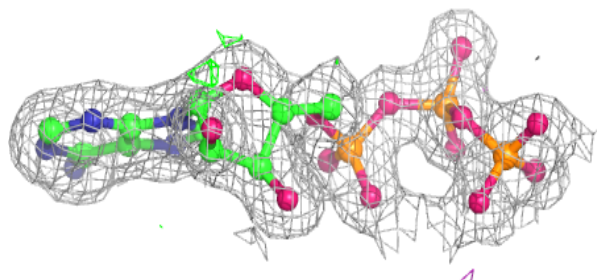
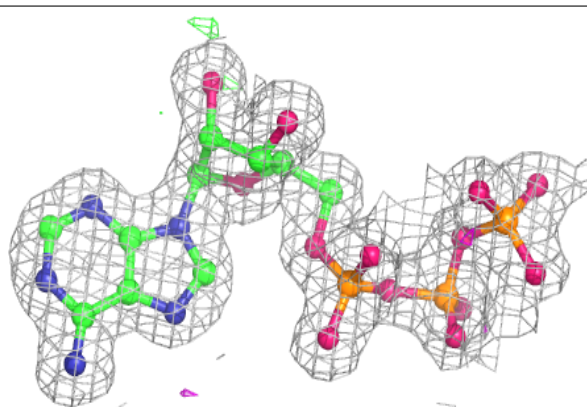
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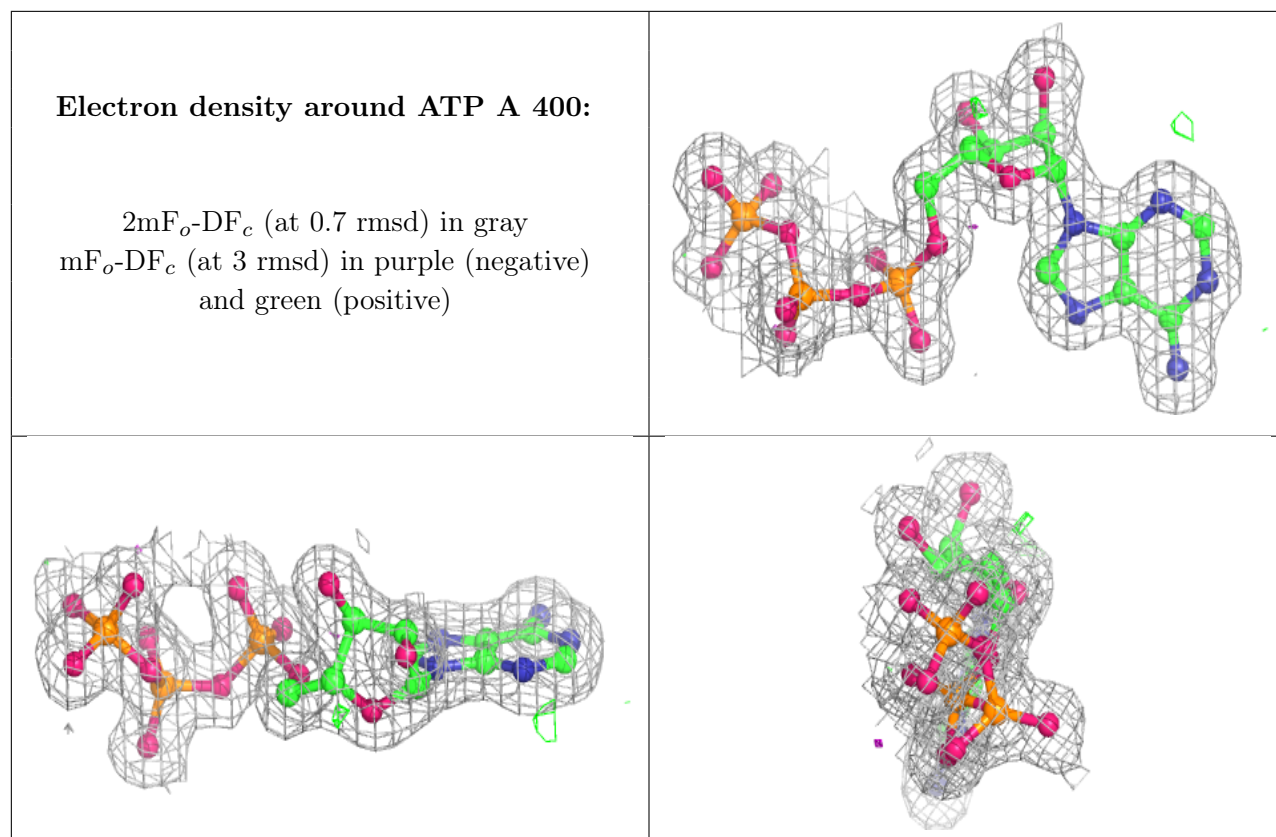
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ATP	B	400	31/31	0.98	0.05	15,19,24,31	0
2	ATP	A	400	31/31	0.98	0.05	13,20,25,29	0
3	MG	B	402	1/1	0.98	0.04	19,19,19,19	0
3	MG	B	401	1/1	0.99	0.03	17,17,17,17	0
3	MG	A	401	1/1	0.99	0.04	17,17,17,17	0

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 ATP B 400:

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.