



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 05:25 PM UTC

PDB ID : 2FGK / pdb_00002fgk
Title : Crystal structure of the ABC-cassette E631Q mutant of HlyB with bound ATP
Authors : Zaitseva, J.; Oswald, C.; Jumpertz, T.; Jenewein, S.; Holland, I.B.; Schmitt, L.
Deposited on : 2005-12-22
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 ⓘ](#)) 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

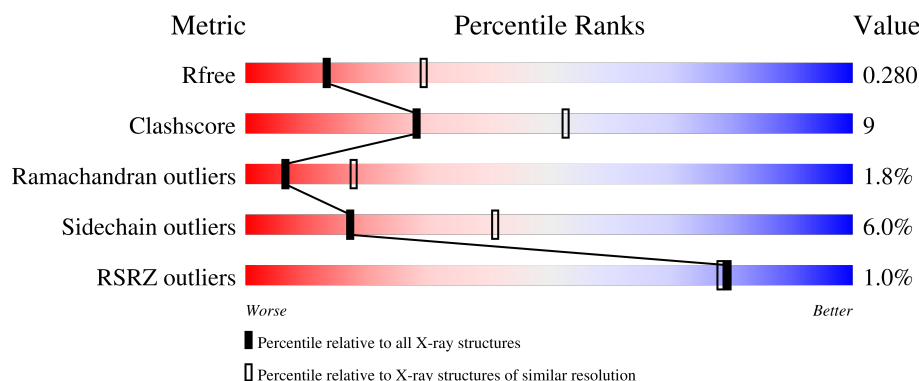
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

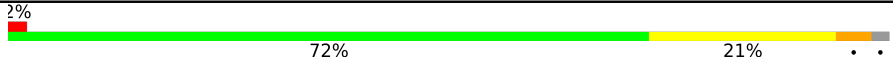
The reported resolution of this entry is 2.70 Å.

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	247	
1	B	247	
1	C	247	
1	D	247	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7848 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 Alpha-hemolysin translocation ATP-binding protein hlyB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	0	0	0
			1893	1194	342	352	5			
1	B	241	Total	C	N	O	S	0	0	0
			1893	1194	342	352	5			
1	C	241	Total	C	N	O	S	0	0	0
			1893	1194	342	352	5			
1	D	241	Total	C	N	O	S	0	0	0
			1893	1194	342	352	5			

There are 28 discrepancies between the modelled and reference sequences:

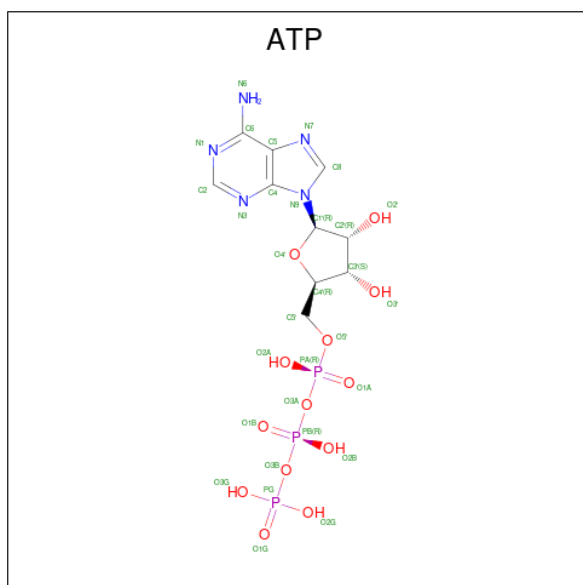
Chain	Residue	Modelled	Actual	Comment	Reference
A	461	HIS	-	expression tag	UNP P08716
A	462	HIS	-	expression tag	UNP P08716
A	463	HIS	-	expression tag	UNP P08716
A	464	HIS	-	expression tag	UNP P08716
A	465	HIS	-	expression tag	UNP P08716
A	466	HIS	-	expression tag	UNP P08716
A	631	GLN	GLU	engineered mutation	UNP P08716
B	461	HIS	-	expression tag	UNP P08716
B	462	HIS	-	expression tag	UNP P08716
B	463	HIS	-	expression tag	UNP P08716
B	464	HIS	-	expression tag	UNP P08716
B	465	HIS	-	expression tag	UNP P08716
B	466	HIS	-	expression tag	UNP P08716
B	631	GLN	GLU	engineered mutation	UNP P08716
C	461	HIS	-	expression tag	UNP P08716
C	462	HIS	-	expression tag	UNP P08716
C	463	HIS	-	expression tag	UNP P08716
C	464	HIS	-	expression tag	UNP P08716
C	465	HIS	-	expression tag	UNP P08716
C	466	HIS	-	expression tag	UNP P08716
C	631	GLN	GLU	engineered mutation	UNP P08716

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Chain	Residue	Modelled	Actual	Comment	Reference
D	461	HIS	-	expression tag	UNP P08716
D	462	HIS	-	expression tag	UNP P08716
D	463	HIS	-	expression tag	UNP P08716
D	464	HIS	-	expression tag	UNP P08716
D	465	HIS	-	expression tag	UNP P08716
D	466	HIS	-	expression tag	UNP P08716
D	631	GLN	GLU	engineered mutation	UNP P08716

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: 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		
2	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is water.

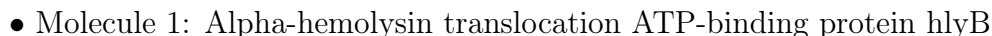
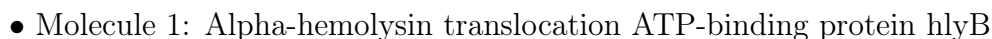
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	36	Total	O	0	0
			36	36		
3	B	44	Total	O	0	0
			44	44		

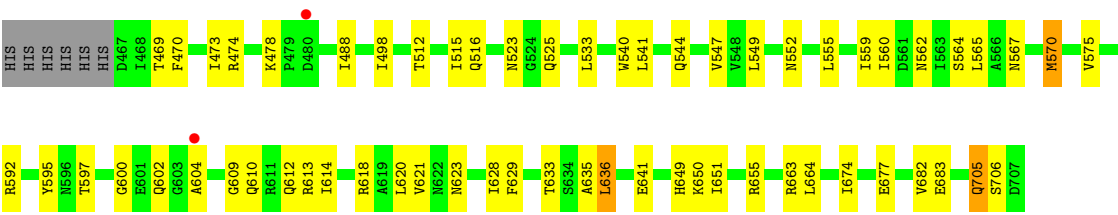
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	32	Total 32	O 32	0	0
3	D	40	Total 40	O 40	0	0

- Molecule 1: Alpha-hemolysin translocation ATP-binding protein hlyB





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	47.14Å 189.26Å 63.48Å 90.00° 111.87° 90.00°	Depositor
Resolution (Å)	20.00 – 2.70 20.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.9 (20.00-2.70) 98.8 (20.00-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.53 (at 2.71Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.223 , 0.280 0.222 , 0.280	Depositor DCC
R_{free} test set	1406 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 66.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.406 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7848	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.71% 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

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	0.95	7/1920 (0.4%)	0.87	0/2589
1	B	0.56	0/1920	0.86	2/2589 (0.1%)
1	C	0.74	8/1920 (0.4%)	0.86	0/2589
1	D	0.55	0/1920	0.83	0/2589
All	All	0.72	15/7680 (0.2%)	0.86	2/10356 (0.0%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	593	GLU	CD-OE2	25.20	1.73	1.25
1	A	557	ARG	NE-CZ	14.00	1.48	1.33
1	A	593	GLU	CD-OE1	10.76	1.45	1.25
1	C	558	SER	CB-OG	9.44	1.61	1.42
1	A	557	ARG	CZ-NH1	8.28	1.44	1.32
1	C	558	SER	C-O	7.60	1.34	1.23
1	A	595	TYR	CE2-CZ	7.18	1.55	1.38
1	C	595	TYR	CE2-CZ	7.11	1.55	1.38
1	A	596	ASN	CG-OD1	7.10	1.37	1.23
1	C	557	ARG	CD-NE	6.93	1.55	1.46
1	C	557	ARG	CZ-NH1	6.89	1.42	1.32
1	C	570	MET	C-O	5.87	1.32	1.23
1	C	570	MET	C-N	5.41	1.41	1.33
1	A	595	TYR	CG-CD2	5.25	1.50	1.39
1	C	589	SER	CB-OG	5.24	1.52	1.42

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	605	GLY	N-CA-C	-8.56	104.42	113.58
1	B	600	GLY	N-CA-C	5.70	119.89	112.25

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	1893	0	1949	33	0
1	B	1893	0	1949	38	0
1	C	1893	0	1949	31	0
1	D	1893	0	1949	39	0
2	A	31	0	12	0	0
2	B	31	0	12	1	0
2	C	31	0	12	0	0
2	D	31	0	12	0	0
3	A	36	0	0	5	0
3	B	44	0	0	2	0
3	C	32	0	0	3	0
3	D	40	0	0	1	0
All	All	7848	0	7844	138	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 9.

All (138) 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:593:GLU:CD	1:A:593:GLU:OE2	1.73	1.30
1:C:571:SER:HB2	1:C:574:LYS:HB2	1.36	1.06
1:B:612:GLN:HE22	1:B:635:ALA:H	1.13	0.90
1:C:560:ILE:HB	3:C:139:HOH:O	1.73	0.88
1:C:612:GLN:HE22	1:C:635:ALA:H	1.27	0.83
1:A:612:GLN:HE22	1:A:635:ALA:H	1.27	0.81
1:A:582:ALA:O	1:A:613:ARG:HG2	1.85	0.77
1:C:571:SER:CB	1:C:574:LYS:HB2	2.15	0.76
1:B:623:ASN:HD22	1:B:655:ARG:HH12	1.36	0.72
1:A:557:ARG:HB2	1:A:561:ASP:OD2	1.89	0.72
1:A:498:ILE:HD11	1:A:674:ILE:HG13	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:600:GLY:H	1:A:604:ALA:H	1.39	0.70
1:A:612:GLN:HE21	1:A:636:LEU:HD13	1.55	0.70
1:D:612:GLN:HE22	1:D:635:ALA:H	1.38	0.70
1:B:512:THR:HA	1:B:515:ILE:HD12	1.74	0.70
1:C:582:ALA:O	1:C:613:ARG:HG2	1.91	0.69
1:C:571:SER:HB2	1:C:574:LYS:CB	2.21	0.68
1:C:498:ILE:HD11	1:C:674:ILE:HG13	1.75	0.67
1:A:555:LEU:H	1:A:562:ASN:HD21	1.42	0.66
1:D:512:THR:HA	1:D:515:ILE:HD12	1.78	0.66
1:B:612:GLN:HE21	1:B:636:LEU:HD13	1.60	0.65
1:D:623:ASN:HD22	1:D:655:ARG:HH12	1.44	0.65
1:A:571:SER:HB2	1:A:574:LYS:HB2	1.78	0.65
1:D:592:ARG:HG2	3:D:9:HOH:O	1.95	0.65
1:C:595:TYR:HB3	3:C:139:HOH:O	1.96	0.64
1:C:612:GLN:HE21	1:C:636:LEU:HD13	1.62	0.64
1:D:609:GLY:O	1:D:613:ARG:HD3	1.98	0.63
1:B:609:GLY:O	1:B:613:ARG:HD3	1.97	0.63
1:D:549:LEU:H	1:D:552:ASN:HD21	1.44	0.63
1:A:570:MET:HE1	1:A:621:VAL:HG13	1.82	0.62
1:B:612:GLN:NE2	1:B:635:ALA:H	1.92	0.62
1:D:623:ASN:ND2	1:D:655:ARG:HH12	2.00	0.60
1:A:561:ASP:HB3	1:C:561:ASP:OD1	2.01	0.59
1:A:558:SER:HB3	1:A:561:ASP:OD1	2.04	0.57
1:D:540:TRP:O	1:D:544:GLN:HG2	2.05	0.57
1:A:584:ALA:HB2	1:A:613:ARG:HB3	1.86	0.56
1:D:612:GLN:HE21	1:D:636:LEU:HD13	1.71	0.56
1:D:567:ASN:ND2	1:D:570:MET:HB2	2.19	0.56
1:D:641:GLU:OE2	1:D:663:ARG:HD2	2.05	0.56
1:B:623:ASN:ND2	1:B:655:ARG:HH12	2.02	0.55
1:D:555:LEU:H	1:D:562:ASN:HD21	1.54	0.55
1:B:540:TRP:O	1:B:544:GLN:HG2	2.08	0.54
1:B:641:GLU:OE2	1:B:663:ARG:HD2	2.07	0.54
1:C:568:PRO:C	1:C:570:MET:N	2.64	0.54
1:D:567:ASN:HD22	1:D:570:MET:HB2	1.73	0.53
1:C:584:ALA:HB2	1:C:613:ARG:HB3	1.89	0.53
1:C:612:GLN:NE2	1:C:635:ALA:H	2.02	0.53
1:A:568:PRO:HD3	1:C:568:PRO:HD3	1.89	0.53
1:B:560:ILE:HG13	1:B:575:VAL:HG11	1.90	0.52
1:D:555:LEU:N	1:D:562:ASN:HD21	2.06	0.52
1:B:512:THR:HB	1:B:628:ILE:HD13	1.92	0.52
1:C:600:GLY:HA3	1:C:604:ALA:HB3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:512:THR:HB	1:D:628:ILE:HD13	1.92	0.52
1:B:682:VAL:HG12	1:B:683:GLU:HG3	1.92	0.51
1:D:682:VAL:HG12	1:D:683:GLU:HG3	1.93	0.50
1:D:549:LEU:N	1:D:552:ASN:HD21	2.09	0.50
1:C:555:LEU:H	1:C:562:ASN:HD21	1.59	0.50
1:A:498:ILE:HD12	3:A:142:HOH:O	2.12	0.49
1:C:703:GLN:HB3	3:C:14:HOH:O	2.12	0.49
1:A:496:GLU:CD	3:A:142:HOH:O	2.55	0.49
1:B:474:ARG:HB3	1:B:523:ASN:HB3	1.95	0.49
1:D:600:GLY:HA3	1:D:604:ALA:O	2.13	0.48
1:B:592:ARG:HG2	3:B:150:HOH:O	2.12	0.48
1:B:550:GLN:HA	3:B:40:HOH:O	2.13	0.48
1:C:568:PRO:C	1:C:570:MET:H	2.22	0.48
1:D:470:PHE:HD1	1:D:473:ILE:HG13	1.78	0.48
1:B:588:ILE:O	1:B:594:GLY:HA2	2.14	0.48
1:D:474:ARG:HB3	1:D:523:ASN:HB3	1.96	0.48
1:D:705:GLN:O	1:D:705:GLN:HG3	2.13	0.48
1:B:555:LEU:H	1:B:562:ASN:HD21	1.62	0.47
1:D:498:ILE:HD11	1:D:674:ILE:HG13	1.96	0.47
1:A:588:ILE:HA	1:A:591:LEU:HD12	1.95	0.47
1:B:705:GLN:O	1:B:705:GLN:HG3	2.13	0.47
1:C:512:THR:HB	1:C:628:ILE:HD13	1.97	0.47
1:A:468:ILE:HB	1:A:492:ILE:HB	1.97	0.47
1:B:567:ASN:ND2	1:B:570:MET:HB2	2.30	0.47
1:D:562:ASN:O	1:D:618:ARG:HG3	2.15	0.47
1:C:558:SER:HA	1:C:598:ILE:HA	1.97	0.47
1:D:633:THR:HA	1:D:636:LEU:HD22	1.96	0.47
1:A:603:GLY:O	1:A:605:GLY:N	2.48	0.46
1:B:470:PHE:HD1	1:B:473:ILE:HG13	1.80	0.46
1:C:623:ASN:ND2	1:C:655:ARG:HH22	2.13	0.46
1:A:512:THR:HB	1:A:628:ILE:HD13	1.96	0.46
1:A:548:VAL:O	1:A:630:ASP:HB3	2.16	0.46
1:A:478:LYS:HB3	1:A:479:PRO:HD2	1.97	0.46
1:B:649:HIS:CE1	1:B:650:LYS:HE2	2.50	0.46
1:B:498:ILE:HD11	1:B:674:ILE:CG1	2.46	0.46
1:C:468:ILE:HB	1:C:492:ILE:HB	1.99	0.45
1:B:588:ILE:HD11	1:B:614:ILE:CD1	2.46	0.45
1:B:498:ILE:HD11	1:B:674:ILE:HG13	1.98	0.45
1:B:516:GLN:CD	1:B:547:VAL:HG21	2.41	0.45
1:B:555:LEU:N	1:B:562:ASN:HD21	2.14	0.45
1:D:560:ILE:HG13	1:D:575:VAL:HG11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:599:VAL:HG11	1:B:606:LEU:HD21	1.98	0.45
1:B:599:VAL:CG1	1:B:606:LEU:HD21	2.47	0.45
1:B:572:VAL:HG12	1:B:576:ILE:HD13	1.99	0.45
1:C:478:LYS:HB3	1:C:479:PRO:HD2	1.98	0.44
1:D:559:ILE:HD12	1:D:597:THR:HB	1.98	0.44
1:A:476:ARG:HD3	1:A:481:SER:O	2.17	0.44
1:A:546:GLY:HA3	1:A:624:PRO:HG3	1.99	0.44
1:A:665:SER:HA	1:A:668:LYS:HG3	1.99	0.44
1:C:546:GLY:HA3	1:C:624:PRO:HG3	1.99	0.44
1:D:498:ILE:HD11	1:D:674:ILE:CG1	2.47	0.44
1:D:549:LEU:H	1:D:552:ASN:ND2	2.14	0.44
1:D:620:LEU:HD21	1:D:651:ILE:HG23	2.00	0.44
1:D:610:GLN:O	1:D:614:ILE:HG12	2.18	0.43
1:D:560:ILE:HB	1:D:595:TYR:HB3	2.00	0.43
1:B:633:THR:HA	1:B:636:LEU:HD22	2.00	0.43
1:C:548:VAL:O	1:C:630:ASP:HB3	2.18	0.43
1:C:476:ARG:HD3	1:C:481:SER:O	2.19	0.43
1:D:533:LEU:HD22	1:D:541:LEU:HD13	2.00	0.43
1:D:612:GLN:NE2	1:D:635:ALA:H	2.12	0.43
1:C:665:SER:HA	1:C:668:LYS:HG3	2.01	0.43
1:A:688:LYS:HE2	3:A:41:HOH:O	2.19	0.43
1:A:556:ASN:C	1:A:557:ARG:HG2	2.44	0.42
1:B:504:SER:HA	2:B:801:ATP:O2G	2.18	0.42
1:B:677:GLU:HB3	1:B:682:VAL:HG21	2.01	0.42
1:D:473:ILE:HB	1:D:488:ILE:HB	2.01	0.42
1:B:560:ILE:HB	1:B:595:TYR:HB3	2.01	0.42
1:B:570:MET:HE2	1:B:570:MET:HB3	1.76	0.42
1:C:610:GLN:O	1:C:614:ILE:HG12	2.18	0.42
1:B:587:PHE:C	1:B:587:PHE:CD1	2.97	0.42
1:A:618:ARG:HD2	3:A:63:HOH:O	2.19	0.42
1:C:567:ASN:HA	1:C:568:PRO:HD3	1.80	0.42
1:B:533:LEU:HD22	1:B:541:LEU:HD13	2.01	0.42
1:D:570:MET:HE2	1:D:570:MET:HB3	1.85	0.42
1:B:473:ILE:HB	1:B:488:ILE:HB	2.02	0.42
1:D:516:GLN:CD	1:D:547:VAL:HG21	2.44	0.42
1:A:557:ARG:HH22	1:C:568:PRO:CB	2.33	0.41
1:A:492:ILE:HG23	3:A:142:HOH:O	2.20	0.41
1:A:584:ALA:HA	1:A:587:PHE:CE2	2.55	0.41
1:C:549:LEU:H	1:C:552:ASN:HD21	1.68	0.41
1:D:649:HIS:CE1	1:D:650:LYS:HE2	2.55	0.41
1:A:587:PHE:CD1	1:A:587:PHE:C	2.98	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:547:VAL:HG22	1:D:628:ILE:HB	2.03	0.41
1:D:677:GLU:HB3	1:D:682:VAL:HG21	2.02	0.41
1:A:629:PHE:CE2	1:A:648:MET:HE1	2.57	0.40
1:B:588:ILE:HD11	1:B:614:ILE:HD12	2.04	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	239/247 (97%)	223 (93%)	10 (4%)	6 (2%)	4	11
1	B	239/247 (97%)	226 (95%)	10 (4%)	3 (1%)	9	25
1	C	239/247 (97%)	222 (93%)	11 (5%)	6 (2%)	4	11
1	D	239/247 (97%)	224 (94%)	13 (5%)	2 (1%)	16	37
All	All	956/988 (97%)	895 (94%)	44 (5%)	17 (2%)	6	18

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	603	GLY
1	C	601	GLU
1	C	571	SER
1	D	565	LEU
1	A	604	ALA
1	C	568	PRO
1	C	594	GLY
1	A	567	ASN
1	A	571	SER
1	A	706	SER
1	B	551	ASP

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Mol	Chain	Res	Type
1	B	601	GLU
1	C	567	ASN
1	C	706	SER
1	B	706	SER
1	D	706	SER
1	A	572	VAL

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	207/213 (97%)	191 (92%)	16 (8%)	12	30
1	B	207/213 (97%)	196 (95%)	11 (5%)	20	46
1	C	207/213 (97%)	195 (94%)	12 (6%)	18	42
1	D	207/213 (97%)	196 (95%)	11 (5%)	20	46
All	All	828/852 (97%)	778 (94%)	50 (6%)	17	41

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	478	LYS
1	A	549	LEU
1	A	552	ASN
1	A	553	VAL
1	A	556	ASN
1	A	561	ASP
1	A	576	ILE
1	A	586	ASP
1	A	599	VAL
1	A	606	LEU
1	A	613	ARG
1	A	623	ASN
1	A	636	LEU
1	A	664	LEU
1	A	682	VAL

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Mol	Chain	Res	Type
1	A	693	GLU
1	B	469	THR
1	B	478	LYS
1	B	525	GLN
1	B	557	ARG
1	B	564	SER
1	B	565	LEU
1	B	621	VAL
1	B	629	PHE
1	B	636	LEU
1	B	664	LEU
1	B	705	GLN
1	C	478	LYS
1	C	556	ASN
1	C	576	ILE
1	C	586	ASP
1	C	602	GLN
1	C	606	LEU
1	C	613	ARG
1	C	618	ARG
1	C	636	LEU
1	C	664	LEU
1	C	682	VAL
1	C	693	GLU
1	D	469	THR
1	D	478	LYS
1	D	525	GLN
1	D	564	SER
1	D	570	MET
1	D	602	GLN
1	D	621	VAL
1	D	629	PHE
1	D	636	LEU
1	D	664	LEU
1	D	705	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	539	ASN
1	A	562	ASN
1	A	567	ASN

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Mol	Chain	Res	Type
1	A	612	GLN
1	A	705	GLN
1	B	487	ASN
1	B	531	HIS
1	B	539	ASN
1	B	562	ASN
1	B	567	ASN
1	B	612	GLN
1	B	623	ASN
1	B	642	HIS
1	B	669	ASN
1	B	705	GLN
1	C	539	ASN
1	C	562	ASN
1	C	612	GLN
1	C	623	ASN
1	C	705	GLN
1	D	487	ASN
1	D	531	HIS
1	D	539	ASN
1	D	552	ASN
1	D	556	ASN
1	D	562	ASN
1	D	567	ASN
1	D	612	GLN
1	D	623	ASN
1	D	642	HIS
1	D	669	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

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	ATP	A	800	-	32,33,33	1.27	4 (12%)	48,52,52	1.87	10 (20%)
2	ATP	B	801	-	32,33,33	1.36	4 (12%)	48,52,52	1.85	9 (18%)
2	ATP	D	803	-	32,33,33	1.62	4 (12%)	48,52,52	1.82	10 (20%)
2	ATP	C	802	-	32,33,33	1.40	6 (18%)	48,52,52	1.82	10 (20%)

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	A	800	-	-	2/22/38/38	0/3/3/3
2	ATP	B	801	-	-	4/22/38/38	0/3/3/3
2	ATP	D	803	-	-	3/22/38/38	0/3/3/3
2	ATP	C	802	-	-	3/22/38/38	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	802	ATP	PA-O3A	4.71	1.64	1.59
2	D	803	ATP	PB-O3B	4.43	1.64	1.59
2	D	803	ATP	PB-O3A	3.98	1.63	1.59
2	B	801	ATP	PG-O1G	3.68	1.61	1.50
2	D	803	ATP	PG-O1G	3.49	1.61	1.50
2	B	801	ATP	C5-N7	-3.31	1.33	1.39
2	A	800	ATP	PG-O1G	3.27	1.60	1.50
2	A	800	ATP	C5-N7	-3.17	1.33	1.39
2	D	803	ATP	C5-N7	-3.07	1.33	1.39
2	B	801	ATP	PB-O3B	3.03	1.62	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	802	ATP	C5-N7	-2.92	1.33	1.39
2	C	802	ATP	PG-O1G	2.37	1.57	1.50
2	A	800	ATP	C8-N9	-2.24	1.33	1.37
2	C	802	ATP	C8-N9	-2.17	1.33	1.37
2	C	802	ATP	PB-O3A	2.13	1.61	1.59
2	A	800	ATP	PB-O3B	2.11	1.61	1.59
2	B	801	ATP	C8-N9	-2.08	1.34	1.37
2	C	802	ATP	C4-N9	-2.06	1.33	1.37

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	ATP	C5-C4-N3	-5.81	118.72	126.72
2	D	803	ATP	C5-C4-N3	-5.74	118.81	126.72
2	A	800	ATP	C5-C4-N3	-5.54	119.08	126.72
2	C	802	ATP	C5-C4-N3	-4.99	119.85	126.72
2	C	802	ATP	N3-C2-N1	-4.63	121.57	128.58
2	A	800	ATP	N3-C2-N1	-4.41	121.91	128.58
2	D	803	ATP	N3-C2-N1	-4.26	122.13	128.58
2	B	801	ATP	N3-C2-N1	-4.22	122.20	128.58
2	A	800	ATP	O3G-PG-O3B	4.06	118.24	104.64
2	B	801	ATP	N3-C4-N9	4.04	134.05	127.17
2	A	800	ATP	N3-C4-N9	3.92	133.84	127.17
2	D	803	ATP	N3-C4-N9	3.85	133.71	127.17
2	B	801	ATP	O3G-PG-O3B	3.67	116.95	104.64
2	C	802	ATP	O3G-PG-O3B	3.67	116.95	104.64
2	A	800	ATP	C2-N3-C4	3.64	120.72	111.83
2	C	802	ATP	N3-C4-N9	3.58	133.26	127.17
2	D	803	ATP	C2-N3-C4	3.56	120.51	111.83
2	B	801	ATP	O4'-C1'-N9	-3.53	101.32	108.09
2	D	803	ATP	O3G-PG-O3B	3.51	116.40	104.64
2	B	801	ATP	C2-N3-C4	3.48	120.34	111.83
2	C	802	ATP	C2-N3-C4	3.44	120.24	111.83
2	A	800	ATP	N9-C8-N7	-3.35	109.19	113.94
2	C	802	ATP	N9-C8-N7	-3.23	109.36	113.94
2	A	800	ATP	C5-N7-C8	3.11	108.33	103.45
2	D	803	ATP	O4'-C1'-N9	-2.98	102.36	108.09
2	D	803	ATP	C5-N7-C8	2.92	108.03	103.45
2	A	800	ATP	C4-C5-N7	-2.90	107.27	110.58
2	B	801	ATP	C5-N7-C8	2.88	107.97	103.45
2	D	803	ATP	C4-C5-N7	-2.86	107.31	110.58
2	B	801	ATP	N9-C8-N7	-2.86	109.88	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	803	ATP	N9-C8-N7	-2.85	109.89	113.94
2	C	802	ATP	C5-N7-C8	2.82	107.89	103.45
2	B	801	ATP	C4-C5-N7	-2.76	107.42	110.58
2	C	802	ATP	C4-C5-N7	-2.72	107.47	110.58
2	C	802	ATP	O4'-C1'-N9	-2.71	102.88	108.09
2	C	802	ATP	C4-N9-C8	2.53	108.39	105.74
2	A	800	ATP	O4'-C1'-N9	-2.45	103.38	108.09
2	A	800	ATP	C4-N9-C8	2.28	108.13	105.74
2	D	803	ATP	O2B-PB-O3B	2.09	112.91	107.27

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	801	ATP	O4'-C4'-C5'-O5'
2	A	800	ATP	PA-O3A-PB-O1B
2	B	801	ATP	C3'-C4'-C5'-O5'
2	D	803	ATP	O4'-C4'-C5'-O5'
2	C	802	ATP	PA-O3A-PB-O1B
2	D	803	ATP	PA-O3A-PB-O2B
2	A	800	ATP	PA-O3A-PB-O2B
2	C	802	ATP	O4'-C4'-C5'-O5'
2	B	801	ATP	PA-O3A-PB-O2B
2	C	802	ATP	PA-O3A-PB-O2B
2	B	801	ATP	PA-O3A-PB-O1B
2	D	803	ATP	PA-O3A-PB-O1B

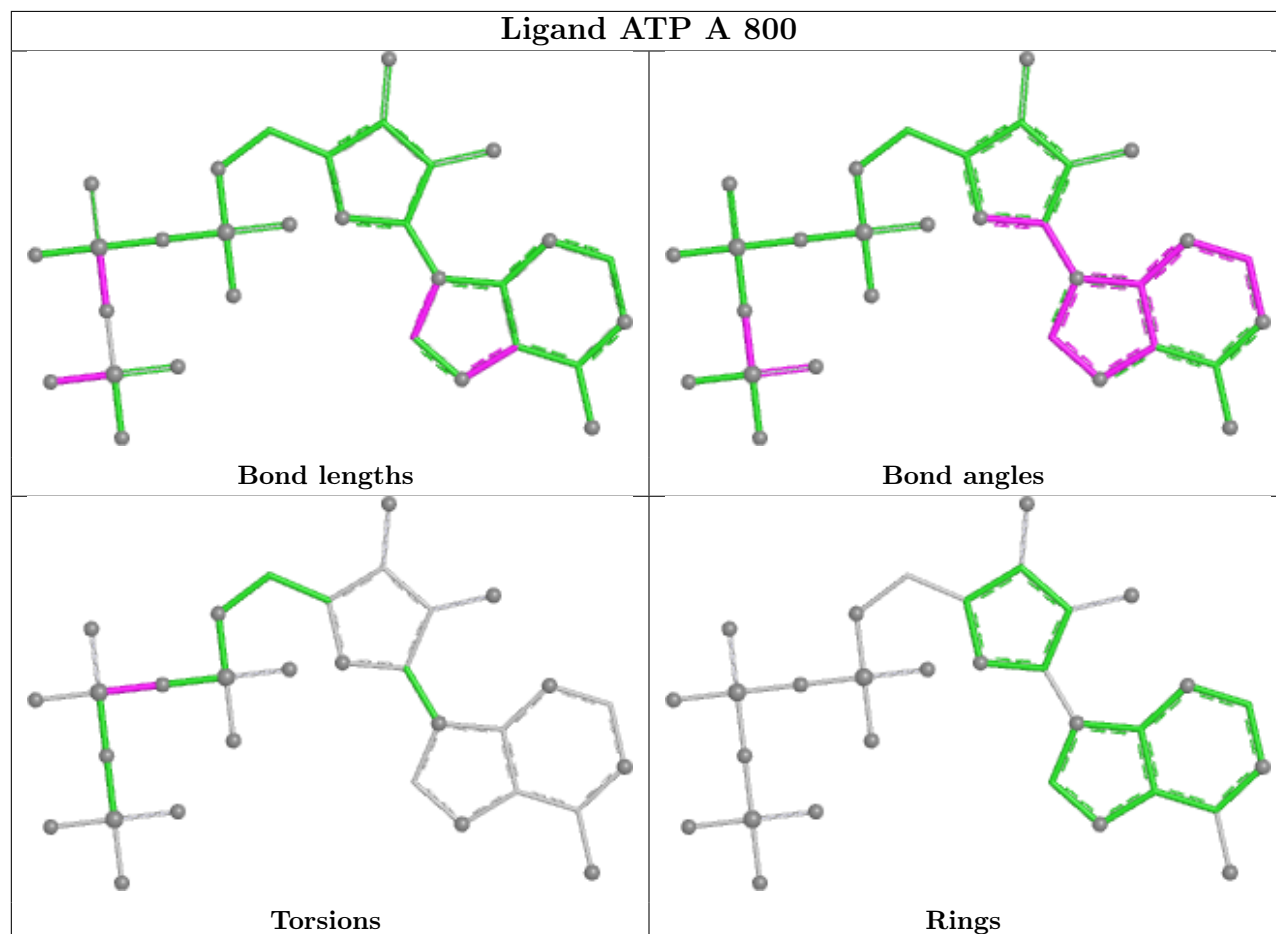
There are no ring outliers.

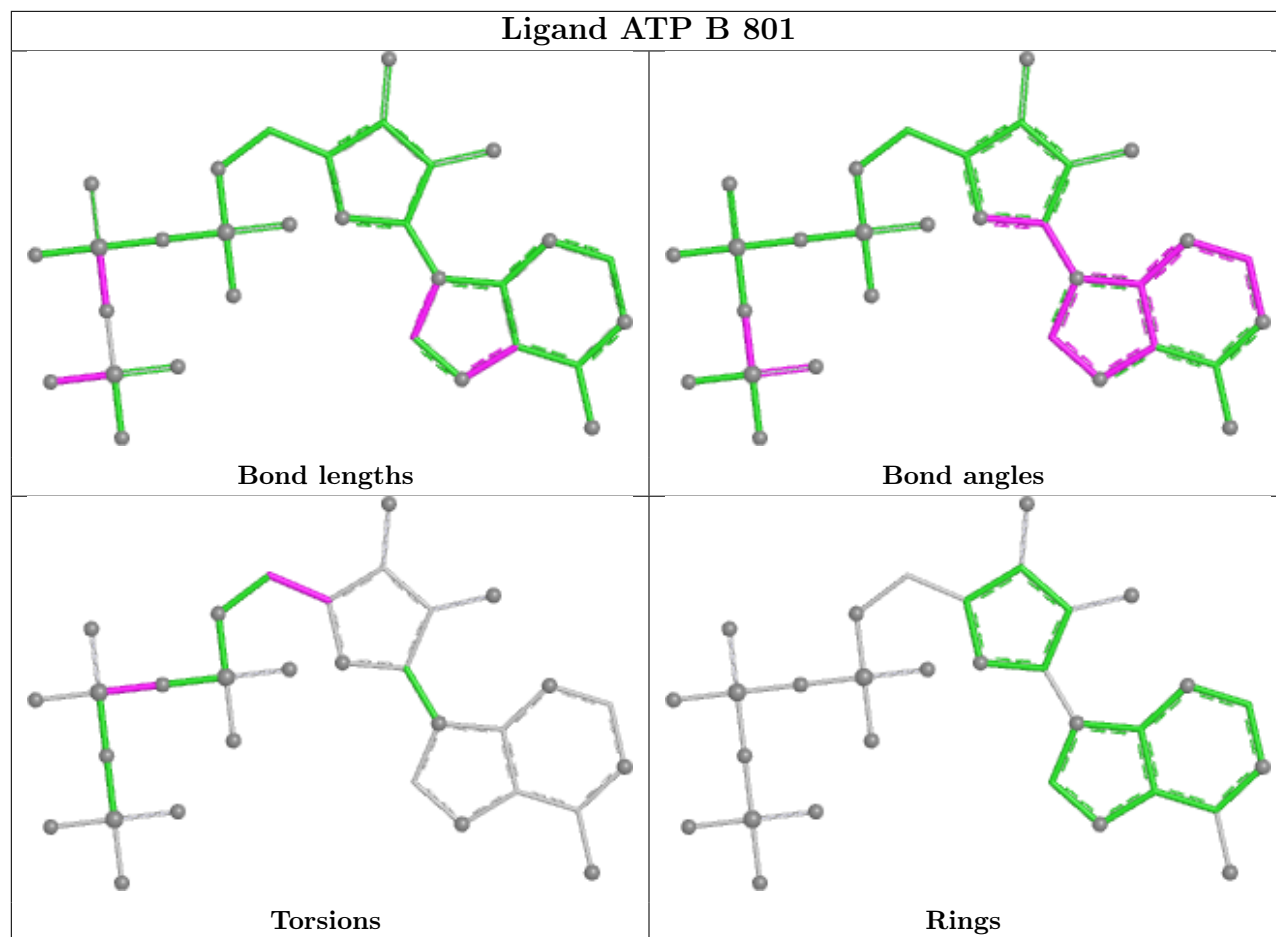
1 monomer is involved in 1 short contact:

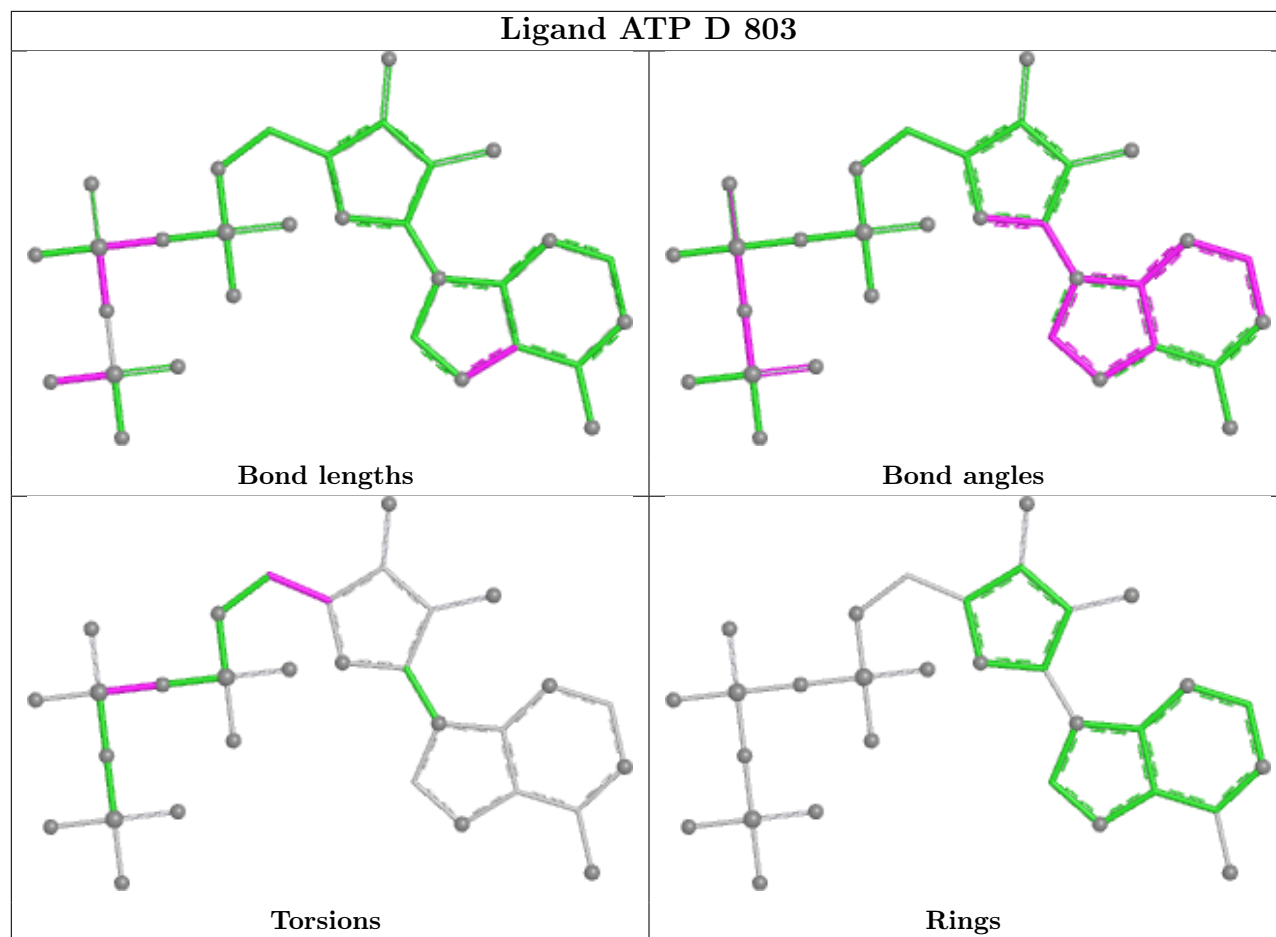
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	ATP	1	0

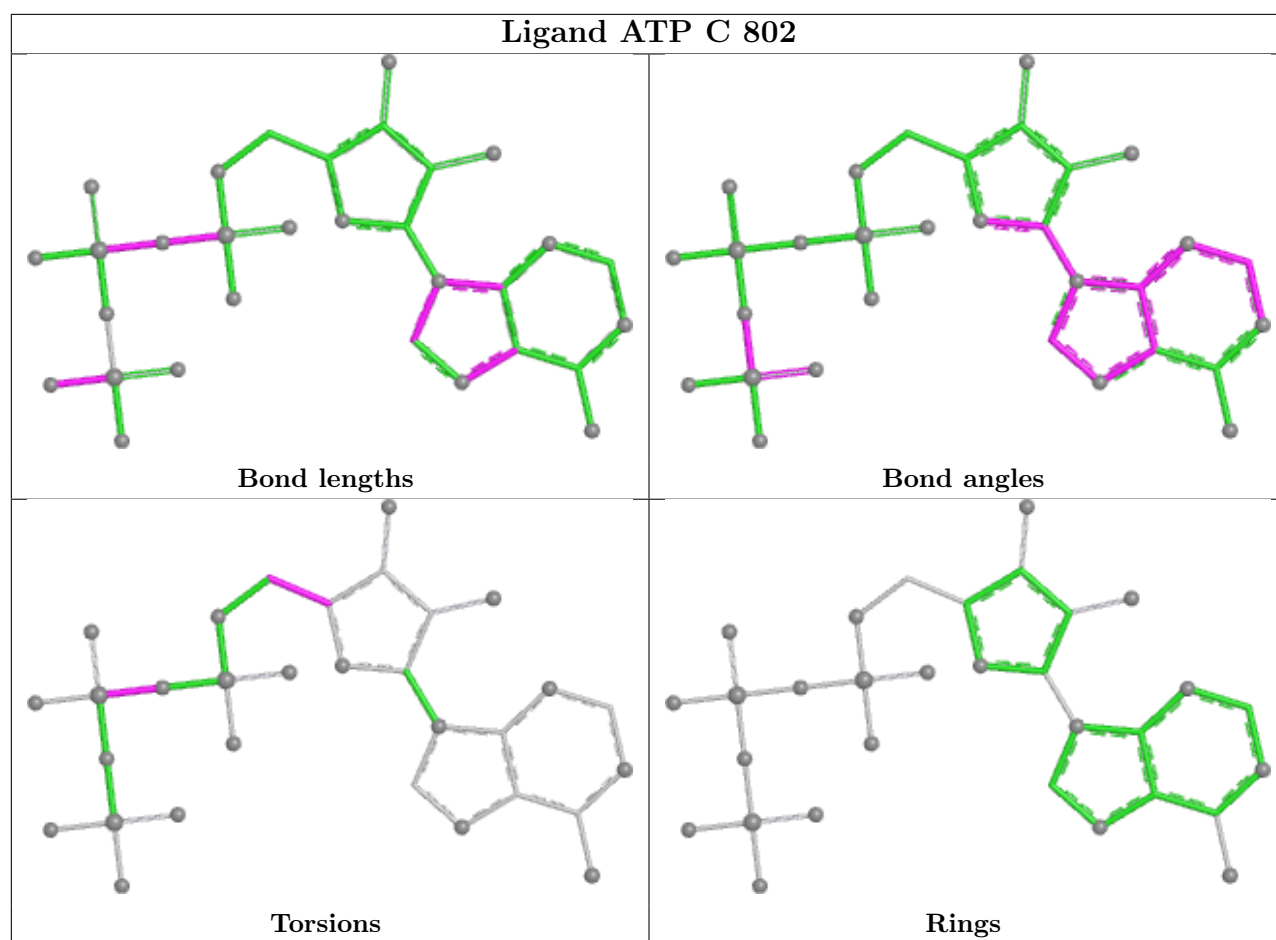
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 [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	241/247 (97%)	-0.24	4 (1%) 69 67	39, 51, 62, 71	0
1	B	241/247 (97%)	-0.68	2 (0%) 82 81	40, 51, 63, 70	0
1	C	241/247 (97%)	-0.43	2 (0%) 82 81	37, 50, 61, 71	0
1	D	241/247 (97%)	-0.63	2 (0%) 82 81	40, 51, 63, 71	0
All	All	964/988 (97%)	-0.49	10 (1%) 79 78	37, 51, 63, 71	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	597	THR	3.5
1	D	604	ALA	2.8
1	B	562	ASN	2.8
1	A	569	GLY	2.5
1	A	605	GLY	2.4
1	D	480	ASP	2.4
1	B	589	SER	2.3
1	A	607	SER	2.3
1	C	507	GLY	2.3
1	C	568	PRO	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](#)

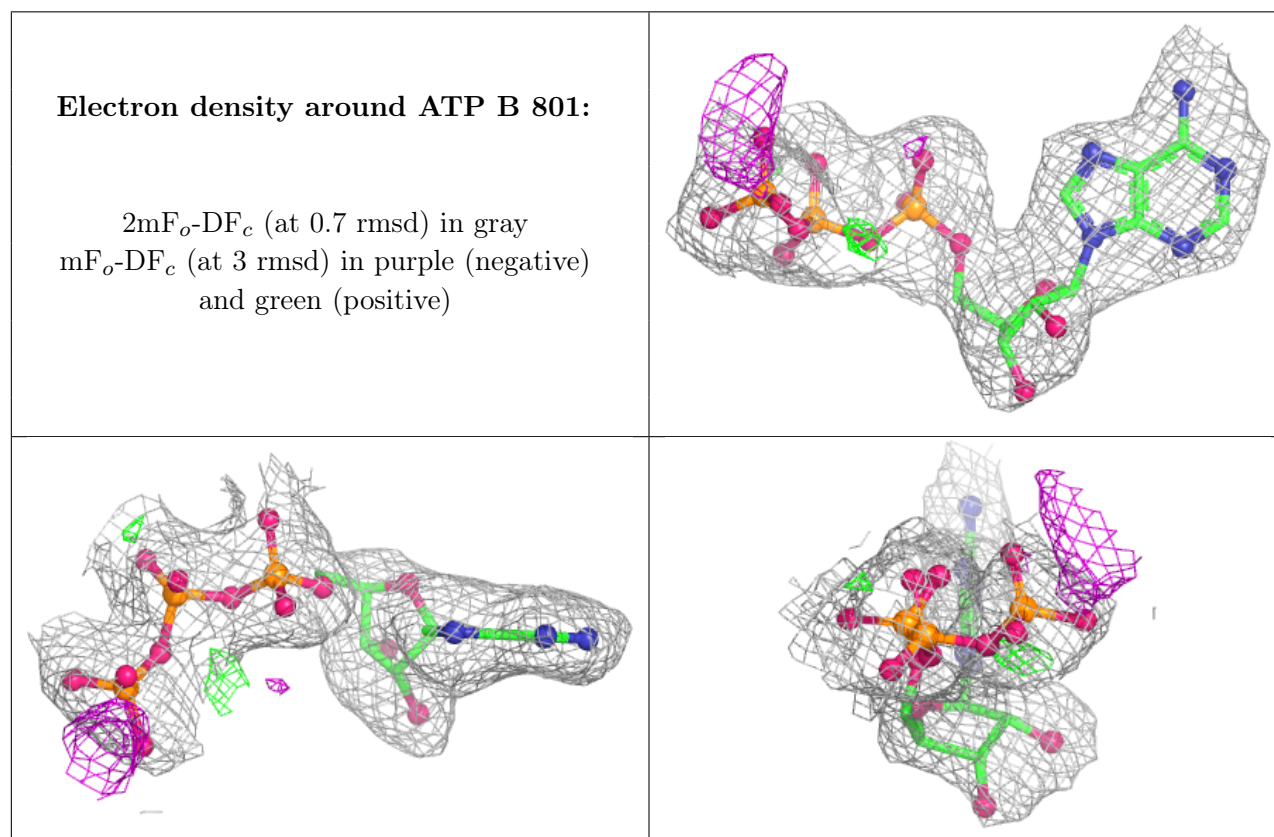
There are no oligosaccharides in this entry.

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.

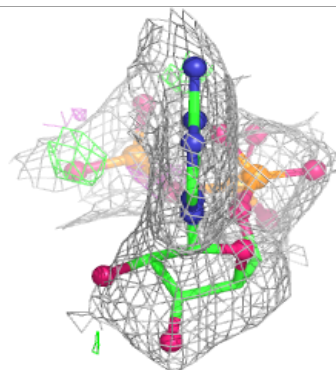
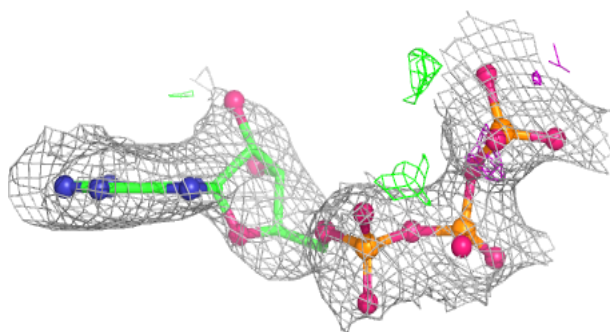
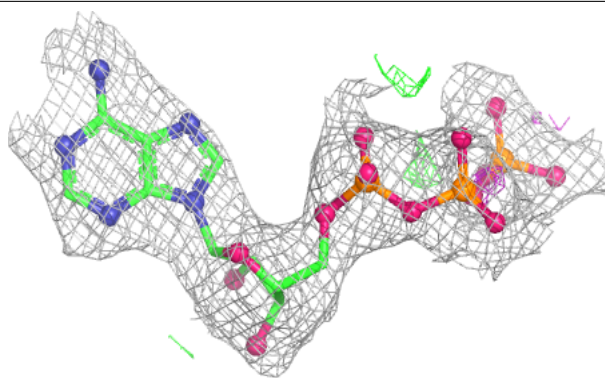
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ATP	B	801	31/31	0.98	0.04	37,42,44,45	0
2	ATP	D	803	31/31	0.98	0.05	38,46,47,50	0
2	ATP	C	802	31/31	0.99	0.04	28,40,43,44	0
2	ATP	A	800	31/31	0.99	0.05	32,41,44,44	0

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.

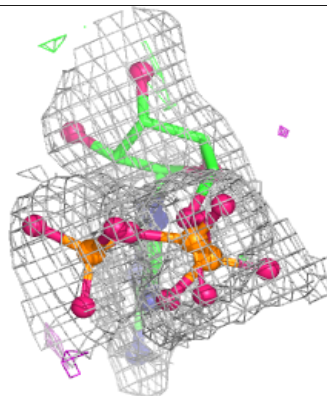
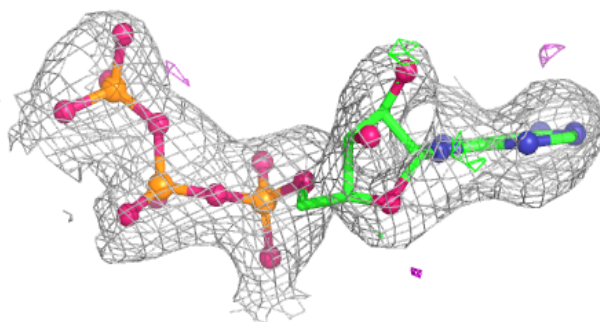
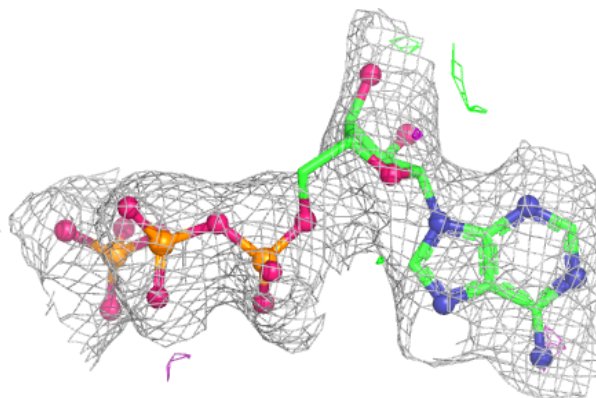


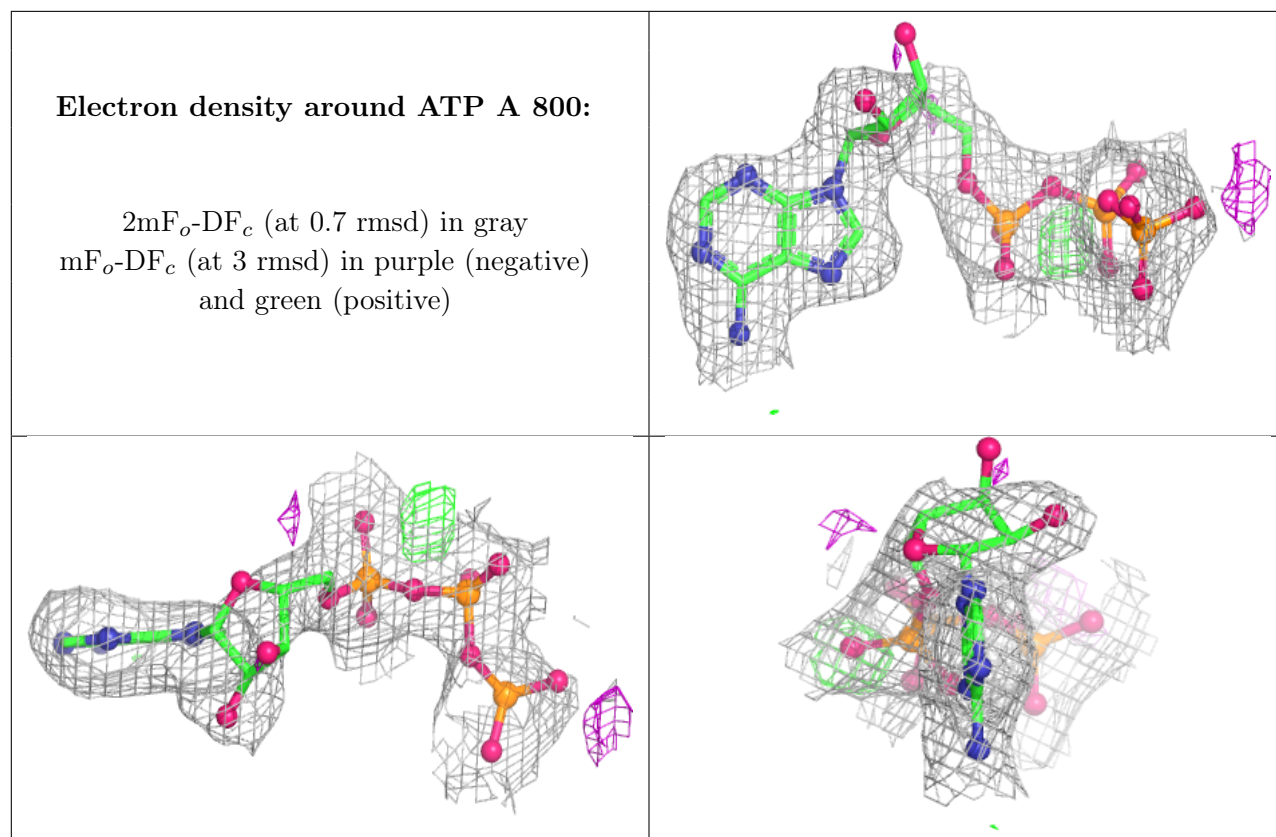
Electron density around ATP D 803:

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

**Electron density around ATP C 802:**

$2mF_o-DF_c$ (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.