



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 01:43 AM UTC

PDB ID : 3FVQ / pdb_00003fvq
Title : Crystal structure of the nucleotide binding domain FbpC complexed with ATP
Authors : Newstead, S.; Bilton, P.; Carpenter, E.P.; Campopiano, D.; Iwata, S.
Deposited on : 2009-01-16
Resolution : 1.90 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

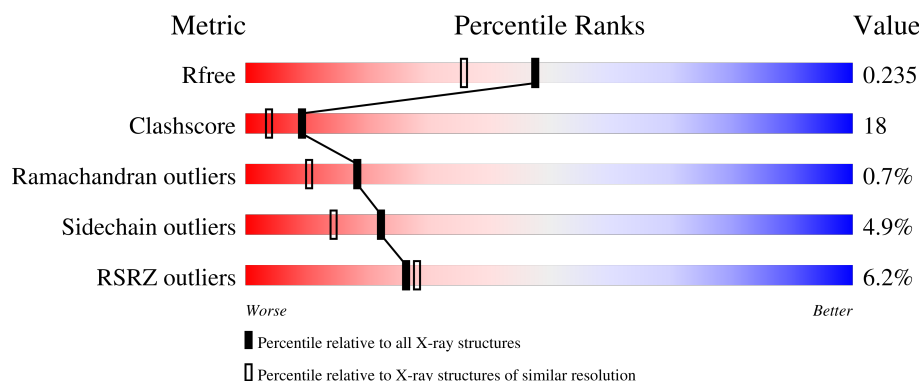
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	359	8% 70% 22% • •
1	B	359	4% 70% 25% • •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5963 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.

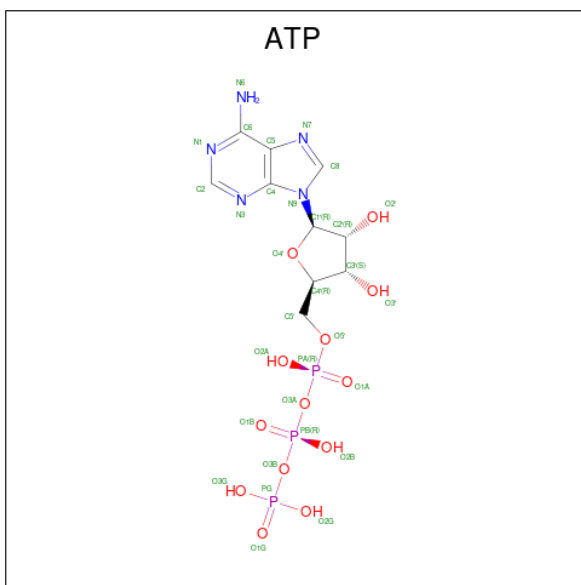
- Molecule 1 is a protein called Fe(3+) ions import ATP-binding protein fbpC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	0	10	0
			2653	1677	477	492	7			
1	B	350	Total	C	N	O	S	0	13	0
			2723	1719	494	504	6			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	147	ALA	VAL	SEE REMARK 999	UNP Q5FA19
A	283	TYR	PHE	SEE REMARK 999	UNP Q5FA19
A	353	GLU	-	expression tag	UNP Q5FA19
A	354	HIS	-	expression tag	UNP Q5FA19
A	355	HIS	-	expression tag	UNP Q5FA19
A	356	HIS	-	expression tag	UNP Q5FA19
A	357	HIS	-	expression tag	UNP Q5FA19
A	358	HIS	-	expression tag	UNP Q5FA19
A	359	HIS	-	expression tag	UNP Q5FA19
B	147	ALA	VAL	SEE REMARK 999	UNP Q5FA19
B	283	TYR	PHE	SEE REMARK 999	UNP Q5FA19
B	353	GLU	-	expression tag	UNP Q5FA19
B	354	HIS	-	expression tag	UNP Q5FA19
B	355	HIS	-	expression tag	UNP Q5FA19
B	356	HIS	-	expression tag	UNP Q5FA19
B	357	HIS	-	expression tag	UNP Q5FA19
B	358	HIS	-	expression tag	UNP Q5FA19
B	359	HIS	-	expression tag	UNP Q5FA19

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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 CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Ca	0	0
			2	2		
3	B	1	Total	Ca	0	0
			1	1		

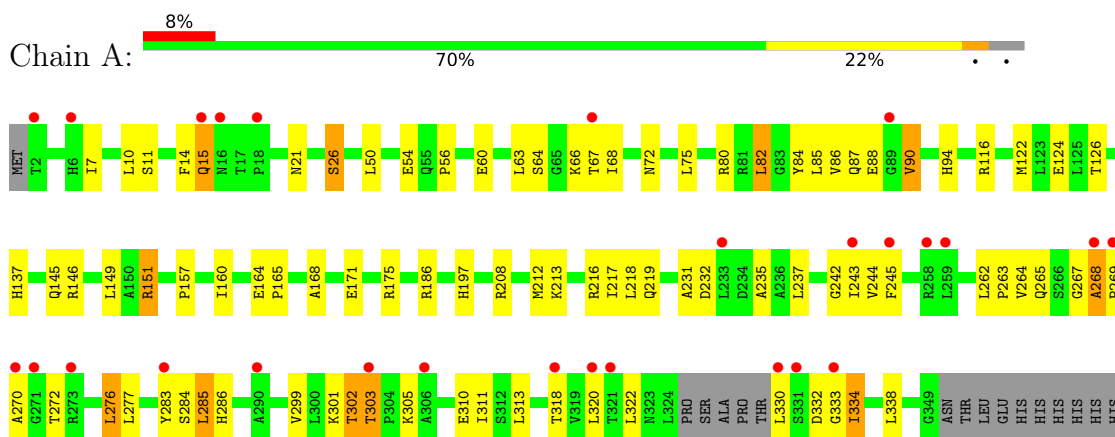
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	243	Total	O	0	0
			243	243		
4	B	279	Total	O	0	0
			279	279		

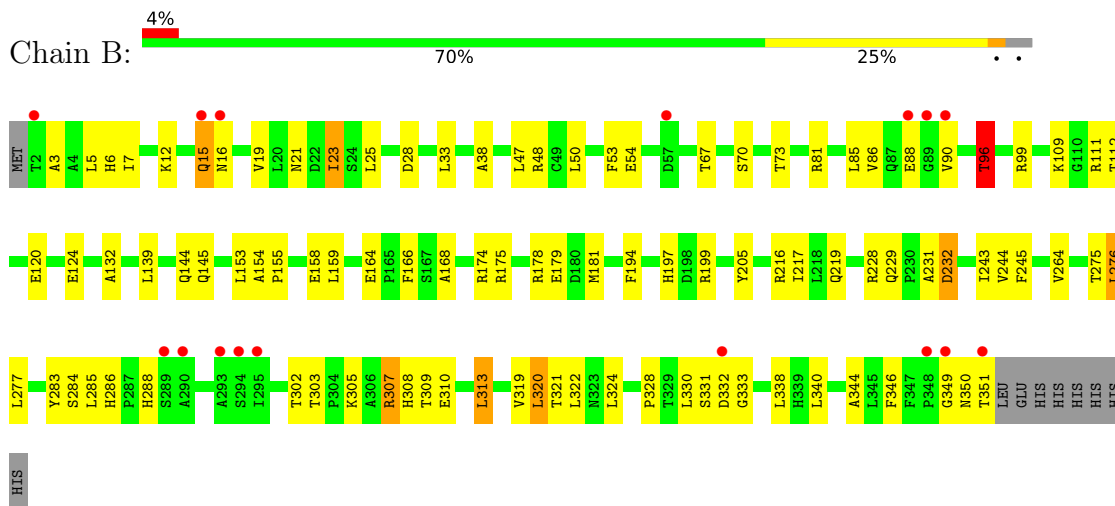
3 Residue-property plots [i](#)

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: Fe(3+) ions import ATP-binding protein fbpC



- Molecule 1: Fe(3+) ions import ATP-binding protein fbpC



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.86Å 89.09Å 149.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	52.27 – 1.90 52.27 – 1.90	Depositor EDS
% Data completeness (in resolution range)	87.4 (52.27-1.90) 87.4 (52.27-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.195 , 0.256 0.185 , 0.235	Depositor DCC
R_{free} test set	2624 reflections (4.41%)	wwPDB-VP
Wilson B-factor (Å ²)	17.1	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5963	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% 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: CA, ATP

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	0.87	0/2731	0.97	2/3702 (0.1%)
1	B	0.96	2/2813 (0.1%)	1.00	3/3819 (0.1%)
All	All	0.92	2/5544 (0.0%)	0.99	5/7521 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	38	ALA	CA-CB	6.52	1.64	1.53
1	B	232	ASP	N-CA	5.58	1.53	1.45

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	23	ILE	CB-CA-C	-7.17	101.83	111.15
1	B	96	THR	N-CA-CB	-5.86	101.89	110.26
1	A	303	THR	CA-C-N	5.55	125.87	119.93
1	A	303	THR	C-N-CA	5.55	125.87	119.93
1	B	90	VAL	N-CA-C	5.31	115.85	108.84

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	2653	0	2711	102	0
1	B	2723	0	2796	110	0
2	A	31	0	12	0	0
2	B	31	0	12	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
4	A	243	0	0	21	0
4	B	279	0	0	13	0
All	All	5963	0	5531	198	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 18.

All (198) 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:276:LEU:HB3	4:A:448:HOH:O	1.39	1.22
1:A:242:GLY:HA3	1:A:277:LEU:HD11	1.30	1.07
1:B:178:ARG:HD2	4:B:448:HOH:O	1.58	1.03
1:B:5:LEU:HD11	1:B:50[B]:LEU:HD11	1.38	1.01
1:B:228:ARG:C	1:B:229[B]:GLN:HE21	1.69	1.01
1:B:145:GLN:HE22	1:B:168:ALA:H	1.10	0.98
1:A:116:ARG:HB2	4:A:545:HOH:O	1.61	0.98
1:B:21:ASN:HB3	4:B:557:HOH:O	1.64	0.96
1:B:112[A]:THR:HG22	4:B:451:HOH:O	1.68	0.94
1:B:5:LEU:HD11	1:B:50[B]:LEU:CD1	1.98	0.93
1:B:303[A]:THR:CG2	1:B:310:GLU:HB3	1.98	0.93
1:B:228:ARG:C	1:B:229[B]:GLN:NE2	2.27	0.93
1:A:303:THR:HG22	1:A:305:LYS:HE3	1.52	0.91
1:A:145:GLN:HE22	1:A:168:ALA:H	1.19	0.90
1:A:151:ARG:HD2	4:A:410:HOH:O	1.71	0.89
1:B:7:ILE:HD11	1:B:50[A]:LEU:HD21	1.55	0.88
1:A:276:LEU:CD2	1:B:344:ALA:HB3	2.04	0.87
1:B:5:LEU:CD1	1:B:50[B]:LEU:HD11	2.05	0.86
1:A:243:ILE:HD12	1:A:283:TYR:HE2	1.37	0.86
1:A:243:ILE:HD12	1:A:283:TYR:CE2	2.11	0.85
1:A:232:ASP:HB3	1:A:235:ALA:H	1.41	0.83
1:A:299:VAL:HG21	1:A:330:LEU:HD21	1.62	0.82
1:B:228:ARG:O	1:B:229[B]:GLN:NE2	2.13	0.82
1:A:264:VAL:HG21	1:B:346:PHE:HB2	1.63	0.81
1:A:242:GLY:CA	1:A:277:LEU:HD11	2.09	0.81
1:A:320:LEU:HG	1:A:322:LEU:HD21	1.62	0.81

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:ALA:HB1	1:A:269:PRO:HD2	1.60	0.81
1:B:70:SER:OG	1:B:73:THR:HG22	1.81	0.81
1:A:276:LEU:HD22	4:A:448:HOH:O	1.81	0.79
1:A:171:GLU:HG3	1:A:175:ARG:HE	1.48	0.78
1:B:85:LEU:HD21	4:B:459:HOH:O	1.84	0.76
1:A:164:GLU:OE1	1:A:197:HIS:HD2	1.69	0.75
1:B:48[B]:ARG:HG2	1:B:53:PHE:HB2	1.69	0.75
1:B:109:LYS:HD3	1:B:111:ARG:NH1	2.02	0.74
1:A:26:SER:O	1:A:208:ARG:NH2	2.20	0.73
1:A:7:ILE:HD11	1:A:50:LEU:HD21	1.69	0.73
1:B:7:ILE:CD1	1:B:50[A]:LEU:HD21	2.18	0.72
1:A:276:LEU:CD2	1:B:344:ALA:CB	2.68	0.71
1:B:349:GLY:O	1:B:351:THR:N	2.24	0.71
1:B:178:ARG:HD3	1:B:205:TYR:CE2	2.25	0.71
1:B:164:GLU:OE1	1:B:197:HIS:HD2	1.73	0.70
1:B:303[A]:THR:HG22	1:B:310:GLU:HB3	1.74	0.70
1:A:116:ARG:CG	4:A:545:HOH:O	2.39	0.69
1:B:67:THR:HG22	1:B:73:THR:HG21	1.75	0.69
1:B:216:ARG:NE	4:B:410:HOH:O	2.25	0.68
1:A:14:PHE:O	1:A:15:GLN:C	2.34	0.68
1:A:334:ILE:HD12	4:A:519:HOH:O	1.91	0.68
1:B:351:THR:HG23	1:B:351:THR:O	1.93	0.68
1:A:94:HIS:HE1	4:A:387:HOH:O	1.76	0.67
1:A:301:LYS:HA	1:A:332:ASP:OD2	1.94	0.67
1:A:268:ALA:HB1	1:A:269:PRO:CD	2.25	0.67
1:B:284:SER:OG	1:B:286:HIS:HE1	1.76	0.67
1:B:219:GLN:NE2	1:B:231:ALA:H	1.92	0.67
1:B:313:LEU:O	1:B:319:VAL:HG13	1.95	0.66
1:B:96:THR:HG21	4:B:477:HOH:O	1.96	0.65
1:B:88:GLU:H	1:B:88:GLU:CD	2.04	0.65
1:B:303[A]:THR:HG23	1:B:310:GLU:HB3	1.74	0.65
1:B:67:THR:CG2	1:B:73:THR:HG21	2.26	0.64
1:B:48[B]:ARG:HG3	1:B:48[B]:ARG:HH11	1.63	0.64
4:A:590:HOH:O	1:B:275:THR:HG23	1.97	0.63
1:A:245:PHE:HZ	1:B:320:LEU:HD22	1.64	0.63
1:B:124:GLU:HG3	4:B:580:HOH:O	1.98	0.62
1:A:82:LEU:O	1:A:82:LEU:HD23	2.00	0.62
1:A:171:GLU:CG	1:A:175:ARG:HE	2.12	0.62
1:A:7:ILE:CD1	1:A:50:LEU:HD21	2.29	0.61
1:A:186:ARG:NH1	4:A:599:HOH:O	2.33	0.61
1:B:228:ARG:HB3	1:B:229[B]:GLN:HE22	1.64	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:GLN:HE22	1:B:231:ALA:H	1.47	0.61
1:B:96:THR:HG22	1:B:99:ARG:H	1.66	0.61
1:A:171:GLU:HG3	1:A:175:ARG:NE	2.16	0.61
1:A:299:VAL:CG2	1:A:330:LEU:HD21	2.30	0.61
1:B:145:GLN:NE2	1:B:168:ALA:H	1.91	0.61
1:A:68:ILE:HD12	1:A:80:ARG:HD3	1.82	0.61
1:A:84:TYR:CE2	1:A:86[B]:VAL:HG12	2.36	0.60
1:A:286:HIS:HD2	4:A:532:HOH:O	1.85	0.60
1:A:219:GLN:HE22	1:A:231:ALA:H	1.49	0.60
1:A:88:GLU:N	1:A:88:GLU:CD	2.60	0.59
1:A:276:LEU:HD21	1:B:344:ALA:HB3	1.83	0.59
1:B:154:ALA:HB3	1:B:155:PRO:HD3	1.85	0.59
1:A:88:GLU:CD	1:A:88:GLU:H	2.10	0.59
1:B:70:SER:OG	1:B:73:THR:CG2	2.50	0.59
1:A:284:SER:OG	1:A:286:HIS:HE1	1.84	0.59
1:A:219:GLN:NE2	1:A:231:ALA:H	2.00	0.59
1:B:275:THR:HG22	4:B:489:HOH:O	2.02	0.58
1:A:303:THR:HG21	4:A:483:HOH:O	2.03	0.58
1:B:330:LEU:C	1:B:332:ASP:H	2.09	0.58
1:B:313:LEU:O	1:B:319:VAL:HA	2.04	0.58
1:B:174:ARG:O	1:B:178:ARG:HG2	2.05	0.57
1:A:216:ARG:HH11	1:A:216:ARG:HG3	1.69	0.57
1:A:302:THR:HG22	1:A:302:THR:O	2.03	0.56
1:A:213:LYS:HB3	1:A:218:LEU:HD13	1.86	0.56
1:B:330:LEU:C	1:B:332:ASP:N	2.61	0.56
1:B:232:ASP:OD1	4:B:543:HOH:O	2.17	0.56
1:A:276:LEU:CD2	4:A:448:HOH:O	2.48	0.56
1:A:84:TYR:CZ	1:A:86[B]:VAL:HG12	2.41	0.55
1:A:116:ARG:CB	4:A:545:HOH:O	2.30	0.55
1:B:175:ARG:HA	1:B:178:ARG:HG2	1.88	0.55
1:B:181:MET:HE3	1:B:194:PHE:HZ	1.72	0.55
1:B:285:LEU:O	1:B:288:HIS:CD2	2.61	0.54
1:A:171:GLU:HG2	1:B:199:ARG:NH1	2.22	0.54
1:A:164:GLU:OE1	1:A:197:HIS:CD2	2.55	0.54
1:A:269:PRO:HG2	1:A:272:THR:OG1	2.08	0.53
1:A:311:ILE:HG22	1:A:313:LEU:HG	1.90	0.53
1:B:109:LYS:HD3	1:B:111:ARG:CZ	2.38	0.53
1:A:87:GLN:O	4:A:565:HOH:O	2.19	0.53
1:A:320:LEU:CG	1:A:322:LEU:HD21	2.37	0.53
1:A:320:LEU:HG	1:A:322:LEU:CD2	2.37	0.53
1:A:66:LYS:HE2	1:A:75:LEU:HD11	1.90	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48[B]:ARG:HH11	1:B:48[B]:ARG:CG	2.21	0.52
1:B:5:LEU:HD11	1:B:50[B]:LEU:HD13	1.89	0.52
1:A:334:ILE:CD1	4:A:519:HOH:O	2.54	0.52
1:A:299:VAL:HG21	1:A:330:LEU:CD2	2.39	0.51
1:A:269:PRO:HD2	1:A:272:THR:OG1	2.10	0.51
1:B:351:THR:O	1:B:351:THR:CG2	2.58	0.51
1:B:330:LEU:O	1:B:332:ASP:N	2.44	0.51
1:A:276:LEU:CG	4:A:448:HOH:O	2.56	0.50
1:A:82:LEU:HD23	1:A:82:LEU:C	2.37	0.50
1:A:145:GLN:NE2	1:A:168:ALA:H	1.99	0.49
1:B:6[A]:HIS:HD2	4:B:369:HOH:O	1.95	0.49
1:A:244:VAL:HG22	4:A:486:HOH:O	2.11	0.49
1:B:124:GLU:OE2	4:B:566:HOH:O	2.19	0.49
1:A:276:LEU:HD22	1:B:344:ALA:CB	2.42	0.49
1:B:139:LEU:HB2	1:B:144:GLN:HG3	1.95	0.49
1:A:126:THR:O	1:A:146:ARG:HD2	2.13	0.48
1:B:86:VAL:CG1	1:B:88:GLU:OE1	2.61	0.48
1:B:96:THR:HG23	1:B:132:ALA:O	2.14	0.48
1:B:81[B]:ARG:HH22	1:B:158:GLU:H	1.61	0.48
1:B:88:GLU:CD	1:B:88:GLU:N	2.69	0.48
1:B:228:ARG:CB	1:B:229[B]:GLN:HE22	2.26	0.48
1:B:276:LEU:C	1:B:276:LEU:HD12	2.38	0.48
1:A:243:ILE:CD1	1:A:283:TYR:HE2	2.16	0.48
1:B:228:ARG:CB	1:B:229[B]:GLN:NE2	2.77	0.48
1:B:120:GLU:O	1:B:124:GLU:HG3	2.14	0.47
1:A:283:TYR:CD2	1:B:322:LEU:HD22	2.49	0.47
1:B:164:GLU:OE1	1:B:197:HIS:CD2	2.63	0.47
1:B:328:PRO:HB2	1:B:330:LEU:HD21	1.95	0.47
1:A:285:LEU:HB2	1:B:324:LEU:CD1	2.43	0.47
1:A:264:VAL:CG2	1:B:346:PHE:HB2	2.39	0.47
1:B:12:LYS:O	1:B:19[B]:VAL:HG22	2.15	0.47
1:A:116:ARG:HG3	4:A:545:HOH:O	2.12	0.46
1:A:212:MET:HG2	1:A:217:ILE:HD13	1.98	0.46
1:B:86:VAL:HG13	1:B:88:GLU:OE1	2.16	0.46
1:B:88:GLU:OE1	1:B:88:GLU:N	2.31	0.46
1:B:15:GLN:HB2	1:B:16:ASN:H	1.45	0.46
1:A:149:LEU:HB2	4:A:480:HOH:O	2.15	0.46
1:B:48[B]:ARG:CG	1:B:48[B]:ARG:NH1	2.78	0.46
1:A:301:LYS:HA	1:A:332:ASP:CG	2.40	0.45
1:A:303:THR:CG2	1:A:305:LYS:HE3	2.37	0.45
1:A:305:LYS:HA	1:A:305:LYS:HD3	1.71	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:ALA:HB3	1:B:159:LEU:HB2	1.99	0.45
1:B:308:HIS:CD2	4:B:441:HOH:O	2.69	0.45
1:A:137:HIS:HD2	4:A:381:HOH:O	2.00	0.45
1:B:243:ILE:HD13	1:B:283:TYR:HE2	1.81	0.44
1:B:307:ARG:O	1:B:307:ARG:HG3	2.18	0.44
1:A:10[A]:LEU:HD11	1:A:56:PRO:HG2	2.00	0.44
1:A:21[B]:ASN:HB2	1:A:216:ARG:HG2	1.99	0.44
1:A:86[A]:VAL:HG12	1:A:88:GLU:OE2	2.17	0.44
1:A:269:PRO:O	1:A:270:ALA:C	2.61	0.44
1:A:285:LEU:HD13	1:B:328:PRO:HG3	1.99	0.44
1:A:216:ARG:HG3	1:A:216:ARG:NH1	2.32	0.44
1:B:7:ILE:HD11	1:B:50[A]:LEU:CD2	2.37	0.44
1:B:302:THR:HG23	1:B:309:THR:HG21	1.99	0.43
1:A:265:GLN:OE1	4:A:531:HOH:O	2.21	0.43
1:B:166:PHE:HZ	1:B:181:MET:HE1	1.84	0.43
1:B:166:PHE:HZ	1:B:181:MET:CE	2.31	0.43
1:B:25:LEU:HD21	1:B:33:LEU:HD21	2.01	0.43
1:B:178:ARG:HG3	1:B:179[B]:GLU:H	1.83	0.43
1:A:262:LEU:HA	1:A:263:PRO:HD3	1.85	0.43
1:A:305:LYS:NZ	1:A:310:GLU:OE1	2.52	0.43
1:B:23:ILE:HD13	1:B:217[A]:ILE:HG13	2.01	0.42
1:A:15:GLN:HE21	1:A:15:GLN:HB3	1.48	0.42
1:A:276:LEU:HD23	1:B:344:ALA:CB	2.47	0.42
1:B:284:SER:OG	1:B:286:HIS:CE1	2.65	0.42
1:A:338:LEU:HD12	1:B:284:SER:O	2.20	0.42
1:B:277:LEU:C	1:B:277:LEU:HD23	2.45	0.42
1:B:244:VAL:HG12	1:B:245:PHE:N	2.33	0.42
1:A:85:LEU:HD23	1:A:165:PRO:HB3	2.02	0.41
1:B:67:THR:HG21	1:B:73:THR:HG21	2.00	0.41
1:A:157:PRO:HG2	1:A:160:ILE:HB	2.02	0.41
1:A:285:LEU:HA	1:A:285:LEU:HD23	1.77	0.41
1:A:67:THR:O	1:A:67:THR:HG23	2.20	0.41
1:B:178:ARG:CD	1:B:205:TYR:CE2	3.01	0.41
1:A:10[B]:LEU:HD23	1:A:11:SER:N	2.36	0.41
1:A:284:SER:O	1:B:338:LEU:HD12	2.21	0.41
1:B:308:HIS:HE1	4:B:474:HOH:O	2.04	0.41
1:B:12:LYS:HE2	1:B:54:GLU:HG3	2.02	0.41
1:B:330:LEU:O	1:B:333:GLY:N	2.39	0.41
1:A:267:GLY:O	1:A:268:ALA:O	2.40	0.40
1:B:313:LEU:HD23	1:B:313:LEU:HA	1.86	0.40
1:A:122:MET:HA	1:A:122:MET:HE2	2.03	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:LEU:HD23	1:A:237:LEU:HA	1.94	0.40
1:A:90:VAL:O	1:A:90:VAL:HG13	2.22	0.40
1:B:81[B]:ARG:NH2	1:B:158:GLU:H	2.18	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	349/359 (97%)	335 (96%)	11 (3%)	3 (1%)	14	6
1	B	361/359 (101%)	348 (96%)	11 (3%)	2 (1%)	21	13
All	All	710/718 (99%)	683 (96%)	22 (3%)	5 (1%)	18	10

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	350	ASN
1	A	90	VAL
1	A	268	ALA
1	A	333	GLY
1	B	331	SER

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	282/288 (98%)	267 (95%)	15 (5%)	20	12
1	B	291/288 (101%)	278 (96%)	13 (4%)	24	17
All	All	573/576 (100%)	545 (95%)	28 (5%)	22	14

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	26	SER
1	A	54[A]	GLU
1	A	54[B]	GLU
1	A	60	GLU
1	A	63	LEU
1	A	64	SER
1	A	72	ASN
1	A	82	LEU
1	A	151	ARG
1	A	276	LEU
1	A	285	LEU
1	A	302	THR
1	A	318	THR
1	A	334	ILE
1	B	15	GLN
1	B	28	ASP
1	B	47	LEU
1	B	96	THR
1	B	153	LEU
1	B	264	VAL
1	B	276	LEU
1	B	305	LYS
1	B	307	ARG
1	B	313	LEU
1	B	320	LEU
1	B	321	THR
1	B	340	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	72	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	94	HIS
1	A	114	GLN
1	A	137	HIS
1	A	145	GLN
1	A	188	ASN
1	A	197	HIS
1	A	214	GLN
1	A	219	GLN
1	A	286	HIS
1	A	288	HIS
1	A	323	ASN
1	B	15	GLN
1	B	21	ASN
1	B	117	GLN
1	B	137	HIS
1	B	145	GLN
1	B	188	ASN
1	B	197	HIS
1	B	204	GLN
1	B	219	GLN
1	B	286	HIS
1	B	308	HIS
1	B	323	ASN

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](#)

Of 5 ligands modelled in this entry, 3 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	360	3	32,33,33	1.58	5 (15%)	48,52,52	1.88	12 (25%)
2	ATP	A	401	3	32,33,33	1.82	8 (25%)	48,52,52	1.85	14 (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	ATP	B	360	3	-	1/22/38/38	0/3/3/3
2	ATP	A	401	3	-	1/22/38/38	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	360	ATP	PB-O3A	5.61	1.65	1.59
2	A	401	ATP	PB-O3A	5.19	1.65	1.59
2	A	401	ATP	C5-C4	4.36	1.46	1.39
2	A	401	ATP	PB-O3B	3.36	1.63	1.59
2	B	360	ATP	C5-C4	3.29	1.45	1.39
2	A	401	ATP	C8-N7	2.82	1.37	1.31
2	A	401	ATP	PA-O3A	2.79	1.62	1.59
2	B	360	ATP	C5-C6	2.64	1.48	1.41
2	A	401	ATP	C5-C6	2.59	1.48	1.41
2	A	401	ATP	C4-N9	-2.48	1.32	1.37
2	B	360	ATP	C5-N7	-2.37	1.34	1.39
2	B	360	ATP	C4-N9	-2.27	1.33	1.37
2	A	401	ATP	C5-N7	-2.03	1.35	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	360	ATP	N3-C2-N1	-4.43	121.87	128.58
2	A	401	ATP	N3-C2-N1	-4.37	121.96	128.58

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	360	ATP	C5-C4-N3	-4.22	120.91	126.72
2	A	401	ATP	C5-C4-N3	-4.22	120.91	126.72
2	B	360	ATP	C4-N9-C8	4.08	110.02	105.74
2	B	360	ATP	N9-C8-N7	-3.87	108.44	113.94
2	B	360	ATP	C4-C5-N7	-3.63	106.44	110.58
2	B	360	ATP	C5-N7-C8	3.60	109.11	103.45
2	A	401	ATP	O2B-PB-O3B	3.57	116.93	107.27
2	B	360	ATP	N3-C4-N9	3.49	133.10	127.17
2	A	401	ATP	C2-N1-C6	3.44	124.38	118.73
2	A	401	ATP	C2-N3-C4	3.26	119.79	111.83
2	A	401	ATP	C4-C5-N7	-3.25	106.86	110.58
2	B	360	ATP	C2-N3-C4	3.22	119.70	111.83
2	A	401	ATP	N3-C4-N9	3.10	132.44	127.17
2	A	401	ATP	C4-N9-C8	3.01	108.90	105.74
2	B	360	ATP	C2-N1-C6	3.01	123.68	118.73
2	B	360	ATP	C6-C5-N7	2.70	137.30	132.09
2	A	401	ATP	C6-C5-N7	2.66	137.23	132.09
2	A	401	ATP	O3A-PB-O1B	-2.55	103.02	110.70
2	B	360	ATP	O2B-PB-O3B	2.44	113.86	107.27
2	B	360	ATP	C4-N9-C1'	-2.40	121.02	126.63
2	A	401	ATP	C4-N9-C1'	-2.35	121.12	126.63
2	A	401	ATP	N9-C8-N7	-2.22	110.78	113.94
2	A	401	ATP	C5-N7-C8	2.14	106.81	103.45
2	A	401	ATP	O2A-PA-O1A	2.04	121.93	112.44

There are no chirality outliers.

All (2) torsion outliers are listed below:

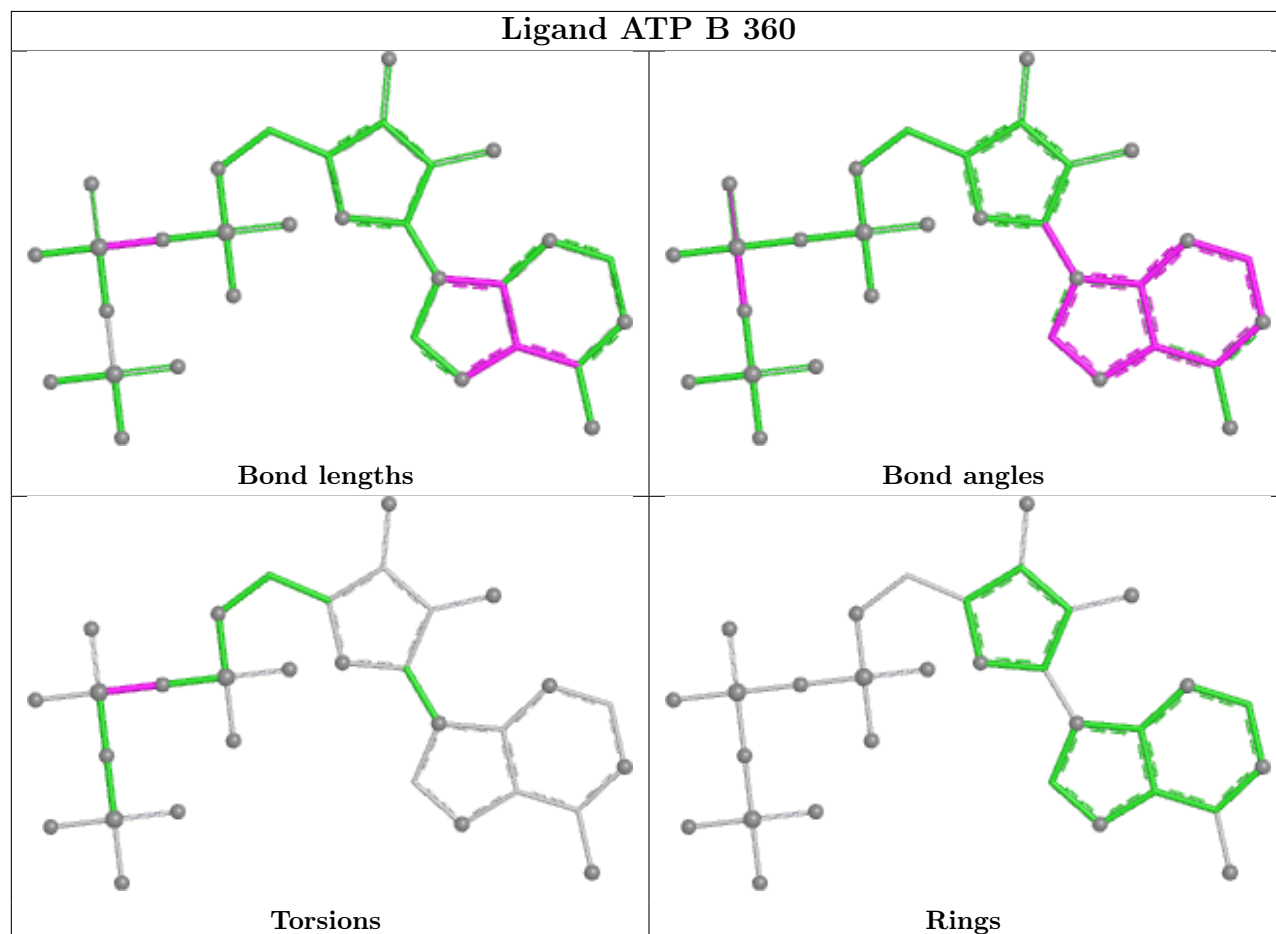
Mol	Chain	Res	Type	Atoms
2	A	401	ATP	PA-O3A-PB-O2B
2	B	360	ATP	PA-O3A-PB-O2B

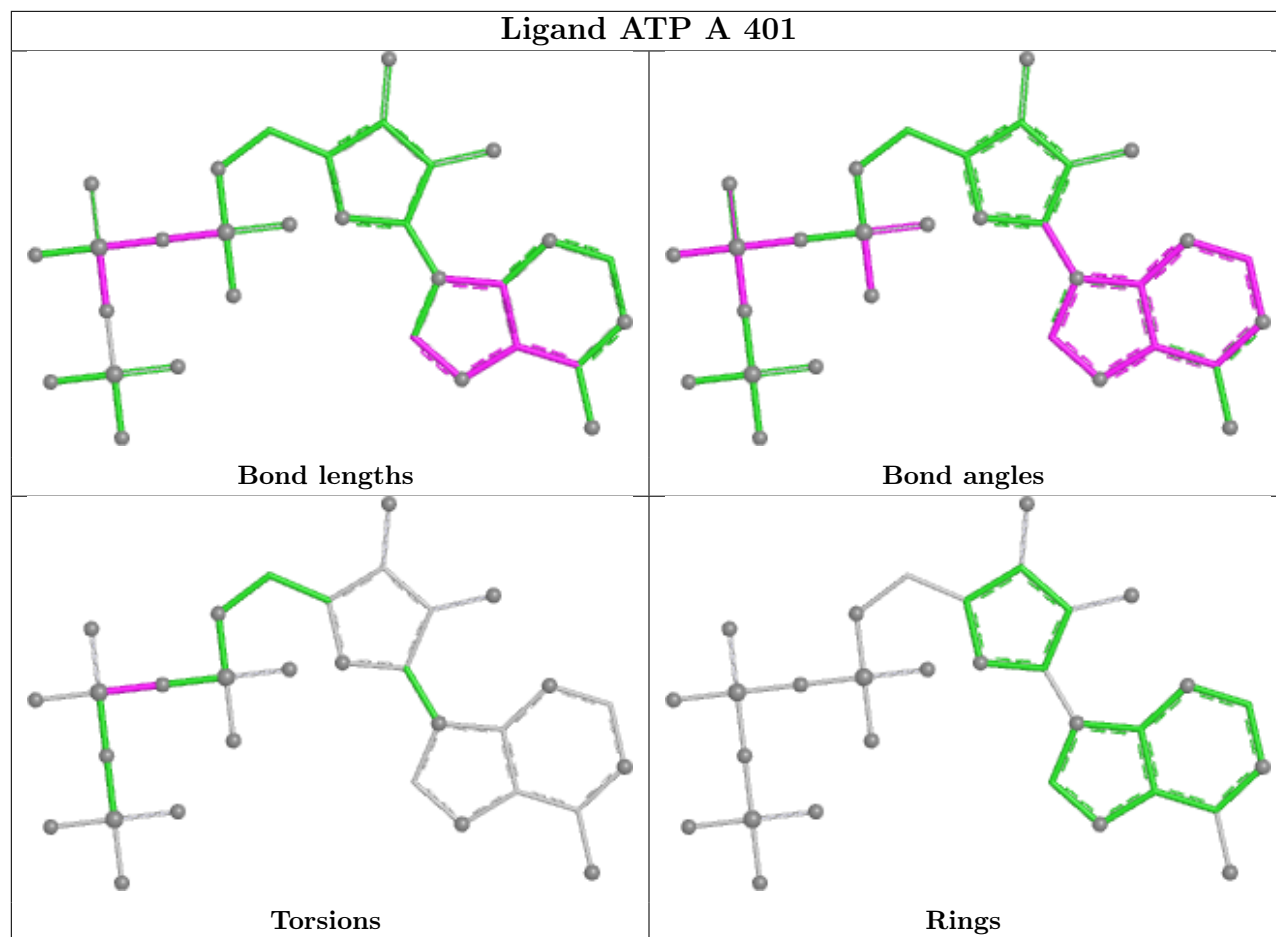
There are no ring outliers.

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 ⓘ

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	343/359 (95%)	0.34	27 (7%) 18 20	3, 20, 39, 47	13 (3%)
1	B	350/359 (97%)	0.12	16 (4%) 37 40	4, 15, 34, 53	13 (3%)
All	All	693/718 (96%)	0.23	43 (6%) 26 28	3, 17, 37, 53	26 (3%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	THR	7.6
1	B	351	THR	5.6
1	B	89	GLY	5.0
1	A	330	LEU	4.3
1	B	90	VAL	3.9
1	A	283	TYR	3.7
1	B	295	ILE	3.7
1	A	15	GLN	3.6
1	A	320	LEU	3.6
1	A	16	ASN	3.5
1	B	290	ALA	3.3
1	B	293	ALA	3.2
1	A	273	ARG	3.1
1	A	271	GLY	3.0
1	A	233	LEU	2.9
1	A	270	ALA	2.8
1	B	294	SER	2.7
1	A	18	PRO	2.6
1	A	245	PHE	2.6
1	A	331	SER	2.6
1	A	318	THR	2.5
1	A	321	THR	2.5
1	B	16	ASN	2.4
1	A	6[A]	HIS	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	289	SER	2.4
1	A	333	GLY	2.4
1	B	348	PRO	2.3
1	A	258	ARG	2.3
1	B	2	THR	2.3
1	A	268	ALA	2.3
1	A	290	ALA	2.3
1	B	349	GLY	2.3
1	A	67	THR	2.3
1	A	269	PRO	2.2
1	A	243	ILE	2.2
1	A	303	THR	2.2
1	B	88	GLU	2.2
1	A	89	GLY	2.2
1	A	259	LEU	2.1
1	B	57	ASP	2.1
1	B	332	ASP	2.1
1	B	15	GLN	2.0
1	A	306	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 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($\text{\AA}^2$)	Q<0.9
3	CA	A	361	1/1	0.92	0.13	46,46,46,46	0
2	ATP	B	360	31/31	0.99	0.04	2,9,13,16	0
2	ATP	A	401	31/31	0.99	0.04	5,14,21,22	0
3	CA	A	360	1/1	1.00	0.01	9,9,9,9	0

Continued on next page...

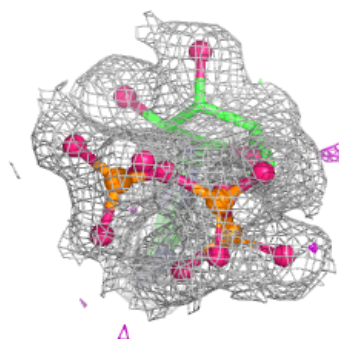
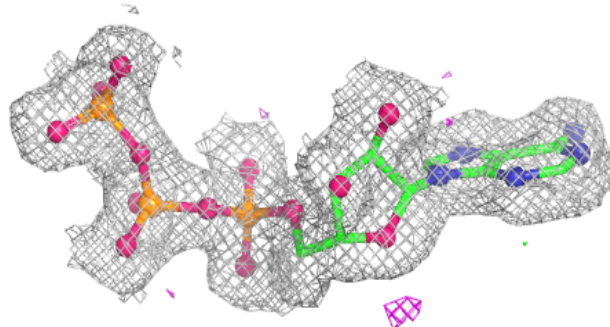
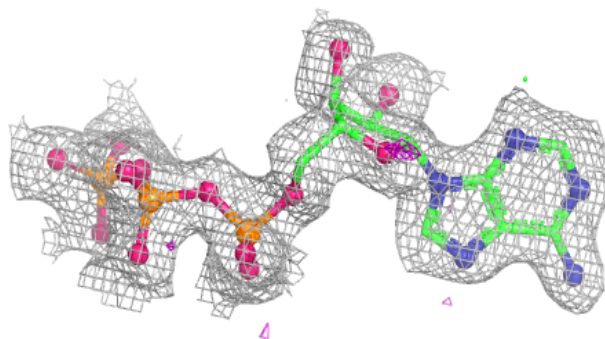
Continued from previous page...

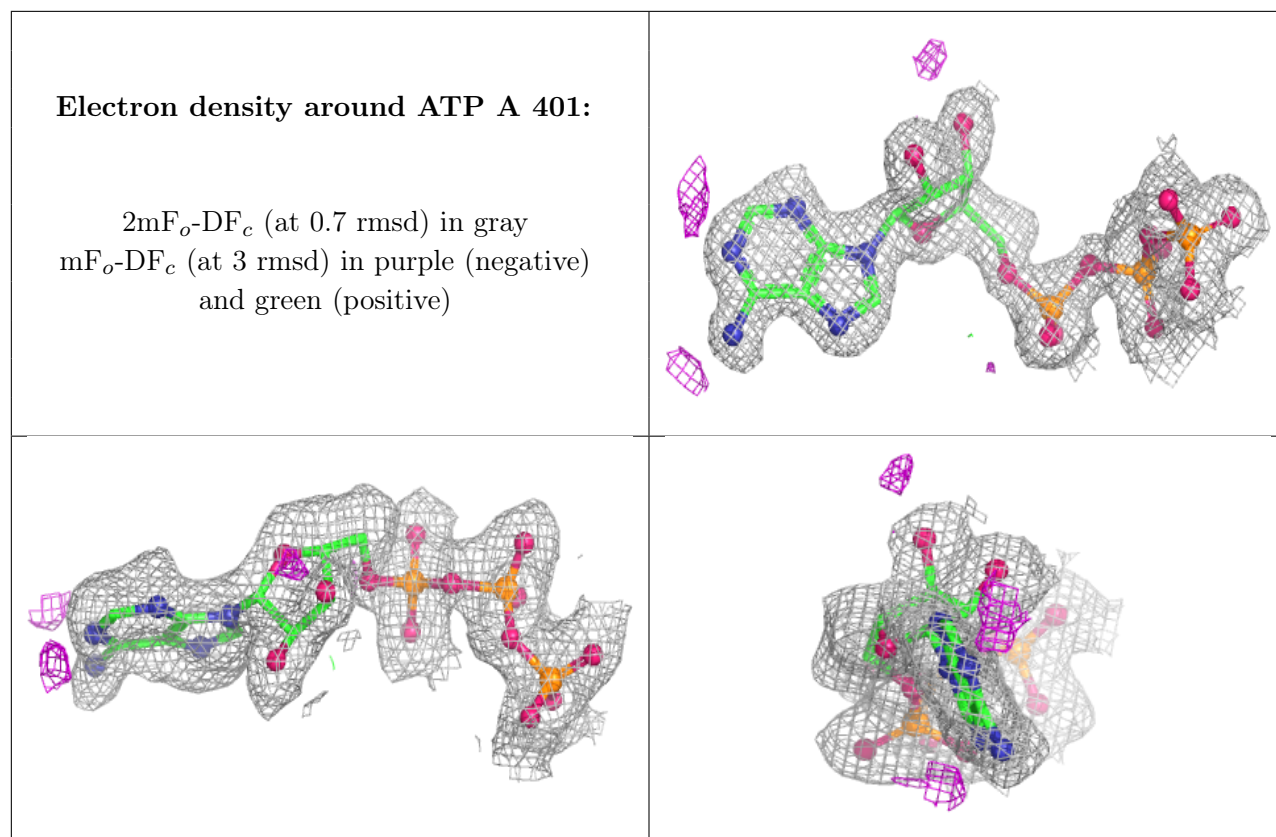
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CA	B	361	1/1	1.00	0.01	5,5,5,5	0

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 360:

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.